# Decision trees[1]

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## 1 Supervised learning

| 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 |
| 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 |
| : | : | : | : | : | : | : | : |
| : | : | : | : | : | : | : | : |
| : | : | : | : | : | : | : | : |
| 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 | europe |
| bad | 5 | medium | medium | medium | medium | 75to78 | europe |

40 Records

Figure 1: Toy set of records (UCI)

[1] slides thanks to Carlos Guestrin@CMU

## 2 Univariate trees for classification

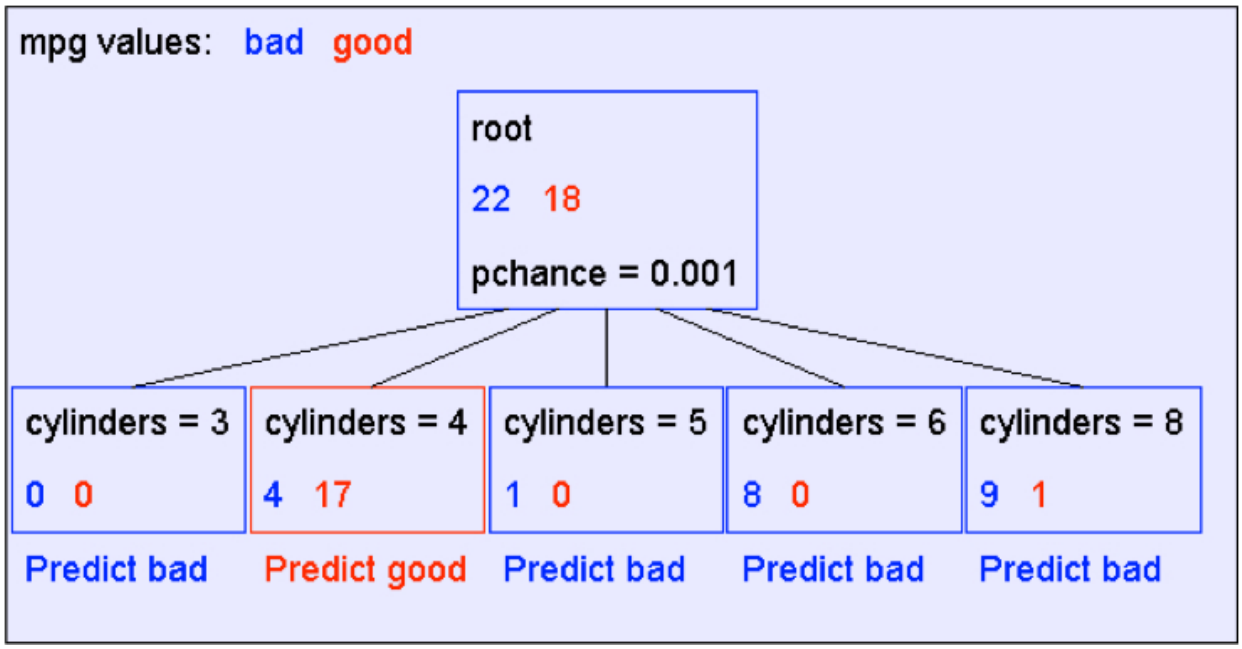


Figure 2: Decision tree, 1 layer.

