Supervised Learning

Regularization

June 26th, 2023

Previously...

We talked about variable selection in the linear model context:

$$\hat{Y}={\hateta}_0+{\hateta}_1x_1+\dots+{\hateta}_px_p$$

Why would we attempt to select a **subset** of the p variables?

- To improve model interpretability
 - Occam's razor: simplest model wins!
 - Eliminating uninformative predictors is obviously a good thing when your goal is to tell the story of how your predictors are associated with your response.
- To improve prediction accuracy
 - Eliminating uninformative predictors can lead to lower variance in the test-set MSE, at the expense of a slight increase in bias

Best subset selection

- Start with the **null model** \mathcal{M}_0 (intercept-only) that has no predictors
 - just predicts the sample mean for each observation
- For $k=1,2,\ldots,p$ (each possible number of predictors)

• Fit **all** $\binom{p}{k} = rac{p!}{k!(p-k)!}$ with exactly k predictors

- $\circ~$ Pick the best (some criteria) among these $\binom{p}{k}$ models, call it \mathcal{M}_k
- $\circ\,$ Best can be up to the user: cross-validation error, highest adjusted R^2 , etc.
- Select a single best model from among $\mathcal{M}_0,\ldots,\mathcal{M}_p$

This is not typically used in research!

• only practical for a smaller number of variables

Remember the bias-variance tradeoff

 $\mathrm{MSE} = \mathrm{(Bias)}^2 + \mathrm{Variance}$

- Introduce bias but decrease variance to improve predictions
- Some questions
 - How do we know that there is a *trade-off* between Bias and Variance?
 - How can we check that it is $(Bias)^2$ and not (Bias) without doing *any* calcs?

Shrinkage methods: Ridge regression

Ridge regression introduces a **shrinkage penalty** $\lambda \geq 0$ by minimizing:

$$\sum_{i}^{n} ig(Y_i - eta_0 - \sum_{j}^{p} eta_j x_{ij}ig)^2 + \lambda \sum_{j}^{p} eta_j^2 = ext{RSS} + \lambda \sum_{j}^{p} eta_j^2$$

or more succinctly we want to minimize:

$$\left\| \mathbf{Y} - \mathbf{X} eta
ight\|_2^2 + \lambda \left\| eta
ight\|_2^2$$

- as λ increases \Rightarrow flexibility of models decreases
 - increases bias, but decreases variance
- for fixed value of λ , ridge regression fits only a single model
 - $\circ\,$ need to use cross-validation to tune λ

Shrinkage methods: Ridge regression

For example: note how the magnitude of the coefficient for Income trends as $\lambda o \infty$



The coefficient **shrinks towards zero**, but never actually reaches it

- Income is always a variable in the learned model, regardless of the value of λ

Shrinkage methods: Lasso regression

Ridge regression **keeps all variables**

But we may believe there is a **sparse** solution

Lasso enables variable selection with λ by minimizing:

$$\sum_{i}^{n}ig(Y_i - eta_0 - \sum_{j}^{p}eta_j X_{ij}ig)^2 + \lambda\sum_{j}^{p}|eta_j| = ext{RSS} + \lambda\sum_{j}^{p}|eta_j|$$

or more succinctly we want to minimize:

$$\left\| \mathbf{Y} - \mathbf{X} eta
ight\|_{2}^{2} + \lambda \left\| eta
ight\|_{1}$$

- Lasso uses an ℓ_1 ("ell 1") penalty
- as λ increases \Rightarrow flexibility of models decreases

• increases bias, but decreases variance

- Can handle the p>n case, i.e. more variables than observations!

Shrinkage methods: Lasso regression

Lasso regression **performs variable selection** yielding **sparse** models



The coefficient shrinks towards and **eventually equals zero** at $\lambda pprox 1000$

- if the optimum value of λ is larger, then Income would NOT be included in the learned model

Which do we use?



Best of both worlds? Elastic net

$$\sum_{i}^{n}\left(Y_{i}-eta_{0}-\sum_{j}^{p}eta_{j}X_{ij}
ight)^{2}+\lambda\left[(1-lpha)\|eta\|_{2}^{2}/2+lpha\|eta\|_{1}
ight]$$

- $||eta||_1$ is the ℓ_1 norm: $||eta||_1 = \sum_j^p |eta_j|$
- $||eta||_2$ is the ℓ_2 , Euclidean, norm: $||eta||_2 = \sqrt{\sum_j^p eta_j^2}$
- Ridge penalty: $\lambda \cdot (1-lpha)/2$
- Lasso penalty: $\lambda \cdot lpha$
- lpha controls the **mixing** between the two types, ranges from 0 to 1
 - $\circ \ lpha = 1$ returns lasso
 - $\circ \ lpha = 0$ return ridge

Caveats to consider...

