

INTRO to DATA SCIENCE

LECTURE 13: ENSEMBLE TECHNIQUES

LAST TIME:

- DIMENSIONALITY REDUCTION**
- PRINCIPAL COMPONENT ANALYSIS**

QUESTIONS?

I. ENSEMBLE TECHNIQUES

II. PROBLEMS IN CLASSIFICATION

III. BAGGING

IV. BOOSTING

V. RANDOM FORESTS

EXERCISE:

VI. RANDOM FORESTS

I. ENSEMBLE TECHNIQUES

Q: What are ensemble techniques?

Q: What are ensemble techniques?

*A: Methods of improving classification accuracy by aggregating predictions over several **base classifiers**.*

Q: What are ensemble techniques?

*A: Methods of improving classification accuracy by aggregating predictions over several **base classifiers**.*

Ensembles are often much more accurate than the base classifiers that compose them.

Q: What are ensemble techniques?

*A: Methods of improving classification accuracy by aggregating predictions over several **base classifiers**.*

Ensembles are often much more accurate than the base classifiers that compose them.

NOTE

Base classifiers and ensemble classifiers are sometimes called *weak learners* and *strong learners*.

In order for an ensemble classifier to outperform a single base classifier, the following conditions must be met:

In order for an ensemble classifier to outperform a single base classifier, the following conditions must be met:

- 1) the bc's must be **accurate**: they must outperform random guessing*

In order for an ensemble classifier to outperform a single base classifier, the following conditions must be met:

- 1) the bc's must be **accurate**: they must outperform random guessing*
- 2) the bc's must be **diverse**: their misclassifications must occur on different training examples*

In order for an ensemble classifier to outperform a single base classifier, the following conditions must be met:

*1) the bc's must be **accurate**: low bias*

*2) the bc's must be **diverse**: uncorrelated*

In order for an ensemble classifier to outperform a single base classifier, the following conditions must be met:

*1) the bc's must be **accurate**: low bias*

*2) the bc's must be **diverse**: uncorrelated*

NOTE

Ideally, we would also like the base classifiers to be *unstable* to variations in the training set.

In other words, *high variance*.

II. PROBLEMS IN CLASSIFICATION

In any supervised learning task, our goal is to make predictions of the true classification function f by learning the classifier h .

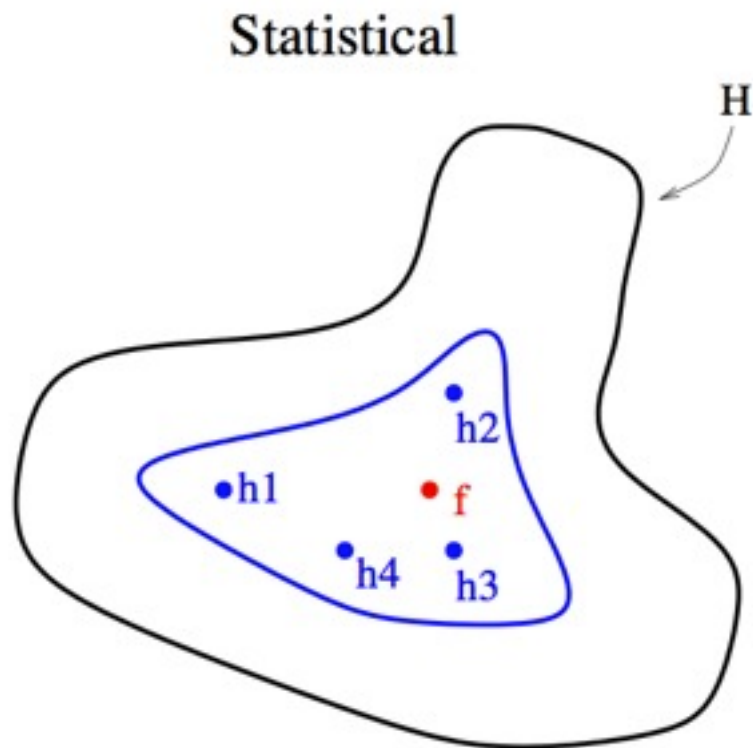
In any supervised learning task, our goal is to make predictions of the true classification function f by learning the classifier h .

There are three main problems that can prevent this:

- statistical problem*
- computational problem*
- representational problem*

If the amount of training data available is small, the base classifier will have difficulty converging to h .

An ensemble classifier can mitigate this problem by “averaging out” base classifier predictions to improve convergence.



