assignment-5

September 15, 2024

1 TASK 1

```
[1]: import pandas as pd
  import numpy as np
  from sklearn.model_selection import train_test_split
  from sklearn.preprocessing import OneHotEncoder, LabelEncoder
  from sklearn.tree import DecisionTreeClassifier
  from sklearn.linear_model import LogisticRegression
  from sklearn.neighbors import KNeighborsClassifier
  from sklearn.metrics import accuracy_score
  import matplotlib.pyplot as plt
```

1.1 Importing dataset

```
[2]: from ucimlrepo import fetch_ucirepo
# fetch dataset
nursery = fetch_ucirepo(id=76)
# data (as pandas dataframes)
X = nursery.data.features
y = nursery.data.targets
```

```
[3]: # Converting 'spec_prior' and 'very_recom' to 'recommend' class to make a 3_ 
class problem

y = y.replace({'spec_prior': 'recommend', 'very_recom' : 'recommend'})
```

```
[4]: # Label Encoding for both X and y
label_encoder = LabelEncoder()
y_label = label_encoder.fit_transform(y)
X_label = pd.DataFrame()
for i in X.columns:
    X_label[i] = label_encoder.fit_transform(X[i])
```

c:\Users\sakth\AppData\Local\Programs\Python\Python312\Lib\sitepackages\sklearn\preprocessing_label.py:114: DataConversionWarning: A columnvector y was passed when a 1d array was expected. Please change the shape of y
to (n_samples,), for example using ravel().

```
y = column_or_1d(y, warn=True)
```

1.2 1. Decision Tree (categorical features)

Here I have used label encoding for both x and y. In other models, I have used one hot encoding for X and label encoding for y.

```
[5]: # Decision Tree model
     def decision_tree(X_train, y_train, X_val, y_val, X_test, y_test):
         param_grid_lr = {'max_depth': [1, 3, 5, 10]}
         best_model = None
         best accuracy = 0
         for i in param_grid_lr['max_depth']:
             mdl_dtc = DecisionTreeClassifier(max_depth=i, random_state=int(100*i))
             mdl_dtc.fit(X_train, y_train)
             y_pred = mdl_dtc.predict(X_val)
             accuracy = accuracy_score(y_val, y_pred)
             if accuracy > best_accuracy:
                 best_accuracy = accuracy
                 best_model = mdl_dtc
         # Print the best model and its best accuracy
         y_pred = best_model.predict(X_test)
         test_accuracy = 100 * accuracy_score(y_test, y_pred)
         return test_accuracy
```

96.6820987654321

```
[7]: # Preprocess data (one-hot encoding)
encoder = OneHotEncoder()
X_encoded = encoder.fit_transform(X)
```

1.3 2. Decision Tree(categorical features in one-hot encoded form)

```
[8]: # Decision Tree (one-hot encoded form) model

def decision_tree_one_hot_encoded_form(X_train, y_train, X_val, y_val, X_test,

□ y_test):

param_grid_lr = {'max_depth': [1, 5, 10, 20, 30]}

best_model = None
```

```
best_accuracy = 0
for i in param_grid_lr['max_depth']:
    mdl_dtc = DecisionTreeClassifier(max_depth=i, random_state=int(100*i))
    mdl_dtc.fit(X_train, y_train)
    y_val_pred = mdl_dtc.predict(X_val)
    accuracy = accuracy_score(y_val, y_val_pred)
    if accuracy > best_accuracy:
        best_accuracy = accuracy
    best_model = mdl_dtc

# Print the best model and its test accuracy
y_pred = best_model.predict(X_test)
test_accuracy = 100 * accuracy_score(y_test, y_pred)
return test_accuracy
```

99.26697530864197

1.4 3. Logistic Regression with L1 regularizer

```
[10]: # Logistic regression model with L1 regularization
      def logistic_regression_with_L1(X_train, y_train, X_val, y_val, X_test, y_test):
          param_grid_lr = {'C': [0.01, 0.1, 1, 10, 100]}
          best_model = None
          best accuracy = 0
          for i in param_grid_lr['C']:
              mdl_lr = LogisticRegression(penalty='12', solver='liblinear', C=i)
              mdl_lr.fit(X_train, y_train)
              y_pred = mdl_lr.predict(X_val)
              accuracy = accuracy_score(y_val, y_pred)
              if accuracy > best_accuracy:
                  best_accuracy = accuracy
                  best_model = mdl_lr
          # Print the best model and its test accuracy
          y_pred = best_model.predict(X_test)
          test_accuracy = 100 * accuracy_score(y_test, y_pred)
          return test_accuracy
```

