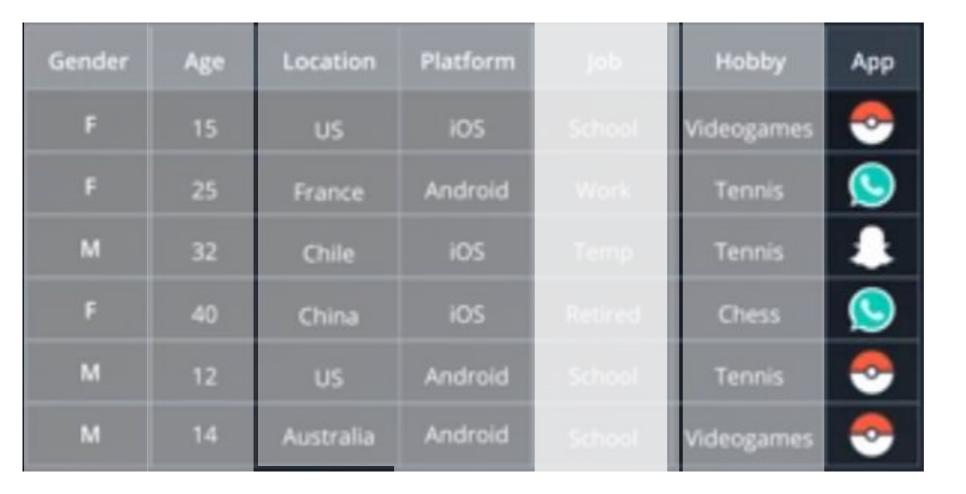
Random Forest

In Random Forest, we grow multiple trees as opposed to a single tree in CART model d. To classify a new object based on attributes, each tree gives a classification and we say the tree "votes" for that class. The forest chooses the classification having the most votes (over all the trees in the forest) and in case of regression, it takes the average of outputs by different trees.













How Random Forest Works?



Random Forest Hyperparameters

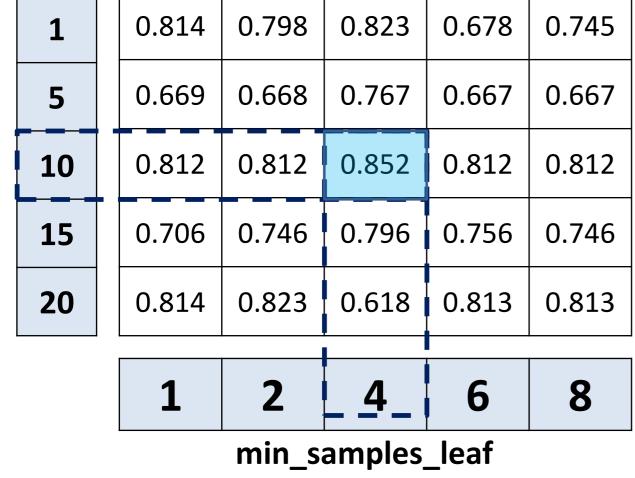
There are 4 hyperparameters required for a Random Forest classifier;

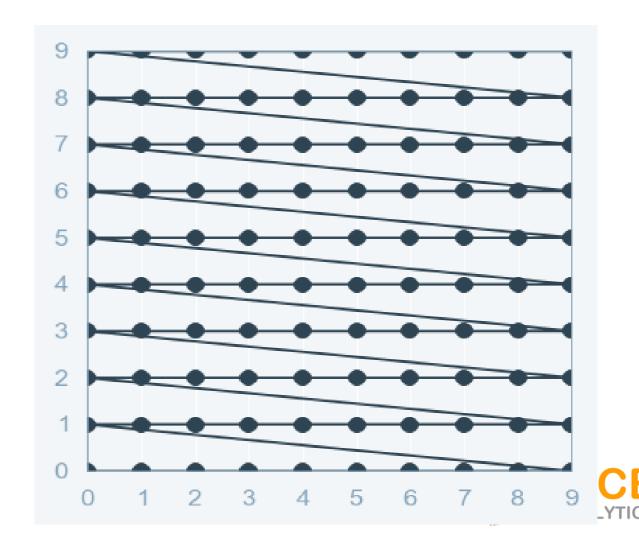
- > The number of trees in the forest (n_estimators).
- ➤ The number of features to consider at each split. By default: square root of total number of features (max_features).
- The maximum depth of a tree i.e. number of nodes (max_depth).
- The minimum number of samples required to be at a leaf node / bottom of a tree (min_samples_leaf).



