

1 Module 20: Ensemble Systems

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- Bagging: Random trees
- Boosting: AdaBoost, Gradient-boosted trees

Individual models must be independent for the ensemble system to work.

$$\text{Variance[ensemble]} = \frac{\text{Variance[individual]}}{N}$$

1.1 Aggregating predictors

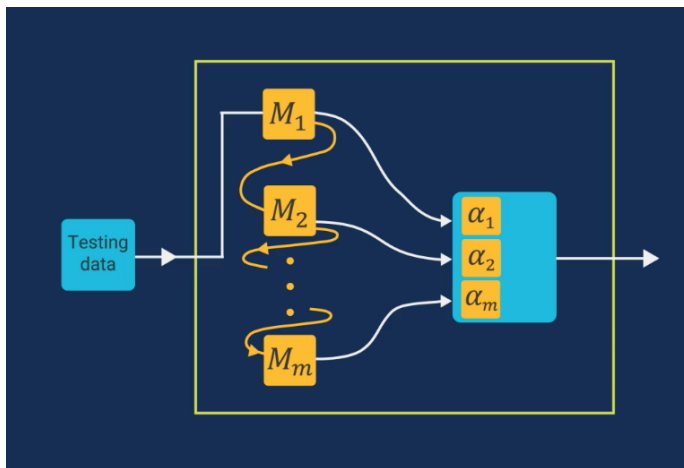
Metamodel = ensemble system

→ assign weights $\alpha_1, \dots, \alpha_m$ to models and fit their influences on final decision.

We find that using the same model in many different variations performs better than multiple different types of classifiers.

- Aggregation for classification
 - Hard voting: Majority vote
 - Soft voting: average the predicted class distributions and choose maximum
- """ for regression
 - Simple average

Boosting: predictors are trained sequentially

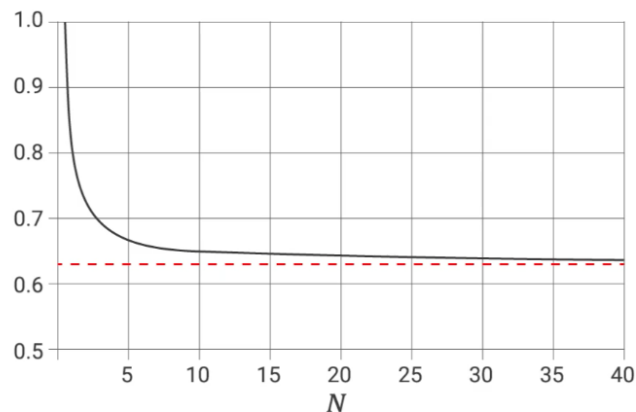


1.2 Bootstrapping & Bagging

High variance \rightarrow too myopically focused on training data.

Bootstrapping (decorrelating models' data) Assume that training data samples from a population \rightarrow sample with replacement (a sampled data point isn't removed from the dataset).

```
indices = np.random.choice( range(ntrain), ntrain )
X_bs = X_train[ind,:]  
y_bs = y_train[ind,:]
```



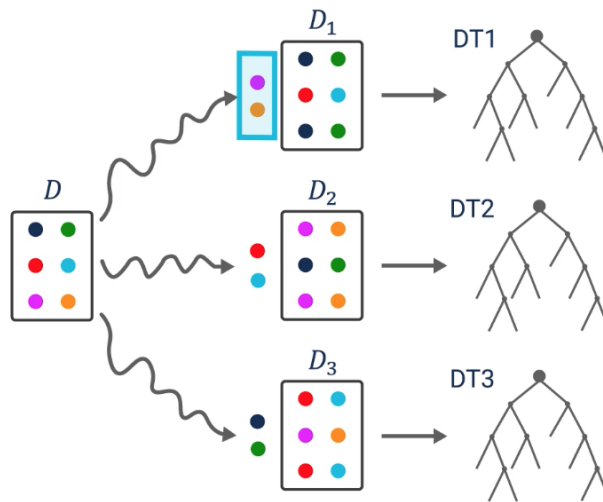
$$\lim_{N \rightarrow \infty} \left(1 - \left(1 - \frac{1}{N} \right)^N \right) = 1 - \frac{1}{e} \cong 0.632$$

We can expect in the lim of \ll training data, we expect for each sample set to be 63% new data and 37% repeated data.

