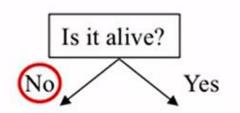
# Decision Trees and Random Forests



# **Decision Tree Example**







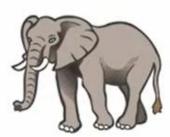
Objects with Alive == No













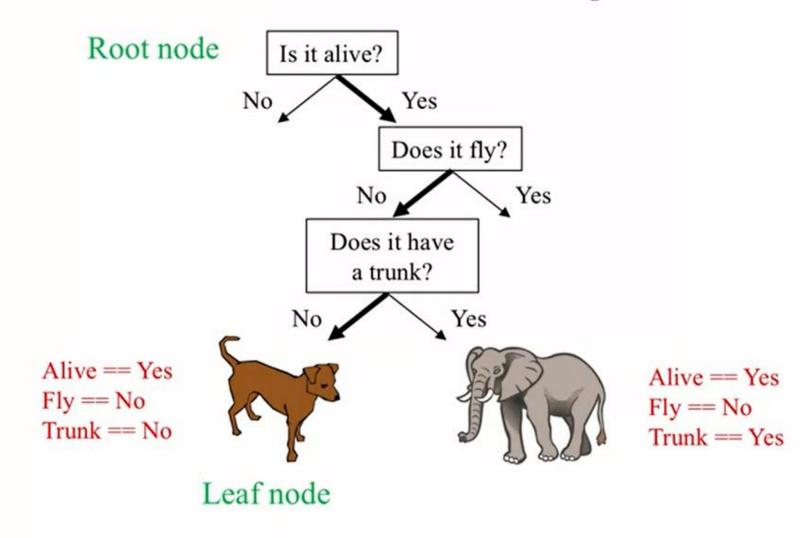


**Decision Tree Example** 





## Decision Tree Example







### The Iris Dataset



150 flowers 3 species 50 examples/species

Iris setosa

Iris versicolor

Iris virginica

#### The Iris Dataset

- Each instance in the dataset represents one of three different species of Iris, a type of flower.
  - setosa
  - Versicolor
  - vrginica
- There are four attributes or features for each instance, that represent measurements of different parts of the flower.
  - The sepal [canak yapragi] length in centimeters,
  - the sepal width in centimeters,
  - petal length in centimeters
  - petal width in centimeters.
- The data set has measurements for 150 flowers with 50 examples of each species.
- The classification task is to predict which of the three species and instances given these measurements.
- The iris data sets features are continuous values. Unlike our simple example involving yes or no questions which were binary features of the object.
- So the rules to be learned will be of the form for example, "Is Petal length greater than 2.3 centimeters?"



