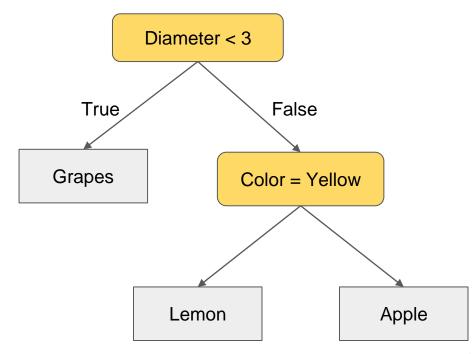
# RANDOM FORESTS



# **DECISION TREE: RECAP**

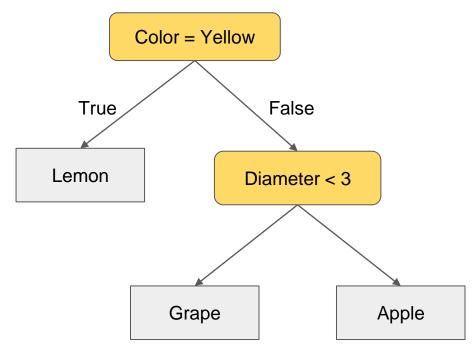
Color	Diameter	Label
Red	3	Apple
Yellow	3	Lemon
Purple	1	Grapes
Red	3	Apple
Yellow	3	Lemon
Purple	1	Grapes





# **ALTERNATE DECISION TREE**

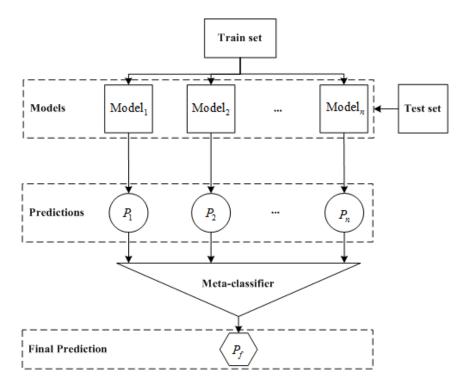
Color	Diameter	Label
Red	3	Apple
Yellow	3	Lemon
Purple	1	Grapes
Red	3	Apple
Yellow	3	Lemon
Purple	1	Grapes





Ensemble Learning is the process of combining multiple models with relatively lower accuracy in order to create a system that eventually produces a high accuracy.

We explicitly use ensemble learning to seek better predictive performance, such as lower error on regression or high accuracy for classification

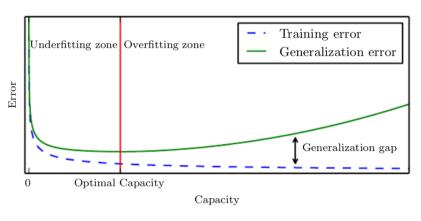




The fundamental principle of the **ensemble model** is that **a group of weak learners come together to form a strong learner, which increases the accuracy of the model.** When we try to **predict the target variable by any machine learning technique**, the **main causes of the difference between the actual and predicted values are noise, variance and bias**. The set reduces these factors (except noise, which is an irreducible error).

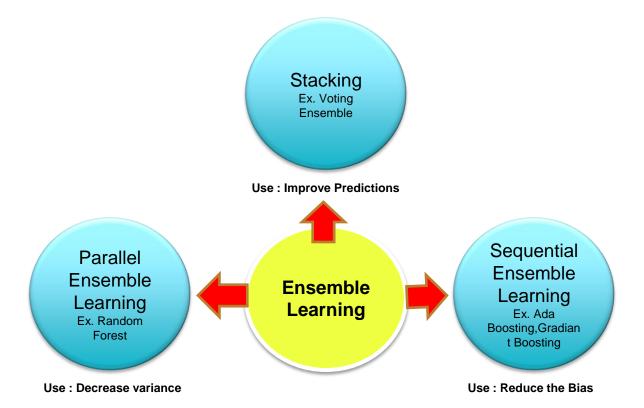


Here, we can see that if we will get the the variance, noise and bias in the raw data, image or any other format of the data. So, our model is going either under-fitting or over-fitting.

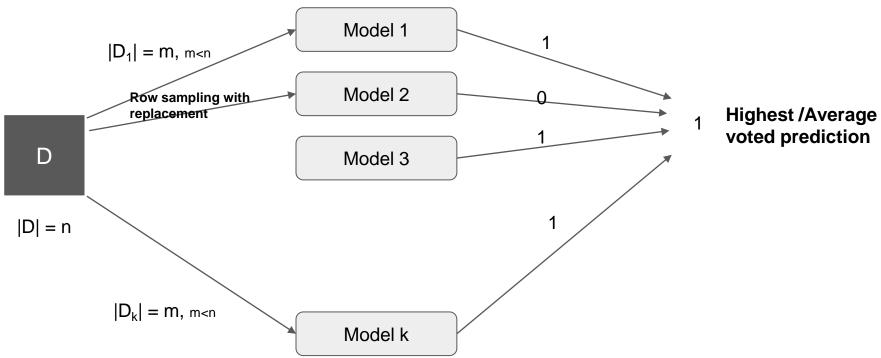


This reason is create big impact on your model directly here the ensemble learning comes in the picture. **Training error** and **generalization error** has gap which is represent as Generalization gap, which is show's that model is **under-fit** or **over-fit**.











