

Decision Trees

Machine Learning – CSE546
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Linear separability

- A dataset is **linearly separable** iff there exists a **separating hyperplane**:
 - Exists \mathbf{w} , such that:
 - $w_0 + \sum_i w_i x_i > 0$; if $\mathbf{x}=\{x_1, \dots, x_k\}$ is a positive example
 - $w_0 + \sum_i w_i x_i < 0$; if $\mathbf{x}=\{x_1, \dots, x_k\}$ is a negative example

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Not linearly separable data

- Some datasets are **not linearly separable!**

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Addressing non-linearly separable data – Option 1, non-linear features

- Choose non-linear features, e.g.,
 - Typical linear features: $w_0 + \sum_i w_i x_i$
 - Example of non-linear features:
 - Degree 2 polynomials, $w_0 + \sum_i w_i x_i + \sum_{ij} w_{ij} x_i x_j$
- Classifier $h_{\mathbf{w}}(\mathbf{x})$ still linear in parameters \mathbf{w}
 - As easy to learn
 - Data is linearly separable in higher dimensional spaces
 - More discussion later this quarter

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Addressing non-linearly separable data – Option 2, non-linear classifier

- Choose a classifier $h_{\mathbf{w}}(\mathbf{x})$ that is non-linear in parameters \mathbf{w} , e.g.,
 - Decision trees, boosting, nearest neighbor, neural networks...
- More general than linear classifiers
- But, can often be harder to learn (non-convex/concave optimization required)
- But, but, often very useful
- (BTW. Later this quarter, we'll see that these options are not that different)

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A small dataset: Miles Per Gallon

Suppose we want to predict MPG

mpg	cylinders	displacement	horsepower	weight	acceleration	modelyear	maker
good	4	low	low	low	high	75to78	asia
bad	6	medium	medium	medium	medium	70to74	america
bad	4	medium	medium	medium	low	75to78	europa
bad	8	high	high	high	low	70to74	america
bad	6	medium	medium	medium	medium	70to74	america
bad	4	low	medium	low	medium	70to74	asia
bad	4	low	medium	low	low	70to74	asia
bad	8	high	high	high	low	75to78	america
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.
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bad	8	high	high	high	low	70to74	america
good	8	high	medium	high	high	79to83	america
bad	8	high	high	high	low	75to78	america
good	4	low	low	low	low	79to83	america
bad	6	medium	medium	medium	high	75to78	america
good	4	medium	low	low	low	79to83	america
good	4	low	low	medium	high	79to83	america
bad	8	high	high	high	low	70to74	america
good	4	low	medium	low	medium	75to78	europa
bad	5	medium	medium	medium	medium	75to78	europa

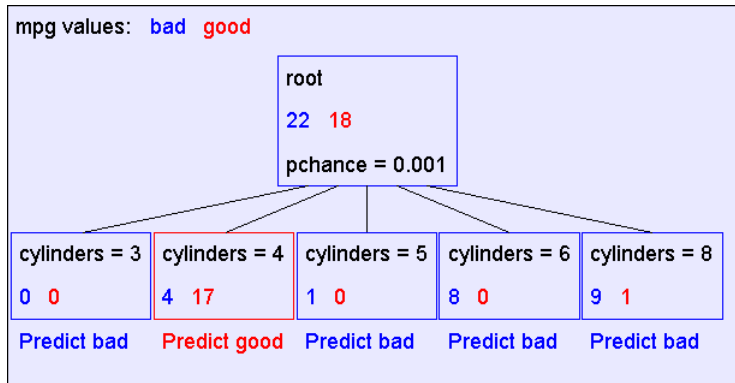
40 training examples

From the UCI repository (thanks to Ross Quinlan)

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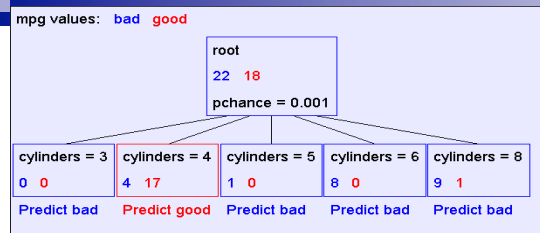
A Decision Stump



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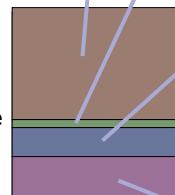
Recursion Step



Take the Original Dataset..



