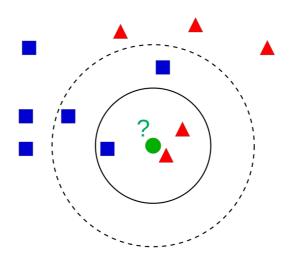


Blockkurs: Introduction to Machine Learning for Psychologists

K Nearest Neighbor



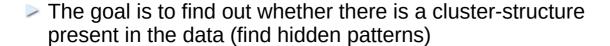
Yannick Rothacher

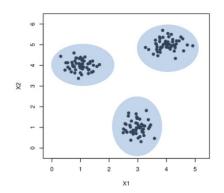
Zürich, FS2025



Recap: k-means clustering

- The k-means algorithm is a partitional clustering method
- Often there is not a "correct" solution in k-means clustering





- Such an approach is generally referred to as unsupervised learning
- In contrast, in supervised learning we try to predict the correct class/value of a new data point
 - > Techniques, which try to predict the correct class of data are referred to as classifiers



Supervised learning: Classification

- Classifiers can be used to predict all kinds of data labels
 - Who will win in the next football-worldcup?
 - Will an operation be successful?
 - Which type of tumor is shown on a medical scan?
- Predictions are always based on training data:

| Hours studied | Score last year | IQ | Exam 4 result |
|------------------|--------------------|-----|-----------------------------|
| 1 | 20 | 101 | failed |
| 20 | 99 | 105 | passed |
| 12 | 54 | 99 | passed |
| | | | |

Target variable

New observation:

Will this person pass or fail?

| Hours | Score last | IQ | Exam |
|---------|------------|----|--------|
| studied | year | | result |
| 25 | 70 | 98 | ? |



Supervised learning: Classification

- The workflow for classifiers is always similar
- First, a model is fitted to training data, for which the true class labels are known
- Afterwards, the model is applied to new observations to predict their class labels
- Thus, we always first need training data to train our model
- ▶ There are different statistical models, which can be used as a classifier
 - e.g. logistic regression is a simple classifier usually applied when the target label has two levels (e.g. predict sex: male/female)

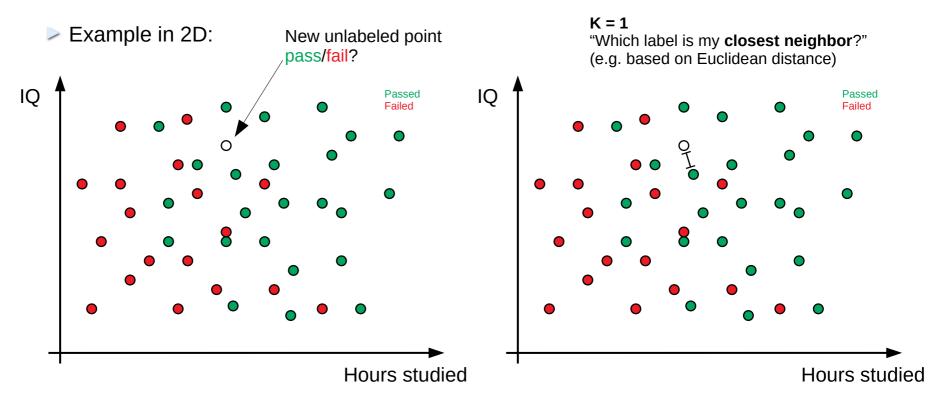
K-nearest neighbor is an easy to understand classifier and, therefore, a good introduction to the topic of classification

| Hours studied | Score last year | IQ | Exam result |
|------------------|--------------------|-----|----------------|
| 1 | 20 | 101 | failed |
| 20 | 99 | 105 | passed |
| 12 | 54 | 99 | passed |
| | | | |



K-nearest neighbor (KNN)

- Not to confuse with k-means algorithm!
- Idea: Classify a new data-point based on the labels of the neighboring training data points

