# Data Mining Lecture 8: Decision Trees

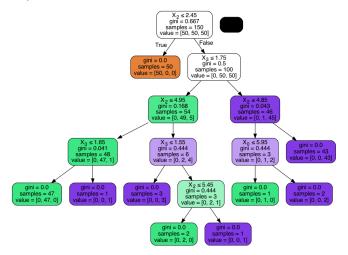
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9<sup>th</sup> March 2023

#### Decision Trees - Introduction

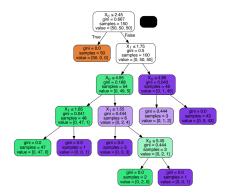
A decision tree is like a flow chart. For example, the iris dataset:



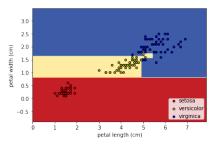
has four features, only three are used here, and one is only used once.

#### Decision Trees - Introduction

#### 2D iris dataset:



Decision surface of decision tree



# Decision Trees - Introduction

Decision Trees can be 'hand crafted' by experts They can also be built up using machine learning techniques

They are **interpretable**, it is easy to see how they made a certain decision

They are used in a wide range of contexts, for example:

- Medicine
- Financial Analysis
- Astronomy

Especially in medicine, the explicit reasoning in decision trees means experts can understand why the algorithm has made its decision.

Each node is a test, each branch is an outcome of the test Can be used for tabulated data with a range of data types, e.g. numerical, categorical ipynb demo

# Decision Trees - Tree Growing Algorithms

There are a good number of algorithms to build decision trees

- CART Classification And Regression Trees
- ► ID3 Iterative Dichotomiser 3
- C4.5 improved ID3
- C5.0 improved C4.5
- CHAID Chi squared Automatic Interaction Detector

The CART algorithm was published by Breiman et al in 1984

- Find best split for each feature minimises impurity measure
- Find feature that minimises impurity the most
- Use the best split on that feature to split the node
- Do the same for each of the leaf nodes

The CART algorithm depends on an impurity measure. It uses Gini impurity

Gini impurity measures how often a randomly chosen element from a set would be incorrectly labelled if it was randomly labelled according to the distribution of labels in the set. The probabilities for each label are summed up.

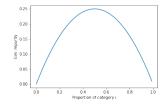
Gini Impurity  $(I_G)$  sums up probability of a mistake for each label:

mistake probability 
$$=\sum_{k
eq i}p_k=1-p_i$$

For J classes:

$$Gini(p) = \sum_{i=1}^{J} p_i \sum_{k \neq i} p_k = \sum_{i=1}^{J} p_i (1 - p_i)$$

It reaches its minimum when all cases in the node fall into a single category



Maximum improvement in impurity found using the equation:

$$Gini(root) - (Gini(Left)\frac{n_L}{n} + Gini(Right)\frac{n_R}{n})$$

Where Gini(root) is the impurity of the node to be split, Gini(Left) and Gini(Right) is the impurity of the left and right branches,  $n_L$  and  $n_R$  are the numbers in left and right branches.

Example:

Car Make	Туре	Colour	Price	Mileage	Bought?
VW	Polo	Grey	£2000	82000	Yes
Ford	Fiesta	Purple	$\pounds 1795$	95000	Yes
Ford	Fiesta	Grey	£1990	90000	No
VW	Golf	Red	£1800	120000	Yes
VW	Polo	Grey	£900	150000	No
Ford	Ka	Yellow	£1400	100000	Yes

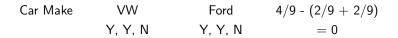
Can go through and calculate best split for each feature.

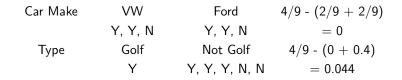
Calculate root node impurity:

$$\textit{Gini(root)} = \frac{1}{3}(1 - \frac{1}{3}) + \frac{2}{3}(1 - \frac{2}{3}) = \frac{2}{9} + \frac{2}{9} = \frac{4}{9}$$

Impurity decrease (or Information gain if using Entropy) is thus:

$$I_{G} = Gini(root) - \left(Gini(Left)\frac{n_{L}}{n} + Gini(Right)\frac{n_{R}}{n}\right)$$





Car Make	VW	Ford	4/9 - (2/9 + 2/9)
	Y, Y, N	Y, Y, N	= 0
Туре	Golf	Not Golf	4/9 - (0 + 0.4)
	Y	Y, Y, Y, N, N	= 0.044
Colour	Grey	Not Grey	4/9 - (2/9 + 0)
	Y, N, N	Y, Y, Y	= 0.222

