

# Model Selection & Regularization

## Lecture 7

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### Recall: Linear Models and Least Squares

$$Y = \beta_0 + \beta_1 X_1 + \beta_2 X_2 + \dots + \beta_p X_p + \epsilon \quad \text{RSS} = \sum_{i=1}^n (y_i - \hat{\beta}_0 - \sum_{j=1}^p \hat{\beta}_{j1} x_{ij})^2$$

Model with all available predictor variables is commonly referred to as **the full model**

#### Issues:

- predictive accuracy
- model interpretability

#### Solutions:

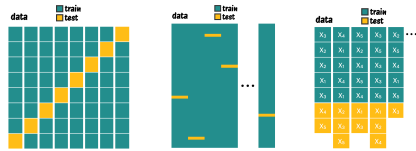
- select **subset** of predictors
- consider **extension to the least squares solution** of full model

### Part I - Variable Subset Selection

## Model Selection Criteria: Validation by Prediction Error

**Last week:** how to use cross validation to choose a set of predictors by directly estimate prediction error using cross-validation techniques

$$\text{e.g. } \text{MSE} = \frac{\text{RSS}}{n} \quad \text{RMSE} = \sqrt{\frac{\text{RSS}}{n}} \quad R^2 = 1 - \frac{\text{RSS}}{\text{TSS}}$$



**Now:** indirectly estimating test performance using an approximation

## Model Selection Criteria

### Four ways to estimate test performance using an approximation

Full model has  $p$  predictors

RSS is the residual sum of squares for model with  $d$  predictors

$\hat{\sigma}^2 = \text{RSS}_p / (n - p - 1)$  is an estimate of the error variance for full model

#### 1. Mallows's $C_p$ criterion:

For a given model with  $d$  (out of the  $p$  available) predictors

$$C_p = \frac{1}{n} (\text{RSS} + 2d\hat{\sigma}^2)$$

we are penalizing models of higher dimensionality (larger  $d$ , greater penalty)

$\Rightarrow$  choose the model which has **minimum**  $C_p$

## Model Selection Criteria

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#### 2. Akaike Information Criterion (AIC)

For linear models: equivalent to Mallows's  $C_p$  (proportional to)

$$AIC = \frac{1}{n\hat{\sigma}^2} (\text{RSS} + 2d\hat{\sigma}^2)$$

we are penalizing models of higher dimensionality (larger  $d$ , greater penalty)

$\Rightarrow$  choose the model which has **minimum**  $AIC$

## Model Selection Criteria

### Four ways to estimate test performance using an approximation

Full model has  $p$  predictors

RSS is the residual sum of squares for model with  $d$  predictors

$\hat{\sigma}^2 = \text{RSS}_p / (n - p - 1)$  is an estimate of the error variance for full model

#### 3. Bayesian Information Criterion (BIC)

For linear models: equivalent to Mallows's  $C_p$  (proportional to)

$$BIC = \frac{1}{n\hat{\sigma}^2} (\text{RSS} + \underbrace{\log(n)d\hat{\sigma}^2}_{\text{heavier penalty}})$$

we are penalizing models of higher dimensionality (larger  $d$ , greater penalty)

$\Rightarrow$  choose the model which has **minimum BIC**

## Model Selection Criteria

### Four ways to estimate test performance using an approximation

Full model has  $p$  predictors

RSS is the residual sum of squares for model with  $d$  predictors

$\hat{\sigma}^2 = \text{RSS}_p / (n - p - 1)$  is an estimate of the error variance for full model

#### 4. Adjusted R-squared value

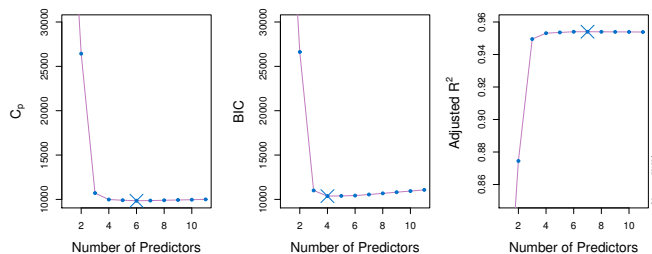
Adjust the regular  $R^2$  by taking into account number of predictors

