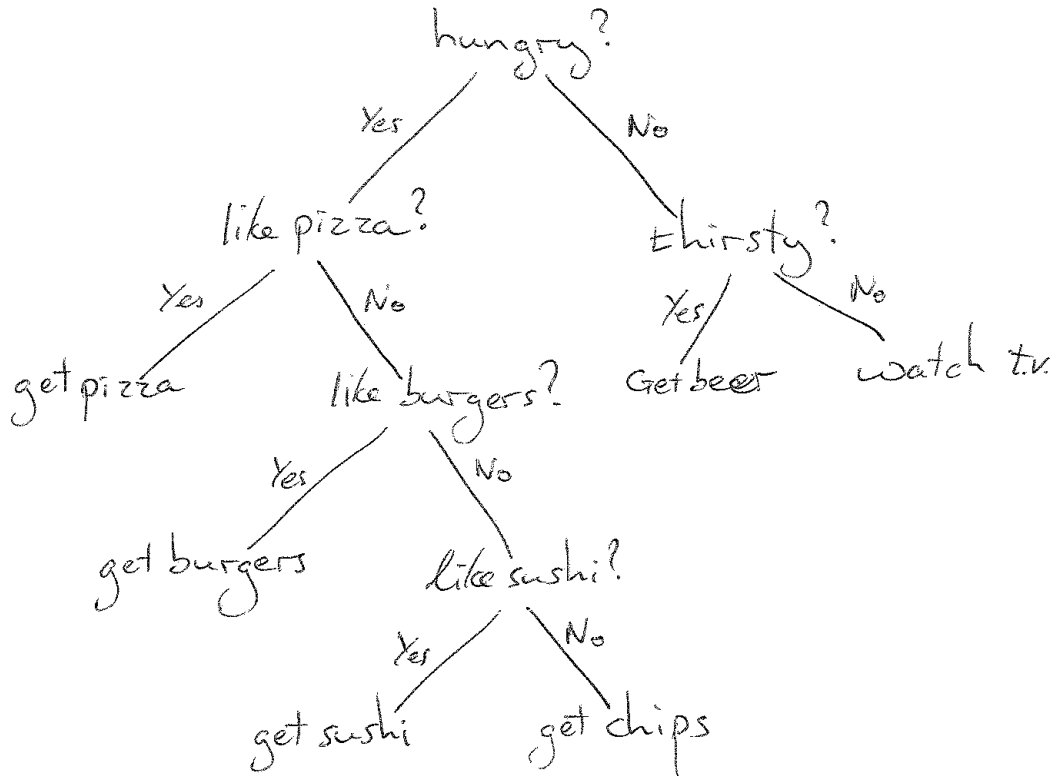


## Adaptive Computation and Machine Learning

## 3. DECISION TREES

A decision tree looks something like the following:



Every node in the tree is labelled by a question about an attribute from the given dataset. Start at the root node and consider the question at that node with respect to the given input. Follow the branch that corresponds to the answer. Continue to traverse the tree in this way until a leaf node is reached which gives us our decision.

We shall describe a way to construct a suitable decision tree for a dataset associated with a classification problem. That is, given a dataset in which every data point has an associated target classification, the objective is to construct a decision tree that predicts the correct class associated with every data point (or for an optimal number of data points).

The main issue when constructing a decision tree is choosing the attribute that should be used at a given node, that is, the ‘question’ to be asked at that node. Any attribute may be chosen at any node, however, we seek to minimise the number of questions required to obtain an outcome, which corresponds to finding a tree of minimum height.

For now, we assume that each attribute in the dataset uses discrete values (e.g., 0/1, T/F, Yes/No, A/B/C, or 0/1/2/3, etc.).

### 3.1. Entropy.

The method that we describe here for choosing an attribute is based on the idea of reducing the uncertainty in the dataset. The (Shannon) entropy of a probability distribution is used; we recall the definition here.

Let  $P$  denote a (discrete) probability distribution, i.e.,  $P = \{p_1, p_2, \dots, p_k\}$  with  $0 \leq p_i \leq 1$  for each  $i$  and  $\sum_{i=1}^k p_i = 1$ .

The **entropy** of  $P$  is:

$$H(P) = - \sum_{i=1}^k p_i \log_2 p_i$$

Examples:

If  $P = \{0.5, 0.5\}$ , then  $H(P) = -0.5 \log_2 0.5 - 0.5 \log_2 0.5 = 1$ .

If  $P = \{0.25, 0.75\}$ , then  $H(P) = -0.25 \log_2 0.25 - 0.75 \log_2 0.75 = 0.811$ .

If  $P = \{\frac{1}{3}, \frac{2}{3}\}$ , then  $H(P) = -\frac{1}{3} \log_2 \frac{1}{3} - \frac{2}{3} \log_2 \frac{2}{3} = 0.918$ .

