



Xi'an Jiaotong-Liverpool University

西交利物浦大学

**INT305 Machine Learning**

**Lecture 7**

**Decision Trees & Bias-Variance Decomposition**

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# Today

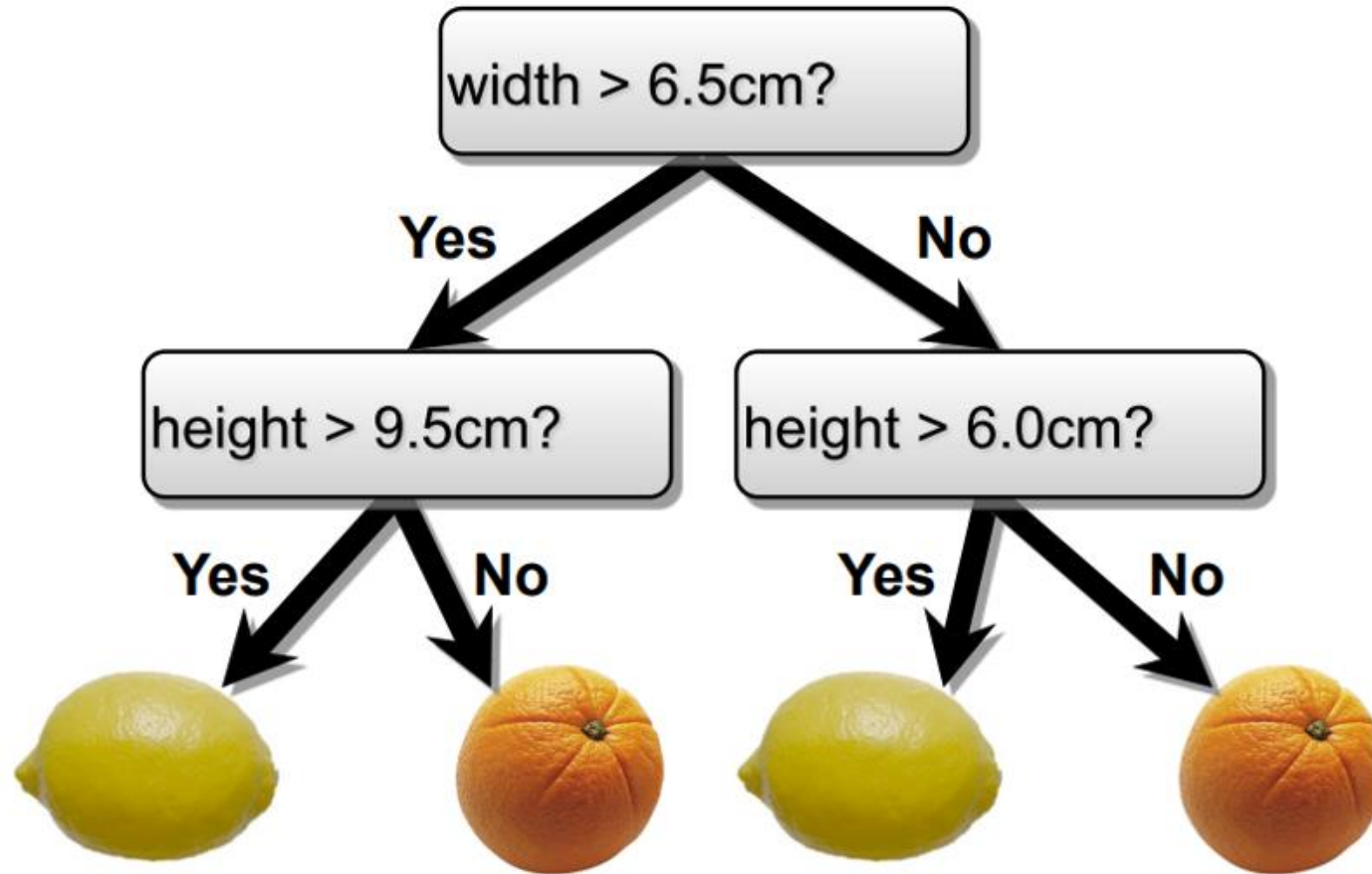
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- Decision Trees
  - Simple but powerful learning algorithm
  - Used widely in Kaggle competitions
  - Let us motivate concepts from information theory (entropy, mutual information, etc.)
- Bias-variance decomposition
  - Let us motivate methods for combining different classifiers.

# Decision Trees

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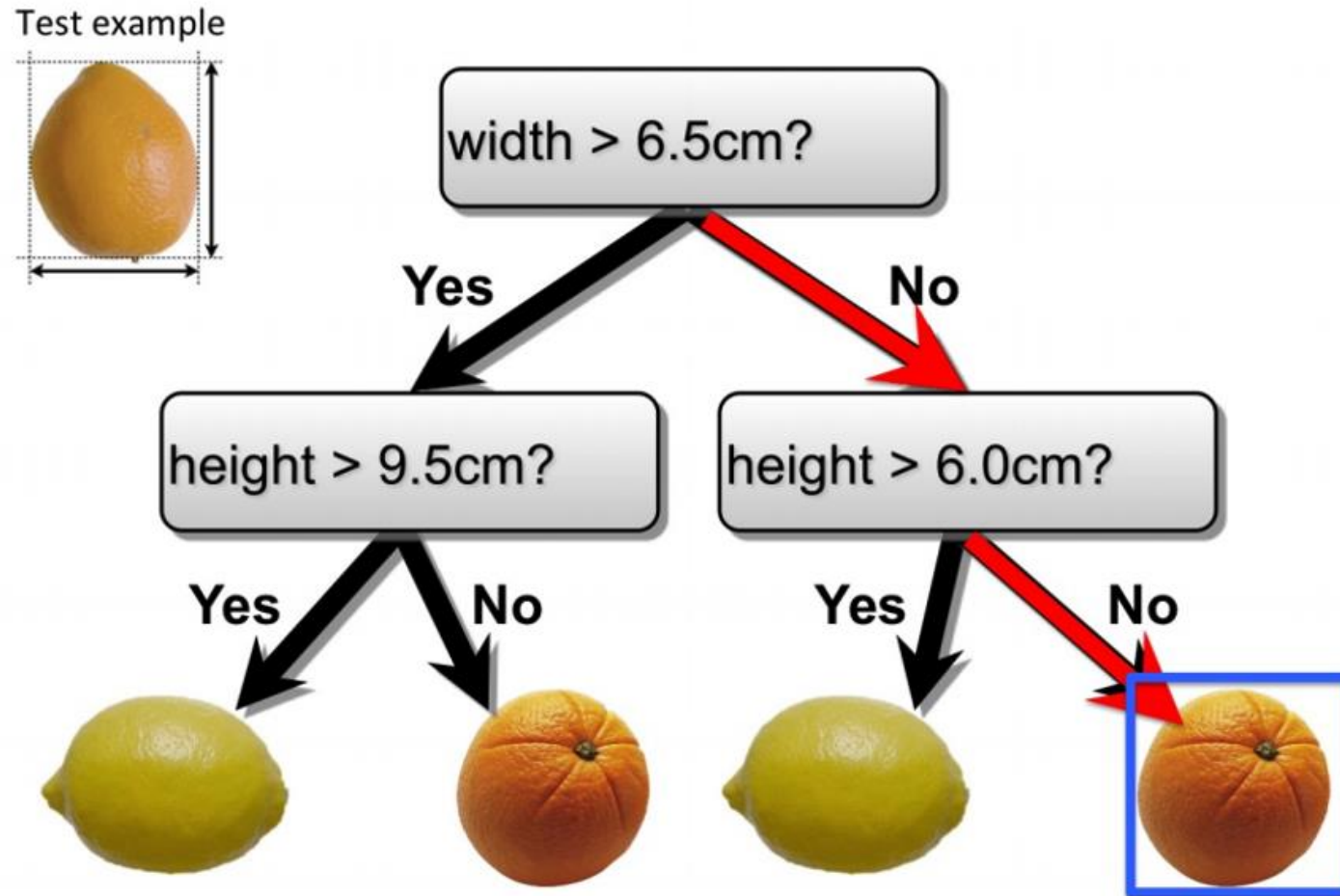
- Make predictions by splitting on features according to a tree structure.



# Decision Trees

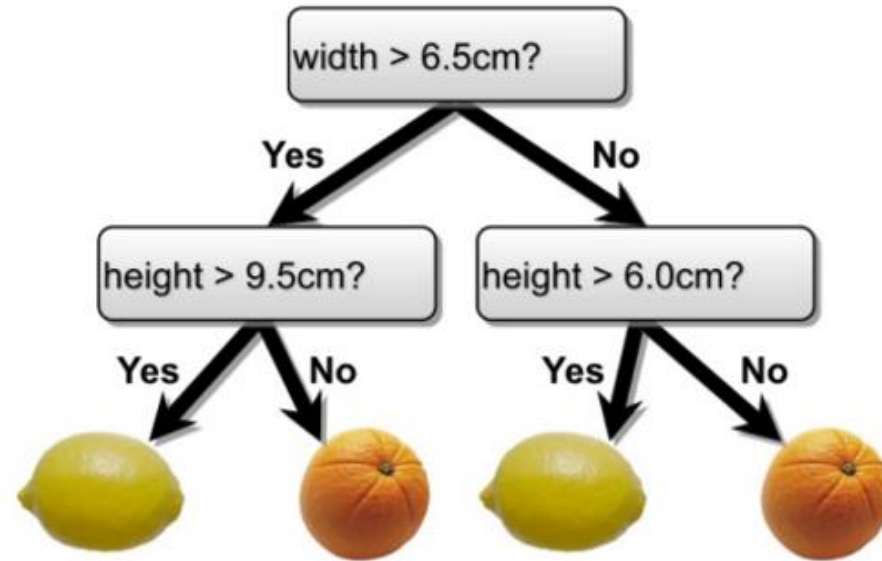
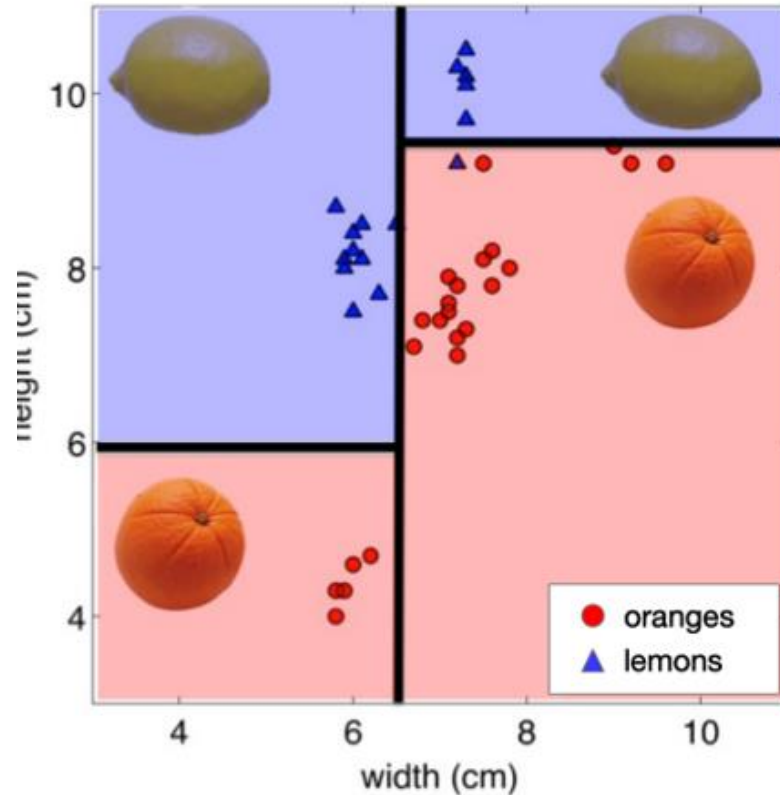
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- Make predictions by splitting on features according to a tree structure.



