

SOSC 5340: Generalized Linear Models

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Feb 23, 2021

Outline

Logistics

Generalized Linear Models

Multinomial and Ordered Logit

Poisson, Negative Binomial, and Zero-inflated Poisson

Model Selection

Bias-Variance Trade-Off

Today's Review

Next week

- Presentation
- Exercise 1 Due (before the class, submit your code and results on Canvas)

Limited Dependent Variable

- Beyond binary outcomes, $Y \in \{0, 1\}$
- Categorical:
 - e.g., major choices;
- Integer (count): $Y \in \{0, 1, 2, \dots\}$
 - e.g., event counts
- Censored: observed Y is in a certain range, but we know in reality they should not be
 - e.g., US census write anyone who report their age > 90 as 90; so in census, age is between $[0, 90]$
- The common problem is that the outcome Y is limited to some regions, not in $(-\infty, \infty)$
 - so economists sometimes call them as **limited dependent variable**

Generalized Linear Model

- To model limited dependent variables, we use **generalized linear model** (GLM)
- GLM looks like:
 - $h(E(Y|X)) = X\beta$
 - or, $E(Y|X) = h^{-1}(X\beta)$
- $h()$ is called **link** function
- Linear regression is a kind of GLM, where $h(X) = X$
- Logistic regression is a kind of GLM, where $h(X) = \text{logit}(X)$
- Other GLM we will learn today choose different $h()$ to model different types of Y

GLM

- In practice, scholars use MLE to make statistical estimation and inference for GLM
- Recall that to use MLE, we need to make assumptions about what $p(Y|X)$ looks like

Estimation and Inference of MLE

- Steps are standard
 1. write down $P(Y|X)$
 2. write down $\log L$: the log-likelihood function
 3. obtain coefficient estimates that maximize log-likelihood
 - and use Hessian matrix to calculate confidence interval

Extending Logistic Regression

- Suppose we have categorical outcome with more than two values
- Sometimes, these categories have no intrinsic orders
 - E.g., majors choices between (Economics = 1, Political Science = 2, Sociology = 3, Public Policy = 4)
- Other times, these categories are **ordinal**
 - E.g., a survey ask whether you think religion deters economic growth, on a 1-7 scale.
 - 1 means strongly disagree, and 7 means strongly agree
 - Order gives more information than pure categories
 - Why not use continuous outcome models?
 - Dont want to assume equal distances between levels
 - Say, moving from 1-4 is different from 4-7
 - Assuming continuous Y does not distinguish these two

Ordered Logit: ordered outcome

- Peter McCullough, *Regression Models for Ordinal Data*, 1980
- Recall that logistic regression assumes a generating process based on latent variables

$$Y^* = X\beta + \epsilon$$

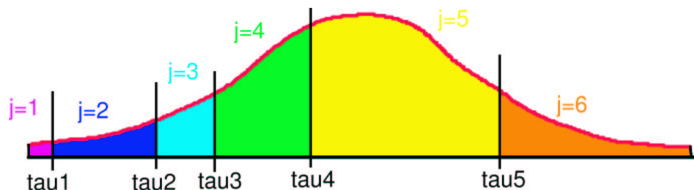
$$Y = \begin{cases} 1 & \text{if } y^* > 0 \\ 0 & \text{otherwise} \end{cases} \quad (1)$$

- Y^* is an unobserved latent variable
- if the latent variable is bigger than a pre-determined **cutoff** (here 0), $Y = 1$
- Otherwise, $Y = 0$

Ordered Logit

- We can borrow the same intuition to derive ordered logit regression, with $J > 2$ ordinal categories
- We create $J - 1$ latent cutoffs

$$Y = \begin{cases} 1 & \text{if } Y^* \leq \tau_1 \\ 2 & \text{if } \tau_1 < Y^* \leq \tau_2 \\ 3 & \text{if } \tau_2 < Y^* \leq \tau_3 \\ \vdots & \\ J & \text{if } \tau_{J-1} \leq Y^* \end{cases} \quad (2)$$



Ordered Logit

- So now the Assumption 3 for ordered logit becomes:

$$Y^* = X\beta + \epsilon$$

$$Y = \begin{cases} 1 & \text{if } Y^* \leq \tau_1 \\ 2 & \text{if } \tau_1 < Y^* \leq \tau_2 \\ 3 & \text{if } \tau_2 < Y^* \leq \tau_3 \\ \vdots & \\ J & \text{if } \tau_{J-1} \leq Y^* \end{cases} \quad (3)$$

