# 10701 Machine Learning

**Boosting** 

### Fighting the bias-variance tradeoff

- Simple (a.k.a. weak) learners are good
  - e.g., naïve Bayes, logistic regression, decision stumps (or shallow decision trees)
  - Low variance, don't usually overfit
- Simple (a.k.a. weak) learners are bad
  - High bias, can't solve hard learning problems

- Can we make all weak learners always good????
  - No!!!
  - But often yes...

# Simplest approach: A "bucket of models"

### Input:

- your top T favorite learners (or tunings)
  - L<sub>1</sub>,....,L<sub>T</sub>
- A dataset D
- Learning algorithm:
  - Use 10-CV to estimate the error of L₁,....,L⊤
  - Pick the best (lowest 10-CV error) learner L\*
  - Train L\* on D and return its hypothesis h\*

### Pros and cons of a "bucket of models"

#### • Pros:

- Simple
- Will give results not much worse than the best of the "base learners"

#### Cons:

- What if there's not a single best learner?
- Other approaches:
  - Vote the hypotheses (how would you weight them?)
  - Combine them some other way?
  - How about *learning* to combine the hypotheses?

# Stacked learners: first attempt

- Input:
  - your top T favorite learners (or tunings)
    - L<sub>1</sub>,...,L<sub>T</sub>
  - A dataset D containing (x,y), ....
- Learning algorithm:
  - Train  $L_1,...,L_T$  on D to get  $h_1,...,h_T$
  - Create a new dataset D' containing (x',y'),....
    - x' is a vector of the T predictions h<sub>1</sub>(x),...,h<sub>T</sub>(x)
    - y is the label y for x
  - Train new classifier on D' to get h' --- which combines the predictions!
- To predict on a new x:
  - Construct x' as before and predict h'(x')

# Pros and cons of stacking

#### • Pros:

- Fairly simple
- Slow, but easy to parallelize

#### Cons:

- What if there's not a single best combination scheme?
- E.g.: for movie recommendation sometimes L1 is best for users with many ratings and L2 is best for users with few ratings.

# Voting (Ensemble Methods)

- Instead of learning a single (weak) classifier, learn many weak classifiers that are good at different parts of the input space
- Output class: (Weighted) vote of each classifier
  - Classifiers that are most "sure" will vote with more conviction
  - Classifiers will be most "sure" about a particular part of the space
  - On average, do better than single classifier!

#### But how do you ???

- force classifiers to learn about different parts of the input space?
- weigh the votes of different classifiers?

### Comments

- Ensembles based on blending/stacking were key approaches used in the netflix competition
  - Winning entries blended many types of classifiers
- Ensembles based on stacking are the main architecture used in Watson
  - Not all of the base classifiers/rankers are learned, however; some are hand-programmed.

### Boosting [Schapire, 1989]

- Idea: given a weak learner, run it multiple times on (reweighted) training data, then let the learned classifiers vote
- On each iteration t.
  - weight each training example by how incorrectly it was classified
  - Learn a hypothesis h<sub>t</sub>
  - A strength for this hypothesis  $\alpha_t$
- Final classifier:
  - A linear combination of the votes of the different classifiers weighted by their strength

- Practically useful
- Theoretically interesting

# Learning from weighted data

- Sometimes not all data points are equal
  - Some data points are more equal than others
- Consider a weighted dataset
  - D(i) weight of *i* th training example ( $\mathbf{x}^i, \mathbf{y}^i$ )
  - Interpretations:
    - *i* th training example counts as D(i) examples
    - If I were to "resample" data, I would get more samples of "heavier" data points
- Now, in all calculations, whenever used, *i* th training example counts as D(i) "examples"
  - e.g., MLE for Naïve Bayes, redefine Count(Y=y) to be weighted count

Given:  $(x_1, y_1), ..., (x_m, y_m)$  where  $x_i \in X, y_i \in Y = \{-1, +1\}$ Initialize  $D_1(i) = 1/m$ .

For t = 1, ..., T:

- Train weak learner using distribution  $D_t$ .
- Getweak classifier  $h_t: X \to \mathbb{R}$ .
- Choose  $\alpha_t \in \mathbb{R}$ .
- Update:

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

where  $Z_t$  is a normalization factor

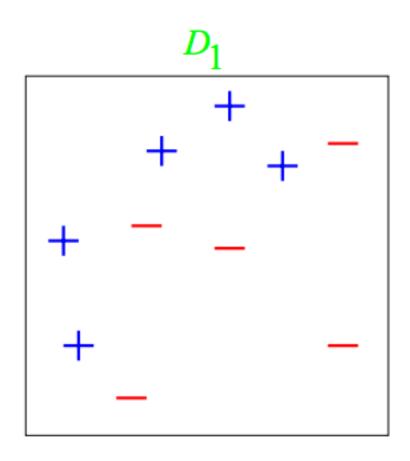
$$Z_t = \sum_{i=1}^m D_t(i) \exp(-\alpha_t y_i h_t(x_i))$$

