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Based on the slides of Vibhav Gogate and David Sontag

Supervised Learning



- Input: labeled training data
 - i.e., data plus desired output
- Assumption: there exists a function f that maps data items x to their correct labels
- Goal: construct an approximation to f

Today



- We've been focusing on linear separators
 - Relatively easy to learn (using standard techniques)
 - Easy to picture, but not clear if data will be separable
- Next two lectures we'll focus on other hypothesis spaces
 - Decision trees
 - Nearest neighbor classification



- Suppose that you go to your doctor with flu-like symptoms
 - How does your doctor determine if you have a flu that requires medical attention?



- Suppose that you go to your doctor with flu-like symptoms
 - How does your doctor determine if you have a flu that requires medical attention?
 - Check a list of symptoms:
 - Do you have a fever over 100.4 degrees Fahrenheit?
 - Do you have a sore throat or a stuffy nose?
 - Do you have a dry cough?

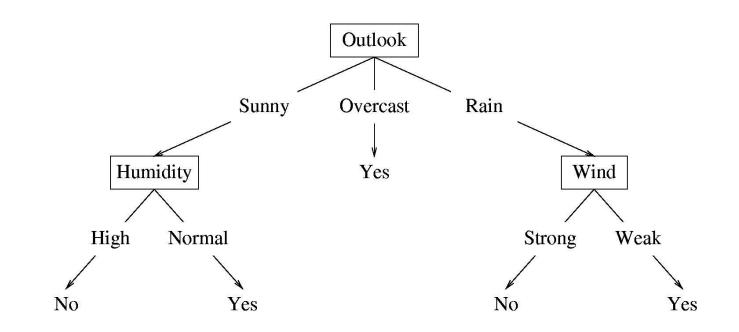


- Just having some symptoms is not enough, you should also not have symptoms that are not consistent with the flu
- For example,
 - If you have a fever over 100.4 degrees Fahrenheit?
 - And you have a sore throat or a stuffy nose?
 - You probably do not have the flu (most likely just a cold)



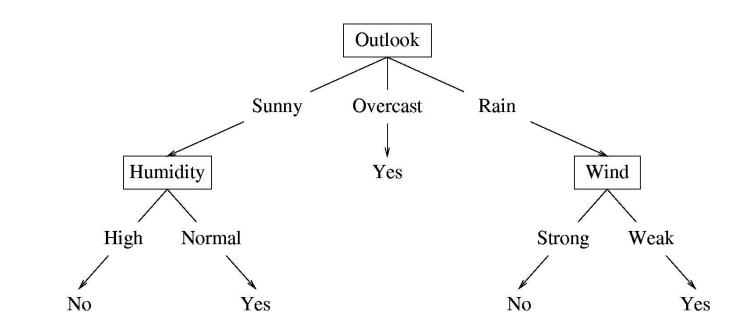
- In other words, your doctor will perform a series of tests and ask a series of questions in order to determine the likelihood of you having a severe case of the flu
- This is a method of coming to a diagnosis (i.e., a classification of your condition)
- We can view this decision making process as a tree





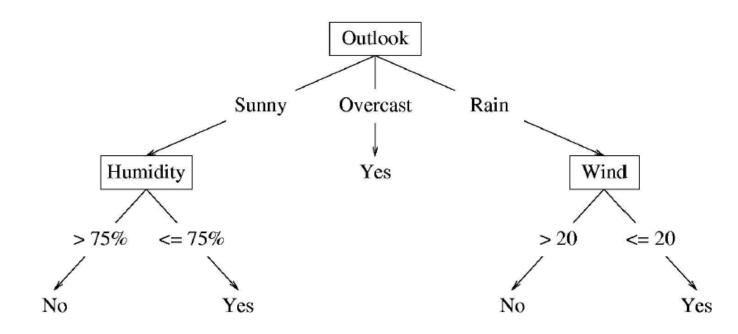
- A tree in which each internal (non-leaf) node tests the value of a particular feature
- Each leaf node specifies a class label (in this case whether or not you should play tennis)