If  $P = \{\frac{1}{3}, \frac{1}{3}, \frac{1}{3}\}$ , then  $H(P) = -\frac{1}{3} \log_2 \frac{1}{3} - \frac{1}{3} \log_2 \frac{1}{3} - \frac{1}{3} \log_2 \frac{1}{3} = 1.585$ .

If  $P = \{0.25, 0.25, 0.5\}$ , then  $H(P) = -0.25 \log_2 0.25 - 0.25 \log_2 0.25 - 0.5 \log_2 0.5 = 1.5$ .

Entropy is a measure of the randomness, or uncertainty, in the system.

For example, take the case  $P = \{p_1, p_2\}$  with  $p_1 + p_2 = 1$ . We can think of this as a probability distribution over two events, where  $p_1$  is the probability of event 1 occurring and  $p_2$  is the probability of event 2 occurring. If  $p_1 = 0.5$  and  $p_2 = 0.5$ , then the entropy (i.e., uncertainty) of the system is at a maximum since event 1 and event 2 are equally likely. If  $p_1 = 0.33$  and  $p_2 = 0.67$ , then the entropy is lower since we know that event 2 is more likely than event 1. If

$p_1 = 0.25$  and  $p_2 = 0.75$ , then the entropy is even lower since event 2 is even more likely than event 1 in this case. If  $p_1 = 0$  and  $p_2 = 1$  then the entropy is 0 since there is a 100% chance that event 2 will occur.

For cases with three events, i.e.,  $P = \{p_1, p_2, p_3\}$ , with  $\sum_{i=1}^3 p_i = 1$ , the maximum entropy occurs when all three events are equally likely, i.e.,  $p_1 = p_2 = p_3 = \frac{1}{3}$ . For any other combination of  $p_1, p_2$  and  $p_3$  the entropy will be less. Note that the entropy for  $\{\frac{1}{3}, \frac{1}{3}, \frac{1}{3}\}$  is greater than the entropy for  $\{\frac{1}{2}, \frac{1}{2}\}$ . Why do you think this is so?

### 3.2. Constructing a decision tree from a dataset.

Consider a dataset  $S$  consisting of an array  $X$  of data points and an array  $T$  of corresponding target values. Let  $F_1, F_2, \dots, F_m$  be the **attributes** of the data points in  $X$ . Attributes are sometimes referred to as **features**.

We seek to construct a decision tree for the given dataset. At each node of the tree a question is asked about one attribute of the data. We must decide which attribute to use at each node. The first step is to choose an attribute for the **root node**, which induces the next layer of nodes and branches. The same method is repeated at every new node until the decision tree is complete.

For each attribute  $F$  define the  $\text{Gain}(S, F)$  as follows:

$$\text{Gain}(S, F) = \text{Entropy}(S) - \frac{1}{|S|} \sum_{f \in \text{values of } F} |S_f| \text{Entropy}(S_f),$$

where  $\text{Entropy}(S)$  denotes the entropy of the probability distribution on the target values of  $S$ , and  $S_f$  is the subset of  $S$  containing of all data points for which attribute  $F$  has value  $f$ .

To determine the attribute to choose at the root node, we calculate  $\text{Gain}(S, F)$  for each attribute  $F$  and choose the attribute for which  $\text{Gain}(S, F)$  is maximum.

Observe that for each attribute  $F$ , the terms  $\text{Entropy}(S)$  and  $\frac{1}{|S|}$  in the formula for  $\text{Gain}(S, F)$  are unaffected by the choice of  $F$ , so it is only necessary to calculate

$$\sum_{f \in \text{values of } F} |S_f| \text{Entropy}(S_f)$$

for each attribute  $F$ . In fact, the attribute that gives the *minimum* value for the above sum will give the maximum Gain.

**Example:** Consider the following dataset on students doing COMS3, where we have combined the data points and targets into one table, for convenience:

$$S : \left[ \begin{array}{c|c|c|c} \text{COMS2} & \text{doing labs?} & \text{doing tuts?} & \text{target} \\ \hline A & N & Y & \text{Pass} \\ C & Y & N & \text{Fail} \\ C & N & Y & \text{Pass} \\ B & Y & Y & \text{Pass} \\ B & N & N & \text{Fail} \\ C & Y & N & \text{Pass} \\ A & N & N & \text{Fail} \\ B & Y & N & \text{Pass} \end{array} \right]$$

In the above dataset, the attributes of the data are: COMS2, doing labs?, and doing tuts?.

Recall that  $\text{Entropy}(S)$  is the entropy of the probability distribution on the targets of the dataset. There are only 2 possible targets: *Pass* and *Fail*. Of the 8 entries, there are 5 *Pass*'s and 3 *Fail*'s. Thus, the probability distribution on the targets is  $P = \{p_1, p_2\}$ , where

$$p_1 = \text{probability of } \text{Pass} = \frac{5}{8},$$

$$p_2 = \text{probability of } \text{Fail} = \frac{3}{8}.$$

$$\text{Thus, } \text{Entropy}(S) = H(P) = -\frac{5}{8} \log_2 \frac{5}{8} - \frac{3}{8} \log_2 \frac{3}{8} = 0.954.$$

To decide which attribute to use at the root node, we calculate  $\text{Gain}(S, \text{COMS2})$ ,  $\text{Gain}(S, \text{doing labs?})$  and  $\text{Gain}(S, \text{doing tuts?})$ .

