



# Linear Regression

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Data Science Decal

Hosted by Machine Learning at Berkeley

## Agenda

Background

Model Estimation

Assumptions

Model Testing

Next Steps

# Background

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	Continuous	Discrete
Supervised	<b>Regression</b>	Classification
Unsupervised	Dimensionality Reduction	Clustering

- Suppose we have data
  - Want to model relationships and make **predictions**
- The data has **continuous** labels ( $y$ )
  - i.e. prices, heights, miles per gallon, etc.
- The data has a set of **explanatory** variables ( $x_i$ )
  - i.e. sales, weights, engine power, etc.
- How does a computer make predictions?

- Regression is one of the most commonly used methods by data scientists
- It is simple, fast, interpretable, and **powerful**
- The techniques we use here are widely applicable
- It is practical!
  - (Physics) Ohm's law, Hooke's law, Charles's law
  - (Economics) Okun's law

- Suppose we have  $p$  predictor variables  $x_1, x_2, \dots, x_p$
- We can approximate  $y$  as a linear function of the  $x_i$ 's:

$$h_{\theta}(x) = \theta_0 + \theta_1 x_1 + \dots + \theta_p x_p$$

- $\theta_i$ 's are the **parameters** (also called **weights**) which we need to estimate
- We introduce  $x_0 = 1$  for simplicity so that:

$$y = \sum_{i=1}^p \theta^T x$$

- Suppose we have the following data about houses:

Price	# of Square Feet	# of Bedrooms
221,900	1180	3
538,000	2570	3
⋮	⋮	⋮
1,225,000	5420	4

- Let's predict the price of a house from the number of square feet it has
- Our linear model has the form:

$$h_{\theta}(sqft) = \theta_0 + \theta_1 sqft$$

# Model Estimation

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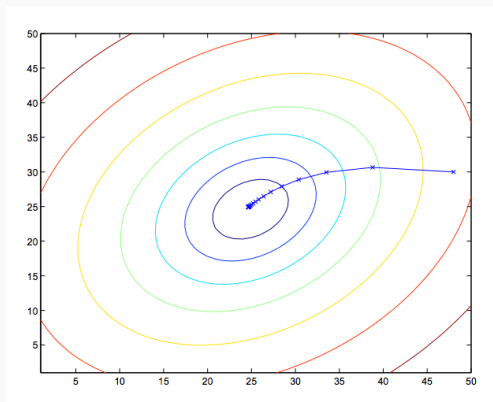


- **Goal:** have  $h_{\theta}(x)$  be as close to  $y$  as possible
- We can translate this goal into mathematics by defining the **cost function**

$$J(\theta) = \frac{1}{2} \sum_{i=1}^n (h_{\theta}(x^{(i)}) - y^{(i)})^2$$

- $J(\theta)$  sums the squared **residuals**
- To have an accurate model, we want to **minimize**  $J(\theta)$

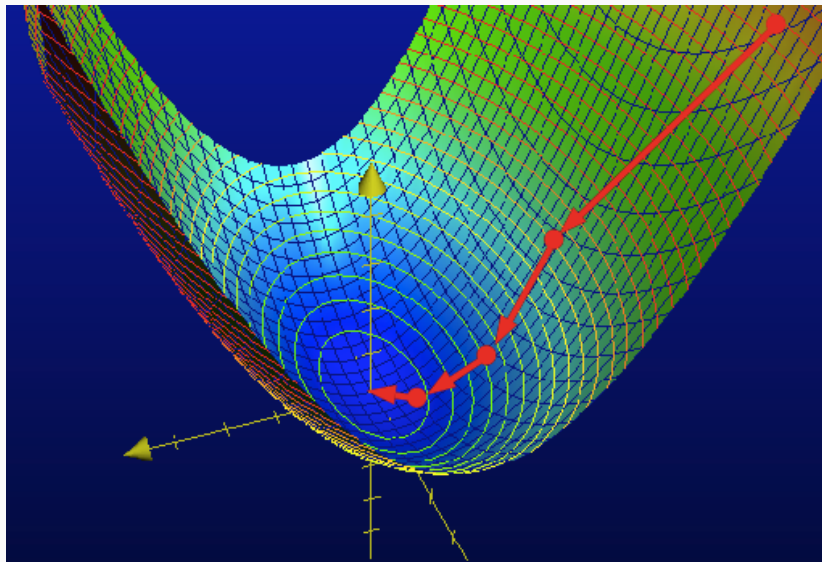
- **Idea:** choose  $\theta$  to minimize  $J(\theta)$
- We can use a search algorithm that follows the scheme:
  - Choose an initial guess for  $\theta$
  - Repeatedly update  $\theta$  to make  $J(\theta)$
  - Keep doing this until  $J(\theta)$  reaches its minimum