# **Decision Tree Splits**

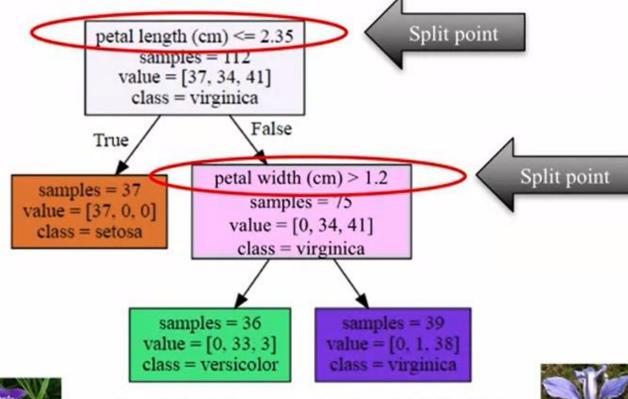


samples at this leaf have: petal length <= 2.35



setosa

versicolor



samples at this leaf have: petal length > 2.35

AND petal width <= 1.2

samples at this leaf have: petal length > 2.35

AND petal width > 1.2



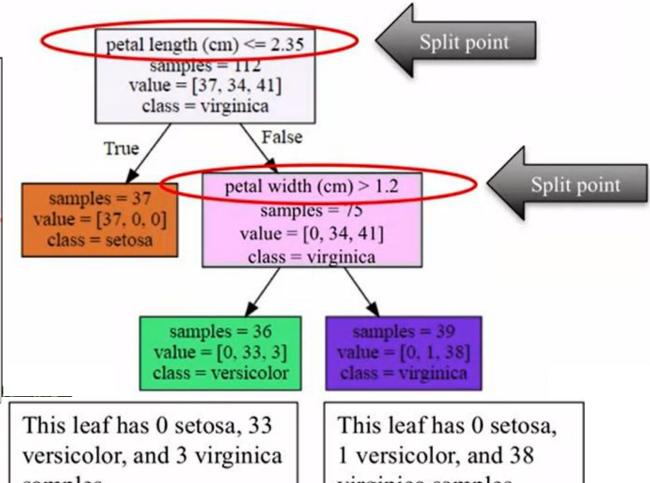
virginica

# **Decision Tree Splits**

The value list gives the number of samples of each class that end up at this leaf node during training.

The iris dataset has 3 classes. so there are three counts.

This leaf has 37 setosa samples, zero versicolor, and zero virginica samples.



samples.

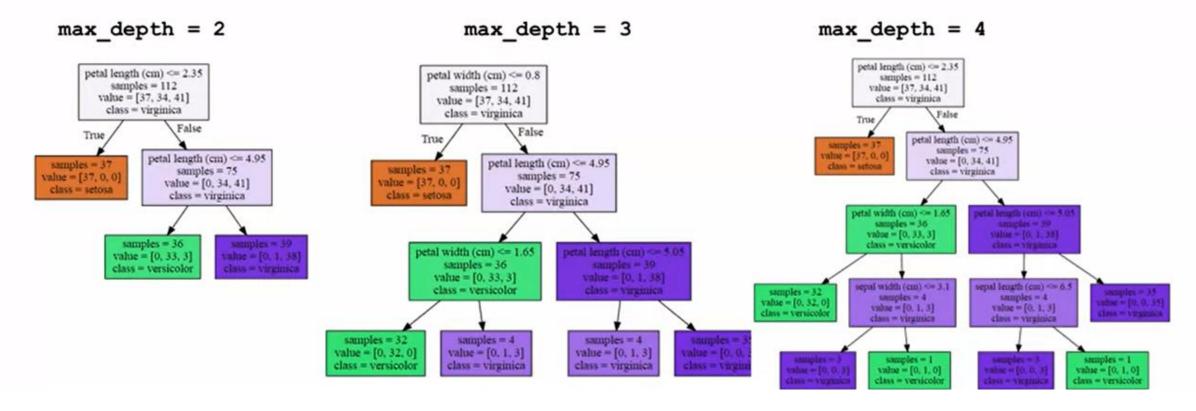
virginica samples.

- So to build the decision tree, the decision tree building algorithm starts by finding the feature that leads to the most informative split.
- For any given split of the data on a particular feature value, even for the best split it's likely that some examples will still be incorrectly classified or need further splitting.
- In the iris data set, if we look at all the flowers with petal length greater than 2.35 centimeters for example, using that split leaves a pool of instances that are a combination of virginica and versicolor that still need to be distinguished further





# Controlling the Model Complexity of Decision Trees



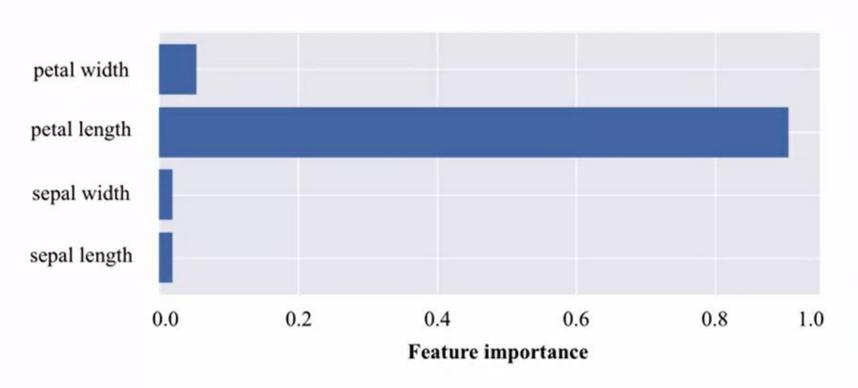


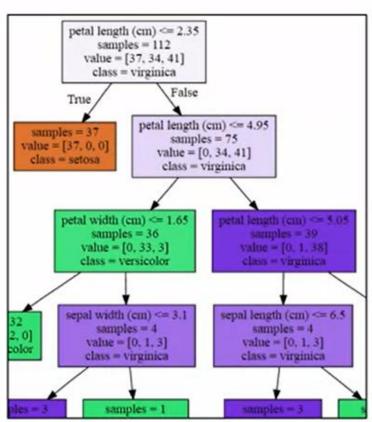
# Feature Importance: How important is a feature to overall prediction accuracy?