And partition it according to the value of the attribute we split on



Examples in which cylinders = 4

Examples in which cylinders = 5

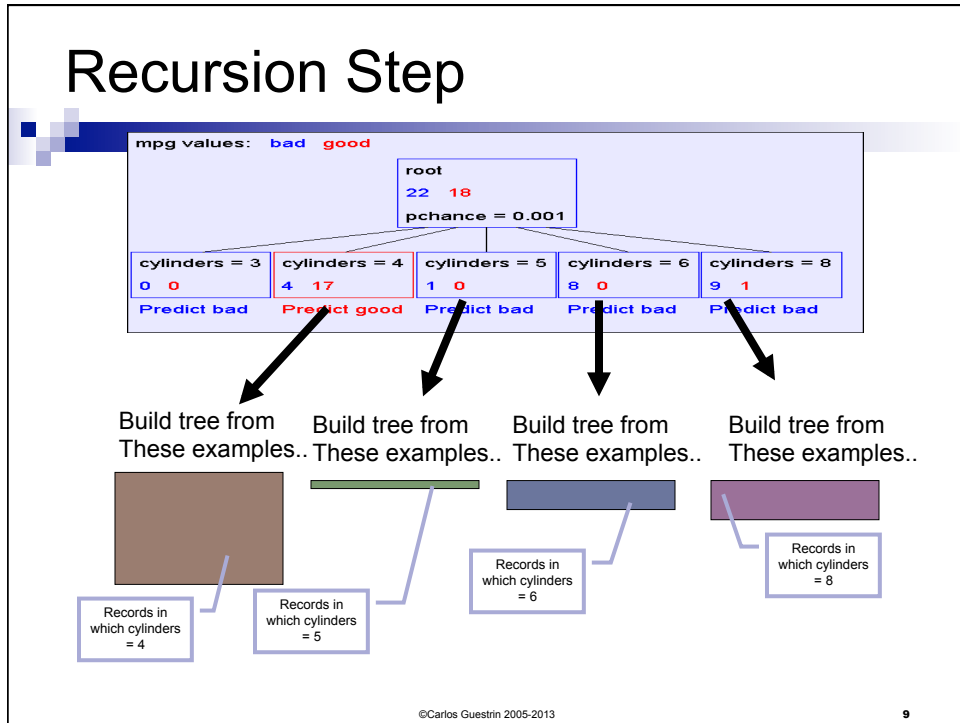
Examples in which cylinders = 6

Examples in which cylinders = 8

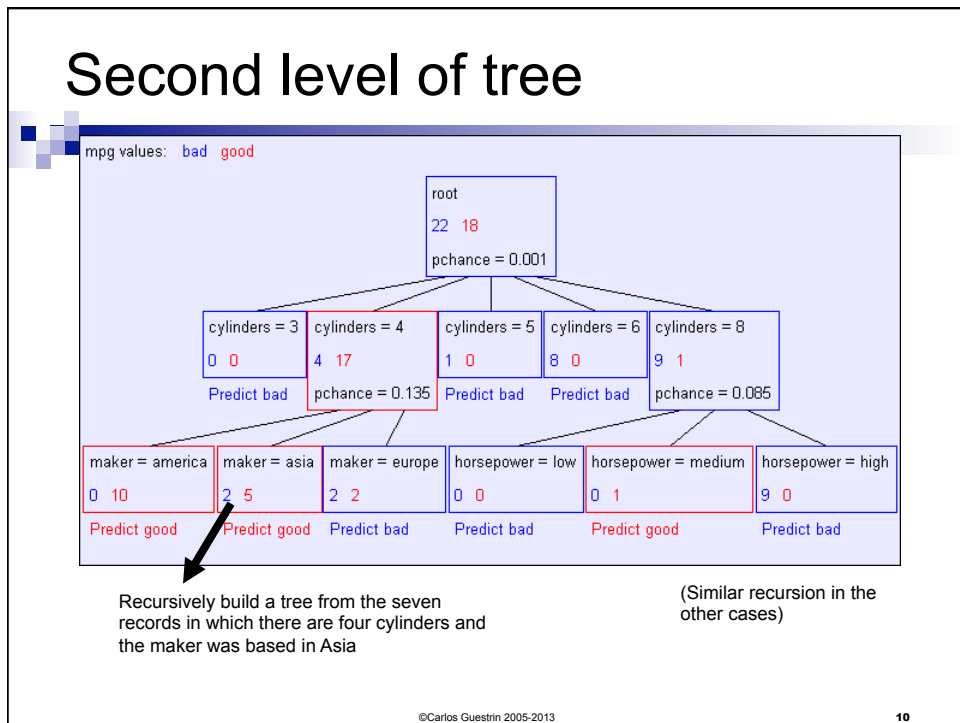
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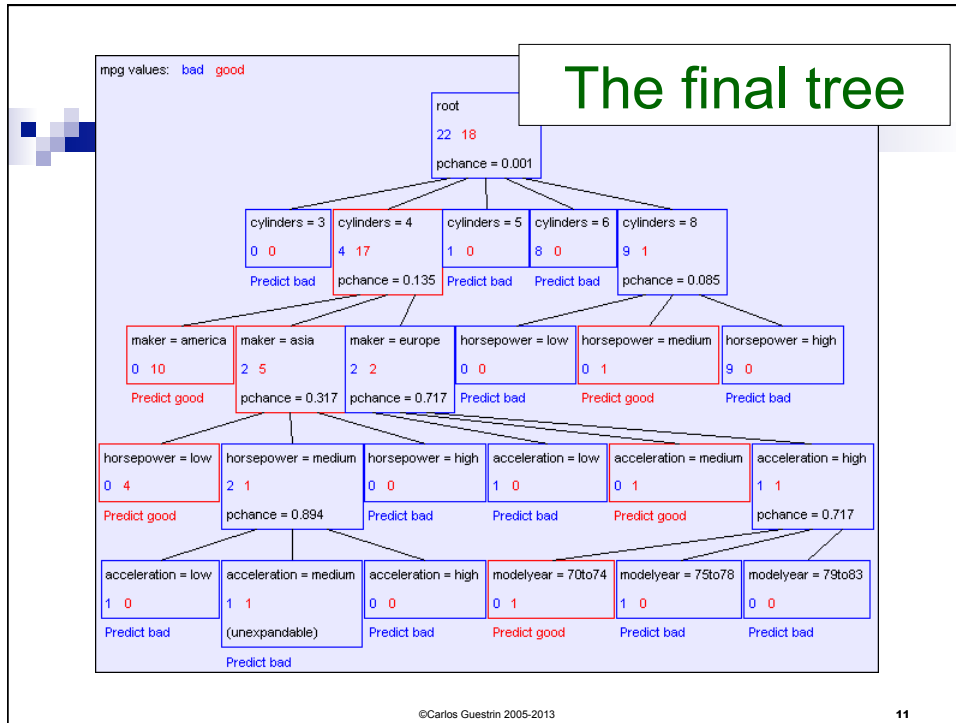
8