- **Note:**  $J(\theta)$  is a convex quadratic function (has nice properties)
- From Math 53: the direction of greatest increase is the same direction of the gradient vector
- **Idea:** let's update  $\theta$  by traversing the opposite direction instead
- This scheme is known as **gradient descent**

$$\theta \leftarrow \theta - \epsilon \nabla_{\theta} J(\theta)$$

- $\epsilon$  is called the **learning rate**



- Let's start with the case where we only have one training example  $(x^{(1)}, y^{(1)})$

$$\begin{aligned}\nabla_{\theta} J(\theta) &= \nabla_{\theta} \frac{1}{2} (h_{\theta}(x^{(1)}) - y^{(1)})^2 \\ &= 2\left(\frac{1}{2}\right) (h_{\theta}(x^{(1)}) - y^{(1)}) \nabla_{\theta} (h_{\theta}(x^{(1)}) - y^{(1)}) \\ &= (h_{\theta}(x^{(1)}) - y^{(1)}) \nabla_{\theta} (\theta^T x^{(1)} - y^{(1)}) \\ &= (h_{\theta}(x^{(1)}) - y^{(1)}) x^{(1)}\end{aligned}$$

- For a single training example, the update rule is:

$$\theta \leftarrow \theta - \epsilon (y^{(1)} - h_{\theta}(x^{(1)})) x^{(1)}$$

- For  $n$  training examples:

$$\theta_j \leftarrow \theta - \epsilon \sum_i^n (h_\theta(x^{(i)}) - y^{(i)}) x_j^{(i)}$$

for  $j = 1, \dots, p$       and  $i = 1, \dots, n$

- This rule is also called the **LMS** update rule ("least mean squares")
- Size of update is proportional to the residual term  $(y^{(i)} - h_\theta(x^{(i)}))$
- If the prediction  $h_\theta(x^{(i)})$  is close the actual  $y^{(i)}$  then the parameters  $\theta$  shouldn't need much changing

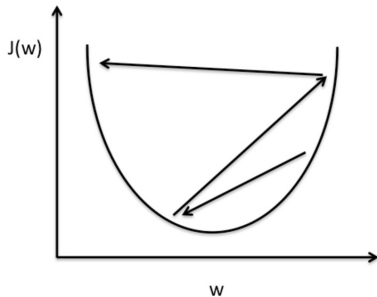
## Stochastic gradient descent for linear regression

While  $J(\theta)$  is not minimized:

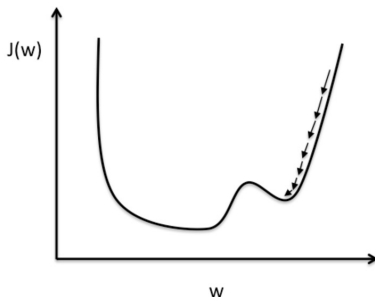
For  $i = 1, \dots, n$ :

$$\theta_j \leftarrow \theta_j - \epsilon(h_{\theta}(x^{(i)}) - y^{(i)})x_j^{(i)} \quad (\text{for each } j)$$

# Choosing the learning rate



**Large learning rate: Overshooting.**



**Small learning rate: Many iterations until convergence and trapping in local minima.**



- SGD is the basis for many optimization algorithms
- Not necessary for linear regression, as there exists a closed form solution
- We can find the optimal  $\theta$  by solving the **normal equations**