Car Make	VW	Ford	4/9 - (2/9 + 2/9)
	Y, Y, N	Y, Y, N	= 0
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	Y, N, N	Y, Y, Y	= 0.222
Price	> 1000	< 1000	4/9 - (0.267+ 0)
	Y, N, Y, Y, Y	Ν	= 0.178

Car Make	VW	Ford	4/9 - (2/9 + 2/9)
	Y, Y, N	Y, Y, N	= 0
Туре	Golf	Not Golf	4/9 - (0 + 0.4)
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	Y, N, N	Y, Y, Y	= 0.222
Price	> 1000	< 1000	4/9 - (0.267+ 0)
	Y, N, Y, Y, Y	Ν	= 0.178

This gives the first split. The same process is repeated for each impure node until all nodes are pure.

The second split has the following items:

Car Make	Туре	Colour	Price	Mileage	Bought?
VW	Polo	Grey	£2000	82000	Yes
Ford	Fiesta	Grey	£1990	90000	No
VW	Polo	Grey	£900	150000	No

The second Split:

 $\begin{array}{cccc} \mbox{Car Make} & \mbox{VW} & \mbox{Ford} & \mbox{4/9 - } (1/3 + 0) \\ \mbox{Y, N} & \mbox{N} & = 0.111 \end{array}$ 

#### The second Split:

Car Make	VW	Ford	4/9 - (1/3 + 0)
	Υ, Ν	Ν	= 0.111
Туре	Polo	Not Polo	4/9 - (1/3 + 0)
	Y, N	Ν	= 0.111

#### The second Split:

Car Make	VW	Ford	4/9 - (1/3 + 0)	
	Υ, Ν	Ν	= 0.111	
Туре	Polo	Not Polo	4/9 - (1/3 + 0)	
	Υ, Ν	Ν	= 0.111	
Mileage	above 85,000	below 85,000	4/9 - (0 + 0)	
	N, N	Y	= 0.444	

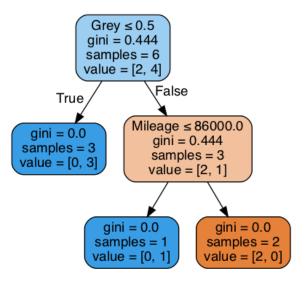
#### The second Split:

Car Make	VW	Ford	4/9 - (1/3 + 0)
	Y, N	Ν	= 0.111
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#### The second Split:

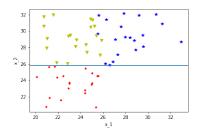
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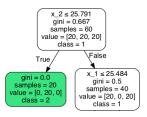
The best splits both remove all impurity so we are done:

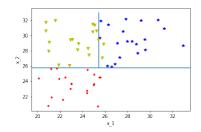


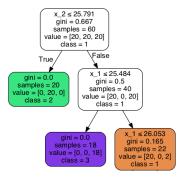
ipynb demo

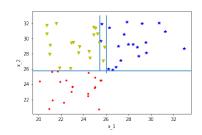
With three classes:

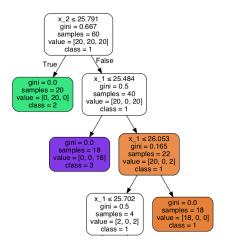


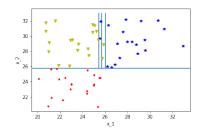












#### x 2 ≤ 25.791 aini = 0.667 samples = 60 value = [20, 20, 20] class = 1False True x 1 ≤ 25.484 gini = 0.0 aini = 0.5 samples = 20samples = 40 value = [0, 20, 0] value = [20, 0, 20] class = 2class = 1 x 1 ≤ 26.053 qini = 0.0gini = 0.165 samples = 18samples = 22value = [0, 0, 18] value = [20, 0, 2] class = 3class = 1x 1 ≤ 25.702 qini = 0.0gini = 0.5 samples = 18 samples = 4value = [18, 0, 0] value = [2, 0, 2] class = 1class = 1gini = 0.0 samples = 2 samples = 2value = [2, 0, 0] value = [0, 0, 2] class = 3class = 1

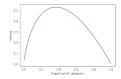
Overfitting ..

