

# Section 8: Classification with R

STA 141A – Fundamentals of Statistical Data Science

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Based on Chapter 4 of ISL book James et al. (2021).

- For more R code examples, see R Markdown files in <https://www.statlearning.com/resources-second-edition>

**1** Why not linear regression?

**2** Logistic regression

- Binary classification
- Multinomial logistic regression

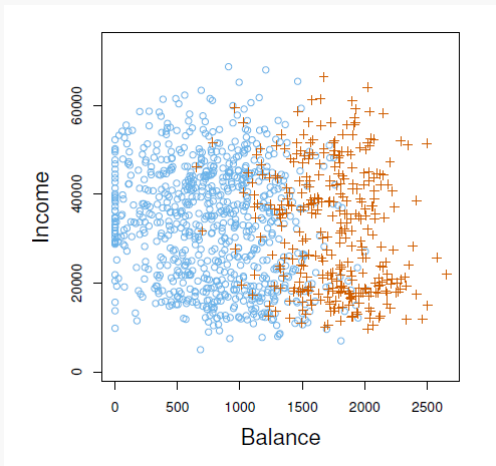
**3** Alternatives to logistic regression

- Linear discriminant analysis for  $p = 1$
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## Example (two categories)



**Figure 1:** Image by James et al. (2021), based on the Default data set in R. The annual incomes and monthly credit card balances of a number of individuals, where the individuals who defaulted on their credit card payments are shown in orange, and those who did not are shown in blue.

What are the predictors and responses in each example?

1. A person arrives at the emergency room with a set of symptoms that could possibly be attributed to one of three medical conditions. Which of these medical conditions does the person have based on the symptoms given?
2. An online banking service must be able to determine whether or not a transaction being performed on the site is fraudulent, on the basis of the user's IP address, past transaction history, and so forth.
3. On the basis of DNA sequence data for a number of patients with and without a given disease, one would like to figure out which DNA mutations are disease-causing and which are not.

**Classification:** the task of predicting *qualitative/categorical* responses

- Each response  $y_i$  is one of finitely many predetermined categories.
- **Classifying** an observation: assigning/predicting that observation to a certain category/class.
- In contrast, regression deals with “continuous” numeric response values.

As in regression, in the classification setting

- We have a set of training observations  $(x_1, y_1), \dots, (x_n, y_n)$  that we can use to build a classifier.
- We want our classifier to perform well not only on the training data, but also on test observations that were not used to train the classifier.

**Why not linear regression?**

## No natural ordering

In example 1 above, a person arrives at the emergency room with a set of symptoms. We would like to treat the person based on three reasonable medical conditions:

Appendicitis, Food poisoning, Gastritis.

- We could code each medical condition  $Y$  as:

$$Y = \begin{cases} 1, & \text{if Appendicitis,} \\ 2, & \text{if Food poisoning,} \\ 3, & \text{if Gastritis.} \end{cases}$$

This coding implies an ordering on the outcomes, insisting that the difference between Appendicitis and Food poisoning is the same as the difference between Food poisoning and Gastritis.

- We could also code:

$$Y = \begin{cases} 1, & \text{if Gastritis,} \\ 2, & \text{if Appendicitis,} \\ 3, & \text{if Food poisoning.} \end{cases}$$

Equally reasonable, but would lead to very different predictions on test observations.

What if categories had a natural ordering, such as **mild**, **moderate**, and **severe**?

- Issue: the distance between **ordinal** categories is generally unknown.
- In general there is no natural way to convert a qualitative response variable with **more than two levels** into a quantitative response that is ready for linear regression.



Can we use linear regression for a **binary** (two levels) response?

- In the Default data set, the two response values can be coded as

$$Y = \begin{cases} 1, & \text{if Default,} \\ 0, & \text{if Not default.} \end{cases}$$

- We could then fit a linear regression to this binary response:

$$\hat{Y} = \hat{\beta}_0 + \hat{\beta}_1 \times \text{Balance} + \hat{\beta}_2 \times \text{Income}$$

and then predict **Default** if  $\hat{Y} > 0.5$  and **Not default** otherwise.