- Recall the equation for a linear model:

$$h_{\theta}(x_i) = \theta_0 + \theta_1 x_{i1} + \cdots + \theta_p x_{ip}$$

- The outcome,  $y$ , which we observe can be thought of as:

$$y_i = h_{\theta}(x_i) + \epsilon_i$$

where  $\epsilon$  is some unobserved error

- We don't know the true  $\theta$  is, so we estimate it with  $\hat{\theta}$
- Our predictions for test points are then

$$\hat{y} = h_{\hat{\theta}}(x)$$

- We can rewrite linear regression as

$$\begin{pmatrix} y_1 \\ y_2 \\ \vdots \\ y_n \end{pmatrix} = \begin{pmatrix} 1 & x_{11} & x_{12} & \dots & x_{1p} \\ 1 & x_{21} & x_{22} & \dots & x_{2p} \\ \vdots & \vdots & \vdots & \ddots & \vdots \\ 1 & x_{n1} & x_{n2} & \dots & x_{np} \end{pmatrix} \begin{pmatrix} \theta_0 \\ \theta_1 \\ \vdots \\ \theta_p \end{pmatrix} + \begin{pmatrix} \epsilon_1 \\ \epsilon_2 \\ \vdots \\ \epsilon_n \end{pmatrix}$$

- In compressed notation:

$$\vec{y} = X\vec{\theta} + \vec{\epsilon}$$

- Here, we're using capital letters to represent matrices, and arrows to represent vectors

- We want our estimate,  $\hat{\theta}$  to be accurate
- We can be accurate by trying to minimize error
- We can be accurate by minimizing our residuals

$$e_i = y_i - \vec{\theta}^T x_i$$

- More mathematically convenient to minimize squared residuals
- That is,

$$\hat{\vec{\theta}} = \operatorname{argmin}_{\vec{\theta}} \sum_{i=1}^n (y_i - (\theta_0 + \theta_1 x_{i1} + \cdots + \theta_p x_{ip}))^2$$

$$\begin{aligned}\hat{\theta} &= \operatorname{argmin}_{\vec{\theta}} \|\vec{y} - X\vec{\theta}\|_2^2 \\ &= \operatorname{argmin}_{\vec{\theta}} (\vec{y} - X\vec{\theta})^T (\vec{y} - X\vec{\theta}) \\ &= \operatorname{argmin}_{\vec{\theta}} \vec{y}^T \vec{y} - 2\vec{y}^T X\vec{\theta} + \vec{\theta}^T X^T X \vec{\theta}\end{aligned}$$

Let  $Q = \|\vec{y} - X\vec{\theta}\|_2^2$

Taking the derivative with respect to the vector  $\vec{\theta}$ :

$$\frac{\partial Q}{\partial \vec{\theta}} = 2X^T X \vec{\theta} - 2X^T \vec{y} = 0$$

$$\boxed{\hat{\theta} = (X^T X)^{-1} X^T \vec{y}}$$

- $\hat{\theta}$  is indeed a minimizer (the second derivative is negative)
- Gauss Markov Theorem:  $\hat{\theta}$  is BLUE (best linear unbiased estimator)
- The residuals are:

$$\vec{e} = (\vec{y} - \hat{\vec{y}}) = (I_{n \times n} - X(X^T X)^{-1} X^T) \vec{y}$$

$$[\hat{\vec{y}} = X\hat{\theta} = X(X^T X)^{-1} X^T Y]$$

- $\hat{\theta}$  is a random variable and thus has variance:

$$\begin{aligned} \text{Var}(\hat{\theta}) &= \text{Var}((X^T X)^{-1} X^T Y) = (X^T X)^{-1} X^T \text{Var}(\vec{y}) X (X^T X)^{-1} \\ &= \sigma^2 (X^T X)^{-1} \end{aligned}$$

$$[\text{Var}(\vec{y}) = \sigma^2 I_{n \times n}]$$

- Recall, once we have our estimate  $\hat{\theta}$ , we can predict new  $x$ 's using:

$$\hat{y}_i = \hat{\theta}_0 + \hat{\theta}_1 x_{i1} + \cdots + \hat{\theta}_p x_{ip}$$

