Outline

• Ensemble classifiers – a motivation exercise
• Combining classifiers into ensembles – general scheme
• Generating random samples by bootstrapping
• Bagging vs. boosting
• Bagging – example classifier
• Random Forests
Consider the following task – we have a binary classification problem and a number of predictors, each with error less than 0.5. Will the resulting majority voting ensemble have an error lower than the single classifiers?

Depends on the accuracy and the diversity of the base learners!

Illustrative example

Particular settings – assume that you have
• 21 classifiers
• each with error $p = 0.3$
• their outputs are statistically independent

Compute the error of the ensemble under these conditions!
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Depends on the accuracy and the diversity of the base learners!
Ensemble classifiers – a motivation exercise

Consider the following task – we have a binary classification problem and a number of predictors, each with error less than 0.5. Will the resulting majority voting ensemble have an error lower than the single classifiers?

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**Illustrative example**

Particular settings – assume that you have

- 21 classifiers
- each with error $p = 0.3$
- their outputs are *statistically independent*

**Compute the error of the ensemble under these conditions!**
Solution of the exercise

How many classifiers will produce error output?
Key idea: The number of them will be binomially distributed! \( \sim \text{Bi}(21, 0.3) \)

\begin{verbatim}
> plot(0:21, dbinom(0:21, 21, 0.3))
> dbinom(11, 21, 0.3)
[1] 0.01764978
> pbinom(10, 21, 0.3)
[1] 0.9736101
\end{verbatim}

Conclusion: Accuracy of the ensemble will be more than 97.3%!
General scheme of combining classifiers

Training data

$\mathcal{L}_1$

$\mathcal{L}_2$

$\mathcal{L}_K$

Combining function

$h(x_i)$

$K$ base learners
$L_1, \ldots, L_K$ that produce $K$ hypotheses $h_1, \ldots, h_K$

Final ensemble prediction
Resampling can be used as a way to produce diversity among base learners

- Distribute the training data into $K$ portions
- Run the learning process to get $K$ different models
- Collect the output of the $K$ models use a combining function to get a final output value
Bootstrapping principle

- New data sets $Data_1, \ldots, Data_K$ are drawn from $Data$ with replacement, each of the same size as the original $Data$, i.e. $n$.

- In the $i$-th step of the iteration, $Data_i$ is used as a training set, while the examples $\{x \mid x \in Data \land x \notin Data_i\}$ form the test set.
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• In the $i$-th step of the iteration, $Data_i$ is used as a training set, while the examples $\{x | x \in Data \land x \notin Data_i\}$ form the test set.

• The probability that we pick an instance is $1/n$, and the probability that we do not pick an instance is $1 - 1/n$. The probability that we do not pick it after $n$ draws is $(1 - 1/n)^n \approx e^{-1} \approx 0.368$.

• It means that for training the system will not use $36.8\%$ of the data, and the error estimate will be pessimistic. So the solution is to repeat the process many times.
Ensemble methods – key ideas

- combining the classification results from different classifiers to produce the final output
- using (un)weighted voting
- different training data – e.g. bootstrapping
- different features
- different values of the relevant parameters
- performance: complementarity $\rightarrow$ potential improvement

Two fundamental approaches

- **Bagging** works by taking a bootstrap sample from the training set
- **Boosting** works by changing the weights on the training set
Are ensembles effective?

Combining multiple learners
- the more complementary the learners are, the more useful their combining is
- the simpliest way to combine multiple learners is voting
- in weighted voting the voters (= base-learners) can have different weights

Unstable learning
- learning algorithm is called unstable if small changes in the training set cause large differences in generated models
- typical unstable algorithm is the decision trees learning
- bagging or boosting techniques are a natural remedy for unstable algorithms
• Bagging is a voting method that uses slightly different training sets (generated by bootstrap) to make different base-learners. Generating complementary base-learners is left to chance and to unstability of the learning method.

