Clustering Human Activity using Inertial Sensors Data

Note:

- Use the next cell to download the data directly, if that didn't work. you can download it manually (available at UCI archive) a copy will also be available on Piazza.
- Don't change the part of the code that labels #Do not change
- Attach this notebook to your answer sheet with all outputs visible.
- make sure you have pytorch, scikit learn, pandas in your environment

```
#### Download the dataset
import urllib.request
import zipfile
import os
dataset url =
"https://archive.ics.uci.edu/static/public/240/human+activity+recognit
ion+using+smartphones.zip"
zip file path = "Dataset.zip"
extracted_downloaded_folder = "Dataset"
extracted data folder = "UCI HAR Dataset"
if not os.path.exists(zip_file_path):
    print("Downloading the dataset...")
    urllib.request.urlretrieve(dataset url, zip file path)
if not os.path.exists(extracted downloaded folder):
    print("Extracting the dataset...")
    with zipfile.ZipFile(zip file path, 'r') as zip ref:
        zip ref.extractall(".")
if not os.path.exists(extracted data folder):
    print("Extracting the dataset...")
    with zipfile.ZipFile(extracted data folder +'.zip', 'r') as
zip ref:
        zip ref.extractall(".")
print("Dataset is ready.")
```

```
Downloading the dataset...
Extracting the dataset...
Extracting the dataset...
Dataset is ready.
```

Load the data into a dataframe

```
# Import necessary libraries
import pandas as pd
import matplotlib.pyplot as plt
# Define paths to data files
train path = "UCI HAR Dataset/train" # TODO
test path = "UCI HAR Dataset/test"
                                     # TODO
activity mapper path = "UCI HAR Dataset/activity labels.txt"
# Load training and testing data
X_train = pd.read_csv(f"{train_path}/X_train.txt",
delim whitespace=True, header=None)
y train = pd.read csv(f"{train path}/y train.txt",
delim whitespace=True, header=None)
X_test, y_test = pd.read_csv(f"{test_path}/X test.txt",
delim whitespace=True, header=None),
pd.read_csv(f"{test_path}/y_test.txt", delim_whitespace=True,
header=None) # TODO
# Display the first 5 rows of the training dataframe
print("First 5 rows of training feature dataframe:")
X train.head() # DO NOT CHANGE
<ipython-input-5-le64e906fa49>:12: FutureWarning: The
'delim whitespace' keyword in pd.read csv is deprecated and will be
removed in a future version. Use ``sep='\s+'`` instead
  X train = pd.read csv(f"{train path}/X train.txt",
delim_whitespace=True, header=None)
<ipython-input-5-le64e906fa49>:13: FutureWarning: The
'delim whitespace' keyword in pd.read csv is deprecated and will be
removed in a future version. Use ``sep='\s+'`` instead
  y train = pd.read_csv(f"{train_path}/y_train.txt",
delim whitespace=True, header=None)
<ipython-input-5-le64e906fa49>:14: FutureWarning: The
'delim whitespace' keyword in pd.read csv is deprecated and will be
removed in a future version. Use ``sep='\s+'`` instead
  X test, y test =
                     pd.read csv(f"{test path}/X test.txt",
delim whitespace=True, header=None),
pd.read csv(f"{test path}/y test.txt", delim whitespace=True,
header=None) # TODO
```

```
First 5 rows of training feature dataframe:
<ipython-input-5-le64e906fa49>:14: FutureWarning: The
'delim whitespace' keyword in pd.read csv is deprecated and will be
removed in a future version. Use ``sep='\s+'`` instead
  X test, y test = pd.read csv(f"{test path}/X test.txt",
delim whitespace=True, header=None),
pd.read csv(f"{test path}/y test.txt", delim whitespace=True,
header=None) # TODO
                            2
                                                          5
                                                                    6
  0.288585 -0.020294 -0.132905 -0.995279 -0.983111 -0.913526 -
0.995112
1 0.278419 -0.016411 -0.123520 -0.998245 -0.975300 -0.960322 -
0.998807
  0.279653 - 0.019467 - 0.113462 - 0.995380 - 0.967187 - 0.978944 -
0.996520
3 0.279174 -0.026201 -0.123283 -0.996091 -0.983403 -0.990675 -
0.997099
4 0.276629 -0.016570 -0.115362 -0.998139 -0.980817 -0.990482 -
0.998321
       7
                 8
                            9
                                           551
                                                     552
                                                               553
554 \
0 -0.983185 -0.923527 -0.934724 ... -0.074323 -0.298676 -0.710304 -
0.112754
1 - 0.974914 - 0.957686 - 0.943068 \dots 0.158075 - 0.595051 - 0.861499
0.053477
2 -0.963668 -0.977469 -0.938692 ... 0.414503 -0.390748 -0.760104 -
0.118559
3 -0.982750 -0.989302 -0.938692 ... 0.404573 -0.117290 -0.482845 -
0.036788
4 -0.979672 -0.990441 -0.942469 ... 0.087753 -0.351471 -0.699205
0.123320
        555
                  556
                                                559
                            557
                                      558
0 0.030400 -0.464761 -0.018446 -0.841247
                                           0.179941 -0.058627
1 -0.007435 -0.732626  0.703511 -0.844788
                                           0.180289 -0.054317
2 0.177899 0.100699 0.808529 -0.848933
                                           0.180637 -0.049118
3 -0.012892
            0.640011 -0.485366 -0.848649
                                           0.181935 -0.047663
4 0.122542 0.693578 -0.615971 -0.847865
                                           0.185151 -0.043892
[5 rows x 561 columns]
```

