

# **Random Forests™**

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# RF motivation

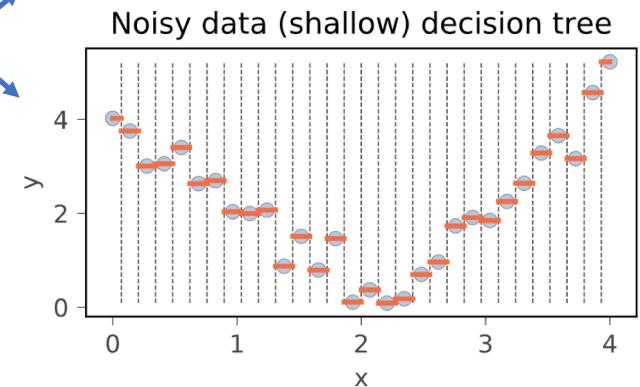
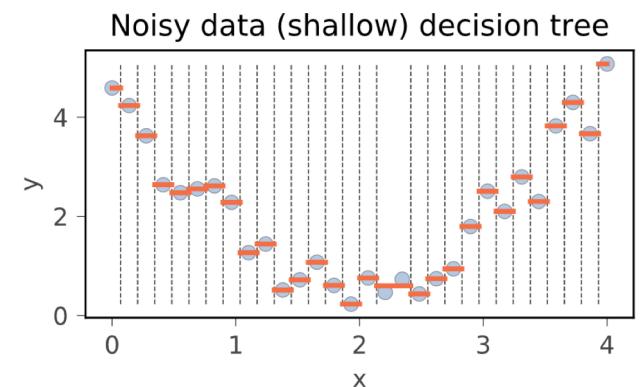
Leo Breiman (1996) introduced bagging  
then Random Forests (2001)

<https://www.stat.berkeley.edu/~breiman/randomforest2001.pdf>

- Decision trees can often get training errors close to zero because we can grow very large trees to partition the feature space into tiny regions with 1 or just a few observations/samples; trees have very **low bias**
- The downside is that decision trees overfit or, using stats nerds terminology, decision trees have **high variance** and don't generalize well
- Goal: keep the low bias, but increase the generality (reduce the variance)

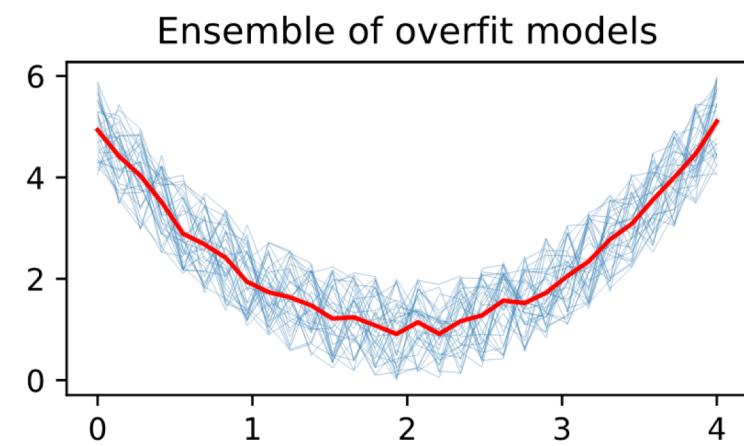
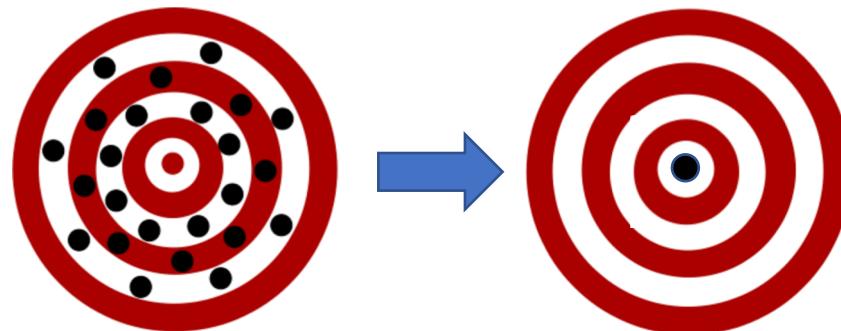
# Variance is the key word here

- Decision trees are sensitive to quirks in the data and so similar but different training sets drawn from the same distribution can yield very different models (parameters)
- (That's the definition of model variance)
- Ex: two i.i.d. X training sets fit by regr tree
- The predictions made by decision trees trained on i.i.d. sets bounce around, but their average approaches the correct value (same for i.d. sets)



