Building Footprint Identification with Airborne LiDAR: A Final Project for FOR 796

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Abstract

This is a test.

1. Introduction

2. Methods

2.1. Building and LiDAR Data

Rasterized building footprint data served as the response data in this study Microsoft (2018). The raw raster contains counts of the number of buildings intersecting each pixel. This data was converted to a Boolean raster where ones represent any buildings present, and zeroes represent no buildings present. This data was chosen for it's availability across the entire country, it's high accuracy ($>$ 99% positive predictive value; Pourpeikari Heris et al. (2020)), and it's 30m resolution.

The raw LiDAR data originates from a single LiDAR acquisition covering portions of Erie, Genesee, and Livingston counties in western New York made available by the New York Stat GIS Program Office (New York Office of Information Technology Services 2019). The raw LiDAR data was then height-normalized and converted into a set of 39 predictors chosen for their prevalence in models of forest structure (Hawbaker et al. 2010; Huang et al. 2019; Pflugmacher et al. 2014). LiDAR data was selected due to it's known ability to characterize three dimensional height-profiles at high-resolution. More specifically, these

forest-structure predictors were chosen for practicality. It would be a great benefit to those mapping forest structure with LiDAR if the same predictors could also be leveraged for building better forest masks.

The LiDAR predictors, in raster stack form, were overlaid with the Boolean building raster to create stack of data where each pixel contained a set of 39 predictors and one building indicator response variable. A stratified random sample was conducted on the raster stack, with the building indicator providing the levels of stratification. 3,500 observations (pixel locations) were selected from each stratum resulting in 7,000 observations for model training and testing. This final dataset was converted to a 7000x40 (rows, columns) data frame. The lidR (Roussel and Auty 2020; Jean-Romain et al. 2020) and raster (Hijmans 2021) packages were used for height-normalization and dataset generation.

Additionally, principle components were derived from the final dataset to remove multicollinearity from the 39 predictors (R Core Team 2021). This alternative dataset consisted of the first seven principle components, as they accounted for $\geq 95\%$ of the information in the predictors. This alternative dataset was a 7000x7 data frame.

2.2. Models

Three candidate classification models were fit to a random 70% (calibration data; n = 3500) of the observations, with the remaining 30% reserved for model performance assessment (holdout data; n = 1500).

The first candidate model was a simple logistic regression model (R Core Team 2021), and was trained on the principle components variant of the calibration data. The second candidate model was a random forest (RF herafter) trained with the ranger R package (Wright and Ziegler 2017) and the calibration dataset. The third candidate model was a stochastic gradient boosting machine (LGB hereafter) trained with the LightGBM R package (Ke et al. 2021) and the

Table 1: Model accuracy metrics computed against 30% holdout partition (n = 1500).

Model	AUC	Overall Accuracy	Sensitivity	Specificity
Logistic	0.87	0.79	0.83	0.75
LGB	0.88	0.80	0.81	0.78
RF	0.96	0.90	0.92	0.89

calibration dataset. The hyperparemeters for both the RF and LGB models were selected using a grid search where each combination of hyperparameters were compared against eachother using the cross-entropy loss function (CEL; Equation (1)) computed from a random five-fold cross-validation with the calibration dataset. CEL is computed as follows:

$$\text{CEL} = \sum_{i=1}^{n} -\log\left(\hat{y_i}\right) \tag{1}$$

Where n is the number of observations in the fold, and $\hat{y_i}$ is the predicted probability of the true class.

Each of the models was assessed against the holdout dataset and compared to one another using overall accuracy, specificity, sensitivity, and AUC. Additionally ROC curves were plotted for each model's results on the holdout set. The caret and pROC R packages were used to compute these accuracy metrics (Kuhn 2021; Robin et al. 2011).