- $\text{Gain}(S, \text{COMS2})$ : attribute COMS2 has 3 values: *A*, *B* and *C*. Thus:

$$\sum_{f \in \text{values of COMS2}} |S_f| \text{Entropy}(S_f) = |S_A| \text{Entropy}(S_A) + |S_B| \text{Entropy}(S_B) + |S_C| \text{Entropy}(S_C).$$

$\text{Entropy}(S_A)$  is calculated on the subset of  $S$  of all data points for which COMS2 is *A*:

$$S_A = \left[ \begin{array}{c|c|c|c} \text{COMS2} & \text{doing labs?} & \text{doing tuts?} & \text{target} \\ \hline A & N & Y & \text{Pass} \\ A & N & N & \text{Fail} \end{array} \right]$$

Note that  $|S_A| = 2$ .  $\text{Entropy}(S_A)$  is calculated on the target column.

The probability distribution on the targets is  $P = \{p_1, p_2\}$ , where

$$p_1 = \text{probability of } \text{Pass} = \frac{1}{2},$$

$p_2 = \text{probability of } Fail = \frac{1}{2}$ .

Thus,  $\text{Entropy}(S_A) = H(P) = -\frac{1}{2} \log_2 \frac{1}{2} - \frac{1}{2} \log_2 \frac{1}{2} = 1$ .

$\text{Entropy}(S_B)$  is calculated on the subset of  $S$  of all data points for which COMS2 is  $B$ :

$$S_B = \left[ \begin{array}{c|c|c|c} \text{COMS2} & \text{doing labs?} & \text{doing tuts?} & \text{target} \\ \hline B & Y & Y & Pass \\ B & N & N & Fail \\ B & Y & N & Pass \end{array} \right]$$

Note that  $|S_B| = 3$ . The probability distribution on the targets is  $P = \{p_1, p_2\}$ , where

$p_1 = \text{probability of } Pass = \frac{2}{3}$ ,

$p_2 = \text{probability of } Fail = \frac{1}{3}$ .

Thus,  $\text{Entropy}(S_B) = H(P) = -\frac{2}{3} \log_2 \frac{2}{3} - \frac{1}{3} \log_2 \frac{1}{3} = 0.918$ .

$\text{Entropy}(S_C)$  is calculated on the subset of  $S$  of all data points for which COMS2 is  $C$ :

$$S_C = \left[ \begin{array}{c|c|c|c} \text{COMS2} & \text{doing labs?} & \text{doing tuts?} & \text{target} \\ \hline C & Y & N & Fail \\ C & N & Y & Pass \\ C & Y & N & Pass \end{array} \right]$$

Note that  $|S_C| = 3$ . The probability distribution on the targets is  $P = \{p_1, p_2\}$ , where

$p_1 = \text{probability of } Pass = \frac{2}{3}$ ,

$p_2 = \text{probability of } Fail = \frac{1}{3}$ .

Thus,  $\text{Entropy}(S_C) = H(P) = -\frac{2}{3} \log_2 \frac{2}{3} - \frac{1}{3} \log_2 \frac{1}{3} = 0.918$ .

From the above calculations,

$$|S_A| \text{Entropy}(S_A) + |S_B| \text{Entropy}(S_B) + |S_C| \text{Entropy}(S_C) = 2(1) + 3(0.918) + 3(0.918) = 7.508$$

$$\text{hence Gain}(S, \text{COMS2}) = 0.954 - \frac{1}{8}(7.508) = 0.016.$$

- $\text{Gain}(S, \text{doing labs?})$ : attribute ‘doing labs?’ has 2 values:  $Y$  and  $N$ . Thus:

$$\sum_{f \in \text{values of 'doing labs?'}} |S_f| \text{Entropy}(S_f) = |S_Y| \text{Entropy}(S_Y) + |S_N| \text{Entropy}(S_N).$$

Entropy( $S_Y$ ) is calculated on the subset of  $S$  of all data points for which ‘doing labs?’ is  $Y$ :

$$S_Y = \left[ \begin{array}{c|c|c|c} \text{COMS2} & \text{doing labs?} & \text{doing tuts?} & \text{target} \\ \hline C & Y & N & \textit{Fail} \\ B & Y & Y & \textit{Pass} \\ C & Y & N & \textit{Pass} \\ B & Y & N & \textit{Pass} \end{array} \right]$$

Note that  $|S_Y| = 4$ . The probability distribution on the targets is  $P = \{p_1, p_2\}$ , where

$p_1 = \text{probability of } \textit{Pass} = \frac{3}{4}$ ,

$p_2 = \text{probability of } \textit{Fail} = \frac{1}{4}$ .