- A number between 0 and 1 assigned to each feature.
- Feature importance of 0 the feature was not used in prediction.
- Feature importance of I the feature predicts the target perfectly.
- All feature importances are normalized to sum to 1.



# Feature Importance Chart





**Decision tree** 



### **Decision Trees: Pros and Cons**

#### Pros:

- Easily visualized and interpreted.
- No feature normalization or scaling typically needed.
- Work well with datasets using a mixture of feature types (continuous, categorical, binary)

#### Cons:

- Even after tuning, decision trees can often still overfit.
- Usually need an ensemble of trees for better generalization performance.



#### Decision Trees: DecisionTreeClassifier Key Parameters

- max\_depth: controls maximum depth (number of split points).
   Most common way to reduce tree complexity and overfitting.
- min\_samples\_leaf: threshold for the minimum # of data instances a leaf can have to avoid further splitting.
- max\_leaf\_nodes: limits total number of leaves in the tree.
- In practice, adjusting only one of these (e.g. max\_depth) is enough to reduce overfitting.

### Ensemble Learning

- A widely used and effective method in machine learning involves creating learning models known as ensembles.
- An ensemble takes multiple individual learning models and combines them to produce an aggregate model that is more powerful than any of its individual learning models alone.
- Why are ensembles effective?
- Well, one reason is that if we have different learning models, although each of them might perform well individually, they'll tend to make different kinds of mistakes on the data set.

### Ensemble Learning

- Because each individual model might overfit to a different part of the data.
- By combining different individual models into an ensemble, we can average out their individual mistakes to reduce the risk of overfitting while maintaining strong prediction performance.
- Random forests are an example of the ensemble idea applied to decision trees.
- Random forests are widely used in practice and achieve very good results on a wide variety of problems.

- An ensemble of trees, not just one tree.
- Widely used, very good results on many problems.
- sklearn.ensemble module:
  - Classification: RandomForestClassifier
  - Regression: RandomForestRegressor
- One decision tree → Prone to overfitting.
- Many decision trees → More stable, better generalization
- Ensemble of trees should be diverse: introduce random variation into tree-building.

The idea is that each of the individual trees in a random forest should do reasonably well at predicting the target values in the training set **but** should also be constructed to be **different** in some way from the other trees in the forest

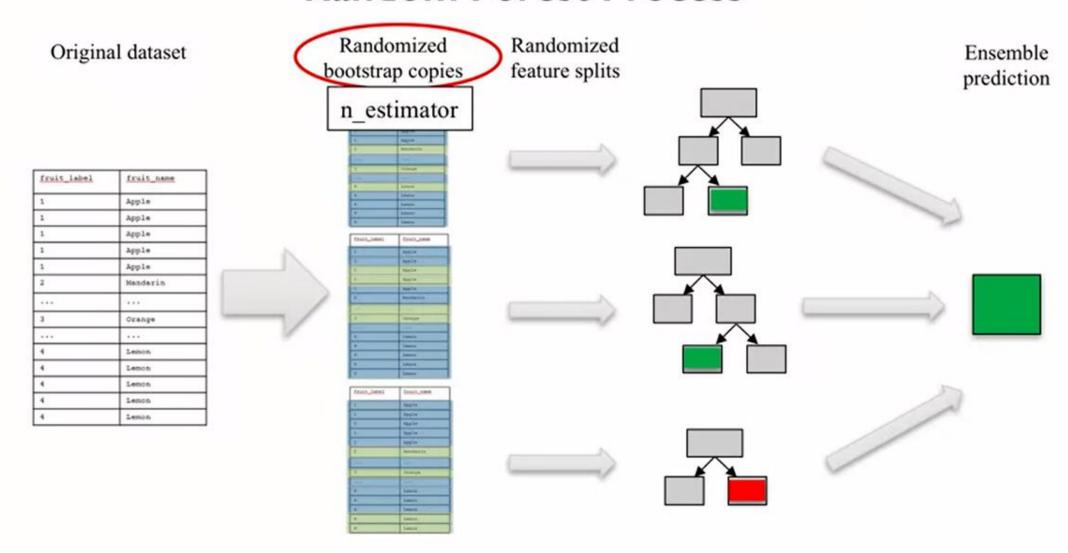
- This difference is accomplished by introducing random variation into the process of building each decision tree.
- This random variation during tree building happens in two ways.
  - First, the data used to build each tree is selected randomly and
  - Second the features chosen in each split tests are also randomly selected
- To create a random forest model you first decide on how many trees to build.
- This is set using the **n\_estimated** parameter.
- Each tree were built from a different random sample of the data called the bootstrap sample.

