Machine Learning (CE 40717) Fall 2024

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Introduction

Introduction

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1 Introduction

Introduction

Condorcet's jury theorem

Ensemble learning
Ensemble methods

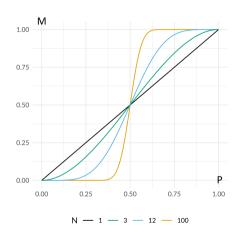
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Condorcet's jury theorem

Introduction

- N voters wish to reach a decision by **major**ity vote.
- Each voter has an independent probability **p** of voting for the correct decision.
- Majority votes for the correct decision with probability M.
- If p > 0.5 and $N \to \infty$, then $M \to 1$
 - How?



Adopted from Wikipedia



Introduction

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- Ensemble learning

Ensemble method

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Strong vs. weak Learners

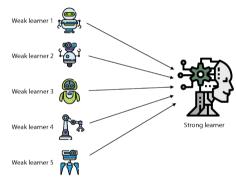
Introduction

- Strong learner: we seek to produce one classifier for which the classification error can be made arbitrarily small.
 - So far we were looking for such methods.
- Weak learner: a classifier which is just better than random guessing (for now this will be our only expectation).

Basic idea

Introduction

- Certain weak learners do well in modeling one aspect of the data, while others do well in modeling another.
- Learn several simple models and combine their outputs to produce the final decision.
- A composite prediction where the final accuracy is better than the accuracy of individual models.



Adopted from [4]

Introduction

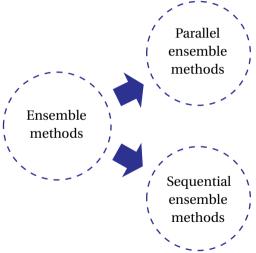
Introduction 000000000

- Ensemble methods



Ensemble Methods

Introduction



- Weak learners are generated in **parallel**.
 - Basic motivation is to use **independence** between the learners.

- Weak learners are generated consecutively.
- Basic motivation is to use **dependence** between the base learners.

What we talk about

Introduction

- Weak or simple learners
 - Low variance: they don't usually overfit
 - **High bias**: they can't learn complex functions
- Bagging (parallel): To decrease the variance
 - Random Forest
- **Boosting** (sequential): To decrease the bias (enhance their capabilities)
 - AdaBoost



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Decision tree (quick review)

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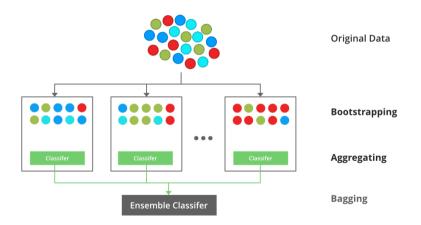


Basic idea

- Bagging = Bootstrap aggregating
- It uses bootstrap resampling to generate different training datasets from the original training dataset.
 - Samples training data uniformly at random with replacement.
- On the training datasets, it trains different weak learners.
- During testing, it **aggregates** the weak learners by uniform averaging or majority voting.
 - Works best with unstable models (high variance models). Why?



Basic idea, Cont.







Algorithm

Algorithm 1 Bagging

- 1: **Input:** *M* (required ensemble size), $D = \{(x^{(1)}, y^{(1)}), \dots, (x^{(N)}, y^{(N)})\}$ (training set)
- 2: **for** t = 1 to M **do**
- 3: Build a dataset D_t by sampling N items randomly with replacement from D \triangleright Bootstrap resampling: like rolling N-face dice N times
- 4: Train a model h_t using D_t and add it to the ensemble
- 5: end for
- 6: $H(x) = \operatorname{sign}\left(\sum_{t=1}^{M} h_t(x)\right)$
 - > Aggregate models by voting for classification or by averaging for regression

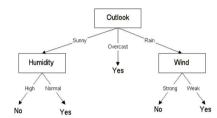


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Structure

- **Terminal nodes** (leaves) represent target variable.
- Each internal node denotes a test on an attribute.