Recursion Step



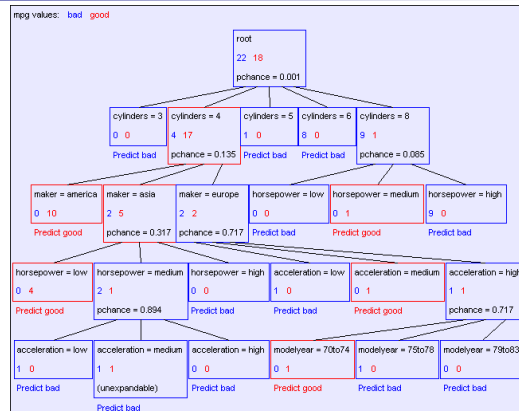
Second level of tree





Classification of a new example

- Classifying a test example – traverse tree and report leaf label



Are all decision trees equal?

- Many trees can represent the same concept
- But, not all trees will have the same size!
 - e.g., $\phi = A \wedge B \vee \neg A \wedge C$ ((A and B) or (not A and C))

Learning decision trees is hard!!!

- Learning the simplest (smallest) decision tree is an NP-complete problem [Hyafil & Rivest '76]
- Resort to a greedy heuristic:
 - Start from empty decision tree
 - Split on **next best attribute (feature)**
 - Recurse

Choosing a good attribute

X_1	X_2	Y
T	T	T
T	F	T
T	T	T
T	F	T
F	T	T
F	F	F
F	T	F
F	F	F

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Measuring uncertainty

- Good split if we are more certain about classification after split
 - Deterministic good (all true or all false)
 - Uniform distribution bad

$P(Y=A) = 1/2$	$P(Y=B) = 1/4$	$P(Y=C) = 1/8$	$P(Y=D) = 1/8$
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$P(Y=A) = 1/4$	$P(Y=B) = 1/4$	$P(Y=C) = 1/4$	$P(Y=D) = 1/4$
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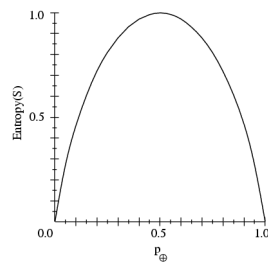
Entropy

Entropy $H(Y)$ of a random variable Y

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

More uncertainty, more entropy!

Information Theory interpretation: $H(Y)$ is the expected number of bits needed to encode a randomly drawn value of Y (under most efficient code)



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Information gain

■ Advantage of attribute – decrease in uncertainty

- Entropy of Y before you split
- Entropy after split
 - Weight by probability of following each branch, i.e., normalized number of records

$$H(Y | X) = - \sum_{j=1}^v P(X = x_j) \sum_{i=1}^k P(Y = y_i | X = x_j) \log_2 P(Y = y_i | X = x_j)$$

X_1	X_2	Y
T	T	T
T	F	T
T	T	T
T	F	T
F	T	T
F	F	F

■ Information gain is difference $IG(X) = H(Y) - H(Y | X)$

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Learning decision trees

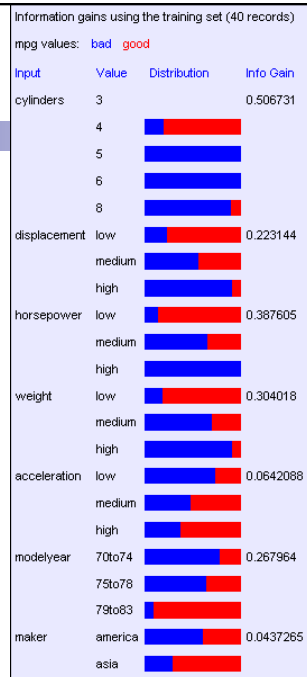
- Start from empty decision tree
- Split on **next best attribute (feature)**
 - Use, for example, information gain to select attribute
 - Split on $\arg \max_i IG(X_i) = \arg \max_i H(Y) - H(Y | X_i)$
- Recurse

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Suppose we want
to predict MPG

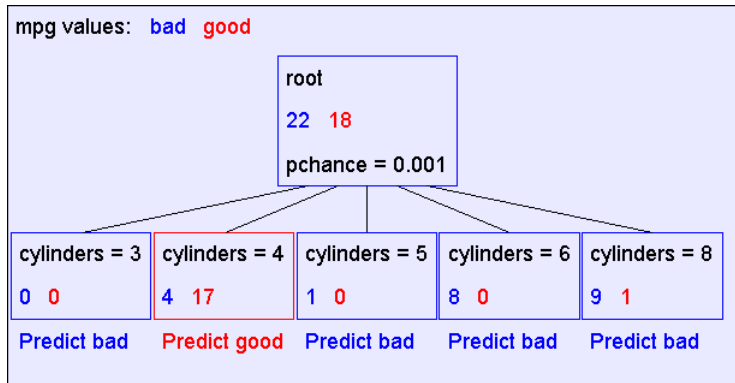
Look at all the
information
gains...