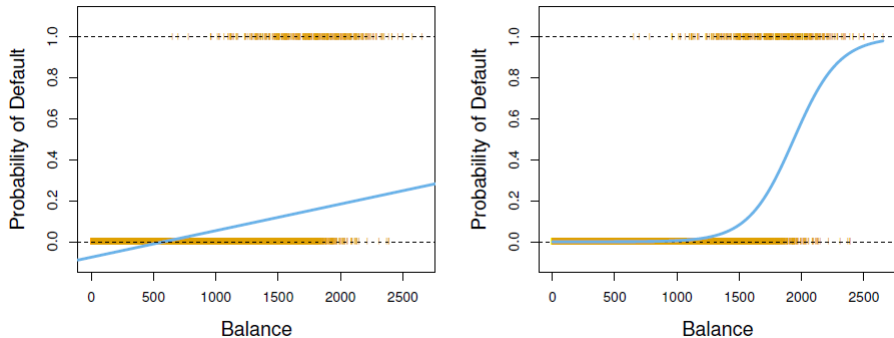
- What if we also want to estimate e.g.,

$$P(\text{Default} \mid \text{Balance} = 4000, \text{Income} = 80000)$$

i.e., the probability of defaulting given certain values of **Balance** and **Income**?

- Issue:  $\hat{Y}$  can be smaller than zero or larger than one.

## Only two levels



**Figure 2:** Image by James et al. (2021), based on the Default data set in R. Left: The estimated probability of default using *linear regression*, where the orange ticks indicate the values "0" for No, and "1" for Yes. Right: Predicted probabilities of default using *logistic regression*, where all probabilities lie between 0 and 1.

■ Let's explore logistic regression.

## Log odds

Let  $P(A)$  be the *probability* that event  $A$  occurs. Then  $P(A) \in [0, 1]$ . Map to  $(-\infty, \infty)$ ?

- The *log odds* of  $A$  occurring is defined as

$$\log \left( \frac{P(A)}{1 - P(A)} \right) \quad (1)$$

which can be a value in  $\mathbb{R}$ . (For this course, assume  $\log$  is the natural logarithm, i.e.,  $\log$  with base  $e$ .) We will also write (1) as *logit*( $P(A)$ ).

# Logistic regression

# Logistic regression

## Binary classification

## Binary classification

Each response belongs to one of two classes, coded as 0 and 1 (e.g., **No** and **Yes**).

- Classification: compute/estimate conditional prob.  $P(Y = k|X)$  for each class  $k$ .
- If only two classes, we only need  $P(Y = 1|X)$ . (Why?)

Suppose we have computed  $P(Y = 1|X)$  for a given value of predictor  $X$ . What class should be assigned to  $X$ ?

- A default decision rule for predictor value  $X$  is to assign:

$$\begin{cases} 1 & \text{if } P(Y = 1|X) > 0.5; \\ 0 & \text{if } P(Y = 1|X) \leq 0.5. \end{cases}$$

- Consequences might not be symmetric. E.g., in court, is it worse give a **guilty** verdict to an innocent person, or give a **not guilty** verdict to guilty person? May want to change the decision rule to assign:

$$\begin{cases} \text{guilty} & \text{if } P(Y = 1|X) > 0.8; \\ \text{not guilty} & \text{if } P(Y = 1|X) \leq 0.8. \end{cases}$$

# Logistic regression

*Logistic regression* models the conditional probability  $P(Y = 1|X)$ .