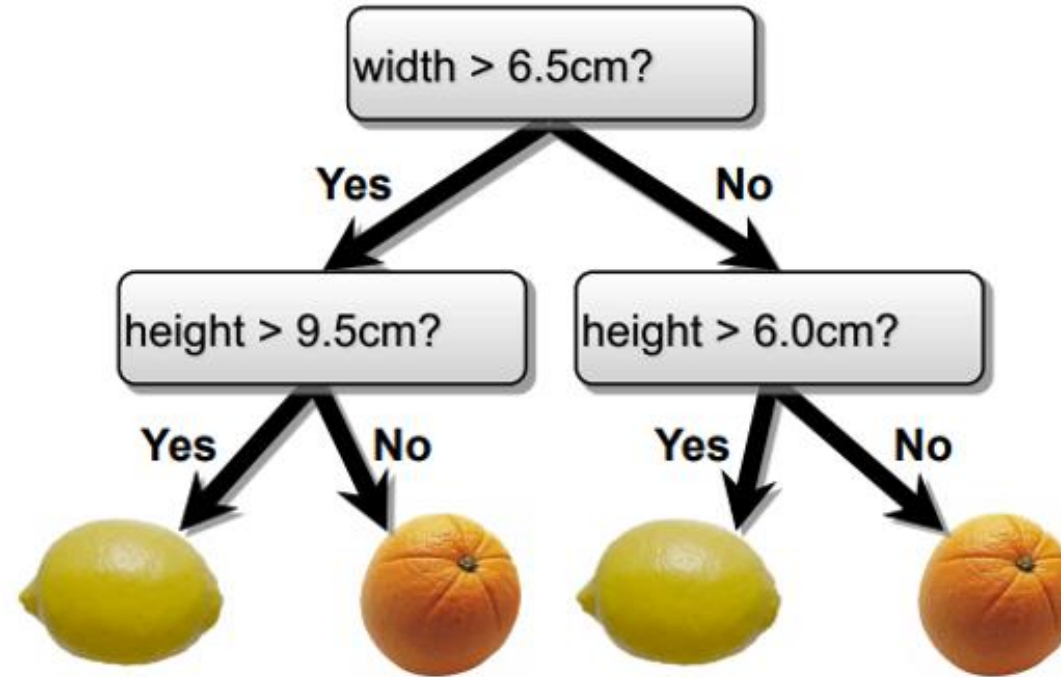
# Decision Trees— Continuous Features

- Split *continuous features* by checking whether that feature is greater than or less than some threshold.
- Decision boundary is made up of axis-aligned planes.



# Decision Trees

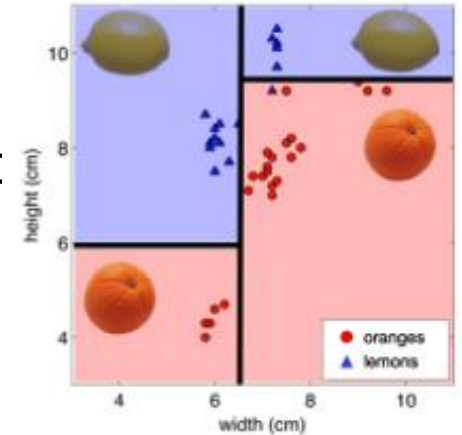
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- Internal nodes test a feature
- Branching is determined by the feature value
- Leaf nodes are outputs (predictions)

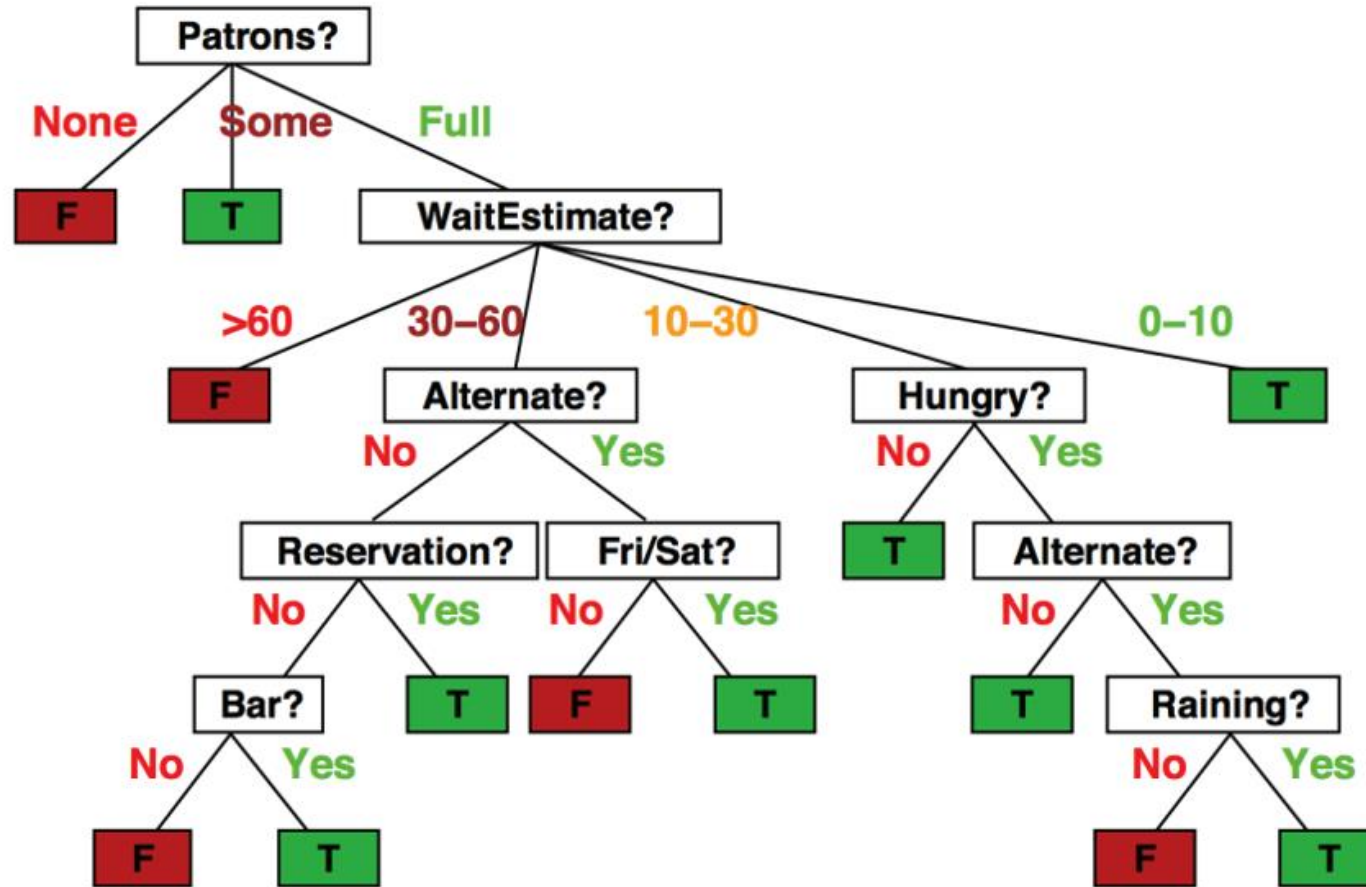
# Decision Trees—Classification and Regression

- Each path from root to a leaf defines a region  $R_m$  of input space
- Let  $\{(x^{(m_1)}, t^{(m_1)}), \dots, (x^{(m_k)}, t^{(m_k)})\}$  be the training examples that fall into  $R_m$
- **Classification tree** (we will focus on this):
  - discrete output
  - leaf value  $y^m$  typically set to the most common value in  $\{t^{(m_1)}, \dots, t^{(m_k)}\}$
- **Regression tree**:
  - Continuous output
  - leaf value  $y^m$  typically set to the mean value in  $\{t^{(m_1)}, \dots, t^{(m_k)}\}$



# Decision Trees—Discrete Features

- Will I wait at this restaurant for dinner?





# Decision Trees—Discrete Features

- Split **discrete features** into a partition of possible values.