- In matrix notation:

$$\hat{\vec{y}} = X\hat{\vec{\theta}}$$

- For a one unit increase in  $x_{ik}$ , we expect  $y_i$  to, **on average** increase by  $\hat{\theta}_k$
- If we take the log of the independent variables, the dependent variable, or both, then the above interpretation changes to involve **percent changes**

# Assumptions

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- Regression is a good summary of data, assuming the data has some key properties
- We need to know what those assumptions are, how to test for them, and what to do when they fall apart

- Linearity
- Normality of errors

$$\epsilon_i \sim N(0, \sigma^2)$$

- Homoscedasticity (constant variance)

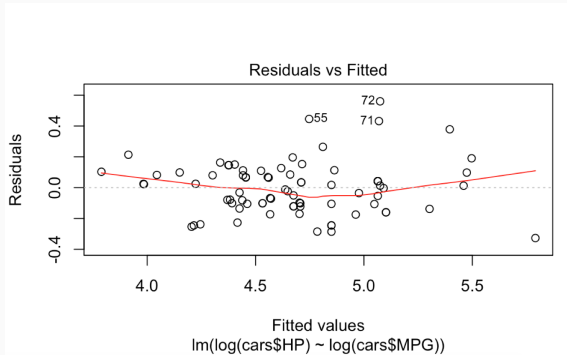
$$\text{Var}(\epsilon_i) = \sigma^2 \neq \sigma^2(x)$$

- Independence of errors

$$\epsilon_i \perp\!\!\!\perp \epsilon_j \quad \forall i \neq j$$

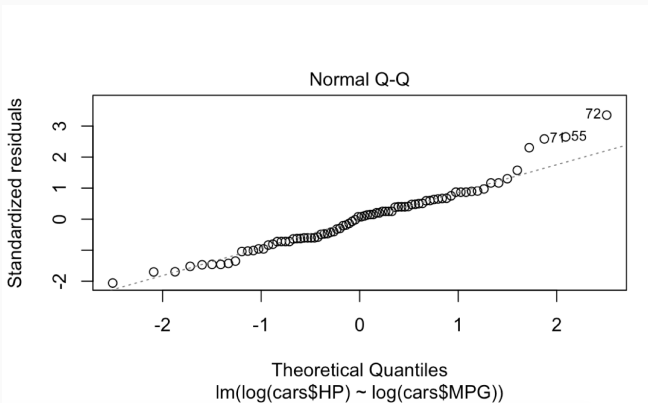
## Linearity

- Scatter plot of Y vs. standardized residuals should have no pattern



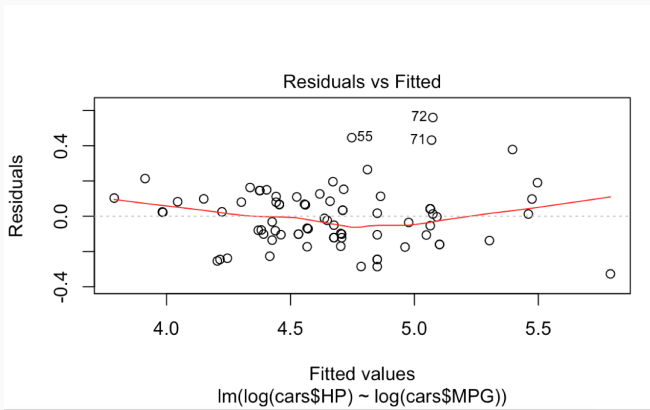
## Normality of errors

- Plot a histogram of the estimated errors (called **residuals**)
- QQplot
- Many tests exist: Kolomogorov-Smirnov, Shapiro-Wilk, ...