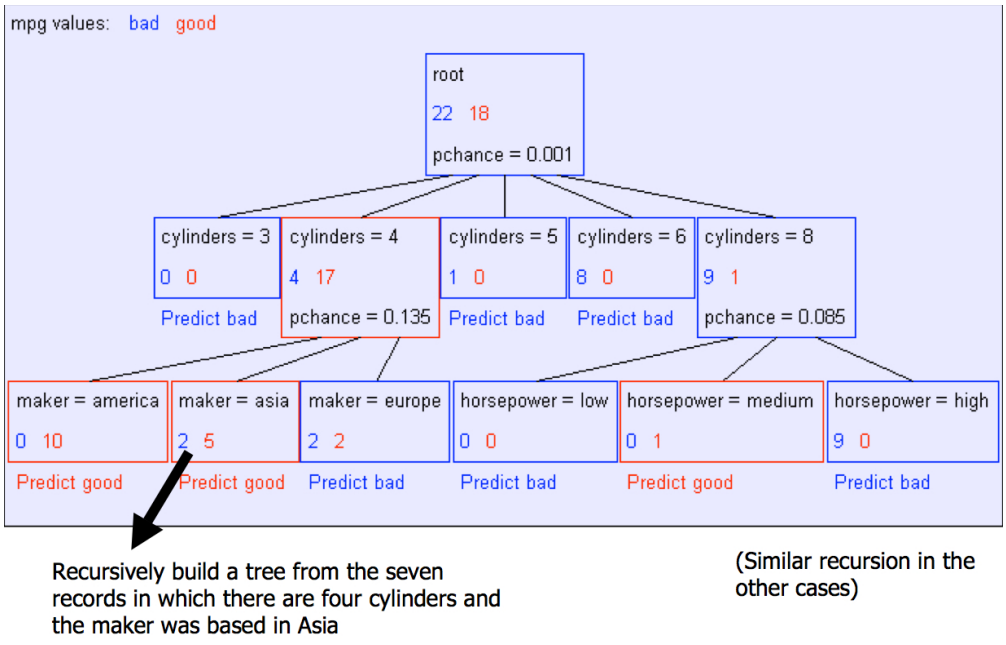


Figure 3: Decision tree, 2 layers.

REMARK: not capable of classifying data not seen in training

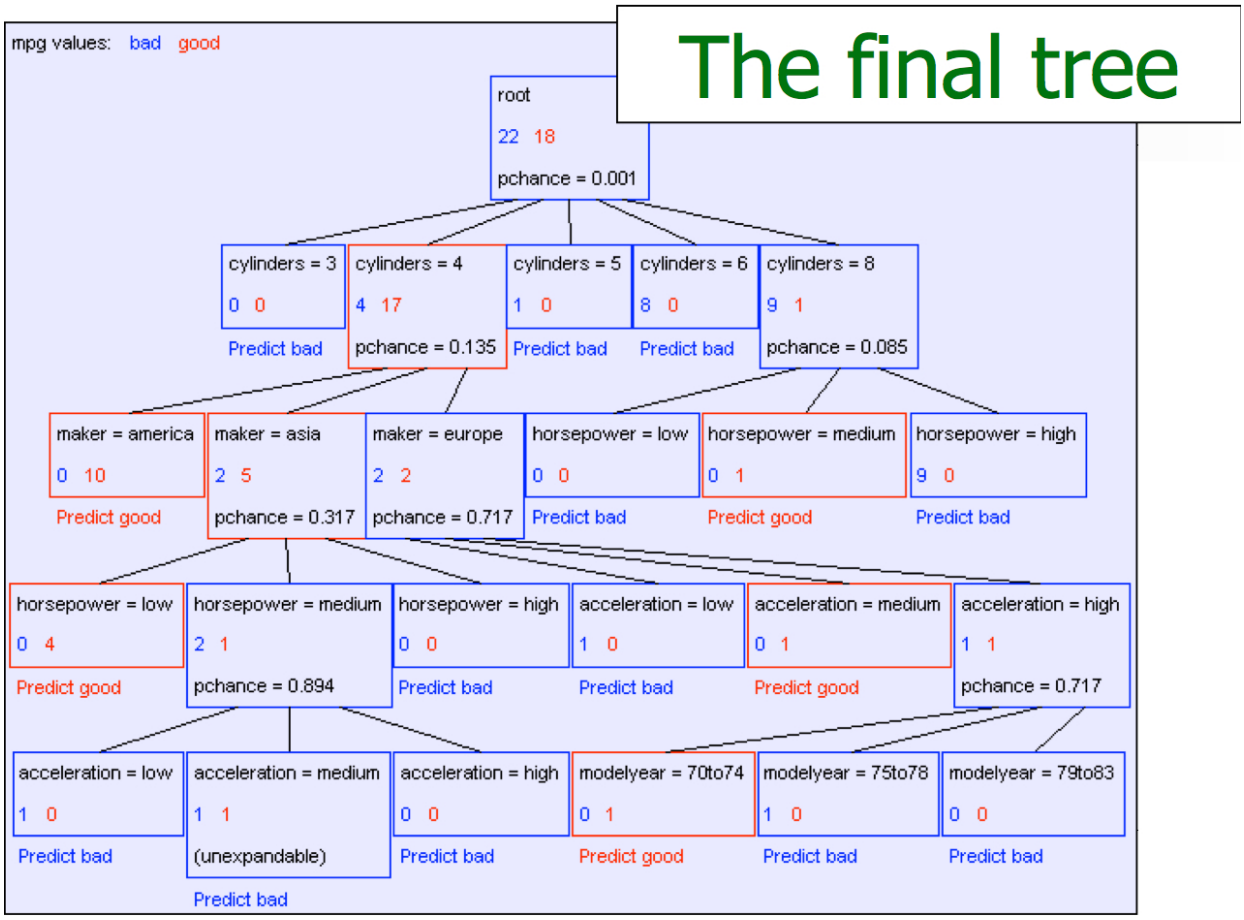


Figure 4: Decision tree, complete

# 3 Tree splitting

Finding the smallest decision tree in NP-complete. Use a heuristic:

- start with an empty decision tree
- split on the best feature.
- recurse

## 3.1 Entropy-based gain

$$H(Y) = \sum_j P(y_j) \log_2(\frac{1}{P(y_j)})$$

**Entropy after split by $X$ feature**

$$H(Y|X) = \sum_i P(x_i) \sum_j P(y_j|x_i) \log_2(\frac{1}{P(y_j|x_i)})$$

**Mutual information (or Information Gain).**

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

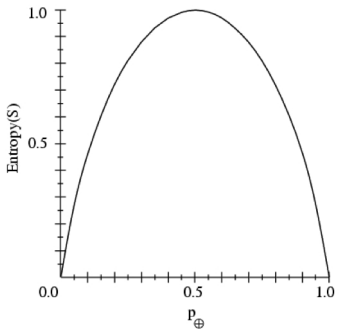


Figure 5: Entropy for 2-valued distribution

At each split we are going to choose the feature that gives the highest information gain.

| $X^1$ | $X^2$ | 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 |

Figure 6: 2 possible features to split by

$$H(Y|X^1) = \frac{1}{2}H(Y|X^1 = T) + \frac{1}{2}H(Y|X^1 = F) = 0 + \frac{1}{2}(\frac{1}{4}\log_2\frac{1}{4} + \frac{3}{4}\log_2\frac{3}{4}) \approx .405$$
$$IG(X^1) = H(Y) - H(Y|X^1) = .954 - .405 = .549$$

$$H(Y|X^2) = \frac{1}{2}H(Y|X^2 = T) + \frac{1}{2}H(Y|X^2 = F) = \frac{1}{2}(\frac{1}{4}\log_2\frac{1}{4} + \frac{3}{4}\log_2\frac{3}{4}) + \frac{1}{2}(\frac{1}{2}\log_2\frac{1}{2} + \frac{1}{2}\log_2\frac{1}{2}) \approx .905$$
$$IG(X^2) = H(Y) - H(Y|X^2) = .954 - .905 = .049$$