Output the final classifier:

$$H(x) = \operatorname{sign}\left(\sum_{t=1}^{T} \alpha_t h_t(x)\right).$$

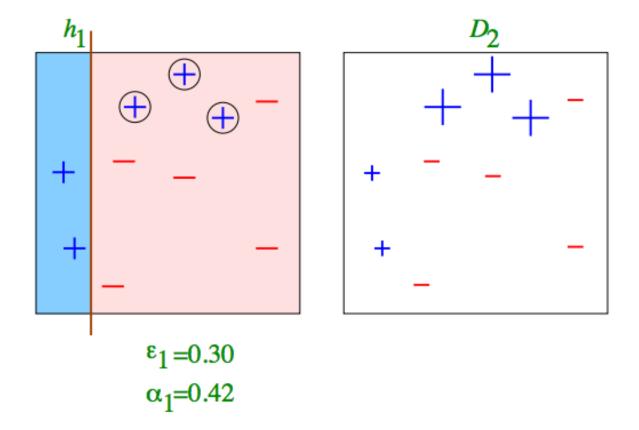
Figure 1: The boosting algorithm AdaBoost.

# Boosting: A toy example



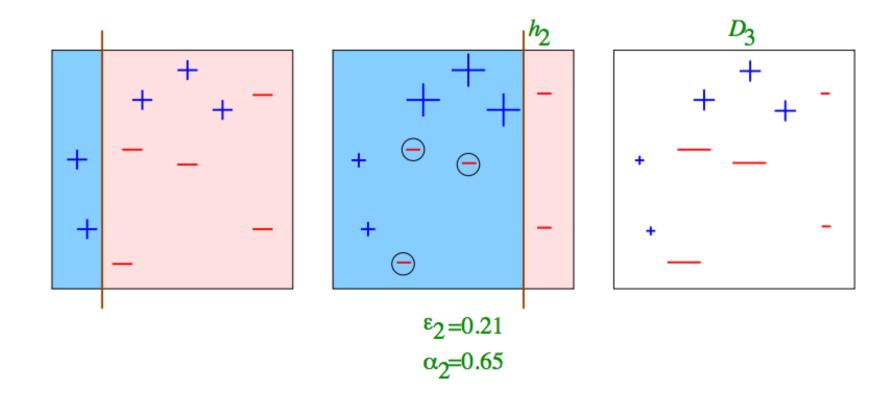
# Boosting: A toy example

### Round 1



# Boosting: A toy example

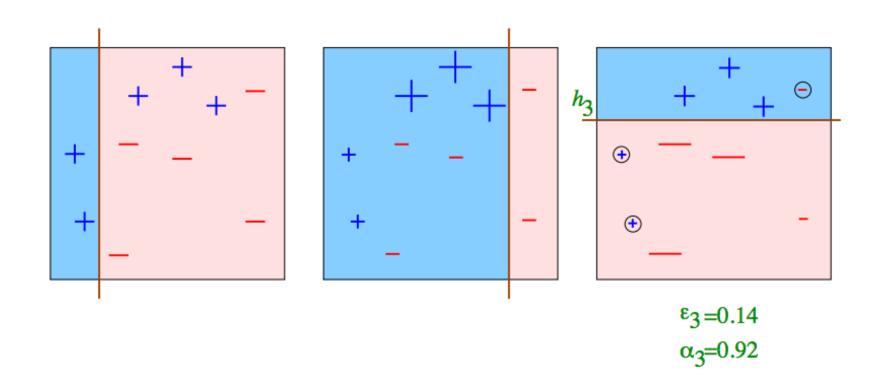
### Round 2



#### Thanks, Rob Schapire

# Boosting: A toy example

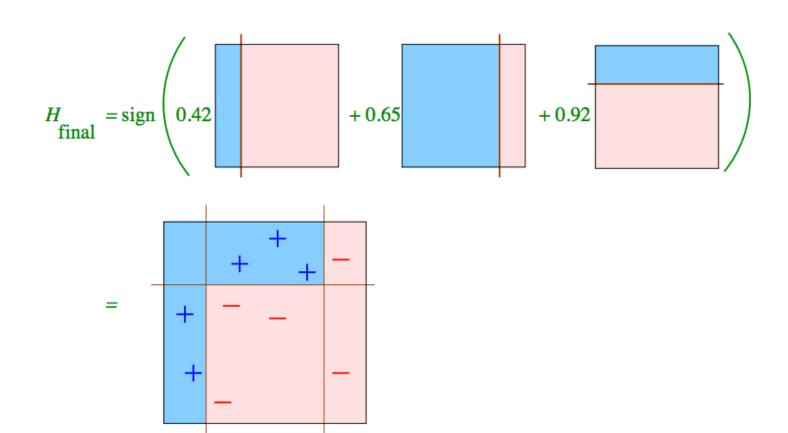
### Round 3



#### Thanks, Rob Schapire

# Boosting: A toy example

#### Final Classifier



[Schapire, 1989]

