

# MODEL SELECTION

Brian Chung

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## LESS THAN PERFECT BOOK RAFFLE

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```
students = ['Doug', 'Dylan', 'Logan', 'Thomas']  
np.random.choice(students)
```

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## FROM LAST TIME

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### **Questions:**

- Other variants of K Means (Why k means or k medians)
- Non-convex data for K Means
- Choosing enough centroids
- How many features to use in k means?

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## **MODEL SELECTION AGENDA**

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**I. LINEAR MODEL SELECTION**

**II. CRITERION**

**III. PROJECT / FREE TIME**

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## **MODEL SELECTION**

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# **I. LINEAR MODEL SELECTION**

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## **CHOOSING FACTORS IN A LINEAR MODEL**

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**Q: In linear regression, what is the purpose of feature selection?**

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## **CHOOSING FACTORS IN A LINEAR MODEL**

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**Q: In linear regression, what is the purpose of feature selection?**

**A: Better prediction accuracy as well as model interpretability**

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## **CHOOSING FACTORS IN A LINEAR MODEL**

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- Subset Selection (**Today**)

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**A:** There are a couple methods, including one such method we’ve already learned

- **Shrinkage (Ridge Regression, LASSO)**
- **Subset Selection (Today)**
- **Dimension Reduction (Future: Principle Components Analysis)**

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## **SUBSET SELECTION**

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**There are a few main methods used in subset selection of features:**

- **Best Subset Selection**
- **Forward Stepwise Selection**
- **Backward Stepwise Selection**

**As well as a few different metrics used in measuring “goodness”:**

- **AIC**
- **BIC**
- **Adjusted  $R^2$**

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## **BEST SUBSET SELECTION**

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**Assume you begin with  $p$  potential features, of which you believe  $k$  of these features are pertinent to your problem.**

**In best subset selection, we iterate through all potential subsets of ( $p$  choose  $k$ ) factors, evaluating them using a common metric (AIC, BIC, Adjusted  $R^2$ )**

**Ultimately, we choose the best set of  $k$  factors that have been tested through cross validation**

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## BEST SUBSET SELECTION – ALGORITHM

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1. Define  $M_0$  to be the *null model*, indicating a model with zero features. The model predicts the response average for every observation
2. For  $k = 1, 2, \dots, p$ :
  1. Fit all  $(p \text{ choose } k)$  models that contain exactly  $k$  features
  2. Pick the model among these  $(p \text{ choose } k)$  models that has the highest adjusted  $R^2$ . Call this model  $M_k$
3. Amongst the models from  $M_0$  to  $M_k$ , choose the single best model with the highest cross validated adjusted  $R^2$ .

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## FORWARD STEPWISE SELECTION

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That's great...except what if you have 100 features? That's  $(100 \text{ choose } 1) + (100 \text{ choose } 2) + (100 \text{ choose } 3) + \dots + (100 \text{ choose } 100)$  different subsets to try. It's computationally extremely unfeasible for large values of  $p$ .

So, what if we try an augmented approach. Let's start with a blank model, and choose the best single feature to add onto this model. Then, let's choose the next best feature to add onto this new model. Almost like cherry picking the best "next additional feature" to add.

Ultimately, we'll have  $M_0, M_1, M_2, \dots, M_k$  models, where each is building upon the last.



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## STEPWISE FORWARD SELECTION – ALGORITHM

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1. Define  $M_0$  to be the *null model*, indicating a model with zero features. The model predicts the response average for every observation
2. For  $k = 1, 2, \dots, p$ :
  1. Fit all models that augment  $M_{k-1}$  with a single additional feature.
  2. Pick the model among these models with the highest adjusted  $R^2$  and call this  $M_k$
3. Amongst the models from  $M_0$  to  $M_k$ , choose the single best model with the highest cross validated adjusted  $R^2$ .

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## BACKWARD STEPWISE SELECTION

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1. Define  $M_p$  to be the *full model*, indicating a model with every feature. The model predicts based on the fitted coefficients.
2. For  $k = p, p-1, p-2, \dots, 1$ :
  1. Fit all models that augment  $M_k$  by removing a single feature.
  2. Pick the model among these models with the highest adjusted  $R^2$  and call this  $M_k$