Tuning by GridSearchCV

- ➤ We try every combination of a preset list of values of the hyperparameters and evaluate the model for each combination.
- Pattern followed here is similar to the grid, where all the values are placed in the form of a matrix
- Once all the combinations are evaluated, the model with the set of parameters which give the top accuracy is considered to be the best.

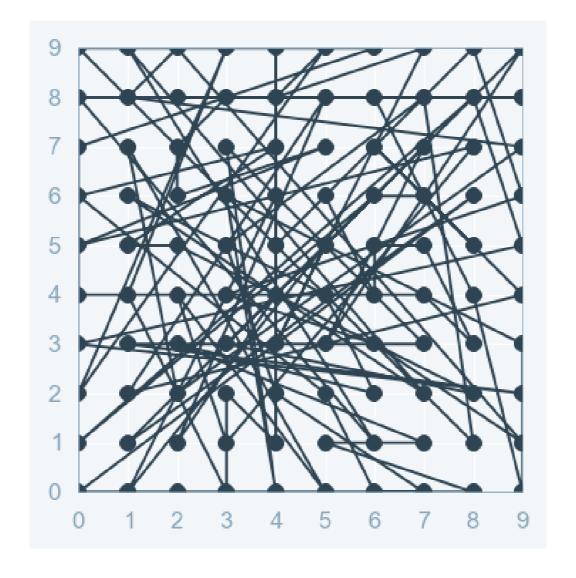




max_depth

Tuning by RandomSearchCV

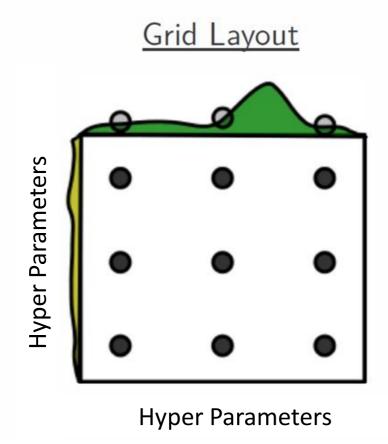
- > Random search is a technique where random combinations of the hyper parameters are used to find the best solution for the built model
- > It tries random combinations of a range of values
- To optimize with random search, the function is evaluated at some number of random configurations in the parameter space

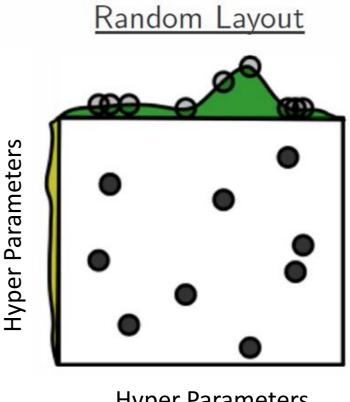




GridSearch vs RandomSearch

- > One of the major drawbacks of grid search is that when it comes to dimensionality, it suffers when the number of hyper parameters grows exponentially.
- > The chances of finding the optimal parameter are comparatively higher in random search because of the random search pattern where the model might end up being trained on the optimized parameters without any aliasing number of random configurations in the parameter space





