In [5]: import pandas as pd df = pd.read csv('tahkeer data cleaned.csv') In [6]: columns = df.columns.tolist() columns.remove("smoking") features x = df[columns]class_y = df["smoking"] In [7]: from sklearn.model_selection import train test split from sklearn.metrics import accuracy score xtrain, xtest, ytrain, ytest = train_test_split(features_x, class_y, test_size=0.30, shuffle=False, tra in size=0.70) In [8]: import numpy as np from sklearn.tree import DecisionTreeClassifier class BaggingClassifier: def init (self, n estimators=100, max depth=None, threshold=0.5): self.n_estimators = n_estimators self.max depth = max depth self.threshold = threshold self.models = [DecisionTreeClassifier(max_depth=self.max_depth) for _ in range(n_estimators)] def fit(self, x, y): for model in self.models: indices = np.random.choice(len(x), len(x), replace=True) x subset, y subset = x.iloc[indices], y.iloc[indices] model.fit(x subset, y subset) return self def predict(self, x): pred = np.zeros((len(x), self.n_estimators)) for i, model in enumerate(self.models): pred[:, i] = model.predict(x) avg predictions = np.mean(pred, axis=1) binary_predictions = (avg_predictions >= self.threshold).astype(int) return binary predictions def get params(self, deep=True): return { 'n_estimators': self.n_estimators, 'max depth': self.max depth, 'threshold': self.threshold, def set params(self, **params): if not params: return self for param, value in params.items(): setattr(self, param, value) return self In [9]: class AdaBoost: def __init__(self, n_estimators=50, max_depth=4): $self.n estimators = n_estimators$ self.alphas = [] self.models = []self.max depth = max depth def fit(self, x, y): x = np.array(x)y = np.array(y)# Get the number of rows and columns n samples, n features = x.shape # Initialize the weights to 1/Nweights = np.ones(n samples) / n samples for in range(self.n estimators): model = DecisionTreeClassifier(max depth=self.max depth) model.fit(x, y, sample weight=weights) y pred = model.predict(x) weighted_error = np.sum(weights[y_pred != y]) if weighted error >= 0.5: break alpha = 0.5 * np.log((1.0 - weighted error) / max(weighted error, 1e-10))self.alphas.append(alpha) self.models.append(model) # Update weights weights *= np.exp(-alpha * y * y_pred) weights /= np.sum(weights) return self def predict(self, x): x = np.array(x)pred = np.zeros(len(x))for alpha, model in zip(self.alphas, self.models): pred += alpha * model.predict(x) return np.sign(pred) def get params(self, deep=True): return { 'n estimators': self.n estimators, 'max depth': self.max depth, def set params(self, **params): if not params: return self for param, value in params.items(): setattr(self, param, value) return self In []: class RandomForestClassifier: def __init__(self, n_estimators=100, max_depth=None, min_samples_split=2, min_samples_leaf=1): self.n estimators = n estimators self.max depth = max depth self.min_samples_split = min samples split self.min_samples_leaf = min_samples_leaf self.models = [] def fit(self, x, y): x = np.array(x)y = np.array(y)n_samples, n_features = x.shape for _ in range(self.n_estimators): # Randomly select a subset of features selected_features = np.random.choice(n_features, size=int(np.sqrt(n_features)), replace=Fal se) x_subset = x[:, selected_features] # Create a decision tree with random features tree = DecisionTreeClassifier(max depth=self.max depth, min_samples_split=self.min_samples_split, min samples leaf=self.min samples leaf tree.fit(x subset, y) self.models.append((tree, selected_features)) return self def predict(self, x): x = np.array(x)pred = np.zeros((x.shape[0], self.n_estimators)) for i, (tree, selected features) in enumerate(self.models): x subset = x[:, selected features] pred[:, i] = tree.predict(x subset) # Use majority voting for the final prediction final_predictions = np.apply_along_axis(lambda x: np.bincount(x.astype(int)).argmax(), axis=1, arr=pred) return final predictions def get params(self, deep=True): return { 'n estimators': self.n estimators, 'max depth': self.max depth, 'min_samples_split': self.min samples split, 'min_samples_leaf': self.min_samples_leaf def set params(self, **params): if not params: return self for param, value in params.items(): setattr(self, param, value) return self In [30]: bagging model = BaggingClassifier(n estimators=100, max depth=60) bagging_model.fit(xtrain, ytrain) bagging_predictions = bagging_model.predict(xtest) bagging_accuracy = accuracy_score(ytest, bagging_predictions) print(f"Bagging Accuracy: {bagging_accuracy}") Bagging Accuracy: 0.7316913595880292 In [31]: | adaboost = AdaBoost(n estimators=300) adaboost.fit(xtrain, ytrain) predictions = adaboost.predict(xtest) accuracy = accuracy score(ytest, predictions) print(f'Boosting Accuracy: {accuracy}') Boosting Accuracy: 0.7187117470769491 In [32]: random forest = RandomForestClassifier(n estimators=150, max depth=3, min samples split=2, min samples leaf=1)random_forest.fit(xtrain, ytrain) predictions = random forest.predict(xtest) accuracy = accuracy_score(ytest, predictions) print(f'Random Forest Accuracy: {accuracy}') Random Forest Accuracy: 0.7062597610907095 **Hyperparameter Tuning** We will use grid search and randomized search methods to choose better hyperparameters In [22]: from sklearn.model_selection import GridSearchCV from sklearn.model selection import RandomizedSearchCV from skopt import BayesSearchCV import multiprocessing n jobs = multiprocessing.cpu count()-1 np.int = intdef print results(search results): best params = search_results.best_params_ best_score = search_results.best_score_ print("Best Parameters:", best params) print("Best Score:", best_score) **Tuning Bagging model** In [23]: params = { 'n_estimators': [100, 200, 250, 300], 'max_depth': [50, 60, 70, 90], 'threshold': [0.3, 0.5, 0.7], bagging = BaggingClassifier() grid_search = GridSearchCV(estimator=bagging, param_grid=params, scoring='accuracy', n_jobs=n_jobs) grid search.fit(xtrain, ytrain) print("Grid search - Bagging: ") print_results(grid_search) Grid search - Bagging: Best Parameters: {'max depth': 50, 'n estimators': 100, 'threshold': 0.5} Best Score: 0.7288794679763679 In [24]: random search = RandomizedSearchCV(estimator=bagging, param distributions=params, n iter=6, scoring='ac curacy', random_state=42, n_jobs=n_jobs) random_search.fit(xtrain, ytrain) print("Random search - Bagging: ") print_results(random_search) Random search - Bagging: Best Parameters: {'threshold': 0.5, 'n_estimators': 100, 'max_depth': 90} Best Score: 0.7281558532037508 In []: bayes optimization = BayesSearchCV(bagging, params, n iter=6, scoring='accuracy', random state=42, n jo bs=n jobs) bayes_optimization.fit(xtrain, ytrain) print("Bayesian Optimization - Bagging: ") print_results(bayes_optimization) **Tuning AdaBoost model** In [15]: params = { 'n estimators': [100, 200, 250, 300, 500, 1000], adaboost = AdaBoost() grid search = GridSearchCV(estimator=adaboost, param grid=params, scoring='accuracy', n jobs=n jobs) grid_search.fit(xtrain, ytrain) print("Grid search - AdaBoost: ") print_results(grid_search) Grid search - AdaBoost: Best Parameters: {'n estimators': 100} Best Score: 0.714262437542956 In [37]: random_search = RandomizedSearchCV(estimator=adaboost, param_distributions=params, n_iter=6, scoring='a ccuracy', random_state=42, n_jobs=n_jobs) random search.fit(xtrain, ytrain) print("Random search - AdaBoost: ") print_results(random_search) Random search - AdaBoost: Best Parameters: {'n_estimators': 100} Best Score: 0.714262437542956 In [18]: bayes optimization = BayesSearchCV(adaboost, params, n iter=6, scoring='accuracy', random state=42, n j obs=n_jobs) bayes optimization.fit(xtrain, ytrain) print("Bayesian Optimization - AdaBoost: ") print_results(bayes_optimization) Bayesian Optimization - AdaBoost: Best Parameters: OrderedDict([('n_estimators', 250)]) Best Score: 0.714262437542956 **Tuning Random Forest model** In [38]: params = { 'n_estimators': [100, 200, 250], 'max depth': [None, 5, 10], 'min_samples_split': [2, 3], 'min_samples_leaf': [1, 2, 4], rf model = RandomForestClassifier() grid_search = GridSearchCV(estimator=rf_model, param_grid=params, scoring='accuracy', n_jobs=n_jobs) grid_search.fit(xtrain, ytrain) print("Grid search - Random Forest: ") print results(grid search) Grid search - Random Forest: Best Parameters: {'max_depth': 10, 'min_samples_leaf': 4, 'min_samples_split': 2, 'n_estimators': 25 Best Score: 0.7214714514997054 In [40]: random_search = RandomizedSearchCV(estimator=rf_model, param_distributions=params, n_iter=6, scoring='a ccuracy', random_state=42, n_jobs=n_jobs) random_search.fit(xtrain, ytrain) print("Random Search - Random Forest: ") print results(random search) Random Search - Random Forest: Best Parameters: {'n_estimators': 250, 'min_samples_split': 2, 'min_samples_leaf': 2, 'max_depth': 1 Best Score: 0.7211367772692194 In [20]: bayes_optimization = BayesSearchCV(rf_model, params, n_iter=6, scoring='accuracy', random_state=42, n_j obs=n jobs) bayes_optimization.fit(xtrain, ytrain) print("Bayesian Optimization - Random Forest: ") print_results(bayes_optimization) Bayesian Optimization - Random Forest: Best Parameters: OrderedDict([('max_depth', 10), ('min_samples_leaf', 2), ('min_samples_split', 3), ('n_estimators', 200)]) Best Score: 0.7210644310501265