$$\text{Adjusted-}R^2 = 1 - \frac{\text{RSS}/(n - d - 1)}{\text{TSS}/(n - 1)}$$

$\Rightarrow$  choose the model which has **maximum Adjusted- $R^2$**

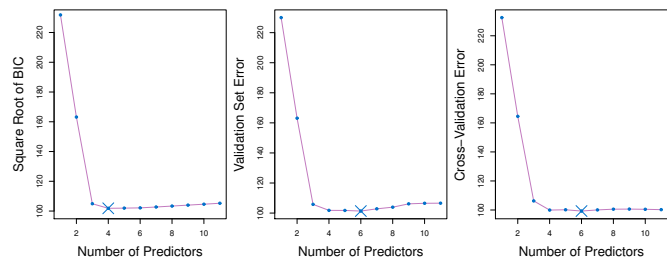
## Model Selection Criteria

### Four ways to estimate test performance using an approximation



## Model Selection Criteria

...and compared to cross validation



## Model Search Methods

### Best Subset Selection

- Let  $M_0$  denote null model which contains no predictors. This model simply predicts the of the response for each observation.
- For  $k = 1, 2, \dots, p$ 
  - Fit all  $\binom{p}{k}$  models that contain exactly  $p$  predictors
  - Pick the best among these  $\binom{p}{k}$  models and call it  $M_k$ .  
Here, best is defined as having the smallest RSS or largest  $R^2$
- Select a single best model from among  $M_0, M_1, \dots, M_p$  using cross validated prediction error,  $C_p$  (AIC), BIC, or Adjusted- $R^2$

requires training  $2^p$  models

#### Example

$p = 3$

$M_0$ : intercept only (null)

$C_1$ :  $X_1$   $X_2$   $X_3$



$M_1$

lowest training RSS within  $C_1$

$C_2$ :  $X_1, X_2$   $X_1, X_3$   $X_2, X_3$



$M_2$

lowest training RSS within  $C_2$

$M_3$ : full model with



## Model Search Methods

### Forward Stepwise Selection

- Let  $M_0$  denote null model which contains no predictors.
- For  $k = 1, 2, \dots, p - 1$ 
  - Consider all  $p - k$  models that augment the predictors in  $M_k$  with one additional predictor
  - Choose the best among these  $p - k$  models and call it  $M_{k+1}$ .  
Here, best is defined as having the smallest RSS or largest  $R^2$
- Select a single best model from among  $M_0, M_1, \dots, M_p$  using cross validated prediction error,  $C_p$  (AIC), BIC, or Adjusted- $R^2$

requires training  $1 + \frac{p(p+1)}{2}$  models

#### Example

$p = 3$

$M_0$ : intercept only (null)

$C_1$ :  $X_1$   $X_2$   $X_3$



$M_1$

lowest training RSS within  $C_1$

$C_2$ :  $X_1, X_2$   $X_2, X_3$



$M_2$

lowest training RSS within  $C_2$

$M_3$ : full model with



# Model Search Methods

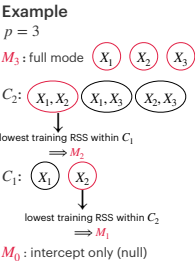
## Backward Stepwise Selection

1. Let  $M_p$  denote full model which all predictors.
2. For  $k = p, p - 1, p - 2, \dots, 1$ 

▸ Consider all  $k$  models that contain all but one of the predictors in  $M_k$ , for a total of  $k - 1$  predictors

▸ Choose the best among these  $k$  models and call it  $M_{k-1}$ .  
Here, *best* is defined as having the smallest RSS or largest  $R^2$
3. Select a single best model from among  $M_0, M_1, \dots, M_p$  using cross validated prediction error,  $C_p$  (*AIC*), *BIC*, or Adjusted- $R^2$

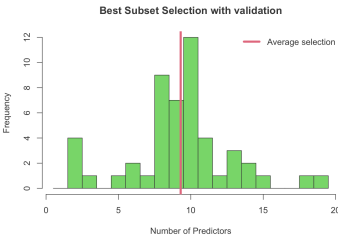
requires training  $1 + \frac{p(p+1)}{2}$  models



# Model Search Methods

## Best Subset Selection

validation approach based on 50 different seeds and storing number of predictors in selected model each time



[plot is made based on the 'hitters' data se used in this week's practical in ISLR2]

## Part II - Shrinkage

## Shrinkage Methods

Before: **Discrete model search methods**

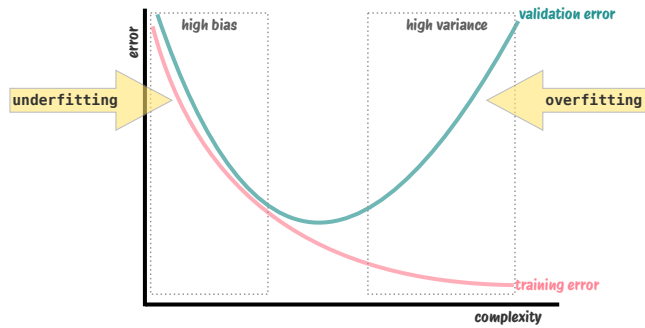
$\underbrace{\text{model fit} + \text{penalty on model dimensionality}}_{\text{RSS}}$