NOTE

The true function f is best approximated as an average of the base classifiers.

source: <http://www.cs.iastate.edu/~jtian/cs573/Papers/Dietterich-ensemble-00.pdf>

Even with sufficient training data, it may still be computationally difficult to find the best classifier h .

For example, if our base classifier is a decision tree, an exhaustive search of the hypothesis space of all possible classifiers is extremely complex (NP-complete).

Even with sufficient training data, it may still be computationally difficult to find the best classifier h .

For example, if our base classifier is a decision tree, an exhaustive search of the hypothesis space of all possible classifiers is extremely complex (NP-complete).

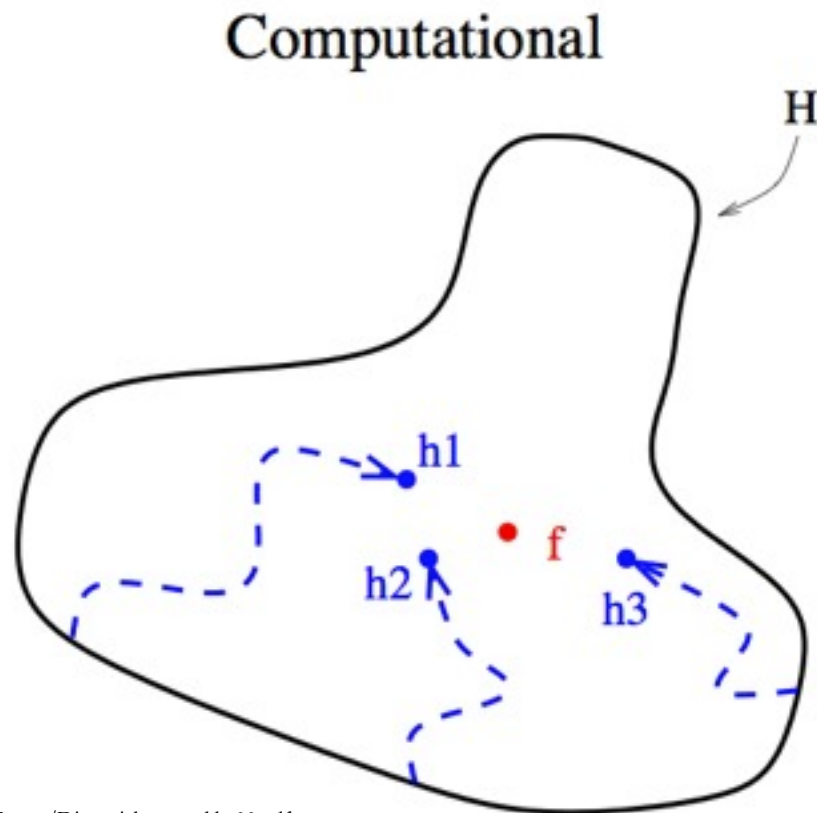
NOTE

Recall that this is why we used a *heuristic algorithm* (greedy search).

Even with sufficient training data, it may still be computationally difficult to find the best classifier h .

For example, if our base classifier is a decision tree, an exhaustive search of the hypothesis space of all possible classifiers is extremely complex (NP-complete).

An ensemble composed of several BC's with different starting points can provide a better approximation to f than any individual BC.

**NOTE**

The true function f is often best approximated by using several starting points to explore the hypothesis space.

source: <http://www.cs.iastate.edu/~jtian/cs573/Papers/Dietterich-ensemble-00.pdf>

Sometimes f cannot be expressed in terms of our hypothesis at all.

