

ECE 295: Lecture 05 Supervised Learning

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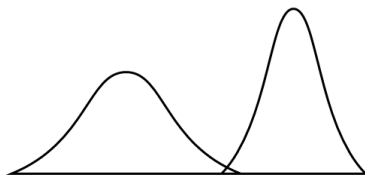


Motivation

Escalator Problem:

- ▶ You study the escalator problem for two airports
- ▶ Repeat the measurements for N days
- ▶ You have two distributions of the sample means
- ▶ I give you a new data point
- ▶ Which class does it belong to?

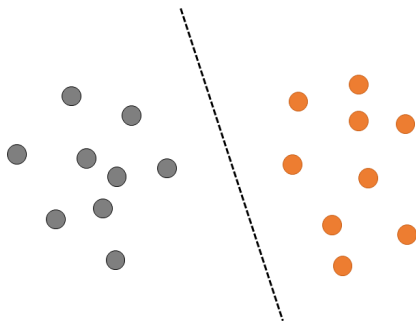
	Indianapolis	Chicago O'Hare
Day 1	$\bar{X}_1 = 10$	$\bar{Y}_1 = 100$
Day 2	$\bar{X}_2 = 11$	$\bar{Y}_2 = 98$
Day 3	$\bar{X}_3 = 10$	$\bar{Y}_3 = 99$
⋮	⋮	⋮
Day N	$\bar{X}_N = 12$	$\bar{Y}_N = 103$



Supervised Learning

Why call **supervised** learning?

- ▶ Labeled ground truth available
- ▶ Build a classifier based on the training data
- ▶ Given a new data point, tell which class does it belong to
- ▶ Can be high-dimensional data points



Supervised Learning Methods

We will talk about two methods

- Naive Bayes**
- ▶ Requires a model, e.g., Gaussian.
 - ▶ Do classification by estimating the likelihood.
 - ▶ High training cost (depending on choice model).
 - ▶ Low testing cost. Likelihood is usually not expensive.

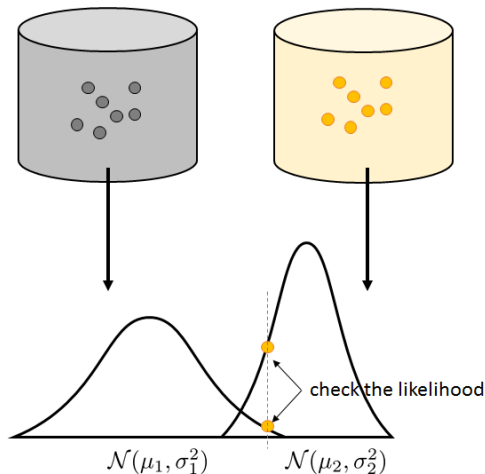
- K Nearest Neighbor**
- ▶ Does not require a model.
 - ▶ Do classification by measuring distance.
 - ▶ No training cost.
 - ▶ High testing cost. You need to measure distance for every testing data point.

There are other methods:

- ▶ Support Vector Machine
- ▶ Neural Networks

Naive Bayes

- ▶ Pick a model. Let's say Gaussian.
- ▶ From the data, estimate the parameters. For Gaussian, estimate μ and σ



Naive Bayes

Recall **Bayes Theorem**: The **posterior** distribution is

$$f_{C|\mathbf{X}}(c | \mathbf{x}) = \frac{f_{\mathbf{X}|C}(\mathbf{x} | c)f_C(c)}{f_{\mathbf{X}}(\mathbf{x})}. \quad (1)$$

- ▶ $f_{\mathbf{X}|C}(\mathbf{x} | C)$: The likelihood of having $\mathbf{X} = \mathbf{x}$ given class $C = c$.
- ▶ $f_C(c)$: The probability of getting class $C = c$.

