10-601
Machine Learning

Classification

Where we are

√

Today

Later

Classification

- Assume we want to teach a computer to distinguish between cats and dogs …








Bayes decision rule

- If we know the conditional probability p(x | y) and class priors p(y) we can determine the appropriate class by using Bayes rule:

$$P(y=i \mid x) = \frac{P(x \mid y=i)P(y=i)}{P(x)} \overset{def}{=} q_i(x)$$

- We can use $q_i(x)$ to select the appropriate class.
- We chose class 0 if $q_0(x) \geq q_1(x)$ and class 1 otherwise
- This is termed the 'Bayes decision rule' and leads to optimal classification.
- However, it is often very hard to compute …

Minimizes our probability of making a mistake

Note that p(x) does not affect our decision

Bayes decision rule

$$P(y=i \mid x) = \frac{P(x \mid y=i)P(y=i)}{P(x)} \stackrel{def}{=} q_i(x)$$

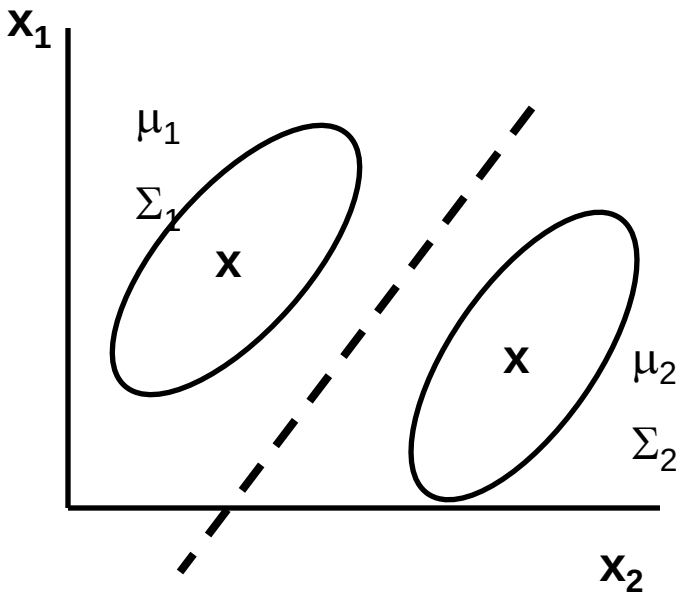
- We can also use the resulting probabilities to determine our confidence in the class assignment by looking at the likelihood ratio:

$$L(x) = \frac{q_0(x)}{q_1(x)}$$

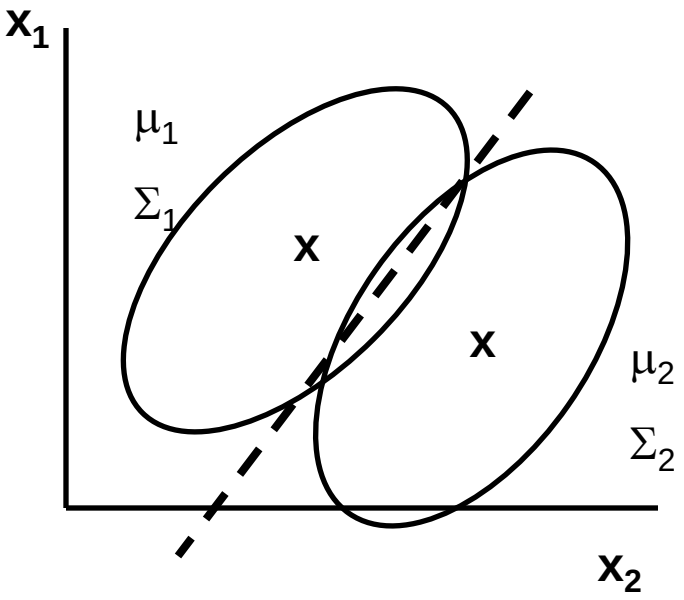
Also known as likelihood ratio test, we will talk more about this later