Training error of final classifier is bounded by:

$$\frac{1}{m} \sum_{i=1}^{m} \delta(H(x_i) \neq y_i) \leq \frac{1}{m} \sum_{i=1}^{m} \exp(-y_i f(x_i))$$

Where 
$$f(x) = \sum_{t} \alpha_t h_t(x)$$
;  $H(x) = sign(f(x))$ 

[Schapire, 1989]

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

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Where 
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#### If we minimize $\prod_t Z_t$ , we minimize our training error

We can tighten this bound greedily, by choosing  $\alpha_t$  and  $h_t$  on each iteration to minimize  $Z_t$ .

$$Z_t = \sum_{i=1}^m D_t(i) \exp(-\alpha_t y_i h_t(x_i))$$

[Schapire, 1989]

We can minimize this bound by choosing  $\alpha_t$  on each iteration to minimize  $Z_t$ 

$$Z_t = \sum_{i=1}^m D_t(i) \exp(-\alpha_t y_i h_t(x_i))$$

Define

$$\epsilon_t = \sum_{i=1}^m D_t(i)\delta(h_t(x_i) \neq y_i)$$

We can show that:

$$Z_{t} = (1 - \varepsilon_{t}) \exp^{-\alpha_{t}} + \varepsilon_{t} \exp^{\alpha_{t}}$$

[Schapire, 1989]

We can minimize this bound by choosing  $\alpha_t$  on each iteration to minimize  $Z_t$ 

$$Z_{t} = (1 - \varepsilon_{t}) \exp^{-\alpha_{t}} + \varepsilon_{t} \exp^{\alpha_{t}}$$

For boolean target function, this is accomplished by [Freund & Schapire '97]:

$$\alpha_t = \frac{1}{2} \ln \left( \frac{1 - \epsilon_t}{\epsilon_t} \right)$$

Where:

$$\epsilon_t = \sum_{i=1}^m D_t(i)\delta(h_t(x_i) \neq y_i)$$

Given:  $(x_1, y_1), \dots, (x_m, y_m)$  where  $x_i \in X, y_i \in Y = \{-1, +1\}$ Initialize  $D_1(i) = 1/m$ .

For t = 1, ..., T:

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- Get base classifier  $h_t: X \to \mathbb{R}$ .
- Choose  $\alpha_t \in \mathbb{R}$ . •
- Update:

$$\alpha_t = \frac{1}{2} \ln \left( \frac{1 - \epsilon_t}{\epsilon_t} \right)$$

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

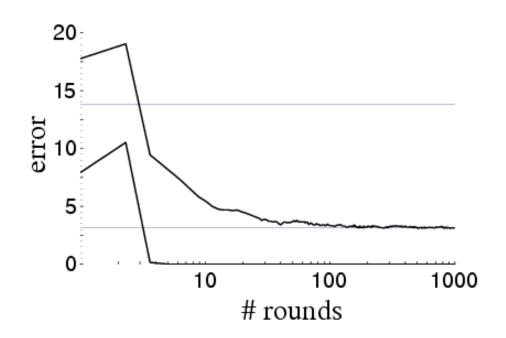
### Strong, weak classifiers

- If each classifier is (at least slightly) better than random
  - $\varepsilon_{t} < 0.5$
- With a few extra steps it can be shown that AdaBoost will achieve zero *training error* (exponentially fast):

$$\frac{1}{m} \sum_{i=1}^{m} \delta(H(x_i) \neq y_i) \leq \prod_{t=1}^{m} Z_t \leq \exp\left(-2\sum_{t=1}^{m} (1/2 - \epsilon_t)^2\right)$$

### Boosting results – Digit recognition

[Schapire, 1989]



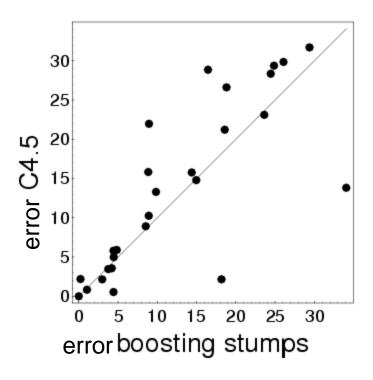
### Boosting often

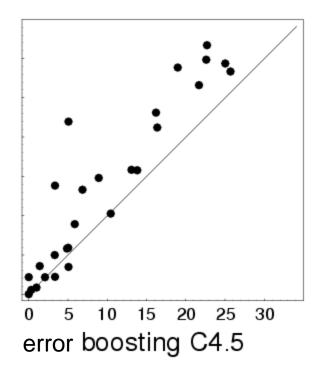
- Robust to overfitting
- Test set error decreases even after training error is zero