- Convert  $p(X) = P(Y = 1|X)$  to log odds, then use linear regression on log odds:

$$\text{logit}(p(X)) = \beta_0 + \beta_1 X_1 + \dots + \beta_p X_p.$$

- The conditional probabilities are then

$$p(X) = \frac{e^{\beta_0 + \beta_1 X_1 + \dots + \beta_p X_p}}{1 + e^{\beta_0 + \beta_1 X_1 + \dots + \beta_p X_p}}. \quad (2)$$

Interpretation:

- Increasing  $X_1$  by one unit changes  $p(X)$  by

$$p(X_1 + 1, X_2, X_3, \dots, X_p) - p(X_1, X_2, X_3, \dots, X_p)$$

which depends on all  $p - 1$  coefficient values and the current predictor values.

- Increasing  $X_1$  by one unit changes the log odds  $\text{logit}(p(X))$  by

$$\text{logit}(p(X_1 + 1, X_2, X_3, \dots, X_p)) - \text{logit}(p(X_1, X_2, X_3, \dots, X_p))$$

which is just  $\beta_1$ .

Estimating the regression coefficients:

- Usually use the method of *maximum likelihood*.
- Details outside scope of this class; we will just use R to compute these estimates.



## Making predictions from estimators $\hat{\beta}_0, \hat{\beta}_1, \dots, \hat{\beta}_p$

To assign 0 or 1 to  $X$ , we can use estimated log odds or estimated  $p(X)$ .

- We can estimate log odds at  $X$  by

$$\hat{\beta}_0 + \hat{\beta}_1 X_1 + \dots + \hat{\beta}_p X_p. \quad (3)$$

- Alternatively, we can estimate the conditional probability  $p(X) = P(Y = 1|X)$  by

$$\hat{p}(X) := \frac{e^{\hat{\beta}_0 + \hat{\beta}_1 X_1 + \dots + \hat{\beta}_p X_p}}{1 + e^{\hat{\beta}_0 + \hat{\beta}_1 X_1 + \dots + \hat{\beta}_p X_p}}.$$

- If this estimated probability is over a certain pre-defined threshold (e.g. 0.25), then we would assign  $X = x$  to the class 1.

### Example

If  $\hat{\beta}_0 = -9.9$  and  $\hat{\beta}_1 = 0.005$ , we predict the probabilities of **default** for individuals with balance  $X = \$1,000$  and  $X = \$2,000$  by

$$\hat{p}(X = 1,000) = \frac{e^{\hat{\beta}_0 + \hat{\beta}_1 X}}{1 + e^{\hat{\beta}_0 + \hat{\beta}_1 X}} = \frac{e^{-9.9 + 0.005 \cdot 1,000}}{1 + e^{-9.9 + 0.005 \cdot 1,000}} \approx 0.007,$$

$$\hat{p}(X = 2,000) = \frac{e^{\hat{\beta}_0 + \hat{\beta}_1 X}}{1 + e^{\hat{\beta}_0 + \hat{\beta}_1 X}} = \frac{e^{-9.9 + 0.005 \cdot 2,000}}{1 + e^{-9.9 + 0.005 \cdot 2,000}} \approx 0.525.$$

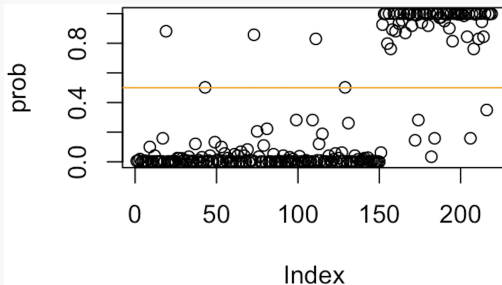
# glm()

We use `glm()` for logistic regression ('glm' stands for *general linear model*).

- Must specify which variables are used, data set, and type of response.
- Must put `family=binomial` to specify a binary response.

```
library(palmerpenguins)
# We work with only Adelie and Chinstrap species (we exclude Gentoo).
peng_binary <- na.omit(penguins[penguins$species != 'Gentoo', ])
logreg <- glm(species ~ bill_length_mm, data=peng_binary, family=binomial)
prob <- predict(logreg, type='response') # 'link' also possible
predicted <- ifelse(prob<.5, 'Adelie', 'Chinstrap')

plot(prob)
abline(a=0.5, b=0, col='orange')
```



# Logistic regression

## Multinomial logistic regression

# Multinomial logistic regression

We sometimes wish to classify *a response variable that has more than two classes*.