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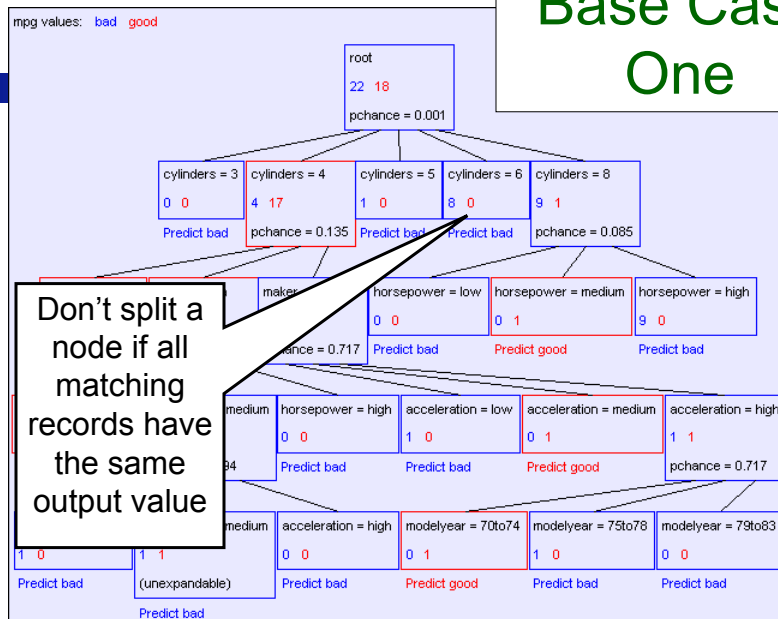
A Decision Stump



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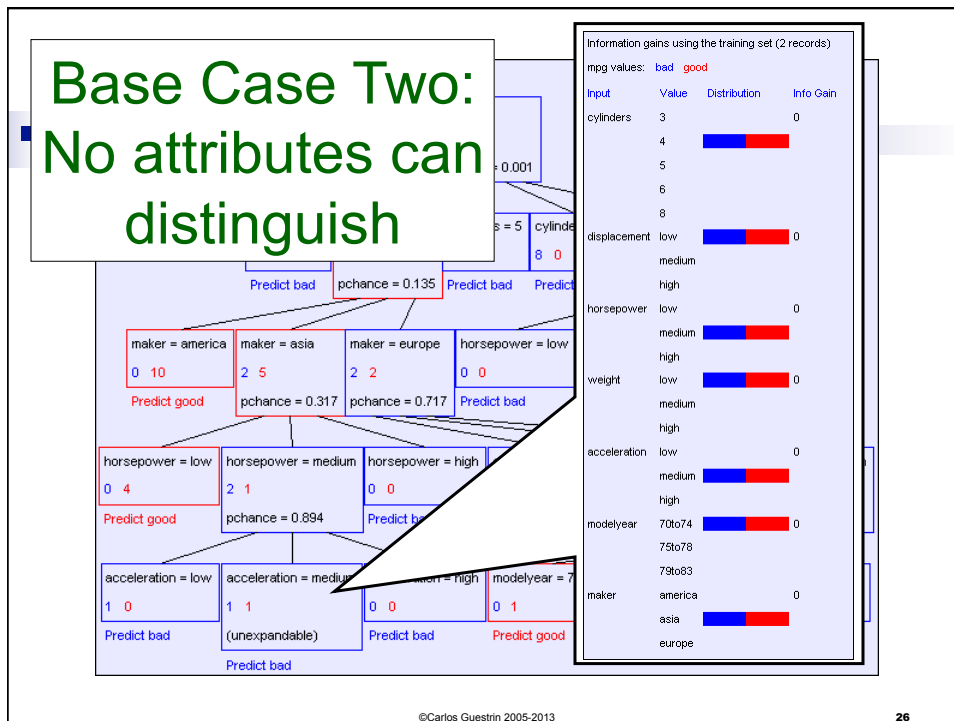
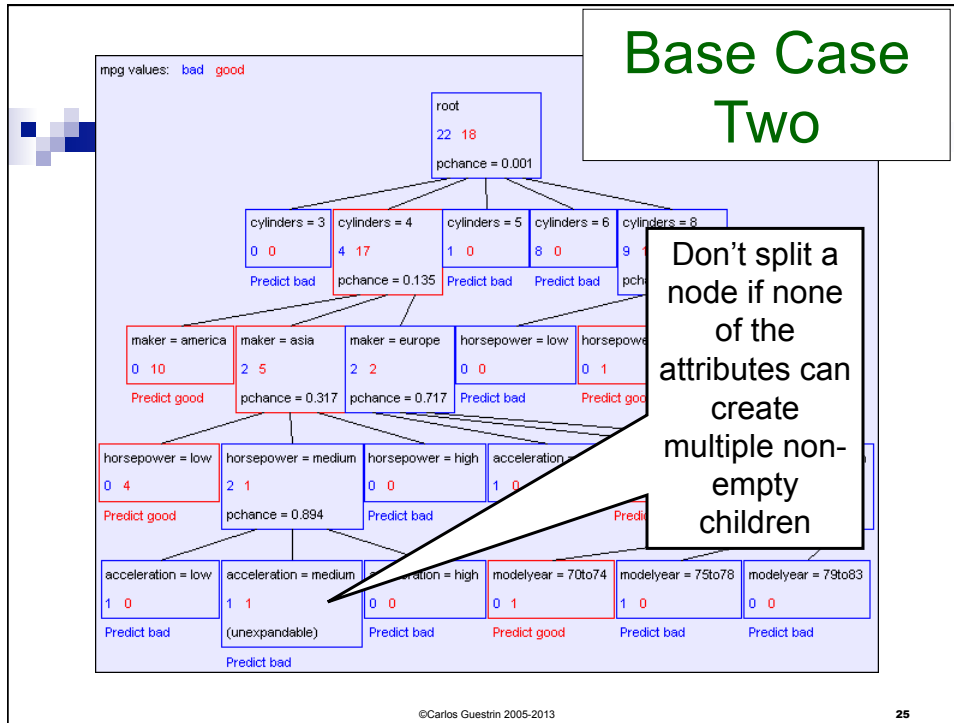
Base Case One