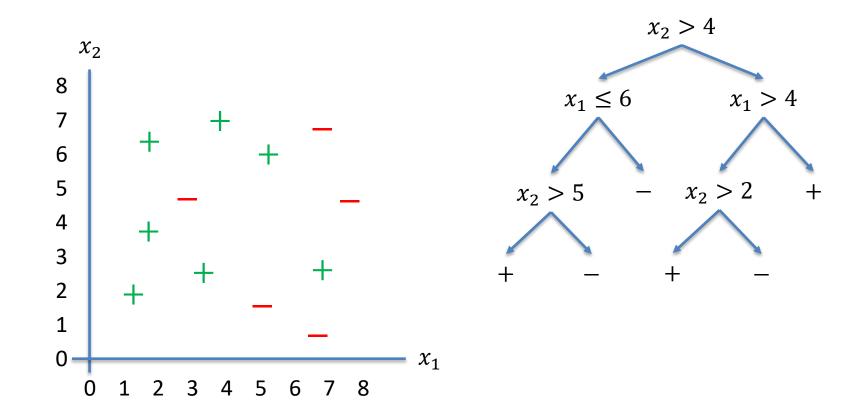
- Features: (Outlook, Humidity, Wind)
- Classification is performed root to leaf
 - The feature vector (Sunny, Normal, Strong) would be classified as a yes instance



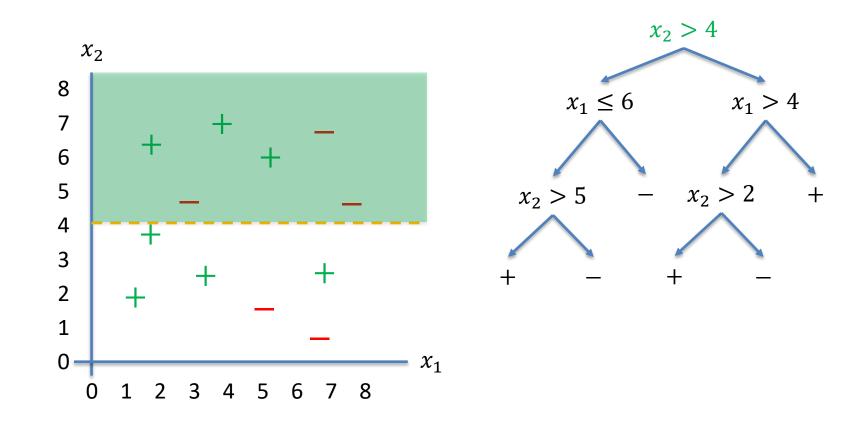


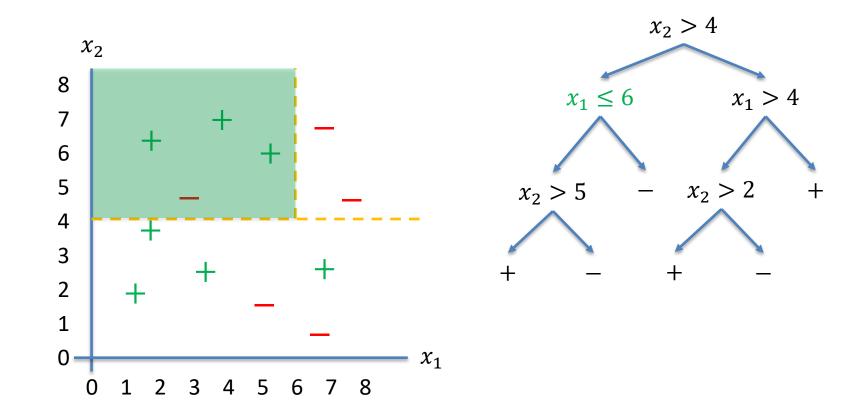
- Can have continuous features too
 - Internal nodes for continuous features correspond to thresholds