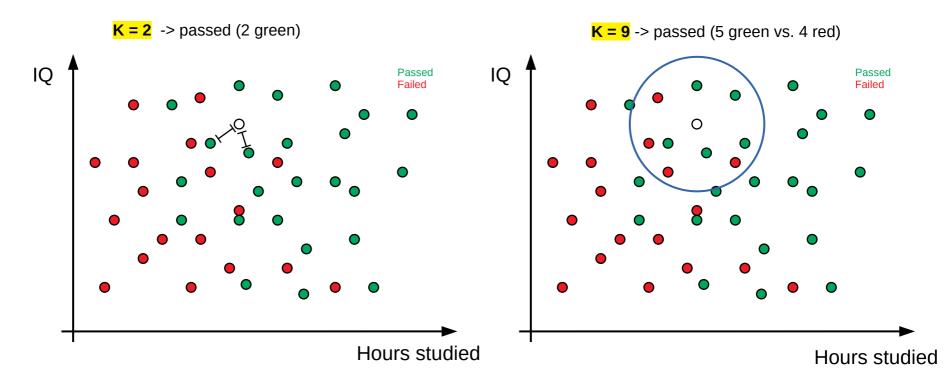


Assign the label "passed" because closest neighbor is "passed" as well



K-nearest neighbor (KNN)

► The parameter k determines how many neighbors are considered (use the majority vote as prediction):



- The prediction may change depending on k
- There is the possibility of ties, what could be done to deal with ties?



KNN - Decision boundaries

► The decision boundaries for new data points depend on how many neighbors are considered:

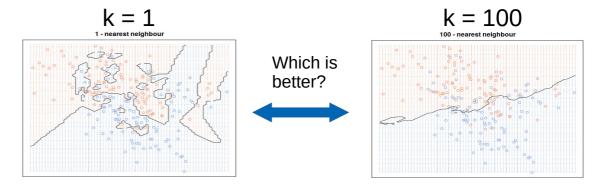


- ► The decision boundary becomes **simpler** with larger k
 - What is the best value of k?



How to choose k

What is the best value of k?



- In order to decide what the best value for k is we want to measure how "well" the classifier is performing
 - How can we do that?
- Simple idea: Check how correctly the classifier is labeling the training data (Training error)
 - Confusion matrices (for k=1 and k=100):

$$k=1$$
 $k=100$ true label true label KNN prediction blue red blue 100 0 blue 66 26 red 0 100 red 34 74

Train- and test-error

Calculate the training error (number of miss-classifications divided by total number of predictions):

- ► Looking at the training error is not that informative. In our example, K=1 forces a 0% training error rate.
- Because we want to use the classifier to predict the class of new observations (not included in the training set), it makes sense to calculate the error rate for unseen data (Test-error)



Train- and test-error

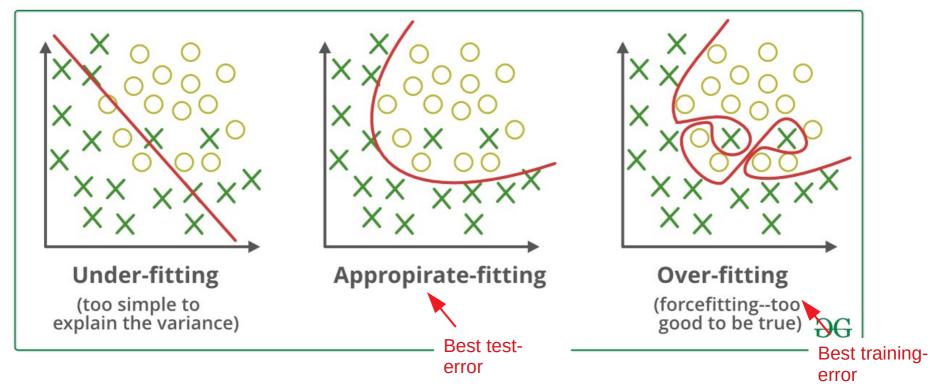
► Test-error for 40 new observations (for k=1 and k=100)

```
k = 1
                                                           k = 100
              true label
                                                          true label
KNN prediction blue red
                                           KNN prediction blue red
          blue
                 12 8
                                                      blue
                                                             13 7
                  3 17
          red
                                                      red
                                                              2 18
                          \rightarrow 11/40 = 27.5\%
                                                                      \rightarrow 9/40 = 22.5%
```

- Now the two classifiers both perform equally well (in this case k=100 even slightly better)
- ▶ The goal is to keep the **test-error** as low as possible
- Most probably, setting k=1 and setting k=100 are both not ideal
 - k=1 is too flexible (over-fitting the training data) and k=100 is perhaps too simple (under-fitting the training data)



