

# Lecture 001

## Statistical learning: Foundations

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Admin

# Admin

## Today

### In-class

- *Course website:* <https://github.com/edrubin/EC524W20/>
- *Resources*
  - [RStudio](#) cheatsheets, books, and tutorials
  - [UO library](#)
  - See course page for more...
- Formalizing statistical learning, notation, goals (and problems)

↻ Eugene R Users Retweeted



**Ryann Crowley**

@ryann\_crowley



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Tweet; h/t: Grant McDermott

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## Upcoming

### Readings

- *Today*
  - ISL Ch1–Ch2
  - **Prediction Policy Problems** by Kleinberg *et al.* (2015)
- *Next*
  - ISL Ch. 3–4

**Problem set** Likely assigned Thursday and due Tuesday.

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- **Regression trees** extend *classification trees* to numerical outcomes (random forests extend, as well).
- **K-means clustering** partitions observations into K groups (clusters) based upon a set of variables.

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$$\mathbf{y} = f(\mathbf{x}_1, \dots, \mathbf{x}_p)$$

that accurately describes  $\mathbf{y}$  given some values of  $\mathbf{x}_1, \dots, \mathbf{x}_p$ .

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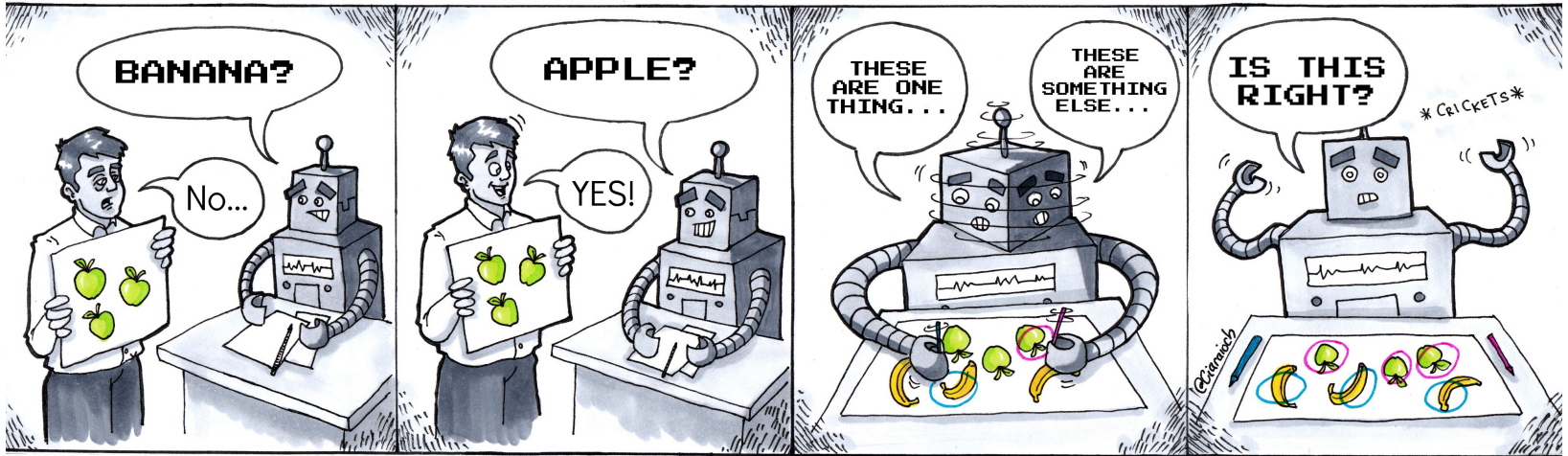
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2. **Unsupervised learning** learns relationships and structure using only **inputs** ( $x_1, \dots, x_p$ ) without any *supervising* output—letting the data "speak for itself."

**Semi-supervised learning** falls somewhere between these supervised and unsupervised learning—generally applied to supervised tasks when labeled **outputs** are incomplete.



## Supervised Learning

## Unsupervised Learning

Source

# Statistical learning

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*E.g.*, race, sex, loan default, hazard, disease, flight status
2. **Regression tasks** in which **y** takes on continuous, numeric values.  
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*Note*<sub>1</sub> The use of *regression* differs from our use of *linear regression*.