The image shows that the bagging example has three steps:

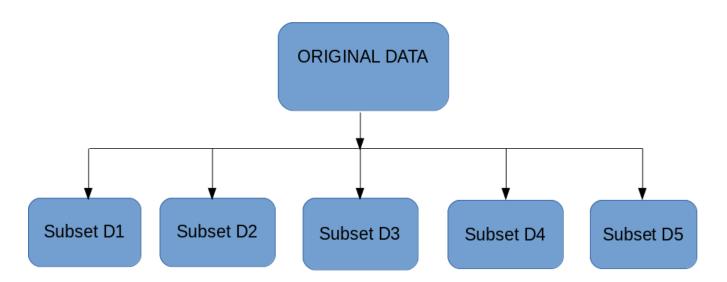
- Bootstrap data(Sampling)
- Aggregation or Model fit
- Combination different model with Result aggregation.
- Bootstrapping is a sampling technique in which we create multiple random sample from our training data-set.



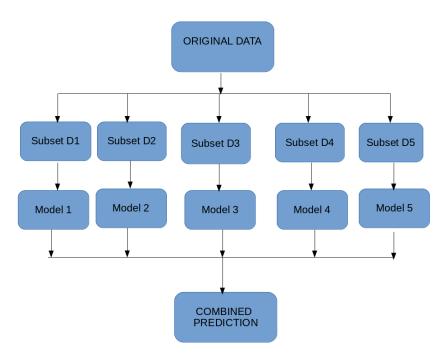
- The idea behind bagging is combining the results of multiple models (for instance, all decision trees) to get a generalized result.
- Bagging uses a sampling technique called Bootstrapping.
- Bootstrapping is a sampling technique in which we create subsets of observations from the original dataset, with replacement.
- Bagging (or Bootstrap Aggregating) technique uses these subsets (bags) to get a fair idea of the distribution (complete set).
- The size of subsets created for bagging may be same or less than the original set.



• Multiple subsets are created from the original dataset, selecting observations with replacement.



- A base model (weak model) is created on each of these subsets.
- The models run in parallel and are independent of each other.
- The final predictions are determined by combining the predictions from all the models.





# **HOW DOES RANDOM FOREST WORK?**

- Decisions trees are very sensitive to the data they are trained on small changes to the training set can result in significantly different tree structures.
- Random forest takes advantage of this by allowing each individual tree to randomly sample from the dataset with replacement, resulting in different trees – Bagging.
- RF consists multiple decision trees which act as base learners. Each
  decision tree is given a subset of random samples from the data set (hence
  the name *random*).



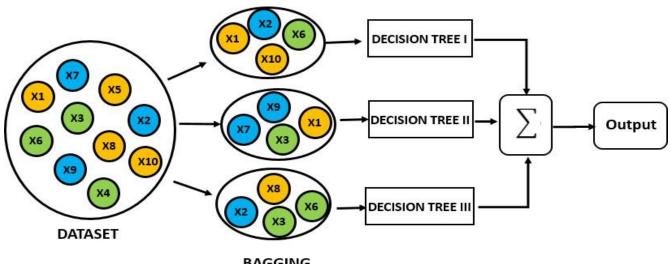
## **HOW DOES RANDOM FOREST WORK?**

- RF algorithm uses an Ensemble method Bagging (Bootstrap Aggregating)
- Then, Random Forest train each base learner (i.e Decision Tree) on a different sample of data and the sampling of data points happens with replacement.



# RANDOM FOREST WORKING

- Consider a training dataset: [X1, X2, X3, ... X10, Y].
- Random forest will create decision trees taking the input from subset using bagging as shown below:





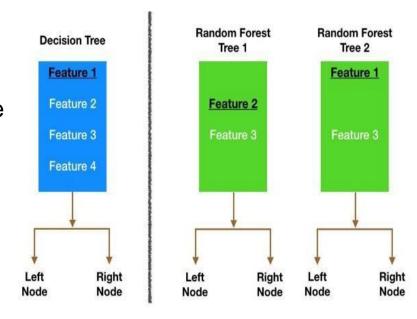
# RANDOM FOREST WORKING

- Note that with bagging we are not sub-setting the training data into smaller chunks and training each tree on a different chunk.
- Rather, if we have a sample of size N or 1000 rows, we are still feeding each tree a training set of size N.
- But instead of the original training data, we take a random sample of size N with replacement.
- For example, if our training data was [10, 11, 12,13, 14, 15, 16] then we might give one of our trees the following list [11, 12,12, 13,13,16, 16].



# RANDOM FOREST WORKING

- Feature Randomness
- In a normal decision tree, we consider the variable with highest gain.
- In contrast, each tree in a random forest can pick only from a random subset of features.
- This forces even more variation amongst the trees in the model and ultimately results in lower correlation across trees and more diversification.





# **HYPER-PARAMETERS RANDOM FOREST:**

- Optimization of RF depends on few inbuilt parameters.
- n\_estimators\* number of decision trees that the algorithm creates. As the number tree increases, the performance increases and the predictions are more stable but it slows down the computation.
- max\_features\* maximum number of features that are considered for splitting a node.
- **n\_jobs** number of jobs to run in parallel. If n\_jobs=1, it uses one processor. If n\_jobs=-1, then the number of jobs is set to the number of cores available.



# **HYPER-PARAMETERS RANDOM FOREST:**

- max\_depth is the maximum depth of the tree. The deeper the tree, the more splits it has and it captures more information about the data.
- **criterion** is the function to measure the quality of a split. Supported criteria are "**gini**" for the Gini impurity and "**entropy**" for the information gain.



## WHY RANDOM FOREST?

- •Reduces Risk of Over fitting
- High Accuracy
- Maintain Accuracy even if more data is missing



# **THAT'S ALL FOLKS**