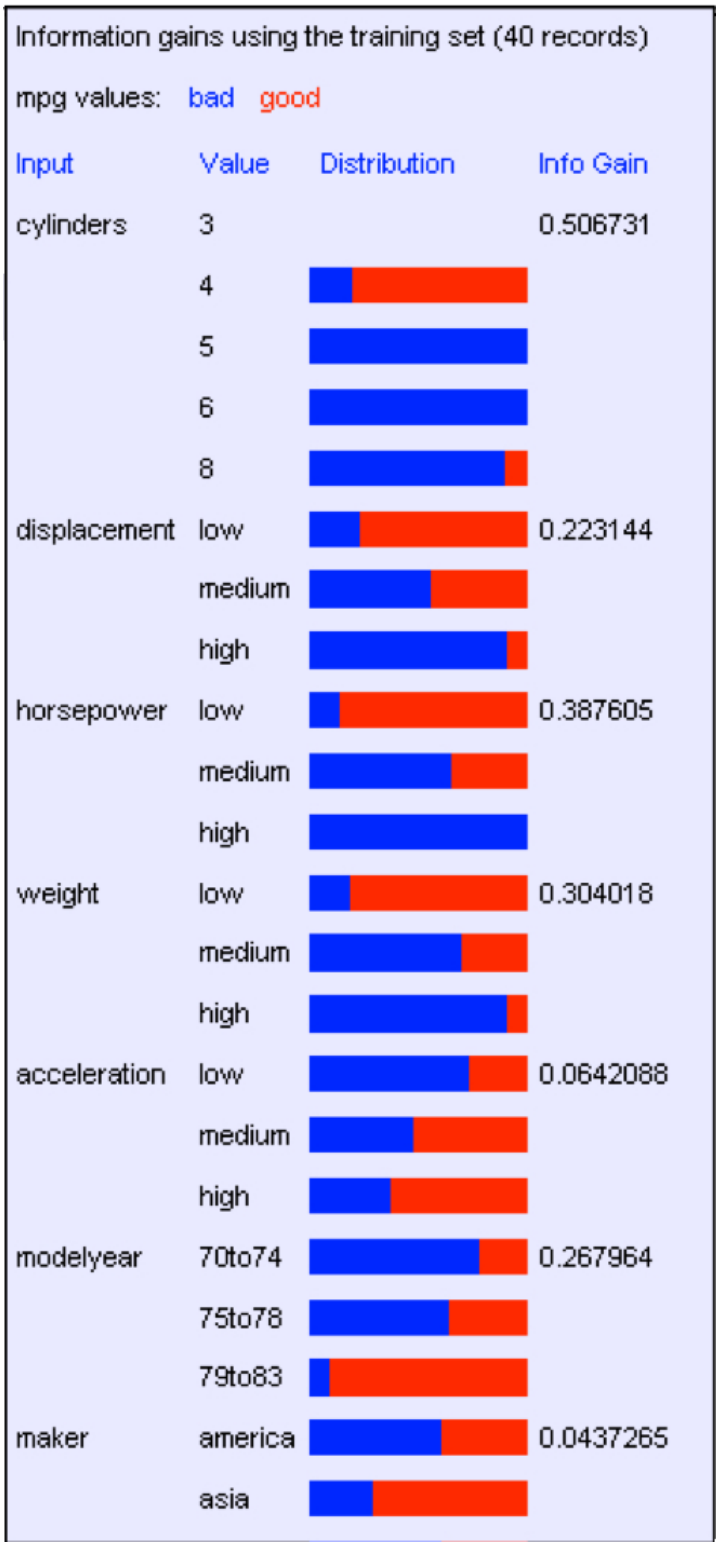


Figure 7: Information gain

# 4 When to stop splitting

- **matching records have the same attribute value**. REMARK: $H(Y) = 0$
- **No attributes can further distinguish records**. REMARK : $H(Y|X) = H(Y)$ for any feature $X$

### 4.0.1 0 information gain case in general

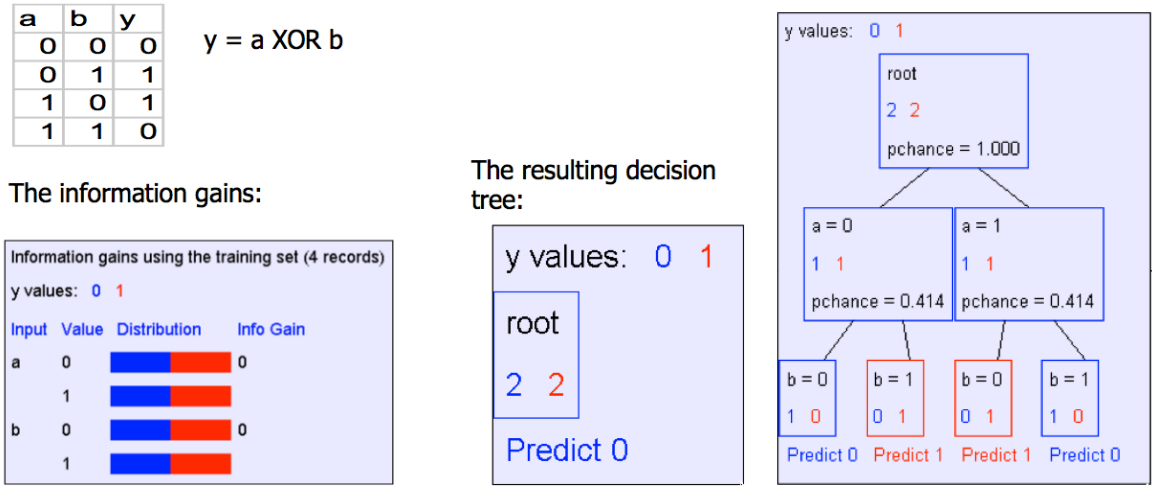


Figure 8: "Do not split" VS "split" when Information gain is 0

# 5 Real-valued inputs

If the input values of $X$ are real (and/or continuous), then splitting branches by feature-values are not feasible. Instead find the best threshold $t$ for the feature $X$.

$$H(Y|X:t) = P(X < t)H(Y|X < t) + P(X \geq t)H(Y|X \geq t)$$
$$IG(X:t) = H(Y) - H(Y|X:t)$$

Find $t$ that maximizes $IG(X:t)$. To do so, a possibility is to consider all $t$ of the form $(x_i + x_{i+1})/2$, where $x_1, x_2, ..., x_n$ are the values of feature $X$ in the training set.

