

MODEL SELECTION AND EVALUATION

Empirical error and overfitting

Evaluation methods

Performance measure

Comparison test

Bias and variance





SUPERVISED LEARNING



WHAT IS SUPERVISED LEARNING?

Supervised

Source: Ben Freundorfer

Doug Rose defines supervised learning as "When a data scientist acts like a tutor for the machine, training it by showing it basic rules and giving it an overall strategy." [5]

- Regression model
- Classification model



SUPERVISED MACHINE LEARNING

- We humans learn from past experiences.
- A computer does not "experience."
 - A computer system learns from data, which represents "past experiences" in an application domain.
- Our focus: learn a target function that can be used to predict the values (labels) of a discrete class attribute, e.g.,
 - high-risk or low risk and approved or not-approved.
- The task is commonly called: supervised learning, classification, or inductive learning.



EXAMPLE APPLICATION

A credit card company receives thousands of applications for new cards. Each application contains information about an applicant,

- age
- annual salary
- outstanding debts
- credit rating
- etc.

<u>Problem</u>: Decide whether an application should be approved, i.e., **classify** applications into two categories, **approved** and **not approved**.

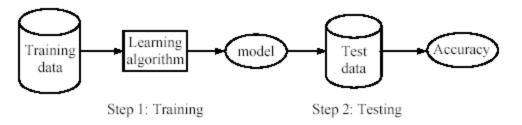


SUPERVISED LEARNING PROCESS: TWO STEPS

Learning or training: Learn a model using the training data (with labels)

Testing: Test the model using unseen test data (without labels) to assess the model accuracy

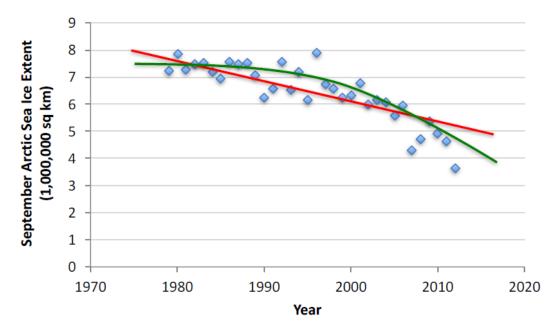
$$Accuracy = \frac{\text{Number of correct classifications}}{\text{Total number of test cases}},$$





REGRESSION MODEL

- Given (x_1, y_1) , (x_2, y_2) , ..., (x_n, y_n)
- Learn a function f(x) to predict y given x
 - -y is real-valued == regression



Data from G. Witt. Journal of Statistics Education, Volume 21, Number 1 (2013)

REGRESSION

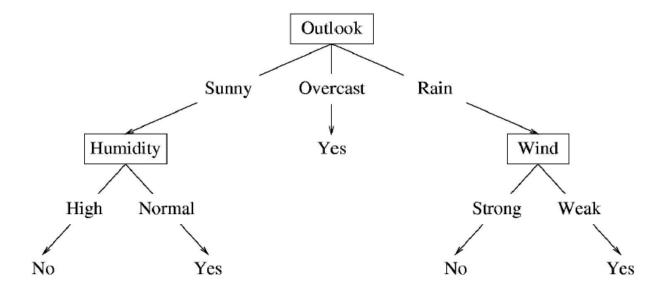
Types of Regression Algorithm:

- Simple Linear Regression
- Multiple Linear Regression
- Polynomial Regression
- Support Vector Regression
- Decision Tree Regression
- Random Forest Regression



DECISION TREE

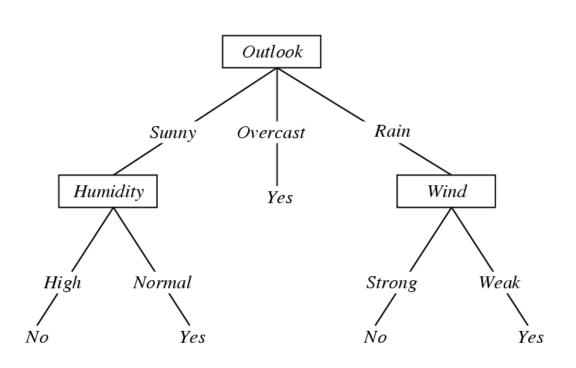
A possible decision tree for the data:



- Each internal node: test one attribute X_i
- Each branch from a node: selects one value for X_i
- Each leaf node: predict Y (or $p(Y \mid \boldsymbol{x} \in \text{leaf})$)

DECISION TREE LEARNING

Decision Tree for PlayTennis

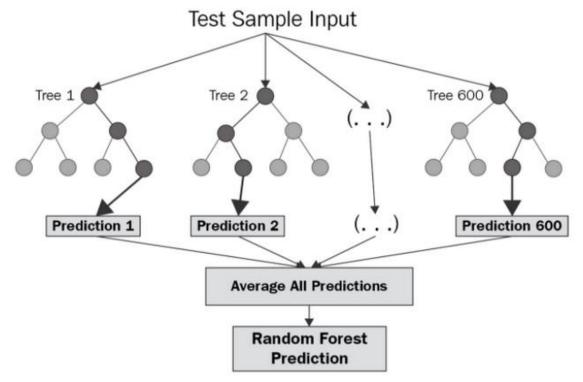


Problem Setting:

- Set of possible instances X
 - each instance x in X is a feature vector
 - e.g., <Humidity=low, Wind=weak, Outlook=rain, Temp=hot>
- Unknown target function $f: X \rightarrow Y$
 - Y is discrete valued
- Set of function hypotheses $H = \{ h \mid h : X \rightarrow Y \}$
- each hypothesis *h* is a decision tree
- trees sorts x to leaf, which assigns y

RANDOM FOREST REGRESSION

Random Forest Regression is a supervised learning algorithm that uses **ensemble learning** method for regression. Ensemble learning method is a technique that combines predictions from multiple machine learning algorithms to make a more accurate prediction than a single model.



Source: Chava Bakshi

RANDOM FOREST REGRESSION

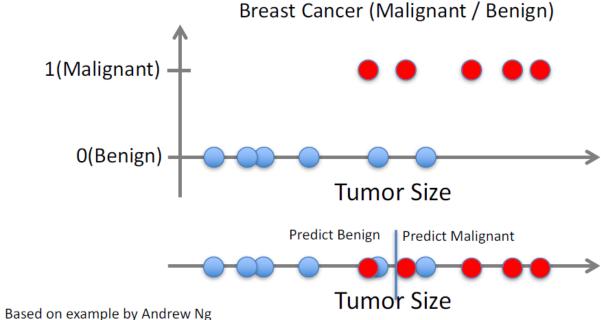
To get a better understanding of the Random Forest algorithm, let's walk through the steps:

- Pick at random k data points from the training set.
- Build a decision tree associated to these k data points.
- Choose the number N of trees you want to build and repeat steps 1 and 2.
- For a new data point, make each one of your N trees predict the value of y for the data point in question and assign the new data point to the average across all of the predicted y values.



CLASSIFICATION MODEL

- Given (x_1, y_1) , (x_2, y_2) , ..., (x_n, y_n)
- Learn a function f(x) to predict y given x
 - -y is categorical == classification



CLASSIFICATION

Types of ML Classification Algorithms:

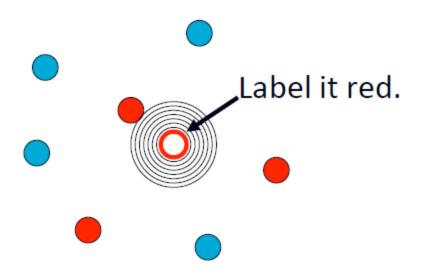
Classification Algorithms can be further divided into the following types:

- Logistic Regression
- K-Nearest Neighbours
- Support Vector Machines
- Kernel SVM
- Naïve Bayes
- Decision Tree Classification
- Random Forest Classification



K-NEAREST NEIGHBOUR

- 1-Nearest Neighbour
- One of the simplest of all machine learning classifiers
- Simple idea: label a new point the same as the closest known point