Example	Input Attributes										Goal
	<i>Alt</i>	<i>Bar</i>	<i>Fri</i>	<i>Hun</i>	<i>Pat</i>	<i>Price</i>	<i>Rain</i>	<i>Res</i>	<i>Type</i>	<i>Est</i>	<i>WillWait</i>
$x_1$	Yes	No	No	Yes	Some	\$\$\$	No	Yes	French	0–10	$y_1 = \text{Yes}$
$x_2$	Yes	No	No	Yes	Full	\$	No	No	Thai	30–60	$y_2 = \text{No}$
$x_3$	No	Yes	No	No	Some	\$	No	No	Burger	0–10	$y_3 = \text{Yes}$
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$x_7$	No	Yes	No	No	None	\$	Yes	No	Burger	0–10	$y_7 = \text{No}$
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$x_{10}$	Yes	Yes	Yes	Yes	Full	\$\$\$	No	Yes	Italian	10–30	$y_{10} = \text{No}$
$x_{11}$	No	No	No	No	None	\$	No	No	Thai	0–10	$y_{11} = \text{No}$
$x_{12}$	Yes	Yes	Yes	Yes	Full	\$	No	No	Burger	30–60	$y_{12} = \text{Yes}$

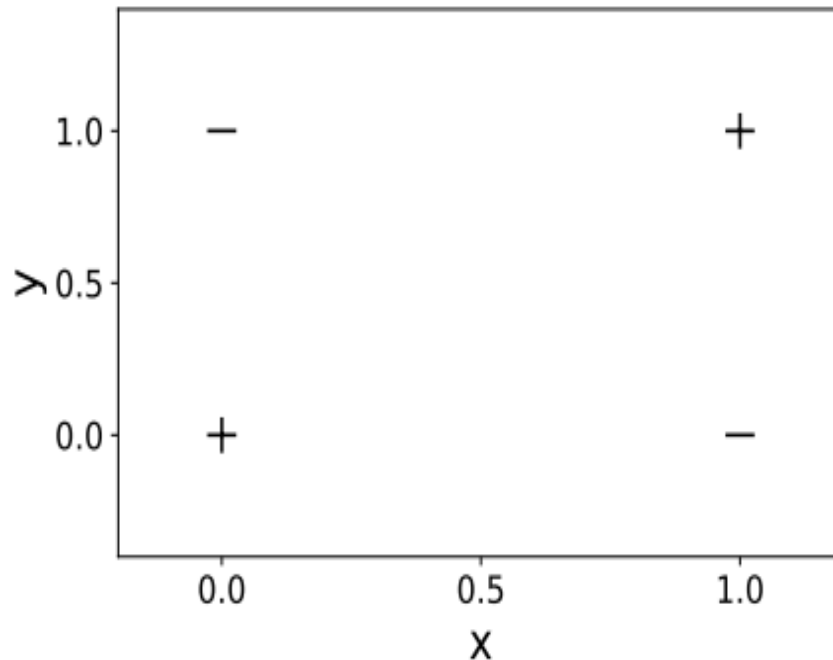
1.	Alternate: whether there is a suitable alternative restaurant nearby.
2.	Bar: whether the restaurant has a comfortable bar area to wait in.
3.	Fri/Sat: true on Fridays and Saturdays.
4.	Hungry: whether we are hungry.
5.	Patrons: how many people are in the restaurant (values are None, Some, and Full).
6.	Price: the restaurant's price range (\$, \$\$, \$\$\$).
7.	Raining: whether it is raining outside.
8.	Reservation: whether we made a reservation.
9.	Type: the kind of restaurant (French, Italian, Thai or Burger).
10.	WaitEstimate: the wait estimated by the host (0-10 minutes, 10-30, 30-60, >60).

Features:

# Attempts

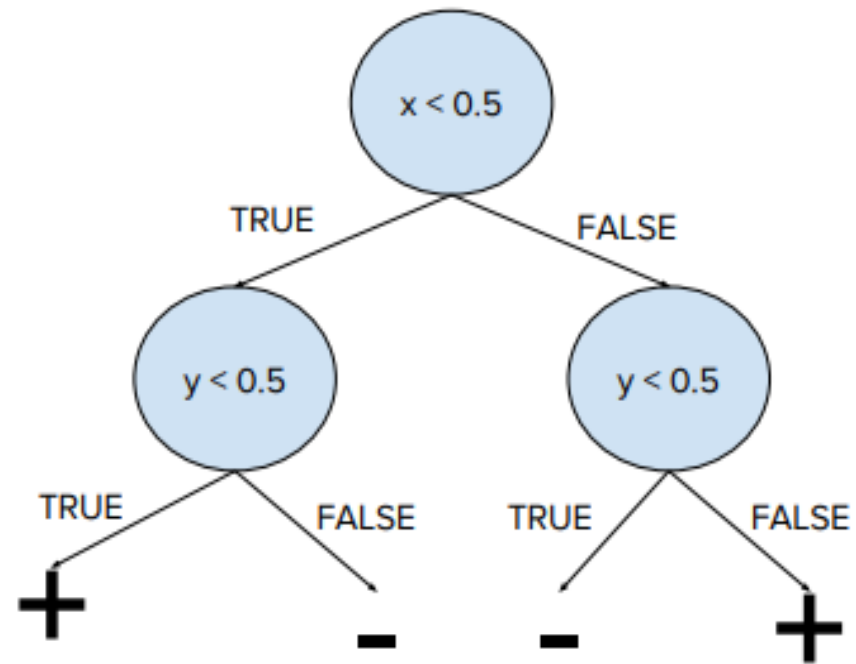
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The drawing below shows a dataset. Each example in the dataset has two inputs features  $x$  and  $y$ , and maybe classified as a positive example (labelled  $+$ ) or a negative example (labelled  $-$ ). Draw a decision tree which correctly classifies each example in the dataset.



# Attempts

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# Learning Decision Trees

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- For any training set we can construct a decision tree that has exactly the one leaf for every training point, but it probably won't generalize.
  - Decision trees are universal function approximators.
- But, finding the smallest decision tree that correctly classifies a training set is NP complete.
  - If you are interested, check: Hyafil & Rivest'76.
- So, how do we construct a useful decision tree?

# Learning Decision Trees

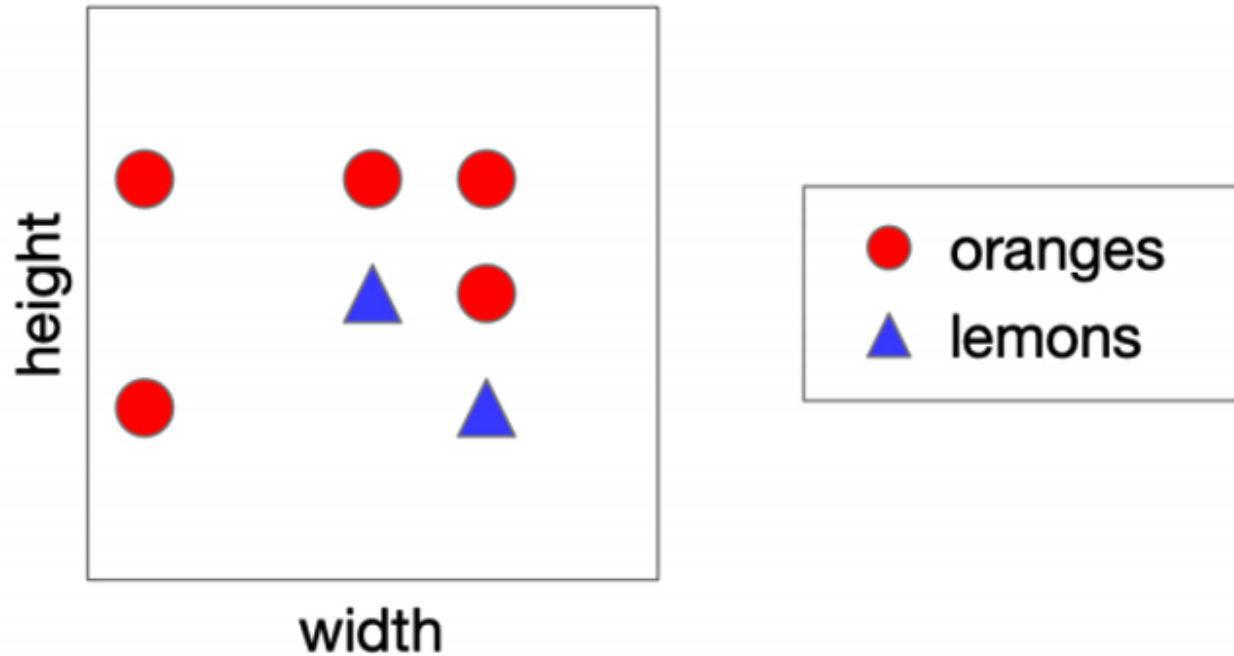
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- Resort to a **greedy heuristic**:
  - Start with the whole training set and an empty decision tree.
  - Pick a feature and candidate split that would most reduce the loss.
  - Split on that feature and recurse on subpartitions.
- Which loss should we use?
  - Let's see if misclassification rate is a good loss.

# Choosing a Good Split

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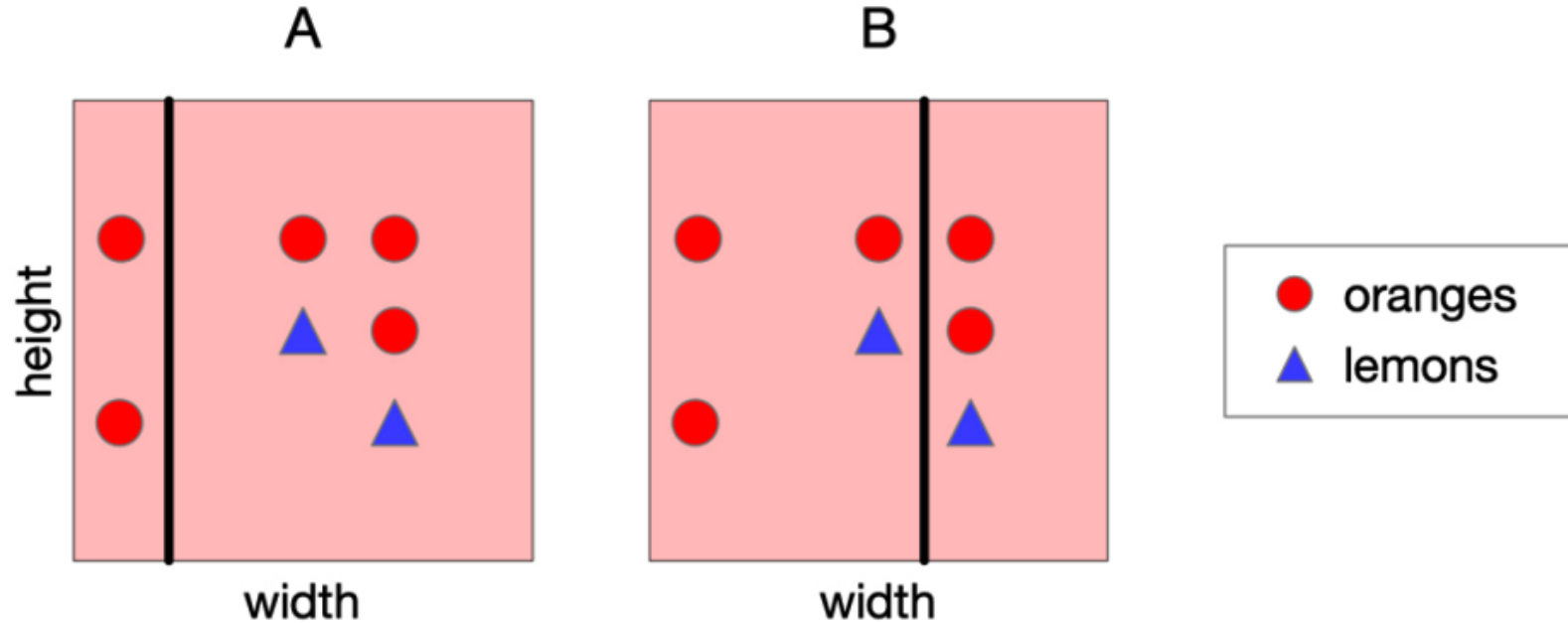
- Consider the following data. Let's split on width.