• Generally, bagging can be combined with any approach to learning.
Simple bagging algorithm

Bootstrap AGGregatING

1. for $i \leftarrow 1$ to $K$ do
2. $Train_i \leftarrow \text{bootstrap}(Data)$
3. $h_i \leftarrow \text{TrainPredictor}(Train_i)$

Combining function

- **Classification:** $h_{\text{final}}(x) = \text{MajorityVote}(h_1(x), h_2(x), \ldots, h_K(x))$
- **Regression:** $h_{\text{final}}(x) = \text{Mean}(h_1(x), h_2(x), \ldots, h_K(x))$
Random Forests

- Random Forest is another example of bagging
- an ensemble method based on decision trees and bagging
- builds a number of *independent* random decision trees and then uses voting
Outline

• Bagging vs. boosting

• Simple boosting trees – the regression case

• Adaptive boosting – classification with AdaBoost
• **Bagging**: each predictor is trained independently

• **Boosting**: each predictor is built on the top of previous predictors trained
  – Like bagging, boosting is also a voting method. In contrast to bagging, boosting actively tries to generate complementary learners by training the next learner on the mistakes of the previous learners.
Boosting combines the outputs of many “weak” classifiers (“rules of thumb”) to produce a powerful “committee.”

**Motivation**

- How to extract rules of thumb that will be the most useful?
- How to combine moderately accurate rules of thumb into a single highly accurate prediction rule?

**Basic idea**

- Boosting is a method that produces a very accurate predictor by combining rough and moderately accurate predictors.
- It is based on the observation that finding many rough predictors (rules of thumb) can be easier than finding a single, highly accurate predictor.
Simple boosting with regression trees

1. Initialization: Set $h(x) = 0$ and $r_i = y_i$ for all $i = 1, \ldots, n$ in the training set.

2. For $b = 1, \ldots, B$, repeat
   (a) Fit a tree $h^b$ with only $d$ splits to the training set $(X, r)$
   (b) Update $h$ by adding the new tree
       
       $$h(x) \leftarrow h(x) + \lambda h^b(x)$$
   (c) Update the residuals
       
       $$r_i \leftarrow r_i - \lambda h^b(x_i)$$

3. Output the boosted model

   $$h(x) = \sum_{b=1}^{B} \lambda h^b(x)$$
• The number of trees $B$

• The shrinkage parameter $\lambda$
  – A small positive number. This controls the rate at which boosting learns. Typical values are 0.01 or 0.001, and the right choice can depend on the problem. Very small $\lambda$ can require using a very large value of $B$ in order to achieve good performance.

• The number $d$ of splits in each tree
  – Trees with just $d = 1$ split are called “stumps”.
**AdaBoost** is a boosting method that repeatedly calls a given weak learner, each time with different distribution over the training data. Then we combine these weak learners into a strong learner.
AdaBoost is a boosting method that repeatedly calls a given weak learner, each time with different distribution over the training data. Then we combine these weak learners into a strong learner.

- originally proposed by Freund and Schapire (1996)

- great success
  - “AdaBoost with trees is the best off-the-shelf classifier in the world.” (Breiman 1998)
  - “Boosting is one of the most powerful learning ideas introduced in the last twenty years.” (Hastie et al, 2009)
Key questions
- How to choose the distribution?
- How to combine the weak predictors into a single predictor?
- How many weak predictors should be trained?

Schapire’s strategy: Change the distribution over the examples in each iteration, feed the resulting sample into the weak learner, and then combine the resulting hypotheses into a voting ensemble, which, in the end, would have a boosted prediction accuracy.
AdaBoost.M1 (Freund and Schapire, 1997) is the most popular boosting algorithm

- Consider a binary classification task with the training data

$$Data = \{\langle x_i, y_i \rangle : x_i \in X, y_i \in \{-1, +1\}, i = 1, \ldots, n\}$$

- We need to define distribution $D$ over $Data$ such that $\sum_{i=1}^{n} D_i = 1$.