scaling the data and PCA

```
from sklearn.preprocessing import StandardScaler
# TODO: Scale X_train
scaler = StandardScaler()
```

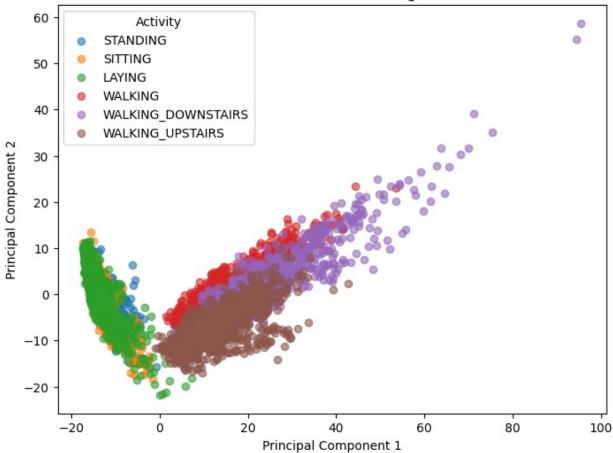
```
X train scaled = scaler.fit transform(X train)
# TODO: Scale X test
X test scaled = scaler.transform(X test)
# Convert scaled arrays back to DataFrames
X_train = pd.DataFrame(X_train_scaled)
X test = pd.DataFrame(X test scaled)
# Add 'Activity' column to create training df and testing df
# TODO: Combine X train and y train into a single DataFrame named
training df.
training df = pd.concat([X train, y train.rename(columns={0:
'Activity'})], axis=1)
# TODO: Combine X test and y test into a single DataFrame named
testing df.
testing df = pd.concat([X test, y test.rename(columns={0:
'Activity'})], axis=1)
# Display the first 5 rows of the training feature dataframe
print("First 5 rows of training feature dataframe:")
training df.head() # DO NOT CHANGE
First 5 rows of training feature dataframe:
                   1
                             2
0 0.200642 -0.063683 -0.419628 -0.868814 -0.939441 -0.737529 -
0.859817
1 0.055948 0.031486 -0.253908 -0.875426 -0.923902 -0.849304 -
0.868531
2 0.073515 -0.043416 -0.076295 -0.869039 -0.907760 -0.893785 -
0.863137
3 0.066696 -0.208422 -0.249712 -0.870626 -0.940022 -0.921805 -
0.864503
4 0.030469 0.027587 -0.109848 -0.875188 -0.934878 -0.921343 -
0.867384
                   8
                             9 ...
                                          552
                                                    553
555 \
0 -0.939019 -0.766437 -0.856036 ... 0.025960 -0.276399 -0.360603
0.062940
1 - 0.921998 - 0.848928 - 0.871359 \dots - 0.897357 - 0.767990 0.133011 -
0.021461
2 -0.898854 -0.896701 -0.863323 ... -0.260878 -0.438316 -0.377840
0.391976
3 -0.938124 -0.925279 -0.863323 ... 0.591045 0.463155 -0.135025 -
```

```
0.033637
4 -0.931789 -0.928028 -0.870260 ... -0.138515 -0.240313 0.340406
0.268486
       556
                 557
                           558
                                     559
                                               560
                                                    Activity
0 -0.778427 -0.026080 -0.687219 0.407946 -0.007568
                                                           5
1 -1.218805 1.484470 -0.694138 0.409117 0.007875
                                                           5
                                                           5
2 0.151207 1.704201 -0.702239 0.410288 0.026502
  1.037851 -1.003019 -0.701684 0.414650 0.031714
                                                           5
4 1.125918 -1.276282 -0.700152 0.425463 0.045225
                                                           5
[5 rows x 562 columns]
# TODO perform PCA on the train data and get the first 2 PC
from sklearn.decomposition import PCA
pca = PCA(n components=2)
X_train_pca = pca.fit_transform(X train) #TODO
```

