UVA CS 6316: Machine Learning

Lecture 18c: More and Extra about Boosting

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Boosting:

- Learners are ordered: Each learner tries to reduce error (residual) on "hard" examples (those misclassified by earlier learners).
- ADABOOST: weight hard samples more;
- GRADIENT BOOST: use residual to train later models. Reduces bias and possibly variance compared to base learners.
- Gradient-boosted decision trees (GBDT) often gives state-of-theart performance on simple classification tasks, e.g. XGBOOST.
- Neural networks are used fairly often with bagging, but rarely with boosting.
- Decision trees work well in both bagging and boosting.

From Stanford CV class

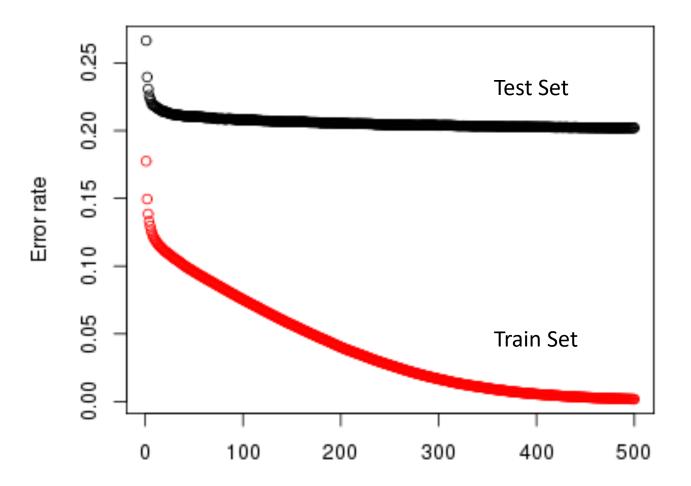
Boosting

- Sequential algorithm where at each step, a weak learner is trained based on the results of the previous learner.
- Two main types:
 - Adaptive Boosting: Reweight datapoints based on performance of last weak learner. Focuses on points where previous learner had trouble. Example: AdaBoost.
 - **Gradient Boosting**: Train new learner on residuals of overall model. Constitutes gradient boosting because approximating the residual and adding to the previous result is essentially a form of gradient descent. Example: XGBoost.

- XGBoost is a very efficient Gradient Boosting Decision Tree implementation with some interesting features:
- **Regularization:** Can use L1 or L2 regularization.
- Handling sparse data: Incorporates a sparsity-aware split finding algorithm to handle different types of sparsity patterns in the data.
- Weighted quantile sketch: Uses distributed weighted quantile sketch algorithm to effectively handle weighted data.
- **Block structure for parallel learning:** Makes use of multiple cores on the CPU, possible because of a block structure in its system design. Block structure enables the data layout to be reused.
- Cache awareness: Allocates internal buffers in each thread, where the gradient statistics can be stored.
- **Out-of-core computing:** Optimizes the available disk space and maximizes its usage when handling huge datasets that do not fit into memory.

XGBoost (an example performance figure)

Error rate Vs. Trees



Number of trees

- Task is to estimate target continuous function F(x). We measure goodness of estimation with loss function L(y, F(x)).
- Gradient boosting assumes that:
- $F(x) = \alpha_0 + \alpha_1 h_1(x) + \dots + \alpha_M h_M(x)$
- Basic Gradient boosting workflow:
 - 1. Initialize $F_0(x) = \alpha_0$
 - 2. Estimate α_m and $h_{m(x)}$ such that:

 $L(y, F_{m-1}(x) + \alpha_m h_m(x)) < L(y, F_{(m-1)}(x))$

- 3. Update $F_m(x) = F_{m-1}(x) + \alpha_m h_m(x)$
- 4. Repeat from 2, M times.

 $L(y, F_{m-1}(x) + \alpha_m h_m(x)) < L(y, F_{(m-1)}(x))$

If we can find a vector r_m that we can plug in here to make this equation true, we can train a basic learner $h_m(x)$ to predict r_m from x!

We are basically searching for a vector that points to the direction that reduces our loss... does that sound familiar?