- Extend the two-class logistic regression approach to the setting of  $K > 2$  classes.
- We will need separate regression coefficients for each of the first  $K - 1$  classes.  
Hence, for any  $x \in \mathbb{R}^p$ , define

$$\alpha_l(x) := \beta_{l,0} + \beta_{l,1}x_1 + \cdots + \beta_{l,p}x_p \quad \text{for any } l = 1, \dots, K - 1. \quad (4)$$

E.g., consider conditions **Appendicitis**, **Food poisoning**, and **Gastritis**.

For  $j = 1, \dots, p$ , consider  $x_j = \text{Severity of symptom } j$  (e.g. how much does head hurt? how nauseated?).

- For **Appendicitis**, we want to define  $\beta_{\text{Appendicitis},j}$  for  $j = 0, 1, \dots, p$ .
- For **Food poisoning**, we want to define  $\beta_{\text{Food poisoning},j}$  for  $j = 0, 1, \dots, p$ .
- We could define similarly for **Gastritis**, but we will see that we won't need to.

# Multinomial logistic regression

For convenience, copy-and-paste Eq. (4) here:

$$\alpha_l(x) := \beta_{l,0} + \beta_{l,1}x_1 + \cdots + \beta_{l,p}x_p \quad \text{for any } l = 1, \dots, K-1.$$

**Multinomial logistic regression** model:

1. Select a class to serve as the baseline; WLOG, select the  $K$ th class for this role.
2. Replace the model (2) with the model

$$P(Y = k \mid X = x) = \begin{cases} \frac{e^{\alpha_k(x)}}{1 + \sum_{l=1}^{K-1} e^{\alpha_l(x)}} & \text{for } k = 1, \dots, K-1, \\ \frac{1}{1 + \sum_{l=1}^{K-1} e^{\alpha_l(x)}} & \text{for } k = K. \end{cases}$$

For  $k = 1, \dots, K-1$ , we have

$$\log \left( \frac{P(Y = k \mid X = x)}{P(Y = K \mid X = x)} \right) = \alpha_k(x)$$

which is linear in the predictors.

Consider classifying ER visits into **Appendicitis**, **Food poisoning**, **Gastritis**.

- Suppose we set **Appendicitis** as the baseline.
- If  $X_j$  increases by one unit, then

$$\log \left( \frac{P(Y = \text{Food poisoning} \mid X = x)}{P(Y = \text{Appendicitis} \mid X = x)} \right)$$

increases by  $\beta_{\text{Food poisoning},j}$ .

- If  $X_j$  increases by one unit, then

$$P(Y = \text{Food poisoning} \mid X = x)$$

increases by a complicated function of all  $p - 1$  coefficient values and the current predictor values.

ISLR2 textbook doesn't have code walkthrough for multinomial logistic regression, so you can find one here:

[https://www.r-bloggers.com/2020/05/  
multinomial-logistic-regression-with-r/](https://www.r-bloggers.com/2020/05/multinomial-logistic-regression-with-r/)

## Alternative coding: softmax coding

In the **softmax coding** (used extensively in some areas of machine learning), rather than selecting a baseline class, we treat all  $K$  classes symmetrically:

$$P(Y = k \mid X = x) = \frac{e^{\alpha_k(x)}}{\sum_{l=1}^K e^{\alpha_l(x)}} \quad \text{for } k = 1, \dots, K$$

- Thus, we estimate coefficients for all  $K$  classes (rather than for just  $K - 1$  classes).
- The log odds ratio between the  $k$ th and  $l$ th classes equals

$$\begin{aligned} \log \left( \frac{P(Y = k \mid X = x)}{P(Y = l \mid X = x)} \right) &= \alpha_k(x) - \alpha_l(x) \\ &= (\beta_{k,0} - \beta_{l,0}) + (\beta_{k,1} - \beta_{l,1})x_1 + \dots + (\beta_{k,p} - \beta_{l,p})x_p. \end{aligned}$$

Example interpretation: if  $X_j$  increases by one unit, then

$$\log \left( \frac{P(Y = \text{Food poisoning} \mid X = x)}{P(Y = \text{Appendicitis} \mid X = x)} \right)$$

increases by  $(\beta_{\text{Food poisoning},j} - \beta_{\text{Appendicitis},j})$ .