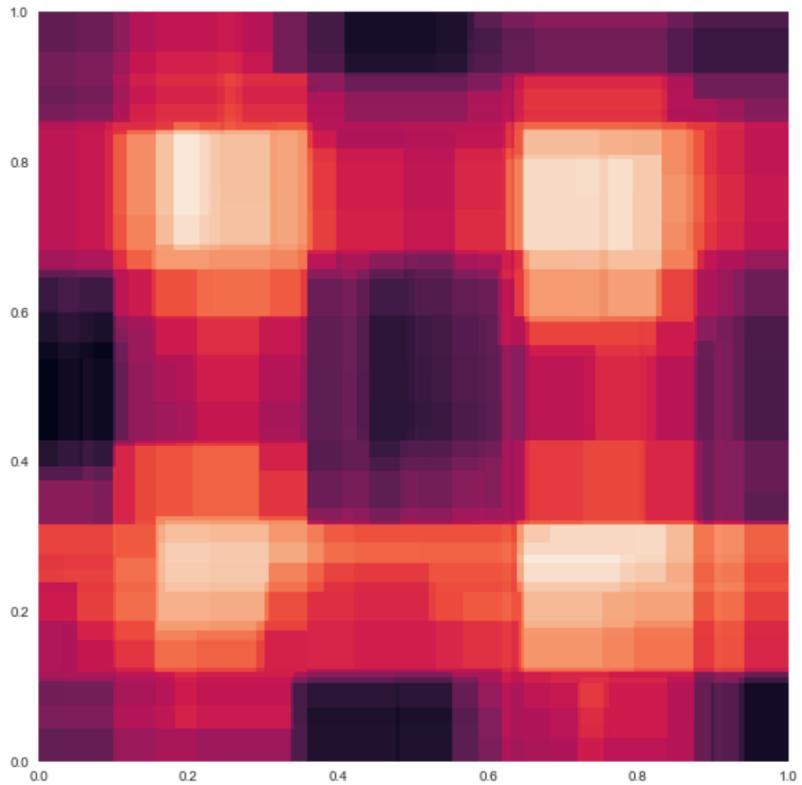
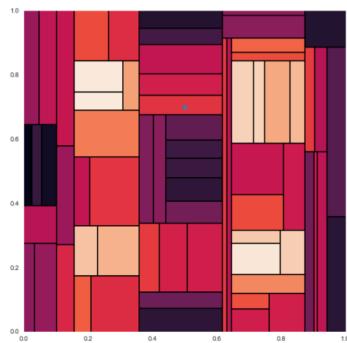
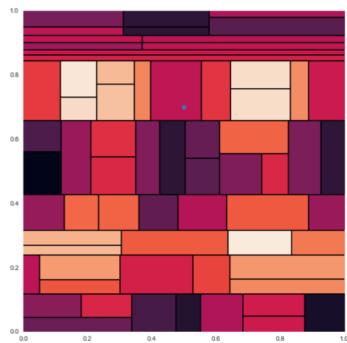
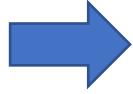
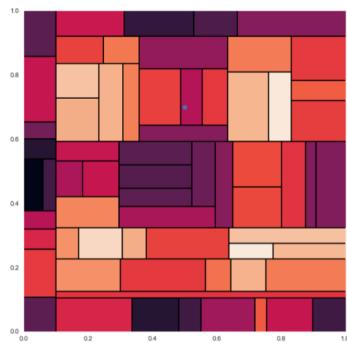
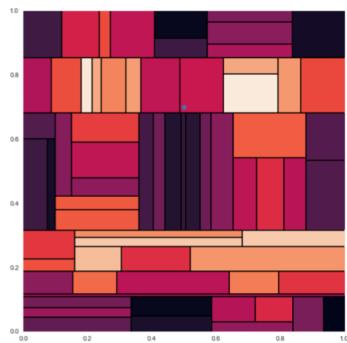
# Ensemble of high-variance decision trees

- Decision trees are inconsistent but are accurate on average across multiple identically distributed (i.d.) training sets
- Pretend we do have multiple i.d. X's and multiple fit models
- Then, we can average predictions from multiple models to get an accurate prediction:



Target image credit: <http://ficscience.blogspot.com/2011/02/technicalities-of-technical-terms.html>

# RF: Overlapping feature space partitions



Images generated by USF MSDS alumnus Tyler White  
<https://gist.github.com/tylerwx51/fc8b316337833c877785222d463a45b0>

# Ensemble's effect on bias and variance

- Train  $T$  trees on  $T$  identically distributed  $X$  data sets (doesn't have to be independent)
- The average of the tree predictions is the same as the expected prediction from any tree trained on one of the  $X$
- Central limit theorem says that if variance of  $T$  i.i.d. random variables is  $\sigma^2$ , the variance of the average is  $\frac{\sigma^2}{T}$
- So, as we add trees, the variance of the prediction (average from ensemble trees) tightens towards true prediction!
- Problem: we only have one  $X$  training data set

