

# Midterm Review Notes

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## Ordinary Linear Regression

### estimate parameters

In OLS, we estimate the parameters by choosing  $\beta_0, \beta_1$  to minimize the sum of squares.

$$\min_{\beta_0, \beta_1} \sum_i (Y_i - (\beta_0 + \beta_1 X_i))^2$$

That is, the sum of squared differences between our prediction and the true values  $Y_i$ . For OLS, closed form solutions exist for the parameter estimates:

$$\hat{\beta}_1 = \frac{\sum_i (X_i - \bar{X})(Y_i - \bar{Y})}{\sum_i (X_i - \bar{X})^2}$$
$$\hat{\beta}_0 = \bar{Y} - \hat{\beta}_1 \bar{X}$$

### estimate variance

The variance of the noise  $\epsilon$  is estimated based on the parameters  $\beta_0, \beta_1$  and the data set  $\{X_i, Y_i\}$ :

$$\hat{Y}_i = \hat{\beta}_0 + \hat{\beta}_1 X_i$$
$$\hat{\sigma}^2 = \frac{\sum_i (\hat{Y}_i - Y_i)^2}{n - 2}$$

Notice the denominator is  $n - 2$  – the “degrees of freedom”, i.e. the number of data points minus the number of estimated parameters. This estimator is unbiased, meaning  $\mathbb{E}(\hat{\sigma}^2) = \sigma^2$ .

### least squares vs MLE

Least squares and MLE estimation for linear regression are very similar; under the assumption of gaussian noise, they are nearly equivalent, with the only difference being the estimation of the sample variance.

$$\hat{Y}_i = \hat{\beta}_0 + \hat{\beta}_1 X_i$$
$$\hat{\sigma}_{MLE}^2 = \frac{\sum_i (\hat{Y}_i - Y_i)^2}{n}$$

Notice the difference between the *MLE* estimator of the sample variance and the least squares estimator of the sample variance. The *MLE* estimator is biased i.e.  $\mathbb{E}(\hat{\sigma}^2) \neq \sigma^2$ ; however, the bias term decays to zero as the number of data points  $n$  increases.

## residuals and their properties

The residuals are defined as  $\varepsilon_i = \hat{y}_i - y_i$ . Of course, the loss is defined as the sum of squared residuals:

$$\sum_i \varepsilon_i = 0$$
$$\sum_i \varepsilon_i \cdot X_i = 0$$

From the above properties, it's very simple to show the following additional properties:

$$\sum_i \hat{Y}_i = \sum Y_i$$
$$\sum_i \varepsilon_i \cdot \hat{Y}_i = 0$$

Typically, we assume the noise is gaussian, in which case the residuals should be roughly normally distributed. We can check this via a q-q plot.

## Confidence intervals vs Prediction Intervals

A confidence interval for  $\beta_1$  is an interval-valued function of the data; intuitively, a confidence interval is supposed to provide a range of reasonable values for the parameter of interest.

- Here's how we construct a confidence interval for  $\beta_1$ :

$$\hat{\beta}_1 + \hat{\sigma} \cdot t\left(\frac{1-\alpha}{2}, n-2\right)$$
$$\hat{\beta}_1 + \hat{\sigma} \cdot t\left(\frac{1+\alpha}{2}, n-2\right)$$

- Interpretation: What is meant by “95%”? The model assume the data is distributed like as follows

$$Y_i = \beta_0 + \beta_1 X_i + \varepsilon_i$$

Confidence intervals are computed with respect to fixed  $X_i$  but varying  $\varepsilon_i$ . So when we refer to a “95% confidence interval for  $\beta_1$ ”, we mean over 100 datasets that share the same  $\{X_i\}$ , roughly 95 of the confidence intervals contain the true parameter  $\beta_1$ .

- Prediction intervals: whereas a confidence interval estimates a range of plausible values of a parameter of interest, e.g.  $E(Y_i|X_i)$ ; a prediction interval, on the other hand, tries to estimate a range of plausible variables for a single out of sample point e.g. the a new  $Y_i$ .
- The difference lies in the fact that a prediction interval has to account for the variance in the distribution of  $Y_i$ .

