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import numpy as np
class KMeans():
    # This function initializes the KMeans class
   def init (self, k = 3, num iter = 1000, order = 2):
        # Set a seed for easy debugging and evaluation
       np.random.seed(42)
       # This variable defines how many clusters to create
       # default is 3
       self.k = k
       # This variable defines how many iterations to recompute centroids
       # default is 1000
       self.num iter = num_iter
       # This variable stores the coordinates of centroids
       self.centers = None
       # This variable defines whether it's K-Means or K-Medians
       # an order of 2 uses Euclidean distance for means
       \# an order of 1 uses Manhattan distance for medians
        # default is 2
       if order == 1 or order == 2:
           self.order = order
       else:
           raise Exception("Unknown Order")
    # This function fits the model with input data (training)
    def fit(self, X):
       # m, n represent the number of rows (observations)
        # and columns (positions in each coordinate)
       m, n = X.shape
       # self.centers are a 2d-array of
        # (number of clusters, number of dimensions of our input data)
       self.centers = np.zeros((self.k, n))
       # self.cluster idx represents the cluster index for each observation
        # which is a 1d-array of (number of observations)
       self.cluster idx = np.zeros(m)
       ##### TODO 1 #####
       # Task: initialize self.centers
        # Instruction:
       # For each dimension (feature) in X, use the 10th percentile and
        # the 90th percentile to form a uniform distribution. Then, we will initialize
        # the values of each center by randomly selecting values from the distributions.
        # Note:
        # This method is by no means the best initialization method. However, we would
        # like you to follow our guidelines in this HW. We will ask you to discuss some better
        # initializaiton methods in the notebook.
        # Hint:
        # 1. np.random.uniform(), np.percentile() might be useful
        # 2. make sure to look over its parameters if you're not sure
       for i in range(n):
           lower, upper = np.percentile(X[:,i], [10, 90])
           self.centers[:,i] = np.random.uniform(lower, upper, self.k)
        ##### END TODO 1 #####
       for i in range(self.num iter):
           # new centers are a 2d-array of
            # (number of clusters, number of dimensions of our input data)
           new_centers = np.zeros((self.k, n))
            ##### TODO 2 ######
            # Task: calculate the distance and create cluster index for each observation
            # Instruction:
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# You should calculate the distance between each observation and each centroid
        # using specified self.order. Then, you should derive the cluster index for
        # each observation based on the minimum distance between an observation and
        # each of the centers.
        # Hint:
        # 1. np.linalg.norm() might be useful, along with parameter axis, ord
        # for that function
        # 2. You can transpose an array using .T at the end
        # 3. np.argmin() might be useful along with parameter axis in finding
        # the desired cluster index of all observations
        # IMPORTANT:
        # Copy-paste this part of your implemented code
        # to the predict function, and return cluster idx in that function
        \texttt{distances} = \texttt{np.array([np.linalg.norm(X - center, ord=self.order, axis=1)} \ \textbf{for} \ \texttt{center} \ \textbf{in} \ \texttt{self.centers])}
        cluster idx = np.argmin(distances.T, axis=1)
        ##### END TODO 2 #####
        ##### TODO 3 ######
        \slash\hspace{-0.4em}\# Task: calculate the coordinates of new_centers based on cluster_idx
        # Instruction:
        # You should assign the coordinates of the new center by calculating
        # mean/median of the coordinates of observations belonging to the same
        # cluster.
        # Hint:
        # 1. np.mean(), np.median() with axis might be helpful
        #####################
        for idx in range(self.k):
           cluster coordinates = X[cluster idx == idx]
            if self.order == 2:
                cluster center = np.mean(cluster coordinates, axis=0)
            elif self.order == 1:
                cluster_center = np.median(cluster_coordinates, axis=0)
            new centers[idx, :] = cluster center
        ##### END TODO 3 #####
        ##### TODO 4 ######
        # Task: determine early stop and update centers and cluster idx
        # Instructions:
        # You should stop tranining as long as cluster index for all
        # observations is the same as the previous iteration
        # Hint:
        # 1. .all() might be helpful
        if (cluster_idx == self.cluster_idx).all():
           print(f"Early Stopped at Iteration {i}")
            return self
        self.centers = new centers
        self.cluster idx = cluster idx
        ##### END TODO 4 #####
    return self
# This function makes predictions with input data
# Copy-paste your code from TODO 2 and return cluster idx
def predict(self, X):
   distances = np.array([np.linalg.norm(X - center, ord=self.order, axis=1) for center in self.centers])
   return np.argmin(distances.T, axis=1)
```