# Choosing a Good Split

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- Recall: classify by majority.

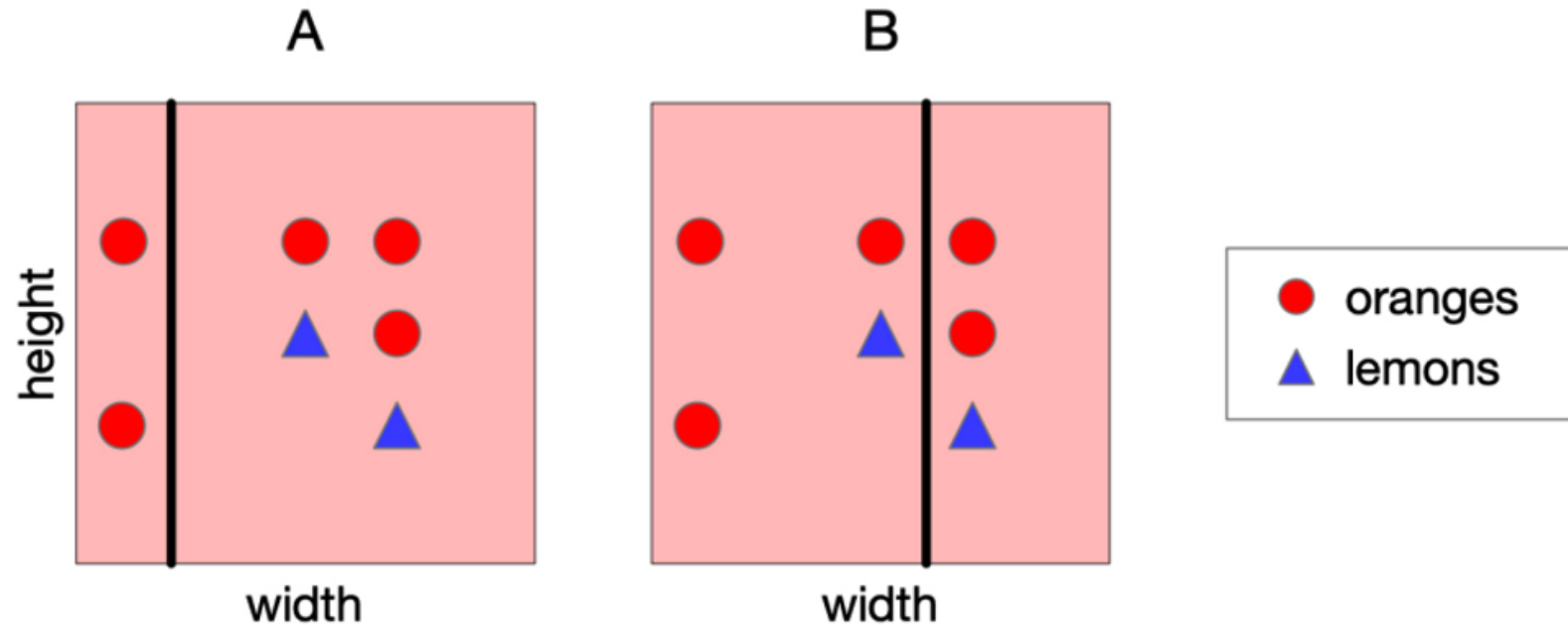


- A and B have the same misclassification rate, so which is the best split? Vote!

# Choosing a Good Split

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- A feels like a better split, because the left-hand region is very certain about whether the fruit is an orange.



- Can we quantify this?



# Choosing a Good Split

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- How can we quantify uncertainty in prediction for a given leaf node?
  - If all examples in leaf have same class: good, low uncertainty
  - If each class has same amount of examples in leaf: bad, high uncertainty
- **Idea:** use counts at leaves to define probability distributions; use a probabilistic notion of uncertainty to decide splits.
- A brief detour through information theory...

# Quantifying Uncertainty

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- The **entropy** of a discrete random variable is a number that quantifies the **uncertainty** inherent in its possible outcomes.
- The mathematical definition of entropy that we give in a few slides may seem arbitrary, but it can be motivated axiomatically.
  - If you're interested, check: *Information Theory* by Robert Ash.
- To explain entropy, consider flipping two different coins...

# We Flip Two Different Coins

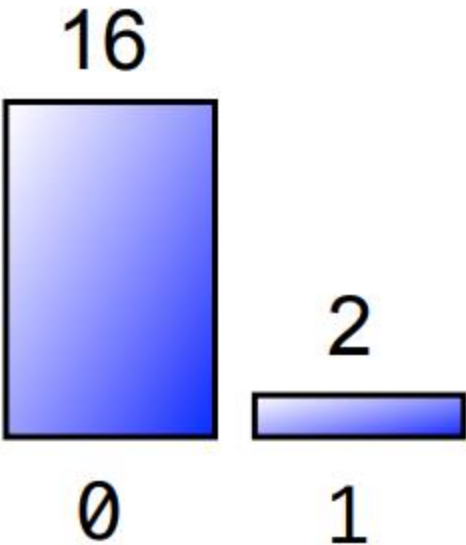
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Sequence 1:

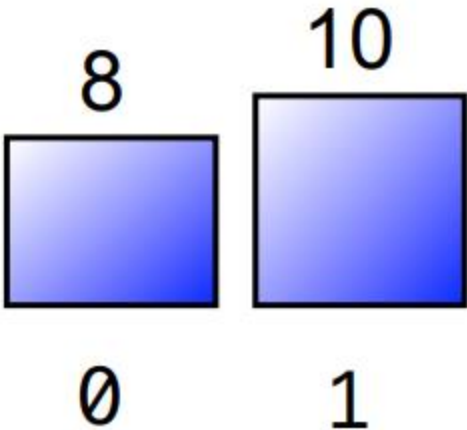
0 0 0 1 0 0 0 0 0 0 0 0 0 0 0 0 1 0 0 ... ?

Sequence 2:

0 1 0 1 0 1 1 1 0 1 0 0 1 1 0 1 0 1 ... ?



versus

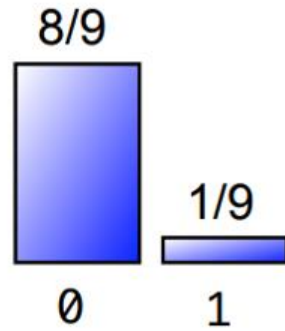


# Quantifying Uncertainty

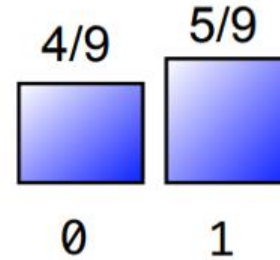
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- The entropy of a loaded coin with probability  $p$  of heads is given by

$$-p\log_2(p) - (1 - p)\log_2(1 - p)$$



$$-\frac{8}{9}\log_2\frac{8}{9} - \frac{1}{9}\log_2\frac{1}{9} \approx \frac{1}{2}$$



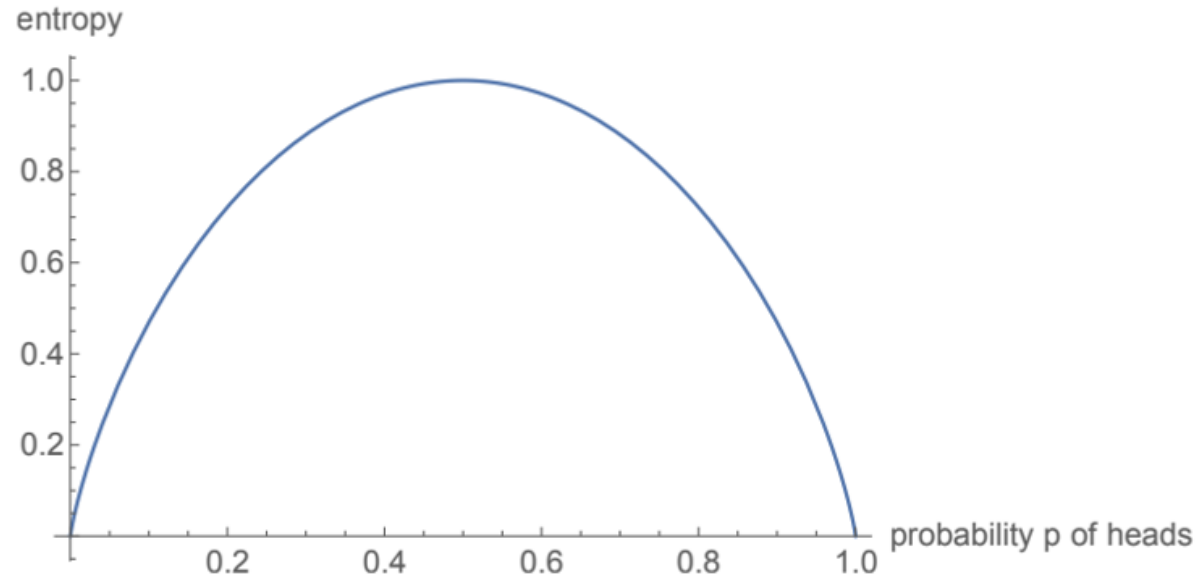
$$-\frac{4}{9}\log_2\frac{4}{9} - \frac{5}{9}\log_2\frac{5}{9} \approx 0.99$$

- Notice: the coin whose outcomes are more certain has a lower entropy
- In the extreme case  $p = 0$  or  $p = 1$ , we were certain of the outcome before observing. So, we gained no certainty by observing it, i.e., entropy is 0.

