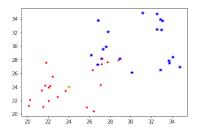
# Data Mining Lecture 9: Nearest Neighbours

Jo Houghton

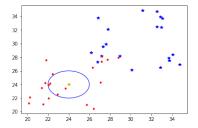
**ECS Southampton** 

March 11, 2019

How would you classify this point?

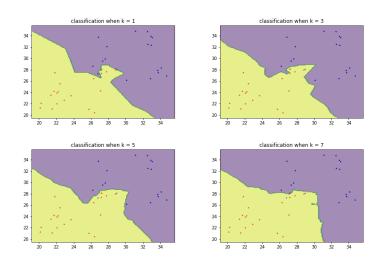


Use the closest samples..

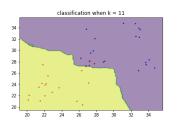


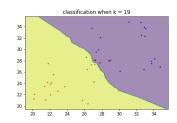
K-Nearest Neighbours: Assigns class based on majority class of closest K neighbours in featurespace

We can get a decision boundary given k: for example:

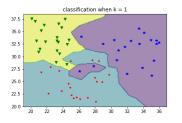


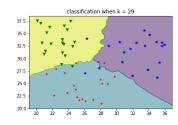
The boundary gets smoother, and generalises better when k is high for example:



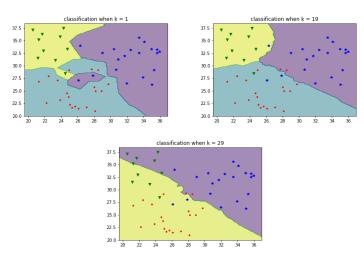


And with multi class classification, equally sized classes:





However, if k is too high, where some classes are less common, they can be missed



Advantages?

#### Advantages?

- ► No assumptions made
- ► No training phase
- ► Simple and easy to implement

#### Problems?

#### Advantages?

- ► No assumptions made
- ► No training phase
- Simple and easy to implement

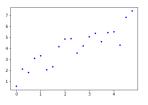
#### Problems?

- Doesn't scale well with lots of data
- Doesn't scale well with many dimensions

# Nearest Neighbours - Regression

KNN can be used to perform regression

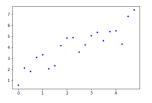
It uses the average value of the k closest data points



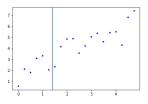
### Nearest Neighbours - Regression

KNN can be used to perform regression

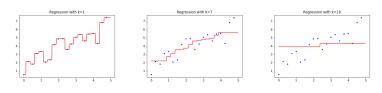
It uses the average value of the k closest data points



So a point at x = 1.4 will have a value  $\approx 2$  if k = 1 - 2

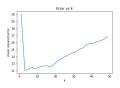


### Nearest Neighbours - Regression



From overfitting to underfitting..

The mean squared errors can be measured for each value of k



Greatest errors at the edges, interpolation easier than extrapolation Tuning k carefully is important - best done using cross validation ipynb Height Weight Age regression demo

Up to now, each value in the k nearest neighbours has been treated equally.

Better: if closer neighbours are more important

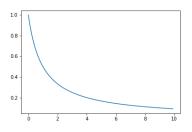
We can use a range of weighting schemes to do this:

- Inverse Weighting
- Subtraction weighting
- Gaussian Weighting

Inverse Weighting:

$$w = \frac{1}{dist + c}$$

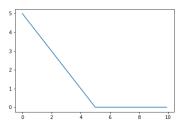
Where c is a constant, avoiding division by zero error if dist = 0



Subtraction Weighting:

$$w = \max(0, c - dist)$$

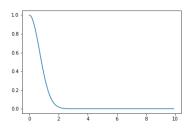
Where c is a constant



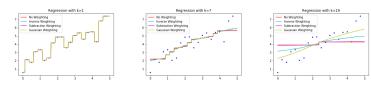
Gaussian Weighting:

$$w = \exp \frac{-dist^2}{c^2}$$

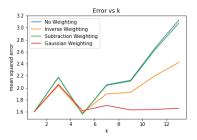
Where c is a constant



Weighted Regression:



Gaussian performs best here, especially with higher values of k



Still has greater errors at the edges of the data, interpolation easier than extrapolation

14 / 31

Again, to chose the best weighting scheme, measure performance using cross validation.