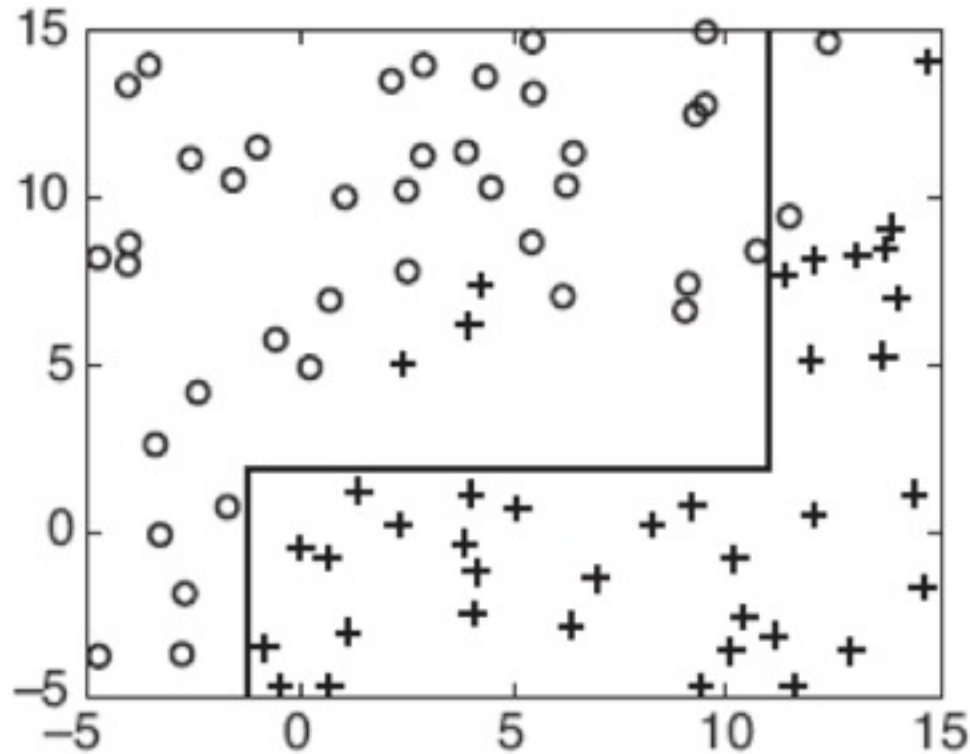
Sometimes f cannot be expressed in terms of our hypothesis at all.

To illustrate this, suppose we use a decision tree as our base classifier.

Sometimes f cannot be expressed in terms of our hypothesis at all.

To illustrate this, suppose we use a decision tree as our base classifier.

A decision tree works by forming a rectilinear partition of the feature space.



NOTE

What is a *rectilinear* decision boundary?

One whose segments are *orthogonal* to the x & y axes.

But what if f is a diagonal line?

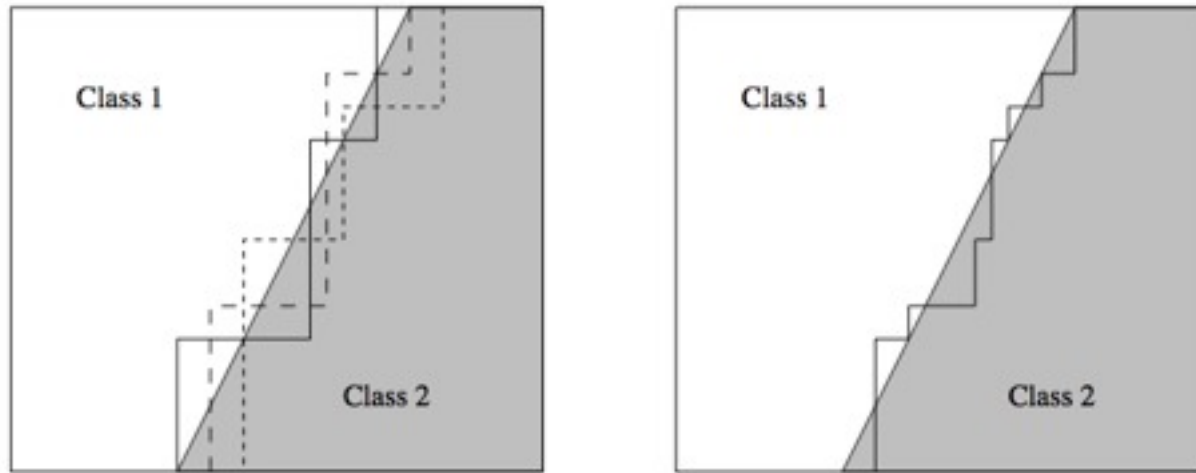
But what if f is a diagonal line?

*Then it cannot be represented by finitely many rectilinear segments,
and therefore the true decision boundary cannot be obtained by a
decision tree classifier.*

But what if f is a diagonal line?

Then it cannot be represented by finitely many rectilinear segments, and therefore the true decision boundary cannot be obtained by a decision tree classifier.