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*Note<sub>2</sub>* Don't get tricked: Not all numbers represent continuous, numerical values—*e.g.*, zip codes, industry codes, social security numbers.<sup>†</sup>

† Q Where would you put responses to 5-item Likert scales?



# Statistical learning

## *Why Learning?*

Q What puts the "learning" in statistical/machine learning?

# Statistical learning

## *Why Learning?*

**Q** What puts the "learning" in statistical/machine learning?

**A** Most learning models/algorithms will **tune model parameters** based upon the observed dataset—learning from the data.

# Notation

Our class will typically follow the notation and definitions of *ISL*.

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- $x_{i,j}$  is observation  $i$  (in  $1, \dots, n$ ) on variable  $j$  (for  $j$  in  $1, \dots, p$ )

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$$\mathbf{X} = \begin{bmatrix} x_{1,1} & x_{1,2} & \cdots & x_{1,p} \\ x_{2,1} & x_{2,2} & \cdots & x_{2,p} \\ \vdots & \vdots & \ddots & \vdots \\ x_{n,1} & x_{n,2} & \cdots & x_{n,p} \end{bmatrix}$$

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Applied to R:

- `dim(x_df)` =  $n p$
- `nrow(x_df)` =  $n$ ; `ncol(x_df)` =  $p$
- `x_df[1,]` ( $i = 1$ ); `x_df[,1]` ( $j = 1$ )

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The **outcome** for our  $i^{\text{th}}$  observation is  $y_i$ . Together the  $n$  observations form

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and our full dataset is composed of  $\left\{ (x_1, y_1), (x_2, y_2), \dots, (x_n, y_n) \right\}$ .

Back to the problem of (supervised) statistical learning...

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You'll have to wait on any real/specific answers...

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## Learning from $\hat{f}$

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Similarly, in causal-inference settings, we don't particularly care about  $\hat{\mathbf{y}}$ .

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*Irreducible* because we defined an error term  $\varepsilon$  unexplained by  $f$ .

*Note* As its name implies, you can't get rid of *irreducible* error—but we can try to get rid of *reducible* errors.



# Statistical learning

## Prediction errors

Why we're stuck with *irreducible* error

$$\begin{aligned} E\left[\{\mathbf{y} - \hat{\mathbf{y}}\}^2\right] &= E\left[\left\{f(\mathbf{X}) + \varepsilon + \hat{f}(\mathbf{X})\right\}^2\right] \\ &= \underbrace{\left[f(\mathbf{X}) - \hat{f}(\mathbf{X})\right]^2}_{\text{Reducible}} + \underbrace{\text{Var}(\varepsilon)}_{\text{Irreducible}} \end{aligned}$$

In less math:

- If  $\varepsilon$  exists, then  $\mathbf{X}$  cannot perfectly explain  $\mathbf{y}$ .
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Thus, to form our **best predictors**, we will **minimize reducible error**.

# Statistical learning

## Which type of $\hat{f}$ ?

Once you have your **inputs** ( $\mathbf{X}$ ) and **output** ( $\mathbf{y}$ ) data, you still need to decide how parametric your  $\hat{f}$  should be.<sup>†</sup>

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**Parametric methods** assume a function typically involve two steps

1. Select a functional form (shape) to represent  $f$
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**Non-parametric methods** avoid explicit assumption about the shape of  $f$ . Attempt to **flexibly fit** the data, while trying to **avoid overfitting**.

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# Statistical learning

## Which type of $\hat{f}$ ?

Methods' parametric assumptions come with tradeoffs.

### **Parametric methods**

- + Simpler to estimate and interpret.
- If assumed functional form is bad, model performance will suffer.

### **Non-parametric methods**

- + Fewer assumptions. More flexibility.
- Lower interpretability. Susceptible to overfitting. Want lots of data.