#### Problems?

- ► Heterogenous Data features with larger ranges have greater effects
- Outliers affect data a good deal, especially for low k
- ► For larger k, less common classes can get ignored
- Distance metric determines similarity usually Euclidean, works badly in high D
- Can use Hamming distance for categorical attributes
- Irrelevant data can force otherwise similar data samples to be far apart
- Computationally expensive if there are lots of data, or highly dimensional data

Curse of dimensionality:

For low dimensions, the number of points on the edge is very low E.g. for a line, the outer 1% of a line is 2% of the line (values at x>0.99, and x<0.1)

Curse of dimensionality:

For low dimensions, the number of points on the edge is very low E.g. for a line, the outer 1% of a line is 2% of the line (values at x>0.99, and x<0.1)

For a square, the outer 1% is  $1-0.98^2=0.0396\approx 4\%$ 

Curse of dimensionality:

For low dimensions, the number of points on the edge is very low E.g. for a line, the outer 1% of a line is 2% of the line (values at x>0.99, and x<0.1)

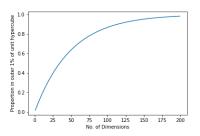
For a square, the outer 1% is  $1-0.98^2=0.0396\approx 4\%$ 

For a cube, the outer 1% is  $1 - 0.98^3 = 0.0588 \approx 6\%$ 

Curse of dimensionality:

For low dimensions, the number of points on the edge is very low E.g. for a line, the outer 1% of a line is 2% of the line (values at x > 0.99, and x < 0.1)

For a square, the outer 1% is  $1-0.98^2=0.0396\approx 4\%$ For a cube, the outer 1% is  $1-0.98^3=0.0588\approx 6\%$ 



This means in higher dimensions, data is nearly always extrapolated

Curse of dimensionality; For low dimensions, the size of a neighbourhood is small.

e.g. for k=10, number of points N=1,000,000In a unit line, the average neighbourhood is  $\frac{10}{10^6}=0.00001$  long

Curse of dimensionality; For low dimensions, the size of a neighbourhood is small.

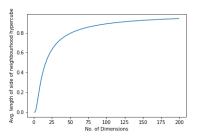
e.g. for k=10, number of points N=1,000,000 In a unit line, the average neighbourhood is  $\frac{10}{10^6}=0.00001$  long In a unit square, the average side length is  $\sqrt{\frac{10}{10^6}}=0.003$  long

Curse of dimensionality; For low dimensions, the size of a neighbourhood is small.

e.g. for k=10, number of points N=1,000,000 In a unit line, the average neighbourhood is  $\frac{10}{10^6}=0.00001$  long In a unit square, the average side length is  $\sqrt{\frac{10}{10^6}}=0.003$  long In a unit cube, the average side length is  $\sqrt[3]{\frac{10}{10^6}}=0.02$  long

Curse of dimensionality; For low dimensions, the size of a neighbourhood is small.

e.g. for k=10, number of points N=1,000,000 In a unit line, the average neighbourhood is  $\frac{10}{10^6}=0.00001$  long In a unit square, the average side length is  $\sqrt{\frac{10}{10^6}}=0.003$  long In a unit cube, the average side length is  $\sqrt[3]{\frac{10}{10^6}}=0.02$  long



This can make it very difficult to work out which are closer, as the distances are nearly all the same

Solutions: For heterogenous data?

Solutions: For heterogenous data?

► For heterogenous data, can *normalise* 

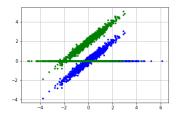
#### Solutions: For heterogenous data?

- For heterogenous data, can normalise
- Better to scale factors for each feature to optimise performance
- Could use this to do feature selection eg. if works best when scale factor = 0, then feature is useless!

Solutions for high D?

Solutions for high D?