# Bagging (Bootstrap aggregation)

- We only have one training set,  $X$ , but can *bootstrap* to get as many as we want from the same distribution
- These bootstrapped sets are i.d. but not independent (some overlap)
- *Bootstrap*: from  $n$  records, randomly select  $n$  with replacement
- Each bootstrapped data set will have about 2/3 of the unique overall records
- *Aggregation*: average the ensemble predictions  
(majority vote for classification)
- Analogy: real estate agents sent out to sample NYC apartments/prices; there will be some overlap but average of all agents' predictions is more accurate than an individual's

# show diff trees

# Problem: bootstrapping is i.d. not i.i.d.

- If just identically distributed and not necessarily independent, with positive pairwise correlation  $\rho$  between trees, then variance of the average for random vars is: (Eqn 15.1 ESLII Hastie *et al*):

$$\rho\sigma^2 + \frac{1 - \rho}{T}\sigma^2$$

- As  $T$ , (num trees) grows, 2<sup>nd</sup> term disappears but correlation of trees would still have variance proportional to  $\rho$  (proportional to  $\sigma^2$ , variance of any single tree predictor)
- ESLII: "...the size of the correlation of pairs of bag trees limits the benefits of averaging."
- So, bagging helps to reduce variance but still has correlation  $\rho > 0$

# What does “correlated trees” mean?

- Model parameters are correlated; “not independent” better term
- **Analogy.** Real estate agents sampling bias: if they visit apartments they observe other agents visit, variance might not shrink with more agents; worst-case: they all visit the same apartments; best case is they choose all apts independently
- **Worst-case.** Imagine replicating single decision tree  $T$  times, variance of average prediction would be same as variance of single tree (i.e., bad strategy); if  $\rho=1$ , variance of avg =  $\sigma^2$
- **Best-case.** if  $\rho=0$ , trees are uncorrelated and variance of average converges to 0 as  $T$  grows

$$\rho\sigma^2 + \frac{1-\rho}{T}\sigma^2$$

# Bagging summary

- Adding bagged trees does not increase **bias**; the avg of tree predictions is same as the expected prediction from any single bootstrapped tree
- Keep in mind, however, that a tree fit to bootstrapped data is only using 2/3 of the data and so each bagged tree will have higher bias than single decision tree fit to entire data set
- Bagging does reduce **variance**, it is averaging predictions from lots of non-identical models afterall
- Degree of tree similarity,  $\rho$ , influences ability to reduce variance to 0
- If  $\rho=1$ , no reduction in variance, but  $\rho=0$  means adding trees reduces variance to 0 asymptotically

$$\rho\sigma^2 + \frac{1-\rho}{T}\sigma^2$$

# De-correlating trees to get Random Forests (RFs)

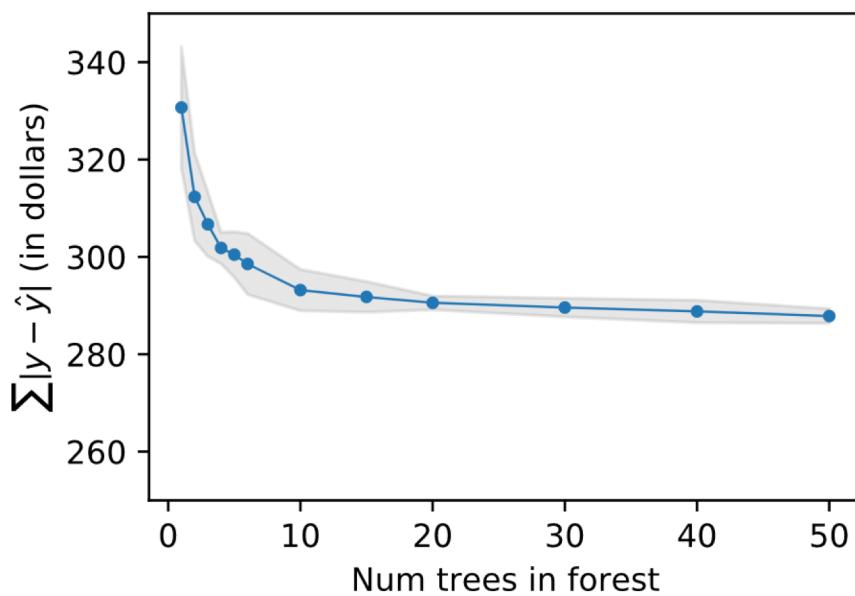
- Bagging overcomes most of the overfitting (Hastie *et al* p589 say  $\rho$  usually small like 0.05)
- Can do a little better de-correlation of the trees to reduce  $\rho$
- Restrict the available features at each split (random selection of  $m$ )
- Choose max features per split,  $m \leq p$ , such as  $\text{sqrt}(p)$
- Imagine one strongly predictive var out of  $p$ , then all trees would be similar; initial root splits, and many others, will likely be same
- Make sure chance of selecting predictive variables ( $m/p$ ) is high enough to find them (See ESLII p596)
- Let validation error be your guide to  $m$  (and other hyperparameters)