# Quantifying Uncertainty

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- Can also think of **entropy** as the expected information content of a random draw from a probability distribution.



- Claude Shannon showed: you cannot store the outcome of a random draw using fewer expected bits than the entropy without losing information.
- So units of entropy are **bits**; a fair coin flip has 1 bit of entropy.

# Entropy

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- More generally, the entropy of a discrete random variable  $Y$  is given by

$$H(Y) = - \sum_{y \in Y} p(y) \log_2 p(y)$$

- “High Entropy”
  - Variable has a uniform like distribution over many outcomes
  - Flat histogram
  - Values sampled from it are less predictable
- “Low Entropy”
  - Distribution is concentrated on only a few outcomes
  - Histogram is concentrated in a few areas
  - Values sampled from it are more predictable

# Entropy

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- Suppose we observe partial information  $X$  about a random variable  $Y$ 
  - For example,  $X = \text{sign}(Y)$ .
- We want to work towards a definition of the expected amount of information that will be conveyed about  $Y$  by observing  $X$ .
  - Or equivalently, the expected reduction in our uncertainty about  $Y$  after observing  $X$ .

# Entropy of a Joint Distribution

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- Example:  $X = \{\text{Raining, Not raining}\}$ ,  $Y = \{\text{Cloudy, Not cloudy}\}$

	Cloudy	Not Cloudy
Raining	24/100	1/100
Not Raining	25/100	50/100

$$\begin{aligned} H(X, Y) &= - \sum_{x \in X} \sum_{y \in Y} p(x, y) \log_2 p(x, y) \\ &= - \frac{24}{100} \log_2 \frac{24}{100} - \frac{1}{100} \log_2 \frac{1}{100} - \frac{25}{100} \log_2 \frac{25}{100} - \frac{50}{100} \log_2 \frac{50}{100} \\ &\approx 1.56 \text{bits} \end{aligned}$$



# Specific Conditional Entropy

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- Example:  $X = \{\text{Raining, Not raining}\}$ ,  $Y = \{\text{Cloudy, Not cloudy}\}$

	Cloudy	Not Cloudy
Raining	24/100	1/100
Not Raining	25/100	50/100

- What is the entropy of cloudiness  $Y$ , **given that it is raining**?

$$\begin{aligned} H(Y|X = x) &= - \sum_{y \in Y} p(y|x) \log_2 p(y|x) \\ &= -\frac{24}{25} \log_2 \frac{24}{25} - \frac{1}{25} \log_2 \frac{1}{25} \\ &\approx 0.24 \text{bits} \end{aligned}$$

- We used:  $p(y|x) = \frac{p(x,y)}{p(x)}$  and  $p(x) = \sum_y p(x, y)$  (sum in a row)

# Conditional Entropy

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	Cloudy	Not Cloudy
Raining	24/100	1/100
Not Raining	25/100	50/100

- The expected conditional entropy:

$$\begin{aligned} H(Y|X) &= \sum_{x \in X} p(x) H(Y|X = x) \\ &= - \sum_{x \in X} \sum_{y \in Y} p(x, y) \log_2 p(y|x) \end{aligned}$$

# Conditional Entropy

---

- Example:  $X = \{\text{Raining, Not raining}\}$ ,  $Y = \{\text{Cloudy, Not cloudy}\}$

	Cloudy	Not Cloudy
Raining	24/100	1/100
Not Raining	25/100	50/100

- What is the entropy of cloudiness  $Y$ , **given the knowledge of whether or not it is raining?**

$$\begin{aligned} H(Y|X) &= \sum_{x \in X} p(x) H(Y|X = x) \\ &= \frac{1}{4} H(\text{cloudiness} | \text{is raining}) + \frac{3}{4} H(\text{cloudiness} | \text{not raining}) \\ &\approx 0.75 \text{bits} \end{aligned}$$

# Conditional Entropy

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- Some useful properties
  - $H$  is always non-negative
  - Chain rule:  $H(X, Y) = H(X|Y) + H(Y) = H(Y|X) + H(X)$
  - If  $X$  and  $Y$  independent, then  $X$  does not affect our uncertainty about  $Y$ :  $H(Y|X) = H(Y)$
  - But knowing  $Y$  makes our knowledge of  $Y$  certain:  $H(Y|Y) = 0$
  - By knowing  $X$ , we can only decrease uncertainty about  $Y$ :  $H(Y|X) \leq H(Y)$

# Information Gain

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	Cloudy	Not Cloudy
Raining	24/100	1/100
Not Raining	25/100	50/100

- How much more certain am I about whether it's cloudy if I'm told whether it is raining?  
My uncertainty in  $Y$  minus my expected uncertainty that would remain in  $Y$  after seeing  $X$ .
- This is called the **information gain**  $IG(Y|X)$  in  $Y$  due to  $X$ , or the **mutual information** of  $Y$  and  $X$

$$IG(Y|X) = H(Y) - H(Y|X) \quad (1)$$

- If  $X$  is completely uninformative about  $Y$ :  $IG(Y|X) = 0$
- If  $X$  is completely informative about  $Y$ :  $IG(Y|X) = H(Y)$

# Revisiting Our Original Example

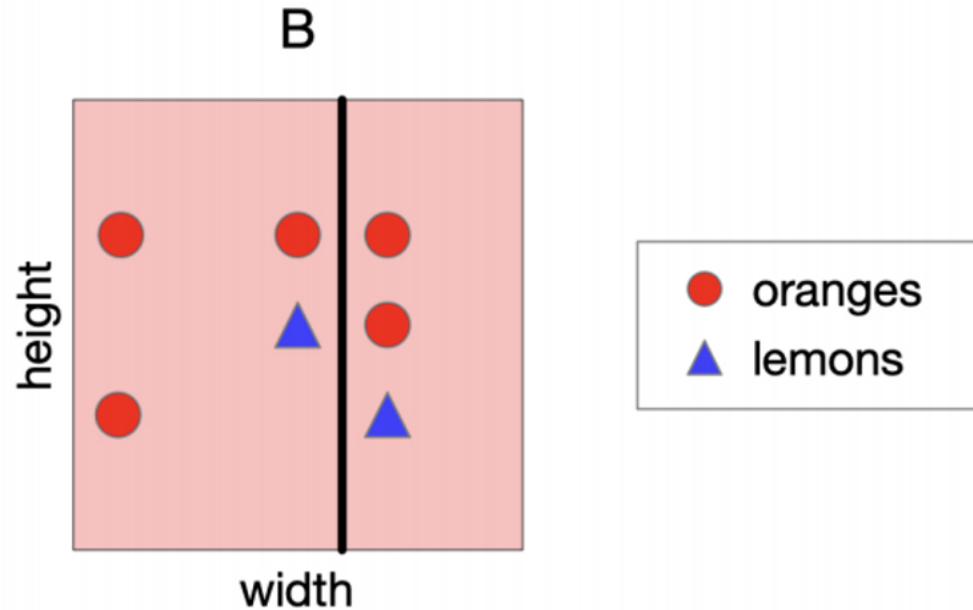
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- Information gain measures the informativeness of a variable, which is exactly what we desire in a decision tree split!
- The information gain of a split: how much information (over the training set) about the class label  $Y$  is gained by knowing which side of a split you're on.

# Revisiting Our Original Example

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- What is the information gain of split B? Not terribly informative...

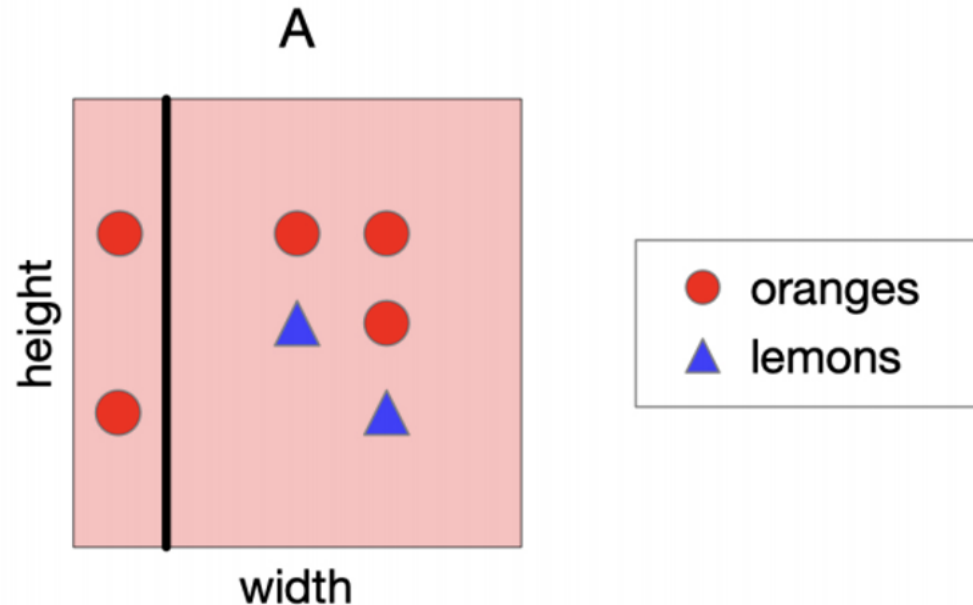


- Root entropy of class outcome:  $H(Y) = -\frac{2}{7}\log_2(\frac{2}{7}) - \frac{5}{7}\log_2(\frac{5}{7}) \approx 0.86$
- Leaf conditional entropy of class outcome:  $H(Y|left) \approx 0.81, H(Y|right) \approx 0.92$
- $IG(split) \approx 0.86 - (\frac{4}{7} \cdot 0.81 + \frac{3}{7} \cdot 0.92) \approx 0.006$

# Revisiting Our Original Example

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- What is the information gain of split A? Very informative!