The Naive Bayes is also called the Maximum-a-Posteriori (MAP) decision. In a two-class classification problem, MAP states that

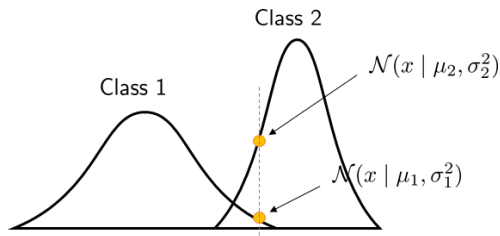
$$f_{C|\mathbf{X}}(1 | \mathbf{x}) \underset{\text{class 0}}{\geq} \underset{\text{class 1}}{1} f_{C|\mathbf{X}}(0 | \mathbf{x}). \quad (2)$$

Example: Single-variable Gaussian

Recall Gaussian:

$$f_{X|C}(x|0) = \frac{1}{\sqrt{2\pi\sigma_0^2}} \exp\left\{-\frac{(x-\mu_0)^2}{2\sigma_0^2}\right\} \stackrel{\text{def}}{=} \mathcal{N}(x|\mu_0, \sigma_0^2)$$

$$f_{X|C}(x|1) = \frac{1}{\sqrt{2\pi\sigma_1^2}} \exp\left\{-\frac{(x-\mu_1)^2}{2\sigma_1^2}\right\} \stackrel{\text{def}}{=} \mathcal{N}(x|\mu_1, \sigma_1^2)$$



Example: Single-variable Gaussian

- ▶ Given a testing data point x
- ▶ Compute likelihoods $\mathcal{N}(x | \mu_1, \sigma_1^2)$ and $\mathcal{N}(x | \mu_0, \sigma_0^2)$
- ▶ The MAP decision rule is

$$\frac{f_{X|C}(x | 1)f_C(1)}{f_X(x)} \underset{\text{class 0}}{\overset{\text{class 1}}{\geq}} \frac{f_{X|C}(x | 0)f_C(0)}{f_X(x)}$$

We can cancel out the denominators:

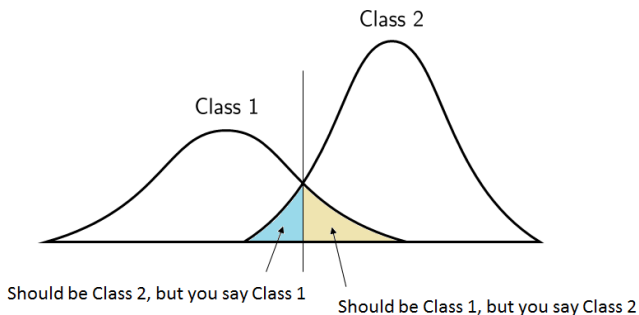
$$f_{X|C}(x | 1)f_C(1) \underset{\text{class 0}}{\overset{\text{class 1}}{\geq}} f_{X|C}(x | 0)f_C(0)$$

Write out the terms explicitly:

$$\mathcal{N}(x | \mu_1, \sigma_1^2)\mathbb{P}[\text{Class 1}] \underset{\text{class 0}}{\overset{\text{class 1}}{\geq}} \mathcal{N}(x | \mu_0, \sigma_0^2)\mathbb{P}[\text{Class 0}]$$

Naive Bayes

- ▶ $\mathbb{P}[\text{Class 1}]$ is the probability that Class 1 shows up
- ▶ $\mathbb{P}[\text{Class 1}]$ usually requires some prior knowledge
- ▶ Naive Bayes tells you “soft-decisions”
- ▶ They are the probabilities that x should belong to Class 1 or 2.
- ▶ Cut off appears when the two Gaussian intersects
- ▶ There are two types of error



Multi-Dimensional Gaussian

What if the data is high-dimensional?

