

Day 1: Overview of Core Concepts

ME314: Introduction to Data Science and Machine Learning

LSE Summer School

21 June 2021

Outline

Overview of Core Concepts

- Supervised Learning

- Unsupervised Learning

Machine Learning

Overview of Core Concepts

Philosophy

- ▶ It is important to understand the ideas behind the various techniques, in order to know how and when to use them.
- ▶ One has to understand the simpler methods first, in order to grasp the more sophisticated ones.
- ▶ It is important to accurately assess the performance of a method, to know how well or how badly it is working (simpler methods often perform as well as fancier ones!).
- ▶ This is an exciting research area, having important applications in science, industry and policy.
- ▶ Machine learning is a fundamental ingredient in the training of a modern **data scientist**.

Two main approaches to machine learning

- ▶ Supervised Learning
- ▶ Unsupervised Learning

The Supervised Learning Problem

Starting point:

- ▶ Outcome measurement Y (also called dependent variable, response, target).
- ▶ Vector of p predictor measurements X (also called inputs, regressors, covariates, features, independent variables).
- ▶ In the **regression problem**, Y is quantitative (e.g price, blood pressure).
- ▶ In the **classification problem**, Y takes values in a finite, unordered set (survived/died, digit 0-9, cancer class of tissue sample).
- ▶ We have training data $(x_1, y_1), \dots, (x_N, y_N)$. These are observations (examples, instances) of these measurements.

Objectives

On the basis of the training data we would like to:

- ▶ Accurately predict unseen test cases.
- ▶ Understand which inputs affect the outcome, and how.
- ▶ Assess the quality of our predictions and inferences.

Unsupervised Learning

- ▶ No outcome variable, just a set of predictors (features) measured on a set of samples.
- ▶ Objective is more fuzzy – find groups of samples that behave similarly, find features that behave similarly, find linear combinations of features with the most variation.
- ▶ Difficult to know how well you are doing.
- ▶ Different from supervised learning, but can be useful as a pre-processing step for supervised learning.

Machine learning and this course

Machine learning

- ▶ **Machine learning** refers to a vast set of tools for *understanding data*.
- ▶ For a quantitative response Y and a set of predictors X :

$$Y = f(X) + \epsilon$$

- ▶ Here, f represents the *systematic* information that X provides about Y .
- ▶ Statistical learning refers to a set of approaches for estimating f .
- ▶ Most of the course we'll spend talking about different ways to estimate f and how to evaluate whether we've done a good job with it.

Where does this course fit in?

- ▶ Supervised versus unsupervised learning.
- ▶ Regression versus classification.
- ▶ No single *best* method. We'll spend a lot of time choosing the most appropriate tool for a given dataset using different measures of the quality of fit. E.g. MSE

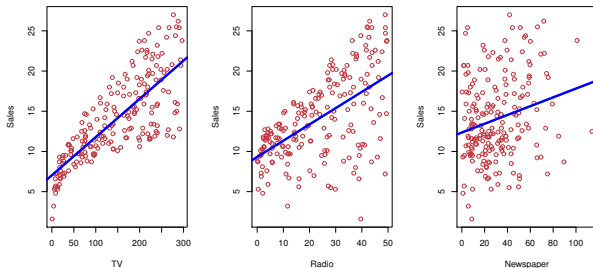
$$MSE_{training} = \frac{1}{n} \sum_{i=1}^n (y_i - \hat{f}(x_i))^2$$

$$MSE_{test} = \text{Ave}(\hat{f}(x_0) - y_0)^2$$

Why should we bother with f ?

1. **Prediction:** $\hat{Y} = \hat{f}(X)$, where \hat{f} is a *black box*.
2. **Inference:** How Y is changing as a function of X .
 - ▶ Depending on whether the ultimate goal is prediction, inference or a mix of both, we may deploy different methods for estimating f .
 - ▶ Also depending on the ultimate goal you may or may not care about evaluating the causal relationship between Y and X .

What is Machine Learning?