Visualize the data

```
# Visualize training data using PCA
# Use the feature decoder to create Activity Name column
# Load activity labels
activity labels = pd.read csv(activity mapper path, header=None,
sep='\s+', names=['id', 'activity_name'])
# Create mapping dictionary {1: "WALKING", 2: "WALKING UPSTAIRS", ...}
activity_mapping = dict(zip(activity_labels['id'],
activity labels['activity_name']))
# Use the mapping to decode the Activities labels
Activity Name = y train[0].map(activity mapping)
# Create a scatter plot using the X train pca and the Activity Names
plt.figure(figsize=(8, 6))
for activity in Activity_Name.unique():
    indices = Activity Name == activity
    plt.scatter(X_train_pca[indices, 0], X train pca[indices, 1],
label=activity, alpha=0.6)
# TODO <--code below-->
plt.title('PCA Visualization of Training Data')
plt.xlabel('Principal Component 1')
plt.ylabel('Principal Component 2')
plt.legend(title='Activity')
plt.show()
```

PCA Visualization of Training Data



Kmeans Clustering and The Optimal Number of Clusters

1. Elbow Method

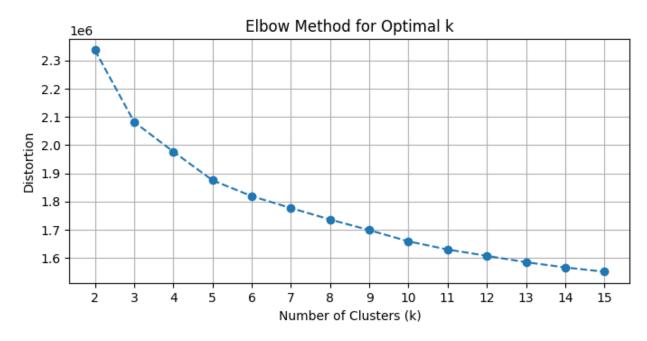
```
from sklearn.cluster import KMeans

# Elbow Method

distortion_values = []
for k in range(2, 16):
    kmeans = KMeans(n_clusters=k, random_state=42, n_init=10)
    kmeans.fit(X_train_scaled)
    distortion_values.append(kmeans.inertia_)

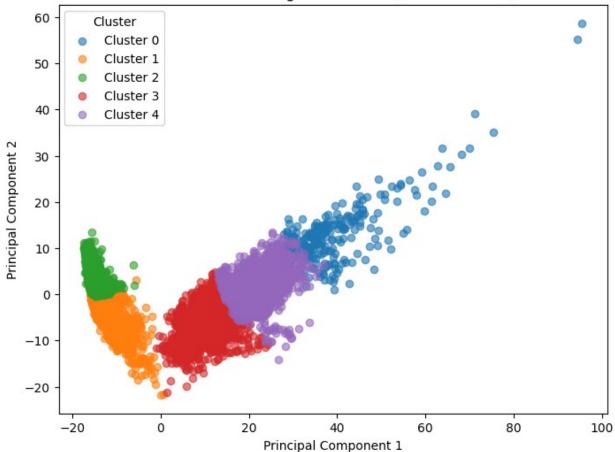
# Plotting the Elbow Method
plt.figure(figsize=(8, 3.5))
plt.plot(range(2, 16), distortion_values, marker='o', linestyle='--')
plt.title('Elbow Method for Optimal k')
plt.xlabel('Number of Clusters (k)')
plt.ylabel('Distortion')
plt.xticks(range(2, 16))
```

```
plt.grid()
plt.show()
```



```
# Choose k based on the elbow method
elbow k = 5
kmeans elbow = KMeans(n clusters=elbow k, random state=42, n init=10)
clusters elbow = kmeans elbow.fit predict(X train scaled)
# PCA for visualization - already computed PCA
X_train_pca_elbow = X_train_pca
# Plotting the clusters
plt.figure(figsize=(8, 6))
for cluster in range(elbow k):
    indices = clusters elbow == cluster
    plt.scatter(X train pca elbow[indices, 0],
X_train_pca_elbow[indices, 1], label=f'Cluster {cluster}', alpha=0.6)
plt.title('K-Means Clustering Visualization (Elbow Method)')
plt.xlabel('Principal Component 1')
plt.ylabel('Principal Component 2')
plt.legend(title='Cluster')
plt.show()
```

K-Means Clustering Visualization (Elbow Method)



Explain distortion with Increasing K

- The optimal number of cluser is 5.
 - Distortion measures the sum of the squared distance between the data points and the assigned clustes this means an increase in the clusters leads to the decrease of the the distortion. This is mainly because increaseing clusters allows the centroing to better fit the data which reduces the distance between points and their nearest centroid. However, as K becomes large the distortion the decreasing rate also diminishes which leads to the elbow point where addding more clustersprovides the diminishing returns in reducing distortion.