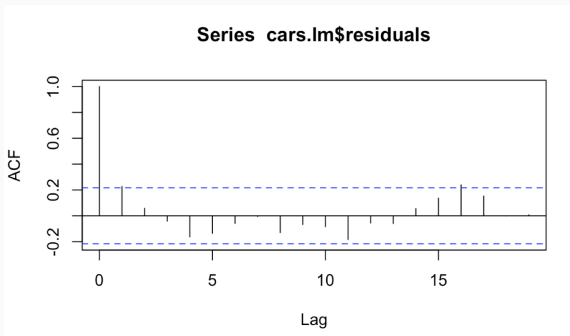
## Homoscedasticity

- Plot of Y vs. residuals should have equal variation across vertical slices
- Tests: Brusch-Pagan, White test, ...



## Independence of errors

- Autocorrelation plots
  - Most of the residuals should fall within the 95% confidence band around 0
- Durban-Watson test



- If the data is nonlinear...
  - Try performing a transformation on the independent or dependent variables such as squaring it, taking the log or square root, or ...
- If the errors are not normal...
  - Often, this isn't a big problem
  - Transformations help here too
  - Maybe subsets of the data are more normal than the overall set
  - Outliers and/or high leverage points may contribute to this issue

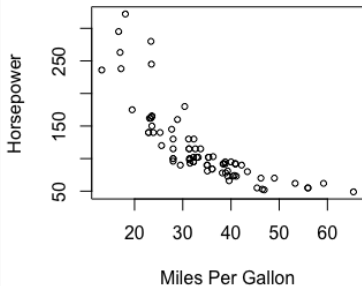
- If the data exhibits heteroscedasticity...
  - Log transformations are helpful
  - Search for and remove outliers/high-leverage points
  - Use a more advanced model (ARCH: auto-regressive conditional heteroscedasticity)
  - Heteroscedasticity may arise from violation of one of the other assumptions
- If the errors are not independent...
  - You have a structural problem in your model
  - Very hard to fix...
  - One way that I am aware of: identify an appropriate ARMA process and fit a generalized least squares model



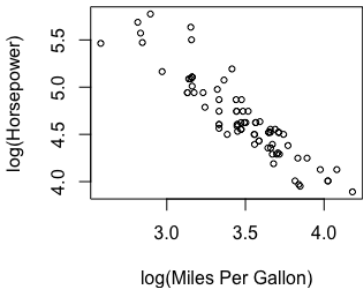
# Example of the beauty of a log transform



**MPG vs HP**



**Log(MPG) vs Log(HP)**



# Model Testing

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Once we have estimated  $\hat{\theta}$ , we have some questions:

- Is  $\theta_i$  significantly different from 0? (Is the variable  $X_i$  relevant?)
- How confident are we about what the true  $\theta$  is?
- How do we know what independent variables to use?

Once we have estimated  $\hat{\theta}$ , we have some questions:

- Is  $\theta_i$  significantly different from 0? (Is the variable  $X_i$  relevant?)
  - Perform some hypothesis tests
  - t-tests, F-tests, etc...  
[https://en.wikipedia.org/wiki/Statistical\\_hypothesis\\_testing](https://en.wikipedia.org/wiki/Statistical_hypothesis_testing)
- How confident are we about what the true  $\beta$  is?
  - Construct a confidence interval (many different kinds)  
[https://en.wikipedia.org/wiki/Confidence\\_interval](https://en.wikipedia.org/wiki/Confidence_interval)
- How do we know what independent variables to use?
  - Let's talk about this one some more

- Before we do any feature selection, we need to make sure to split our dataset into a **training set** and a **validation set**
- Greedy forwards selection
- Greedy backwards selection
- Other search algorithms...
- Many different "goodness" metrics exist to compare models:
  - $R^2$  (want more), MSE (want less), AIC and BIC (want less), ...
  - MSE (mean squared error):

$$MSE = \frac{1}{n} \sum_{i=1}^n (Y_i - \hat{Y}_i)^2$$

## Next Steps

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- **Classification problems:** logistic regression, support vector machines
- **Non-linearity:** kernel smoothing, splines and generalized additive models; nearest neighbor methods
- **Interactions:** tree-based methods, bagging, random forests and boosting (also capture non-linearities)
- **Regularized fitting:** ridge regression and lasso

- Polynomial transformations
- Basis expansions
- Dummy coding of categorical inputs
- Time series models
- Hierarchical modeling
- Causal inference
- Spatial models



Questions?