- ▶ Shown are Sales vs TV, Radio and Newspaper, with a blue linear-regression line fit separately to each.
- ▶ We can predict Sales using a model

$$Sales \approx f(TV, Radio, Newspaper)$$

Notation

- ▶ Here Sales is a *response* or *target* that we wish to predict. We generically refer to the response as Y .
- ▶ TV is a *feature*, or *input*, or *predictor*; we name it X_1 .
- ▶ Likewise name Radio as X_2 , and so on.
- ▶ We can refer to the input vector collectively as

$$X = (X_1, X_2, X_3)$$

- ▶ Now we write our model as

$$Y = f(X) + \epsilon$$

where ϵ captures measurement errors and other discrepancies.

What is $f(X)$ good for?

- ▶ With a good f we can make predictions of Y at new points $X = x$.
- ▶ We can understand which components of $X = (X_1, X_2, \dots, X_p)$ are important in explaining Y , and which are irrelevant. For example, Seniority and Years of Education have a big impact on Income, but Marital Status typically does not.
- ▶ Depending on the complexity of f , we may be able to understand how each component X_j of X affects Y .

- ▶ Is there an ideal $f(X)$? In particular, what is a good value for $f(X)$ at any selected value of X , say $X = 4$?
- ▶ There can be many Y values at $X = 4$. A good value is

$$f(4) = E(Y|X = 4)$$

- ▶ $E(Y|X = 4)$ means expected value (average) of Y given $X = 4$.
- ▶ This ideal $f(x) = E(Y|X = x)$ is called the regression function.

The regression function $f(x)$

- ▶ Is also defined for vector X ; e.g.
 $f(x) = f(x_1, x_2, x_3) = E(Y|X_1 = x_1, X_2 = x_2, X_3 = x_3)$
- ▶ Is the **ideal** or **optimal** predictor of Y with regard to mean-squared prediction error: $f(x) = E(Y|X = x)$ is the function that minimizes $E[(Y - g(X))^2|X = x]$ over all functions g at all points $X = x$.
- ▶ $\epsilon = Y - f(x)$ is the **irreducible** error – i.e. even if we knew $f(x)$, we would still make errors in prediction, since at each $X = x$ there is typically a distribution of possible Y values.
- ▶ For any estimate $\hat{f}(x)$ of $f(x)$, we have

$$E[(Y - \hat{f}(X))^2|X = x] = \underbrace{[f(x) - \hat{f}(x)]^2}_{\text{Reducible}} + \underbrace{\text{Var}(\epsilon)}_{\text{Irreducible}}$$

How to estimate f

- ▶ Typically we have few if any data points with $X = 4$ exactly.
- ▶ So we cannot compute $E(Y|X = x)$!
- ▶ Relax the definition and let

$$\hat{f}(x) = \text{Ave}(Y|X \in N(x))$$

where $N(x)$ is some **neighborhood** of x .

- ▶ Nearest neighbor averaging can be pretty good for small p – i.e. $p \leq 4$ and large-ish N .
- ▶ We will discuss smoother versions, such as kernel and spline smoothing later in the course.
- ▶ Nearest neighbor methods can be lousy when p is large. Reason: the **curse of dimensionality**. Nearest neighbors tend to be far away in high dimensions.
 - ▶ We need to get a reasonable fraction of the N values of y_i to average to bring the variance down – e.g. 10%.
 - ▶ A 10% neighborhood in high dimensions need no longer be local, so we lose the spirit of estimating $E(Y|X = x)$ by local averaging.

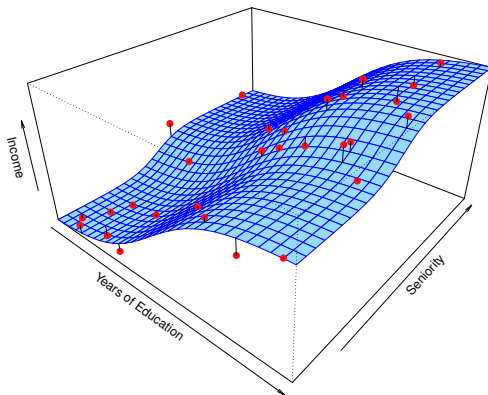
Parametric and structured models

The **linear** model is an important example of a parametric model:

$$f_L(X) = \beta_0 + \beta_1 X_1 + \beta_2 X_2 + \cdots + \beta_p X_p.$$

- ▶ A linear model is specified in terms of $p + 1$ parameters $\beta_0, \beta_1, \beta_2, \dots, \beta_p$.
- ▶ We estimate the parameters by fitting the model to training data.
- ▶ Although it is **almost never correct**, a linear model often serves as a good and interpretable approximation to the unknown true function $f(X)$.