Bayes decision rule: Example

Normal Gaussians

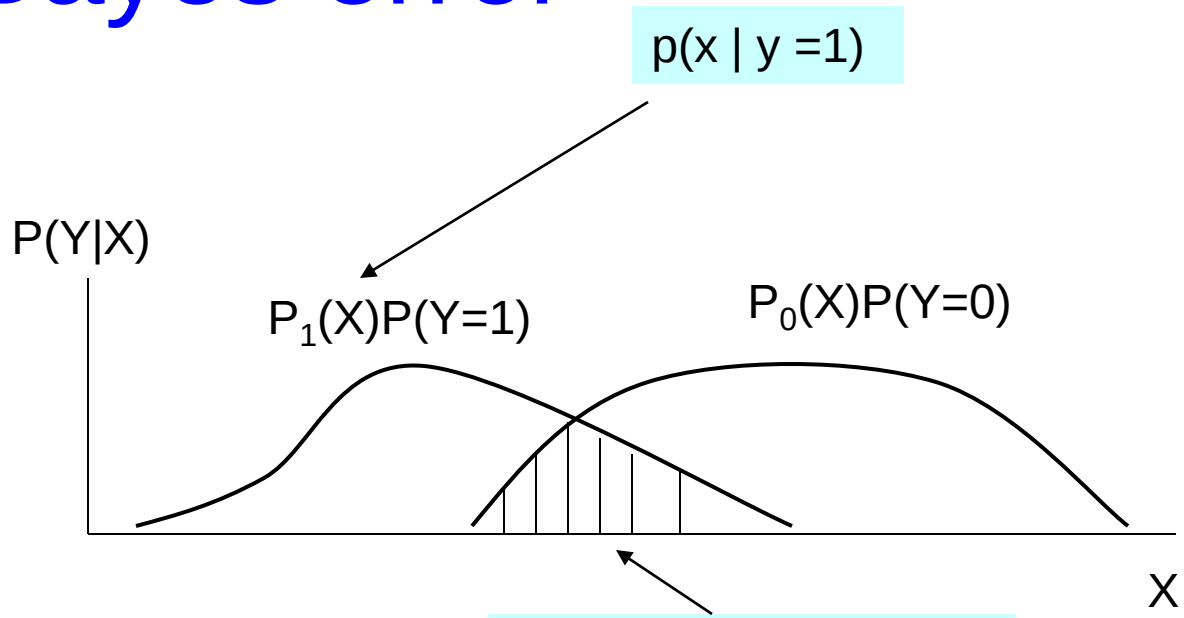


Normal Gaussians



Bayes error

- For the Bayes decision rule we can calculate the probability of an error
- This is the probability that we assign a sample to the wrong class, also known as the **risk**

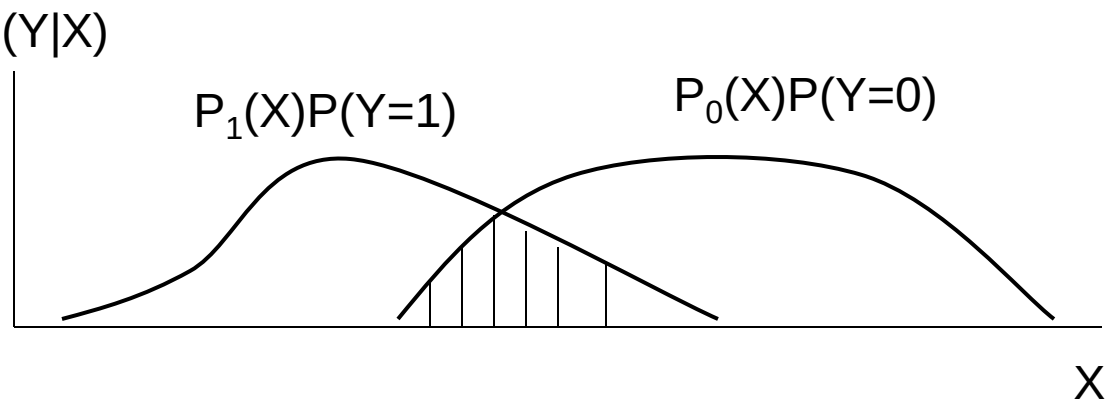


- The risk for sample x is:

$$R(x) = \min\{P_1(x)P(y=1), P_0(x)P(y=0)\}$$

Risk can be used to determine a ‘reject’ region

Bayes error



• The probability that we assign a sample to the wrong class, is known as the **risk**

• The risk for sample x is:

$R(x) = \min\{P_1(x)P(y=1), P_0(x)P(y=0)\}$

• We can also compute the expected risk (the risk for the entire range of values of x):

Assuming all values equally likely

$$E[r(x)] = \int_x r(x)p(x)dx$$

$$= \int_x \min\{p_1(x)p(y=1), p_0(x)p(y=0)\}dx$$

$$= p(y=0)\int_{L_1} p_0(x)dx + p(y=1)\int_{L_0} p_1(x)dx$$

L_1 is the region where we assign instances to class 1

Loss function

- The risk value we computed assumes that both errors (assigning instances of class 1 to class 0 and vice versa) are equally harmful.
- However, this is not always the case.
- Why?
- In general our goal is to minimize loss, often defined by a loss function: $L_{0,1}(x)$ which is the penalty we pay when assigning instances of class 0 to class 1

$$E[L] = L_{0,1}\, p(y=0) \int_{L_1} p_0(x)dx + L_{1,0}\, p(y=1) \int_{L_0} p_1(x)\, p(y=1)dx$$

Types of classifiers

- We can divide the large variety of classification approaches into roughly two main types

1. Instance based classifiers
   - Use observation directly (no models)
   - e.g. K nearest neighbors

2. Generative:
   - build a generative statistical model
   - e.g., Bayesian networks

3. Discriminative
   - directly estimate a decision rule/boundary
   - e.g., decision tree

Classification

- Assume we want to teach a computer to distinguish between cats and dogs …

Several steps:

1. feature transformation
2. Model / classifier specification
3. Model / classifier estimation (with regularization)
4. feature selection

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How do we encode the picture? A collection of pixels? Do we use the entire image or a subset? …

Classification

- Assume we want to teach a computer to distinguish between cats and dogs …

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What type of classifier should we use?