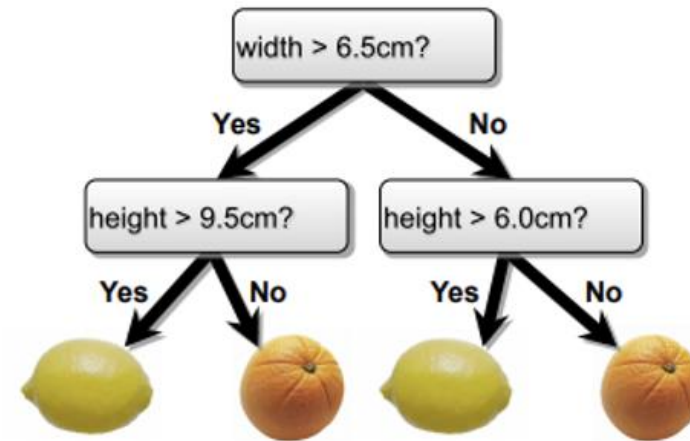
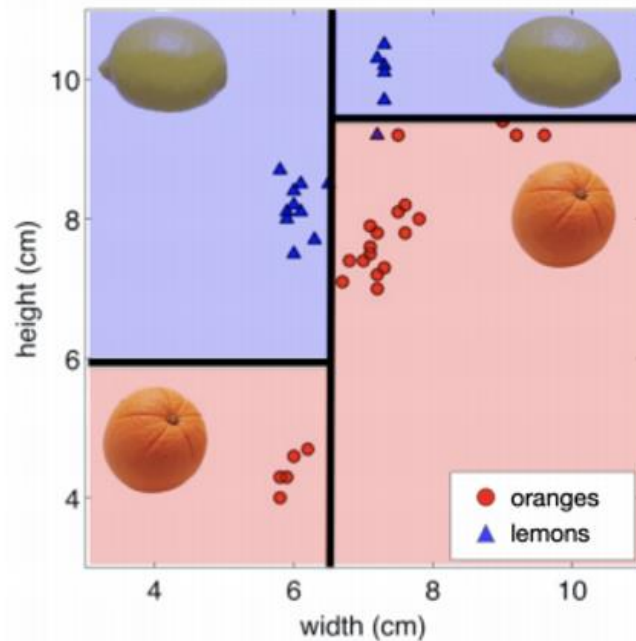


- Root entropy of class outcome:  $H(Y) = -\frac{2}{7}\log_2(\frac{2}{7}) - \frac{5}{7}\log_2(\frac{5}{7}) \approx 0.86$
- Leaf conditional entropy of class outcome:  $H(Y|left) \approx 0, H(Y|right) \approx 0.97$
- $IG(split) \approx 0.86 - (\frac{2}{7} \cdot 0 + \frac{5}{7} \cdot 0.97) \approx 0.17!$



# Constructing Decision Trees

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- At each level, one must choose:
  1. Which feature to split.
  2. Possibly where to split it.
- Choose them based on how much information we would gain from the decision! (choose feature that gives the highest gain)

# Decision Tree Construction Algorithm

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- Simple, greedy, recursive approach, builds up tree node-by-node
  1. Pick a feature to split at a non-terminal node
  2. Split examples into groups based on feature value
  3. For each group:
    - If no examples – return majority from parent
    - Else if all examples in same class – return class
    - Else loop to step 1
- Terminates when all leaves contain only examples in the same class or are empty.

# Back to Our Example

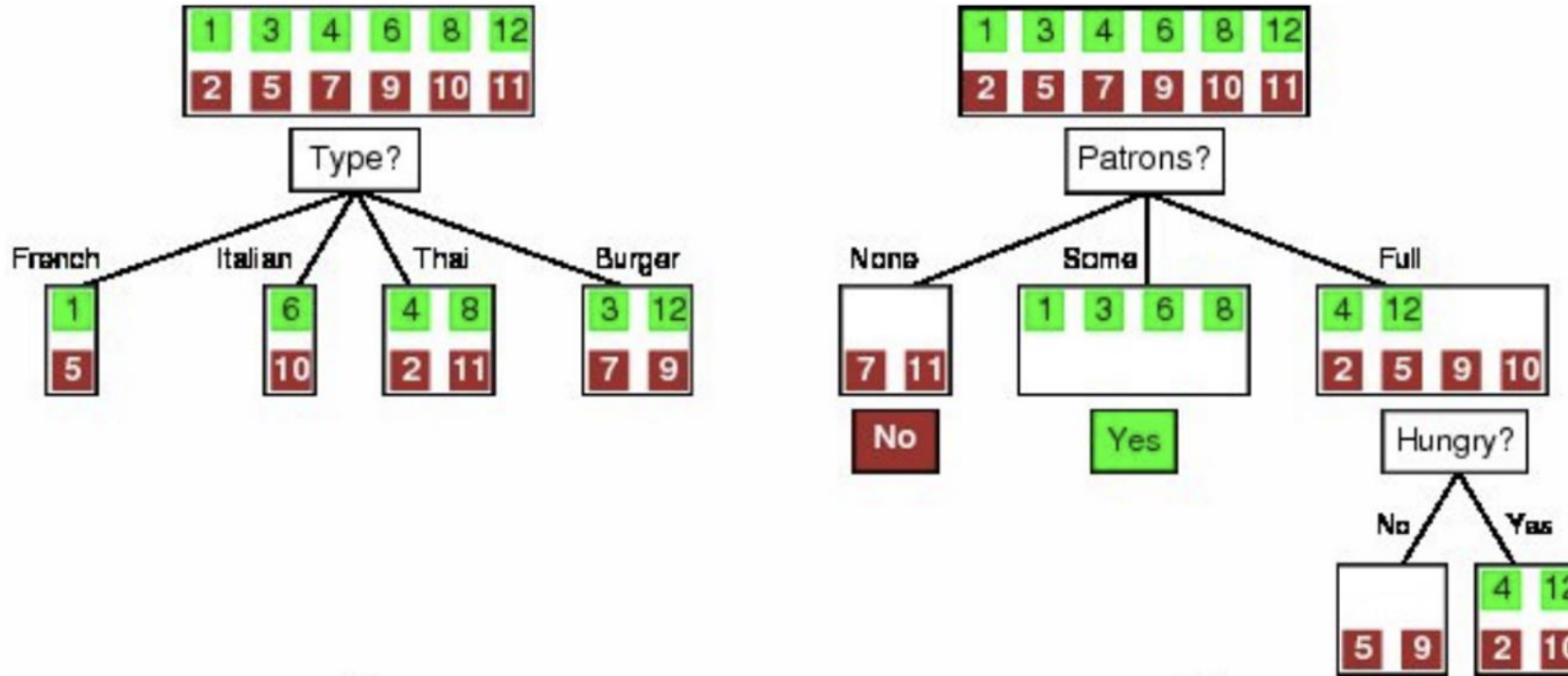
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1. Alternate: whether there is a suitable alternative restaurant nearby.
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7. Raining: whether it is raining outside.
8. Reservation: whether we made a reservation.
9. Type: the kind of restaurant (French, Italian, Thai or Burger).
10. WaitEstimate: the wait estimated by the host (0-10 minutes, 10-30, 30-60, >60).

Features:

[from: Russell & Norvig]

# Feature Selection



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

$$IG(type) = 1 - \left[ \frac{2}{12} H(Y|Fr.) + \frac{2}{12} H(Y|It.) + \frac{4}{12} H(Y|Thai) + \frac{4}{12} H(Y|Bur.) \right] = 0$$

$$IG(Patrons) = 1 - \left[ \frac{2}{12} H(0,1) + \frac{4}{12} H(1,0) + \frac{6}{12} H\left(\frac{2}{6}, \frac{4}{6}\right) \right] \approx 0.541$$

# What Makes a Good Tree?

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- Not too small: need to handle important but possibly subtle distinctions in data
- Not too big:
  - Computational efficiency (avoid redundant, spurious attributes)
  - Avoid over-fitting training examples
  - Human interpretability
- “Occam’s Razor”: find the simplest hypothesis that fits the observations
  - Useful principle, but hard to formalize (how to define simplicity?)
  - See Domingos, 1999, “The role of Occam’s razor in knowledge discovery”
- We desire small trees with informative nodes near the root

# Decision Tree Miscellany

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- Problems:
  - You have exponentially less data at lower levels
  - Too big of a tree can overfit the data
  - Greedy algorithms don't necessarily yield the global optimum
- Handling continuous attributes
  - Split based on a threshold, chosen to maximize information gain

# Comparison to some other classifiers

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## Advantages of decision trees over KNNs and neural nets

- Simple to deal with discrete features, missing values, and poorly scaled data
- Fast at test time
- More interpretable

## Advantages of KNNs over decision trees

- Few hyperparameters
- Can incorporate interesting distance measures (e.g. shape contexts)

## Advantages of neural nets over decision trees

- Able to handle attributes/features that interact in very complex ways (e.g. pixels)