91.62808641975309

1.5 4. k-Nearest Neighbors

```
[12]: # k-Nearest Neighbors model
      def k_nearest_neighbors(X_train, y_train, X_val, y_val, X_test, y_test):
          param_grid_lr = {'k': [3, 5, 7, 9, 11]}
          best_model = None
          best_accuracy = 0
          for i in param grid lr['k']:
              mdl_knn = KNeighborsClassifier(n_neighbors=i)
              mdl_knn.fit(X_train, y_train)
              y_pred = mdl_knn.predict(X_val)
              accuracy = accuracy score(y val, y pred)
              if accuracy > best_accuracy:
                  best_accuracy = accuracy
                  best_model = mdl_knn
          # Print the best model and its test accuracy
          y_pred = best_model.predict(X_test)
          test_accuracy = 100 * accuracy_score(y_test, y_pred)
          return test_accuracy
```

95.25462962962963

1.6 Performance of all 4 models calculated

Accuracy of Decision Tree model: 96.6820987654321

Accuracy of Decision Tree model with one-hot encoding: 99.26697530864197

Accuracy of Logistic Regression with L1 regularization model: 91.62808641975309

Accuracy of k-Nearest Neighbors model: 95.25462962963

To compute varaince and mean, I have repeated the task 5 times for each model and stored all the accuracies in a dictionary

```
[15]: # Creating a dictionary to store each values of accuracy for each model
     accuracy = {'accuracy_dst_1': [], 'accuracy_dst_2': [], 'accuracy_lr_l1': [], u
     for i in range(5):
        # Data split in label encoded form
        X_train_full, X_test, y_train_full, y_test = train_test_split(X_label,__
      X_train, X_val, y_train, y_val = train_test_split(X_train_full,_
      accuracy['accuracy dst 1'].append(decision_tree(X_train, y_train, X_val,_

y_val, X_test, y_test))
        # Data split in one-hot encoded form
        X_train_full, X_test, y_train_full, y_test = train_test_split(X_encoded,_

y_label, test_size=0.2, stratify=y, random_state=i)
        X_train, X_val, y_train, y_val = train_test_split(X_train_full,_
      wy_train_full, test_size=0.25, stratify=y_train_full, random_state=i)
        accuracy['accuracy_dst_2'].
      →append(decision_tree_one_hot_encoded_form(X_train, y_train, X_val, y_val, __
      →X_test, y_test))
        accuracy['accuracy lr l1'].append(logistic regression with L1(X train, __
      accuracy['accuracy_knn'].append(k_nearest_neighbors(X_train, y_train, u
      →X_val, y_val, X_test, y_test))
     print(accuracy)
```

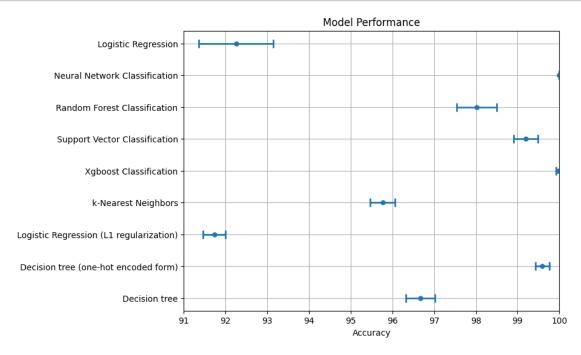
{'accuracy_dst_1': [96.72067901234568, 96.10339506172839, 96.5277777777779, 97.14506172839506, 96.83641975308642], 'accuracy_dst_2': [99.8070987654321, 99.5756172839506, 99.5756172839506, 99.3055555555555556, 99.69135802469135],

```
'accuracy_lr_11': [92.2067901234568, 91.666666666666, 91.7824074074074, 91.43518518518519, 91.55092592592592], 'accuracy_knn': [96.06481481481481, 96.02623456790124, 95.5246913580247, 95.9104938271605, 95.2932098765432]}
```