Classification

- Assume we want to teach a computer to distinguish between cats and dogs …




Several steps:

1. feature transformation
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3. Model / classifier estimation (with regularization)
4. feature selection

How do we learn the parameters of our classifier? Do we have enough examples to learn a good model?

Classification

- Assume we want to teach a computer to distinguish between cats and dogs …




Several steps:

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3. Model / classifier estimation (with regularization)
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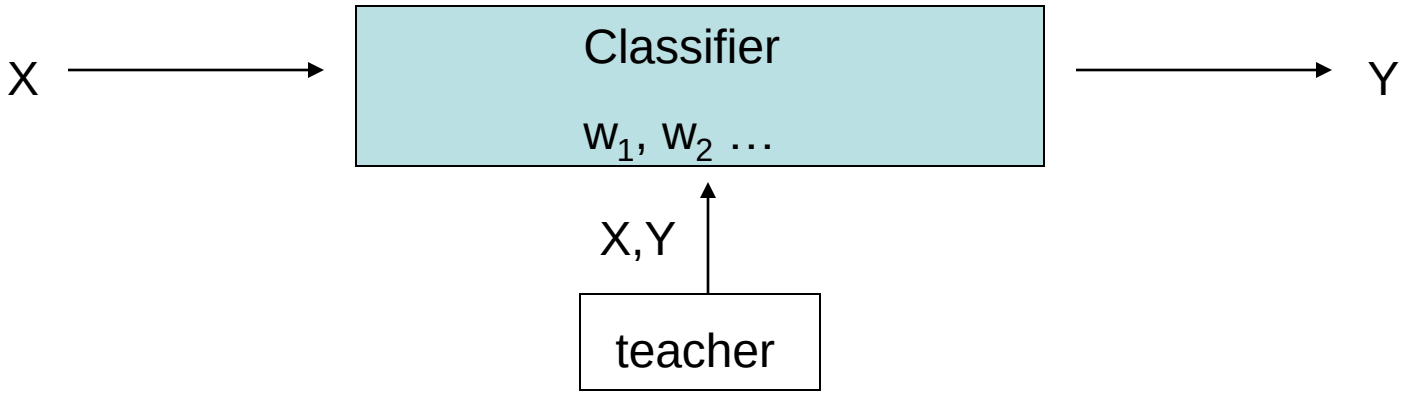
Do we really need all the features? Can we use a smaller number and still achieve the same (or better) results?

Supervised learning

- Classification is one of the key components of 'supervised learning'
- Unlike other learning paradigms, in supervised learning the teacher (us) provides the algorithm with the solutions to some of the instances and the goal is to generalize so that a model / method can be used to determine the labels of the unobserved samples



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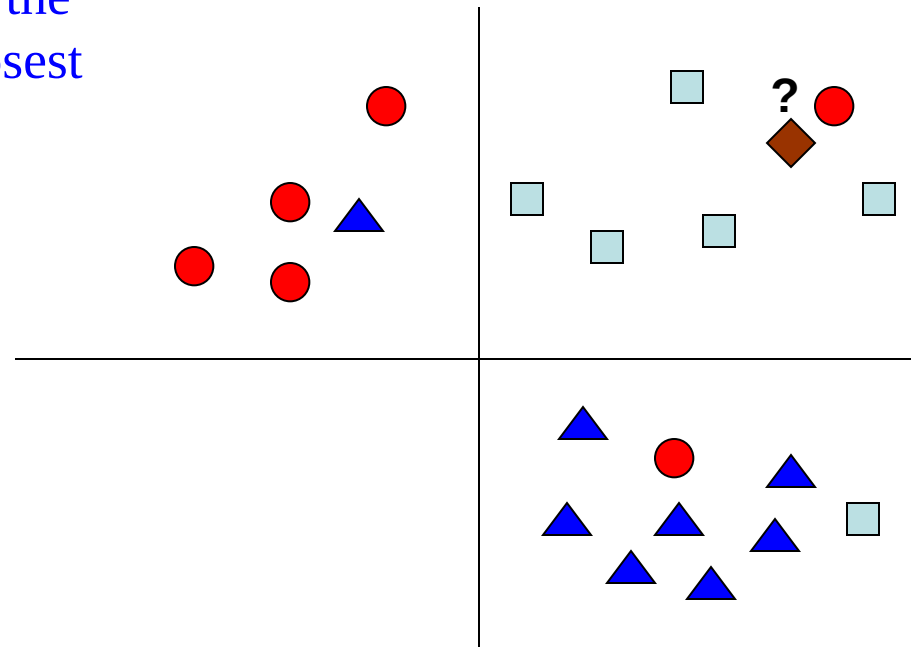
K nearest neighbors