- Assumption: a weak classifier $h_t$ has the property

$$\text{error}_D(h_t) < 1/2.$$
Adaboost (Adaptive Boosting) — key idea

Classifiers are trained on weighted versions of the original training data set, and then combined to produce a final prediction

- Training examples $\rightarrow h_1(x)$
- Weighted examples $\rightarrow h_2(x)$
- Weighted examples $\rightarrow h_3(x)$
- Weighted examples $\rightarrow h_M(x)$
Adaboost (Adaptive Boosting) — key idea

Classifiers are trained on weighted versions of the original training data set, and then combined to produce a final prediction

- Training examples $\rightarrow h_1(x)$
- Weighted examples $\rightarrow h_2(x)$
- Weighted examples $\rightarrow h_3(x)$
- \ldots
- Weighted examples $\rightarrow h_M(x)$

**Final hypothesis** $h(x) = \text{sign} \sum_{t=1}^{M} \alpha_t h_t(x)$, where $\alpha_t$ are computed by the boosting algorithm, and weight the contribution of each respective $h_t$
AdaBoost – iterative algorithm

- Initialize the training distribution \( D_1(i) = 1/n \) for \( i = 1, \ldots, n \)
- At each step \( t \)
  - Learn \( h_t \) using \( D_t \): find the weak classifier \( h_t \) with the minimum weighted sample error \( \text{error}_{D_t}(h_t) = \sum_{i=1}^{n} D_t(i) \delta(h(x_i) \neq y_i) \)
  - Set weight \( \alpha_t \) of \( h_t \) based on the sample error
    \[
    \alpha_t = \frac{1}{2} \ln \left( \frac{1 - \text{error}_{D_t}(h_t)}{\text{error}_{D_t}(h_t)} \right)
    \]
  - Update the training distribution
    \[
    D_{t+1} = D_t e^{-\alpha_t y_i h_t(x_i)} / Z_t \quad \text{where } Z_t \text{ is a normalization factor}
    \]
- Stop when impossible to find a weak classifier being better than chance
- Output the final classifier \( h_{\text{final}}(x) = \text{sign} \sum_{t=1}^{T} \alpha_t h_t(x) \)
AdaBoost – training data weighting

**Constructing** $D_t$

- On each round, the weights of incorrectly classified instances are increased so that the algorithm is forced to focus on the hard training examples.

- $D_1(i) = 1/n$ for $i = 1, \ldots, n$

- given $D_t$ and $h_t$ (i.e. update $D_t$):

\[
D_{t+1}(i) = \frac{D_t(i)}{Z_t} \cdot \begin{cases} 
  e^{-\alpha_t} & \text{if } y_i = h_t(x_i) \\
  e^{\alpha_t} & \text{if } y_i \neq h_t(x_i) 
\end{cases} = \frac{D_t(i)}{Z_t} e^{-\alpha_t y_i h_t(x_i)},
\]

where $Z_t$ is normalization constant $Z_t = \sum_i D_t(i) e^{-\alpha_t y_i h_t(x_i)}$

- $\alpha_t$ measures the importance that is assigned to $h_t$

As the iterations proceed, examples that are difficult to classify correctly receive ever-increasing influence
Weights of the base learners $\alpha_t$

- $\text{error}_{D_t}(h_t) < \frac{1}{2} \Rightarrow \alpha_t > 0$

- the smaller the error, the bigger the weight of the (weak) base learner

- the bigger the weight, the more impact on the (strong) resulting classifier

$$\text{error}_{D_t}(h_1) < \text{error}_{D_t}(h_2) \Rightarrow \alpha_1 > \alpha_2$$

- $D_{t+1} = \frac{1}{Z_t} D_t e^{-\alpha_t y_i h_t(x_i)}$

  The weights of correctly classified instances are reduced while weights of misclassified instances are increased.
Multiclass problem – generalization of the two-class case

- Assume classification task where \( Y = \{ y_1, \ldots, y_k \} \)

\[ h_t : X \to Y, \]

\[ D_{t+1}(i) = \frac{D_t(i)}{Z_t} \cdot \begin{cases} e^{-\alpha_t} & \text{if } y_i = h_t(x_i) \\ e^{\alpha_t} & \text{if } y_i \neq h_t(x_i) \end{cases} \]

\[ h_{\text{final}}(x) = \arg \max_{y \in Y} \sum_{\{t \mid h_t(x) = y\}} \alpha_t. \]
Ensembles, bagging, boosting – general principles

Simple bagging algorithm

Random Forests

Boosting with regression trees

AdaBoost

Practical use of bagging and boosting methods in R