## **Alternatives to logistic regression**

Recall: logistic regression directly models  $P(Y = k|X = x)$  for binary responses by using the logistic function.

Pros:

- We can see impact of e.g., a unit increase in  $X_j$  on log odds.
- Assumptions are relatively loose: independence of errors, a linear relationship between the logit and independent variables, and absence of severe multicollinearity among independent variables.

Some issues:

- When there are big differences between two classes, the parameter estimates in the logistic regression model are unstable.
- If the sample size is small, other approaches can be more accurate than logistic regression.

## Using Bayes' theorem

We can instead use *Bayes' theorem* to obtain the *posterior* probability  $P(Y = k \mid X = x)$ :

$$p_k(x) := P(Y = k \mid X = x) = \frac{\pi_k f_k(x)}{\sum_{\ell=1}^K \pi_\ell f_\ell(x)}, \quad (5)$$

- $\pi_k := P(Y = k)$  is the overall or *prior* probability that a randomly chosen observation comes from the  $k$ th class.
  - ▶ Here  $\pi$  is just a variable name — not the same as  $\pi = 3.14159 \dots$
- $f_k$  is the predictor's PMF/PDF given that the response is from the  $k$ th class.
  - ▶ PMF case:  $f_k(x) = P(X = x \mid Y = k)$ .

We typically don't know either  $\pi_k$  or  $f_k$ .

- Can estimate  $\pi_k$  by the proportion of observed elements in the  $k$ th class.
  - ▶ E.g. if there are 3, 2, 5 elements in the classes 1, 2, 3, respectively, then the estimated probabilities are  $\hat{\pi}_1 = \frac{3}{10}$ ,  $\hat{\pi}_2 = \frac{2}{10}$ ,  $\hat{\pi}_3 = \frac{5}{10}$ .
- Estimating  $f_k$  is more challenging — typically requires a huge amount of data unless *strong simplifying assumptions* are made.
  - ▶ Bias-variance tradeoff: the assumptions will reduce variance (hopefully by a lot) at the cost of introducing some bias.
  - ▶ Two approaches will be discussed in the next few slides.

# Alternatives to logistic regression

Linear discriminant analysis for  $p = 1$

## Assumptions of $f_k$ in LDA for $p = 1$

The *linear discriminant analysis (LDA)* approach estimates  $f_k$  by assuming:

1.  $f_k$  is a normal/Gaussian PDF, i.e. for all  $x$  holds

$$f_k(x) = \frac{1}{\sqrt{2\pi}\sigma_k} \exp \left\{ -\frac{1}{2\sigma_k^2} (x - \mu_k)^2 \right\}. \quad (6)$$

- ▶ This class will focus on  $p = 1$ .
  - ▶ But if  $p > 1$ , replace  $\mu_k$  with  $p$ -tuple and replace  $\sigma_k$  with  $p \times p$  covariance matrix  $\Sigma_k$ .
2. Same variance parameter across all  $K$  classes:  $\sigma_1^2 = \sigma_2^2 = \dots = \sigma_K^2 = \sigma^2$ .
    - ▶ The *quadratic discriminant analysis (QDA)* relaxes this *equivariance* assumption, but we won't discuss QDA in this class.