Outlook	Temperature	Humidity	Wind	Played football(yes/no
Sunny	Hot	High	Weak	No
Sunny	Hot	High	Strong	No
Overcast	Hot	High	Weak	Yes
Rain	Mild	High	Weak	Yes
Rain	Cool	Normal	Weak	Yes
Rain	Cool	Normal	Strong	No
Overcast	Cool	Normal	Strong	Yes
Sunny	Mild	High	Weak	No
Sunny	Cool	Normal	Weak	Yes
Rain	Mild	Normal	Weak	Yes
Sunny	Mild	Normal	Strong	Yes
Overcast	Mild	High	Strong	Yes
Overcast	Hot	Normal	Weak	Yes
Rain	Mild	High	Strong	No

Adopted from Medium



Learning

- Learning an optimal decision tree is NP-Complete.
 - Instead, we use a **greedy search** based on a heuristic.
 - We can't guarantee to return the globally-optimal decision tree.
- The most common strategy for DT learning is a greedy top-down approach.
- Tree is constructed by splitting samples into subsets based on an **attribute value test** in a recursive manner.

Algorithm

Algorithm 2 Constructing DT

```
1: procedure FINDTREE(S, A)
                                                                  \triangleright Input: S (samples), A (attributes)
        if A is empty or all labels in S are the same then
            status ← leaf
 3:
            class \leftarrow most common class in S
 4:
        else
 5:
            status ← internal
 6:
            a \leftarrow \text{bestAttribute}(S, A)
                                                                              ➤ The attribute value test
            LeftNode \leftarrow FindTree(S(a = 1), A \setminus \{a\})
 8:
 9:
            RightNode \leftarrow FindTree(S(a = 0), A \setminus \{a\})
        end if
10:
11: end procedure
```

Which attribute is the best?

• Entropy measures the uncertainty in a specific distribution.

$$H(X) = -\sum_{x_i \in x} P(x_i) \log P(x_i)$$

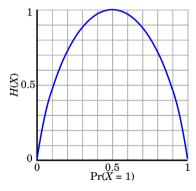
Information Gain (IG)

$$Gain(S, A) = H_S(Y) - \sum_{v \in Values(A)} \frac{|S_v|}{|S|} H_{S_V}(Y)$$

A: variable used to split samples

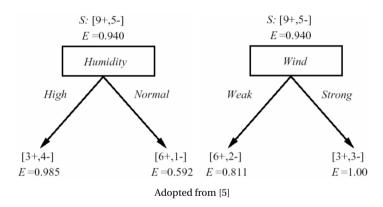
Y: target variable

S: samples



Adopted from Wikipedia

Example



$$Gain(S, Humidity) = 0.940 - (7/14)0.985 - (7/14)0.592 = 0.151$$

Gain(S, Humidity) = 0.940 - (8/14)0.811 - (6/14)1.0 = 0.48



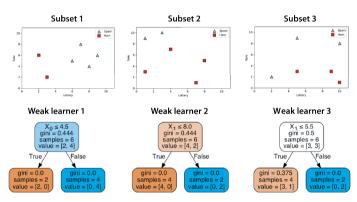
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Bagging on decision trees?

Why decision trees?

- Interpretable
- Robust to outliers
- Low bias
- High variance



Adopted from [4]

Perfect candidates

- Why are **DTs** perfect candidates for ensembles?
 - Consider averaging many (nearly) **unbiased** tree estimators.
 - Bias remains similar, but variance is reduced.
- Remember Bagging?
 - Train many trees on bootstrapped data, then average the outputs.



Algorithm

Algorithm 3 Random Forest

- 1: **Input:** *T* (number of trees), *m* (number of variables used to split each node)
- 2: **for** t = 1 to T **do**
- 3: Draw a bootstrap dataset
- 4: Learn a tree on this dataset
 - ightharpoonup Select $m{m}$ features randomly out of $m{d}$ features as candidates before splitting
- 5: end for
- 6: Output:

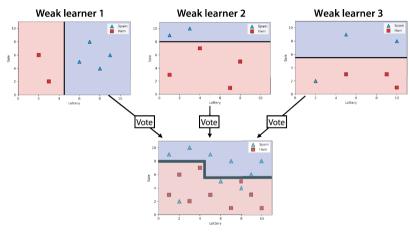
 $ightharpoonup Usually: m \le \sqrt{d}$

- 7: Regression: average of the outputs
- 8: Classification: majority voting



n Bagging **Random Forest** Boosting AdaBoost Comparison References

Example



Strong learner (random forest)

Adopted from [4]



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Problems with bagging

- Bagging created a diversity of **weak learners** by creating random datasets.
 - Examples: Decision stumps (shallow decision trees), Logistic regression, ...
- Did we have full control over the usefulness of the weak learners?
 - The **diversity** or **complementarity** of the weak learners is not controlled in any way, it is left to chance and to the instability of the models.