Thus,  $\text{Entropy}(S_Y) = H(P) = -\frac{3}{4} \log_2 \frac{3}{4} - \frac{1}{4} \log_2 \frac{1}{4} = 0.811$ .

Entropy( $S_N$ ) is calculated on the subset of  $S$  of all data points for which ‘doing labs?’ is  $N$ :

$$S_N = \left[ \begin{array}{c|c|c|c} \text{COMS2} & \text{doing labs?} & \text{doing tuts?} & \text{target} \\ \hline A & N & Y & \textit{Pass} \\ C & N & Y & \textit{Pass} \\ B & N & N & \textit{Fail} \\ A & N & N & \textit{Fail} \end{array} \right]$$

Note that  $|S_N| = 4$ . The probability distribution on the targets is  $P = \{p_1, p_2\}$ , where

$p_1 = \text{probability of } \textit{Pass} = \frac{1}{2}$ ,

$p_2 = \text{probability of } \textit{Fail} = \frac{1}{2}$ .

Thus,  $\text{Entropy}(S_N) = H(P) = -\frac{1}{2} \log_2 \frac{1}{2} - \frac{1}{2} \log_2 \frac{1}{2} = 1$ .

From the above calculations,

$$|S_Y| \text{Entropy}(S_Y) + |S_N| \text{Entropy}(S_N) = 4(0.811) + 4(1) = 7.244$$

hence  $\text{Gain}(S, \text{doing labs?}) = 0.954 - \frac{1}{8}(7.244) = 0.0485$ .

•  $\text{Gain}(S, \text{doing tuts?})$ : attribute ‘doing tuts?’ has 2 values:  $Y$  and  $N$ . Thus:

$$\sum_{f \in \text{values of 'doing tuts?'}} |S_f| \text{Entropy}(S_f) = |S_Y| \text{Entropy}(S_Y) + |S_N| \text{Entropy}(S_N).$$

Entropy( $S_Y$ ) is calculated on the subset of  $S$  of all data points for which ‘doing tuts?’ is  $Y$ :

$$S_Y = \left[ \begin{array}{c|c|c|c} \text{COMS2} & \text{doing labs?} & \text{doing tuts?} & \text{target} \\ \hline A & N & Y & Pass \\ C & N & Y & Pass \\ B & Y & Y & Pass \end{array} \right]$$

Note that  $|S_Y| = 3$ . The probability distribution on the targets is  $P = \{p_1, p_2\}$ , where

$p_1$  = probability of  $Pass = 1$ ,

$p_2$  = probability of  $Fail = 0$ .

Thus, Entropy( $S_Y$ ) =  $H(P) = -\log_2 1 - 0 \log_2 0 = 0$ . (Note that we always take  $0 \log_2 0 = 0$ .)

Entropy( $S_N$ ) is calculated on the subset of  $S$  of all data points for which ‘doing labs?’ is  $N$ :

$$S_N = \left[ \begin{array}{c|c|c|c} \text{COMS2} & \text{doing labs?} & \text{doing tuts?} & \text{target} \\ \hline C & Y & N & Fail \\ B & N & N & Fail \\ C & Y & N & Pass \\ A & N & N & Fail \\ B & Y & N & Pass \end{array} \right]$$

Note that  $|S_N| = 5$ . The probability distribution on the targets is  $P = \{p_1, p_2\}$ , where

$p_1$  = probability of  $Pass = \frac{2}{5}$ ,

$p_2$  = probability of  $Fail = \frac{3}{5}$ .

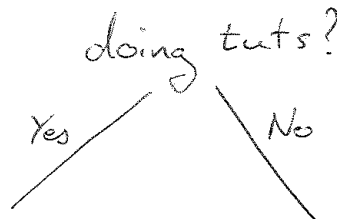
Thus, Entropy( $S_N$ ) =  $H(P) = -\frac{2}{5} \log_2 \frac{2}{5} - \frac{3}{5} \log_2 \frac{3}{5} = 0.971$ .

From the above calculations,

$$|S_Y| \text{Entropy}(S_Y) + |S_N| \text{Entropy}(S_N) = 3(0) + 5(0.971) = 4.855$$

hence  $\text{Gain}(S, \text{doing tuts?}) = 0.954 - \frac{1}{8}(4.855) = 0.347$ .

Thus, the maximum gain is obtained from using the attribute ‘doing tuts?’ at the root node, so the decision tree starts as follows:



Next, the process is repeated on the  $Y$  branch of the tree and also on the  $N$  branch of the tree. On the  $Y$  branch, the dataset is restricted to the data points in  $S$  for which ‘doing tuts?’ has value  $Y$ , i.e.,

COMS2	doing labs?	doing tuts?	target
$A$	$N$	$Y$	$Pass$
$C$	$N$	$Y$	$Pass$
$B$	$Y$	$Y$	$Pass$

On the above dataset, the target is always  $Pass$  hence we can make this node a leaf node labeled with the decision  $Pass$ .