The distribution for an single out of sample  $\hat{Y}_i$  would be gaussian with the following mean and variance:

$$E[\hat{Y}_i] = \beta_0 + \beta_1 X_i$$
$$\sigma^2[\hat{Y}_i] = \hat{\sigma}^2 \left[ 1 + \frac{1}{n} + \frac{(X_i - \bar{X})^2}{\sum_j (X_j - \bar{X})^2} \right]$$

Notice the difference between the above and the *confidence* interval for  $Y_i$ :

$$E[\hat{Y}_i] = \beta_0 + \beta_1 X_i$$
$$\sigma^2[\hat{Y}_i] = \hat{\sigma}^2 \left[ \frac{1}{n} + \frac{(X_i - \bar{X})^2}{\sum_j (X_j - \bar{X})^2} \right]$$

## Diagnostics (R squared)

$$R^2 = 1 - \frac{RSS}{TSS} = 1 - \frac{\sum_i (\hat{y}_i - y_i)^2}{\sum_i (\bar{y} - y_i)^2}$$

## Multiple Variables

If we have more than one predictor, we can still find closed form solutions for all our coefficients. The form is now:

$$X_i = (1, x_{i,1}, x_{i,2}, \dots, x_{i,k})$$
$$\hat{\beta} = (X^T X)^{-1} X^T Y$$

Keep in mind that adding a variable to a regression can change ALL of the regression coefficients.

## Qualitative

- Qualitative variables (e.g. species, race) can be represented by indicators i.e. a variable that takes value 1 if a condition is met and zero otherwise.
- Keep in mind R chooses a reference point, meaning if we have  $k$  categories in a variable, R will create  $k - 1$  indicator variables and the  $k$ th category will be represented by all the others taking a value of zero.

## AIC

- The AIC, short for Akaike-Information Criterion, is a measure of fit of a model that balances goodness of fit (i.e. log likelihood) against the number of variables in the model  $q$

$$L(\beta) = \Pi_i P(y_i | x_i, \beta)$$
$$\text{AIC}(\text{lmod}) = 2 \cdot \log(L(\hat{\beta})) + q$$

- This is a very general and powerful way of comparing models.

## Bernoulli/Binomial Data

A word on terminology. Make sure not to confuse bernoulli with binomial! They're closely related but different concepts.

- A bernoulli random variable models a binary outcome, e.g. “yes” or “no”, “success” or “failure”, “male” or “female” etc. It takes values 0 or 1. A bernoulli random variable has one parameter – the “success probability”  $p$ . The distribution is:

$$\text{bernoulli}(p) = p^k (1 - p)^{1-k}$$

- A binomial random variable models the number of successes in  $m$  independent and identically distributed bernoulli random variables. It has parameters  $m$  the total number of trials,  $p$  the success probability of any given trial.
- A binomial random variable can take values from 0 to  $m$ , inclusive.

$$P(Y = k) = \binom{m}{k} \cdot p^k \cdot (1 - p)^{m-k}$$

- Binomial data is data where the response variable is a binomial random variable. For example, the input might be economic data at the county-level, and the response might be the number of people in the county who voted democrat in the most recent election (out of the total number of people in the county).

- Note in binomial regression we need to know both the parameter  $m_i$  and the number of successes  $k_i$  for each data point  $i$ .
- Bernoulli data can be thought of as binomial data where the  $m = 1$

## link functions

The problem with using basic linear regression to model binomial and bernoulli data is that the linear predictor  $\sum_i \beta_i * x_i$  is unbounded – it isn't guaranteed to fall in  $0, 1$  or  $0, m$ .

## predictions

- Pass the odds through the inverse-logit function to get the predicted probability  $p_i$

## odds, probability

- Odds is a different way of parametrizing of probability. The odds of an event with probability  $p$  is  $p/1 - p$ .
- The odds concept is useful in thinking about bernoulli, binomial regression because the the linear predictor is the log of the odds:

$$\log\left(\frac{p_i}{1 - p_i}\right) = \beta_0 + \beta_1 x_1 + \dots + \beta_k x_k$$

## risk ratio vs odds ratio

- An odds ratio is a ratio of the odds of two events  $odds(A)/odds(B)$ .
- Because the linear predictor models the log odds for each data point, the success odds of data point  $i$  are:

$$odds_i | X_i = e^{\beta_0} \prod_k e^{\beta_k X_{i,k}}$$

- The odds ratio of success given two levels of predictor  $X_{i,k}$ , say  $s, t$  is

$$odds\_ratio_i(X_i = s, X_i = t) = e^{\beta_k(s-t)}$$

- A risk ratio is the ratio probabilities given two events

$$risk\_ratio(A, B) = P(A)/P(B)$$

## goodness of fit - We can assess goodness of fit with a calibration curve, which plots *empirical<sub>p</sub>probability* against *pred<sub>p</sub>probability*. Empirical probabilities are determined via binning.

- Ideally the curve would follow the diagonal  $y = x$

## Diagnostics (Pearson Residuals)

## different scoring functions

## confusion matrix and properties (sensitivity, specificity, PPV, accuracy, NPV)

## comparing nested models

## overdispersion

## f statistics

## quasibinomial