Basic idea

- We would expect a better performance if the weak learners also complemented each other.
 - They would have "expertise" on different subsets of the dataset.
 - So they would work better on different subsets.
- The basic idea of boosting is to generate a series of weak learners which complement each other.
 - For this, we will force each learner to focus on the mistakes of the previous learner.



Bagging Random Forest **Boosting** AdaBoost Comparison References

Basic idea, Cont.



Adopted from GeeksForGeeks



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Algorithm

- Try to combine many simple weak classifiers (in sequence) to find a single strong classifier.
 - Each component is a simple binary ±1 classifier
 - Voted combinations of component classifiers

$$H_m(\mathbf{x}) = \alpha_1 h(\mathbf{x}; \boldsymbol{\theta}_1) + \dots + \alpha_m h(\mathbf{x}; \boldsymbol{\theta}_m)$$

• To simplify notations: $h(x; \theta_i) = h_i(x)$

$$H_m(\mathbf{x}) = \alpha_1 h_1(\mathbf{x}) + \dots + \alpha_m h_m(\mathbf{x})$$

• **Prediction**: $\hat{y} = \text{sign}(H_m(x))$

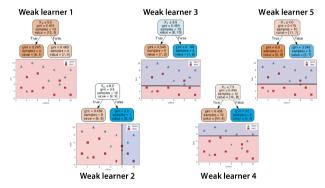


Candidate for $h_i(x)$

Decision stumps

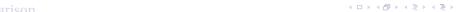
 Each classifier is based on a single feature of x(e.g., x_k):

$$h(\mathbf{x}; \boldsymbol{\theta}) = \operatorname{sign}(w_1 \mathbf{x}_k - w_0)$$
$$\boldsymbol{\theta} = \{k, w_1, w_0\}$$



Adopted from [4]

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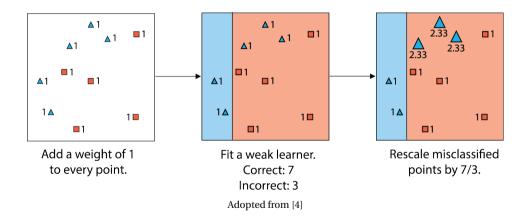


Basic idea

- Sequential production of classifiers
 - Iteratively add the classifier whose addition will be most helpful.
- Represent the important of each sample by assigning weights to them.
 - Correct classification \implies smaller weights
 - Misclassified samples ⇒ larger weights
- Each classifier is **dependent** on the previous ones.
 - Focuses on the **previous ones' error**.

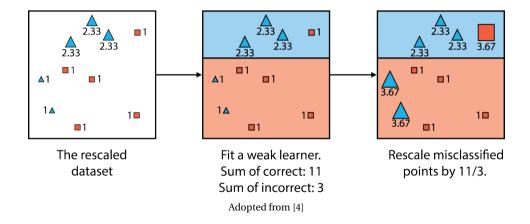


Example



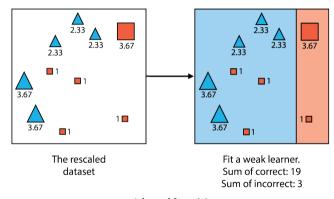


Example, Cont.





Example, Cont.

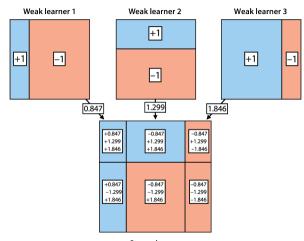


Adopted from [4]



n Bagging Random Forest Boosting **AdaBoost** Comparison References

Example, Cont.

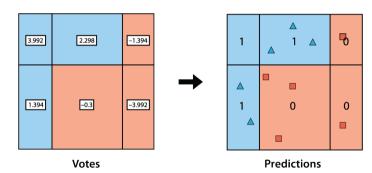


Strong learner

Adopted from [4]



Example, Cont.