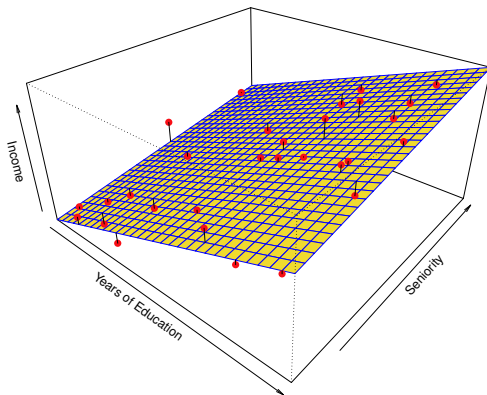
Simulated example



Red points are simulated values for income from the model

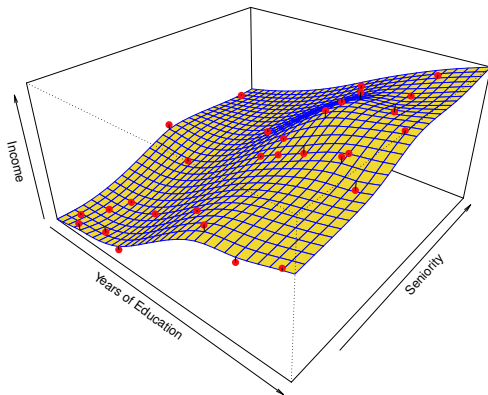
$$income = f(education, seniority) + \epsilon$$

f is the blue surface.

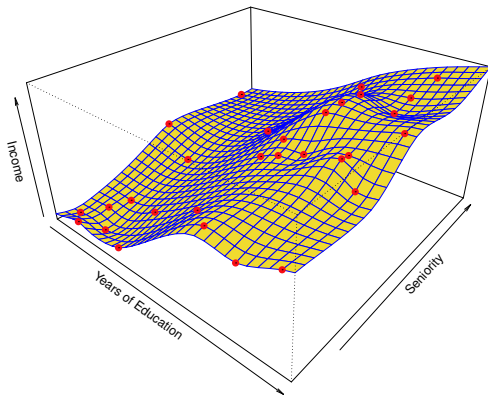


Linear regression model fit to the simulated data.

$$\hat{f}_L(\text{education}, \text{seniority}) = \hat{\beta}_0 + \hat{\beta}_1 \times \text{education} + \hat{\beta}_2 \times \text{seniority}$$



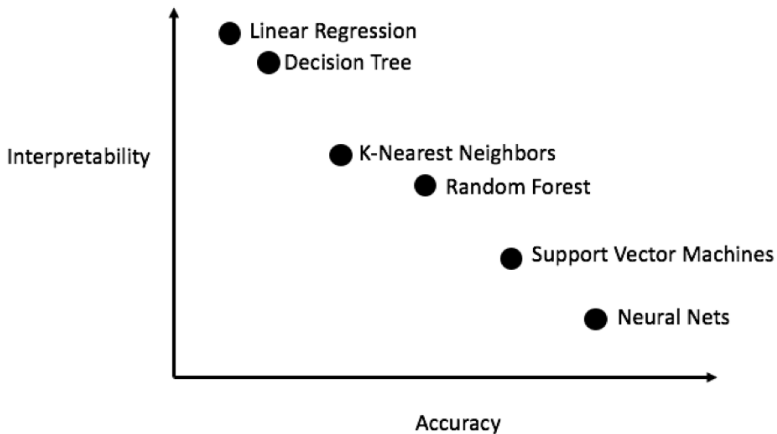
- ▶ More flexible regression model $\hat{f}_S(\text{education}, \text{seniority})$ fit to the simulated data.
- ▶ Here we use a technique called a **thin-plate spline** to fit a flexible surface.
- ▶ We control the roughness of the fit.



- ▶ Even more flexible spline regression model $\hat{f}_S(\text{education}, \text{seniority})$ fit to the simulated data.
- ▶ Here the fitted model makes no errors on the training data!
- ▶ Also known as **overfitting**.