Definition (High-dimensional Gaussian)

A d -dimensional **Gaussian** has a PDF

$$\mathcal{N}(\mathbf{x} \mid \boldsymbol{\mu}, \boldsymbol{\Sigma}) = \frac{1}{\sqrt{(2\pi)^d |\boldsymbol{\Sigma}|}} \exp \left\{ -\frac{1}{2} (\mathbf{x} - \boldsymbol{\mu})^T \boldsymbol{\Sigma}^{-1} (\mathbf{x} - \boldsymbol{\mu}) \right\}.$$

- ▶ $\boldsymbol{\mu} \in \mathbb{R}^d$ is the mean vector
- ▶ $\boldsymbol{\Sigma} \in \mathbb{R}^{d \times d}$ is the covariance matrix

The mean vector and the covariance matrix can be computed using commands

mean, cov

Caution: Be careful about the transpose of the data matrix.

Classification

- ▶ Given a testing dataset $\mathbf{y}_1, \dots, \mathbf{y}_N$.
- ▶ Assume $\mathbb{P}[\text{Class 1}] = \mathbb{P}[\text{Class 0}] = \frac{1}{2}$.

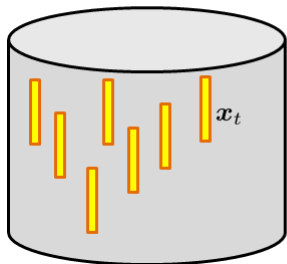
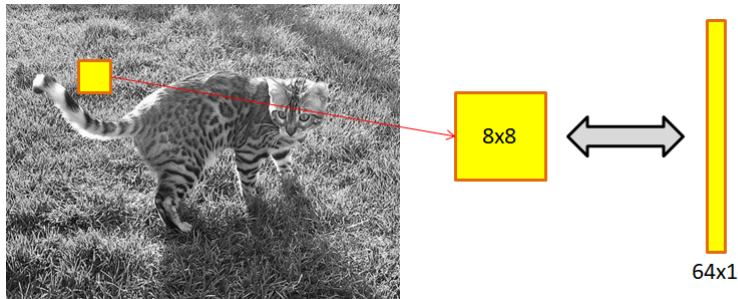
The Naive Bayes (i.e., the MAP) decision is

$$\mathcal{N}(\mathbf{y}_i | \boldsymbol{\mu}_1, \boldsymbol{\Sigma}_1) \underset{\text{class 0}}{\overset{\text{class 1}}{\geq}} \mathcal{N}(\mathbf{y}_i | \boldsymbol{\mu}_0, \boldsymbol{\Sigma}_0),$$

This is equivalent to

$$\frac{1}{\sqrt{(2\pi)^d |\boldsymbol{\Sigma}_1|}} \exp \left\{ -\frac{1}{2} (\mathbf{x} - \boldsymbol{\mu}_1)^T \boldsymbol{\Sigma}_1^{-1} (\mathbf{x} - \boldsymbol{\mu}_1) \right\} \\ \underset{\text{class 0}}{\overset{\text{class 1}}{\geq}} \frac{1}{\sqrt{(2\pi)^d |\boldsymbol{\Sigma}_0|}} \exp \left\{ -\frac{1}{2} (\mathbf{x} - \boldsymbol{\mu}_0)^T \boldsymbol{\Sigma}_0^{-1} (\mathbf{x} - \boldsymbol{\mu}_0) \right\} \quad (3)$$

Homework 5



$$\boldsymbol{\mu} = \frac{1}{T} \sum_{t=1}^T \mathbf{x}_t \quad \in \mathbb{R}^{64 \times 1}$$

$$\boldsymbol{\Sigma} = \frac{1}{T} \sum_{t=1}^T (\mathbf{x}_t - \boldsymbol{\mu})(\mathbf{x}_t - \boldsymbol{\mu})^T \quad \in \mathbb{R}^{64 \times 64}$$

Naive Bayes

Pros:

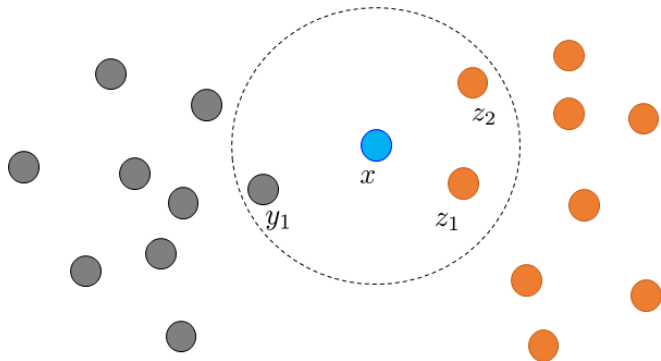
- ▶ You have a model
- ▶ More interpretable
- ▶ Usually cheap to compute the likelihood
- ▶ Robust against outliers
- ▶ Good for missing data