On the  $N$  branch, the dataset is restricted to the data points for which ‘doing tuts?’ has value  $N$ , i.e.,

COMS2	doing labs?	doing tuts?	target
$C$	$Y$	$N$	$Fail$
$B$	$N$	$N$	$Fail$
$C$	$Y$	$N$	$Pass$
$A$	$N$	$N$	$Fail$
$B$	$Y$	$N$	$Pass$

On the above dataset, there are both  $Pass$  and  $Fail$  targets, so another attribute is required at this node. That is, a question can be asked about either: COMS2 or doing labs?. Both  $\text{Gain}(S, \text{COMS2})$  and  $\text{Gain}(S, \text{doing labs?})$  are calculated for this dataset and whichever attribute gives maximum Gain is used as the attribute for this node. This process continues until all branches of the tree end in a leaf node, in which case the tree is complete.

A leaf node is created in two ways. Firstly, if all the data points in the current dataset have the same target, then a leaf node is created and labeled with that target. Secondly, if the current dataset has no more attributes left to test, then a leaf node is created. In this case, it may happen that the targets are not all the same. Then the label chosen for the leaf node is the most common target in the current dataset. (It may happen that there is a tie for the most common, and then the label must be chosen some other way, or the label can indicate a split decision. Another possibility is that the dataset is empty at this node, which can also be indicated by some special label.)



### 3.3. ID3 Algorithm.

The method for constructing a decision tree from a dataset using the entropy measure described above is called the ID3 algorithm. An outline of the algorithm is given below.

#### ID3 ALGORITHM

Given a dataset with attributes  $F_1, \dots, F_m$ , which are all discrete (i.e., each attribute has a finite set of possible values) and discrete classification values for targets, a decision tree is built as follows:

Initially, **assign**  $S$  to be the whole dataset.

- if** all data points in  $S$  have the same target value, create a leaf node and label it with that target value,
- else if** there are no attributes left to test, create a leaf node and label it with the target value that is the most common in  $S$ ,
- else** find the feature  $\hat{F}$  that maximises the information gain
  - for** each value  $f$  of  $\hat{F}$ , add a new branch and node and calculate  $S_f$ 
    - by selecting only those data points with  $\hat{F}$ -value equal to  $f$
    - remove attribute  $\hat{F}$ ,
    - recursively call the algorithm on dataset  $S_f$ .

## EXERCISES

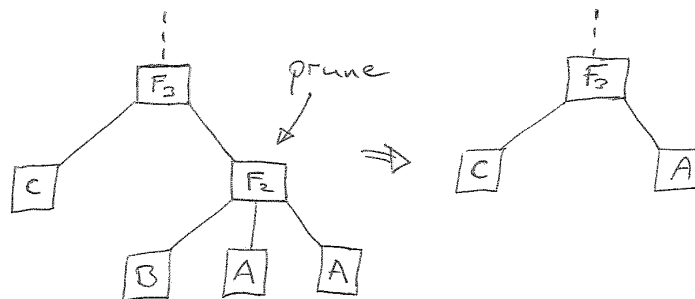
- (1) Complete the decision tree in the above example.
- (2) Construct the decision tree for the following dataset using the entropy method described above:

$S :$

$F_1$	$F_2$	$F_3$	target
$T$	0	$B$	$Yes$
$F$	2	$A$	$No$
$F$	1	$C$	$Yes$
$T$	2	$B$	$Yes$
$F$	0	$A$	$No$
$F$	2	$C$	$No$
$F$	0	$A$	$Yes$
$T$	1	$B$	$No$
$T$	0	$B$	$Yes$
$F$	1	$C$	$Yes$
$T$	2	$A$	$No$
$T$	2	$C$	$No$

### 3.4. Pruning a decision tree.

It is very likely that overfitting will occur when constructing a decision tree using the ID3 algorithm. To avoid overfitting, the method of **pruning** a tree can be used. This means replacing any subtree with a leaf node and giving the node a label that is the most common target of all the datapoints that would be decided in this subtree. That is, consider every data point in the dataset whose prediction would be obtained by following a path that ends up in a leaf within this subtree. Of those data points choose the most common target.



To decide where to prune a tree, a validation set is used (as in the case of neural networks). The initial dataset is split into three subsets: training, validation and test datasets. (The usual splitting rules apply: approximately: 50%, 25%, 25% or 60%, 20%, 20%).

The initial tree, say  $T$ , is built using the ID3 ALGORITHM on the training dataset. Using tree  $T$ , calculate the **error<sub>V</sub>** on the validation set as follows: Each data point from the validation dataset is input into the existing tree and a class prediction is obtained. If the prediction differs from the target, this is a misclassification. The validation set error on  $T$  is then:

$$\text{error}_V(T) = \frac{\text{number of misclassifications}}{\text{number of data points}}$$

Choose any subtree that you are considering pruning and let  $T'$  be the pruned tree. As above, calculate the validation set error on the pruned tree,  $\text{error}_V(T')$ . If  $\text{error}_V(T') < \text{error}_V(T)$  then replace the tree  $T$  by the pruned tree  $T'$ ; otherwise keep the original tree  $T$ .