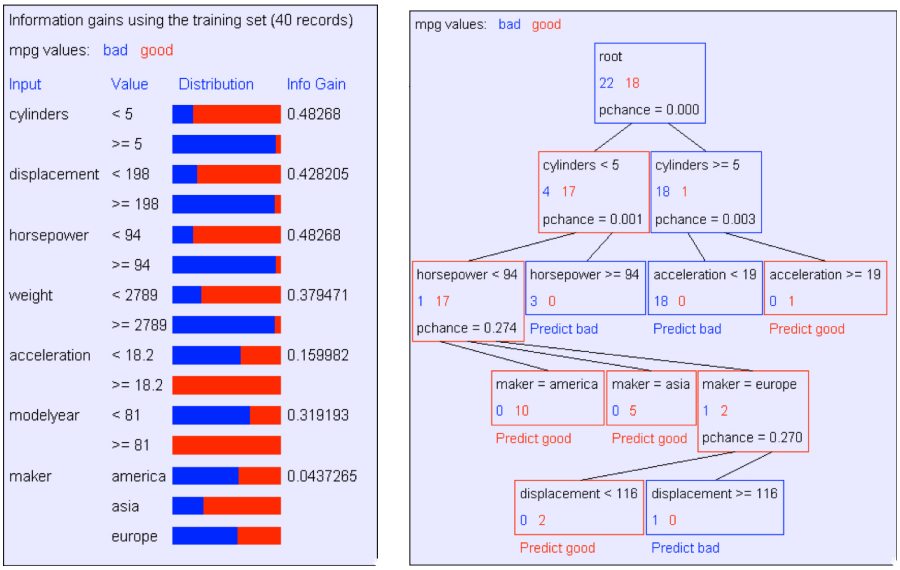


Figure 9: Real-valued information gain and decision tree

# 6 Regression trees

Lets say that fro each node $m$, $\chi_m$ is the set of datapoints reaching that node.

**Estimate a predicted value per tree node**

$$g_m = \frac{\sum_{t \in \chi_m} y_t}{|\chi_m|}$$

**Calculate mean square error**

$$E_m = \frac{\sum_{t \in \chi_m} (y_t - g_m)^2}{|\chi_m|}$$

How to choose the next split. If $E_m < \theta$, then stop splitting. Otherwise choose the split that realizes the maximum drop in error for all all brances. Say we are considering feature $X$ with branches $x_1, x_2, ..., x_k$, and lets call $\chi_{mj}$ the subset of $\chi_m$ for which $X = x_j$.

$$g_{mj} = \frac{\sum_{t \in \chi_{mj}} y_t}{|\chi_{mj}|}$$

$$E'_m(X) = \frac{\sum_j \sum_{t \in \chi_{mj}} (y_t - g_{mj})^2}{|\chi_m|}$$

We shall choose $X$ such that $E'_m(X)$ is minimized, or the drop in error is maximized.