- And the error ϵ follows a standard logistic distribution (the same as logistic regression)

Ordered Logit vs Linear Regression

- It may be easier to change from “very unlikely” (1) to “unlikely” (2), but it is more difficult to change from “unlikely” to “neutral” (3)
- For linear regression
 - It takes the same amount of changes in X to turn Y from 1 to 2 versus Y from 2 to 3
 - Linear regression does not capture this difference
- For ordered logit
 - Y changing from 1 to 2 means latent Y^* changes from below τ_1 to (τ_1, τ_2)
 - Y changing from 2 to 3 means latent Y^* changes from (τ_1, τ_2) to (τ_2, τ_3)
 - It often requires a different amount a change in X to move Y from 1 to 2 versus from 2 to 3. That's what we want to capture

Ordered Logit

- For MLE, we have to explicitly write down $P(Y|X)$

$$\begin{aligned}
 P(Y = 1|X) &= \Pr(\beta X + \epsilon \leq \tau_1|X) \\
 &= P(\epsilon \leq \tau_1 - \beta X|X) \\
 &= F(\tau_1 - \beta X), \text{ (definition of cumulative probability } F) \\
 &= \text{logit}^{-1}(\tau_1 - \beta X)
 \end{aligned} \tag{4}$$

$$\begin{aligned}
 P(Y = 2|X) &= \Pr(\tau_1 < \beta X + \epsilon \leq \tau_2|X) \\
 &= \Pr(\tau_1 - \beta X < \epsilon \leq \tau_2 - \beta X|X) \\
 &= F(\tau_2 - \beta X) - F(\tau_1 - \beta X) \\
 &= \text{logit}^{-1}(\tau_2 - \beta X) - \text{logit}^{-1}(\tau_1 - \beta X)
 \end{aligned} \tag{5}$$

And so on and so forth, for j up to $J - 1$

Ordered Logit

The last category J

$$\begin{aligned}
 P(Y = J|X) &= P(\tau_{J-1} \leq \beta X + \epsilon|X) \\
 &= P(\epsilon \geq \tau_{J-1} - \beta X|X) \\
 &= 1 - P(\epsilon < \tau_{J-1} - \beta X) & (6) \\
 &= 1 - F(\tau_{J-1} - \beta X) \\
 &= 1 - \text{logit}^{-1}(\tau_{J-1} - \beta X)
 \end{aligned}$$

- We have written down $P(Y|X)$ for every possible value of Y .
- Now we can use MLE to estimate parameters
- Now, there are regression coefficients β , as well as cutoffs τ
- Statistical software will return estimates for both

Ordered Logit

- What do the cutoffs τ mean?
- Recall that $P(Y = 1|X) = \text{logit}^{-1}(\tau_1 - \beta X)$
- And $P(Y = 2|X) = \text{logit}^{-1}(\tau_2 - \beta X) - \text{logit}^{-1}(\tau_1 - \beta X)$
- We add then together:

$$P(Y = 1|X) + P(Y = 2|X) = P(Y \leq 2|X) = \text{logit}^{-1}(\tau_2 - \beta X) \quad (7)$$

- And take the logit:

$$\text{logit}(P(Y \leq 2)) = \tau_2 - \beta X \quad (8)$$

- The rest is similar

$$\text{logit}(P(Y \leq j)) = \tau_j - \beta X$$

- In this way, τ looks like intercepts in normal regressions; so some other software (R) call them intercepts

Multinomial Logit: categorical outcome

- Multinomial logit: for categorical outcomes that have **no intrinsic** order
- We extend logistic regression in a different way
- Y has J levels, from 0 to $J - 1$
- For logistic regression, $P(Y = 1|X) = \text{logit}^{-1}X\beta = \frac{\exp(X\beta)}{1 + \exp(X\beta)}$
- For multinomial logit, we make similar assumptions about $P(Y = j|X)$

$$P(Y = j|X) = \text{logit}^{-1}X\beta_j = \frac{\exp(X\beta_j)}{1 + \sum_{j=1}^J \exp(X\beta_j)} \quad (9)$$

- And for reference group, its

$$P(Y = 0|X) = \text{logit}^{-1}X\beta_j = \frac{1}{1 + \sum_{j=1}^J \exp(X\beta_j)} \quad (10)$$