- Bootstrap samples are commonly used in statistics and machine learning.
- If your training set has N instances or samples in total, a bootstrap sample of size N is created by just repeatedly picking one of the N dataset rows at random with replacement, that is, allowing for the possibility of picking the same row again at each selection.
- You repeat this random selection process N times.
- The resulting bootstrap sample has N rows just like the original training set but with possibly some rows from the original dataset missing and others occurring multiple times just due to the nature of the random selection with replacement





#### **Random Forest Process**







### Random Forest Process: Bootstrap Samples

Bootstrap sample 1

Bootstrap	sample 2
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Bootstra	n	sam	n]	le	3
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fruit_label	fruit_name
1	Apple
2	Mandarin
3	Orange
***	
4	Lemon

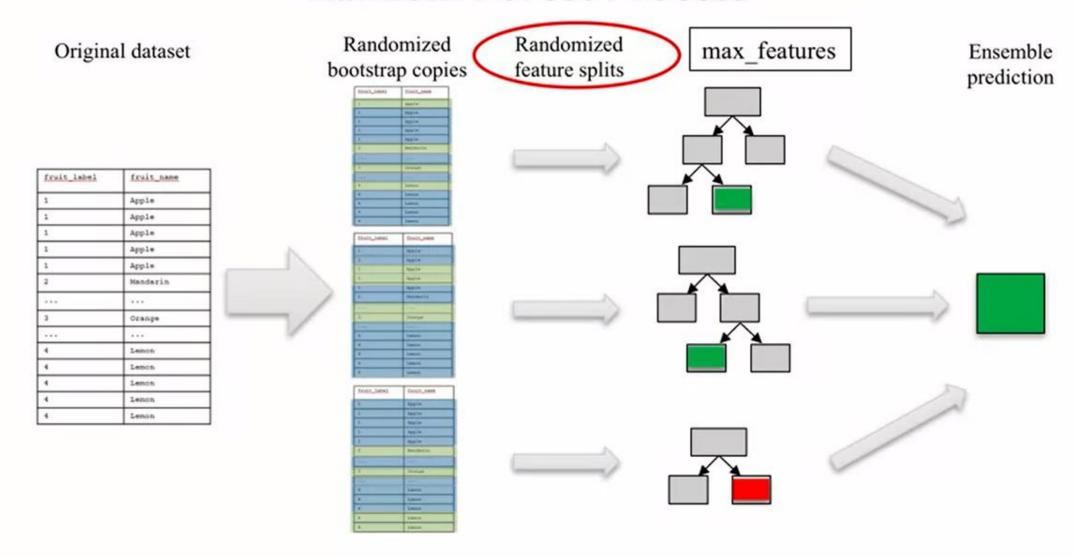
fruit_label	fruit_name
1	Apple
2	Mandarin
***	***
3	Orange
4	Lemon

fruit_label	fruit_name
1	Apple
2	Mandarin
3	Orange
***	
4	Lemon





#### **Random Forest Process**



### Random Forest max\_features Parameter

- Learning is quite sensitive to max\_features.
- Setting max\_features = 1 leads to forests with diverse, more complex trees.
- Setting max\_features = <close to number of features> will lead to similar forests with simpler trees.



