

# Ensemble

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# Which Classifier/Model to Choose?

- Possible strategies:
- Go from simplest model to more complex model until you obtain desired accuracy
- Discover a new model if the existing ones do not work for you
- Combine all (simple) models

# Common Strategy: Bagging

(Bootstrap Aggregating)

Originally designed for combining multiple models,  
to improve classification “stability” [Leo Breiman, 94]

Uses random training datasets  
(sampled from one dataset)

<http://statistics.about.com/od/Applications/a/What-Is-Bootstrapping.htm>

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Consider the data set  $S = \{(X_i, Y_i)\}_{i=1, \dots, n}$

- Pick a sample  $S^*$  with **replacement** of size  $n$   
( $S^*$  called a “bootstrap sample”)

$$S \rightarrow x = \begin{bmatrix} 1 & 2 & 3 & 4 \\ 9 & 10 & 11 & 12 \\ 20 & 21 & 22 & 23 \\ 5 & 6 & 7 & 8 \end{bmatrix} y = \begin{bmatrix} 1 \\ 1 \\ -1 \\ -1 \end{bmatrix}$$

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## (Bootstrap Aggregating)

Consider the data set  $S = \{(X_i, Y_i)\}_{i=1, \dots, n}$

- Pick a sample  $S^*$  with **replacement** of size  $n$   
( $S^*$  called a “bootstrap sample”)
- Train on  $S^*$  to get a classifier  $f^*$
- Repeat above steps  $B$  times to get  $f_1, f_2, \dots, f_B$
- Final classifier  $f(x) = \text{majority}\{f_b(x)\}_{j=1, \dots, B}$

# Common Strategy: Bagging

Why would bagging work?

- Combining multiple classifiers reduces the variance of the final classifier

When would this be useful?

- We have a classifier with high variance

# Bagging decision trees

Consider the data set  $S$

- Pick a sample  $S^*$  with replacement of size  $n$
- Grow a decision tree  $T_b$
- Repeat  $B$  times to get  $T_1, \dots, T_B$
- The final classifier will be

$$f(x) = \text{majority}\{f_{T_b}(x)\}_{b=1, \dots, B}$$

# Random Forests

Almost identical to bagging decision trees,  
except we introduce some randomness:

- Randomly pick  $m$  of the  $d$  available attributes, at every split when growing the tree  
(i.e.,  $d - m$  attributes ignored)

Bagged **random** decision trees =  
**Random forests**

# What are our Hyper-Parameters in Random Forest

$m$  = *Number of randomly chosen attributes*

Usual values for  $m = \sqrt{d}, 1, 10$

$d$  is number of dimensions or features or attributes

How to optimize  $m$ ?      Cross-Validation

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$B$  = *Number of models or decision trees in Random Forest*

Keep adding trees until training error stabilizes (reaches to a plateau)

# Important points about random forests

## Algorithm (hyper) parameters

- Size/#nodes of each tree
  - as in when building a decision tree
- May randomly pick an attribute, and may even randomly pick the split point!
  - Significantly simplifies implementation and increases training speed
    - PERT - Perfect Random Tree Ensembles  
<http://www.interfacesymposia.org/I01/I2001Proceedings/ACutler/ACutler.pdf>
    - Extremely randomized trees  
<http://orbi.ulg.be/bitstream/2268/9357/1/geurts-mlj-advance.pdf>