### **Boosting: Experimental Results**

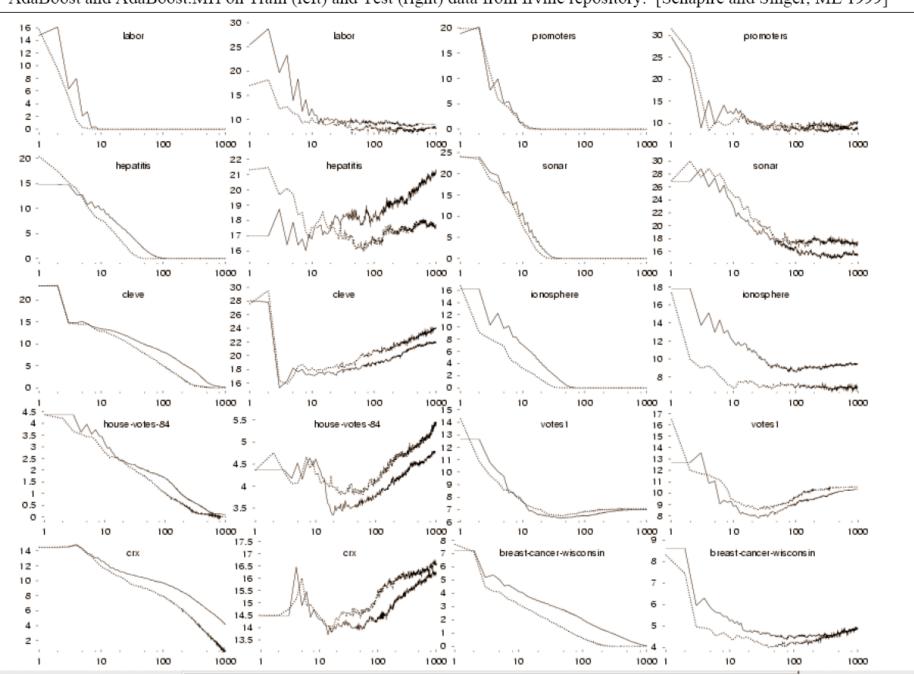
[Freund & Schapire, 1996]

Comparison of C4.5, Boosting C4.5, Boosting decision stumps (depth 1 trees), 27 benchmark datasets





AdaBoost and AdaBoost.MH on Train (left) and Test (right) data from Irvine repository. [Schapire and Singer, ML 1999]



### Random forest

- A collection of decision trees
- For each tree we select a subset of the attributes (recommended square root of |A|) and build tree using just these attributes

 An input sample is sing majority voting **TAP** GeneExpress Y2H GeneExpress Domain GeneExpress **GOProcess** HMS PCI N SynExpress **HMS-PCI** Y2H ProteinExpress GOLocalization GeneOccur GeneExpress **ProteinExpress** 

### What you need to know about Boosting

- Combine weak classifiers to obtain very strong classifier
  - Weak classifier slightly better than random on training data
  - Resulting very strong classifier can eventually provide zero training error
- AdaBoost algorithm
- Most popular application of Boosting:
  - Boosted decision stumps!
  - Very simple to implement, very effective classifier

### **Boosting and Logistic Regression**

Logistic regression assumes:

$$P(Y = 1|X) = \frac{1}{1 + \exp(f(x))}$$

And tries to maximize data likelihood:

$$P(\mathcal{D}|H) = \prod_{i=1}^{m} \frac{1}{1 + \exp(-y_i f(x_i))}$$

Equivalent to minimizing log loss

$$\sum_{i=1}^{m} \ln(1 + \exp(-y_i f(x_i)))$$

### **Boosting and Logistic Regression**

Logistic regression equivalent to minimizing log loss

$$\sum_{i=1}^{m} \ln(1 + \exp(-y_i f(x_i)))$$

Boosting minimizes similar loss function!!

$$\frac{1}{m}\sum_{i} \exp(-y_i f(x_i)) = \prod_{t} Z_t$$

Both smooth approximations of 0/1 loss!

# Logistic regression and Boosting

#### Logistic regression:

Minimize loss fn

$$\sum_{i=1}^{m} \ln(1 + \exp(-y_i f(x_i)))$$

Define

$$f(x) = \sum_{j} w_j x_j$$

where  $x_i$  predefined

#### **Boosting:**

Minimize loss fn

$$\sum_{i=1}^{m} \exp(-y_i f(x_i))$$

• Define  $f(x) = \sum_{t} \alpha_t h_t(x)$ 

where  $h_t(x_i)$  defined dynamically to fit data

(not a linear classifier)

• Weights  $\alpha_j$  learned incrementally