Multinomial Logit

- For all levels except the reference group, it has its own regression coefficients
- Say we have 7 categories and 4 predictors (each of them is continuous), then in total we will have $6 * 5 = 30$ coefficients
 - $6 = 7 - 1$
 - $5 = 4 + 1$ (plus intercepts)
- Also because we know what $P(Y = j|X)$ looks like for every possible value of Y , we can use MLE to estimate β_j

Interpreting multinomial logit

- Based on the assumptions of multinomial, it is easy to see:

$$\frac{P(Y = j|X)}{P(Y = 0|X)} = \exp(X\beta_j) \quad (11)$$

- Therefore, one unit increase in X leads to $\exp(\beta_j)$ increase in odds ratio of $Y = j$ occurring, relative to $Y = 0$

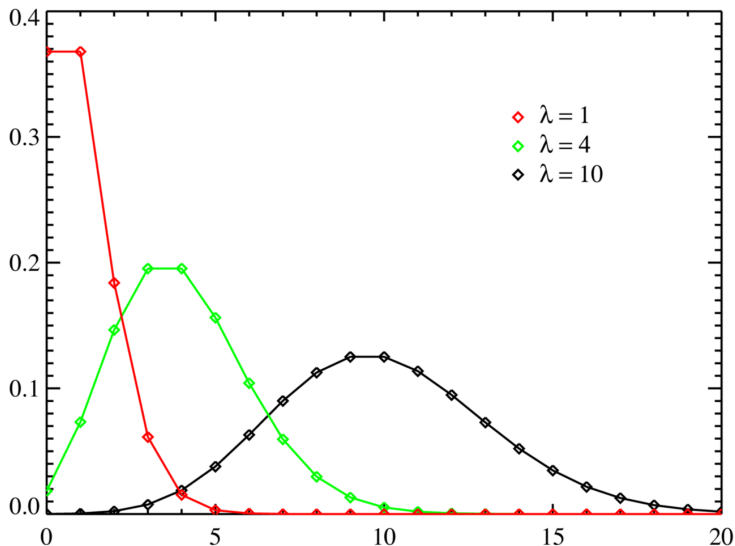
Poisson Distribution

- Example: Y is event count
 - e.g., number of times each person visit a physician)
 - Number of new born / decease in a country
 - Usually small counts are more likely than large counts
- Key difference: Y are **non-negative integers**; in linear regression Y is assumed to be continuous variable between $(-\infty, \infty)$
- Event count usually follows Poisson distribution

$$Pr(X = k) = \frac{\tau^k e^{-\tau}}{k!}$$

- $k! = k(k-1)(k-2) \cdots 1$ is factorial
- Property: $E(X) = V(X) = \tau$

Poisson Distribution



Poisson Regression

- The conditional probability $P(Y|X)$ is assumed to be distributed according to Poisson:

$$P(Y = y|X) = \frac{\exp(-\tau) \tau^y}{y!}, \quad y = 0, 1, 2, \dots \quad (12)$$

$$\tau = \exp(X\beta)$$

- And the conditional expectation $E(Y|X)$ is given by:

$$E(Y|X) = \tau = \exp(X\beta) \quad (13)$$

Poisson Regression (cont'd)

- Why don't we explicitly write $E(\epsilon) = 0$ and $E(\epsilon X) = 0$ as in the Assumption 1 and 2 of linear, logistic and probit regressions?
 - Hint: our assumption of the form of $P(Y|X)$ is very strong
 - It directly gives what $E(Y|X)$ should look like
 - And $E(\epsilon) = 0$ and $E(\epsilon X) = 0$ are essentially the property of $\epsilon = Y - E(Y|X)$
 - So in many textbooks, when introducing generalized linear models, they will omit Assumptions 1 and 2, since it is implied by the assumption of the function form of $P(Y|X)$
- Poisson assumption implies that the data is **heteroskedastic**:

$$\begin{aligned}
 V(\epsilon|X) &= V(Y - E(Y|X)|X) \\
 &= V(Y|X) \\
 &= \exp(X\beta)
 \end{aligned}
 \tag{14}$$