Don't split a node if all matching records have the same output value

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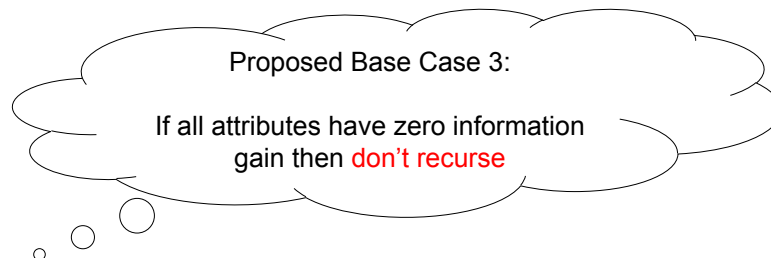


Base Cases

- Base Case One: If all records in current data subset have the same output then **don't recurse**
- Base Case Two: If all records have exactly the same set of input attributes then **don't recurse**

Base Cases: An idea

- Base Case One: If all records in current data subset have the same output then **don't recurse**
- Base Case Two: If all records have exactly the same set of input attributes then **don't recurse**



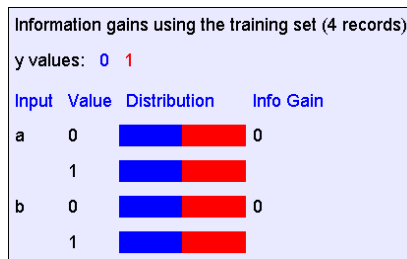
•*Is this a good idea?*

The problem with Base Case 3

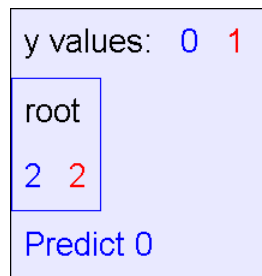
a	b	y
0	0	0
0	1	1
1	0	1
1	1	0

$Y = A \text{ XOR } B$

The information gains:



The resulting bad decision tree:

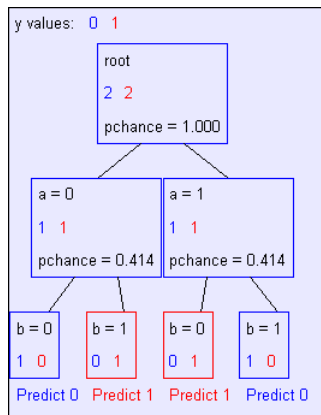


If we omit Base Case 3:

a	b	y
0	0	0
0	1	1
1	0	1
1	1	0

$y = a \text{ XOR } b$

The resulting decision tree:



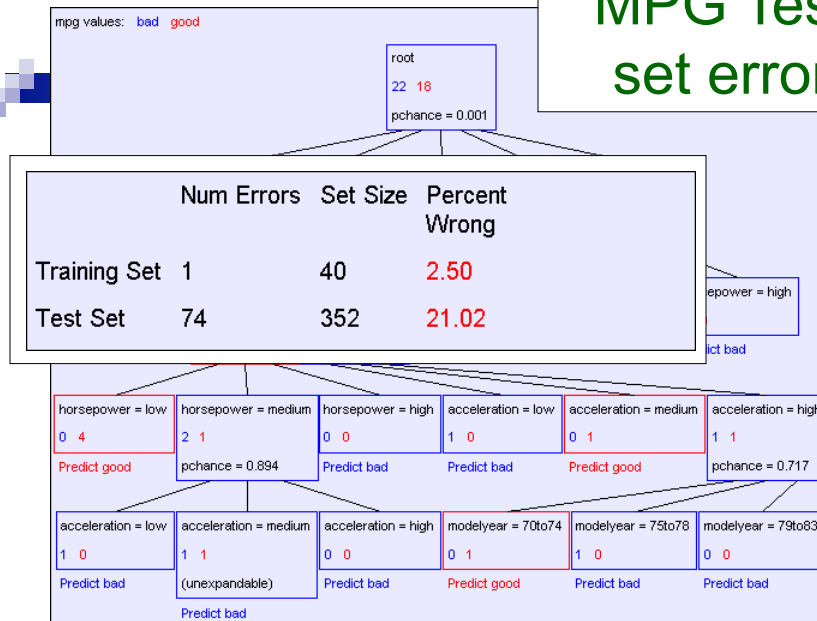
Basic Decision Tree Building Summarized

BuildTree(DataSet, Output)

- If all output values are the same in DataSet, return a leaf node that says “predict this unique output”
- If all input values are the same, return a leaf node that says “predict the majority output”
- Else find attribute X with highest Info Gain
- Suppose X has n_x distinct values (i.e. X has arity n_x).
 - Create and return a non-leaf node with n_x children.
 - The i 'th child should be built by calling
BuildTree(DS_i , Output)

Where DS_i built consists of all those records in DataSet for which $X = i$ th distinct value of X.

MPG Test set error



MPG Test set error

mpg values: bad good

root
22 18
pchance = 0.001

	Num Errors	Set Size	Percent Wrong
Training Set	1	40	2.50
Test Set	74	352	21.02

The test set error is much worse than the training set error...
...why?

horsepower = low horsepower = medium horsepower = high acceleration = low acceleration = medium acceleration = high

Predict bad (unexpandable) Predict bad Predict good Predict bad Predict bad

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Decision trees & Learning Bias

mpg	cylinders	displacement	horsepower	weight	acceleration	modelyear/maker
good	4	low	low	low	high	75to78 asia
bad	6	medium	medium	medium	medium	70to74 america
bad	4	medium	medium	medium	low	75to78 europe
bad	8	high	high	high	low	70to74 america
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bad	4	low	medium	low	medium	70to74 asia
bad	4	low	medium	low	low	70to74 asia
bad	8	high	high	high	low	75to78 america
:	:	:	:	:	:	:
:	:	:	:	:	:	:
:	:	:	:	:	:	:
:	:	:	:	:	:	:
bad	8	high	high	high	low	70to74 america
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bad	8	high	high	high	low	70to74 america
good	4	low	medium	low	medium	75to78 america
bad	5	medium	medium	medium	medium	75to78 europe

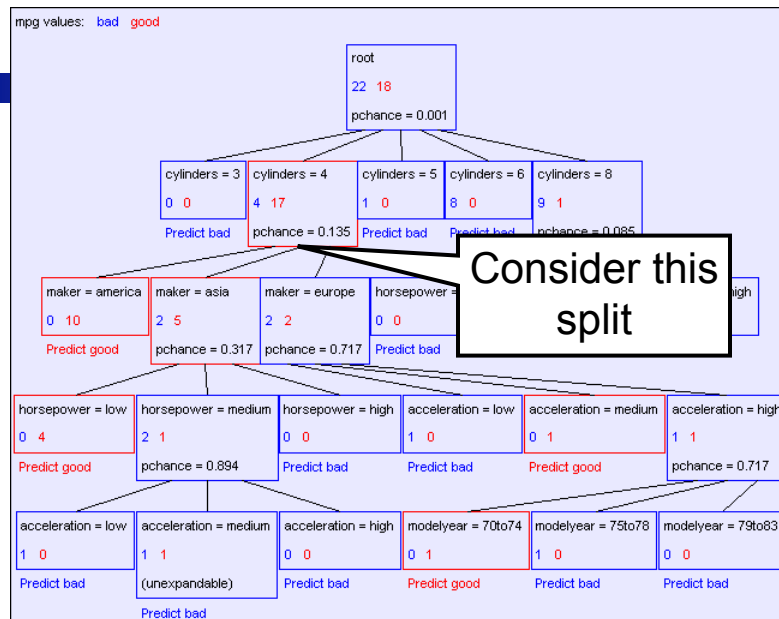
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Decision trees will overfit

- Standard decision trees have no learning bias
 - Training set error is always zero!
 - (If there is no label noise)
 - Lots of variance
 - Will definitely overfit!!!
 - Must bias towards simpler trees
- Many strategies for picking simpler trees:
 - Fixed depth
 - Fixed number of leaves
 - Or something smarter...

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A chi-square test

mpg values: bad good

maker	bad	good	H(mpg maker)
america	0	10	0
asia	2	5	0.863121
europa	2	2	1

$H(\text{mpg}) = 0.702467$ $H(\text{mpg}|\text{maker}) = 0.478183$
 $IG(\text{mpg}|\text{maker}) = 0.224284$

- Suppose that MPG was completely uncorrelated with maker.
- What is the chance we'd have seen data of at least this apparent level of association anyway?

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A chi-square test

mpg values: bad good

maker	bad	good	H(mpg maker)
america	0	10	0
asia	2	5	0.863121
europa	2	2	1

$H(\text{mpg}) = 0.702467$ $H(\text{mpg}|\text{maker}) = 0.478183$
 $IG(\text{mpg}|\text{maker}) = 0.224284$

- Suppose that mpg was completely uncorrelated with maker.
- What is the chance we'd have seen data of at least this apparent level of association anyway?