Adopted from [4]

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Algorithm

Algorithm 4 AdaBoost

1: Initialize data weight $w_1^{(i)} = \frac{1}{N}$ for all N samples

2: **for**
$$m = 1$$
 to M **do**

3:
$$J_m = \sum_{i=1}^N w_m^{(i)} \times I(y^{(i)} \neq h_m(\mathbf{x}^{(i)}))$$

 \triangleright Find $h_m(\mathbf{x})$ by minimizing the weighted error

3:
$$J_{m} = \sum_{i=1}^{N} w_{m}^{(i)} \times I(y^{(i)} \neq h_{m}(\mathbf{x}^{(i)}))$$
4:
$$\epsilon_{m} = \frac{\sum_{i=1}^{N} w_{m}^{(i)} \times I(y^{(i)} \neq h_{m}(\mathbf{x}^{(i)}))}{\sum_{i=1}^{N} w_{m}^{(i)}}$$

 \triangleright Find the weighted error of $h_m(x)$

5:
$$\alpha_{m} = \ln \left(\frac{1 - \epsilon_{m}}{\epsilon_{m}} \right)^{-1}$$
6:
$$w_{m+1}^{(i)} = w_{m}^{(i)} e^{\alpha_{m} I(y^{(i)} \neq h_{m}(\boldsymbol{x}^{(i)}))}$$

> Assign votes based on the error

> Update normalized weights

7: end for

8: Combined classifier: $\hat{y} = \text{sign}(H_M(\mathbf{x})), H_M(\mathbf{x}) = \sum_{m=1}^{M} \alpha_m h_m(\mathbf{x})$



Notations & conditions

- $w_m^{(i)}$: weighting coefficient of data point *i* in iteration *m*
- α_m : weighting coefficient of *m*-th base classifier in the final ensemble
 - ϵ_m : weighted error rate of *m*-th classifier

- Only when $h_m(x)$ with $\epsilon_m < 0.5$ (better than chance) is found, boosting continues.
 - Condorcet's jury theorem?

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Loss function

- We need a loss function for the combination
 - To Determine the new component, $h(x;\theta)$
 - And how many votes it should receive, α

$$H_m(\mathbf{x}) = \alpha_1 h_1(\mathbf{x}) + \dots + \alpha_m h_m(\mathbf{x})$$

- Many options for the loss function
 - AdaBoost is equivalent to using the following **exponential loss**

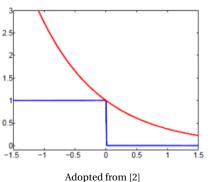
$$\operatorname{Loss}(y,\hat{y}) = e^{-y \times H_m(x)}$$

$$\hat{y} = \text{sign}(H_m(\mathbf{x}))$$



Why the exponential loss?

- Differentiable approximation (bound) of the 0/1 loss
 - Easy to optimize
 - Optimizing an upper bound on classification error.



Calculations

• Consider adding the *m*-th component:

$$H_m(\mathbf{x}) = \frac{1}{2} [\alpha_1 h_1(\mathbf{x}) + \dots, + \alpha_m h_m(\mathbf{x})]$$

For a cleaner form later

$$E = \sum_{i=1}^{N} e^{-y^{(i)} H_{m}(\mathbf{x}^{(i)})} = \sum_{i=1}^{N} e^{-y^{(i)} [H_{m-1}(\mathbf{x}^{(i)}) + \frac{1}{2} \alpha_{m} h_{m}(\mathbf{x}^{(i)})]}$$

$$= \sum_{i=1}^{N} e^{-y^{(i)} H_{m-1}(\mathbf{x}^{(i)})} e^{-\frac{1}{2} \alpha_{m} y^{(i)} h_{m}(\mathbf{x}^{(i)})} = \sum_{i=1}^{N} \underbrace{w_{m}^{(i)}}_{e^{-y^{(i)} H_{m-1}(\mathbf{x}^{(i)})}} e^{-\frac{1}{2} \alpha_{m} y^{(i)} h_{m}(\mathbf{x}^{(i)})}$$

Suppose it is fixed at stage m

Should be optimized at stage m by seeking $h_m(x)$ and α_m

Weighted exponential loss

$$E = \sum_{i=1}^{N} w_m^{(i)} e^{-\frac{1}{2}\alpha_m y^{(i)} h_m(\mathbf{x}^{(i)})}$$

- Sequentially adds a new component trained on reweighted training samples.
- $w_m^{(i)}$: history of classification of $x^{(i)}$ by H_{m-1}
 - Loss weighted towards mistakes
- Iteration *m* optimization:
 - Choose the new component, $h_m = h(\mathbf{x}; \boldsymbol{\theta}_m)$
 - And the vote that optimizes the weighted exponential loss, α_m .