- For either ridge, lasso, or elastic net: you should standardize your data
- Common convention: within each column, compute then subtract off the sample mean, and compute the divide off the sample standard deviation:

$$ilde{x}_{ij} = rac{x_{ij} - ar{x}_j}{s_{x,j}}$$

- glmnet package does this by default and reports coefficients on the original scale
- λ and α are tuning parameters
- Have to select appropriate values based on test data / cross-validation
- When using glmnet, the cv.glmnet() function will perform the cross-validation for you

Example data: NFL teams summary

Created dataset using nflfastR summarizing NFL team performances from 1999 to 2020

```
library(tidyverse)
nfl_teams_data <- read_csv("https://shorturl.at/uwAV2")
nfl_model_data <- nfl_teams_data %>%
    mutate(score_diff = points_scored - points_allowed) %>%
    # Only use rows with air yards
    filter(season >= 2006) %>%
    dplyr::select(-wins, -losses, -ties, -points_scored, -points_allowed, -season, -team)
```

Introduction to glmnet

We will use the glmnet package for ridge, lasso, and elastic net

library(glmnet)

• could use the model.matrix() function (which converts factors to 0-1 dummy variables!)

```
model_x <- nfl_model_data %>%
    dplyr::select(-score_diff) %>%
    as.matrix()
model_y <- nfl_model_data$score_diff
# model_x <- model.matrix(score_diff ~ ., nfl_model_data)[, -1]</pre>
```

Initial model with lm()

- What do the initial regression coefficients look like?
- Use broom to tidy model output for plotting

```
init_reg_fit <- lm(score_diff ~ ., nfl_model_
library(broom)
tidy(init_reg_fit) %>%
  mutate(coef_sign = as.factor(sign(estimate))
        term = fct_reorder(term, estimate))
  ggplot(aes(x = term, y = estimate, fill = c
  geom_bar(stat = "identity", color = "white"
  scale_fill_manual(values = c("darkred", "da
            guide = FALSE) +
  coord_flip() + theme_bw()
```

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defense success rate run -								
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term

Ridge regression example

Perform ridge regression using glmnet with alpha = 0 (more on that later)

By default it standardizes your predictors and fits model across a range of λ values (can plot these!)

init_ridge_fit <- glmnet(model_x, model_y, alpha = 0)
plot(init_ridge_fit, xvar = "lambda")</pre>



Log Lambda

Ridge regression example

We use cross-validation to select λ with cv.glmnet() which uses 10-folds by default

• specify ridge regression with alpha = 0

fit_ridge_cv <- cv.glmnet(model_x, model_y, alpha = 0)
plot(fit_ridge_cv)</pre>



Tidy ridge regression



Tidy ridge regression



Ridge regression coefficients

Coefficients using the 1 standard error rule λ



Lasso regression example

Similar syntax to ridge but specify alpha = 1:



Lasso regression example

Number of non-zero predictors by λ

```
tidy_lasso_cv <- tidy(fit_lasso_cv)
tidy_lasso_cv %>%
ggplot(aes(x = lambda, y = nzero)) +
geom_line() +
geom_vline(xintercept = fit_lasso_cv$lambda
geom_vline(xintercept = fit_lasso_cv$lambda
linetype = "dashed", color = "re
scale_x_log10() + theme_bw()
```

Reduction in variables using **1 standard error rule** λ



Lasso regression example

Coefficients using the 1 standard error rule λ

```
tidy_lasso_coef %>%
  filter(lambda == fit_lasso_cv$lambda.1se) %
  mutate(coef_sign = as.factor(sign(estimate)
            term = fct_reorder(term, estimate))
  ggplot(aes(x = term, y = estimate,
            fill = coef_sign)) +
  geom_bar(stat = "identity", color = "white"
  scale_fill_manual(values = c("darkred", "da
            guide = FALSE) +
  coord_flip() +
  theme_bw()
```



Elastic net example

Need to tune both λ and lpha - can do so manually with our own folds

```
set.seed(2020)
fold_id <- sample(rep(1:10, length.out = nrow(model_x)))</pre>
```

Then use cross-validation with these folds for different candidate alpha values:

```
cv_en_25 <- cv.glmnet(model_x, model_y, foldid = fold_id, alpha = .25)
cv_en_50 <- cv.glmnet(model_x, model_y, foldid = fold_id, alpha = .5)
cv_ridge <- cv.glmnet(model_x, model_y, foldid = fold_id, alpha = 0)
cv_lasso <- cv.glmnet(model_x, model_y, foldid = fold_id, alpha = 1)</pre>
```

Can see which one had the lowest CV error among its candidate λ values:

which.min(c(min(cv_en_25\$cvm), min(cv_en_50\$cvm), min(cv_ridge\$cvm), min(cv_lasso\$cvm)))

[1] 2

Elastic net example

Can view same type of summary

• More relaxed than lasso for variable entry



Comparison of models based on holdout performance

```
set.seed(2020)
nfl_model_data <- nfl_model_data %>% mutate(test_fold = sample(rep(1:5, length.out = n())))
holdout predictions <-
 map_dfr(unique(nfl_model_data$test_fold),
          function(holdout) {
            # Separate test and training data:
            test data <- nfl model data %>% filter(test fold == holdout)
            train data <- nfl model data %>% filter(test fold != holdout)
            # Repeat for matrices
            test x <- as.matrix(dplyr::select(test data, -score diff))</pre>
            train_x <- as.matrix(dplyr::select(train_data, -score_diff))</pre>
            # Train models:
            lm model <- lm(score diff ~ ., data = train data)</pre>
            ridge_model <- cv.glmnet(train_x, train_data$score_diff, alpha = 0)</pre>
            lasso_model <- cv.glmnet(train_x, train_data$score_diff, alpha = 1)</pre>
            en_model <- cv.glmnet(train_x, train_data$score_diff, alpha = .5)</pre>
            # Return tibble of holdout results:
            tibble(lm_preds = predict(lm_model, newdata = test_data),
                   ridge_preds = as.numeric(predict(ridge_model, newx = test_x)),
                   lasso_preds = as.numeric(predict(lasso_model, newx = test_x)),
                   en_preds = as.numeric(predict(en_model, newx = test_x)),
                   test_actual = test_data$score_diff, test_fold = holdout)
          2)
```

```
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```

Predictions compared to lm?

Compute RMSE across folds with std error intervals



In this case lm actually "beat" regularization, but within intervals