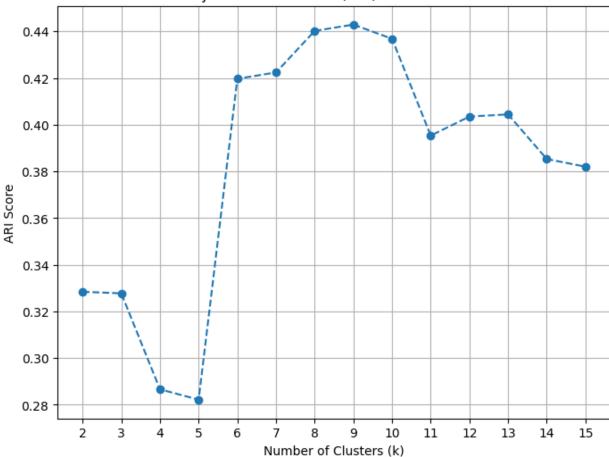
2. Adjusted Rand Index (ARI)

```
from sklearn.metrics import adjusted_rand_score
# 2. Adjusted Rand Index (ARI)
ari_scores = []
for k in range(2, 16):
    kmeans = KMeans(n_clusters=k, random_state=42)
    clusters = kmeans.fit_predict(X_train_scaled)
```

```
ari = adjusted rand score(y train[0], clusters)
    ari scores.append(ari)
# Plotting ARI Scores
plt.figure(figsize=(8, 6))
plt.plot(range(2, 16), ari scores, marker='o', linestyle='--')
plt.title('Adjusted Rand Index (ARI) for Different k')
plt.xlabel('Number of Clusters (k)')
plt.ylabel('ARI Score')
plt.xticks(range(2, 16))
plt.grid()
plt.show()
/usr/local/lib/python3.10/dist-packages/sklearn/cluster/
kmeans.py:870: FutureWarning: The default value of `n init` will
change from 10 to 'auto' in 1.4. Set the value of `n init` explicitly
to suppress the warning
 warnings.warn(
/usr/local/lib/python3.10/dist-packages/sklearn/cluster/ kmeans.py:870
: FutureWarning: The default value of `n_init` will change from 10 to
'auto' in 1.4. Set the value of `n_init` explicitly to suppress the
warning
 warnings.warn(
/usr/local/lib/python3.10/dist-packages/sklearn/cluster/_kmeans.py:870
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/usr/local/lib/python3.10/dist-packages/sklearn/cluster/ kmeans.py:870
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'auto' in 1.4. Set the value of `n init` explicitly to suppress the
warning
 warnings.warn(
/usr/local/lib/python3.10/dist-packages/sklearn/cluster/ kmeans.py:870
: FutureWarning: The default value of `n_init` will change from 10 to
```

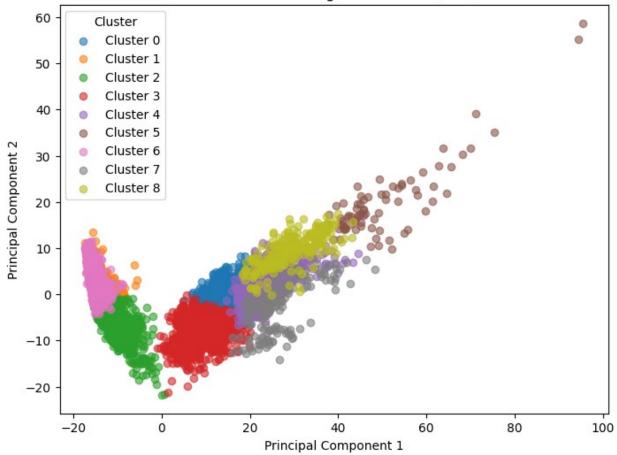
```
'auto' in 1.4. Set the value of `n init` explicitly to suppress the
warning
 warnings.warn(
/usr/local/lib/python3.10/dist-packages/sklearn/cluster/ kmeans.py:870
: FutureWarning: The default value of `n init` will change from 10 to
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/usr/local/lib/python3.10/dist-packages/sklearn/cluster/_kmeans.py:870
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 warnings.warn(
/usr/local/lib/python3.10/dist-packages/sklearn/cluster/ kmeans.py:870
: FutureWarning: The default value of `n_init` will change from 10 to
'auto' in 1.4. Set the value of `n init` explicitly to suppress the
warning
 warnings.warn(
```





```
# Choose k based on ARI
best ari k = 9 # Based on ARI scores
kmeans ari = KMeans(n clusters=best ari k, random state=42, n init=10)
clusters ari = kmeans ari.fit predict(X train scaled)
# PCA for visualization
X_train_pca_ari = X_train_pca # PCA already computed
# Plotting the clusters
plt.figure(figsize=(8, 6))
for cluster in range(best ari k):
    indices = clusters ari == cluster
    plt.scatter(X train pca ari[indices, 0], X train pca ari[indices,
1], label=f'Cluster {cluster}', alpha=0.6)
plt.title('K-Means Clustering Visualization (ARI)')
plt.xlabel('Principal Component 1')
plt.vlabel('Principal Component 2')
plt.legend(title='Cluster')
plt.show()
```