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Minimizing loss: finding h_m

$$E = \sum_{i=1}^{N} w_{m}^{(i)} e^{-\frac{1}{2}\alpha_{m}y^{(i)}h_{m}(\mathbf{x}^{(i)})}$$

$$= e^{\frac{-\alpha_{m}}{2}} \left(\sum_{y^{(i)} = h_{m}(\mathbf{x}^{(i)})} w_{m}^{(i)} \right) + e^{\frac{\alpha_{m}}{2}} \left(\sum_{y^{(i)} \neq h_{m}(\mathbf{x}^{(i)})} w_{m}^{(i)} \right)$$

$$= (e^{\frac{\alpha_{m}}{2}} - e^{\frac{-\alpha_{m}}{2}}) \left(\sum_{y^{(i)} \neq h_{m}(\mathbf{x}^{(i)})} w_{m}^{(i)} \right) + e^{\frac{-\alpha_{m}}{2}} \left(\sum_{i=1}^{N} w_{m}^{(i)} \right)$$

$$J_{m} = \sum_{i=1}^{N} w_{m}^{(i)} \times I \left(y^{(i)} \neq h_{m}(\mathbf{x}^{(i)}) \right)$$

Find $h_m(x)$ that minimizes J_m

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Minimizing loss: finding α_m

$$\begin{split} \frac{\partial E}{\partial \alpha_m} &= 0 \\ \implies \frac{1}{2} (e^{\frac{\alpha_m}{2}} + e^{\frac{-\alpha_m}{2}}) \left(\sum_{y^{(i)} \neq h_m(\mathbf{x}^{(i)})} w_m^{(i)} \right) - \frac{1}{2} e^{\frac{-\alpha_m}{2}} \left(\sum_{i=1}^N w_m^{(i)} \right) = 0 \\ \implies \frac{e^{\frac{-\alpha_m}{2}}}{(e^{\frac{\alpha_m}{2}} + e^{\frac{-\alpha_m}{2}})} &= \frac{\sum_{y^{(i)} \neq h_m(\mathbf{x}^{(i)})} w_m^{(i)}}{\sum_{i=1}^N w_m^{(i)}} \end{split}$$

$$\epsilon_m = \frac{\sum_{i=1}^{N} w_m^{(i)} I\left(y^{(i)} \neq h_m(\boldsymbol{x}^{(i)})\right)}{\sum_{i=1}^{N} w_m^{(i)}}, \quad \alpha_m = \ln\left(\frac{1 - \epsilon_m}{\epsilon_m}\right)$$



Updating weights

• Updating weights in AdaBoost algorithm:

$$w_i^{m+1} = w_i^m e^{-\frac{1}{2}\alpha_m y^{(i)} h_m(\mathbf{x}^{(i)})}$$

$$\xrightarrow{y^{(i)}h_m(\mathbf{x}^{(i)})=1-2I(y^{(i)}\neq h_m(\mathbf{x}^{(i)}))} w_i^{m+1} = w_i^m e^{-\frac{1}{2}\alpha_m} e^{\alpha_m I(y^{(i)}\neq h_m(\mathbf{x}^{(i)}))}$$

Independent of i and can be ignored

$$\implies w_i^{m+1} = w_i^m e^{\alpha_m I\left(y^{(i)} \neq h_m(\boldsymbol{x}^{(i)})\right)}$$

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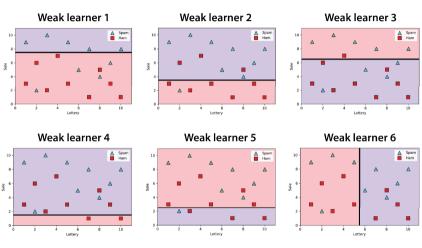
Algorithm 5 AdaBoost Summary