Some trade-offs

- ▶ Prediction accuracy versus interpretability.
 - ▶ Linear models are easy to interpret; thin-plate splines are not.
- ▶ Good fit versus over-fit or under-fit.
 - ▶ How do we know when the fit is just right?
- ▶ Parsimony versus black-box.
 - ▶ We often prefer a simpler model involving fewer variables over a black-box predictor involving them all.



Source: (<https://medium.com/ansaro-blog/interpreting-machine-learning-models-1234d735d6c9>)

Assessing Model Accuracy

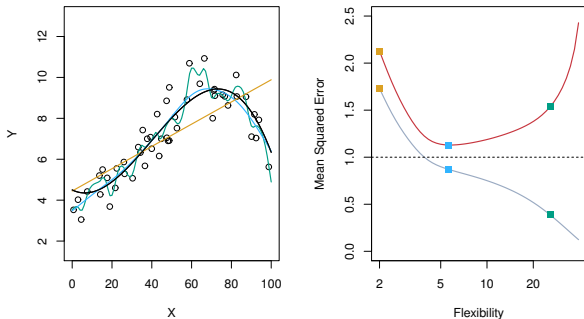
- ▶ Suppose we fit a model $\hat{f}(x)$ to some training data $Tr = \{x_i, y_i\}_1^N$, and we wish to see how well it performs.
- ▶ We could compute the average squared prediction error over Tr :

$$MSE_{Tr} = Ave_{i \in Tr} [y_i - \hat{f}(x_i)]^2$$

This may be biased toward more overfit models.

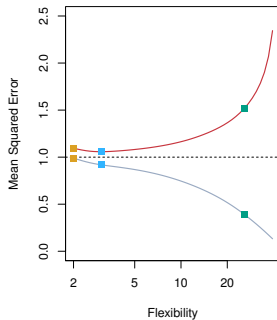
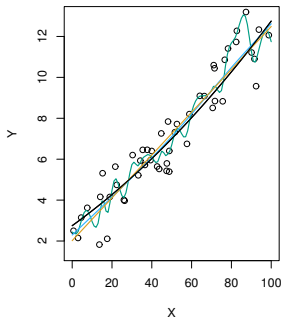
- ▶ Instead we should, if possible, compute it using fresh **test** data $Te = \{x_i, y_i\}_1^M$:

$$MSE_{Te} = Ave_{i \in Te} [y_i - \hat{f}(x_i)]^2$$

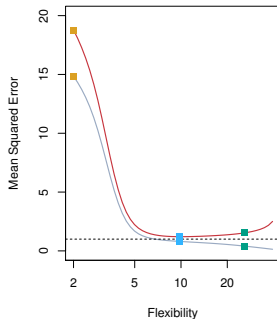
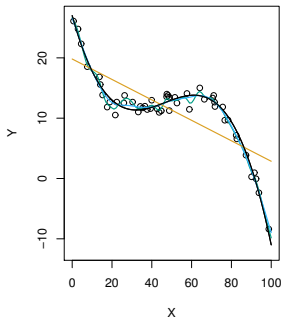


Data simulated from f , shown in black. Three estimates of f are shown: the linear regression line (orange curve), and two smoothing splines.

- ▶ Black curve is truth.
- ▶ Red curve on right is MSE_{Te} , grey curve is MSE_{Tr} .
- ▶ Orange, blue and green curves/squares correspond to fits of different flexibility.



- ▶ The setup as before, using a different true f that is much closer to linear. In this setting, linear regression provides a very good fit to the data.
- ▶ Here the truth is smoother, so the smoother fit and linear model do really well.



- ▶ Setup as above, using a different f that is far from linear.
- ▶ In this setting, linear regression provides a very poor fit to the data.
- ▶ Here the truth is wiggly and the noise is low, so the more flexible fits do the best.