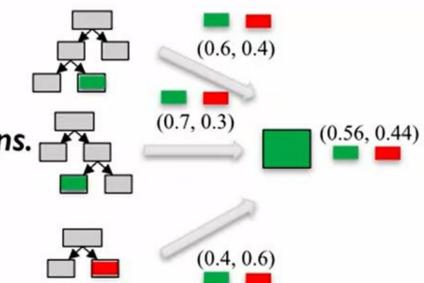


# Prediction Using Random Forests

I. Make a prediction for every tree in the forest.

#### 2. Combine individual predictions

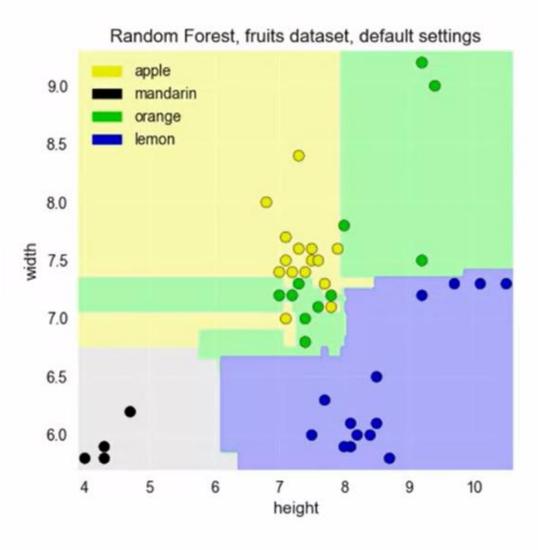
- Regression: mean of individual tree predictions.
- Classification:
  - Each tree gives probability for each class.
  - Probabilities averaged across trees.
  - Predict the class with highest probability.







### Random Forest: Fruit Dataset





### Random Forest: Pros and Cons

#### Pros:

- Widely used, excellent prediction performance on many problems.
- Doesn't require careful normalization of features or extensive parameter tuning.
- Like decision trees, handles a mixture of feature types.
- Easily parallelized across multiple CPUs.

#### Cons:

- The resulting models are often difficult for humans to interpret.
- Like decision trees, random forests may not be a good choice for very highdimensional tasks (e.g. text classifiers) compared to fast, accurate linear models.

# Random Forests: RandomForestClassifier Key Parameters

- n\_estimators: number of trees to use in ensemble (default: 10).
  - Should be larger for larger datasets to reduce overfitting (but uses more computation).
- max\_features: has a strong effect on performance. Influences the diversity of trees in the forest.
  - Default works well in practice, but adjusting may lead to some further gains.
- max\_depth: controls the depth of each tree (default: None. Splits until all leaves are pure).
- n\_jobs: How many cores to use in parallel during training.
- Choose a fixed setting for the random\_state parameter if you need reproducible results.

- **N\_estimators** sets the number of trees to use.
  - The default value for n\_estimators is 10 and increasing this number for larger data sets is almost certainly a good idea since ensembles that can average over more trees will reduce overfitting.
  - Just bear in mind that increasing the number of trees in the model will also increase the computational cost of training.
  - You'll use more time and more memory.
  - So in practice you'll want to choose the parameters that make best use of the resources available on your system.
- The max\_features parameter has a strong effect on performance.
- It has a large influence on how diverse the random trees in the forest are.
- Typically, the default setting of max features,
  - for classification is the square root of the total number of features and
  - for regression is the log base two of the total number of features

- The max depth parameter controls the depth of each tree in the ensemble.
- The default setting for this is none, in other words, the nodes in a tree will continue to be split until *all leaves* contain the same class or have fewer samples than the minimum sample split parameter value, which is two by default.
- Most systems now have a multi-core processor and so you can use the **end jobs** parameter to tell the random forest algorithm how many cores to use in parallel to train the model.
- Generally, you can expect something close to a linear speed up.
- So, for example, if you have four cores, the training will be four times as fast as if you just used
  one.
- If you set end jobs to -1 it will use all the cores on your system and setting end jobs to a number that's more than the number of cores on your system won't have any additional effect.
- Finally, given the random nature of random forests, if you want reproducible results it's
  especially important to choose a fixed setting for the random state parameter.