3. Amongst the models from  $M_0$  to  $M_k$ , choose the single best model with the highest cross validated adjusted  $R^2$ .

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## **FORWARD OR BACKWARDS?**

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**Forward and backward selection will often give similar but slightly different results.**

**Personally, I choose forward selection when I want the simplest possible model, and choose backwards selection when I believe the 'true' model is close to what I have right now**

**The main issue with either stepwise selection methods is that of optimality. We use a heuristic to choose features, but these may not be the most ideal set of features**

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## **MODEL SELECTION**

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**I. LINEAR MODEL SELECTION**

**II. CRITERION**

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## OTHER SELECTION CRITERION

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Recall that the *adjusted  $R^2$*  penalizes the goodness of fit for adding many features or variables.

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However, for models that are fit using maximum likelihood (this include linear regression as well as logistic regression), you can also perform model selection by minimizing the Akaike Information Criterion (AIC) or the Bayesian Information Criterion (BIC).

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However, for models that are fit using maximum likelihood (this include linear regression as well as logistic regression), you can also perform model selection by minimizing the Akaike Information Criterion (AIC) or the Bayesian Information Criterion (BIC).

Information criterion essentially measure the "loss of information" when moving from some true model "f" to a model "g". While we do not know the true model "f", we can compare two different models  $g_1$  and  $g_2$  using AIC or BIC.

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## AKAIKE INFORMATION CRITERION

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The AIC is defined as -2 times the log likelihood of the model plus 2 times the number of features

Given two models with different AIC values, the preferred model is the one with a lower AIC

$$\text{AIC}(\mathbf{M}) = -2 \ln(L) + 2k$$



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$$AIC(M) = -2 \ln(L) + 2k$$

AIC(M): Akaike Information Criterion for model M  
ln(L): Log likelihood for model and data  
k: number of features

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## BAYESIAN INFORMATION CRITERION

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The BIC is defined as -2 times the log likelihood of the model plus the natural log of the number of observations times the number of features

Given two models with different BIC values, the preferred model is the one with a lower BIC

$$\text{BIC}(M) = -2 \ln(L) + \ln(N) * k$$

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## BAYESIAN INFORMATION CRITERION

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The BIC is defined as -2 times the log likelihood of the model plus the natural log of the number of observations times the number of features

Given two models with different BIC values, the preferred model is the one with a lower BIC

$$\text{BIC}(M) = -2 \ln(L) + \ln(N) * k$$

BIC(M): Bayesian Information Criterion for model M

$\ln(L)$ : Log likelihood for model and data

$\ln(N)$ : log of number of observations

k: number of features

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## AIC VS BIC

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The equations are similar, but are born from different justifications.

- AIC believes that all models are incorrect, and only looks for the best approximation.
- BIC believes that the “true model” is a simplification of the current set of features.
- Generally, AIC tends to favor overfitting, while BIC tends to favor underfitting

$$\text{AIC}(\mathbf{M}) = -2 \ln(L) + 2k$$

$$\text{BIC}(\mathbf{M}) = -2 \ln(L) + \ln(N) * k$$

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## SO WHICH DO WE USE IN MODEL SELECTION?

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Generally speaking, when using AIC vs BIC for stepwise selection, both should yield similar results.

*Statsmodels* provides all the metrics including AIC and BIC.

Let's look at a revised example of Forward Stepwise Selection using the AIC (you can replace AIC with BIC). Just remember that now, instead of maximizing adjusted  $R^2$  we want to minimize AIC or BIC.

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## STEPWISE FORWARD SELECTION – ALGORITHM

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2. For  $k = 1, 2, \dots, p$ :
  1. Fit all models that augment  $M_{k-1}$  with a single additional feature.
  2. Pick the model among these models with the **lowest AIC** and call this  $M_k$
3. Amongst the models from  $M_0$  to  $M_k$ , choose the single best model with the **lowest cross validated AIC**.

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## **MODEL SELECTION**

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**I. LINEAR MODEL SELECTION**

**II. CRITERION**

**III. PROJECT / REVIEW / ANYTHING**

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**THAT'S IT!**

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▸ Exit Tickets: DAT1 - Lesson 10 - Model Selection