3. Results

4. Discussion

- RF was superior to all the others
- Logistic still rather good. Makes you wonder if the marginal benefits of ML here are worth the effort, time, understandability

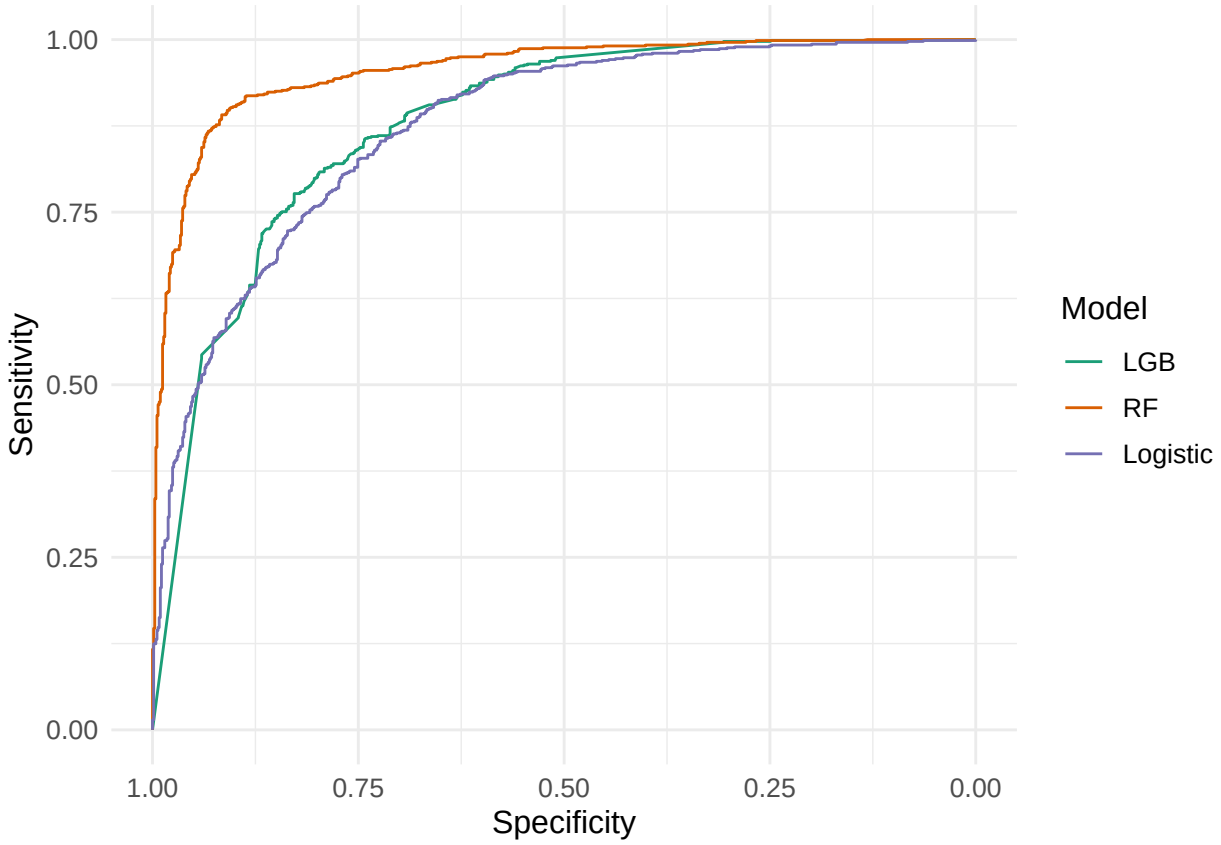

Figure 1: ROC curves showing sensitivity against specificity for each model.

- What could have been done better?
  - ensembling methods?
  - wider ranges of hyperparameters
  - larger inclusion than just 95% of the information for principle components and logistic regression.
  - Timing. Response from 2018. LiDAR from 2019
- What might these models be used for
  - Since the predictors herein are primarily used for predicting forest structure - these models make a relatively easy way to extract man-made structures away from LiDAR-based maps of forest structure like biomass.
  - There are questions about transferrability of these data from this region and this lidar coverages to others from different locations and times.

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