$$\rho\sigma^2 + \frac{1 - \rho}{T}\sigma^2$$

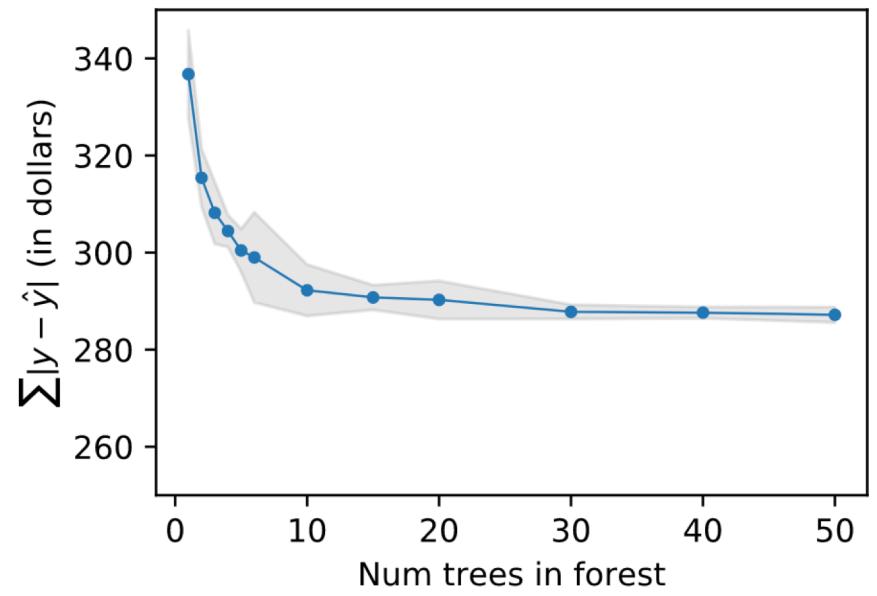
# Compare bagged trees vs RF (Rent)

Very little difference with just p=4 features

Rent bagged forest error vs num trees  
5 trials, 41055 training obs., 7245 test  
4 max features



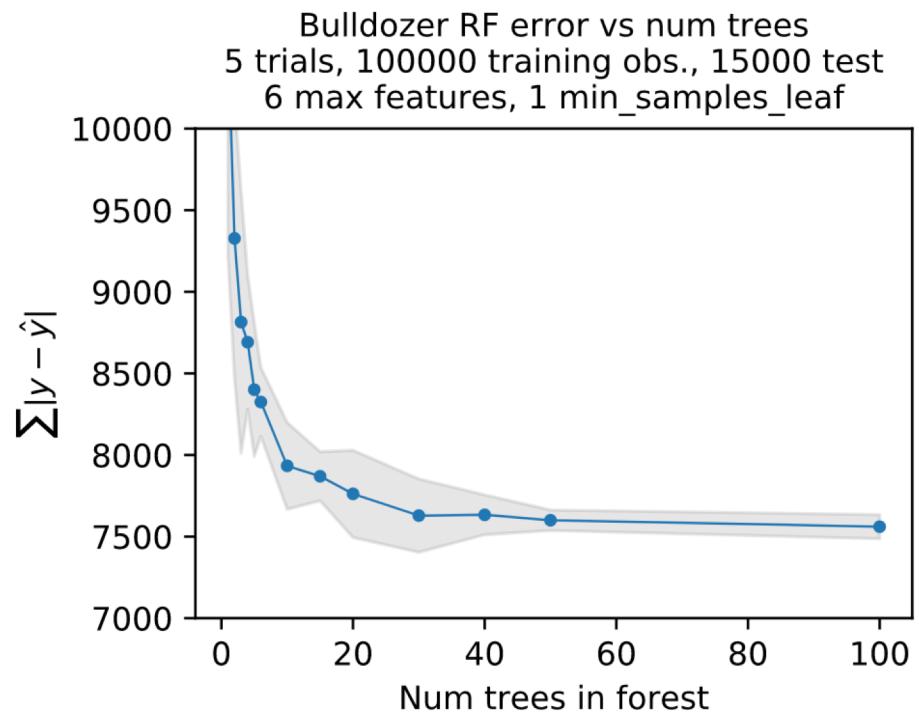
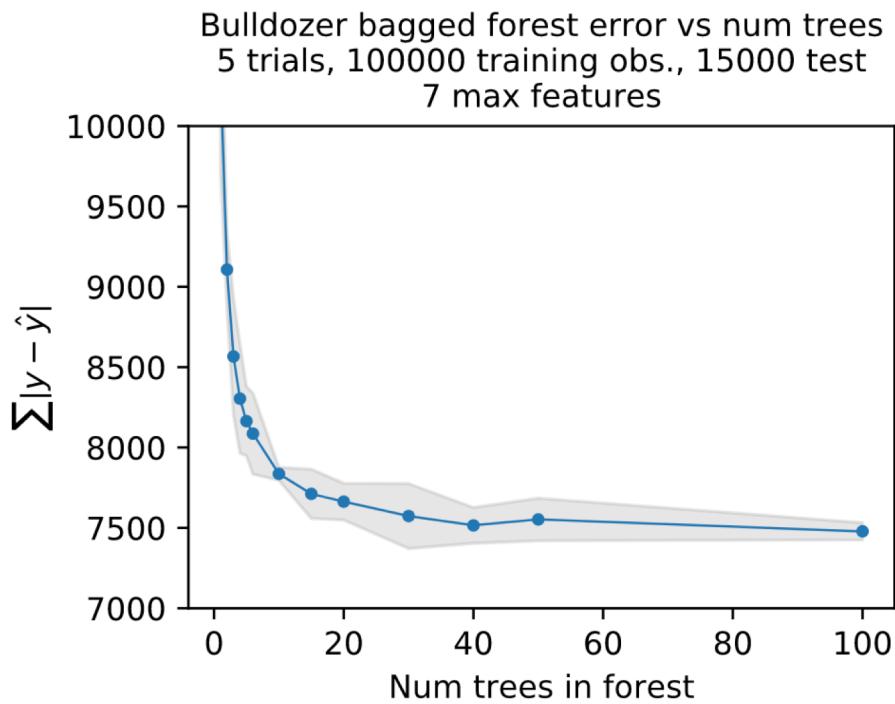
Rent RF error vs num trees  
5 trials, 41055 training obs., 7245 test  
1 max features