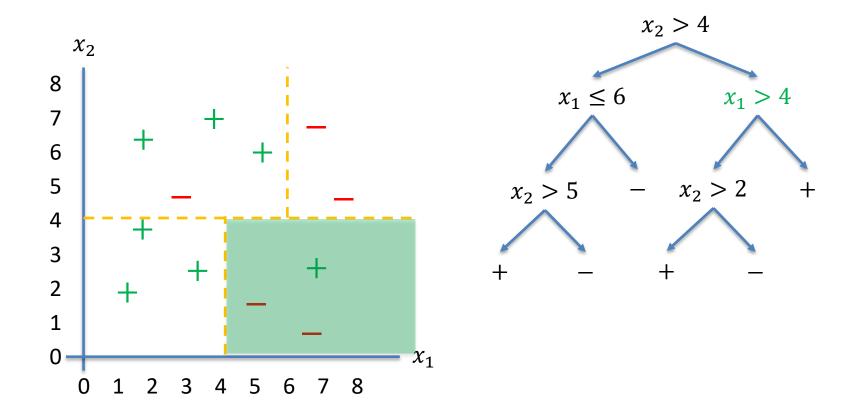




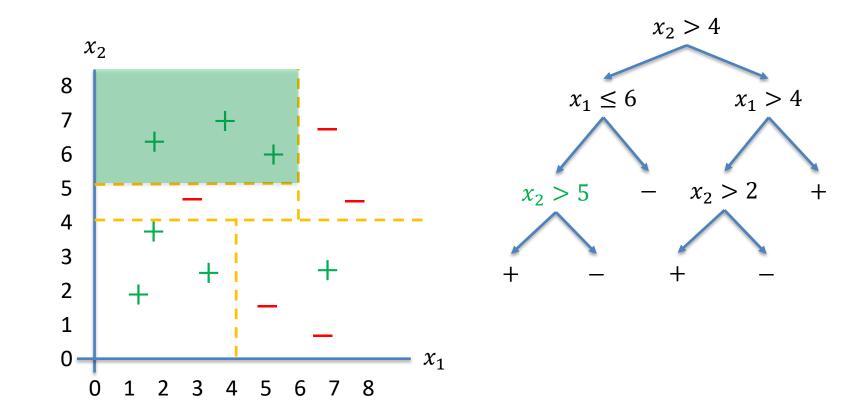




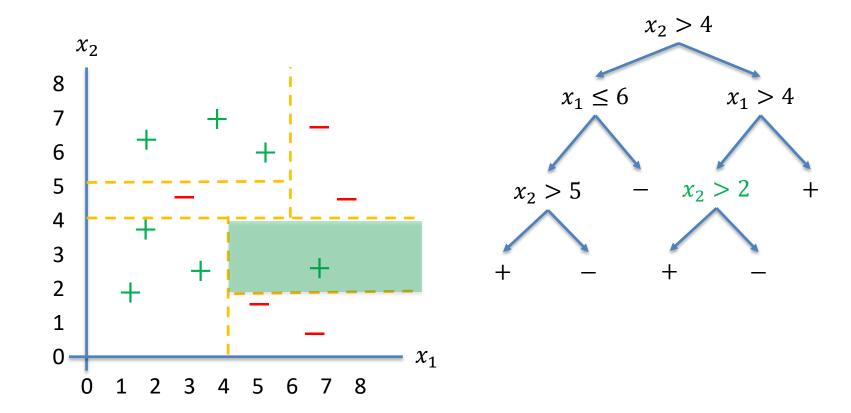






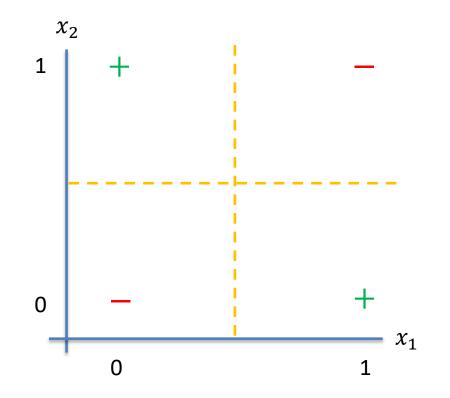








• Worst case decision tree may require exponentially (in the dimension of the data) many nodes



Decision Tree Learning



- Basic decision tree building algorithm:
 - Pick some feature/attribute
 - Partition the data based on the value of this attribute
 - Recurse over each new partition

Decision Tree Learning



- Basic decision tree building algorithm:
 - Pick some feature/attribute (how to pick the "best"?)
 - Partition the data based on the value of this attribute
 - Recurse over each new partition (when to stop?)