#### Decision Trees - ID3

Similar to CART, Iterative Dichotomy 3 (ID3) minimises entropy instead of Gini impurity Entropy:

$$H(S) = \sum_{x \in X} -p(x) \log_2 p(x)$$

Where S is the data set, X is the set of classes in S, p(x) is the proportion of the number of elements in class x to the number of elements in set S



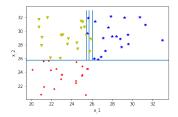
# Information Gain is measured for a split along each possible attribute $\boldsymbol{A}$

$$I_G(S,A) = H(S) - \sum_{x \in X} \frac{|S_A|}{|S|} H(S_A)$$

ID3 is very similar to CART, though doesn't technically support numerical values

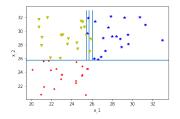
# Decision Trees - Pruning

Overfitting is a serious problem with Decision Trees



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Overfitting is a serious problem with Decision Trees



.. are four splits really required here?

The trees it creates are too complex.

One solution is pruning

This can be done in a variety of ways, including:

- Reduced Error Pruning
- Entropy Based Merging

# Decision Trees - Reduced Error Pruning

Growing a decision tree fully, then removing branches without reducing predictive accuracy, measured using a *validation set*.

- Start at leaf nodes
- Look up branches at last decision split
- replace with a leaf node predicting the majority class
- If validation set classification accuracy is not affected, then keep the change

This is a simple and fast algorithm that can simplify over complex decision trees

Grow a decision tree fully, then

- Chose a pair of leaf nodes with the same parent
- What is the entropy gain from merging them?
- If lower than a threshold, merge nodes

This doesn't require additional data

ID3 ignores missing data, CART generally puts them to the node that has the largest number of the same category

- > You can assign a branch specifically to an unknown value
- You can assign it to the branch with the most of the same target value
- You can weight each branch using the known factors and put it in the most similar branch

# Decision Trees - Regression

CART - Classification and Regression Trees How do we use decision trees for regression? i.e. to give numerical values rather than a classification.

Could use classification, but using each numerical value as a class..

Problems?

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Problems?

- How would you generalise?
- Loses all meaning of ordering, or similarity

Solution?

## Decision Trees - Regression

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Solution?

Use Variance instead of Gini or Entropy

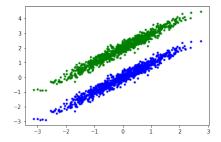
Maximise Variance Gain:

- Split on the feature values that give maximum gain in variance
- Should make similar numbers group together
- ▶ I. e. lower numbers on one side, higher on the other

#### Decision Trees - In General

There are problems with Decision Trees:

- Finding an optimal Tree is NP-complete
- They overfit, so don't generalise well Hence need to prune
- Information Gain is biased to features with more categories
- Splits are axis aligned..



Bagging: Bootstrap aggregating

Uniformly sample initial dataset with replacement in to m subsets For example:

if data set has 5 samples,  $(s_1, s_2, s_3, s_4, s_5)$  make a whole bunch of similar data sets:

 $\blacktriangleright$  ( $s_5$ ,  $s_2$ ,  $s_2$ ,  $s_1$ ,  $s_5$ )

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Train a different decision tree on each set

To classify, apply each classifier and chose the correct one by majority vote

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This improves generalisation, as decreases variance without increasing bias

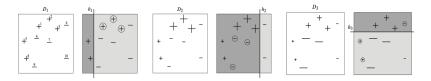
Boosting - Kearns and Valiant (1988): "Can a set of weak learners create a single strong learner?"

We make a weighted sum of very weak learners - so long as they all learn different things then it works! AdaBoost:

- Train a weak learner on one feature
- See what it does well on
- Weight the remaining data more
- repeat

This makes a series of weak learners that have learned how to use different features to discriminate between classes.

#### AdaBoost:



From Schapire and Freund, 2012) Shortcomings are identified by high weight data points

Gradient boosting with trees - Friedman (1999): Generalise Adaboost to Gradient boosting to handle any loss function

Shortcomings are where the residuals are larger

So fit a tree to the residuals:

• 
$$x_1, y_q - f(x_1)$$
  
•  $x_2, y_q - f(x_1)$   
•  $x_3, y_q - f(x_1)$   
•  $\vdots$ 

$$\blacktriangleright x_n, y_q - f(x_n)$$

more detail available at

http://www.chengli.io/tutorials/gradient\_boosting.pdf

Random Forests

Apply bagging but when learning the tree for each subset, chose the split by searching over a random sample of the features

Reduces overfitting

Decision Trees - Summary

Advantages:

- Interpretability
- Ability to work with numerical and categorical features
- Good with mixed tabular data

Disadvantages:

- Might not scale effectively for lots of classes
- Features that interact are problematic

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