STA 35C: Statistical Data Science III

Lecture 16: Linear Model Selection – Regularization Methods

Dogyoon Song

Spring 2025, UC Davis

Agenda

- Last time: Model selection via subset selection
 - Best subset selection: identify relevant predictors among many
 - Stepwise selection: A computationally more tractable alternative (greedy alg)
- Today: Regularization
 - Overview: what regularization is & why it can help
 - Two main examples in linear regression
 - Ridge regression
 - The lasso (least absolute shrinkage and selection operator)

Recap: Subset selection

Best subset selection:

- Exhaustively search all 2^p subsets
- Pick the best model for each size k, then choose among them (via R_{adj}^2 or CV)
- Feasible only for smaller p (high computational cost)

Forward stepwise selection:

- Greedy approach: add one predictor at a time
- Complexity: $\mathcal{O}(p^2)$ instead of 2^p
- May miss the global optimum if local decisions are suboptimal

Backward stepwise selection:

- Greedy approach: remove one predictor at a time
- Similar pros/cons as forward stepwise

Recap: Subset selection – Comparison of search paths

An illustrative example with p = 4:

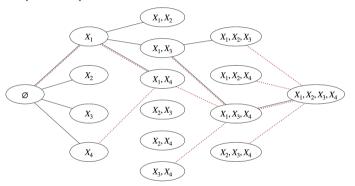


Figure: Illustration of forward stepwise (black solid path; successively adding $X_1 \to X_4 \to X_3 \to X_2$) and backward stepwise (red dashed path; removing $X_2 \to X_3 \to X_4 \to X_1$). Best subset selection checks all 2^p possibilities; all three can yield different outcomes.

For more R examples, see the discussion section slides on Canvas

Recap: Subset selection – Summary

• Summary:

- Goal: identify a relevant subset of predictors
- **Procedure**: evaluate subsets (all or partial), then pick best via R_{adj}^2 , CV, etc.
- BSS is exhaustive but expensive; stepwise is faster
- Typically refit the final chosen "best" subset with least squares

• Advantages:

- Direct variable selection: some $\beta_j = 0$ (excluded)
- Straightforward implementation and intuitive interpretation

• Disadvantages:

- Even stepwise can be costly if p is very large
- Instability: small changes in the data can alter the chosen "best" subset
- \Rightarrow **Regularization** can handle large p, offering stable estimates without discrete exclusion

Regularization: What and why?

Recall least squares: find parameters $\hat{\beta}_0, \dots \hat{\beta}_p$ that minimize

RSS =
$$\sum_{i=1}^{n} (y_i - \hat{y}_i)^2$$
 where $\hat{y}_i = \hat{\beta}_0 + \hat{\beta}_1 x_{i1} + \dots + \hat{\beta}_{ip} x_p$

- Challenges:
 - If p is large or data are noisy, least squares solutions can be unstable
 - Overfitting (huge variance in $\hat{\beta}$) or no unique solution if p > n
- Idea: modify the objective by adding a penalty on β_j 's to stabilize fitting
 - Balance data fidelity vs. "simplicity" (by favoring smaller β_j)
 - This approach is called *regularization* (or *shrinkage*)

We will learn two prominent regularization techniques for linear regression:

- Ridge regression (ℓ_2 penalty)
- Lasso (ℓ_1 penalty)

Ridge regression: 1) Formulation

Ridge regression: Find $\hat{\beta}_0, \dots \hat{\beta}_p$ that minimize

$$\underbrace{\sum_{i=1}^{n} \left(y_i - \beta_0 - \sum_{j=1}^{p} \beta_j x_{ij} \right)^2}_{\text{RSS}} + \lambda \underbrace{\sum_{j=1}^{p} \beta_j^2}_{\text{penalty}}$$

- $\lambda \geq 0$ is a tuning parameter
- Each λ yields a different set of coefficient estimates $\hat{\beta}^R_{\lambda}$