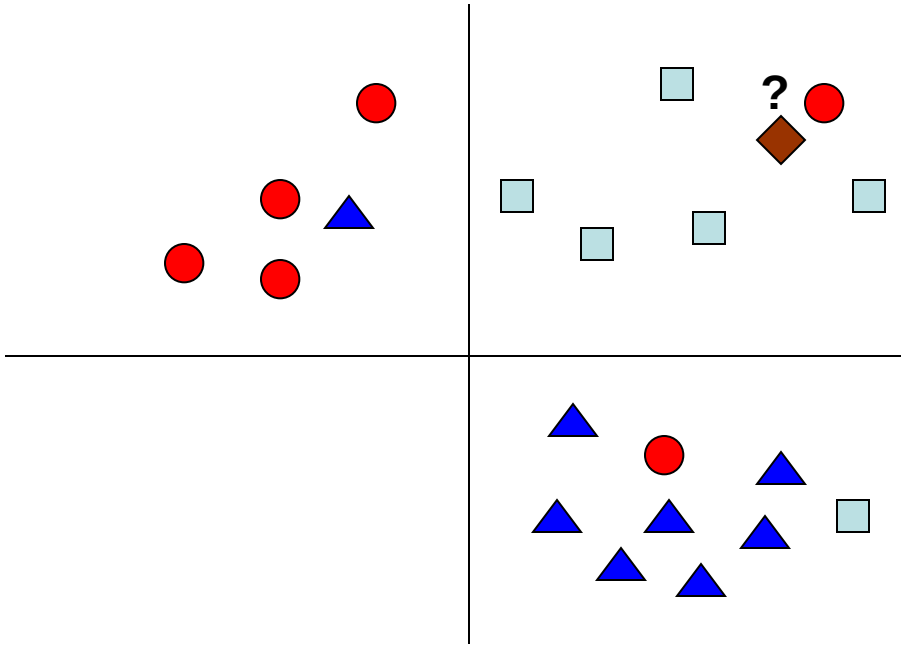
K nearest neighbors (KNN)

- A simple, yet surprisingly efficient algorithm
- Requires the definition of a distance function or similarity measures between samples
- Select the class based on the majority vote in the k closest points



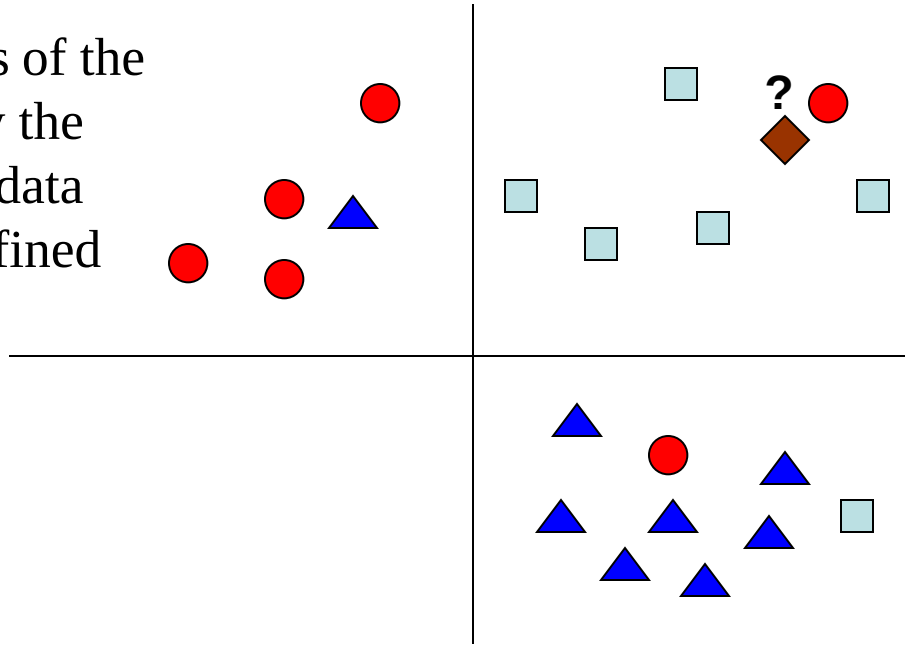
K nearest neighbors (KNN)

- Need to determine an appropriates value for k
- What happens if we chose k=1?
- What if k=3?