Poisson and Log-Linear model

- Poisson regressions:

$$E(Y|X) = \tau = \exp(X\beta)$$

- An alternative way is to take **log** at both side of the equation

$$\log E(Y|X) = \log(\tau) = X\beta$$

- It means that the **link function** of Poisson regression is **log**
- Sociologists and demographers call $\log E(Y|X) = \log(\tau) = X\beta$ as **log-linear** model

MLE for Poisson regression

1. $P(Y = y|X) = \frac{\exp(-\tau)\tau^y}{y!}; \tau = \exp(X\beta)$

2. Likelihood is: $L = \prod_{i=1}^N \frac{\exp(-\tau_i)\tau_i^{y_i}}{y_i!}$

- and log-likelihood is :

$$\sum_{i=1}^n y_i X_i' \beta - \exp(X_i' \beta) - \log y_i!$$

3. try to maximize by setting the derivative to be 0

$$\frac{\partial l}{\partial \beta} = \sum_{i=1}^n (y_i - \exp(X_i' \beta)) X_i = 0$$

- There is no closed-form solution, unfortunately. Numerical optimization is required.

Interpretation of Poisson Regression

- In log-linear model format:

$$\log E(Y|X) = \log(\tau) = X\beta$$

- One unit increase in X leads to β increase of the average of y in its **log scale**
- In Poisson regression format:

$$E(Y|X) = \exp(X\beta)$$

- One unit increase in X leads to $\exp(\beta) - 1$ increase in Y
- One unit increase in X multiplies the mean of Y by a factor $\exp(\beta)$
- The ratio between the new Y and old Y is $\exp(\beta)$, on average

Over-dispersion of Count Data

- Poisson regression assumes that $P(Y|X)$ follows a Poisson distribution
- Recall that Poisson distribution assumes that the mean and the variance is the same
- Sometimes we have data whose variance is bigger than mean
- E.g., Long, J. Scott. 1990. *The Origins of Sex Differences in Science*. Social Forces. 68(3):1297-1316.
- The outcome is the number of published articles by a Ph.D. student in biochemistry
- The mean number of articles is 1.69 and the variance is 3.71, a bit more than twice the mean.
- Why? There are always super-starts :) and people who publish nothing : (

Zero-inflated Poisson Regression

- One common situation of over-dispersion: there are a lot of zeros in the outcome Y and a few big values, which boosts the variance of outcome
- Example: civil war as outcome.
- Zero-inflated Poisson Regression is designed to address this issue
- It assumes that data has two generating processes
 1. With probability $1 - \lambda$, the data is generated according to Poisson with mean τ
 2. With probability λ , we generate excess zeros.
- The final conditional probability is

$$P(Y = y|X) = \lambda + (1 - \lambda) \frac{\exp(-\tau) \tau^y}{y!}$$

Zero-inflated Poisson Regression (cont'd)

- With the assumptions in the previous slide
- $E(Y|X) = (1 - \lambda)\tau$
- $V(Y|X) = (1 - \lambda)\tau(1 + \tau\lambda)$
- V is bigger than E , of a ratio of $1 + \tau\lambda$
- Essentially, zero-inflated Poisson regression is the mix of two regressions:
 - One Poisson regression, with prob $1 - \lambda$
 - One logistic regressions (0 and all others), with prob λ
 - Each regression has its own coefficients
- So it is a more complex model than negative binomial regression, which adds only one additional parameter

Negative binomial regression

- Another way to deal with over-dispersion: choose a different functional form about $P(Y|X)$

$$P(Y = y|X) = \frac{\Gamma(\alpha + y)}{y! \Gamma(\alpha) (\tau + \alpha)^{\alpha+y}} \quad (15)$$

- And $\tau = \exp(X\beta)$
- Γ is Gamma function, an extension of factorial
- With this more complex parametric assumption
- $E(Y|X) = \tau$ (similar to Poisson regression)
- $V(Y|X) = \tau(1 + \frac{1}{\alpha}\tau)$
- Positive α ensures that variance is bigger than the mean

Other count data model

- Zero truncated regressions
 - Say, the outcome of the length of stay in a hospital, which is at least 1 day
 - Zero-truncated Poisson:
 - Remove the probability $P(y = 0)$ because it's not possible)
 - Re-scale the rest of the probability distribution to make it sums to 1

How do we choose between models?