- 1: **for** i = 1 to N **do**
- 2: Initialize the data weight $w_1^{(i)} = \frac{1}{N}$
- 3: end for
- 4: **for** m = 1 to M **do**
- 5: Find a classifier $h_m(x)$ by minimzing the weighted error function
- 6: Find the normalized weighted error of $h_m(\mathbf{x})$ as ϵ_m
- 7: Compute the new component weight as α_m
- 8: Update example weights for the next iteration $w_{m+1}^{(i)}$
- 9: end for
- 10: Combined classifier $\hat{y} = \text{sign}(H_M(\mathbf{x}))$ where $H_M(\mathbf{x}) = \sum_{m=1}^{M} \alpha_m h_m(\mathbf{x})$



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Example

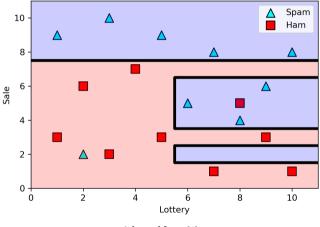






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Example, Cont.



Adopted from [4]



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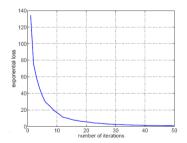


Exponential loss properties

• In each boosting iteration, assuming we can find $h(x; \hat{\theta}_m)$ whose weighted error is better than chance.

$$H_m(x) = \frac{1}{2} [\hat{\alpha}_1 h(\boldsymbol{x}; \hat{\boldsymbol{\theta}}_1) + \dots + \hat{\alpha}_m h(\boldsymbol{x}; \hat{\boldsymbol{\theta}}_m)]$$

• Thus, **lower exponential loss** over training data is guaranteed.

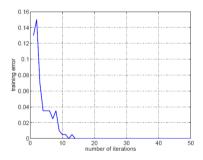


Adopted from [6]



Training error properties

• Boosting iterations typically **decrease** the **training error** of $H_m(\mathbf{x})$ over training examples.



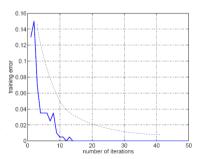
Adopted from [6]



Training error properties, Cont.

• Training error has to go down exponentially fast if the weighted error of each h_m is strictly better than chance (i.e., $\epsilon_m < 0.5$)

$$E_{\text{train}}(H_M) \le \prod_{m=1}^{M} 2\sqrt{\epsilon_m(1-\epsilon_m)}$$



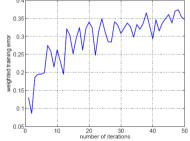
Adopted from [6]



Weighted error properties

• Weighted error of each new component classifier tends to increase as a function of boosting iterations.

$$\epsilon_{m} = \frac{\sum_{i=1}^{N} w_{m}^{(i)} I(y^{(i)} \neq h_{m}(x^{(i)}))}{\sum_{i=1}^{N} w_{m}^{(i)}}$$

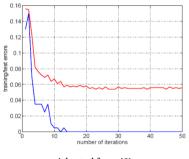


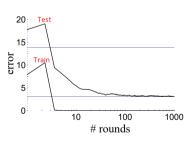
Adopted from [6]



Test error properties

- **Test error** can still **decrease** after training error is flat (even zero).
- But, is it robust to overfitting?
 - May easily overfit in the presence of labeling noise or overlap of classes.





Adopted from [6]

Adopted from [3]



Typical behavior

- Exponential loss goes strictly down.
- Training error of H goes down.
- Weighted error ϵ_m goes $\mathbf{up} \implies \text{votes } \alpha_m$ go \mathbf{down} .
- **Test error decreases** even after a flat training error.



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Bagging vs. Boosting

Bagging

- Uses bootstrap sampling to construct several training sets from the training set and then aggregates the learners trained (in parallel) on these datasets.
- **Reduces the variance** of high variance learners (e.g., decision trees)
- Boosting
 - Combine many weak classifiers in sequence to find a single strong classifier (by putting emphasize on the misclassified samples in each iteration).
 - There is empirical evidence that it reduces both the bias and the variance.
 - It seems that bias is mostly reduced in earlier iterations, while variance in later ones.



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- **6** Comparison
- References



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