K-Means Clustering Visualization (ARI)



ARI with increase of k

The optimal was 9.

- ARI measures the similarity between clustering results and the ground truth labels. As the number of clusters increase the following trend is observeds.
 - Plateau: Beyond the optimal K, the Ari scrore declines as the clusters become too granular, this leads reduction of alignment with the ground truth.
 - Fluctuations: As K increases, the ARI fluctuares due to creation of more clustres that do not correspond well to the true data distributions or overfitting.

Prototype Selection using K-means Clustering.

1. Random Selection

```
import numpy as np
import pandas as pd
from sklearn.linear_model import LogisticRegression
from sklearn.metrics import accuracy_score
```

```
def random prototype selection(X, y, n samples):
    Selects a random subset from the data. Train a logistic regression
model
    on the selected data.
   Args:
        X (pd.DataFrame): The input features.
        y (pd.Series): The target labels.
        n samples(int): The number of samples to select from each
class.
    Returns:
        tuple: A tuple containing the selected features (X selected)
and labels (y selected).
    X selected = []
    y selected = []
    # Iterate over each unique class
    for label in np.unique(y):
        # Select random samples for the current class
        indices = np.random.choice(np.where(y == label)[0], n samples,
replace=False)
        X selected.append(X[indices])
        y selected.append(y[indices])
    # Concatenate the selected samples
    X selected = np.vstack(X selected)
    y selected = np.concatenate(y selected)
    return X selected, y selected
n repetitions = 10
accuracies = []
n \text{ samples} = 120
# Calculate the accuracy for the randomly selected prototypes over 10
experiments
for in range(n repetitions):
    X selected, y selected =
random prototype selection(X train scaled, y train[0].values,
n samples)
    # Train Logistic Regression model
    logistic regression = LogisticRegression(random state=42,
\max iter=1000)
    logistic regression.fit(X selected, y selected)
```

```
# Make predictions and calculate accuracy
y_pred = logistic_regression.predict(X_test_scaled)
accuracy = accuracy_score(y_test[0].values, y_pred)
accuracies.append(accuracy)

average_accuracy = np.mean(accuracies)
print(f"Average Accuracy with Random Selection over {n_repetitions}
repetitions: {average_accuracy:.4f}")

Average Accuracy with Random Selection over 10 repetitions: 0.9192
```