Now: **Continuous model search method** (also faster)

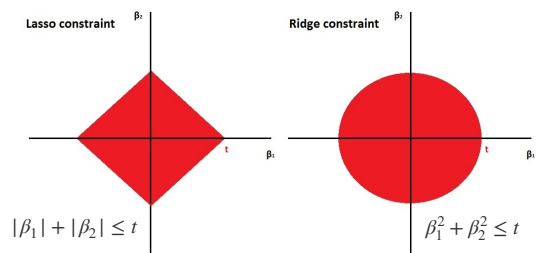
$\underbrace{\text{model fit} + \text{penalty on size of coefficients}}_{\text{RSS}}$

this is called **penalized** or **regularized regression**

## Bias Variance Trade-Off



## Ridge and Lasso Regression: The Constraints



# Ridge Regression

Least Squares produces estimates by minimizing

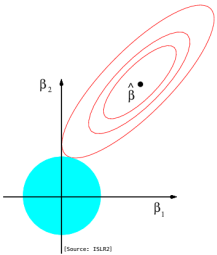
$$RSS = \sum_{i=1}^n (y_i - \hat{\beta}_0 - \sum_{j=1}^p \hat{\beta}_j x_{ij})^2$$

Ridge regression instead minimizes

$$\underbrace{\sum_{i=1}^n (y_i - \hat{\beta}_0 - \sum_{j=1}^p \hat{\beta}_j x_{ij})^2}_{\text{model fit}} + \underbrace{\lambda \sum_{j=1}^p \beta_j^2}_{\text{penalty}} = RSS + \lambda \sum_{j=1}^p \beta_j^2$$

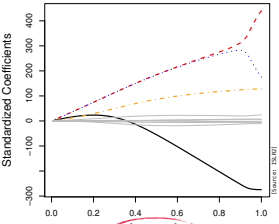
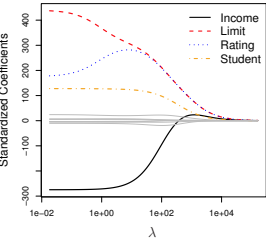
where  $\lambda \geq 0$  is the tuning parameter controlling trade off between model fit and size of coefficients ( $\lambda \rightarrow \infty, \hat{\beta}_j \rightarrow 0$ )

ridge uses  $\ell_2$  penalty



# Ridge Regression

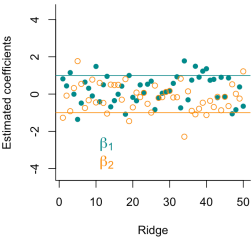
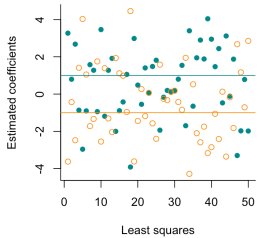
Regularization Paths



$\ell_2 \text{ norm} = \|\beta\|_2 = \sqrt{\sum_{j=1}^p \beta_j^2}$

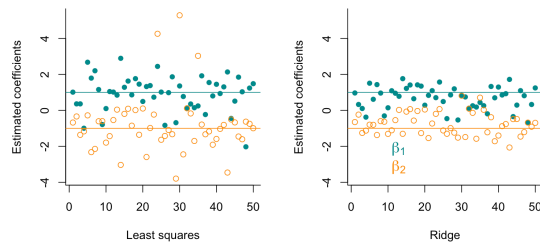
# Ridge Regression

Advantage 1: Multicollinearity (a simulation study)



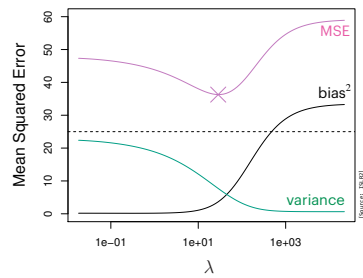
## Ridge Regression

Advantage 2: When  $p$  is close to  $n$  (a simulation study)



## Ridge Regression

Bias-Variance Trade Off



## Lasso Regression

Least Absolute Shrinkage and Selection Operator

Least Squares produces estimates by minimizing

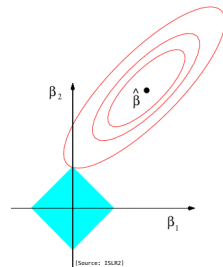
$$RSS = \sum_{i=1}^n (y_i - \hat{\beta}_0 - \sum_{j=1}^p \hat{\beta}_j x_{ij})^2$$