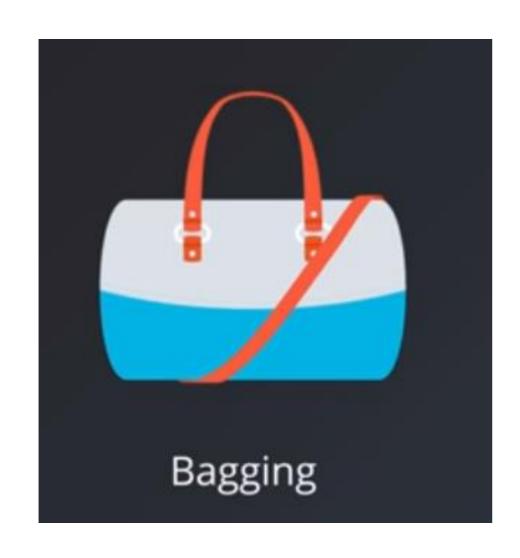




Let's Practice



ENSEMBLE METHODS



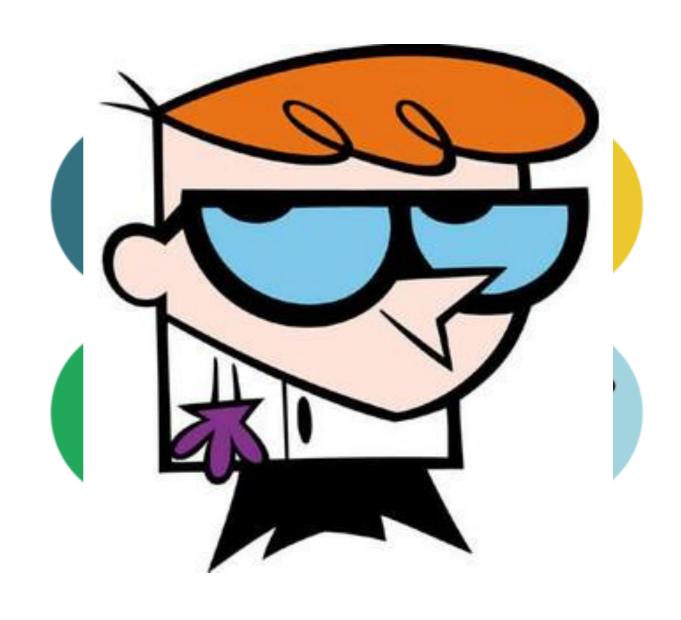


BOOTSRAP AGGREGATION



BAGGING





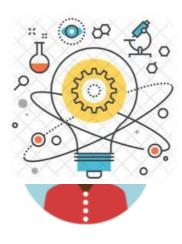


BOOSTING















ENSEMBLE METHODS

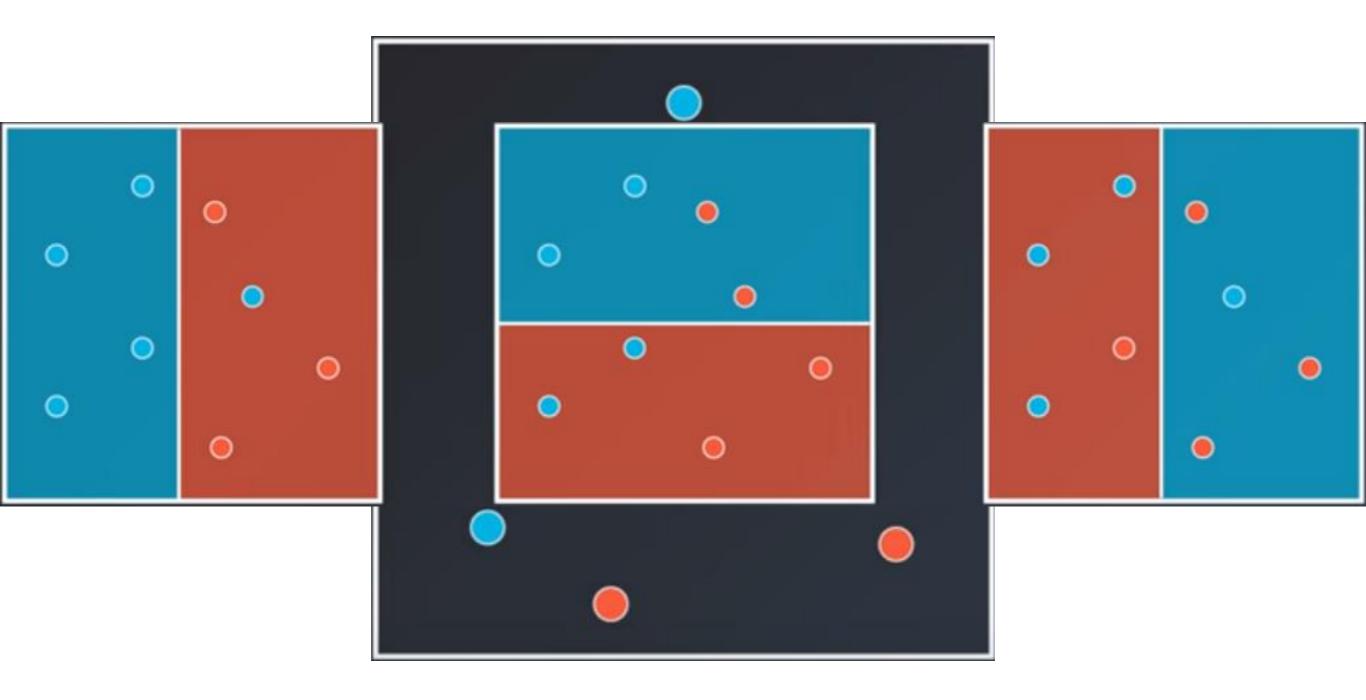


WEAK LEARNERS





Bagging Explained





Bagging Explained