2. K-means Clustering by Class

```
# 2. K-means Clustering by Class
def kmeans prototype selection(X, y, n prototypes per class):
    Selects prototypes using K-means clustering for each class.
   Args:
       X (pd.DataFrame): The input features.
        y (pd.Series): The target labels.
        n prototypes per class (int): The number of prototypes to
select from each class.
    Returns:
        pd.DataFrame: The selected prototypes.
        pd.Series: The selected labels.
    #Initialize lists to store selected prototypes and labels
    X selected = [] # List to store selected feature subsets for each
class
    y selected = [] # List to store selected labels for each class
    # Step 1: Iterate over each unique class label in the target
labels
    for label in np.unique(y):
        # Filter data points belonging to the current class
        class data = X.loc[y.values.flatten() == label]
        # Step 2: Cluster the points using K-means with k =
n prototypes per class
        kmeans = KMeans(n clusters=n prototypes per class,
random state=42)
        kmeans.fit(class data)
        # Step 3: Find the closest points to each centroid
        centroids = kmeans.cluster centers
```

```
for centroid in centroids:
            distances = np.linalg.norm(class data.values - centroid,
axis=1)
            closest point idx = np.argmin(distances)
X selected.append(class data.iloc[closest point idx].values)
            y selected.append(label)
    # Convert lists to numpy arrays
    X selected = np.array(X selected)
    y selected = np.array(y selected)
    return X selected, y selected
# Select prototypes using K-means
y_train = y_train.rename(columns={0: 'Activity'})
X train selected kmeans, y train selected kmeans =
kmeans prototype selection(X train, y train['Activity'], 20)
# Train Logistic Regression model
logistic regression kmeans = LogisticRegression(random state=42,
max iter=1000)
logistic regression kmeans.fit(X train selected kmeans,
y train selected kmeans)
# Make predictions and calculate accuracy
y pred kmeans = logistic regression kmeans.predict(X test)
accuracy kmeans = accuracy score(y test, y pred kmeans)
print(f"Accuracy with K-means Selection: {accuracy kmeans:.4f}")
/usr/local/lib/python3.10/dist-packages/sklearn/cluster/
kmeans.py:870: FutureWarning: The default value of `n init` will
change from 10 to 'auto' in 1.4. Set the value of `n init` explicitly
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 warnings.warn(
/usr/local/lib/python3.10/dist-packages/sklearn/cluster/ kmeans.py:870
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 warnings.warn(
/usr/local/lib/python3.10/dist-packages/sklearn/cluster/ kmeans.py:870
```

```
: FutureWarning: The default value of `n_init` will change from 10 to 'auto' in 1.4. Set the value of `n_init` explicitly to suppress the warning
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: FutureWarning: The default value of `n_init` will change from 10 to 'auto' in 1.4. Set the value of `n_init` explicitly to suppress the warning
  warnings.warn(

Accuracy with K-means Selection: 0.8931
```

Comparizon of the Random selection and the K-meands clusering by class model Accuracy

Random Selection: Average Accuracy with Random Selection over 10 repetitions: 0.9192.

Using K-means clustering by Class: Accuracy with K-means Selection: 0.8931

The Random prototype selection had a higher Accuracy of 92.23% compared to the K-means clustering that had 90.57%, this could be becaue Random seelction relies on randomly selected prototypes that may capture diverse sample from each class while K-means clustering selects the prototype based on clusters that ensure that the selected samples are representative of the cluser centroids however, this may miss some outlier or diverse samples. And since the later focuess more on centroirds rather than capturing the full variablity of the data this may be the reason of the less performance of the models compared to the reandom selection model.

Autoencoder for Features Learning.

####1. Data Preparation:

```
import glob
import numpy as np

# Load data with proper tensor formatting
def load_inertial_data(path):
    files = glob.glob(path)
    data_dict = {}
    for f in files:
        name = f.split('/')[-1][:-4]
        # Read as numpy array and convert to float32
        data_dict[name] = pd.read_csv(f, sep='\s+', header=None).values.astype(np.float32)
    return data_dict
```

```
# Load training data
train data = load inertial data("UCI HAR Dataset/train/Inertial
Signals/*.txt")
train labels = pd.read csv("UCI HAR Dataset/train/y train.txt",
header=None)[0].values
# Load Test data
test data = load inertial data("UCI HAR Dataset/test/Inertial
Signals/*.txt")
test labels = pd.read csv("UCI HAR Dataset/test/y test.txt",
header=None)[0].values
print(f"Train Data Dictionary keys: {list(train data.keys())}")
print(f"For each sensor the Data shape:
{train data['body acc x train'].shape}")
Train Data Dictionary keys: ['body_acc_z_train', 'body_gyro_z_train',
'total_acc_z_train', 'body_acc_y_train', 'total_acc_y_train', 'body_acc_x_train', 'total_acc_x_train', 'body_gyro_x_train',
'body gyro y train']
For each sensor the Data shape: (7352, 128)
from torch.utils.data import Dataset, DataLoader
import torch
# Create PyTorch Dataset
class SensorsDataset(Dataset):
    def __init__(self, data_dict, labels):
        # Stack all signals along the feature dimension Shape:
(num samples, 128, num features)
        self.data = torch.tensor(
            np.stack([data dict[key] for key in data dict.keys()],
axis=-1), dtype=torch.float32
        ) # Shape: (num samples, 128, num features)
        self.labels = torch.tensor(labels, dtype=torch.long)
    def len_(self):
        return len(self.data)
    def __getitem__(self, idx):
        return self.data[idx], self.labels[idx]
# Create dataset and dataloader
train dataset = SensorsDataset(train data, train labels)
train loader = DataLoader(train dataset, batch size=32, shuffle=True)
# Verify shapes
sample, label = next(iter(train loader))
```

```
print(f"Input shape: {sample.shape}") # Should be (batch_size, 128,
9)
print(f"Label shape: {label.shape}") # Should be (batch_size)

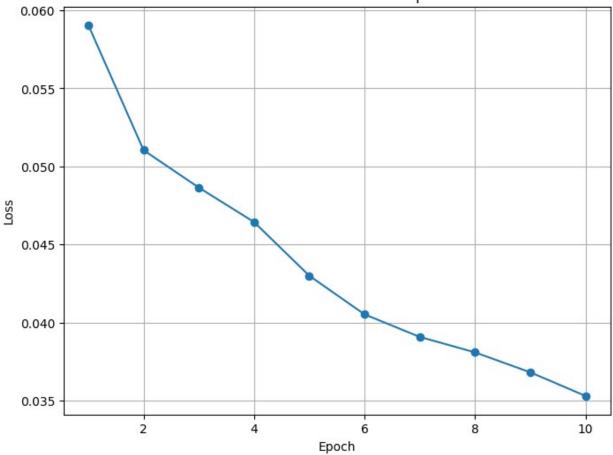
Input shape: torch.Size([32, 128, 9])
Label shape: torch.Size([32])
```

