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# Step 1: Import necessary libraries
                                     # For loading and handling the
import pandas as pd
dataset
                                     # For numerical operations
import numpy as np
from sklearn.model selection import train test split # To split the
from sklearn.ensemble import RandomForestClassifier # Our base
model
                                                     # To evaluate
from sklearn.metrics import accuracy score
from sklearn.preprocessing import LabelEncoder
                                                     # To convert
strings to numbers
# Step 2: Load the dataset
data = pd.read csv('/content/healthcare-dataset-stroke-data.csv')
# Step 3: Drop rows with missing values
data = data.dropna() # Removes rows that have missing (NaN) values
# Step 4: Encode categorical columns
label encoders = {}
categorical cols = ['gender', 'ever married', 'work type',
'Residence type', 'smoking status']
for col in categorical cols:
   le = LabelEncoder()
   data[col] = le.fit transform(data[col]) # Convert text to numbers
   label encoders[col] = le
# Step 5: Define features (X) and label (y)
X = data.drop(columns=['id', 'stroke']) # Features: all columns
except 'id' and 'stroke'
y = data['stroke']
                                        # Target variable: stroke (0
or 1)
# Step 6: Split data into labeled and unlabeled
# Assume only 10% is labeled initially
X_labeled, X_unlabeled, y_labeled, y_unlabeled = train test split(
   X, y, test size=0.90, stratify=y, random state=42
# Step 7: Train model on labeled data
model = RandomForestClassifier(random state=42)
model.fit(X labeled, y labeled) # Train only on the small labeled
part
RandomForestClassifier(random state=42)
# Step 8: Predict probabilities on the unlabeled data
probs = model.predict proba(X unlabeled) # Predict class
probabilities
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# Step 9: Select confident predictions (confidence > 0.9)
confident indices = []
pseudo labels = []
for i, prob in enumerate(probs):
    if max(prob) > 0.9:
                                                # If model is very
confident
        confident indices.append(i)
                                                # Save index
        pseudo labels.append(np.argmax(prob)) # Save predicted label
# Step 10: Add confident pseudo-labeled data to the labeled set
X confident = X unlabeled.iloc[confident indices]
y confident = np.array(pseudo labels)
# Combine with original labeled data
X labeled = pd.concat([X labeled, X confident])
y labeled = np.concatenate([y labeled, y confident])
# Remove used unlabeled samples
X unlabeled = X unlabeled.drop(X unlabeled.index[confident indices])
# Step 11: Retrain model on the updated labeled dataset
model.fit(X labeled, y labeled)
RandomForestClassifier(random state=42)
# Step 12: Evaluate on a hold-out test set (optional, for accuracy
# Let's split a separate test set from the original data
X train, X test, y train, y test = train test split(
    X, y, test size=0.2, stratify=y, random state=42
)
model.fit(X train, y train)
y pred = model.predict(X test)
print("Accuracy on test data:", accuracy_score(y_test, y_pred))
Accuracy on test data: 0.9500978473581213
import matplotlib.pyplot as plt
# Prepare to repeat self-training for graph
X labeled iter = X labeled.copy()
y labeled_iter = y_labeled.copy()
X unlabeled iter = X unlabeled.copy()
# Lists to store data for plotting
accuracy list = []
pseudo_count_list = []
iterations = []
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# Fixed test set for evaluation
X train eval, X test eval, y train eval, y test eval =
train_test split(
   X, y, test size=0.2, stratify=y, random state=42
# Self-training loop with graph data collection
for iteration in range(1, 11): # up to 10 rounds
   model = RandomForestClassifier(random state=42)
   model.fit(X labeled iter, y labeled iter)
   y pred = model.predict(X test eval)
   accuracy = accuracy score(y test eval, y pred)
   # Predict on unlabeled
   probs = model.predict proba(X unlabeled iter)
    confident indices = []
   pseudo labels = []
   for i, prob in enumerate(probs):
        if max(prob) > 0.9: # confidence threshold
           confident indices.append(i)
           pseudo labels.append(np.argmax(prob))
   if not confident indices:
        break # Stop if no more confident predictions
   # Save graph info
   accuracy list.append(accuracy)
    pseudo count list.append(len(confident indices))
   iterations.append(iteration)
   # Add confident pseudo-labeled data
   X confident = X unlabeled iter.iloc[confident indices]
   y confident = np.array(pseudo labels)
   X labeled iter = pd.concat([X labeled iter, X confident])
   y labeled iter = np.concatenate([y labeled iter, y confident])
   X unlabeled iter =
X unlabeled iter.drop(X unlabeled iter.index[confident indices])
plt.figure(figsize=(10, 6))
plt.plot(iterations, accuracy list, marker='o', color='blue',
label='Test Accuracy')
plt.bar(iterations, pseudo count list, alpha=0.3, color='orange',
label='Pseudo-labeled added')
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plt.title("Self-Training Learning: Accuracy & Pseudo-Labels Over
Iterations")
plt.xlabel("Iteration")
plt.ylabel("Accuracy / Pseudo-Labels")
plt.legend()
plt.grid(True)
plt.tight_layout()
plt.show()
```