By using a particular kind of chi-square test, the answer is 7.2%

(Such simple hypothesis tests are very easy to compute, unfortunately, not enough time to cover in the lecture, but see readings...)

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Using Chi-squared to avoid overfitting

- Build the full decision tree as before
- But when you can grow it no more, start to prune:
 - Beginning at the bottom of the tree, delete splits in which $p_{chance} > MaxPchance$
 - Continue working your way up until there are no more prunable nodes

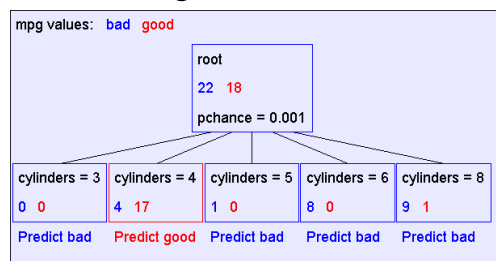
MaxPchance is a magic parameter you must specify to the decision tree, indicating your willingness to risk fitting noise

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Pruning example

- With *MaxPchance* = 0.1, you will see the following MPG decision tree:



Note the improved test set accuracy compared with the unpruned tree

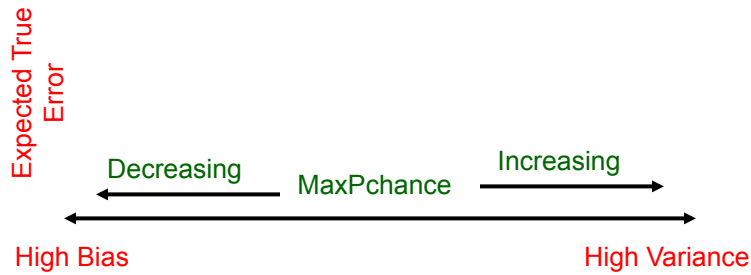
	Num Errors	Set Size	Percent Wrong
Training Set	5	40	12.50
Test Set	56	352	15.91

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MaxPchance

- Technical note MaxPchance is a regularization parameter that helps us bias towards simpler models



Real-Valued inputs

- What should we do if some of the inputs are real-valued?

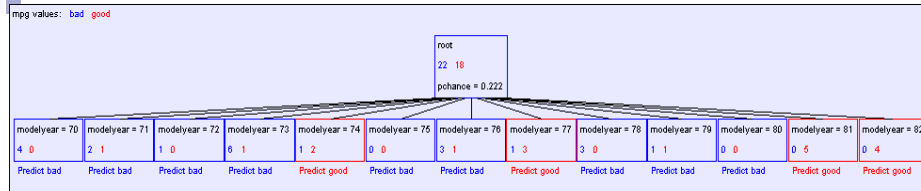
mpg	cylinders	displacemen	horsepower	weight	acceleration	modelyear	maker
good	4	97	75	2265	18.2	77	asia
bad	6	199	90	2648	15	70	america
bad	4	121	110	2600	12.8	77	europa
bad	8	350	175	4100	13	73	america
bad	6	198	95	3102	16.5	74	america
bad	4	108	94	2379	16.5	73	asia
bad	4	113	95	2228	14	71	asia
bad	8	302	139	3570	12.8	78	america
:	:	:	:	:	:	:	:
:	:	:	:	:	:	:	:
:	:	:	:	:	:	:	:
:	:	:	:	:	:	:	:
good	4	120	79	2625	18.6	82	america
bad	8	455	225	4425	10	70	america
good	4	107	86	2464	15.5	76	europa
bad	5	131	103	2830	15.9	78	europa

Infinite number of possible split values!!!

Finite dataset, only finite number of relevant splits!

Idea One: Branch on each possible real value

“One branch for each numeric value” idea:



Hopeless: with such high branching factor will shatter the dataset and overfit

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Threshold splits

- Binary tree, split on attribute X_i
 - One branch: $X_i < t$
 - Other branch: $X_i \geq t$

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Choosing threshold split

- Binary tree, split on attribute X_i
 - One branch: $X_i < t$
 - Other branch: $X_i \geq t$
- Search through possible values of t
 - Seems hard!!!
- But only finite number of t 's are important
 - Sort data according to X_i into $\{x_1, \dots, x_m\}$
 - Consider split points of the form $x_a + (x_{a+1} - x_a)/2$