- Let us use our example of number of published articles by Ph.D. biochemists
- We can choose between three models:
 - Poisson regression
 - Negative binomial regression
 - Zero-inflated Poisson regression
- Decide whether or not to use Poisson regression is relative easier: (Cameron and Trivedi, "Regression-based tests for overdispersion in the Poisson model", *Journal of Econometrics*, 1990)
- Assume $E(Y|X) = \tau$, then
- Null Hypothesis: $V(Y|X) = E(Y|X) = \tau$
- Alternative Hypothesis: $V(Y|X) = \tau + c\tau$
- Cameron and Trivedi's overdispersion test just seeks to examine whether $c = 0$
- (For R users: `dispersiontest` in AER package)

Use Likelihood for Hypothesis Testing

- But how can we compare negative binomial regression vs zero-inflated Poisson regression?
- We can compare **Likelihood** among similar models to choose the best one
- Intuition:
 - Likelihood L represents the joint probability that we observe the entire data, given our parameters
 - Assume we have two models
 - A better model should have larger likelihood

Likelihood Ratio Test

- Define Likelihood Ratio Test Statistics D as:

$$\begin{aligned}
 D &= -2 \log \frac{L_{\text{null}}}{L_{\text{alternative}}} \\
 &= 2(\log L_{\text{alternative}} - \log L_{\text{null}})
 \end{aligned}
 \tag{16}$$

- For comparing models, null model is often the simpler model, and alternative model is often the more complex model
- Null Hypothesis: $D = 0$
- Alternative Hypothesis: $D > 0$
- The bigger the D , the more evidence for the alternative model

Likelihood Ratio Test (cont'd)

- Wilk's Theorem (1938): D has an χ^2 -distribution, with degrees of freedom equal to the difference in number of parameters between alternative model and the null model, if the null model is **nested** within the alternative model
- Nested basically means that the null model can be viewed as a simple case of the alternative model
 - e.g., null is logistic regression with 5 variables; alternative adds another variable
 - null is Poisson; alternative is negative binomial or zero-inflated Poisson
- For non-nested models, Wilk's Theorem does not hold; we need something else (shortly)

Likelihood Ratio Test (cont'd)

- How do express Wilk's Theorem in the p-value language?
 - Say we get a $D = 12$, and the degree of freedom is 2
 - Definition: the probability of obtaining a test statistics that equals to D or higher is approximately $p \iff$ p-value is p
 - $P(D < 12, d.f. = 2) = 0.9975$
 - in R, just type `pchisq(12, 2)`, which is the cumulative probability distribution of D
 - It means that the probability of observing a D smaller than 12 is 0.9975
 - So the probability we observe a D equal to or larger than 12 is $1 - 0.9975 = 0.0025$, which is our **p-value**)

Bias-Variance Trade-Off and Likelihood Ratio Test

- But, a more complex model (adding more parameters) usually can predict more accurately and thus often always have larger likelihood
- AIC: Akaike information criterion (named after Hirotugu Akaike, 1974); reaching balance between predictive power and model complexity
- k is the number of parameters in a model

$$AIC = 2k - 2 \log L \quad (17)$$

Bias-variance trade-off

- AIC wants to balance predictive power and model complexity
- This is a fundamental idea in machine learning
- The idea is called **bias-variance trade-off**
- Recall:
 - We use $g(X)$ to predict Y ;
 - Among all possible $g(X)$, $E(Y|X)$ is the best predictor of Y because it minimizes mean squared error $E[(Y - g(X))^2]$
 - and $\hat{g}(X)$ is estimator of $g(X)$ based on sample data
- Now we compare the mean squared error between Y and its empirical prediction $\hat{g}(X)$, $E[(Y - \hat{g}(X))^2]$

Bias Variance Decomposition

$$E[(Y - \hat{g}(X))^2] = V[Y - E(Y|X)] + [\hat{g}(X) - E(Y|X)]^2 + V(\hat{g}(x))$$

= variance of irreducible error + (prediction bias)² +
prediction variance

- Variance of irreducible error: this only relates to your data;
 - They are irreducible as long as you have selected X ; it will be large if X has nothing to do with Y .
- Prediction bias: relating to your model
 - How current estimator $\hat{g}(X)$ differs from the best predictor $E(Y|X)$
 - OLS is often **bad** at approximating $E(Y|X)$
- Prediction Variance: relating to your model
 - It roughly indicates how varied your predictions can be
 - OLS actually has small prediction variance