2. Autoencoder Implementation

```
import torch.nn as nn
import torch.optim as optim
# 2. Autoencoder Implementation
class TimeSeriesAE(nn.Module):
    def __init__(self, input_size=9, hidden_size=64, encoding_dim=64):
        super().__init__()
        # Encoder
        self.encoder = nn.GRU(input size, hidden size,
batch first=True, bidirectional=True)
        self.enc fc = nn.Linear(hidden size * 2, encoding dim)
        # Decoder
        self.dec fc = nn.Linear(encoding dim, hidden size * 2)
        self.decoder = nn.GRU(hidden size * 2, hidden size,
batch first=True, bidirectional=True)
        self.output layer = torch.nn.Linear(hidden size * 2,
input size)
    def forward(self, x):
        , hidden = self.encoder(x)
        hidden = torch.cat([hidden[-2], hidden[-1]], dim=1) # Combine
bidirectional
        encoded = self.enc fc(hidden)
        # Decoding
        decoded = self.dec fc(encoded).unsqueeze(1).repeat(1,
x.size(1), 1)
        out, = self.decoder(decoded)
        reconstructed = self.output layer(out)
        return reconstructed, encoded
# Instantiate the model
input size = 9 # Number of features
hidden size = 64
model = TimeSeriesAE(input size, hidden size)
criterion = nn.MSELoss()
optimizer = torch.optim.Adam(model.parameters(), lr=0.001)
```

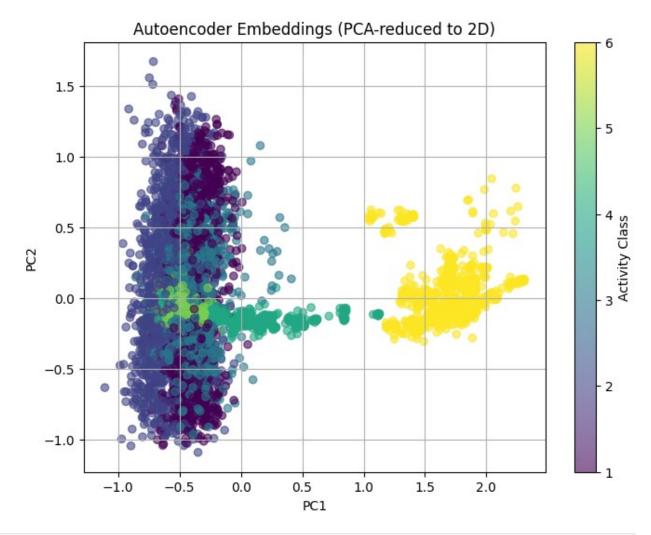
```
# Train loop for the autoencoder
loss history = []
num epochs = 10
for epoch in range(num epochs):
    model.train()
    total_loss = 0
    for batch_X, _ in train_loader:
    optimizer.zero_grad()
        recon, \_ = model(batch X)
        loss = criterion(recon, batch X)
        loss.backward()
        optimizer.step()
        total loss += loss.item()
    avg_loss = total_loss / len(train_loader)
    loss history.append(avg loss)
    print(f"Epoch {epoch+1}/{num epochs}, Loss: {avg loss:.4f}")
# Plotting the reconstruction loss vs epoch
plt.figure(figsize=(8, 6))
plt.plot(range(1, num epochs + 1), loss history, marker='o')
plt.title("Reconstruction Loss Over Epochs")
plt.xlabel("Epoch")
plt.vlabel("Loss")
plt.grid(True)
plt.show()
Epoch 1/10, Loss: 0.0591
Epoch 2/10, Loss: 0.0510
Epoch 3/10, Loss: 0.0486
Epoch 4/10, Loss: 0.0464
Epoch 5/10, Loss: 0.0430
Epoch 6/10, Loss: 0.0405
Epoch 7/10, Loss: 0.0391
Epoch 8/10, Loss: 0.0381
Epoch 9/10, Loss: 0.0368
Epoch 10/10, Loss: 0.0353
```