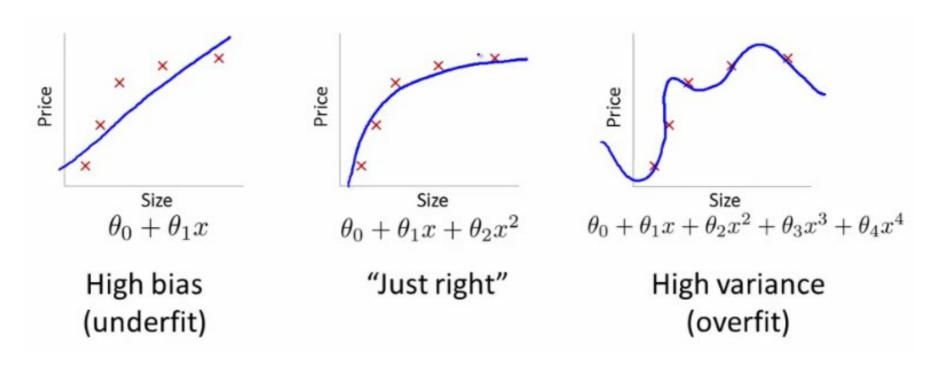
Over- vs. under-fitting



- Imagine you apply repeatedly an over- or under-fitting model to collected data:
 - Under-fitting will show high bias and low variance
 - Over-fitting will show low bias and high variance



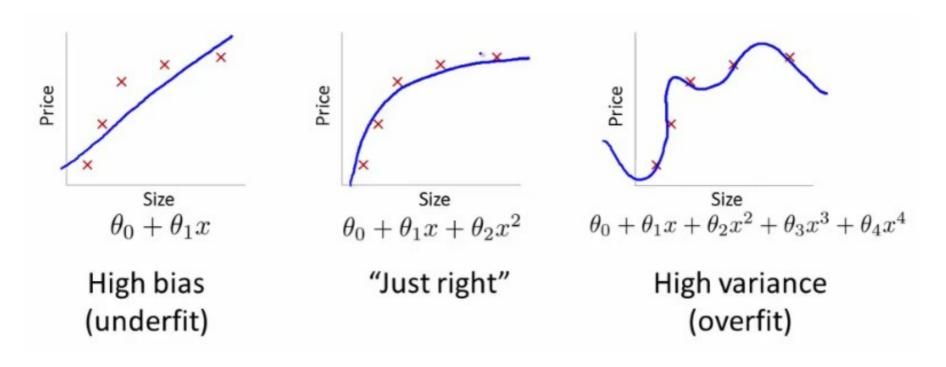
Over- vs. under-fitting (in regression)



- Imagine you apply repeatedly an over- or under-fitting model to collected data:
 - Under-fitting will show high bias and low variance
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Over- vs. under-fitting (in regression)



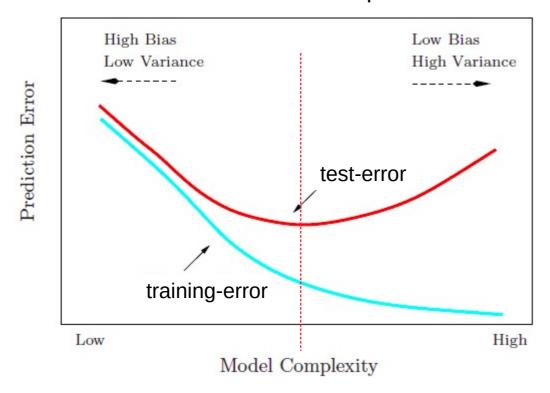
- Imagine you apply repeatedly an over- or under-fitting model to collected data:
 - Under-fitting will show high bias and low variance
 - Over-fitting will show low bias and high variance

 \rightarrow show in R



Where is the optimum between over- and under-fitting?