Remarks:

- No penalty on β_0 (the intercept)
- As $\lambda \to 0$, ridge regression \to standard least squares
- As λ grows, $\beta_{1:p} = (\beta_1, \dots, \beta_p)$ shrinks toward 0
 - Reduces variance of $\beta_{1:p}$ but increases bias

Ridge regression: 2) Effects of scaling

The least squares coefficients are scale equivariant:

- Multiplying X_j by constant c scales $\hat{\beta}_j$ by 1/c
- Regardless of scaling, $X_i \hat{\beta}_i$ remains the same, not affecting other coefficients

Ridge regression is *not* scale-equivariant:

- Rescaling one predictor can affect others through the penalty term
- Hence, $X_j \hat{\beta}_{j,\lambda}^R$ depends on both λ and predictor scaling

Therefore, it is recommended to standardize predictors before ridge via

$$ilde{x}_{ij} = rac{x_{ij}}{s_{ij}}$$
 where $s_{ij}^2 = rac{1}{n} \sum_{i=1}^n (x_{ij} - ar{x}_j)^2$ and $ar{x}_j = rac{1}{n} \sum_{i=1}^n x_{ij}$

Ridge regression: 3) Credit dataset example

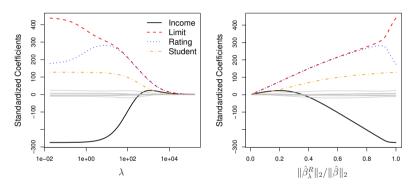


Figure: Standardized ridge coefficients for Credit, plotted vs. λ and $\|\hat{\beta}_{\lambda}^{R}\|_{2}/\|\hat{\beta}\|_{2}$ [JWHT21, Figure 6.4]. (Note: $\|v\|_{2} = \sqrt{v_{1}^{2} + \dots + v_{p}^{2}}$.)

- As λ increases, all β_i shrink toward 0 (none exactly zero though)
- If variables are correlated, ridge shrinks them together in a "group" manner

Ridge regression: 4) Advantages over least squares

Ridge's advantage rooted in **Bias-variance tradeoff**:

- $\lambda = 0$: no bias but high variance
- Larger λ : more bias but lower variance

When ridge can be most helpful:

- If least squares has high variance (e.g. $p \approx n$ or predictors are collinear)
- If data are noisy and β_i can fluctuate a lot
- Ridge still works even if p > n, producing a unique solution

Computational advantages of ridge:

- No need for 2^p model fits (as in best subset); for any λ , ridge requires only a single fit
- Indeed, we can compute solutions for all λ at near the same cost as one OLS fit

Ridge regression: 5) Visualization of bias-variance tradeoff

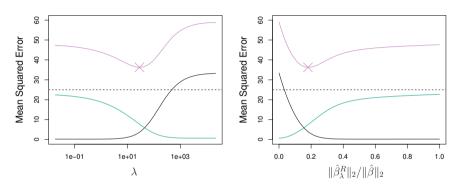


Figure: Bias-variance tradeoff in ridge for a simulated data set (p = 45, n = 50). Shown are squared bias (**black**), variance (**green**), and test MSE (**purple**) vs. λ [JWHT21, Figure 6.5].

- At $\lambda = 0$, there is no bias but high variance
- Increasing λ significantly reduces variance at the cost of slightly higher bias
- Eventually, added bias overtakes the benefit of reduced variance

Ridge regression: 6) Example with a toy dataset (R script)

