Saarland University

The Elements of Stastical Learning

Assignement 4

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Prove that for linear and polynomial least squares regression, the LOOCV estimate for the test MSE can be calculated using the following formula:

$$CV_{(n)} = \frac{1}{n} \sum_{i=1}^{n} \left(\frac{y_i - \hat{y}_i}{1 - h_i} \right)^2$$
 (1.1)

Where h_i is the leverage (3.37, ISLR p98)

$$h_i = \frac{1}{n} + \frac{(x_i - \bar{x})^2}{\sum_{i'=1}^n (x_i' - \bar{x})^2}$$
 (1.2)

We first have this equation, that can take long if n is big because it has to fit every model.

$$MSE_i = (y_i - \hat{y_i})^2$$

$$CV_{(n)} = \frac{1}{n} \sum_{i=1}^{n} MSE_i$$

So knowing that $\hat{y} = Hy$ and as it is a Leave One Out cross validation, we fit n times the model with one element out. So we have:

$$H = X(X^{T}X)^{-1}X^{T}$$

$$H^{-i} = X_{-i}(X_{-i}^{T}X_{-i})^{-1}X_{-i}^{T}$$

The hat matrix with all the data and with one out, respectively Then we have

$$\hat{y}_i = x_i^T [X(X^T X)^{-1} X^T] y$$

$$\hat{y}_{-i} = x_i^T [X_{-i} (X_{-i}^T X_{-i})^{-1} X_{-i}^T] y_{-i}$$

The fitted values at x_i when using all the data points and when leaving one out. We can then do the following:

$$\hat{y}_{-i} = \sum_{i \neq j} H_{ij} y_j + H_{ii} \hat{y}_{-i}$$

$$\hat{y}_{-i} = \sum_{j}^{m} H_{ij} y_j - H_{ii} y_i + H_{ii} \hat{y}_{-i}$$

We have $\sum_{j=1}^{m} H_{ij}y_j = \hat{y}_i$ so:

$$\hat{y}_{-i} = \hat{y}_i - H_{ii}y_i + H_{ii}\hat{y}_{-i}$$

We substitute \hat{y}_{-i} in the prediction error:

$$y_{i} - \hat{y}_{-i} = y_{i} - (\hat{y}_{i} - H_{ii}y_{i} + H_{ii}\hat{y}_{-i})$$
$$y_{i} - H_{ii}y_{i} - \hat{y}_{-i} - H_{ii}\hat{y}_{-i} = y_{i} - \hat{y}_{i}$$
$$y_{i} - \hat{y}_{-i} = \frac{y_{i} - \hat{y}_{i}}{1 - H_{ii}}$$

Taking the Mean Square Error leads to equation 1.1.

1. Ridge regression is done by minimizing the RSS with a quadratic penalty term:

$$minimize(y - X\beta)^T(y - X\beta) + \lambda \beta^T \beta$$

to show that the solutions take on the form:

$$\hat{\beta}^{ridge} = (X^T X + \lambda y I)^{-1} X^T y$$

First we expand the equation to:

$$y^T y - y^T X \beta - y X^T \beta^T + X^T \beta^T X \beta + \lambda \beta^T \beta$$

which simplifies to:

$$y^T y - y X^T \beta^T - y X^T \beta^T + X^T \beta^T X \beta + \lambda \beta^T \beta$$

and:

$$y^Ty - 2yX^T\beta^T + X^T\beta^T X\beta + \lambda\beta^T\beta$$

We take the first derivative with respect to β , which gives us:

$$0 - 2X^T y + 2(XX^T)\beta + 2\lambda\beta$$

setting this to zero this can be simplified to:

$$2(XX^T)\beta + 2\lambda\beta = 2X^Ty$$

and further simplified to:

$$((XX^T) + \lambda I)\beta = X^T y$$

where I is the p x p identity matrix (added to the matrix math works out correctly). Solving for β gives:

$$\hat{\beta}^{ridge} = (X^T X + \lambda y I)^{-1} X^T y$$

Assume a scenario in which the number of observations equals the number of features (n = p) and X is the $n \times n$ identity matrix. Furthermore, assume that we perform regression without an intercept. In this setting, lasso simplifies to:

$$\underset{\beta}{\text{minimize}} \sum_{j=1}^{p} (y_j - \beta_j)^2 + \lambda \sum_{j=1}^{p} |\beta_j|$$
(3.1)