# Comparison to some other classifiers

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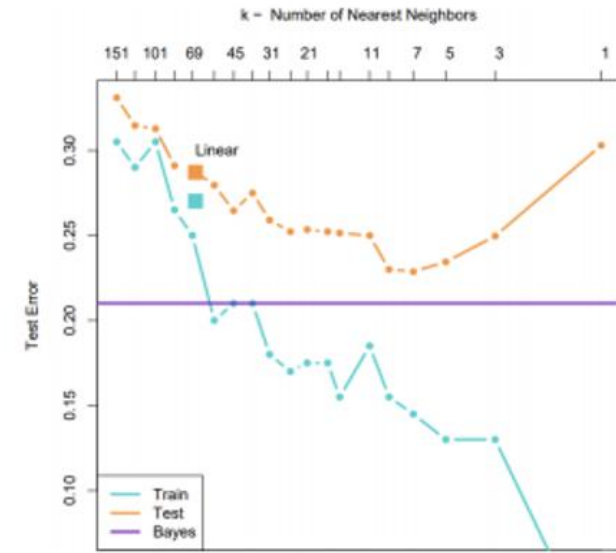
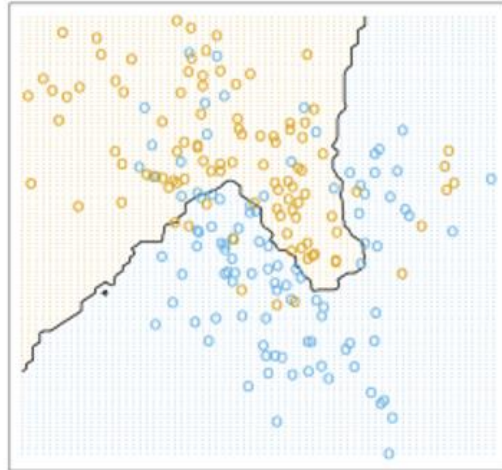
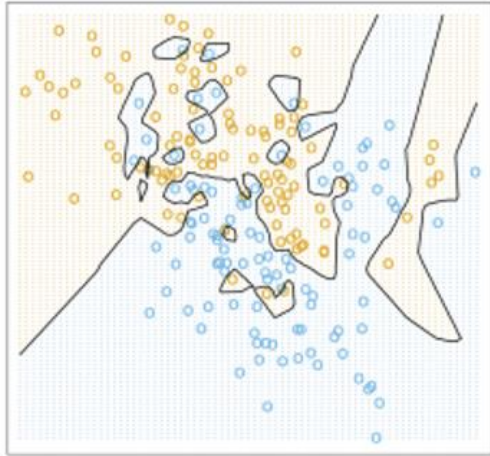
- We've seen many classification algorithms.
- We can combine multiple classifiers into an **ensemble**, which is a set of predictors whose individual decisions are combined in some way to classify new examples
  - E.g., (possibly weighted) majority vote
- For this to be nontrivial, the classifiers must differ somehow, e.g.
  - Different algorithm
  - Different choice of hyperparameters
  - Trained on different data
  - Trained with different weighting of the training examples
- Next lecture, we will study some specific ensembling techniques.



- 
- Today, we deepen our understanding of generalization through a bias-variance decomposition.
    - This will help us understand ensembling methods.

# Bias-Variance Decomposition

- Recall that overly simple models underfit the data, and overly complex models overfit.

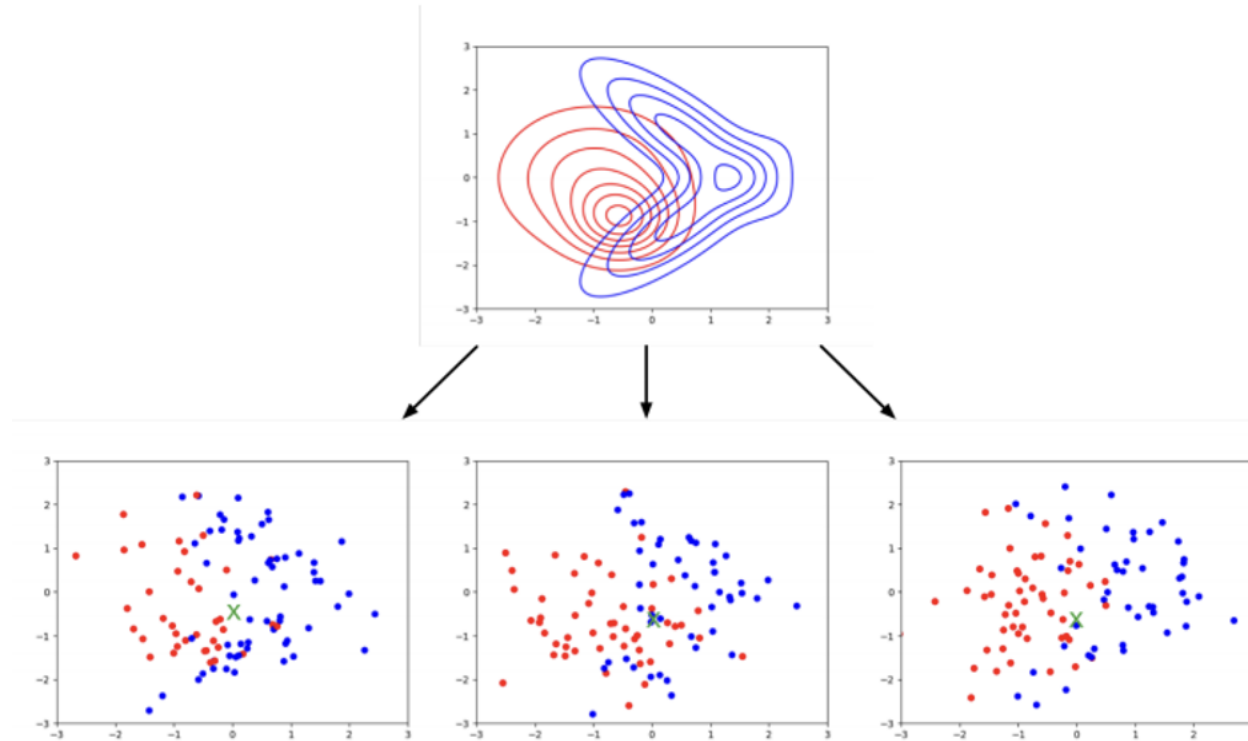


- We can quantify this effect in terms of the [bias/variance decomposition](#).
  - Bias and variance of what?

# Bias-Variance Decomposition: Basic Setup

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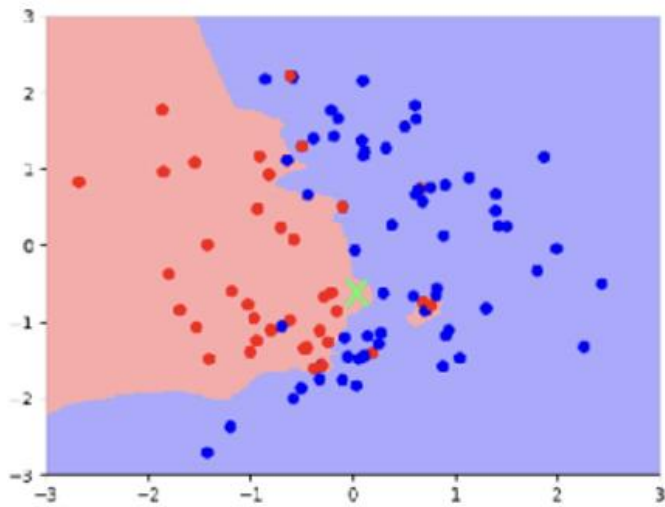
- Suppose the training set  $\mathcal{D}$  consists of pairs  $(\mathbf{x}_i, t_i)$  sampled **independent and identically distributed (i.i.d)** from a single **data generating distribution**  $p_{\text{sample}}$ .
- Pick a fixed query point  $\mathbf{x}$  (denoted with a green x).
- Consider an experiment where we sample lots of training sets independently from  $p_{\text{sample}}$ .



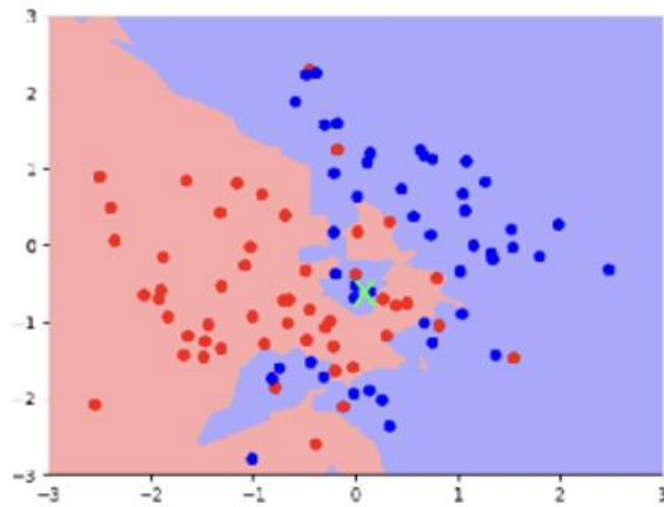
# Bias-Variance Decomposition: Basic Setup

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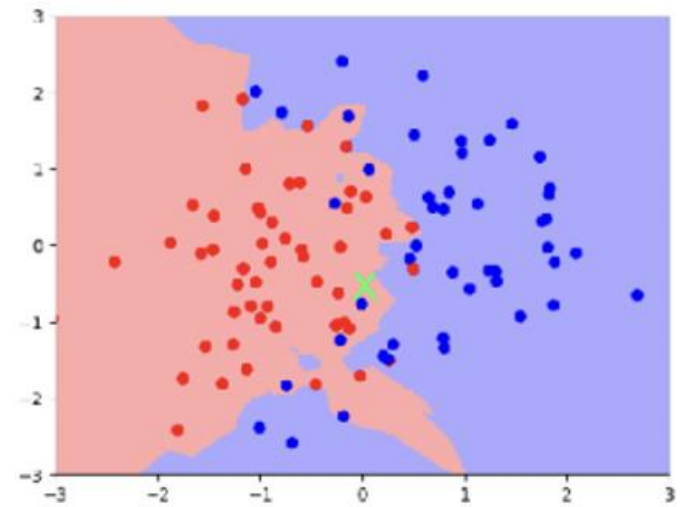
- Let's run our learning algorithm on each training set, and compute its prediction  $y$  at the query point  $\mathbf{x}$ .
- we can view  $y$  as a random variable, where the randomness comes from the choice of training set.
- The classification accuracy is determined by the distribution of  $y$ .