We'll focus on the discrete case first (i.e., each feature takes a value in some finite set)



• What functions can be represented by decision trees?

• Are decision trees unique?



- What functions can be represented by decision trees?
 - Every function of +/- can be represented by a sufficiently complicated decision tree
- Are decision trees unique?
 - No, many different decision trees are possible for the same set of labels

Choosing the Best Attribute

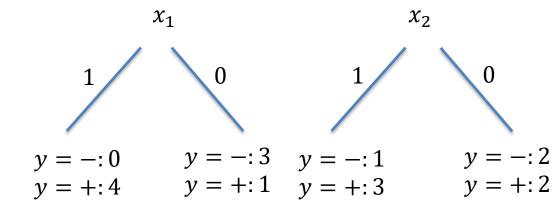


- Because the complexity of storage and classification increases with the size of the tree, should prefer smaller trees
 - Simplest models that explain the data are usually preferred over more complicated ones
 - Finding the smallest tree is an NP-hard problem
 - Instead, use a greedy heuristic based approach to pick the best attribute at each stage



 $x_1, x_2 \in \{0, 1\}$

Which attribute should you split on?

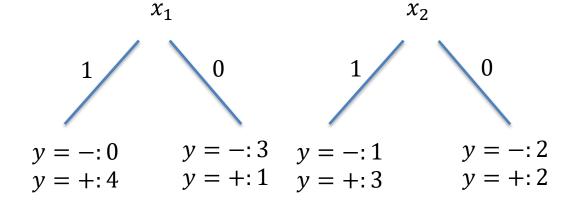


<i>x</i> ₁	<i>x</i> ₂	у
1	1	+
1	0	+
1	1	+
1	0	+
0	1	+
0	0	—
0	1	_
0	0	



 $x_1, x_2 \in \{0, 1\}$

Which attribute should you split on?



Can think of these counts as probability distributions over the labels: if x = 1, the probability that y = + is equal to 1

<i>x</i> ₁	<i>x</i> ₂	у
1	1	+
1	0	+
1	1	+
1	0	+
0	1	+
0	0	—
0	1	—
0	0	_



- The selected attribute is a good split if we are more "certain" about the classification after the split
 - If each partition with respect to the chosen attribute has a distinct class label, we are completely certain about the classification after partitioning
 - If the class labels are evenly divided between the partitions, the split isn't very good (we are very uncertain about the label for each partition)
 - What about other situations? How do you measure the uncertainty of a random process?

Discrete Probability



- Sample space specifies the set of possible outcomes
 - For example, $\Omega = \{H, T\}$ would be the set of possible outcomes of a coin flip
- Each element $\omega \in \Omega$ is associated with a number $p(\omega) \in [0,1]$ called a **probability**

$$\sum_{\omega\in\Omega}p(\omega)=1$$

• For example, a biased coin might have p(H) = .6 and p(T) = .4

Discrete Probability



- An event is a subset of the sample space
 - Let Ω = {1, 2, 3, 4, 5, 6} be the 6 possible outcomes of a dice role
 - A = {1, 5, 6} ⊆ Ω would be the event that the dice roll comes up as a one, five, or six
- The probability of an event is just the sum of all of the outcomes that it contains

•
$$p(A) = p(1) + p(5) + p(6)$$

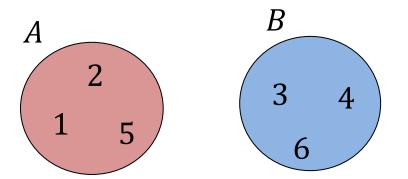


• Two events A and B are **independent** if

 $p(A \cap B) = p(A)P(B)$

Let's suppose that we have a fair die: $p(1) = \dots = p(6) = 1/6$

If $A = \{1, 2, 5\}$ and $B = \{3, 4, 6\}$ are A and B indpendent?



