

Random Forest

Random Forest is an ensemble learning method that combines many decision trees. The math behind it comes from bootstrap, ~~features~~ sampling, ~~features~~ randomness, and majority voting. These ideas reduce variance and create a stronger, more stable model.

1. Bootstrap Sampling (Bagging)

Each tree is trained on a random sample of the dataset created by sampling with replacement.

If original dataset is:

$$D = \{(x_1, y_1), (x_2, y_2), \dots, (x_N, y_N)\}$$

A bootstrap sample D_b is generated by:

$$D_b = \{(x_{i_1}, y_{i_1}), \dots, (x_{i_m}, y_{i_m})\}$$

where each i_k is sampled from $\{1, \dots, N\}$ with replacement.

Some points repeat, some disappear - this makes every tree see a different version of the world.

2. Feature Subsampling

When a tree tries to split a node, Random Forest restricts the possible features to a random subset.

If there are F total features, each ~~split~~ split sees only:

$$K = \sqrt{F} \quad (\text{for classification})$$

This creates diversity among trees because each tree uses different features at split.

3. Decision Tree Split Mathematics

Each tree uses the same math as a normal decision tree.

Entropy:

$$H(S) = - \sum_c P(c) \log_2 P(c)$$

Weighted Child Entropy:

$$H_{\text{split}} = \frac{|S_L|}{|S|} H(S_L) + \frac{|S_R|}{|S|} H(S_R)$$

Information Gain:

$$IG = H(S) - H_{\text{split}}$$

Every tree finds splits that maximize Information Gain.

Because every tree sees different samples and features, their splits differ naturally.

4. Ensemble Prediction (Majority Vote)

Once all trees make a prediction, the forest combines them.

For classification:

If T trees produce predictions:

$$h_1(x), h_2(x), \dots, h_T(x)$$

The final prediction is the most common class:

$$H_{\text{forest}}(x) = \text{mode}(h_1(x), h_2(x), \dots, h_T(x))$$

For regression:

Take the average:

$$H_{\text{forest}}(x) = \frac{1}{T} \sum_{i=1}^T h_i(x)$$

5. Why Random Forest Works (Mathematical Insight)

A single decision tree has low bias but high variance - it overfits.

Bagging reduces variance by averaging many noisy predictions.

Variance decreases roughly as:

$$\text{Var}_{\text{ensemble}} \approx \frac{\text{Var}_{\text{tree}}}{T}$$

(assuming trees are not too correlated)
Feature subsampling reduces correlation between trees, making this reduction stronger.
So even though each individual tree is a "weak learner", the forest becomes a strong model.

6. Out-of-Bag Error

Because each tree sees about 63% of the data (due to sampling with replacement), the remaining 37% can be used to estimate the model's performance without a separate test set.

This works because the probability a point is not chosen in a bootstrap sample is:

$$\left(1 - \frac{1}{N}\right)^N \approx e^{-1} \approx 0.37$$