- Dimensionality reduction.. Care! Some aren't suitable (e.g. MDS, SOM)
- ► Also.. PCA:



A random direction could be better!

Johnson Lindenstrauss lemma:

if points in a vector space are of high enough dimensionality,
they may be projected into a lower dimensional space in a way
which approximately preserves the distances between the
points, this basis can be generated randomly

More solutions for high D?

More solutions for high D?

► Use different metric:

#### More solutions for high D?

- Use different metric:
  - ► Hamming distance for categorical attributes
  - BM25 or TF-IDF for text data
  - Minkowski distance (p-norm) generalisation of Euclidean distance
  - Kullback Liebler Divergence for histograms

#### Solutions for lots of data?

► Need to quickly find the nearest neighbour to a particular point in a highly dimensional space

#### Solutions for lots of data?

- Need to quickly find the nearest neighbour to a particular point in a highly dimensional space
  - Could index points in a tree structure?
  - Could hash the points?
  - Could break up the space

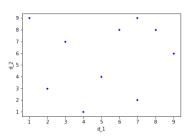
K-D trees are binary tree structures that partition the space along an axis-aligned hyperplane

- Chose random dimension
- Divide along median value
- Repeat until depth limit reached or certain number of items in each leaf

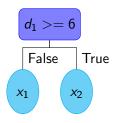
#### For a simple dataset:

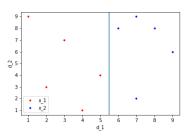
Tree:

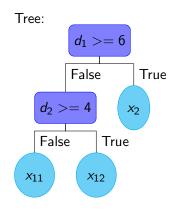


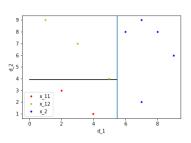


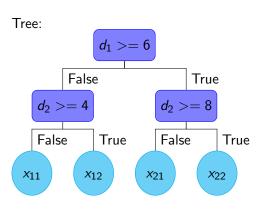
Tree:

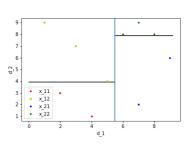






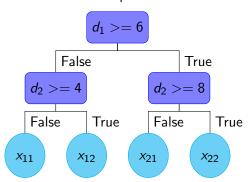


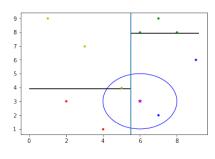




### To Classify: (6, 3)

- Go to correct part of tree and search subspace
- ► If the border is closer than k-neighbours in the subspace:
  - ► Go back up tree and search





#### Problems?

- Doesn't scale well to high dimensions
- Often need to search much of the tree
- Need many more examples than there are dimensions, at least  $2^n$
- There are approximate versions, not guaranteed exact answer but do scale
  - Based on ensembles of trees with a randomised split dimension

# Nearest Neighbours - LSH

Locality Sensitive Hashing

Makes hash codes that are similar for similar vectors

- ▶ Similar items map to the same buckets with high probability
- number of buckets much smaller than number of data samples
- Aims to maximise the probability of a collision for similar items

### Nearest Neighbours - LSH

#### Accomplished by:

- ▶ Chose random hyperplanes  $(h_1, h_2, ..., h_k)$
- ► Each hyperplane with split the space in to 2 regions
- $\triangleright$  : the space will be sliced in to  $2^k$  regions (buckets)
- Compare new point only to training points in the same region
- ▶ Repeat with different random hyperplanes  $(h_1, h_2, ..., h_k)$

Gives low complexity ,  $\approx O(d \log n)$ , as compare new data to only  $\frac{n}{2^k}$ 

# Nearest Neighbours - Summary

KNN can be used for regression as well as classification

- Using weighting can improve performance
- Poor performance with large data sets
- Can use K-D trees to help overcome these issues
  - ▶ Still can have issues with highly dimensional data
  - Often not much improvement in performance
- Curse of dimensionality
  - Affects neighbourhood size
  - Affects amount of extrapolation
- Can use dimensionality reduction to help (but be careful!)
- Fast approximate Nearest Neighbourhood methods LSH

Also: Final presentation does not need to be for full coursework, it is to show what you have done so far