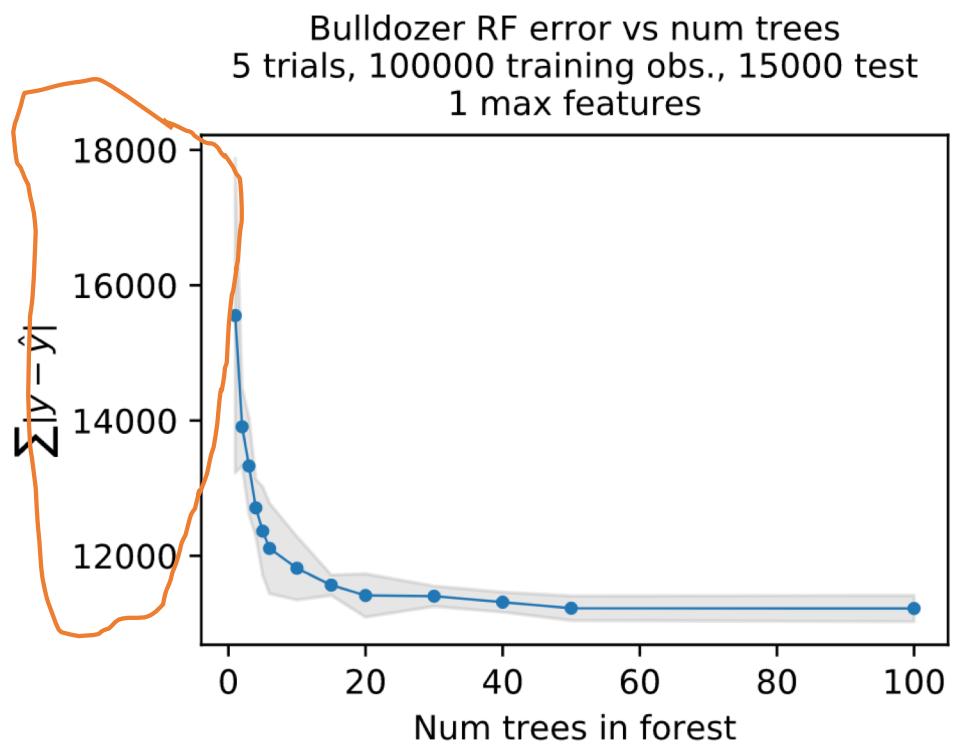
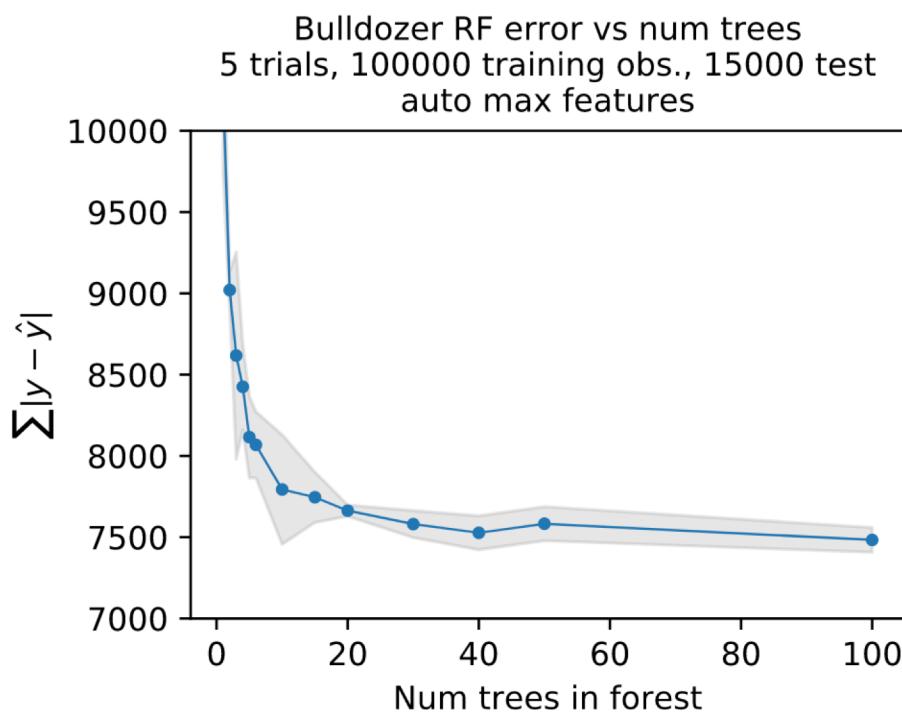
Some data sets will be more sensitive to max features than others