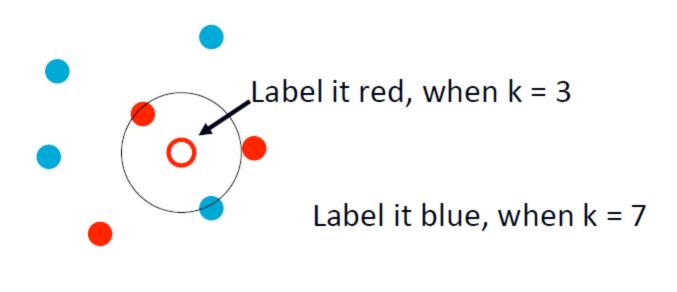
K-NEAREST NEIGHBOUR

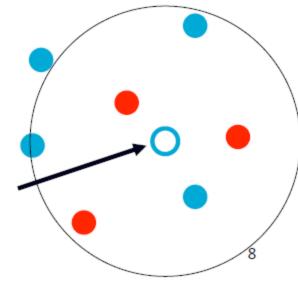
- Four Aspects of an Instance---Based Learner:
- 1. A distance metric
- 2. How many nearby neighbours to look at?
- 3. A weighting function (optional)
- 4. How to fit with the local points?



K-NEAREST NEIGHBOUR

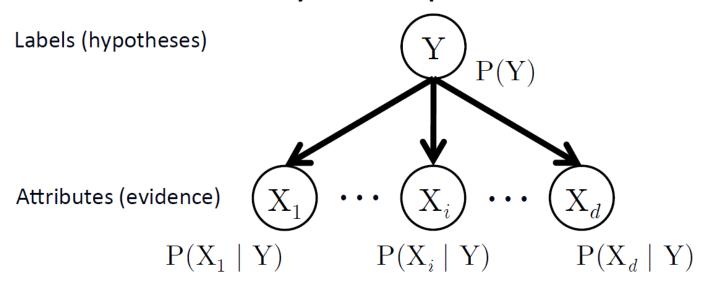
- Generalizes 1---NN to smooth away noise in the labels
- A new point is now assigned the most frequent label of its *k* nearest neighbours







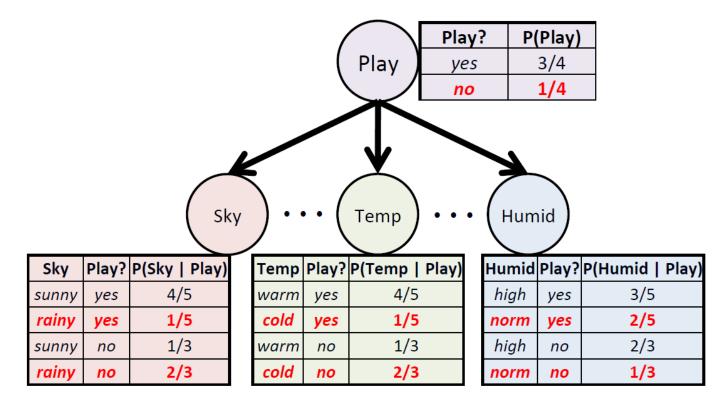
The Naïve Bayes Graphical Model



- Nodes denote random variables
- Edges denote dependency
- Each node has an associated conditional probability table (CPT), conditioned upon its parents



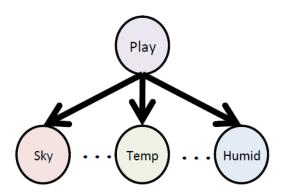
Example NB Graphical Model



• Some redundancies in CPTs that can be eliminated



Example Using NB for Classification



| Play? | P(Play) |
|-------|---------|
| yes | 3/4 |
| no | 1/4 |

| Temp | Play? | P(Temp Play) |
|------|-------|----------------|
| warm | yes | 4/5 |
| cold | yes | 1/5 |
| warm | no | 1/3 |
| cold | no | 2/3 |

| Sky | Play? | P(Sky Play) |
|-------|-------|---------------|
| sunny | yes | 4/5 |
| rainy | yes | 1/5 |
| sunny | no | 1/3 |
| rainy | no | 2/3 |

$$h(\mathbf{x}) = \underset{y_k}{\operatorname{arg\,max}} \log P(Y = y_k) + \sum_{j=1}^{d} \log P(X_j = x_j \mid Y = y_k)$$

Goal: Predict label for x = (rainy, warm, normal)



• Advantages:

- Fast to train (single scan through data)
- Fast to classify
- Not sensitive to irrelevant features
- Handles real and discrete data
- Handles streaming data well

• Disadvantages:

• Assumes independence of features



REFERENCES