Lasso regression instead minimizes

$$\underbrace{\sum_{i=1}^n (y_i - \hat{\beta}_0 - \sum_{j=1}^p \hat{\beta}_j x_{ij})^2}_{\text{model fit}} + \underbrace{\lambda \sum_{j=1}^p |\beta_j|}_{\text{penalty}} = RSS + \lambda \sum_{j=1}^p |\beta_j|$$

where  $\lambda \geq 0$  is the tuning parameter controlling trade off between model fit and size of coefficients ( $\lambda \rightarrow \infty, \hat{\beta}_j = 0$ )

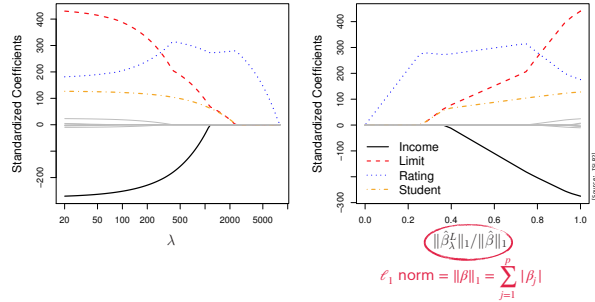
lasso uses  $\ell_1$  penalty





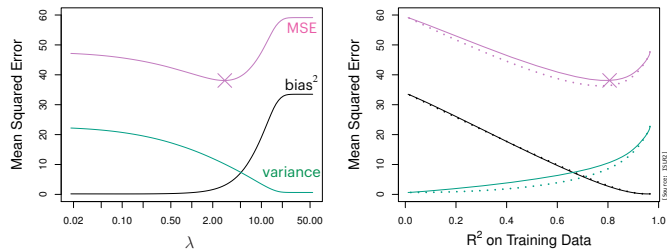
## Lasso Regression

### Regularization Paths



## Lasso Regression

### Bias-Variance Trade Off



## Ridge vs. Lasso Regression

- Both ridge and lasso are convex optimization
- The ridge solution exists in closed form
- Lasso does not have closed form solution, but very efficient optimization algorithms exist

### When to choose which?

- When the actual data-generating mechanism is **sparse** lasso has the advantage
- When the actual data-generating mechanism is **dense** ridge has the advantage

**Sparse mechanisms:** Few predictors are relevant to the response → good setting for lasso regression

**Dense mechanisms:** A lot of predictors are relevant to the response → good setting for ridge regression

- Also depends on:
  - Signal strength (the magnitude of the effects of the relevant variables)
  - The correlation structure among predictors
  - Sample size  $n$  vs. number of predictors  $p$

## Ridge vs. Lasso Regression

### Ridge

- + Reduces Multicollinearity
- + Continuous Shrinking
- + Stable Solutions
- + Computationally Efficient

- No variable selection
- Interpretability
- Sensitive to scale

### Lasso

- + Variable selection
- + Sparse models
- + Improves interpretability
- + Particularly useful for when  $p > n$

- Collinearity issues
- Bias in coefficients ( $\ell_1$  penalty is harsher)
- Computationally intensive

## $\lambda$ Tuning

### • K-fold Cross Validation

1. Choose the number of folds  $K$
2. Split the data accordingly into training and testing sets.
3. Define a grid of values for  $\lambda$
4. For each  $\lambda$ , calculate the validation MSE within each fold
5. For each  $\lambda$ , calculate the overall cross-validation MSE
6. Locate under which  $\lambda$  cross-validation MSE is minimized, i.e. `minimum_cv  $\lambda$`

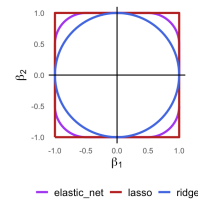
- Packages such as `glmnet` do this automatically



## Hybrid Approach: Elastic Nets

$$\text{RSS} + \underbrace{\lambda_1 \sum_{j=1}^p \beta_j^2}_{\text{"ridge"}} + \underbrace{\lambda_2 \sum_{j=1}^p |\beta_j|}_{\text{"lasso"}}$$

$\lambda_1$  and  $\lambda_2$  are regularization parameters controlling the strength of the penalties



- Helps stabilize the solution when predictors are correlated
- Shrinks some coefficients to zero, enabling feature selection
- Particularly useful for high-dimensional datasets with correlated predictors

## Part III- Dimensionality Reduction

another strategy which aims to reduce dimensionality **before** applying LS

create  $q$  transformed variables which are linear combinations of the original predictors ( $q < p$ )

we return to this during our PCA lecture...

## Part IV- Transformations: next week!

extensions to the regression model when the best straight line doesn't quite work!

[illegible]