# Compare bagged trees vs RF (bulldozer)

For this data, using all p=7 is best

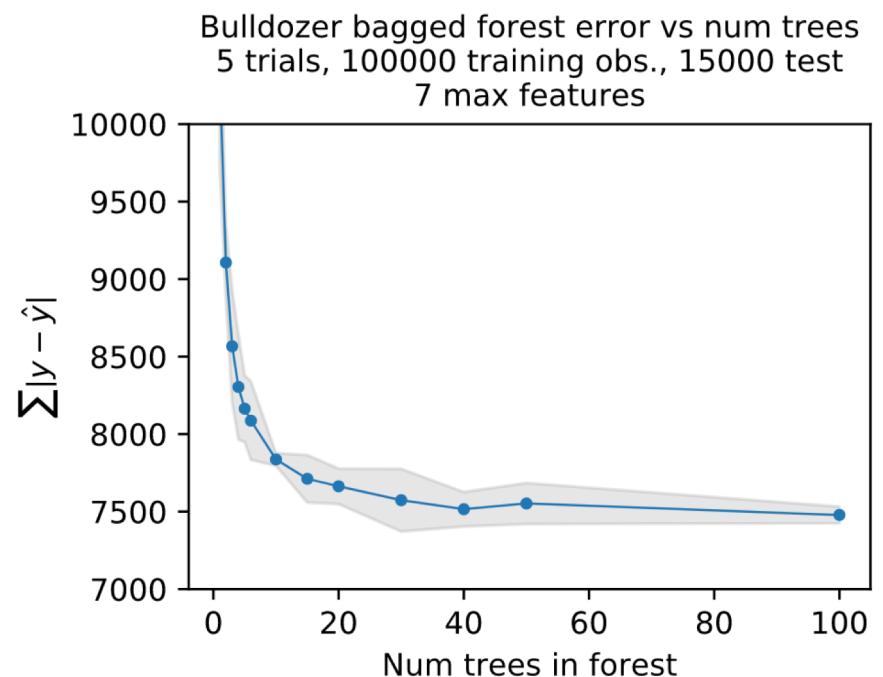


# If max\_features too low, high bias



# Effect of forest size on bias

- Why does accuracy improve greatly (initially) as we add trees?
  - Each tree sees only 2/3 of data so adding bootstrapped trees increases use of training data
- Why does accuracy asymptotically approach a minimum instead of continual improvement?
  - With enough trees, ensemble sees 100% of the training data; it's approaching the bias of single decision tree in ideal world where it's not overfit



# Properties (Breiman 2001)

- p4 “*Random forests do not overfit as more trees are added*” **Why?**
  - Adding more trees actually REDUCES variance, and overfitting==variance
  - New trees get averaged in so each new tree has less individual effect
- p7 “*It's relatively robust to outliers and noise*” **Why?**
  - Outliers get shunted to their own leaf since doing so reducing loss function, particularly if squared-error is used; noise vars are predictive so not chosen as split vars
- p10 Bagging helps more the more unstable the model. **Why?**
  - Averaging is a smoothing operator, squeezing predictions to centroid
  - If model is low variance already, there is no point in bagging

