

Regression

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Data are the set of inputs and outputs, $\mathcal{D} = \{(x_i, y_i)\}_{i=1}^n$



In *linear regression*, the goal is to predict y from x using a linear function



Examples of linear regression:

- given a child's age and gender, what is his/her height?
- given unemployment, inflation, number of wars, and economic growth, what will the president's approval rating be?
- given a browsing history, how long will a user stay on a page?



Multiple Covariates

Often, we have a vector of inputs where each represents a different *feature* of the data

$$\mathbf{x} = (x_1, \dots, x_p)$$

The function fitted to the response is a linear combination of the covariates

$$f(\mathbf{x}) = \beta_0 + \sum_{j=1}^{p} \beta_j x_j$$

Multiple Covariates

- Often, it is convenient to represent **x** as $(1, x_1, \dots, x_p)$
- In this case **x** is a vector, and so is β (we'll represent them in bold face)
- This is the dot product between these two vectors
- This then becomes a sum (this should be familiar!)

$$\beta \mathbf{x} = \beta_0 + \sum_{j=1}^{p} \beta_j x_j$$

Hyperplanes: Linear Functions in Multiple Dimensions

Hyperplane



Covariates

- Do not need to be raw value of x₁, x₂,...
- Can be any feature or function of the data:
 - Transformations like $x_2 = \log(x_1)$ or $x_2 = \cos(x_1)$
 - Basis expansions like $x_2 = x_1^2$, $x_3 = x_1^3$, $x_4 = x_1^4$, etc
 - Indicators of events like $x_2 = 1_{\{-1 \le x_1 \le 1\}}$
 - Interactions between variables like $x_3 = x_1 x_2$
- Because of its simplicity and flexibility, it is one of the most widely implemented regression techniques

Fitting a Linear Regression



Idea: minimize the Euclidean distance between data and fitted line

$$RSS(\beta) = \frac{1}{2} \sum_{i=1}^{n} (y_i - \beta \mathbf{x}_i)^2$$

How to Find β

- Use calculus to find the value of β that minimizes the RSS
- The optimal value is

$$\hat{\beta} = \frac{\sum_{i=1}^{n} y_i x_i}{\sum_{i=1}^{n} x_i^2}$$



- After finding $\hat{\beta}$, we would like to predict an output value for a new set of covariates
- We just find the point on the line that corresponds to the new input:

$$\hat{y} = \beta_0 + \beta_1 x \tag{1}$$



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$$\hat{y} = 1.0 + 0.5x$$
 (1)



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$$\hat{y} = 1.0 + 0.5 * 5$$
 (1)

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$$\hat{y} = 3.5 \tag{1}$$

Probabilistic Interpretation

- Our analysis so far has not included any probabilities
- Linear regression does have a probabilisitc (probability model-based) interpretation



Probabilistic Interpretation

 Linear regression assumes that response values have a Gaussian distribution around the linear mean function,

$$Y_i | \mathbf{x}_i, \beta \sim N(\mathbf{x}_i \beta, \sigma^2)$$

This is a discriminative model, where inputs x are not modeled



Minimizing RSS is equivalent to maximizing conditional likelihood



We will predict the time that we will have to wait to see the next eruption given the duration of the current eruption



We can plot our data and make a function for new predictions

```
> # Plot a line on the data
>
   abline (fit.lm, col="red", lwd=3)
>
>
 # Make a function for prediction
>
   fit.lm$coefficients[1]
(Intercept)
     33.4744
> fit.lm$coefficients[2]
eruptions
  10.72964
> faithful.fit <- function(x) fit.lm$coefficients[1] +</pre>
fit.lm$coefficients[2] *x
> x.pred <- c(2.0, 2.7, 3.8, 4.9)</pre>
> faithful.fit(x.pred)
[1] 54.93368 62.44443 74.24703 86.04964
```



Multivariate Linear Regression

Example: p = 1, have 2 points



- Have p+1 or fewer points, line hits all (or p with mean 0 data)
- $\geq p + 1$ (but still close to that number), line goes *close* to all points

Noise, Bias, Variance Tradeoff



- Noise: Lower bound on performance
- Bias: Error as a result as choosing the wrong model
- Variance: Variation due to training sample and randomization

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- No model is perfect
- More complex models are more susceptible to errors due to variance

Multivariate Linear Regression

Why linear regression:

- has few parameters to estimate (p)
- really restrictive model-low variance, higher bias



- should be good for data with few observations, large number of covariates...
- ... but we can't use it in this situation

Multivariate Linear Regression

Idea: if we have a large number of covariates compared to observations, say n < 2p, best to estimate most coefficients as 0!