## Bayes decision boundary

With these assumptions, we plug the PDF (6) into the posterior probability (5) to get

$$p_k(x) = \frac{\pi_k \frac{1}{\sqrt{2\pi}\sigma} \exp\left\{-\frac{1}{2\sigma^2}(x - \mu_k)^2\right\}}{\sum_{\ell=1}^K \pi_\ell \frac{1}{\sqrt{2\pi}\sigma} \exp\left\{-\frac{1}{2\sigma^2}(x - \mu_\ell)^2\right\}}. \quad (7)$$

- Recall: **Bayes classifier** assigns observation with predictor  $x$  to class

$$\arg \max_{k \in \{1, 2, \dots, K\}} p_k(x)$$

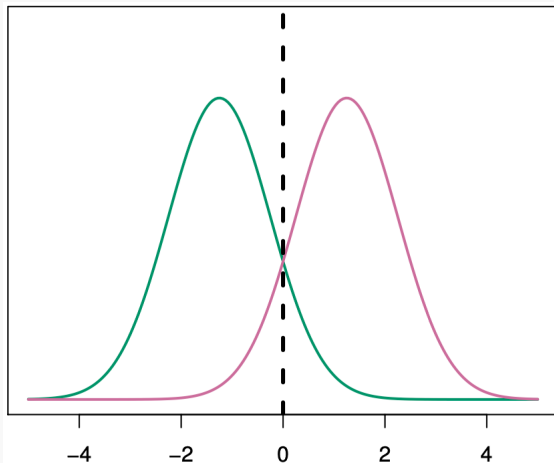
- This class is equivalent (take the log of (7)) to the class  $\arg \max_{k \in \{1, 2, \dots, K\}} \delta_k(x)$  where

$$\delta_k(x) := x \cdot \frac{\mu_k}{\sigma^2} - \frac{\mu_k^2}{2\sigma^2} + \log(\pi_k). \quad (8)$$

Here  $\delta_k$  is called a **discriminant function**.

- If  $K = 2$ , classifier assigns  $x$  to class 1 if  $\delta_1(x) > \delta_2(x)$ , to class 2 otherwise.
- If  $p = 1$ , the **Bayes decision boundary** is the value  $x$  for which  $\delta_1(x) = \delta_2(x)$ 
  - What does this inequality  $\delta_1(x) > \delta_2(x)$  and boundary simplify to if  $\pi_1 = \pi_2$ ?

## Example for decision boundary



**Figure 3:** Image by James et al. (2021). Two PDFs of normal distributions with means  $\mu_1 = -1.25$  and  $\mu_2 = 1.25$ , respectively, and variance  $\sigma^2 = 1$ . The dashed vertical line represents the Bayes decision boundary, so we assign the observation to class 1 if  $x$  is left of the line, and to class 2 otherwise.

## The LDA method for $p = 1$

In the plot above, we can calculate the Bayes classifier because we know values for all parameters  $\pi_1, \dots, \pi_K, \mu_1, \dots, \mu_K, \sigma^2$ .

- In practice, we must estimate these parameters to apply the Bayes classifier.
- Consider the estimates

$$\hat{\pi}_k = \frac{n_k}{n}, \quad \hat{\mu}_k = \frac{1}{n_k} \sum_{i:y_i=k} x_i, \quad \hat{\sigma}^2 = \frac{1}{n-K} \sum_{k=1}^K \sum_{i:y_i=k} (x_i - \hat{\mu}_k)^2,$$

- ▶  $\hat{\pi}_k$  — proportion of all training observations from  $k$ th class.
- ▶  $\hat{\mu}_k$  — average of all training observations from  $k$ th class.
- ▶  $\hat{\sigma}^2$  — weighted average of sample variances for each class.