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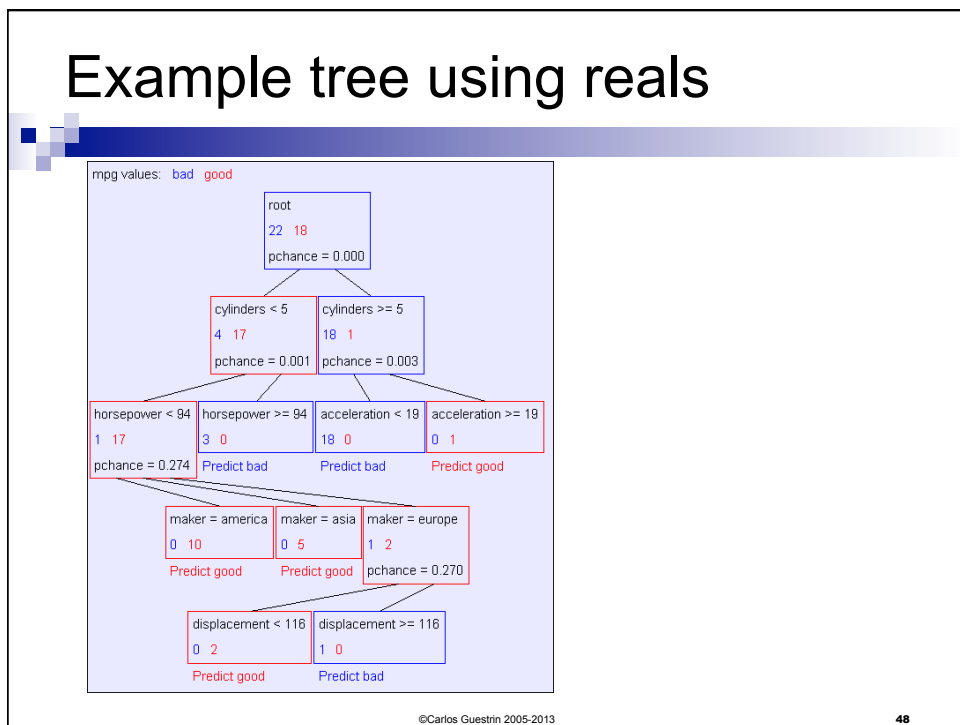
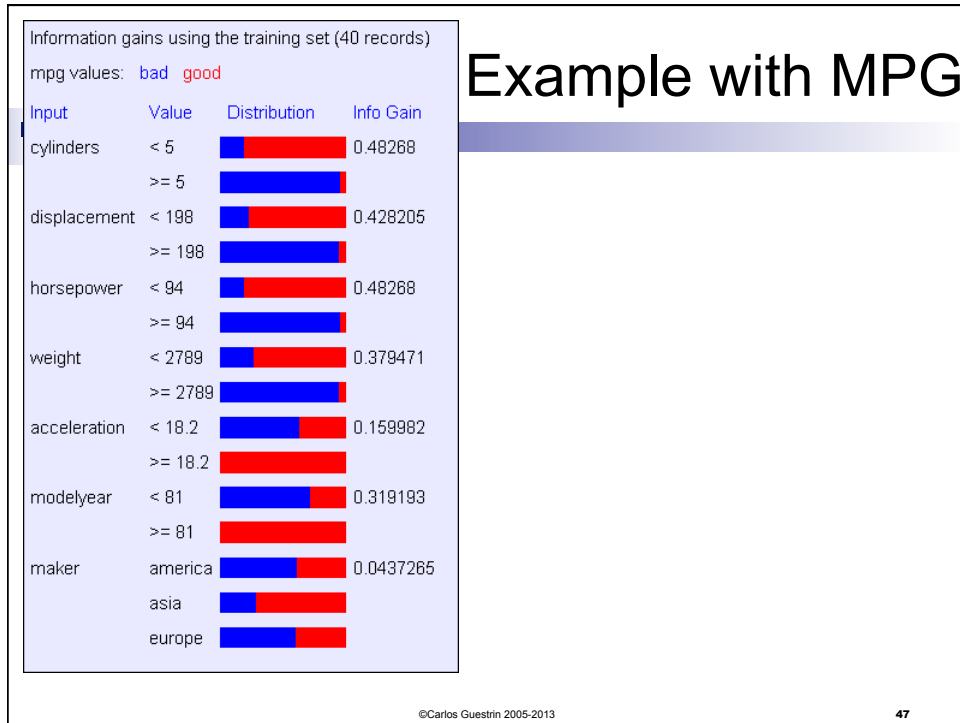
A better idea: thresholded splits

- Suppose X_i is real valued
- Define $IG(Y|X_i:t)$ as $H(Y) - H(Y|X_i:t)$
- Define $H(Y|X_i:t) =$

$$H(Y|X_i < t) P(X_i < t) + H(Y|X_i \geq t) P(X_i \geq t)$$
 - $IG(Y|X_i:t)$ is the information gain for predicting Y if all you know is whether X_i is greater than or less than t
- Then define $IG^*(Y|X_i) = \max_t IG(Y|X_i:t)$
- For each real-valued attribute, use $IG^*(Y|X_i)$ for assessing its suitability as a split
- Note, may split on an attribute multiple times, with different thresholds

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What you need to know about decision trees

- Decision trees are one of the most popular data mining tools
 - Easy to understand
 - Easy to implement
 - Easy to use
 - Computationally cheap (to solve heuristically)
- Information gain to select attributes (ID3, C4.5,...)
- Presented for classification, can be used for regression and density estimation too
- Decision trees will overfit!!!
 - Zero bias classifier ! Lots of variance
 - Must use tricks to find “simple trees”, e.g.,
 - Fixed depth/Early stopping
 - Pruning
 - Hypothesis testing

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Acknowledgements

- Some of the material in the decision trees presentation is courtesy of Andrew Moore, from his excellent collection of ML tutorials:
 - <http://www.cs.cmu.edu/~awm/tutorials>

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