- Problem setup:
 - n = 5, p = 2
 - Ridge vs. least squares (at $\lambda = 1$)
- Goal: See how ridge regression shrinks coefficients $\hat{\beta}$ compared to standard least squares

```
# Toy data: 5 obs, 2 predictors, no intercept

df <- data.frame(
    x1 = c(1,2,3,4,5),
    x2 = c(2,1,3,1,2),
    y = c(2,2.5,6,4,6.5)
)
```

```
# OLS fit (no intercept => '-1')
ols_fit \leftarrow lm(y \sim x1 + x2 - 1, data = df)
cat("OLS Coeffs:\n", coef(ols fit), "\n")
install.packages("glmnet") # if not installed
library(glmnet)
# Prepare X.u for almnet
X <- as.matrix(df[,c("x1","x2")])</pre>
v \leftarrow df$v
# Ridge fit with lambda=1, no intercept
     penalization
ridge_fit <- glmnet(
 X, y, alpha=0, lambda=1,
 intercept=FALSE, standardize=TRUE
cat("Ridge Coeffs (lambda=1):\n", as.matrix(
     coef(ridge fit)). "\n")
```

Ridge regression: 7) Selecting λ

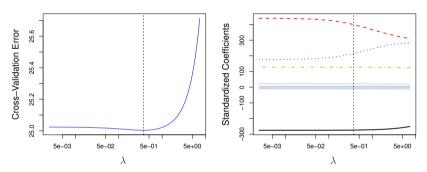


Figure: Left: CV errors for ridge regression on the Credit dataset with various values of λ . Right: Ridge regression coefficient estimates. The vertical dashed lines indicate the λ selected by CV [JWHT21, Figure 6.12].

- ullet Here, the chosen λ is relatively small, meaning minimal shrinkage relative to least square
- ullet The error curve's dip is not very pronounced, suggesting a broad range of λ values yield similar performance

Pop-up quiz #1: Ridge regression

Which statement below is false about ridge regression?

(Hint: Think carefully. Which one might not always hold?)

- A) It often lowers variance compared to least squares, at the cost of higher bias.
- B) All ridge coefficients are strictly smaller in magnitude than the least squares estimates for each predictor.
- C) The tuning parameter λ strongly influences ridge performance, often chosen via cross-validation.
- D) Ridge tends to shrink correlated predictors together.

Answer: (B) is false.

Brief explanation:

- While ridge regression shrinks the overall magnitude of the coefficient vector, it does not guarantee every individual coefficient is always smaller than its least squares counterpart.
- In some cases, ridge can redistribute weights among correlated predictors, causing some estimates to exceed their OLS magnitudes even though the sum of squares $\|\beta\|_2^2$ is reduced.

Ridge regression: 8) Summary

Ridge regression formulation:

- Add a penalty term $\sum_{j=1}^{p} \beta_j^2$ to the least squares objective, scaled by $\lambda \geq 0$
- Shrinks β_i more strongly as λ grows, stabilizing estimates

Bias-variance tradeoff:

- Larger $\lambda \Rightarrow$ higher bias but lower variance
- Especially helpful when OLS has high variance (e.g. large p, or p > n)

• Computation:

 \bullet Efficient to solve for all λ at roughly the cost of one OLS fit

Limitation:

ullet Coefficients seldom reach exactly zero \Longrightarrow no direct variable selection

The lasso: 1) Formulation

The lasso: Find $\hat{\beta}_0, \dots, \hat{\beta}_p$ that minimize

$$\sum_{i=1}^{n} \left(y_i - \beta_0 - \sum_{j=1}^{p} \beta_j x_{ij} \right)^2 + \lambda \underbrace{\sum_{j=1}^{p} |\beta_j|}_{\text{penalty}}$$

- $\lambda \ge 0$ is a tuning parameter
- Each choice of λ gives a different set of Lasso estimates \hat{eta}_{λ}^L

Remarks:

- No penalty on β_0 (the intercept)
- As $\lambda \to 0$, lasso \to standard least squares
- ullet As λ grows, many eta_j shrink toward zero, and some exactly become 0

Unlike ridge, the lasso can yield exact zero estimates \implies variable selection!