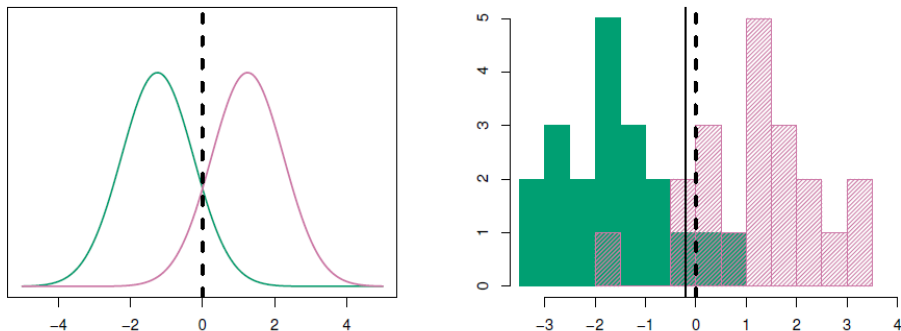
- Plug these estimates into (8) to get the **LDA discriminant function**

$$\hat{\delta}_k(x) = x \cdot \frac{\hat{\mu}_k}{\hat{\sigma}^2} - \frac{\hat{\mu}_k^2}{2\hat{\sigma}^2} + \log(\hat{\pi}_k), \quad (9)$$

which is linear in  $x$ , hence the “linear” in LDA.



## Example for decision boundary



**Figure 4:** Image by James et al. (2021). Left: Two PDFs of normal distributions with means  $\mu_1 = -1.25$  and  $\mu_2 = 1.25$ , respectively, and variance  $\sigma^2 = 1$ . The dashed vertical line represents the Bayes decision boundary, so we assign the observation to class 1 if  $x < 0$  and class 2 otherwise. Right: 20 observations were drawn from each of the two classes, and are shown as histograms. The Bayes decision boundary is shown as a dashed vertical line, and the solid vertical line represents the LDA decision boundary estimated from the training data.

## Calculation example

For  $K = 2$  classes (class "0" and "1"), we have the five data points

$$(9, 1), (8, 0), (6, 0), (7, 1), (4, 0)$$

and want to calculate the LDA discrimination function (9) for  $k = 0$  and  $k = 1$ .

# Alternatives to logistic regression

## Naive Bayes

# Naive Bayes classifier

The *naive Bayes classifier* assumes for each class  $k = 1, \dots, K$  that:

Within the  $k$ th class, the  $p$  predictors are independent.

Mathematically, this means that for each class  $k$ :

$$f_k(x) = f_{k,1}(x_1) \times f_{k,2}(x_2) \times \dots \times f_{k,p}(x_p) \quad (10)$$

where  $f_{kj}$  is the PDF/PMF of the  $j$ th predictor for observations in the  $k$ th class.

- Plug (10) into (5) to get the posterior probability

$$p_k(x) = P(Y = k|X = x) = \frac{f_{k,1}(x_1) \cdot f_{k,2}(x_2) \cdot f_{k,p}(x_p) \cdot \pi_k}{\sum_{\ell=1}^K f_{\ell,1}(x_1) \cdot f_{\ell,2}(x_2) \cdot f_{\ell,p}(x_p) \cdot \pi_{\ell}}. \quad (11)$$

- The *naive Bayes classifier* will assign an observation with predictor  $x$  to the class that maximizes posterior probability (11).
- The independence assumption, though often unrealistic, produces decent results, especially when  $n$  is too small to effectively estimate the joint distribution  $f_k$ .

## Estimation approach

Estimating the posterior probability (11) requires estimating the univariate density functions  $f_{k,j}$  for all classes  $k = 1, \dots, K$  and all predictors  $j = 1, \dots, p$ . Some options:

- If  $X_j$  is quantitative, we can assume  $X_j|Y = k \sim \mathcal{N}(\mu_{k,j}, \sigma_{k,j}^2)$  (as in LDA) or use a nonparametric approach.
- If  $X_j$  is qualitative, we could count the proportion of training observations for the  $j$ th predictor corresponding to each class  $k$ .
  - E.g. suppose we want to predict whether a student studies more than 10 hours per week based on their major (so  $p = 1$ ). We survey 100 people:

	Math major	Art major	Poli Sci major
Study > 10 hr /wk	20	25	5
Study $\leq$ 10 hr /wk	15	25	10

We use proportions (i.e., divide each cell by column sum) to estimate the “true” PMFs  $f_{k,j}$  (each column is a PMF):

	Math major	Art major	Poli Sci major
Study > 10 hr /wk	0.6	0.55	0.35
Study $\leq$ 10 hr /wk	0.4	0.45	0.65

## Errors in classification

## Confusion matrix

In classification, observations can be assigned to the wrong class.