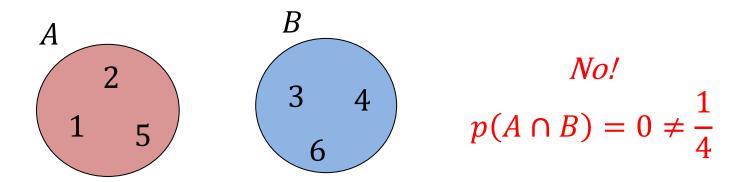


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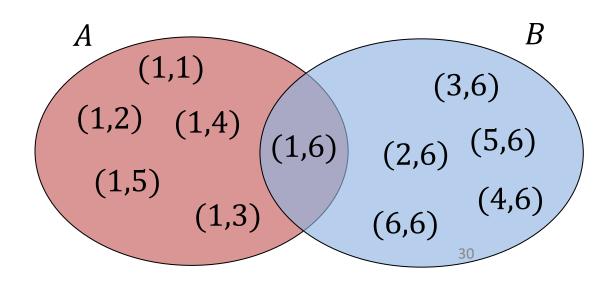
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Independence



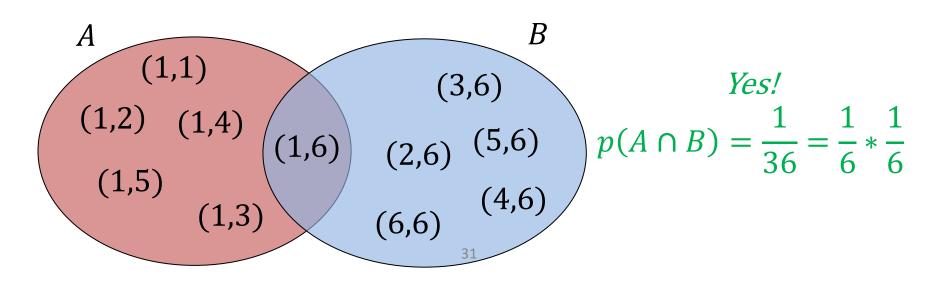
- Now, suppose that Ω = {(1,1), (1,2), ..., (6,6)} is the set of all possible rolls of two unbiased dice
- Let A = {(1,1), (1,2), (1,3), ..., (1,6)} be the event that the first die is a one and let B = {(1,6), (2,6), ..., (6,6)} be the event that the second die is a six
- Are A and B independent?



Independence



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The conditional probability of an event A given an event B with p(B) > 0 is defined to be

$$p(A|B) = \frac{p(A \cap B)}{P(B)}$$

- This is the probability of the event A ∩ B over the sample space
 Ω' = B
- Some properties:
 - $\sum_{\omega \in \Omega} p(\omega|B) = 1$
 - If A and B are independent, then p(A|B) = p(A)

Discrete Random Variables



- A discrete random variable, X, is a function from the state space
 Ω into a discrete space D
 - For each $x \in D$,

$$p(X = x) \equiv p(\{\omega \in \Omega : X(\omega) = x\})$$

is the probability that X takes the value x

• p(X) defines a probability distribution

•
$$\sum_{x \in D} p(X = x) = 1$$

• Random variables partition the state space into disjoint events

Example: Pair of Dice



- Let Ω be the set of all possible outcomes of rolling a pair of dice
- Let p be the uniform probability distribution over all possible outcomes in $\boldsymbol{\Omega}$
- Let $X(\omega)$ be equal to the sum of the value showing on the pair of dice in the outcome ω
 - p(X = 2) = ?
 - p(X = 8) = ?