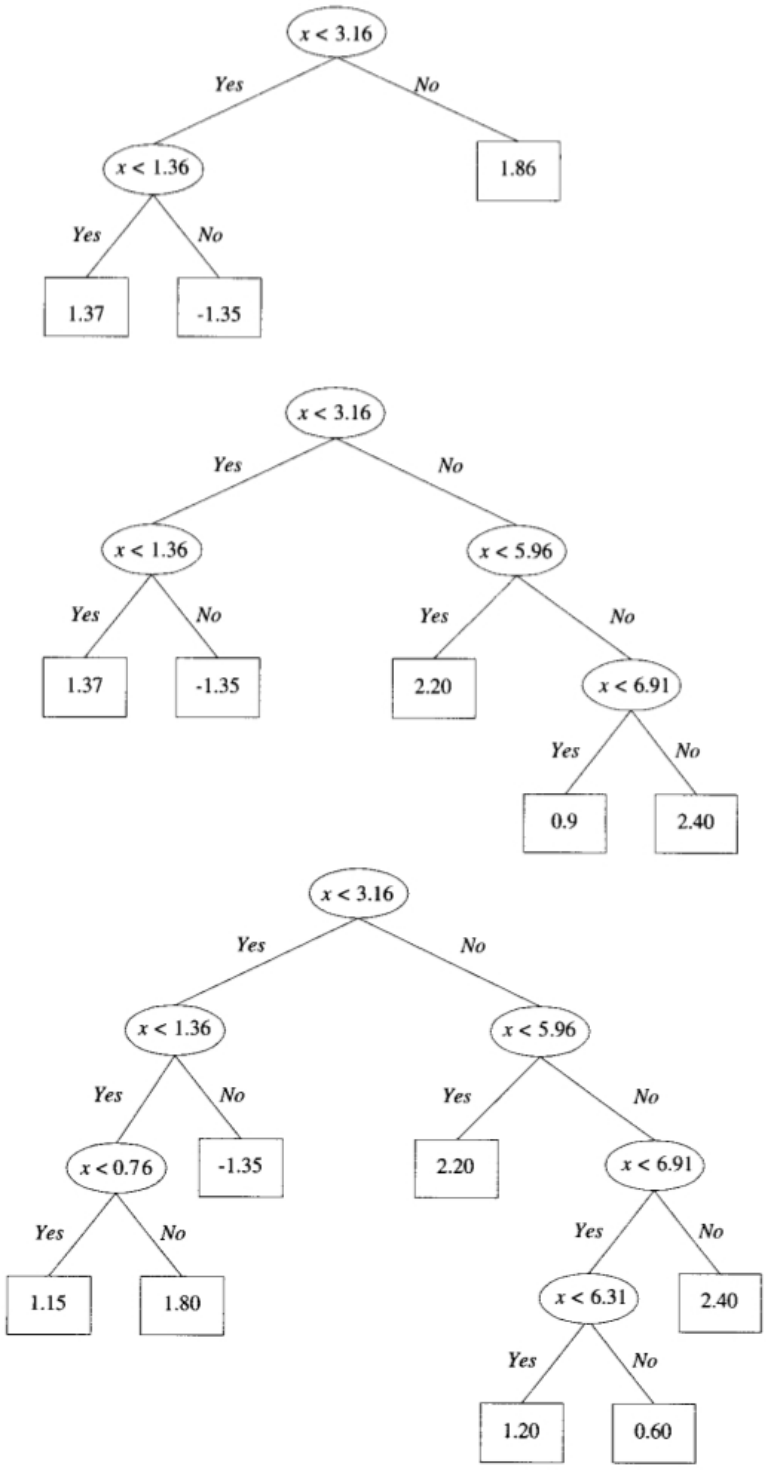


Figure 10: Regression tree

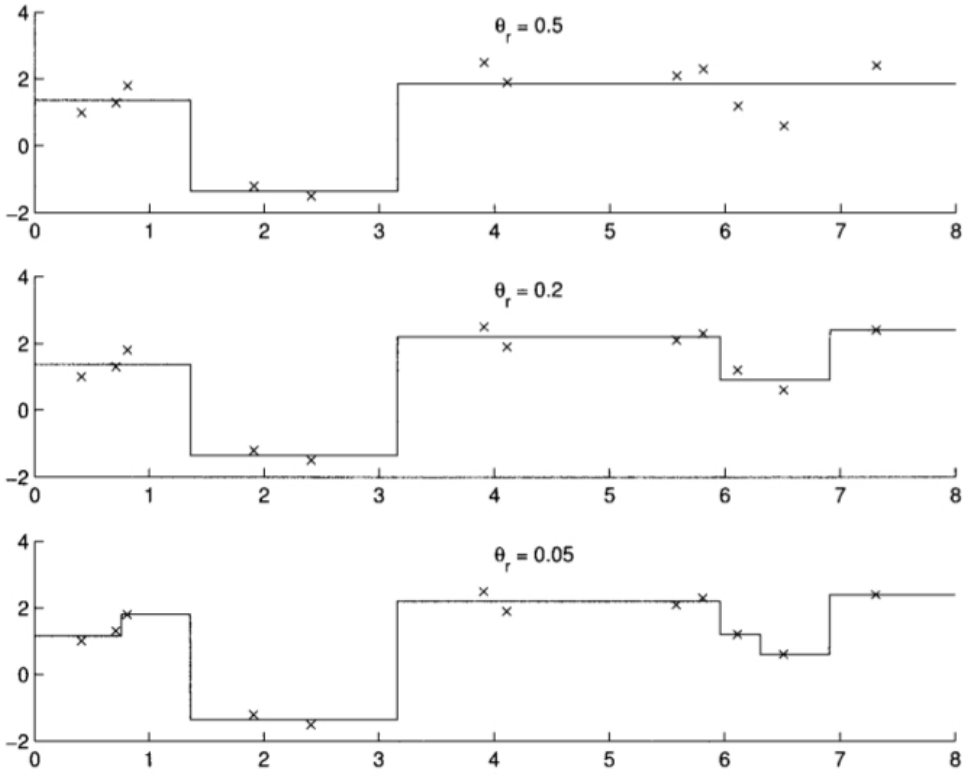


Figure 11: Regression fit

# 7 Pruning

If a tree is "too small", the model does not capture all structure of data, or it *underfits*. if the tree is too big, it captures structure that is too local and it cannot be generalized (*overfits*). Pruning helps heuristically to find the appropriate tree size.

**Pre-pruning** If a tree node contains less that, say, 5% of the training set, stop splitting (even if there are features with positive information gain).

**Post-pruning** Grow the tree until all positive information gains are used for splitting; then find the overfitting subtrees and merge them together. To do so, we need a pruning set (separate from testing or validation sets): if merging subtrees does not increase the classification error on the pruning set (by more than $\epsilon$), then we merge the subtrees.

# 8 Rules extraction

Go over the branches of the tree and write down the splits. For example, for the tree in figure 9, some rules are:

```
IF (CYlinders<5) AND (horsepower<94) AND (maker= asia) THEN "predict good"