Test-error should be as small as possible



- Dashed, red line shows the ideal model flexibility
 - Everything to the right of dashed line is over-fitting
 - Everything to the left of dashed line is under-fitting



Outlook: Cross-validation

- It is important to estimate the test-error of a classifier to judge its performance
- Ideally, one uses a training data set to fit the model and separate test data set to assess its performance
 - Often a separate test data set is not available
- In the next topic we see what methods are used to estimate the test-error using **one single data set** (cross-validation)



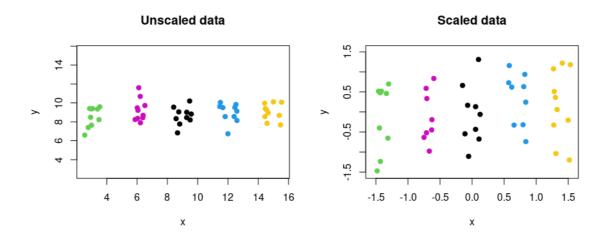
KNN in R

```
library(class)
knn.pred <- knn(train = iris[,1:4], # use the iris data as training data
                cl = iris$Species, # correct labels of training data
                test = iris[,1:4], # knn always needs a specified test set
                k = 5
                                     # parameter k
# Confusion matrix:
confT <- table(knn.pred, iris$Species)</pre>
confT
knn.pred setosa versicolor virginica
                                                              Because we do not have a
  setosa
                 50
                                                              test set, we reuse the
 versicolor 0
                            47
                                                              training data as test data
  Virginica
                                      48
# Training error:
missT <- confT
diag(missT) <- 0 # turn diagonal values to zeros</pre>
missCount <- sum(missT) # count the miss-classifications (5)
trainErr <- missCount/sum(confT)</pre>
trainErr
[1] 0.03333333
```



Standardization in KNN

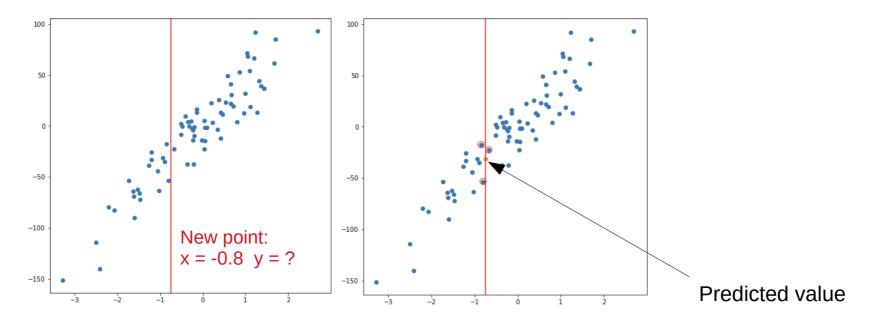
- Similar to the case of PCA it often makes sense to standardize the data before applying KNN classification
 - Why could standardization be important?
- ► Example: If one variable has a much larger variance than all the others (e.g. ranging from 0 100'000, other variables: 0 0.1), it will be much more important in determining the Euclidean distance between two data points.
 - Other variables are neglected!
- Through standardization, all variables have equal variance
 - In R: use the scale() function to standardize numeric vectors





KNN for regression

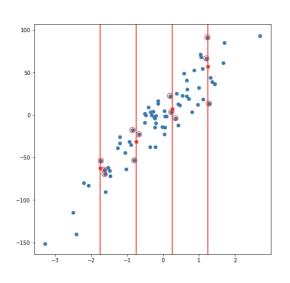
- KNN cannot only be used for classification
 - Can also predict numerical values
 - Instead of taking the majority vote of neighbors, take e.g. the **mean of k neighbors**

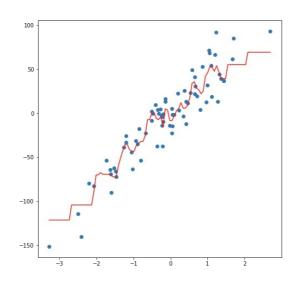


- ► E.g. univariate data (figures above): Use the average of **k=3** closest neighbors (with respect to x value) as a prediction for y value
 - Can use the distances as weights for the mean

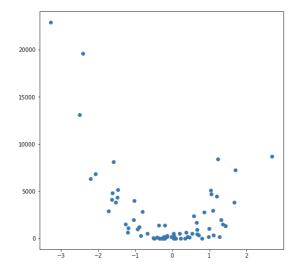


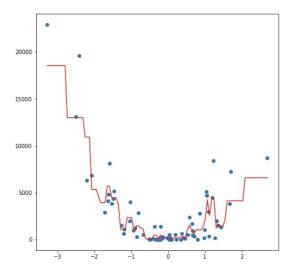
KNN for regression





- KNN for linear relation
- Result is not a clean line but still general pattern is captured





- Advantage of KNN for regression:
 - Simple algorithm which can capture non-linearity quite well