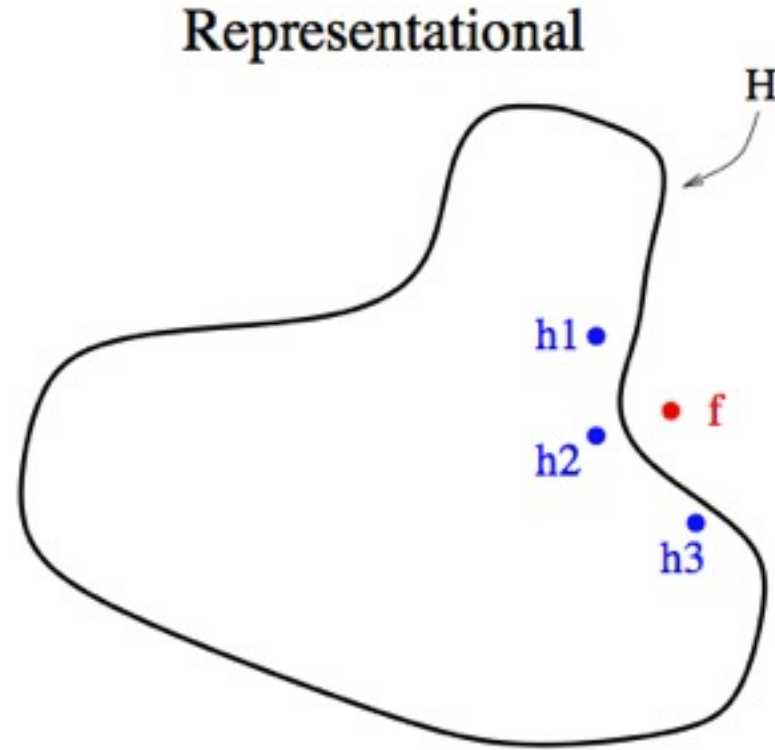
However, it may be still be possible to approximate f or even to expand the space of representable functions using ensemble methods.



NOTE

An ensemble of decision trees can approximate a diagonal decision boundary.

Fig. 4. The left figure shows the true diagonal decision boundary and three staircase approximations to it (of the kind that are created by decision tree algorithms). The right figure shows the voted decision boundary, which is a much better approximation to the diagonal boundary.



NOTE

Ensemble classifiers can be effective even if the true decision boundary lies outside the hypothesis space.

Q: How do you create an ensemble classifier?

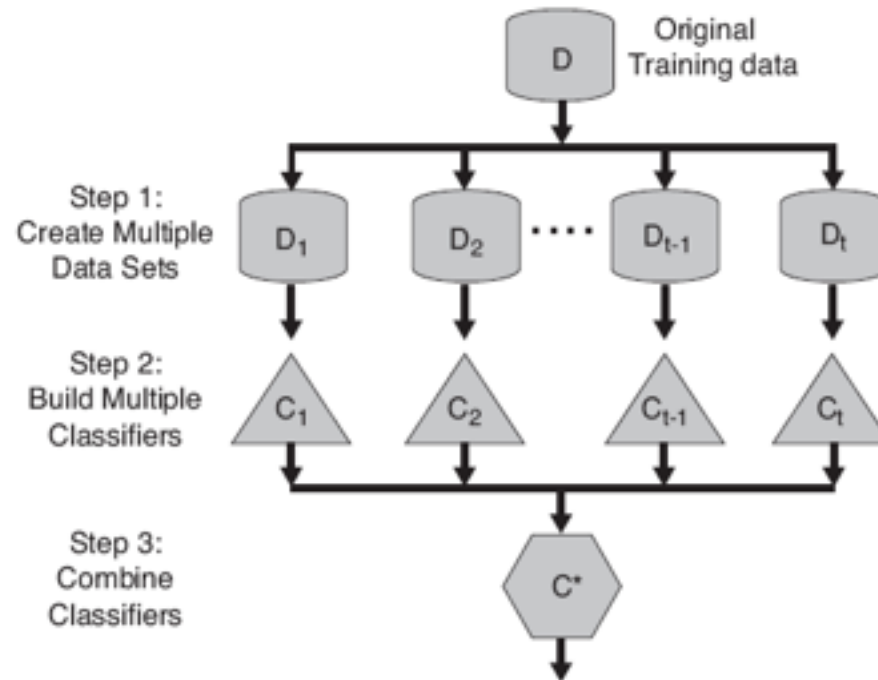


Figure 5.31. A logical view of the ensemble learning method.

Q: How do you generate several base classifiers?

Q: How do you generate several base classifiers?

A: There are several ways to do this:

- manipulating the training set*
- manipulating the output labels*
- manipulating the learning algorithm itself*

We will talk about a few examples of each of these.

III. BAGGING

Bagging (*bootstrap aggregating*) is a method that involves manipulating the training set by **resampling**.

Bagging (*bootstrap aggregating*) is a method that involves manipulating the training set by **resampling**.

We learn k base classifiers on k different samples of training data.

*These samples are independently created by resampling the training data using uniform weights (eg, a uniform **sampling distribution**).*

Bagging (*bootstrap aggregating*) is a method that involves manipulating the training set by **resampling**.

We learn k base classifiers on k different samples of training data.