The above process can be repeated until there are no ways left to prune the tree that reduce the validation set error (or until the height of the tree is at a specified level). After pruning is complete, the test dataset is used to evaluate the tree.

### 3.5. Gini Impurity.

An alternative to using Entropy for constructing a decision tree for a given dataset is the Gini Impurity. For a probability distribution  $P = \{p_1, p_2, \dots, p_k\}$ , the **Gini Impurity** of  $P$  is defined as:

$$\text{Gini}(P) = 1 - \sum_{i=1}^k p_i^2.$$

For example, if  $P = \{\frac{1}{4}, \frac{3}{4}\}$ , then  $\text{Gini}(P) = 1 - ((\frac{1}{4})^2 + (\frac{3}{4})^2) = 0.375$ .

The minimum Gini Impurity occurs if the distribution is of the form  $P = \{0, \dots, 0, 1, 0, \dots, 0\}$ , in which case  $\text{Gini}(P) = 0$ . This occurs if all data points in the set have the same target. Such a set is called **pure**. As the target values become more mixed, the impurity increases.

The maximum Gini Impurity occurs if the distribution is of the form  $P = \{\frac{1}{k}, \frac{1}{k}, \dots, \frac{1}{k}\}$ , in which case  $\text{Gini}(P) = 1 - \sum_{i=1}^k (\frac{1}{k})^2 = 1 - \frac{k}{k^2} = \frac{k-1}{k}$ . In this case, as  $k$  increases, the Gini Impurity tends to 1.

For a dataset  $S$ , the Gini Impurity of  $S$ , denoted  $\text{Gini}(S)$ , is defined on its set of targets (as in the Entropy case) by finding the probabilities  $p_1, \dots, p_k$  for each of the possible target classes and obtaining the Gini Impurity for this distribution. For example, consider a dataset that looks like:

$$S : \left[ \begin{array}{c|c|c|c|c} F_1 & F_2 & \dots & F_m & \text{target} \\ \hline \vdots & \vdots & \vdots & \vdots & A \\ \vdots & \vdots & \vdots & \vdots & B \\ \vdots & \vdots & \vdots & \vdots & A \\ \vdots & \vdots & \vdots & \vdots & C \\ \vdots & \vdots & \vdots & \vdots & B \\ \vdots & \vdots & \vdots & \vdots & A \\ \vdots & \vdots & \vdots & \vdots & C \end{array} \right]$$

The probability distribution on the targets of  $S$  is  $P = \{p_1, p_2, p_3\}$ , where  $p_1 = P(A) = \frac{3}{7}$ ,  $p_2 = P(B) = \frac{2}{7}$ ,  $p_3 = P(C) = \frac{2}{7}$ . Then  $\text{Gini}(S) = 1 - ((\frac{3}{7})^2 + (\frac{2}{7})^2 + (\frac{2}{7})^2) = 0.653$ .

Constructing a decision tree using Gini Impurity is very similar to using Entropy. To choose an attribute for a given node, calculate the following Gain for each attribute  $F$ :

$$\text{Gain}(S, F) = \text{Gini}(S) - \frac{1}{|S|} \sum_{f \in \text{values of } F} |S_f| \text{Gini}(S_f).$$

Then, as in the ID3 algorithm, a branch is created for each value  $f$  of  $F$  and the algorithm recurses on each of the sets  $S_f$  until leaf nodes are created.

### 3.6. Dealing with continuous variables.

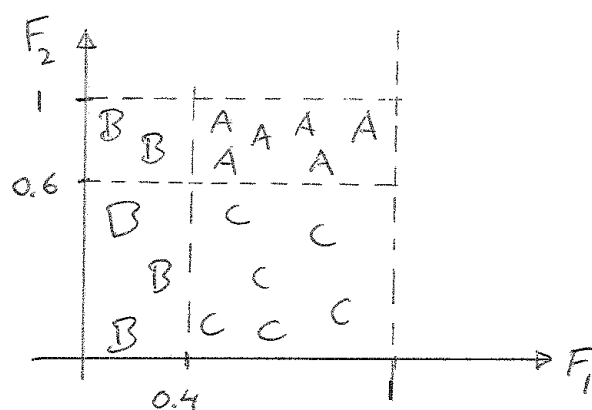
Suppose that a dataset has attributes that are not discrete, but continuous, i.e., they can take any real value in some range, as in the following example:

$$S : \left[ \begin{array}{c|c|c} F_1 & F_2 & \text{target} \\ \hline 0.75 & 0.92 & A \\ 0.51 & 0.43 & B \\ 0.89 & 0.14 & C \\ 0.72 & 0.25 & A \\ \vdots & \vdots & \vdots \end{array} \right]$$