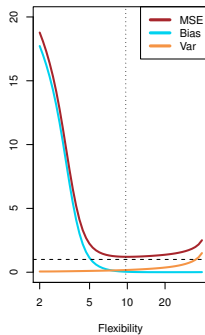
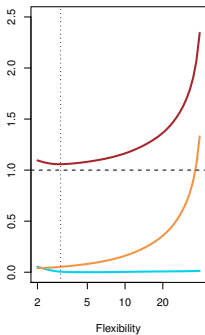
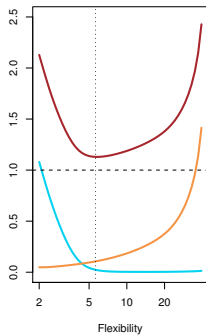
Bias-Variance Trade-off

- ▶ Suppose we have fit a model $f(x)$ to some training data Tr , and let (x_0, y_0) be a test observation drawn from the population.
- ▶ If the true model is $Y = f(X) + \epsilon$ (with $f(x) = E(Y|X = x)$), then

$$E(y_0 - \hat{f}(x_0)) = \text{Var}(\hat{f}(x_0)) + [\text{Bias}(\hat{f}(x_0))]^2 + \text{Var}(\epsilon).$$

- ▶ The expectation averages over the variability of y_0 as well as the variability in Tr . Note that $\text{Bias}(\hat{f}(x_0)) = E[\hat{f}(x_0)] - f(x_0)$.
- ▶ Typically as the **flexibility** of \hat{f} increases, its variance increases, and its bias decreases.
- ▶ So choosing the flexibility based on average test error amounts to a **bias-variance trade-off**.

Bias-variance trade-off for the three examples



Classification Problems

Here the response variable Y is **qualitative** – e.g. email is one of $\mathcal{C} = (\textit{spam}, \textit{ham})$ ($\textit{ham} = \textit{goodmail}$), digitclass is one of $\mathcal{C} = \{0, 1, \dots, 9\}$. Our goals are to:

- ▶ Build a classifier $\mathcal{C}(X)$ that assigns a class label from \mathcal{C} to a future unlabeled observation X .
- ▶ Assess the uncertainty in each classification.
- ▶ Understand the roles of the different predictors among $X = (X_1, X_2, \dots, X_p)$.

- ▶ Is there an ideal $\mathcal{C}(X)$? Suppose the K elements in \mathcal{C} are numbered $1, 2, \dots, K$. Let

$$p_k(x) = \Pr(Y = k | X = x), k = 1, 2, \dots, K.$$

- ▶ These are the **conditional class probabilities** at x . Then the **Bayes optimal** classifier at x is

$$\mathcal{C}(x) = j \text{ if } p_j(x) = \max\{p_1(x), p_2(x), \dots, p_K(x)\}$$

- ▶ Nearest-neighbor averaging can be used as before.
- ▶ Also breaks down as dimension grows.
- ▶ However, the impact on $\hat{\mathcal{C}}(x)$ is less than on $\hat{p}_k(x)$, $k = 1, \dots, K$.

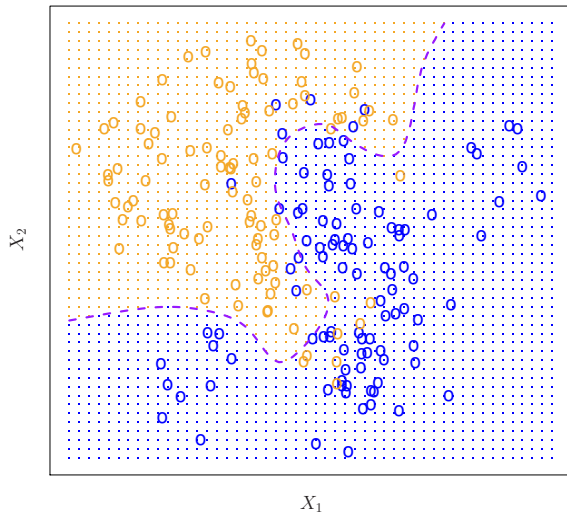
Classification: some details

- ▶ Typically we measure the performance of $\hat{\mathcal{C}}(x)$ using the misclassification error rate:

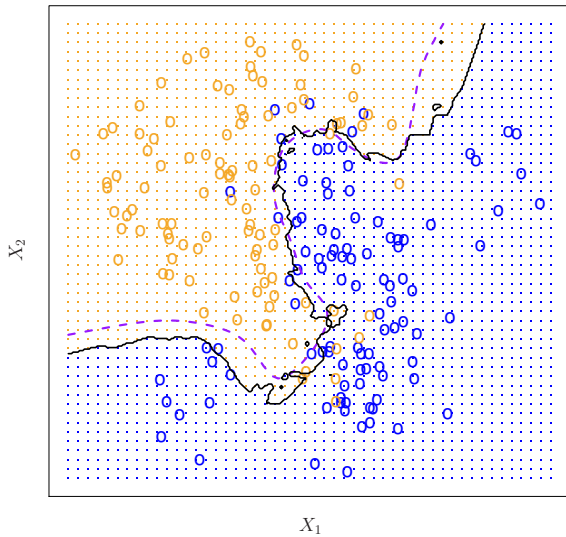
$$Err_{Te} = Ave_{i \in Te} \mathcal{I}[y_i \neq \hat{\mathcal{C}}(x_i)]$$

- ▶ The Bayes classifier (using the true $p_k(x)$) has smallest error (in the population).
- ▶ Support-vector machines build structured models for $\mathcal{C}(x)$.
- ▶ We will also build structured models for representing the $p_k(x)$. For example, logistic regression, generalized additive models.

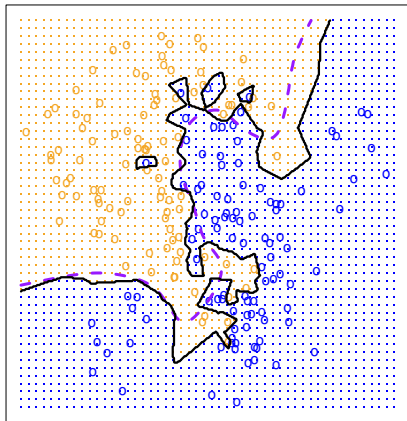
Example: K-nearest neighbors in two dimensions



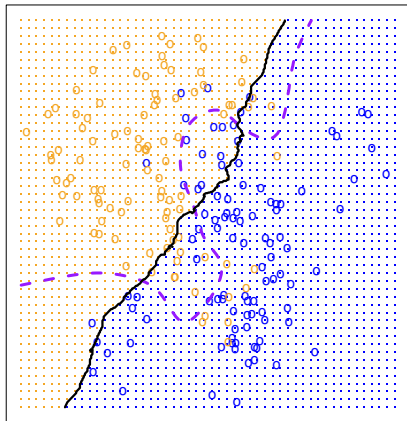
KNN: K=10

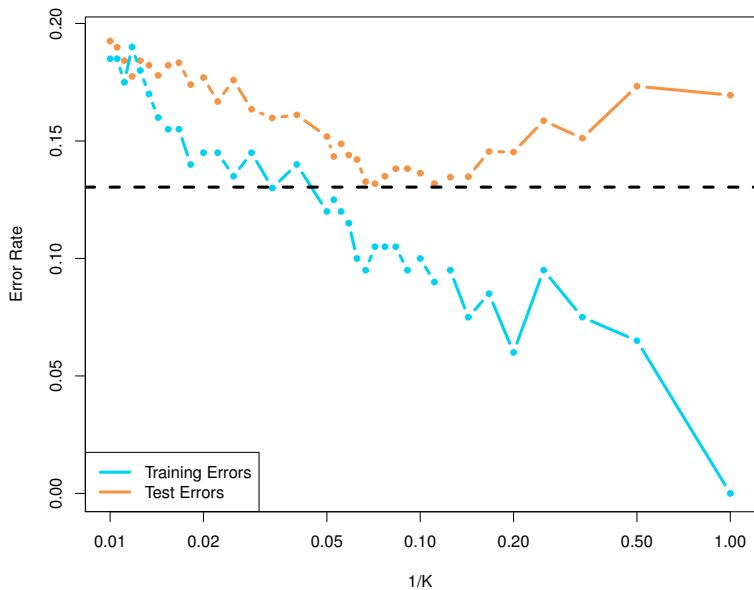


KNN: $K=1$



KNN: $K=100$





Prediction versus explanation

- ▶ Social science: The goal is typically *explanation*
- ▶ Data science: The goal is frequently *prediction*, or data exploration
- ▶ Many of the same methods are used for both objectives