Cons:

- ▶ You need to choose a model
- ▶ Your model may not work — It may not describe the data accurately
- ▶ Decision boundary could be over-simplified
- ▶ You need to have prior knowledge

k-Nearest Neighbor

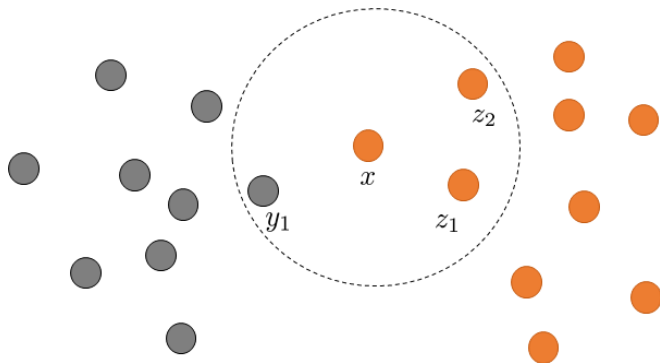
- ▶ Start with two labeled clusters
- ▶ Give me a new data point x
- ▶ Draw a circle around x



k-NN

k-Nearest Neighbor

- ▶ Grow the circle until you find k data points, e.g., $k = 3$
- ▶ Count how many are in Class 1 and Class 2
- ▶ If Class 1 is more, then assign x to Class 1



Distance

- ▶ For 1D data points, we can set

$$D(x, y) = (x - y)^2, \quad \text{or} \quad D(x, y) = |x - y|$$

- ▶ For high dimensional data points, we can set

$$D(\mathbf{x}, \mathbf{y}) = \sum_{i=1}^d (x_i - y_i)^2, \quad \text{or} \quad D(\mathbf{x}, \mathbf{y}) = \sum_{i=1}^d |x_i - y_i|$$

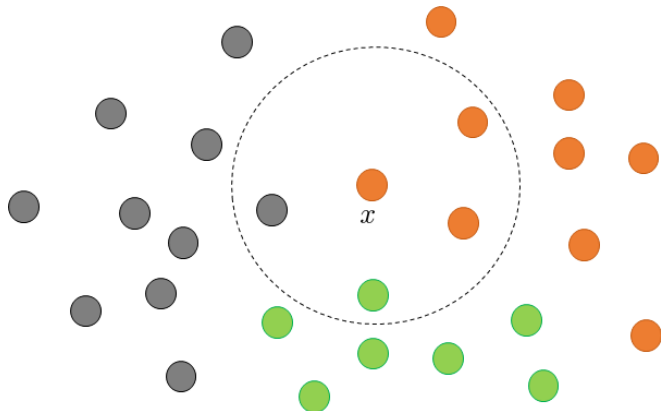
How large K should be?

- ▶ Depends on your problem
- ▶ Small datasets, k should be small

k-NN

Multiple Classes

- ▶ Say there are M classes
- ▶ Then k cannot be a multiple of M
- ▶ Otherwise there will be tie



kNN

Pros:

- ▶ No need to have a model
- ▶ Can have arbitrary decision boundary
- ▶ Could be efficient if dataset is small

Cons:

- ▶ Very expensive if dataset is large
- ▶ Not as interpretable as Naive Bayes
- ▶ Need to tune k
- ▶ Doesn't handle missing data

Summary

- ▶ Supervised learning: Ground truth available
- ▶ Two methods in this lecture
- ▶ Naive Bayes: Require a model, but fast and interpretable
- ▶ kNN: Does not require a model, but slow
- ▶ There are many other supervised learning methods