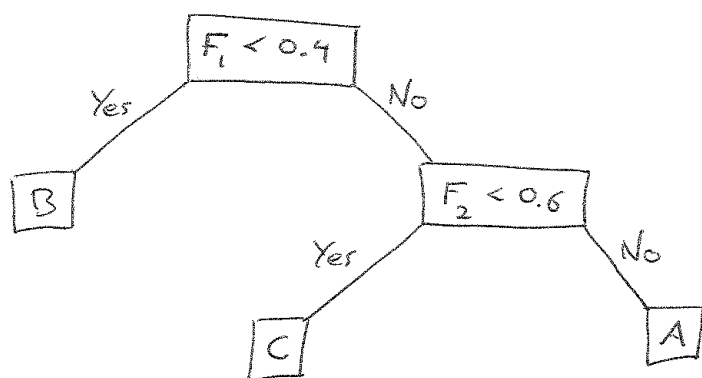
To build a decision tree for this dataset, the types of questions that are asked about attributes are of the form:  $F_1 < 0.6$ , or  $F_2 < 0.3$ , for example. The values 0.6 and 0.3 must be chosen somehow - they are called **split-points**.

There are different ways of choosing a split-point. One way is to consider a number of possible split-points at regular intervals, say at 0.1, 0.2, 0.3, 0.4, 0.5, 0.6, 0.7, 0.8, 0.9. For each of these possible split-points, calculate the Gain (using the Entropy method) for that feature for each split point. Do this for each continuous attribute and also calculate the Gain for any discrete-valued attributes. Then choose the attribute/split point that gives the maximum Gain.

For example, consider a dataset that has two continuous attributes  $F_1$  and  $F_2$  in the range  $[0, 1]$  and targets in the set  $\{A, B, C\}$  that is illustrated in the following diagram:



A good decision tree for the above data set would be:



When dealing with continuous-valued attributes, an attribute may be used more than once as a node question. For example, we may ask ' $F_1 < 0.4$ ' at the root and then later on ' $F_1 < 0.2$ '. (In the discrete case, asking about the same attribute a second time is not useful since the data has been split into subsets where that attribute has same value.)

For the above reason, and also because of the need to search for good split points, it is often better to use random forests for datasets whose attributes are mainly continuous-valued (see below).

### 3.7. Regression trees.

A decision tree can also be constructed for a regression problem, that is, a dataset in which the targets are real values (as opposed to a classification problem where the targets are discrete values). An example of such a dataset is the following:

$$S = \begin{bmatrix} F_1 & F_2 & \dots & F_m & \text{target} \\ \hline \vdots & \vdots & \vdots & \vdots & 2.7 \\ \vdots & \vdots & \vdots & \vdots & 2.35 \\ \vdots & \vdots & \vdots & \vdots & 1.98 \\ \vdots & \vdots & \vdots & \vdots & -0.33 \\ \vdots & \vdots & \vdots & \vdots & -1.05 \\ \vdots & \vdots & \vdots & \vdots & 0.77 \\ \vdots & \vdots & \vdots & \vdots & 0.5 \end{bmatrix}$$

In the case of a regression problem, the Entropy and Gini Impurity measures are not appropriate. Instead, the mean square error is used as a measure.

If the column of target values for dataset  $S$  is  $\begin{bmatrix} t_1 \\ \vdots \\ t_N \end{bmatrix}$  then the **mean square error** of  $S$ , denoted  $MSE(S)$ , is obtained by first calculating the mean of the targets:

$$\text{mean of } S = \mu_S = \frac{1}{N} \sum_{i=1}^N t_i$$

and then computing

$$MSE(S) = \frac{1}{N} \sum_{i=1}^N (t_i - \mu_S)^2.$$

For example, for the above data set, first calculate the mean:

$$\mu_S = \frac{1}{7}(2.7 + 2.35 + 1.98 - 0.33 - 1.05 + 0.77 + 0.5) = 0.989.$$

Then

$$\begin{aligned}
 MSE(S) &= \frac{1}{7} \left( (2.7 - \mu_S)^2 + (2.35 - \mu_S)^2 + (1.98 - \mu_S)^2 + (-0.33 - \mu_S)^2 \right. \\
 &\quad \left. + (-1.05 - \mu_S)^2 + (0.77 - \mu_S)^2 + (0.5 - \mu_S)^2 \right) \\
 &= 1.707.
 \end{aligned}$$

If all the target values in the dataset are close together, then  $MSE$  is small, and the more spread out the target values are, the greater  $MSE$  becomes. If the values are all the same then  $MSE$  is 0.