Bagging: "bootstrap aggregation" Performed by our choice of aggregation technique.

```
y_ensemble = scipy.stats.mode(y_pred).mode[0]
```

Out-of-bag evaluation: withhold a subset of data per trained ensemble classifier, which is used as the testing set.



```
model = BaggingClassifier(
    DecisionTreeClassifier(),
    n_estimators = 500, # number of models in ensemble
    oob_score = True # Out of bag score
)
model.fit(X,y)
model.oob_score_
```

1.3 Random Forests

Ordinarily, DTrees choose to split on features that produce the greatest change in entropy. We can simply require that the `max_features > 1` such that we search along otherwise neglected branches. → increase diversity of ensemble

```
from sklearn.ensemble import RandomForestClassifier
```

```

model = RandomForestClassifier( n_estimators = 30, max_features = 10, oob_score =
    ↪ True )
model.fit( X,y )
model.oob_score_

```

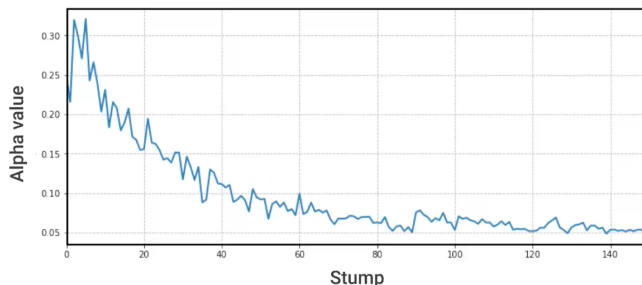
1.4 Boosting

Reduces bias of a set of weak learners (without increasing variance as much) → Combine a large number of weak models in a clever way to create a strong model.

Typically employ very shallow (one-node) decision trees (decision stumps) i.e. cut the dataset with a single slice.

Adaboost algorithm

- Initialize, $s=0$
 - for all samples $i = 1, \dots, N$ and stumps s : $w_s^i = \frac{1}{N}$
 - == the effort a stump should put into correctly classifying i
 - to begin with, all points are weighted equally
- Loop as log derived:
 - Create a stump to correctly classify samples for at least half of the total weight $S_s \cdot \text{fit}(x_1 y_1, W_s)$; if < 0.5 , just flip the inequality
 - misclassification score $\epsilon_s = \sum_{\text{misclassified}} W_s^i < 0.5$ as the system is at least better than random chance
 - influence coefficient $\alpha_s = \frac{1}{2} \log \frac{1-\epsilon_s}{\epsilon_s}$ use misclassification score to determine influence of particular stump on the ensemble
- Allow influence coefficient to update weights
 - $w_{s+1}^i = \begin{cases} w_s^i e^{\alpha_s} & i \text{ was misclassified} \\ w_s^i e^{-\alpha_s} & \text{otherwise} \end{cases}$



```

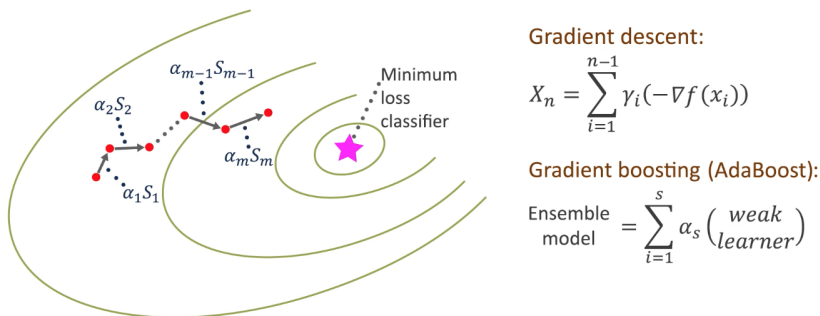
from sklearn.ensemble import AdaBoostClassifier
from sklearn.tree import DecisionTreeClassifier

model = AdaBoostClassifier(DecisionTreeClassifier(max_depth=1))
model.fit(X,y)

```

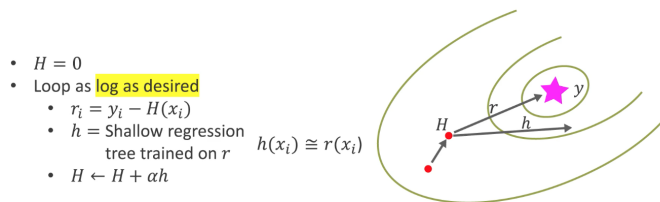
Algorithm is based on **gradient boosting**. Boosting algorithms are not easily overfitted.

GRADIENT BOOSTING



Gradient Boosting Trees Use trees for base model BUT: use squared loss as loss fcn.

- $H = 0$
- Loop as log
 - $r_i = y_i - H(X_i)$
 - h = Shallow regression tree (depth ≈ 4) trained on $r \rightarrow h(x_i) \approx r(x_i)$
 - $H \leftarrow H + \alpha h$



```

from sklearn.ensemble import GradientBoostingRegressor
GradientBoostingRegressor(n_estimators = 1000, max_depth = 1)

```