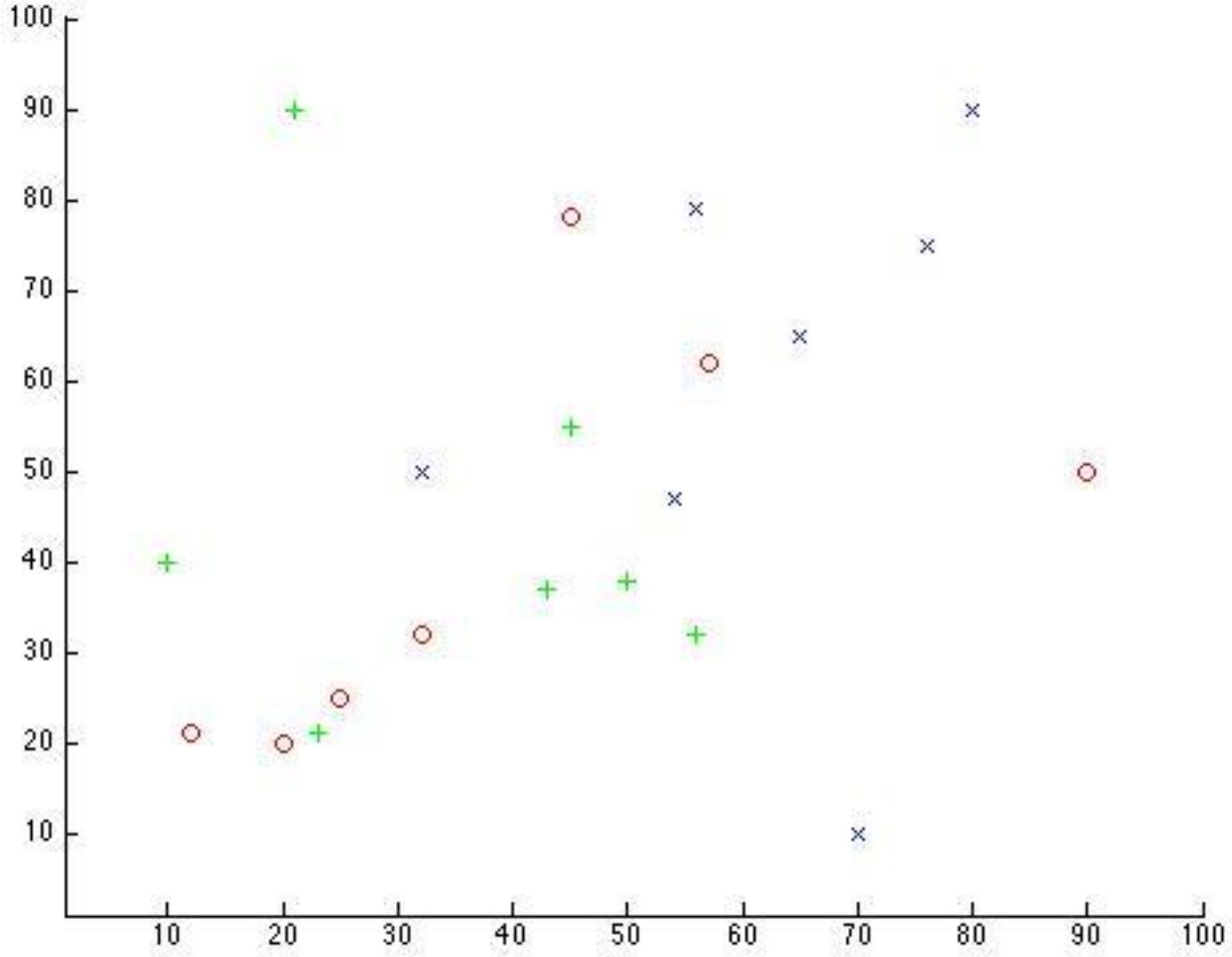


K nearest neighbors (KNN)

- Choice of k influences the 'smoothness' of the resulting classifier
- In that sense it is similar to a kernel methods (discussed later in the course)
- However, the smoothness of the function is determined by the actual distribution of the data (p(x)) and not by a predefined parameter.