The algorithm for building the decision tree is the same as in the earlier cases, except that  $MSE$  is used instead of Entropy or Gini Impurity. For each attribute  $F$ , calculate the Gain as follows:

$$Gain(S, F) = MSE(S) - \frac{1}{|S|} \sum_{f \in \text{values of } F} |S_f| MSE(S_f).$$

Using the observation that  $MSE(S_f) = \frac{1}{|S_f|} \sum_{i \in S_f} (t_i - \mu_{S_f})^2$ , the above simplifies to:

$$Gain(S, F) = MSE(S) - \frac{1}{|S|} \sum_{f \in \text{values of } F} \sum_{i \in S_f} (t_i - \mu_{S_f})^2.$$

Next, select the attribute  $F$  for which the Gain is maximum.

Then, as in the ID3 algorithm, a branch is created for each value  $f$  of  $F$  and the algorithm recurses on each of the sets  $S_f$  until leaf nodes are created. When a leaf node is created, the label it is given is the average of all the values in the current dataset.

The above algorithms for using trees for classification and regression are often referred to as CART algorithms (Classification and Regression Trees).

### 3.8. Random Forests.

If one tree is not good enough, sometimes it is useful to create a number of trees for a given dataset. This is especially the case when dealing with continuous-valued attributes.

The trees are created randomly: at each node an attribute is chosen at random. If the attribute is continuous-valued, then the split-point can also be randomly selected. Usually the height of the tree is restricted so as to prevent over-training. The leaf nodes are labelled by the most

common target in the dataset at that node, or by the mean of the targets in the regression case. A collection of randomly created trees is called a **random forest**.

Once a random forest has been created, it can be used as a classifier. Given some data point, feed it into each tree in the forest and get a prediction from that tree. In the case of classification, choose the class that occurs most often in the set of predictions. In the case of regression, take the average value of the set of predictions.

## EXERCISES

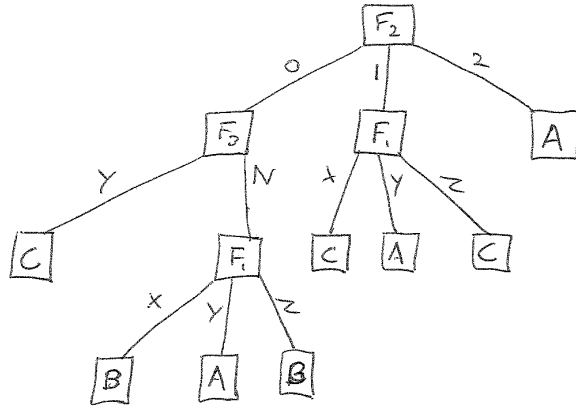
- (1) Construct the decision tree for the following dataset using
  - (a) the Entropy method and
  - (b) the Gini Impurity method.

$$S : \left[ \begin{array}{c|c|c|c} F_1 & F_2 & F_3 & \text{target} \\ \hline a & 0 & Y & A \\ b & 0 & N & C \\ c & 1 & Y & B \\ b & 1 & Y & B \\ c & 0 & N & A \\ a & 0 & N & C \\ c & 1 & N & B \\ a & 1 & Y & A \\ b & 0 & Y & B \\ a & 1 & N & C \end{array} \right]$$

- (c) Find the predictions from the trees obtained in (a) and (b) for the following inputs:
  - (i)  $(b, 1, N)$
  - (ii)  $(c, 1, Y)$ .



- (2) Suppose that the decision tree below has been constructed using some training dataset.



Let the following dataset be the validation set.

Validation set :

$F_1$	$F_2$	$F_3$	target
X	2	Y	A
Y	0	N	B
Z	1	Y	C
Z	1	N	B
Y	2	N	A
Z	0	Y	C
X	2	N	A
Y	1	Y	A
Z	2	Y	C
X	0	N	C
Y	0	Y	B
X	0	N	B

- (a) Compute the validation set error for the above decision tree.
- (b) For each (non-leaf) node in the tree, prune the node and see if the validation set error decreases on the pruned tree.
- (3) Suppose a decision tree has been constructed from some dataset. Then it can be used to make predictions for new data points that are not part of the original dataset. Suppose a data point has one of its attribute values missing. Describe how a prediction could still be made in this case. For example, for the tree in Question 2, suppose the input is:  $[Z, ?, Y]$ . What prediction could be obtained from the tree?

- (4) A regression tree is to be constructed for the dataset below using  $MSE$  on the targets.

The attributes  $F_1$  and  $F_2$  can take any real value between 0 and 1.

Determine what question should be asked at the root node of the tree.

When choosing a split-point value for attributes  $F_1$  and  $F_2$ , just consider the values 0.25, 0.5 and 0.75.

$$S = \begin{array}{c|c|c|c} F_1 & F_2 & F_3 & \text{target} \\ \hline 0.8 & 0.2 & a & 2.7 \\ 0.7 & 0.1 & b & 2.3 \\ 0.2 & 0.9 & c & 1.7 \\ 0.3 & 0.9 & b & 2.1 \\ 0.3 & 0.6 & a & 2.0 \\ 0.1 & 0.4 & c & 1.3 \\ 0.4 & 0.8 & b & 2.9 \\ 0.1 & 0.3 & c & 1.4 \end{array}$$