Lecture 08 Decision Trees II

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Yale

Updated office hours:

- ► Taylor Arnold Wednesdays, 16:00 17:00, HH 24
- ▶ Elena Khusainova Tuesdays, 13:00-15:00, HH 24, Basement
- ▶ Yu Lu Wednesdays, 19:00-20:30, HH 24, Basement
- ▶ Jason Klusowski Thursdays, 19:00-20:30, HH 24

► We will give you full credit either way you approach this, but your solution for Question #4 will be much more interesting if you weight the MSE of the census tracts by their population

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- ► As well, many of the R and Python scripts did not actually run correctly; a common problem being that they threw errors when they could not find global variables or local files

Problem Set 1, cont.

- ► Paying attention to these details is very important!
- ► In almost all cases, you got full credit for the implementation and data analysis questions as long as you wrote *something* and your code ran without any errors
- Future problem sets will likely have fewer 'parts' to them, but the actually content will be graded more strictly, so you do not want to lose points on silly things
- ► Test your code!
- ► Remember:
 - No late submissions (unless you have a true emergency, and for undergraduates a Deans excuse)
 - ▶ We do drop the lowest grade from the final assignment

► Today:

- Finish decision trees
- ▶ Walk through an analysis of trees for classification and regression
- Introduction to SVMs (time permitting)

Gradient boosted trees

Last time, we looked at decision trees and random forests. There is one additional variant on these ideas which has been shown to produce extremely powerful predictions. Called gradient boosted trees.

The algorithm for this can conceptually be described as follows:

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- 5. Calculate decision tree $T^{(3)}$ on a randomly sampled set of the residuals $r^{(2)}$, and calculate the predicted values $\hat{r}^{(2)}$ for the entire dataset.

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- 3. Fit a new decision tree $T^{(2)}$ on a random resampling of the residuals $r^{(1)}$, and calculate the predicted values $\hat{r}^{(1)}$ for the entire dataset.
- 4. Now, set $\hat{y}^{(2)}$ to be $\hat{y}^{(1)} + \hat{r}^{(1)}$ and set the new residuals to be $r^{(2)} = y \hat{y}^{(2)}$.
- 5. Calculate decision tree $T^{(3)}$ on a randomly sampled set of the residuals $r^{(2)}$, and calculate the predicted values $\hat{r}^{(2)}$ for the entire dataset.
- 6. Repeat as needed until you have predictions $\hat{y}^{(K)}$.

The actual algorithm is modified it two ways:

- 1. The predicted values in the terminal nodes of the decision tree $T^{(j)}$ for any j are not determined just by the sampled set of data, but rather chosen to minimize the actual loss of using the $\hat{y}^{(j)}$.
- 2. The values for the predicted residuals $\hat{r}^{(j)}$ is 'shrunk' towards zero by a factor of $0 < \rho \leq 1$. This value is called the learning rate.

Due to the learning rate factor, it makes sense to center the input variables before applying this algorithm (this is usually done by the software).