IF (cylinders>=5) AND (acceleration<19) THEN "predict bad"
....
```

**Rules extraction directly from data**. Also based on Information gain, but it traverses the data DFS instead of BFS.

# 9 Multivariate tree

In a multivariate tree, the splitting criteria can be a functional of more than one feature. For example, at the root we can have the following split:

$$\texttt{cylinders} * 20 + \texttt{horsepower} < 180$$

More generally, a binary linear multivariate node $m$ split can look like

$$w^1 x^1 + w^2 x^2 + ...w^d x^d + w^0 > 0$$

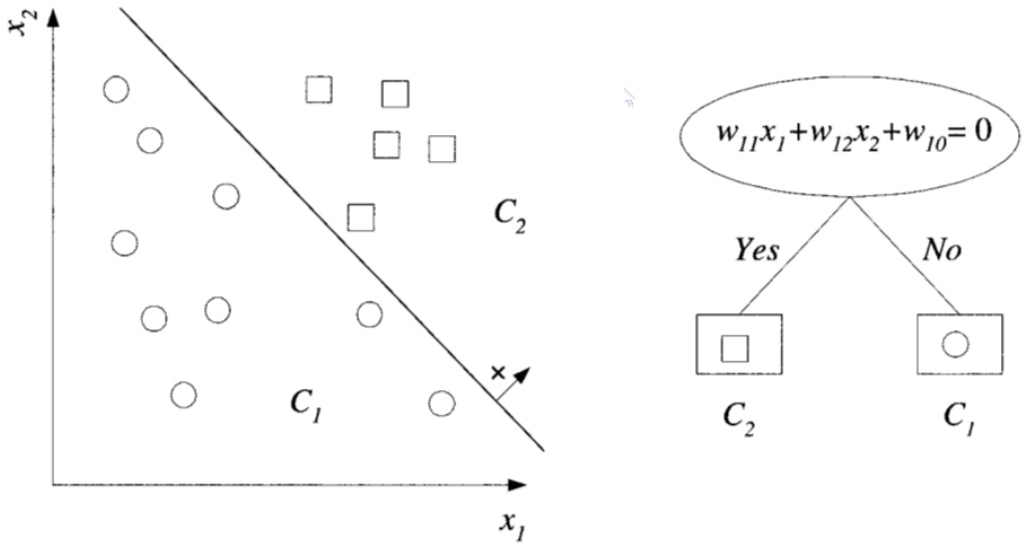


Figure 12: Multivariate split

Such splits can be extremely powerful (if data is linearly separable, a single split at root can create a perfect classification); even more complex splits can be obtained using nonlinear split functions.

However, finding a good multivariate split is not anymore a matter of brute force: there are $2^d\binom{N}{d}$ possible splits (or hyperplanes). Later on in the course we will discuss linear classification and how good hyperplanes can be obtained without an exhaustive search.