Modern High Dimensional Linear Regression

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Analysis of Variance (ANOVA)

$$\begin{array}{c|c} \mathsf{POPULATION} & \mathsf{SAMPLE} \\ Y = \beta'X + \varepsilon, \ E \varepsilon X = 0 \\ EY^2 = E(\beta'X)^2 + E \varepsilon^2 \\ \mathit{MSE}_{pop} = E \varepsilon^2 \\ R_{pop}^2 := \frac{E(\beta'X)^2}{EY^2} = \\ 1 - \frac{E \varepsilon^2}{EY^2} \in [0,1] \end{array} \tag{1} \begin{array}{c} \mathsf{SAMPLE} \\ Y_i = \hat{\beta}'X_i + \hat{\varepsilon}_i \\ \mathbb{E}_n Y_i^2 = \mathbb{E}_n (\hat{\beta}'X_i)^2 + \mathbb{E}_n \hat{\varepsilon}_i^2 \\ \mathit{MSE}_{sample} = \mathbb{E}_n \hat{\varepsilon}_i^2 \\ R_{sample}^2 := \frac{\mathbb{E}_n (\hat{\beta}'X_i)^2}{\mathbb{E}_n Y_i^2} = \\ 1 - \frac{\mathbb{E}_n \hat{\varepsilon}_i^2}{\mathbb{E}_n Y_i^2} \in [0,1] \end{array}$$

By law of large numbers when p/n is small and n is large:

$$\mathbb{E}_n Y_i^2 \approx E Y^2, \ \mathbb{E}_n (\hat{\beta}' X_i)^2 \approx E (\beta' X)^2, \ \mathbb{E}_n \hat{\varepsilon}_i^2 \approx E \varepsilon^2$$
 $R_{sample}^2 \approx R_{pop}^2 \ \text{and} \ MSE_{sample} \approx MSE_{pop}$

(3)

Overfitting: What happens when p/n is not small

When p/n is not small, the discrepancy between the in-sample and out-of-sample measures of fit can be substantial. Let's check the next example :

$$X \sim N(0, I_p)$$
 and $Y \sim N(0, 1)$, $\beta' X = 0$, $R_{pop}^2 = 0$
if $p = n$, then R_{sample}^2 is $1 \gg 0$
if $p = \frac{n}{2}$, then R_{sample}^2 is about $0.5 \gg 0$
if $p = \frac{n}{20}$, then R_{sample}^2 is about 0.05

Better measures of out-of-sample predictive ability are the "adjusted" R^2 and MSE.

$$MSE_{adjusted} = \frac{n}{n-p} \mathbb{E}_n \hat{\varepsilon}_i^2, \quad R_{adjusted}^2 := 1 - \frac{n}{n-p} \frac{\mathbb{E}_n \hat{\varepsilon}_i^2}{\mathbb{E}_n Y_i^2}$$
 (5)

Measuring Predictive Ability by Sample Splitting

To measure out-of-sample performance: **Data splitting**. The idea can be summarized in two parts:

- 1. Use a random part of data, called the **training sample**, for estimating/training the prediction rule.
- 2. Use the other part, called the **testing sample**, to evaluate the quality of the prediction rule, recording out-of-sample mean squared error and R^2 .

Generic Evaluation of Prediction Rules by Sample-Splitting

- 1. Randomly partition the data into training and testing samples. Suppose we use *n* observations for training and *m* for testing/validation.
- 2. Use the training sample to compute a prediction rule $\hat{f}(X)$, for example, $\hat{f}(X) = \beta' X$.
- 3. Let *V* denote the indexes of the observations in the test sample. Then the out-of-sample/test mean squared error is

$$MES_{test} = \frac{1}{m} \sum_{k \in V} (Y_k - \hat{f}(X_k))^2$$
 (6)

and the out-of-sample/test R^2 is

$$R_{test}^2 = 1 - \frac{MSE_{test}}{\frac{1}{m} \sum_{k \in V} Y_k^2}$$
 (7)

1. Lets imagine we have the next equation

$$Y = \underbrace{\left[\beta_1 D + \beta_2' W\right]}_{Predicted \ value} + \underbrace{\varepsilon}_{error}$$
(8)

How does the predicted value of Y change if D increases by a unit while W remains unchanged?