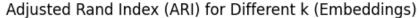
```
# 3. Embedding Extraction and Visualization
ae loader = DataLoader(train dataset, batch size=32, shuffle=False)
# Extract embeddings for the training data
model.eval()
embeddings = []
train labels = []
with torch.no_grad():
    for batch_X, labels in ae_loader:
        _, encoded = model(batch X)
        embeddings.append(encoded.numpy())
        train labels.extend(labels.numpy())
embeddings = np.concatenate(embeddings, axis=0)
train labels = np.array(train labels)
# Reduce embeddings to 2D using PCA for visualization
from sklearn.decomposition import PCA
pca = PCA(n components=2)
embeddings\overline{2}d = pca.fit_transform(embeddings)
```

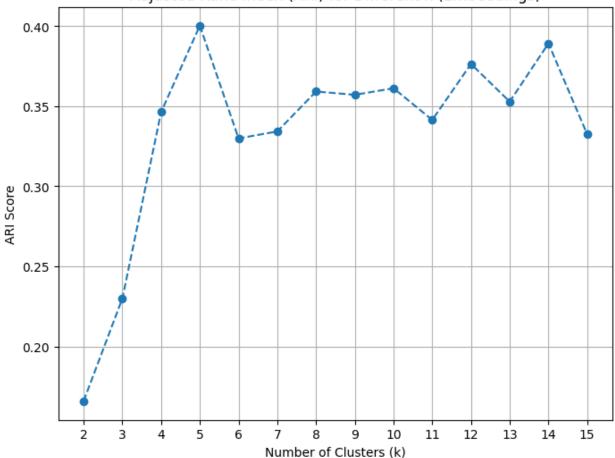
```
# Create a scatter plot of the 2D embeddings
plt.figure(figsize=(8, 6))
activities = np.unique(train_labels)
scatter = plt.scatter(embeddings_2d[:, 0], embeddings_2d[:, 1],
c=train_labels, cmap='viridis', alpha=0.6)
plt.colorbar(scatter, label='Activity Class')
plt.title('Autoencoder Embeddings (PCA-reduced to 2D)')
plt.xlabel("PC1")
plt.ylabel("PC2")
plt.grid(True)
plt.show()
```



```
from sklearn.metrics import adjusted_rand_score
# Adjusted Rand Index (ARI) for the embeddings
ari_scores = []
```

```
for k in range(2, 16):
    kmeans = KMeans(n clusters=k, random state=42, n init=10)
    clusters = kmeans.fit_predict(embeddings) # Use embeddings for
clusterina
    ari = adjusted rand score(train labels, clusters) # Compare with
ground truth labels
    ari scores.append(ari)
# Plotting ARI Scores
plt.figure(figsize=(8, 6))
plt.plot(range(2, 16), ari_scores, marker='o', linestyle='--')
plt.title('Adjusted Rand Index (ARI) for Different k (Embeddings)')
plt.xlabel('Number of Clusters (k)')
plt.ylabel('ARI Score')
plt.xticks(range(2, 16))
plt.grid()
plt.show()
```





```
# Choose k based on ARI
best embedd ari k = 5 # Based on ARI scores or other criteria
kmeans_ari = KMeans(n_clusters=best_embedd_ari_k, random_state=42,
n init=10
clusters ari = kmeans ari.fit predict(embeddings) # Use embeddings
for clustering
# PCA for visualization
pca = PCA(n components=2) # Reduce to 2 dimensions for visualization
X_train_pca_ari = pca.fit transform(embeddings)
# Plotting the clusters
plt.figure(figsize=(8, 6))
for cluster in range(best embedd ari k):
    indices = clusters ari == cluster
   plt.scatter(X_train_pca_ari[indices, 0], X_train_pca_ari[indices,
1], label=f'Cluster {cluster}', alpha=0.6)
plt.title('K-Means Clustering Visualization (Embeddings)')
plt.xlabel('Principal Component 1')
plt.ylabel('Principal Component 2')
plt.legend(title='Cluster')
plt.show()
```

