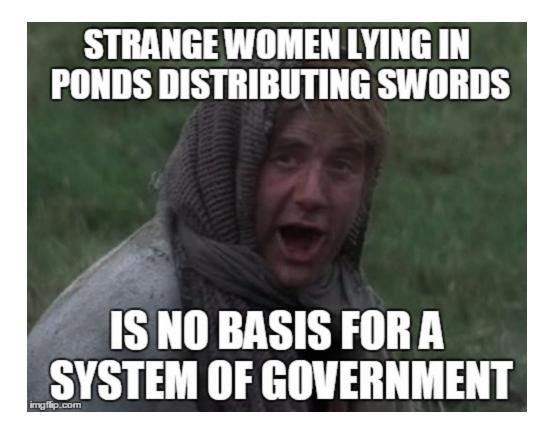
Random Forests



Machine learning algorithms are vulnerable to overfitting with respect to in-sample patterns, reducing our ability to accurately model out-of-sample

Who chose the next Roman emperor? Who chose the next king of France or England?

Who chose the next Roman emperor? Who chose the next king of France or England?

The current emperor or king (mostly)

Why is this a bad idea?

- A single person (algorithm?) making a decision can easily make an error of judgement.
- If we choose a Nero, we end up with Rome in flames

How do most developed countries now choose leaders?

They vote!

Why?

- NOT because we care if everyone has their opinion heard (see history of voting rights)
- Because large groups of people, when their opinions are averaged, make good better choices

My students averaged 55% on their final exam

On the other hand, if a student (in the same class!) had chosen the most popular response to each question based on the responses of their classmates would have scored ~85%

In aggregate, weak learners can become strong learners

This principle also applies to statistical learning algorithms.

Aggregating "weak" algorithms can lead to a "strong" algorithm.

Collections of learning algorithms are called **ensembles**.

Ensemble Flavors - First Flavor

Bagging

Bagging (Bootstrap Aggregation) is a simple way to start creating an ensemble model.

Standard Model:

$$\hat{f}(x) = f^*(x)$$

All training data is used to generate our single best estimate of the true functional form, f(x).

Bagging

$$\hat{f}_{bag}(x) = rac{1}{B} \sum_{b=1}^{B} f_b^*(x)$$

In bagging, each estimate utilizes a bootstrap (random) sample of the training data

The bagged estimate is then based on the weighted average of all of the models

Tree Problems

One drawback to bagging can be illustrated by thinking about how decision trees are generated

- 1. Find the biggest information gain
- 2. Split the tree
- 3. On each branch, find the next best information gain
- 4. Split again
- 5. Repeat 3 and 4 until stopping rule is reached (depth, purity, number of observations, etc.)

Random Forests

Using bagging on decision trees in a situation where one variable is clearly superior to other inputs, the data itself will almost never allow a model to explore other inputs.

- The most informative input will mask the other options (always be chosen)
- Each tree in the bagging algorithm is highly correlated with the other trees
 - Permits overfitting, and reduces predictive power

Random Forests

How can we alleviate this tendency?

- Restrict the inputs that the tree is allowed to choose from
- Bootstrap the sample
- Aggregate the forecasts to make a single, more accurate, prediction.

Restricting Inputs

When a classification tree looks for maximum information gain, it searches across **all** available inputs.

Trees in a random forest are restricted to a random subset of inputs at each branching:

- Typically, \sqrt{k} (where k is the number of available parameters) inputs are provided at each branch
- When a new branch occurs, a new random subset of inputs is provided
- This is repeated for all branches on all trees

Making a Classification

Once each tree in a random forest has been grown, we can use the trees to create a decision rule based on a vote by the classifiers:

- Each tree classifies an observation
- Whichever class receives the most votes (has highest predicted probability of being the true observed class)
 "wins," and is assigned as the predicted class for the observation

MNIST Dataset

MNIST Handwriting Recognition Data

- Contains digits 0-9
- Full Dataset is 60,000 training observations, 10,000 testing observations
- 28 x 28 pixel images, (784 factors)
- Numbers are centered in the image
- Goal is to predict the written digit based on the image
- Great learning dataset

Implementing Ensembles

We will implement Random Forests, Bagging, and Boosting using the scikit-learn module in Python, as we did for Decision Trees

```
import pandas as pd
import numpy as np
from sklearn.model_selection import train_test_split
from sklearn.metrics import accuracy_score
```

Prepare the Data

Baseline Decision Tree Classifier

Resulting in:

```
The decision tree has an accuracy of: 0.5795555555555555
```

Bagging Ensemble (of decision trees)

Resulting in:

```
The bagging algorithm has an accuracy of: 0.8011111111111
```

Random Forest Classifier

```
from sklearn.ensemble import RandomForestClassifier
# Generate the random forest model
forest = RandomForestClassifier(n_estimators=100,
        n jobs = -1, random state=42)
# Fit the model to the training data
fclf = forest.fit(x, y)
# Make predictions
fpred = fclf.predict(xt)
# Print the accuracy score of the fitted model
print("The random forest has an accuracy of : %s\n"
        % str(accuracy_score(fpred, yt)))
```

Resulting in:

Summary of Results

1. Decision Tree: 57.9%

2. Bagging Algorithm: 80.1%

3. Random Forest: 84.4%

Lab Time!