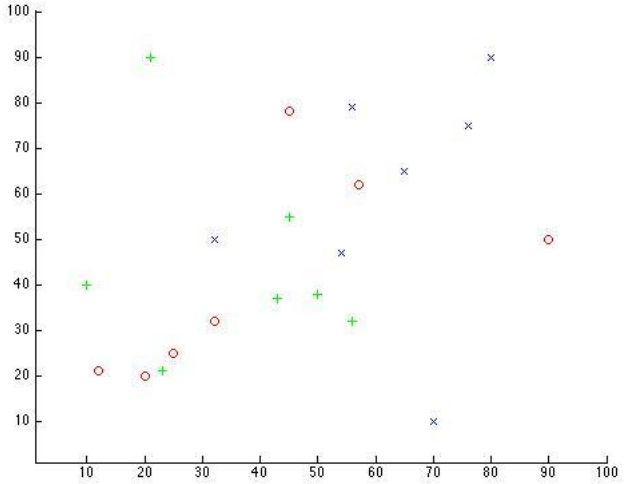
The effect of increasing k



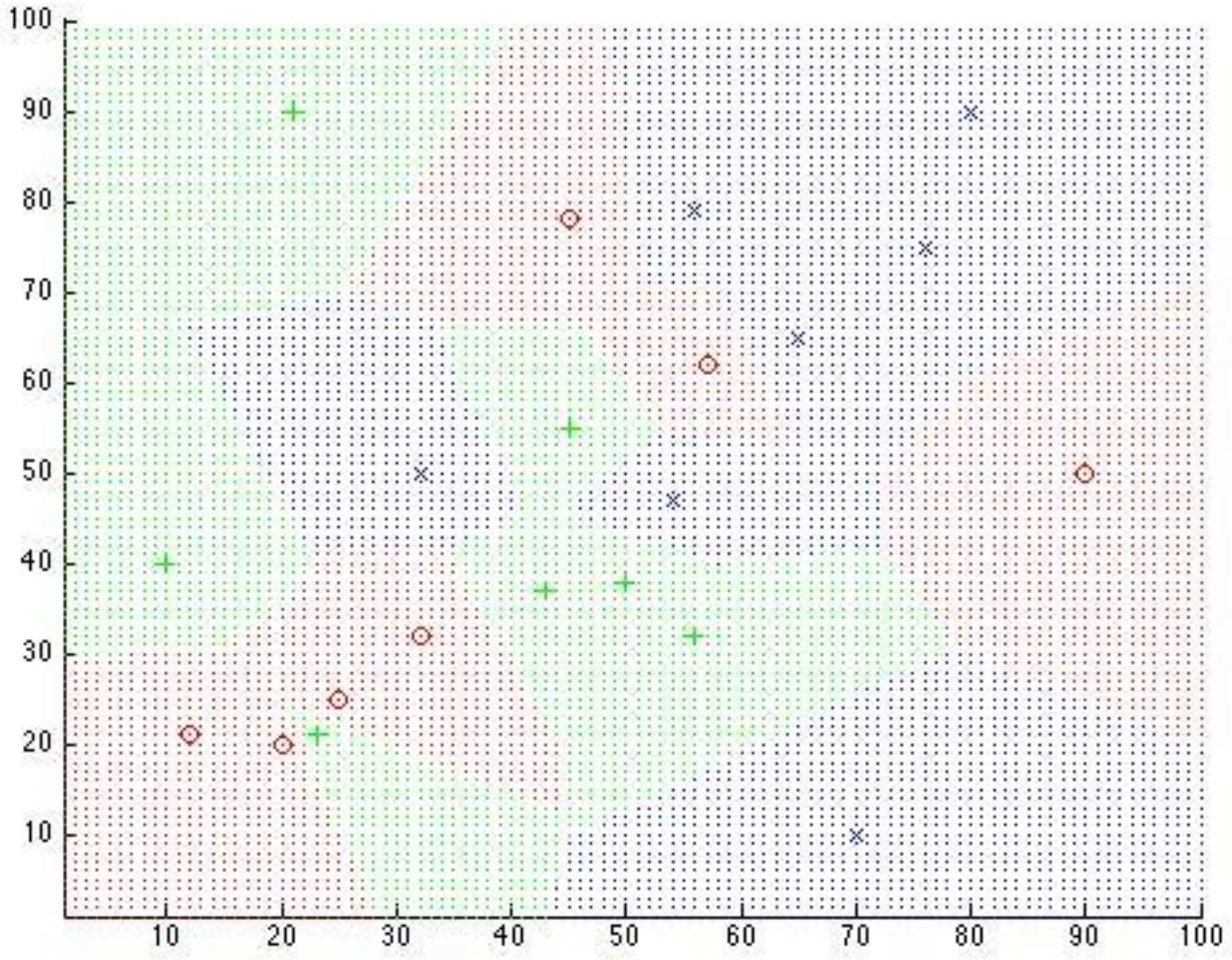
The effect of increasing k

We will be using Euclidian distance to determine what the k nearest neighbors are:

$$d(x,x') = \sqrt{\sum_i (x_i - x_i')^2}$$



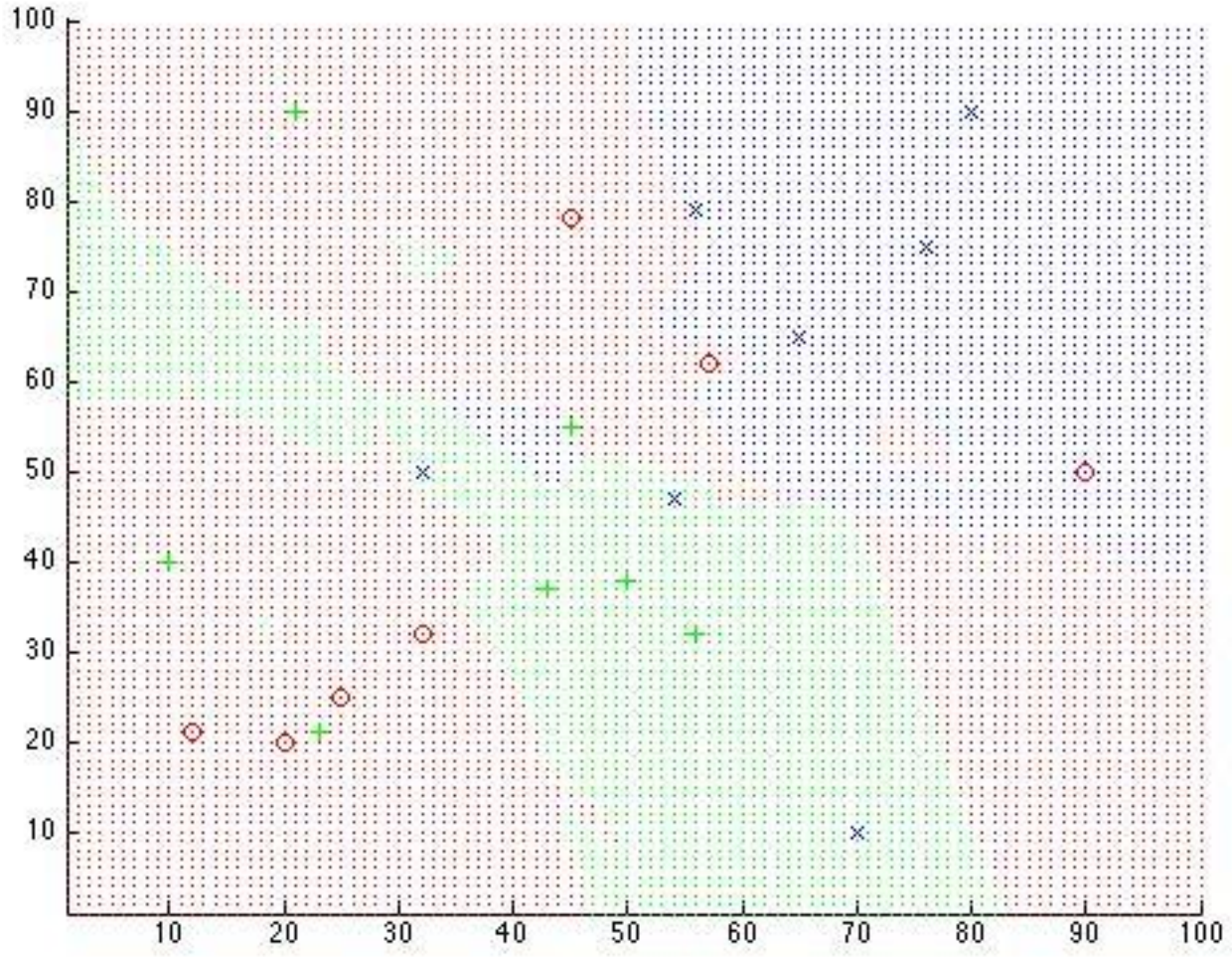
KNN with k=1



KNN with k=3

Ties are broken using the order:

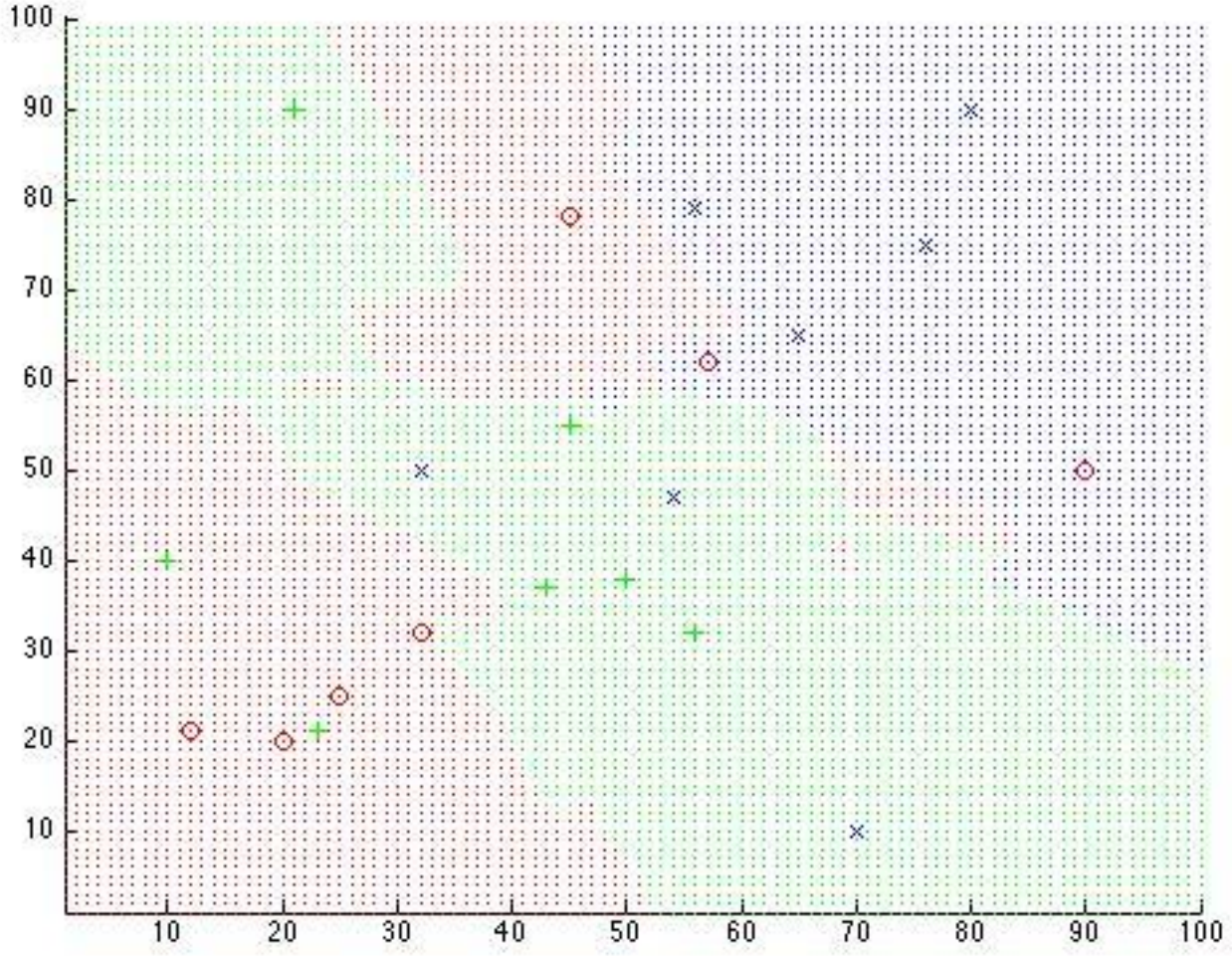
Red , Green, Blue



KNN with k=5

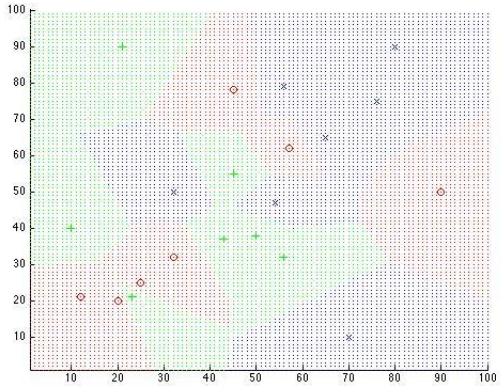
Ties are broken using the order:

Red , Green, Blue

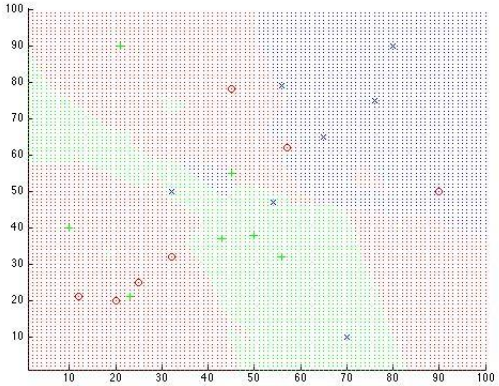


Comparisons of different k's

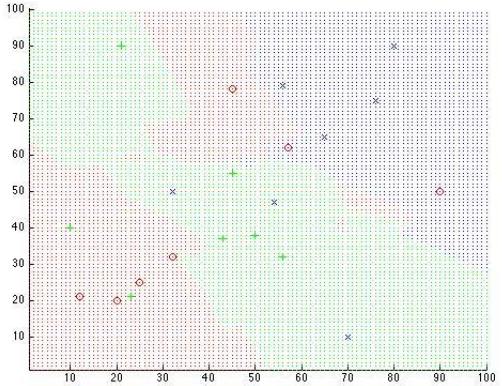
K = 1



K = 3



K = 5



A probabilistic interpretation of KNN

- The decision rule of KNN can be viewed using a probabilistic interpretation
- What KNN is trying to do is approximate the Bayes decision rule on a subset of the data
- To do that we need to compute certain properties including the conditional probability of the data given the class ($p(x|y)$), the prior probability of each class ($p(y)$) and the marginal probability of the data ($p(x)$)
- These properties would be computed for some small region around our sample and the size of that region will be *dependent on the distribution of the test samples*

Computing probabilities for KNN

- Let V be the volume of the m dimensional ball around x containing the k nearest neighbors for x (where m is the number of features in x).
- Then we can write

$$p(x)V = P = \frac{K}{N} \qquad p(x) = \frac{K}{NV} \qquad p(x \mid y=1) = \frac{K_1}{N_1 V} \qquad p(y=1) = \frac{N_1}{N}$$

- Using Bayes rule we get:

$$p(y=1 \mid x) = \frac{p(x \mid y=1)p(y=1)}{p(x)} = \frac{K_1}{K}$$

P - probability of points in region

N - total number of samples

V - volume of selected ball

K - number of nearest neighbors

N_1 - total number of samples from class 1

K_1 - number of samples from class 1 in K

Computing probabilities for KNN

N - total number of samples

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- Using Bayes rule we get:

$$p(y=1 \mid x) = \frac{p(x \mid y=1)p(y=1)}{p(x)} = \frac{K_1}{K}$$

Using Bayes decision rule we will chose the class with the highest probability, which in this case is the class with the highest number of samples in K

Important points

- Optimal decision using Bayes rule
- Types of classifiers
- Effect of values of k on knn classifiers
- Probabilistic interpretation of knn