Gradient descent!

- By solving a simple 1D optimization problem, we could also find the optimal α_m for each step, by computing:
 - $\alpha_m = argmin_{\gamma}L(y, F_{m-1}(x) + \gamma h_m(x))$
- This gives us an updated Gradient Boosting algorithm:
 - 1. Initialize $F_0(x) = \alpha_0$
 - 2. Compute negative gradient per observation: $r_{m_i} = -\frac{\partial L(y_i, F_{m-1}(x_i))}{\partial F_{m-1}(x_i)}$
 - 3. Train base learner $h_m(x)$ on predicting the gradients r_{m_i}
 - 4. Compute α_m with line search strategy
 - 5. Update $F_m(x) = F_{m-1}(x) + \alpha_m h_m(x)$
 - 6. Repeat from 2, M times.

- Where do the residuals come in?
- If we consider Mean Squared Error as our loss function, the perobservation gradient is:

•
$$\frac{\partial L(y_i, F_m(x_i))}{\partial F_m(x_i)} = \frac{\partial \left(\frac{1}{2n} \sum_i (y_i - F_m(x_i))^2\right)}{\partial F_m(x_i)} = \frac{\partial \left(\frac{1}{2} (y_i - F_m(x_i))^2\right)}{\partial F_m(x_i)} = y_i - F_m(x_i)$$

• The derivation we found before works with any loss function.

Gradient Tree Boosting

• When dealing with decision trees, we can take the concept further by selecting a specific α_m for each of the tree's regions. The output of a tree is:

•
$$h_m(x) = \sum^{J_m} b_{jm} \mathbf{1}_{R_{jm}}(x)$$

• The model update rule becomes:

•
$$F_m(x) = F_{m-1}(x) + \sum_{j=1}^{J_m} \alpha_{jm} \mathbf{1}_{R_{jm}(x)}$$

• $\alpha_{jm} = argmin_{\gamma} \sum_{x_i \in R_{jm}} L(y_i, F_{m-1}(x_i) + \gamma)$

J_m: Disjoint regions partitioned by the tree

R_{jm}: Number of leaves

- Three main forms of gradient boosting are supported:
- Gradient Boosting algorithm, as we defined above.
- Stochastic Gradient Boosting with sub-sampling at the row, column and column per split levels.
 - Random procedure where we subsample observations and features
- **Regularized Gradient Boosting** with both L1 and L2 regularization.
 - add a regularization term to the loss function that we are optimizing: $L_R(y, F(x)) = L(y, F(x)) + \Omega(F)$ Where $\Omega(F) = \gamma T + \frac{1}{2}\lambda ||w||^2$

T: Number of leaves

W: Leaf weights: prediction of each leaf

- Remember, we still want to find the tree structure that minimizes our loss, which means best score structure. Doing this for all possible tree structures is unfeasible.
- A greedy algorithm that starts from a single leaf and iteratively adds branches to the tree is used instead.

- XGBoost adds multiple other important advancements that make it state of the art in several industrial applications.
- In practice:
- Can take a while to run if you don't set the n_jobs parameter correctly
- Defining the eta parameter (analogous to learning rate) and max_depth is crucial to obtain good performance.
- Alpha parameter controls L1 regularization, can be increased on high dimensionality problems to increase run time.

- General approach to parameter tuning:
- Cross-validate learning rate.
- Determine the optimum number of trees for this learning rate. XGBoost can perform cross-validation at each boosting iteration for this, with the "cv" function.
- **Tune tree-specific parameters** (max_depth, min_child_weight, gamma, subsample, colsample_bytree) for chosen learning rate and number of trees.
- Tune **regularization parameters** (lambda, alpha).

LGBM

- Stands for Light Gradient Boosted Machines. It is a library for training GBMs developed by Microsoft, and it competes with XGBoost.
- Extremely efficient implementation.
- Usually much faster than XGBoost with low hit on accuracy.
- Main contributions are two novel techniques to speed up split analysis: Gradient based one-side sampling and Exclusive Feature Building.
- Leaf-wise tree growth vs level-wise tree growth of XGBoost.