# Properties continued

- RFs are scale and range insensitive in features and target y **Why?**
  - Comparing feature values in decision nodes not doing math on them
  - Computing mean or mode of y to predict
- ESLII p596 “*Classifiers are less sensitive to variance [than regressors]*” **Why?**
  - (not sure haha) I believe it has something to do with mode vs mean (mode is same until a threshold whereas mean is influenced by any value added, unless it is also the mean)

# Bootstrapping vs subsampling

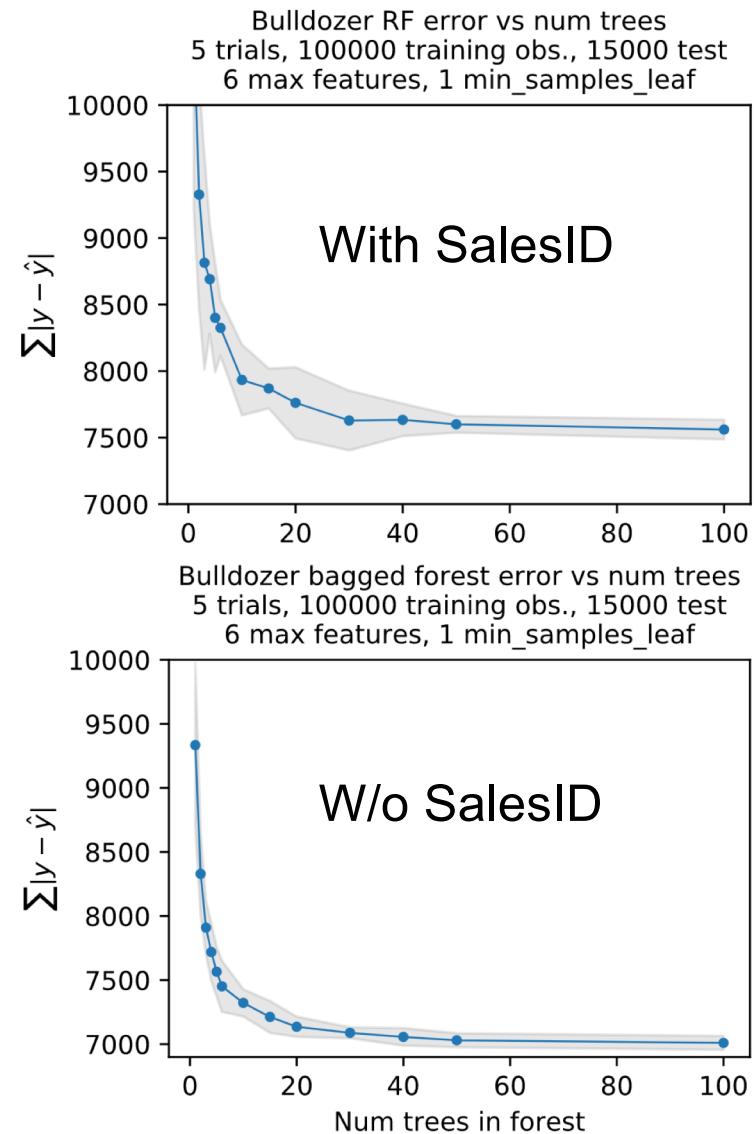
- Bootstrapping is sampling with replacement vs subsampling w/o replacement
-  Friedman and Hall (2000): subsampling also works, showing that training trees with  $n/2$  subsamples is similar in bias/variance to bagging <http://statweb.stanford.edu/~jhf/ftp/bag.pdf>
- Smaller training set is a big win in terms of speed
- Using even smaller fractions of  $n$  improve generality (reduce variance) because trees are less correlated (they work on different data chunks); note that each tree would become more biased as  $n$  subsample size decreases

# RF Tuning strategy

- Good news: very little tuning needed
- Start with maybe 20 trees and work upwards til validation error stops getting better; or just pick 100
- Sklearn uses `max_features=1.0` (all features) by default; try dropping this using  $\text{sqrt}(p)$ ,  $\log(p)$ , or similar; ESLII suggests  $p/3$  for regression and  $\text{sqrt}(p)$  for classification
- Try adjusting min samples per leaf: 1, 3, 5, 10, 25, 100
- Goal: minimize validation error
- Can also try grid search, but I never bother;  
Start with num trees, then tune the others

# Feature engineering beats model tuning

- SalesID: unique record ID, and is never seen again in future predictions
- Is that useful for prediction? No
- Does the model think it's useful? Yes
- Model is overfit not on noise but on falsely-predictive feature
  - Could be that sales ID correlates with inflation or change in type of models sold in auction creates “trend” in sale prices
- A case where using LESS data improves the model a lot (\$500 diff)
- Dropping useless features also often gives a small bump