$y = \text{red}$



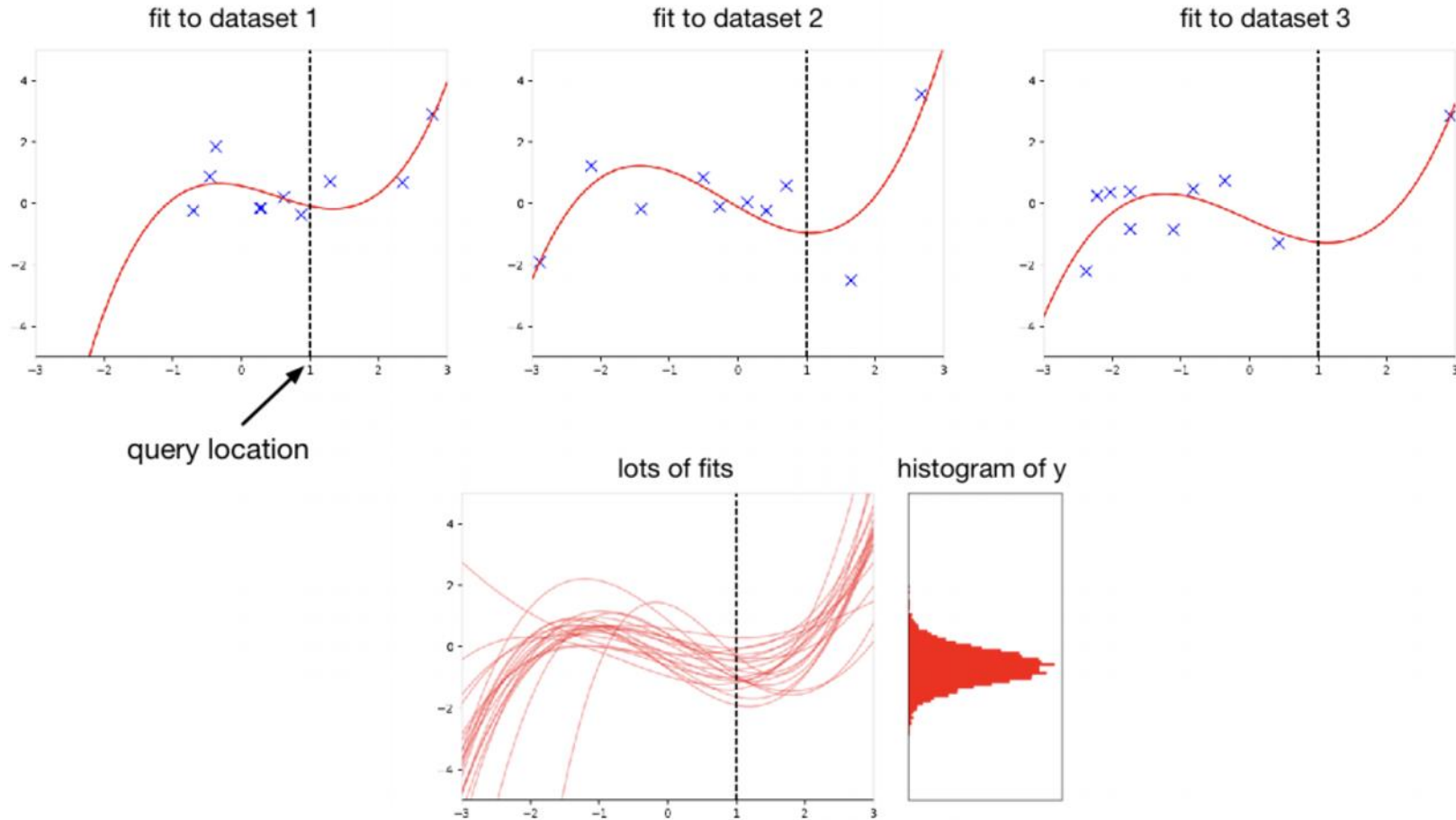
$y = \text{blue}$



$y = \text{red}$

# Bias-Variance Decomposition: Basic Setup

Here is the analogous setup for regression:



Since  $y$  is a random variable, we can talk about its expectation, variance, etc.

# Bias-Variance Decomposition: Basic Setup

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- Recap of basic setup:
  - Fix a query point  $\mathbf{x}$ .
  - Repeat:
    - Sample a random training dataset  $\mathcal{D}$  i.i.d. from the data generating distribution  $p_{\text{sample}}$ .
    - Run the learning algorithm on  $\mathcal{D}$  to get a prediction  $y$  at  $\mathbf{x}$ .
    - Sample the (true) target from the conditional distribution  $p(t|\mathbf{x})$ .
    - Compute the loss  $L(y, t)$ .
- Notice:  $y$  is independent of  $t$ .
- This gives a distribution over the loss at  $\mathbf{x}$ , with expectation  $\mathbb{E}[L(y, t)|\mathbf{x}]$ .
- For each query point  $\mathbf{x}$ , the expected loss is different. We are interested in minimizing the expectation of this with respect to  $\mathbf{x} \sim p_{\text{sample}}$ .

# Bayes Optimality

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- For now, focus on squared error loss,  $L(y, t) = \frac{1}{2} (y - t)^2$ .
- A first step: suppose we knew the conditional distribution  $p(t|\mathbf{x})$ . What value  $y$  should we predict?
  - Here, we are treating  $t$  as a random variable and choosing  $y$ .
- **Claim:**  $y_* = \mathbb{E}[t|\mathbf{x}]$  is the best possible prediction.
- **Proof:**

$$\begin{aligned}\mathbb{E}[(y - t)^2|\mathbf{x}] &= \mathbb{E}[y^2 - 2yt + t^2|\mathbf{x}] \\ &= y^2 - 2y\mathbb{E}[t|\mathbf{x}] + \mathbb{E}[t^2|\mathbf{x}] \\ &= y^2 - 2y\mathbb{E}[t|\mathbf{x}] + \mathbb{E}[t|\mathbf{x}]^2 + \text{Var}[t|\mathbf{x}] \\ &= y^2 - 2yy_* + y_*^2 + \text{Var}[t|\mathbf{x}] \\ &= (y - y_*)^2 + \text{Var}[t|\mathbf{x}]\end{aligned}$$

# Bayes Optimality

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$$\mathbb{E}[(y - t)^2 | \mathbf{x}] = (y - y_*)^2 + \text{Var}[t | \mathbf{x}]$$

- The first term is nonnegative, and can be made 0 by setting  $y = y_*$ .
- The second term corresponds to the inherent unpredictability, or **noise**, of the targets, and is called the **Bayes error**.
  - This is the best we can ever hope to do with any learning algorithm. An algorithm that achieves it is **Bayes optimal**.
  - Notice that this term doesn't depend on  $y$ .
- This process of choosing a single value  $y_*$  based on  $p(t | \mathbf{x})$  is an example of **decision theory**.



# Bayes Optimality

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- Now return to treating  $y$  as a random variable (where the randomness comes from the choice of dataset).
- We can decompose out the expected loss (suppressing the conditioning on  $\mathbf{x}$  for clarity):

$$\begin{aligned}\mathbb{E}[(y - t)^2] &= \mathbb{E}[(y - y_\star)^2] + \text{Var}(t) \\ &= \mathbb{E}[y_\star^2 - 2y_\star y + y^2] + \text{Var}(t) \\ &= y_\star^2 - 2y_\star \mathbb{E}[y] + \mathbb{E}[y^2] + \text{Var}(t) \\ &= y_\star^2 - 2y_\star \mathbb{E}[y] + \mathbb{E}[y]^2 + \text{Var}(y) + \text{Var}(t) \\ &= \underbrace{(y_\star - \mathbb{E}[y])^2}_{\text{bias}} + \underbrace{\text{Var}(y)}_{\text{variance}} + \underbrace{\text{Var}(t)}_{\text{Bayes error}}\end{aligned}$$

# Bayes Optimality

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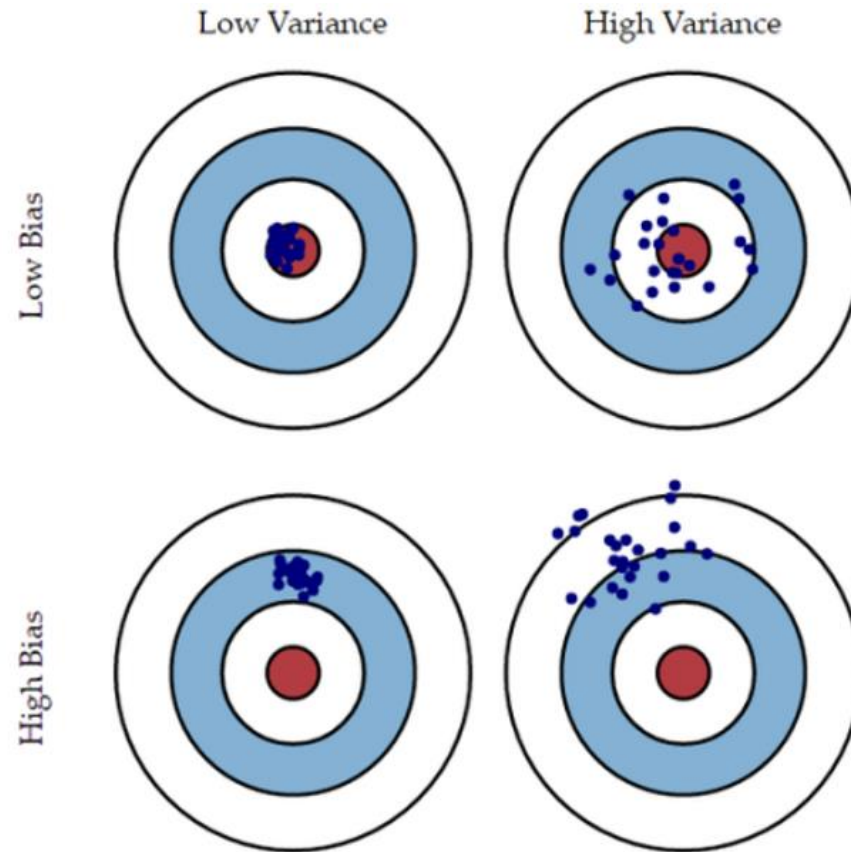
$$\mathbb{E}[(y - t)^2] = \underbrace{(y_\star - \mathbb{E}[y])^2}_{\text{bias}} + \underbrace{\text{Var}(y)}_{\text{variance}} + \underbrace{\text{Var}(t)}_{\text{Bayes error}}$$

- We just split the expected loss into three terms:
  - **bias**: how wrong the expected prediction is (corresponds to underfitting)
  - **variance**: the amount of variability in the predictions (corresponds to overfitting)
  - Bayes error: the inherent unpredictability of the targets
- Even though this analysis only applies to squared error, we often loosely use “bias” and “variance” as synonyms for “underfitting” and “overfitting”.

# Bias and Variance

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- Throwing darts = predictions for each draw of a dataset



- Be careful, what doesn't this capture?
  - We average over points  $x$  from the data distribution.