

Data Mining and Machine Learning

ID3 and Regression

Gergely Horváth

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Outline

1 Decision trees

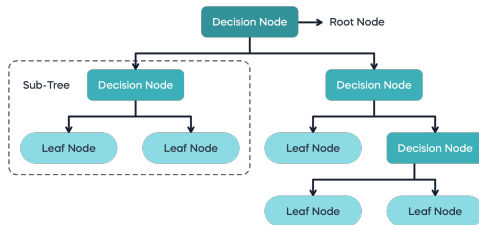
- ID3
- ID3

2 Regression

Decision trees

Classification and regression trees (categorical & numerical data handling)

- Splits dataset into small subsets
- Final result: tree with
 - root node
 - decision nodes: branches → possible values for the attribute
 - leaf nodes: represents a classification
- Which feature splits the data better (which is the best attribute)?



ID3 – Iterative Dichotomiser 3

- Core algorithm for building decision trees
 - top-down, greedy search to test each attribute at every node of the tree
- Which is the best attribute?
 - the one which will result in the smallest tree
 - choose the attribute that produces the “purest” nodes
 - information gain (IG)
 - information before splitting – information after splitting
 - is used to construct a tree
- Entropy: a measure of randomness
 - unbiased coin toss (head and tail is equally likely): $E = 1$
 - biased (2 head): $E = 0$
 - ID3 uses entropy to calculate the homogeneity of a sample
 - $E(p_1, p_2, \dots, p_N) = -p_1 \log(p_1) - p_2 \log(p_2) - \dots - p_N \log(p_N)$

ID3 – Iterative Dichotomiser 3

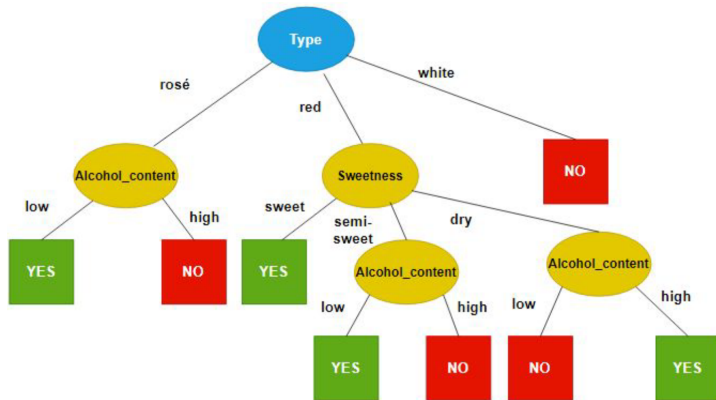
General recipe:

- 1 Compute the overall entropy of the class distribution
- 2 Choose a feature and compute the entropy of the class distribution for each unique value
- 3 Calculate a weighted sum of the unique values' entropy (weight is its relative overall presence in the examined dataset)
- 4 Subtract this value from the overall entropy to receive the information gain
- 5 Repeat the previous 3 steps for every attribute and choose the one with the highest information gain
- 6 Make the first split with the best performing attribute
- 7 Split the data with respect to the chosen feature
- 8 Where the subsets are uniform in class values, that class value should be assigned to that node which will make it a leaf node, otherwise, the previous steps should be repeated for the assigned subsets

ID3 – Iterative Dichotomiser 3

Alcohol_content	Sweetness	Type	(Year)	Popular
low	sweet	rosé	2012	yes
low	dry	red	2009	no
low	semi-sweet	red	2008	yes
high	sweet	rosé	2013	no
low	dry	white	2013	no
low	sweet	white	2006	no
high	semi-sweet	red	2011	no
high	sweet	red	2007	yes
high	dry	red	2005	yes

ID3 – Iterative Dichotomiser 3



Regression – I.

The "reality":

$$y = \sum_i^n (p_i \cdot x_i) + \epsilon$$

The model:

$$\hat{y} = \sum_i^n (p_i \cdot x_i)$$

Let us have matrix notations:

$$Y = \begin{bmatrix} y^{(1)} \\ y^{(2)} \\ \vdots \\ y^{(m)} \end{bmatrix}$$

$$X = \begin{bmatrix} x^{(1)} \\ x^{(2)} \\ \vdots \\ x^{(m)} \end{bmatrix}$$

Where $Y \in \mathbb{R}^{m \times 1}$, $X \in \mathbb{R}^{m \times n}$ and $p \in \mathbb{R}^{n \times 1}$.

Regression – II.

Notation repetition:

$$Y = X \cdot p + \epsilon$$

$$\hat{Y} = X \cdot p$$

The next step is to define the error made by the model (in this case the – arguably – simplest convex function will be used):

$$L(D, p) = \|Y - \hat{Y}\|_2 = \|Y - (X \cdot p)\|_2$$

The error is expressed with a 2-norm, so the next step is to minimize the loss, to find the global minimum of this function. → This is going to be easy and trivial since we have a convex function.

Logistic regression will be a model, where our linear regression model is being fed to a sigmoid function:

$$\text{sigmoid}(\text{model}) = \frac{1}{1 + e^{-\text{model}}}$$