Example: Pair of Dice



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$$p(X=2) = \frac{1}{36}$$

•
$$p(X = 8) = ?$$

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$$p(X=2) = \frac{1}{36}$$

•
$$p(X = 8) = \frac{5}{36}$$

Discrete Random Variables

• We can have vectors of random variables as well

 $X(\omega) = [X_1(\omega), \dots, X_n(\omega)]$

• The joint distribution is $p(X_1 = x_1, ..., X_n = x_n)$ is

$$p(X_1 = x_1 \cap \dots \cap X_n = x_n)$$

 $p(x_1,\ldots,x_n)$

typically written as

Because
$$X_i = x_i$$
 is an event, all of the same rules from basic probability apply



Entropy

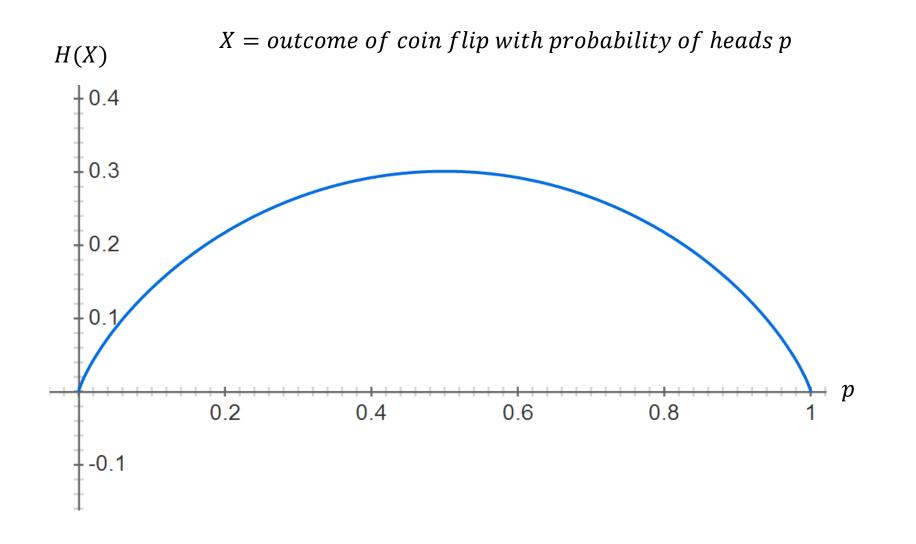


 A standard way to measure uncertainty of a random variable is to use the entropy

$$H(Y) = -\sum_{Y=y} p(Y=y) \log p(Y=y)$$

- Entropy is maximized for uniform distributions
- Entropy is minimized for distributions that place all their probability on a single outcome

Entropy of a Coin Flip



Conditional Entropy



• We can also compute the entropy of a random variable conditioned on a different random variable

$$H(Y|X) = -\sum_{x} p(X = x) \sum_{y} p(Y = y|X = x) \log p(Y = y|X = x)$$

- This is called the conditional entropy
- This is the amount of information needed to quantify the random variable *Y* given the random variable *X*



 Using entropy to measure uncertainty, we can greedily select an attribute that guarantees the largest expected decrease in entropy (with respect to the empirical partitions)

$$IG(X) = H(Y) - H(Y|X)$$

- Called information gain
- Larger information gain corresponds to less uncertainty about *Y* given *X*
 - Note that $H(Y|X) \leq H(Y)$

Decision Tree Learning

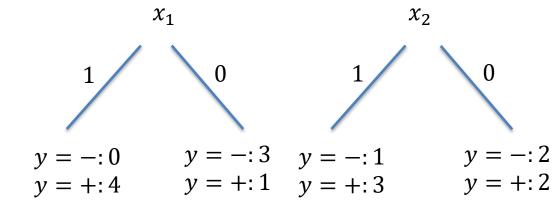


- Basic decision tree building algorithm:
 - Pick the feature/attribute with the highest information gain
 - Partition the data based on the value of this attribute
 - Recurse over each new partition



 $x_1, x_2 \in \{0, 1\}$

Which attribute should you split on?



<i>x</i> ₁	<i>x</i> ₂	у
1	1	+
1	0	+
1	1	+
1	0	+
0	1	+
0	0	
0	1	_
0	0	_

What is the information gain in each case?