Show that the lasso estimates take the form:

$$\hat{\beta}_{j}^{lasso} = \begin{cases} y_{j} - \frac{\lambda}{2}, & \text{if } y_{j} > \frac{\lambda}{2}; \\ y_{j} + \frac{\lambda}{2}, & \text{if } y_{j} < -\frac{\lambda}{2}; \\ 0, & \text{if } |y_{j}| \leq \frac{\lambda}{2}; \end{cases}$$
(3.2)

For equation 3.1 to be a minimum, we must derive it and equal it to zero:

$$\frac{d}{d\beta_j} \sum_{j=1}^p (y_j - \beta_j)^2 + \frac{d}{d\beta_j} \lambda \sum_{j=1}^p |\beta_j| = 0$$
(3.3)

We can rewrite equation 3.1 with the matrix notation like following:

$$F(\beta) = (y - \beta)^T (y - \beta) + \lambda \beta = yy^T - y\beta^T - y^T\beta + \beta\beta^T + \lambda\beta$$
$$\frac{dF}{d\beta} = -2y + 2\beta + \lambda$$
$$\beta = y - \frac{\lambda}{2}$$

To make β negative y has to be smaller than $\frac{\lambda}{2}$ To make β positive y has to be greater than $\frac{\lambda}{2}$ To make β equals 0, |y| has to be equal to $\frac{\lambda}{2}$

- 1. See the attached R code for the data normalization and splits.
- 2. We applied the best subset selection on the training set. Figure 1 shows the curves for R^2 , adjusted R^2 , C_p and BIC as a function of the number of predictors. As expected R^2 increases as the number of predictors increases. During training we try to estimate the model coefficients such that the RSS is as small as possible, and adding more predictors means the training error will get smaller. However, this does not mean the test error will get smaller. Since our real goal is to reduce test error, RSS and R^2 are not good measures of a "best" model. Adjusted R^2 and C_p both suggest that models with 7 predictors would be the best model, since these models have the highest adjusted \mathbb{R}^2 and lowest \mathbb{C}_p . These two also show that a model with 6 predictors will achieve about the same performance. BIC replaces the penalty term used by C_P , $2d\hat{\sigma}^2$, with $log(n)d\hat{\sigma}^2$, where n is the number of observations. Since log(n) > 2 for any n > 7, BIC places a heavier penalty of models with more predictors. This can be observed in Figure 1, where the BIC statistic suggests a model with 2-4 predictors is best. Which model should we choose, a model with 6-7 predictors as indicated by the C_P statistic (and corroborated by the R^2 statistic, even though it's not a well motivated in statistical theory as the other statistics), or a model with 2-4 predictors as indicated by the BIC statistic? We'll choose the simplest "best" model, which is two predictor model indicated by the BIC statistic. Figure 2 shows the the selected features used for the best model for the range of predictors. In our selected model the features used are lcavol and lweight. The training error for this model is 0.553. The **test error** is [].

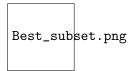


Figure 1: RSq, adjusted RSq, Cp and BIC values versus the number of predictors in best subset selection.



Figure 2: Selected features for best model with predictors from BIC statistic.

3. Figure 3 shows the results of the ridge regression fit on the training data, specifically the values of the coefficients in relation to λ . The results are as expected. On the left hand side λ is almost zero, therefore the ridge coefficient estimates are essentially the same as the least squares estimates. As lambda increases the coefficient estimates shrink to zero. On the right hand side of the plot, when lambda is large, all of the estimates are zero.



Figure 3: Ridge regression coefficients in relation to lambda.

- 4. See code for 5-fold cross-validation for the ridge regression model. For this model the **training error** in MSE is 0.09645, and the **test error** is 0.49327.
- 5. Figure 4 shows the plot of the lasso coefficient estimates plotted against λ . As expected, on the far left, when λ is zero, the lasso gives the least squares fit, and when lambda is large the coefficient estimates shrink to zero. For the ridge regression the coefficients all smoothly shrink to zero, however for the lasso coefficients most of them sharply shrink to zero as lambda increases.



Figure 4: Lasso coefficients in relation to lambda.

- 6. See the code for 5-fold cross-validation for the lasso model. For this model the **training error** is 0.09423, and the **test error** is 0.45305. [How many features?? Compare coef.]
- 7. For the linear regression model on all the features the **training error** is [] and the **test error** is [].