# Extremely randomized trees (Geurts *et al* 2006)

- The variable/value pair is highly sensitive to the training set, and responsible much of error rate
- “*The optimal cut-point was shown to depend very strongly on the particular learning sample used...this cut-point variance appeared to be responsible for a significant part of the error rates of tree-based methods.*” <https://link.springer.com/article/10.1007/s10994-006-6226-1>
- Geurts wondered if more randomness could reduce variance further
- Pick random split value in  $\min(X[:,j]) \dots \max(X[:,j])$ , ignoring y!
- Like RF, select  $m \leq p$  variables and choose var/value with lowest loss
- Fits using entire X training set, not bootstrap and not subsample (trying to reduce bias)
- Our use of just 11 (not n) X candidate values in the project is similar (an effort to reduce variance and increase speed)

# The RF algorithms

# Fitting RFs

**Algorithm:**  $RFpredict_{regr}(\{T_1..T_{ntrees}\}, x)$

**Let**  $leaves = \{leaf(T_i.root, x) \forall i = 1..ntrees\}$  (*set of leaves reached by x*)

$$nobs = \sum_{i=1}^{ntrees} |leaves_i|$$

$$ysum = \sum_{i=1}^{ntrees} \sum_{y \in leaves_i} y$$

**return**  $\frac{1}{nobs} ysum$  (*weighted avg*)

For regression, pass in loss = MSE or stddev

For classifier, pass in loss = gini

# Fitting a single tree in RF

```
1 Algorithm:  $RFdtreefit(X, y, loss, max\_features, min\_samples\_leaf)$ 
2 if  $|X| < min\_samples\_leaf$  then return Leaf( $y$ )
3  $col, split = RFbestsplit(X, y, loss, max\_features)$ 
4 if  $col = -1$  then return Leaf( $y$ )
5  $lchild = fit(X[X \leq split], y[X \leq split])$ 
6  $rchild = fit(X[X > split], y[X > split])$ 
7 return  $DecisionNode(col, split, lchild, rchild)$ 
```

Same as decision tree except  
we pass `max_features` to  
`RFbestsplit()`

# Finding best split in decision node in RF

```
1 Algorithm: RFbestsplit( $X$ ,  $y$ , loss, max_features)
2   best = ( $col = -1$ ,  $split = -1$ ,  $loss = loss(y)$ )
3   vars = pick max_features variables from all  $p$  ← Only diff with decision tree
4   for  $j \in vars$  do
5     candidates = pick 11 values from  $X_{-,j}$ 
6     foreach  $split \in candidates$  do
7        $yl = y[X \leq split]$ 
8        $yr = y[X > split]$ 
9       if  $|yl| = 0$  or  $|yr| = 0$  then continue
10       $l = \frac{loss(yl) + loss(yr)}{2}$  ← Should pick midpoint between
11      if  $l = 0$  then return  $col$ ,  $split$ 
12      if  $l < best[loss]$  then  $best = (col, split, l)$ 
13    end
14  end
15  return  $col$ ,  $split$ 
```

Pick, say, 11 not all possible X values. We get better generality and code is much faster!

Only diff with decision tree

Should pick midpoint between split value and next smallest X



# Simplest RF prediction (ESLII p588)

- But doesn't use all information to make best prediction
- Should use weighted averages / votes

Regression:  $\hat{f}_{\text{rf}}^B(x) = \frac{1}{B} \sum_{b=1}^B T_b(x).$

Classification: Let  $\hat{C}_b(x)$  be the class prediction of the  $b$ th random-forest tree. Then  $\hat{C}_{\text{rf}}^B(x) = \text{majority vote } \{\hat{C}_b(x)\}_1^B.$

# RF prediction

Weighted average of y values among the leaves reached by running x down each tree

**Algorithm:**  $RFpredict_{regr}(\{T_1..T_{ntrees}\}, x)$

**Let**  $leaves = \{leaf(T_i.root, x) \forall i = 1..ntrees\}$

$nobs = \sum_{i=1}^{ntrees} |leaves_i|$

$ysum = \sum_{i=1}^{ntrees} \sum_{y \in leaves_i} y$

**return**  $\frac{1}{nobs} ysum$

Weight each leaf's vote by  $|leaf|$  among the leaves reached by running x down each tree

**Algorithm:**  $RFpredict_{class}(\{T_1..T_{ntrees}\}, x)$

$counts[k] = 0 \forall \text{ classes } k$

**foreach**  $i = 1..ntrees$  **do**

$leaf = leaf(T_i.root, x)$      (*leaf reached by x*)

$y' = \text{mode}(y \in leaf)$      (*most common y in leaf*)

$counts[y'] += |leaf|$      (*track count of leaf modes*)

**end**

**return**  $\text{argmax}(counts)$