- In binary classification, two mistakes are: *false positives* and *false negatives*.
- Examples: not default vs default, cancer vs no cancer, spam vs not spam.
- A *confusion matrix* displays both error types.

		<i>True class</i>		
		– or Null	+ or Non-null	Total
<i>Predicted class</i>	– or Null	True Neg. (TN)	False Neg. (FN)	N*
	+ or Non-null	False Pos. (FP)	True Pos. (TP)	P*
	Total	N	P	

Name	Definition	Synonyms
False Pos. rate	FP/N	Type I error, 1–Specificity
True Pos. rate	TP/P	1–Type II error, power, sensitivity, recall
Pos. Pred. value	TP/P*	Precision, 1–false discovery proportion
Neg. Pred. value	TN/N*	

**Figure 5:** Tables by James et al. (2021). A confusion matrix compares the LDA predictions to the true default statuses for the 10,000 training observations in the Default data set, using a modified threshold value that predicts default for any individuals whose posterior default probability exceeds 20 %.

# Confusion matrix

```
# Using peng_binary and predicted from earlier slide
pb_species <- factor(peng_binary$species, levels=c('Adelie', 'Chinstrap'))
table(pb_species, predicted)
```

```
> table(pb_species, predicted)
      predicted
pb_species  Adelie Chinstrap
Adelie      141         5
Chinstrap    6        62
```

```
# pb_species line is not necessary, but what happens if we instead did:
table(peng_binary$species, predicted)
```

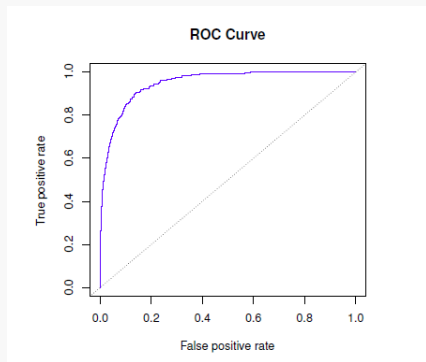


Recall the earlier “default” decision rule for binary responses: assign  $x$  to Yes if

$$P(\text{default} = \text{Yes} | X = x) > 0.5.$$

- This rule weights both types of mistakes (FN and FP) the same.
- But sometimes we care more about lowering false negatives. E.g., a credit card company trying to detect a fraudulent charge.
- Can lower the threshold from 0.5 to e.g., 0.2.
- What happens to TP rate and FP rate as threshold decreases?

The *ROC curve* simultaneously displays both types of errors for all thresholds.



**Figure 6:** Image by James et al. (2021). An *ROC curve* for LDA classifier on Default data. Dotted line represents “no information” classifier, i.e., one that doesn’t use predictors.

- ROC curve is parameterized by the possible threshold values.
- Overall performance of a classifier, summarized over all possible thresholds, is given by the *area under the ROC curve (AUC)*.
- The larger the AUC, the better the classifier.

# **Comparison of classification methods**

*Analytical* (or mathematical) comparison:

- Classifiers with a linear decision boundary are special cases of naive Bayes. Hence, LDA is a special case of naive Bayes. (This is not obvious.)
- No method uniformly dominates others: The appropriate model depends on the predictor's distribution in each class as well as  $n$  and  $p$ .
- $K$ -nearest neighbors (KNN) is a nonparametric approach. Hence, one can expect that it dominates naive Bayes and LDA when the true decision boundary is highly non-linear. However, KNN requires many observations relative to the number of predictors to perform well.

For an *empirical* (or data-based) comparison, see Section 4.5.2 of ISLR2 textbook.