These samples are independently created by resampling the data using uniform weights (eg, a uniform sampling distribution).

NOTE

Each training sample is the same size as the original training set.

Bagging (*bootstrap aggregating*) is a method that involves manipulating the training set by **resampling**.

We learn k base classifiers on k different samples of training data.

These samples are independently created by resampling the data using uniform weights (eg, a uniform sampling distribution).

NOTE

NOTE

Resampling means that some training records may appear in a sample more than once, or even not at all.

Bagging (*bootstrap aggregating*) is a method that involves manipulating the training set by **resampling**.

We learn k base classifiers on k different samples of training data.

*These samples are independently created by resampling the training data using uniform weights (eg, a uniform **sampling distribution**).*

The final prediction is made by taking a majority vote across bc's.

Bagging reduces the variance in our generalization error by aggregating multiple base classifiers together (provided they satisfy our earlier requirements).

Bagging reduces the variance in our generalization error by aggregating multiple base classifiers together (provided they satisfy our earlier requirements).

If the base classifier is stable, then the ensemble error is primarily due to bc bias, and bagging may not be effective.

Bagging reduces the variance in our generalization error by aggregating multiple base classifiers together (provided they satisfy our earlier requirements).

If the base classifier is stable, then the ensemble error is primarily due to bc bias, and bagging may not be effective.

Since each sample of training data is equally likely, bagging is not very susceptible to overfitting with noisy data.

IV. BOOSTING

Boosting is an iterative procedure that adaptively changes the sampling distribution of training records at each iteration.

Boosting is an iterative procedure that adaptively changes the sampling distribution of training records at each iteration.

The first iteration uses uniform weights (like bagging). In subsequent iterations, the weights are adjusted to emphasize records that were misclassified in previous iterations.

Boosting is an iterative procedure that adaptively changes the sampling distribution of training records at each iteration.

The first iteration uses uniform weights (like bagging). In subsequent iterations, the weights are adjusted to emphasize records that were misclassified in previous iterations.

The final prediction is constructed by a weighted vote (where the weights for a bc depends on its training error).

Boosting is an iterative procedure that adaptively changes the sampling distribution of training records at each iteration.

The first iteration uses uniform weights (like bagging). In subsequent iterations, the weights are adjusted to emphasize records that were misclassified in previous iterations.

The final prediction is constructed by a weighted vote (where the weights for a bc depends on its training error).

NOTE

The bc's focus more and more closely on records that are difficult to classify as the sequence of iterations progresses.

Thus the bc's are faced with progressively more difficult learning problems.

Like in bagging, sampling is done with replacement, and as a result some records may not appear in a given training sample.

Like in bagging, sampling is done with replacement, and as a result some records may not appear in a given training sample.

These omitted records will likely be misclassified, and given greater weight in subsequent iterations once the sampling distribution is updated.

Like in bagging, sampling is done with replacement, and as a result some records may not appear in a given training sample.

These omitted records will likely be misclassified, and given greater weight in subsequent iterations once the sampling distribution is updated.

So even if a record is left out at one stage, it will be emphasized later.

Updating the sampling distribution and forming an ensemble prediction leads to a nonlinear combination of the base classifiers.

Updating the sampling distribution and forming an ensemble prediction leads to a nonlinear combination of the base classifiers.

By explicitly trying to optimize the weighted ensemble vote, boosting attacks the representation problem head-on.

INTRO TO DATA SCIENCE

V. RANDOM FORESTS

Tuesday, October 15, 2013

A random forest is an ensemble of decision trees where each base classifier is grown using a random effect.

A random forest is an ensemble of decision trees where each base classifier is grown using a random effect.

One way to do this is to randomly choose one of the top k features to split each node.

A random forest is an ensemble of decision trees where each base classifier is grown using a random effect.

One way to do this is to randomly choose one of the top k features to split each node.

For a small number of features, we can also create linear combinations of features and select splits from the enhanced feature set (Forest-RC).

A random forest is an ensemble of decision trees where each base classifier is grown using a random effect.

One way to do this is to randomly choose one of the top k features to split each node.

For a small number of features, we can also create linear combinations of features and select splits from the enhanced feature set (Forest-RC).

Or, we can select splitting features completely at random (Forest-RI).

Random forests are about as accurate as AdaBoost, more robust to noise, and can also have better runtime than other ensemble methods (since the feature space is reduced in some cases).

Wisdom of Crowds

IBM's WATSON

Nate Silver's election models

Kaggle competitions

INTRO TO DATA SCIENCE

EX: ENSEMBLE METHODS IN SCIKIT-LEARN

Tuesday, October 15, 2013