

Foundations of Machine Learning

AI2000 and AI5000

FoML-37
Model Combination

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So far in FoML

- Intro to ML and Probability refresher
- MLE, MAP, and fully Bayesian treatment
- Supervised learning
 - a. Linear Regression with basis functions
 - b. Bias-Variance Decomposition
 - c. Decision Theory - three broad classification strategies
 - d. Neural Networks
- Unsupervised learning
 - a. K-Means, Hierarchical, and GMM for clustering
- Kernelizing linear Models
 - a. Dual representation, Kernel trick, SVM (max-margin classifier)
- Tree-based Methods



For today

- Model combination



Single vs Multiple models

- Combining multiple models (often) → improved performance

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Single vs Multiple models

- Combining multiple models (often) → improved performance
 - E.g., train L different models and use the average of the predictions made by each model
- Such combinations of models → Committees

Model combination - variants

- Boosting
 - Training multiple models in sequence
 - Error function used to train a models depends on the performance of the previous model



Model combination - variants

- Select one of the models to make the prediction
 - Choice of the model is a function of the input
 - Different models are responsible for making predictions in different regions

Model combination - variants

- Select one of the models to make the prediction
 - Choice of the model is a function of the input
 - Different models are responsible for making predictions in different regions
- E.g., decision trees
 - Selection process is a sequence of binary selections

Bayesian Model Averaging vs. Model combination



Model combination

- E.g., density estimation using a mixture of Gaussians (GMM)
- Several Gaussian components are combined probabilistically
 - Binary latent variable z is responsible for generating x

Model combination

$$p(\mathbf{x}, \mathbf{z})$$

$$p(\mathbf{x}) = \sum_{\mathbf{z}} p(\mathbf{x}, \mathbf{z}).$$

$$p(\mathbf{x}) = \sum_{k=1}^K \pi_k \mathcal{N}(\mathbf{x} | \boldsymbol{\mu}_k, \boldsymbol{\Sigma}_k)$$

$$p(\mathbf{X}) = \prod_{n=1}^N p(\mathbf{x}_n) = \prod_{n=1}^N \left[\sum_{\mathbf{z}_n} p(\mathbf{x}_n, \mathbf{z}_n) \right].$$

Each data sample has a corresponding latent variable



Bayesian Model Averaging

- Several different models indexed by h and prior $p(h)$
 - E.g., GMM or mixture of Cauchy distributions

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Bayesian Model Averaging

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One model is responsible for generating the whole data, $p(h)$ captures our uncertainty as to which model that is



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Decision Trees

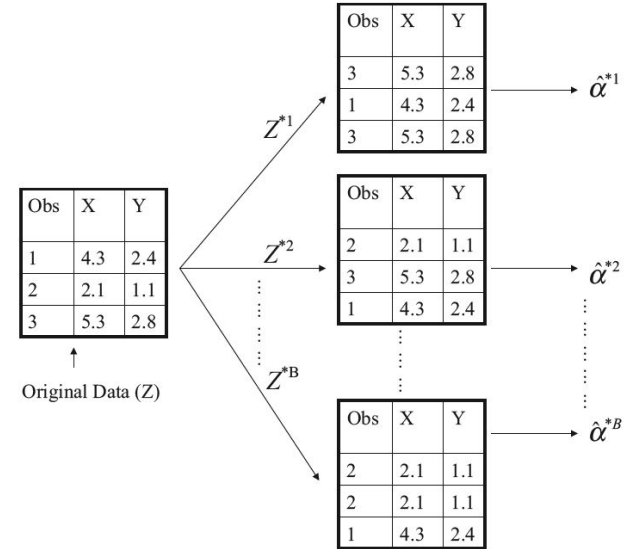
- Suffer from high variance
 - Different splits of training data → quite different results
- Random Forests, and Boosting reduce the variance
 - These are general purpose procedures

Bagging



Bootstrap

- Creates multiple datasets sampled with replacement
- Used to quantify the uncertainty associated with a given estimator



Bootstrap

- *Averaging a set of observations reduces the variance*
- Take many training sets, train separate models and average the resulting predictions

Bagging

- Compute B different models using B separate training sets

$$\hat{f}_{\text{bag}}(x) = \frac{1}{B} \sum_{b=1}^B \hat{f}^{*b}(x).$$



Bagging

- Useful for decision trees (improves predictions)
- B trees are trained on the bootstrapped datasets
 - Trees are grown deep without pruning
 - High variance and low bias
 - Aggregating → low variance

Bagging

- Prediction aggregation
 - Average for regression
 - Majority voting for classification



Random Forests



Random Forests

- Improvement over bagged trees
 - Via decorrelating them



Random Forests

- Similar to bagging, we build several trees
- When building trees
 - During a split, a random subset of predictors are chosen as candidates
 - Instead of all the 'p' predictors, only a random sample of 'm' ($\sim \sqrt{p}$) are allowed to conduct split

Random Forests

- Suppose one strong predictor and multiple moderate predictors are present in the data
- Bagging → most trees use the strong predictor at the top



Random Forests

- → Most of them will be similar → predictions will be correlated
- Averaging doesn't lead to a large reduction in variance



Random Forests

- RF overcome this by forcing each split to use a subset of predictors
- Majority of the splits do not consider the strong predictor
- → decorrelating the trees

Boosting



Boosting

- Bagging → multiple copies → trees are learned independently
- Boosting → Trees are grown sequentially
 - each tree is grown using information from previously grown trees

Boosting

- Does not involve bootstrap sampling
- instead each tree is fit on a modified version of the original data set

Boosting

- Given the current model, we fit a decision tree to the residuals from the model.
- Fit a tree using the current residuals, rather than the outcome Y , as the response.

Boosting for Regression Trees

1. Set $\hat{f}(x) = 0$ and $r_i = y_i$ for all i in the training set.
2. For $b = 1, 2, \dots, B$, repeat:
 - (a) Fit a tree \hat{f}^b with d splits ($d + 1$ terminal nodes) to the training data (X, r) .
 - (b) Update \hat{f} by adding in a shrunk version of the new tree:

$$\hat{f}(x) \leftarrow \hat{f}(x) + \lambda \hat{f}^b(x). \quad (8.10)$$

- (c) Update the residuals,

$$r_i \leftarrow r_i - \lambda \hat{f}^b(x_i). \quad (8.11)$$

3. Output the boosted model,

$$\hat{f}(x) = \sum_{b=1}^B \lambda \hat{f}^b(x). \quad (8.12)$$



Rough



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