2. Partialling-out operation: procedure that takes a random variable V and creates a "residual" \widetilde{V} by subtracting the part of V that is linearly predicted by:

$$\widetilde{V} = V - \gamma'_{VW}W, \quad \gamma_{VW} = \arg\min_{\gamma} E(V - \gamma'W)^2$$
 (9)

3. We can show that

$$Y = V + U \implies \widetilde{Y} = \widetilde{V} + \widetilde{U}$$
 (10)

1. Partialling-out to both sides of our regression equation $Y = \underbrace{[\beta_1 D + \beta_2' W]}_{Predicted\ value} + \underbrace{\varepsilon}_{error}$, we get:

$$\widetilde{\mathbf{Y}} = \beta_1 \widetilde{\mathbf{D}} + \beta_2' \widetilde{\mathbf{W}} + \widetilde{\varepsilon} \tag{11}$$

which simplifies to:

$$\widetilde{Y} = \beta_1 \widetilde{D} + \varepsilon, E[\varepsilon \widetilde{D}] = 0$$
 (12)

- $\beta_2'W$ $\widetilde{W}=0$
- $-\widetilde{\varepsilon} = \varepsilon$
- $E[\varepsilon \widetilde{D}] = 0$; since \widetilde{D} is a linear function of X = (D, W)
- 2. What we found in (16) are the Normal Equations for the population regression of \widetilde{Y} on \widetilde{D} .

Theorem (Frisch-Waugh-Lovell, FWL)

The population linear regression coefficient β_1 can be recovered from the population linear regression of \widetilde{Y} on \widetilde{D} .

$$\beta_1 = \arg\min_{b_1} E(\widetilde{Y} - \beta_1 \widetilde{D})^2 = (E\widetilde{D}^2)^{-1} E\widetilde{D}\widetilde{Y}, \ E\widetilde{D}^2 > 0$$
 (13)

Theorem (Frisch-Waugh-Lovell, FWL)

Follow the next algorithm:

- Regress Y on W , obtain residuals $arepsilon_1$
- Regress D on W , obtain residuals ε_2
- Regress ε_1 on ε_2 , obtain OLS estimates β_1

Theorem (Frisch-Waugh-Lovell, FWL)

In otherwords, β_1 can be interpreted as a (univariate) linear regression coefficient in the linear regression of **residualized** Y **on residualized** D, where the residuals are defined by **partialling-out the linear effect of** W **from** Y **and** D.

Theorem (Frisch-Waugh-Lovell, FWL-IN SAMPLE)

When we work with the sample, we mimic the partiallingout in the population.

$$\hat{\beta}_1 = \arg\min_{b_1} \mathbb{E}(\check{Y}_i - \beta_1 \check{D}_i)^2 = (\mathbb{E} \check{D}_i^2)^{-1} \mathbb{E} \check{D}_i \check{Y}_i$$
(14)

where V_i denote the residual left after predicting V_i with controls W_i in the sample:

$$\ddot{V}_i = V_i - \hat{\gamma}'_{VW} W_i, \quad \hat{\gamma}_{VW} = \arg\min_{\gamma} \mathbb{E}_n (V_i - \gamma' W_i)^2$$
(15)

- 1. Frisch-Waugh-Lovell-Python
- 2. Frisch-Waugh-Lovell-Julia

Regression in a High-Dimensional Setting / LASSO

Lasso constructs $\hat{\beta}$ as the solution of the following penalized least squares problem:

$$\min_{b \in \mathbb{R}^p} \sum_{i} (Y_i - b'X_i)^2 + \lambda \cdot \sum_{j=1}^p |b_j| \tag{16}$$

- 1. The first term is *n* times the sample mean square error
- 2. The second term is a penalty term, which penalizes the size of coefficients b_j by their absolute values times the penalty level λ . A crucial point is the choice of the penalization parameter λ .
- 3. A theoretically valid choice is (Belloni Chernozhukov, 2013)

$$\lambda = 2.c\widehat{\sigma}\sqrt{2nlog(2p/\gamma)}, \ \widehat{\sigma} \approx \sigma = \sqrt{E \in ^2}$$
 (17)

4. Another good way to pick penalty level is by cross-validation (Chetverikov et al, 2020)

Contours of the error and constraint functions for the lasso

Figure 1: Lasso optimization with two coefficients.