 $x_1, x_2 \in \{0, 1\}$

Which attribute should you split on?	<i>x</i> ₁	<i>x</i> ₂	У		
x_1 x_2	1	1	+		
	1	0	+		
	1	1	+		
	1	0	+		
y = -:0 $y = -:3$ $y = -:1$ $y = -:2y = +:4$ $y = +:1$ $y = +:3$ $y = +:2$	0	1	+		
y = 1.4 $y = 1.1$ $y = 1.5$ $y = 1.2$	0	0	—		
	0	1	—		
$H(Y) = -\frac{5}{8}\log\frac{5}{8} - \frac{3}{8}\log\frac{3}{8}$	0	0	_		
$H(Y) = -\frac{1}{8} \log \frac{1}{8} - \frac{1}{8} \log \frac{1}{8}$ $H(Y X_1) = .5[-0 \log 0 - 1 \log 1] + .5[75 \log .7525 \log .25]$ $H(Y X_2) = .5[5 \log .55 \log .5] + .5[75 \log .7525 \log .25]$					

 $H(Y) - H(Y|X_1) - H(Y) + H(Y|X_2) = -.5 \log .5 > 0$ Should split on x_1



- If the current set is "pure" (i.e., has a single label in the output), stop
- If you run out of attributes to recurse on, even if the current data set isn't pure, stop and use a majority vote
- If a partition contains no data points, use the majority vote at its parent in the tree
- If a partition contains no data items, nothing to recurse on
- For fixed depth decision trees, the final label is determined by majority vote



- For continuous attributes, use threshold splits
 - Split the tree into $x_k < t$ and $x_k \ge t$
 - Can split on the same attribute multiple times on the same path down the tree
- How to pick the threshold *t*?

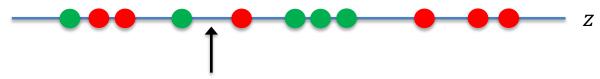


- For continuous attributes, use threshold splits
 - Split the tree into $x_k < t$ and $x_k \ge t$
 - Can split on the same attribute multiple times on the same path down the tree
- How to pick the threshold *t*?
 - Try every possible *t*

How many possible *t* are there?



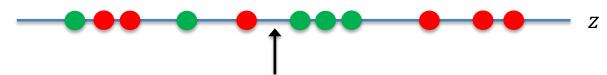
• Sort the data according to the k^{th} attribute: $z_1 > z_2 > \cdots > z_n$



• Only a finite number of thresholds make sense

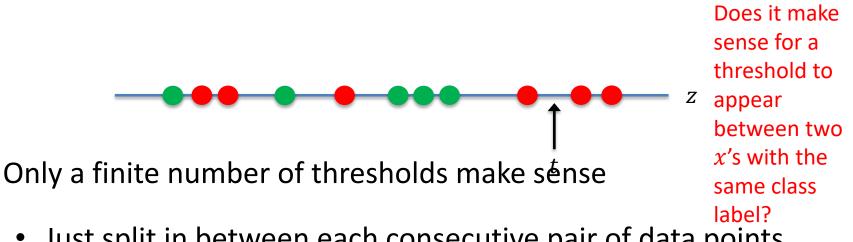


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 - Just split in between each consecutive pair of data points (e.g., splits of the form $t = \frac{z_i + z_{i+1}}{2}$)

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• Just split in between each consecutive pair of data points (e.g., splits of the form $t = \frac{z_i + z_{i+1}}{2}$)

- Compute the information gain of each threshold
- Let *X*: *t* denote splitting with threshold *t* and compute

$$H(Y|X:t) = -p(X < t) \sum_{y} p(Y = y|X < t) \log p(Y = y|X < t) + -p(X \ge t) \sum_{y} p(Y = y|X \ge t) \log p(Y = y|X \ge t)$$

 In the learning algorithm, maximize over all attributes and all possible thresholds of the real-valued attributes

$$\max_{t} H(Y) - H(Y|X:t), \text{ for real-valued } X$$
$$H(Y) - H(Y|X), \text{ for discrete } X$$





- Because of speed/ease of implementation, decision trees are quite popular
 - Can be used for regression too
- Decision trees will **always** overfit!
 - It is always possible to obtain zero training error on the input data with a deep enough tree (if there is no noise in the labels)
 - Solution?