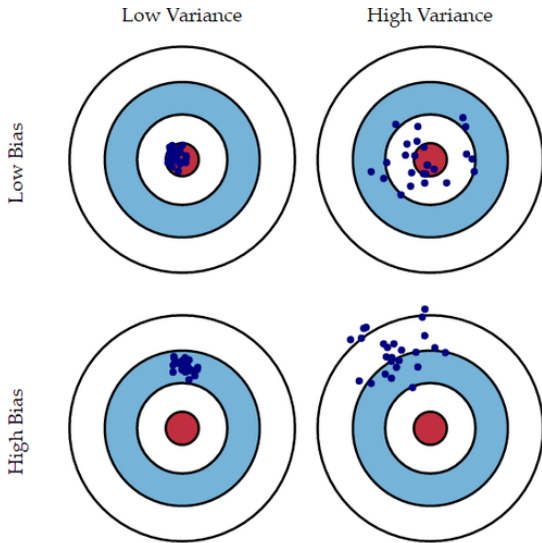
Bias Variance Decomposition (cont'd)

- Variance of Estimator: $V(\hat{g}(x))$
 - It is not the variance of estimated parameters $\hat{V}(\hat{\beta})$; it's the variance of your predicted values!
 - One intuition: the population has 10000 individuals, and each time you sample 100 individuals, and fit an OLS regression.
 - These OLS fitted lines would not vary a lot.
 - But if you use a very complex model, each time predictions can change a lot; thus prediction variances can be high

Bias Variance Trade-off

- To reduce irreducible error: find more predictive X
- The other two quantities relate to your model (estimator):
 - Simple models (like OLS) have large estimator bias, but small estimator variance
 - Complex models have small estimator bias, but large estimator variance

Bias vs Variance (illustration)

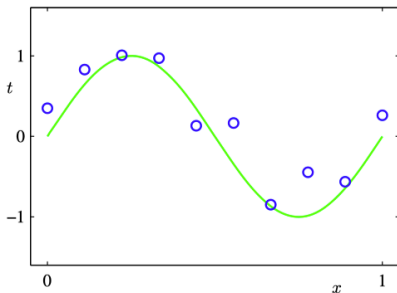


Bias Variance Trade-off (example)

- We have a linear regression with only one variable X , but we add higher order terms

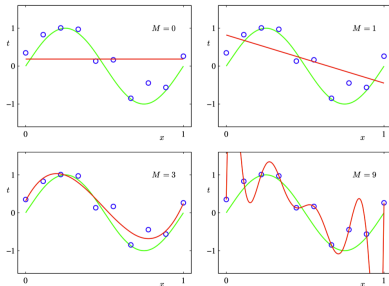
$$Y = \beta_0 + \beta_1 X + \beta_2 X^2 + \cdots + \beta_M X^M$$

- The true M is 3; simulate some data
- Then we try different OLS models by adding more and more high-order terms



Bias Variance Trade-off (cont'd)

- $M = 1$, fits the data very bad (high prediction bias)
- $M = 9$, fits the data so well (small prediction bias), but it is highly sensitive to small changes in observations
 - The prediction on new data can be very bad
 - This is known as **over-fitting**
- $M = 3$, it achieves a good balance between prediction bias and variance
 - And it actually is the correct M



Bias Variance Trade-Off

- Simple model predicts the data very bad (high prediction bias)
- Complex model predicts the data too well (low prediction bias), but it has high estimation variance and is does not generalize well
 - If social science research care about policy implications, generalizability is important.
- Ideal predictive models should balance the prediction bias and variances
- And this principle has been used in many statistics/machine learning applications
 - AIC is one example
 - We will see more next week

Today's Review

Type of Y	Regression to use
Continuous	linear
Binary	logit/probit
Categorical	multinomial logit / ordered logit
Count (integer)	Poisson, negative binomial and zero-inflated

Recommended Readings

- There are many other GLMs (e.g., censored outcome).
- GLM
 - <https://data.princeton.edu/wws509>, Generalized Linear Models course by Germán Rodríguez
 - Powers, Daniel, and Yu Xie. *Statistical methods for categorical data analysis*. Emerald Group Publishing, 2008.
- Machine Learning:
 - Hastie, T., Tibshirani, R., & Friedman, J. (2009). *The elements of statistical learning*. Springer
 - <https://web.stanford.edu/~hastie/ElemStatLearn/>
 - Bias-variance decomposition is discussed in Chapter 2