The lasso: 2) Credit dataset example

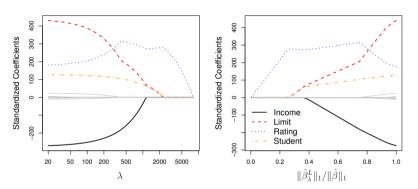


Figure: Standardized lasso coefficients for Credit, plotted vs. λ and $\|\hat{\beta}_{\lambda}^{R}\|_{2}/\|\hat{\beta}\|_{2}$ [JWHT21, Figure 6.6].

- ullet Lasso can force some coefficients to zero as λ increases
- ullet Achieves variable selection directly (predictors with $\hat{eta}_i^L=0$ are excluded)

The lasso: 3) Example with a toy dataset (R script)

- Setup:
 - n = 5, p = 2
 - Compare lasso vs. OLS at $\lambda = 1$
- Objective: See how lasso can shrink coefficients to zero

```
# Toy data: 5 obs, 2 predictors, no intercept

df <- data.frame(

x1 = c(1,2,3,4,5),

x2 = c(2,1,3,1,2),

y = c(2,2.5,6,4,6.5)

)
```

```
# OLS fit (no intercept => '-1')
ols_fit <- lm(y \sim x1 + x2 - 1, data=df)
cat("OLS Coeffs:\n", coef(ols fit), "\n")
# If needed: install.packages("qlmnet")
library(glmnet)
# Prepare X, y
X \leftarrow as.matrix(df[, c("x1", "x2")])
v \leftarrow df$v
# Lasso with alpha=1, lambda=0.1
lasso_fit <- glmnet(X, y,</pre>
                   alpha=1,
                   lambda=1,
                   intercept=FALSE,
                   standardize=TRUE)
cat("Lasso Coeffs (lambda=0.1):\n".
   as.matrix(coef(lasso fit)),"\n")
```

The lasso: 4) Selecting λ

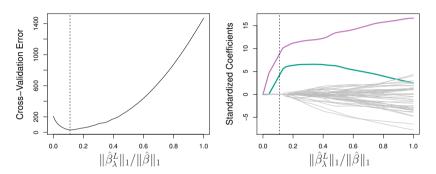


Figure: **Left:** Ten-fold cross-validation MSE for the lasso, applied to the sparse simulated data. **Right:** The corresponding lasso coefficient estimates, with the two *signal* variables in color and the *noise* variables in gray. The vertical dashed line indicates the fit that minimizes the cross-validation error [JWHT21, Figure 6.13].

- The lasso cleanly separates two signal variables from noise variables
- In contrast, standard least squares (far right, with $\|\hat{\beta}_{\lambda}^{L}\|_{1}/\|\hat{\beta}\|_{1}=1$) only identifies the purple variable without discarding the noise predictors

The lasso: 5) Summary

Lasso formulation:

- Penalty term $\sum_{i=1}^{p} |\beta_i|$ scaled by λ
- As λ grows, some β_i become exactly zero \implies variable selection

Lasso advantages:

- Encourages a sparse model for easier interpretability
- Slightly more complex than ridge, but fairly efficient to solve

Ridge vs. lasso:

- Lasso can set some coefficients to exact zeros, while ridge never does
- Ridge tends to be more stable especially when predictors are highly correlated
- Both typically tune λ via cross-validation

Wrap-up & Takeaways

Subset selection:

- Great for small p, but can be expensive or unstable for large p
- Final model is refit by least squares on the chosen subset

Regularization:

- **Ridge**: ℓ_2 penalty shrinks all coefficients toward zero, good if many have modest nonzero effects
- The lasso: ℓ_1 penalty can set some coefficients exactly to zero (variable selection)
- ullet Tuning λ typically via cross-validation for both

Next lecture:

- Geometric intuition for ridge vs. lasso
- Transition to multiple hypothesis testing

References



Gareth James, Daniela Witten, Trevor Hastie, and Robert Tibshirani.

An Introduction to Statistical Learning: with Applications in R, volume 112 of Springer Texts in Statistics.

Springer, New York, NY, 2nd edition, 2021.