- not enough info to determine all coefficients
- try to estimate ones with strong signal
- set everything else to 0 (or close)

Coefficients of 0 may not be a bad assumption...

If we have 1,000s of coefficients, are they all equally important?

Gene Expression

Example: microarray gene expression data

- gene expression: want to measure the level at which information in a gene is used in the synthesis of a functional gene product (usually protein)
- can use gene expression data to determine subtype of cancer (e.g. which type of Lymphoma B?) or predict recurrence, survival time, etc
- problem: thousands of genes, hundreds of patients, p > n!

Intuition: only a handful of genes should affect outcomes

Gene Expression



- gene expression levels are continuous values
- data: observation *i* is gene expression levels from patient *i*, attached to outcome for patient (survival time)
- covariates: expression levels for p genes

Gene Expression



- collinearity: does it matter which gene is selected for prediction? No!
- overfitting: now fitting p' non-0 coefficients to n observations with p' << n means less fitting of noise

Regularization:

- still minimize the RSS
- place a *penalty* on large values for β₁, ..., β_p (why not β₀? can always easily estimate mean)
- add this penalty to the objective function
- solve for $\hat{\beta}$!

New objective function:

$$\hat{\beta} = \arg\min_{\beta} \frac{1}{2} \sum_{i=1}^{n} (y_i - \mathbf{x}_i \beta)^2 + \lambda \sum_{j=1}^{p} \operatorname{penalty}(\beta_j)$$

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Regularization: what is a good penalty function?

Same as penalties used to fit errors:

Ridge regression (squared penalty):

$$\hat{\beta}^{\text{Ridge}} = \arg\min_{\beta} \frac{1}{2} \sum_{i=1}^{n} (y_i - \mathbf{x}_i \beta)^2 + \lambda \sum_{j=1}^{p} \beta_j^2$$

Lasso regression (absolute value penalty):

$$\hat{\beta}^{Lasso} = \arg\min_{\beta} \frac{1}{2} \sum_{i=1}^{n} (y_i - \mathbf{x}_i \beta)^2 + \lambda \sum_{j=1}^{p} |\beta_j|$$

Comparing Ridge and Lasso

	Ridge	Lasso
Objective	$\frac{1}{2}\sum_{i=1}^{n}(y_{i}-\mathbf{x}_{i}\beta)^{2}+\lambda\sum_{i=0}^{p}\beta_{i}^{2}$	$\frac{1}{2}\sum_{i=1}^{n}(\mathbf{y}_{i}-\mathbf{x}_{i}\boldsymbol{\beta})^{2}+\lambda\sum_{j=0}^{p} \boldsymbol{\beta}_{j} $
Estimator	$(\mathbf{X}^T\mathbf{X} + \lambda I)^{-1}\mathbf{X}^T\mathbf{y}^T$	not closed form
Coefficients	most close to 0	most exactly 0
Stability	robust to changes in X , y	not robust to changes in X , y

Regularized linear regression is fantastic for low signal datasets or those with p >> n

- Ridge: good when many coefficients affect value but not large (gene expression)
- Lasso: good when you want an *interpretable* estimator

Choosing λ

Both Ridge and Lasso have a tunable parameter, λ

use cross validation to find best λ

$$\hat{\lambda} = \arg\min_{\lambda} \sum_{i=1}^{n} (y_i - \mathbf{x}_i \hat{\beta}_{-i,\lambda})^2$$

- try out many values
- see how well it works on "development" data

Regression

- Workhorse technique of data analysis
- Fundamental tool that we saw before ("Logistic Regression")
- Important to understand interpretation of regression parameters