**Example:** Let's start with a pretty funky, nonlinear function.

**Truth:** The (nonlinear)  $f(\mathbf{X})$  that we hope to recover.



**The sample:**  $n = 70$  randomly drawn observations for  $\mathbf{y} = f(\mathbf{x}_1, \mathbf{x}_2) + \varepsilon$

**Estimated linear-regression model:**  $\hat{y} = \hat{\beta}_0 + \hat{\beta}_1 \mathbf{x}_1 + \hat{\beta}_2 \mathbf{x}_2 + \hat{\beta}_3 \mathbf{x}_1 \mathbf{x}_2$

**Prediction error** from our fitted linear regression model

**k-nearest neighbors** (kNN) using  $k=5$  (a *non-parametric* method)

**k-nearest neighbors** (kNN) using  $k=10$  (notice increased smoothness)

**k-nearest neighbors** (kNN) using  $k=1$  (notice decreased smoothness)

**Prediction error** from our fitted kNN (k=5) model

**Prediction error** from our fitted kNN (k=10) model



**Prediction error** from our fitted kNN ( $k=1$ ) model

Recall **Prediction error** from our fitted linear regression model

# Model accuracy

## Questions

1. Which of the methods was the most flexible? Inflexible?
2. Why do you think kNN with  $k=1$  had such low prediction error?
3. How could we (better) assess model/predictive performance?
4. Why would we ever want to choose a less flexible model?

# Model accuracy

## Measurement

You probably will not be surprised to know that there is no one-size-fits-all solution in statistical learning.

Q How do we choose between competing models?

# Model accuracy

## Measurement

You probably will not be surprised to know that there is no one-size-fits-all solution in statistical learning.

**Q** How do we choose between competing models?

**A** We're a few steps away, but before we do anything, we need a way to **define model performance**.

# Model accuracy

## Subtlety

Defining performance can actually be quite tricky...

*Regression setting, 1* Which do you prefer?

1. Lots of little errors and a few really large errors.
2. Medium-sized errors for everyone.

*Regression setting, 2* Is a 1-unit error (e.g., \$1,000) equally bad for everyone?

# Model accuracy

## Subtlety

Defining performance can actually be quite tricky...

*Classification setting, 1* Which is worse?

1. False positive (e.g., incorrectly diagnosing cancer)
2. False negative (e.g., missing cancer)

*Classification setting, 2* Which is more important?

1. True positive (e.g., correct diagnosis of cancer)
2. True negative (e.g., correct diagnosis of "no cancer")

# Model accuracy

## MSE

**Mean squared error (MSE)** is the most common<sup>†</sup> way to measure model performance in a regression setting.

$$\text{MSE} = \frac{1}{n} \sum_{i=1}^n \left[ y_i - \hat{f}(x_i) \right]^2$$

Recall:  $y_i - \hat{f}(x_i) = y_i - \hat{y}_i$  is our prediction error.

<sup>†</sup> *Most common* does not mean best—it just means lots of people use it.



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## MSE

**Mean squared error (MSE)** is the most common<sup>†</sup> way to measure model performance in a regression setting.

$$\text{MSE} = \frac{1}{n} \sum_{i=1}^n \left[ y_i - \hat{f}(x_i) \right]^2$$

Recall:  $y_i - \hat{f}(x_i) = y_i - \hat{y}_i$  is our prediction error.

Two notes about MSE

1. MSE will be (relatively) very small when **prediction error** is nearly zero.
2. MSE **penalizes** big errors more than little errors (the squared part).

<sup>†</sup> *Most common* does not mean best—it just means lots of people use it.

# Model accuracy

## Training or testing?

Low MSE (accurate performance) on the data that trained the model isn't actually impressive—maybe the model is just overfitting our data.<sup>†</sup>

*What we want:* How well does the model perform **on data it has never seen?**

<sup>†</sup> Recall the kNN performance for  $k=1$ .

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This introduces an important distinction:

1. **Training data:** The observations  $(y_i, x_i)$  used to **train** our model  $\hat{f}$ .
2. **Testing data:** The observations  $(y_0, x_0)$  that our model has yet to see—and which we can use to evaluate the performance of  $\hat{f}$ .

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**Real goal: Low test-sample MSE** (not the training MSE from before).

<sup>†</sup> Recall the kNN performance for  $k=1$ .

**Next time:** model performance, the variance-bias tradeoff, and kNN

# Sources

These notes draw upon

- [An Introduction to Statistical Learning \(ISL\)](#)  
James, Witten, Hastie, and Tibshirani
- [Python Data Science Handbook](#)  
Jake VanderPlas

I pulled the comic from [Twitter](#).

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