Intuition: The j-th component $\hat{\beta}_j$ of the lasso estimator $\hat{\beta}$ is set to zero if the marginal predictive benefit of changing $\hat{\beta}_j$ away from zero is smaller than the marginal increase in penalty:

$$\widehat{\beta}_j = 0 \quad \text{if} \quad \left| \partial_{b_j} \sum_i (Y_i - \widehat{\beta}' X_i)^2 \right| < \lambda$$
 (18)

OLS post Lasso

We can then **use the Lasso-selected set of regressors** to refit the model by least squares. This method is called the "least squares post Lasso" or simply **post-Lasso**

$$\widetilde{\beta} \in \arg\min_{\beta \in \mathbb{R}^p} \sum_i (Y_i - X_i' \beta)^2 : \beta_j = 0 \text{ if } \widehat{\beta}_j = 0 \text{ for each } j$$
 (19)

Under approximate sparsity Lasso and Post-Lasso will approximate the best linear predictor well. This means that they won't overfit the data, and we can use the sample and adjusted R^2 and MSE to assess out-of-sample predictive performance. Of course, it is always a good idea to verify the out-of-sample predictive performance by using sample splitting

How to select λ

Big lambdas tend to result in a lot of shrinkage and sparsity, as $\lambda->0$ our solution approaches the OLS solution

Two ways to select λ

- Select model with lowest AIC/BIC/other plug-in criterion. This uses no out-of-sample information for selection but is fast.
- Cross-validate by testing on our hold-out test sample. Variants of cross-validation are most commonly used.

k-fold cross-validation

The next algorithm is taken from Ivan Rudik' Lectures Note in Dynamic Optimization(Cornell, Fall 2021)

Figure 2: k-fold cross-validation

k-fold cross-validation

In k-fold cross-validation we do the following:

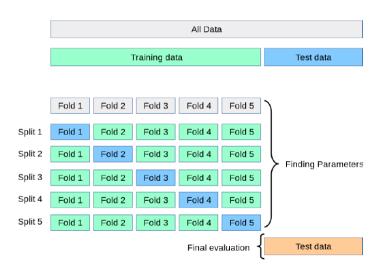
- Create a grid of λ s
- For each λ :
 - \circ Split data into k mutually-exclusive folds of about equal size, usually choose k=5.10
 - \circ For $i = 1, \ldots, k$
 - fit the model using all folds but fold i
 - Predict out-of-sample on fold j
 - \circ Compute average mean squared prediction error across the k folds:

$$ar{Q}(\lambda) = rac{1}{k} \sum_{i=1}^k \sum_{i \in ext{fold } i} ig(y_i - (lpha_0 + x_i'eta) ig)^2 + \lambda ||eta||_1$$

• Choose $\hat{\lambda}_{min} = argmin_{\lambda}\bar{Q}(\lambda)$ or to avoid modest overfitting choose the largest λ such that $\bar{Q}(\lambda) \leq \hat{\lambda}_{min} + \sigma_{\hat{\lambda}}$. (1 standard deviation rule)

k-fold cross-validation

Figure 3: k-fold cross-validation



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Practical Value of CEF

- 1. Randomly split your sample with N observations into a training set with N_{tr} and a test set with N_{te} observations.
- 2. Estimate different CEF functions in the training sample $\hat{m}(X)$.
- 3. Calculate and compare their out-of-sample Mean Squared Error (MSE) in the test sample:

$$MSE_{te} = \frac{1}{N_{te}} \sum_{i=1}^{N_{te}} (Y_i - \hat{m}(X_i))^2$$

4. Pick the function with the lowest MSE.