Next, finding the mean and variance of these individual model accuracies

```
[16]: mean_accuracy = {k: np.mean(v) for k, v in accuracy.items()}
      print(mean_accuracy)
      variance_accuracy = {k: np.var(v) for k, v in accuracy.items()}
      print(variance_accuracy)
     {'accuracy_dst_1': np.float64(96.666666666666), 'accuracy_dst_2':
     np.float64(99.59104938271604), 'accuracy lr l1': np.float64(91.72839506172839),
     'accuracy_knn': np.float64(95.76388888888889)}
     {'accuracy dst 1': np.float64(0.11943206066148507), 'accuracy dst 2':
     np.float64(0.027744436823654968), 'accuracy_lr_l1':
     np.float64(0.07061137783874538), 'accuracy_knn':
     np.float64(0.09192577351013673)}
[17]: # Adding accuracies of other models to the dictionary
      mean_accuracy['accuracy_xgb'] = 99.969
      mean_accuracy['accuracy_svc'] = 99.198
      mean_accuracy['accuracy_rfc'] = 98.025
      mean_accuracy['accuracy_nn'] = 100
      mean_accuracy['accuracy_lr'] = 92.253
      variance_accuracy['accuracy_xgb'] = 0.00216225
      variance accuracy['accuracy svc'] = 0.085849
      variance_accuracy['accuracy_rfc'] = 0.22896225
      variance accuracy['accuracy nn'] = 0.0
      variance accuracy['accuracy lr'] = 0.801025
```

1.7 Visualisation of all 9 methods



2 TASK 2

2.1 PART 1

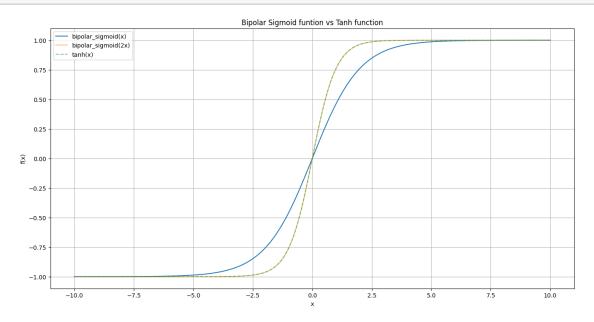
```
[19]: import numpy as np
import matplotlib.pyplot as plt

# Define sigmoid function
def sigmoid(x):
    return 1 / (1 + np.exp(-x))
```

```
# Define bipolar_sigmoid function as a function of sigmoid function
def bipolar_sigmoid(x):
    return 2 * sigmoid(x) - 1
```

2.2 PART 2

```
[20]: # Defining tanh function
      def tanh_fn(x):
          return np.tanh(x)
      # Create x values
      x = np.linspace(-10, 10, 1000)
      # Plotting tanh and bipolar_sigmoid
      plt.figure(figsize=(16, 8))
      plt.plot(x, bipolar_sigmoid(x), label=f'bipolar_sigmoid(x)')
      plt.plot(x, bipolar_sigmoid(2*x), label=f'bipolar_sigmoid(2x)', alpha = 0.5)
      plt.plot(x, tanh_fn(x), '--', label=f'tanh(x)', alpha=0.6)
      plt.title('Bipolar Sigmoid funtion vs Tanh function')
      plt.xlabel('x')
      plt.ylabel('f(x)')
      plt.legend()
      plt.grid(True)
      plt.show()
```



From the graph above, we can see that tanh(x) is equivalent to bipolar_sigmoid(2*x)

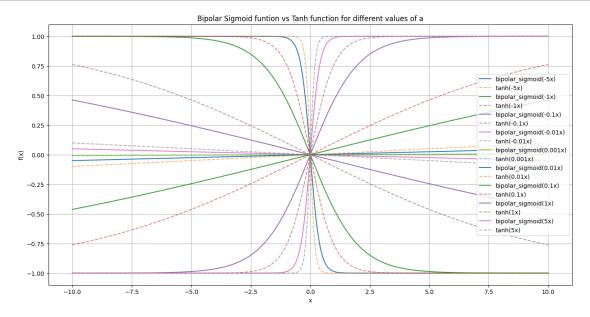
2.3 PART 3

```
[21]: # Define parameter 'a' values
a_values = [-5, -1, -0.1, -0.01, 0.001, 0.01, 0.1, 1, 5]

# Plot bipolar_sigmoid(ax) and tanh(ax) for different values of a
plt.figure(figsize=(16, 8))

for a in a_values:
    plt.plot(x, bipolar_sigmoid(a * x), label=f'bipolar_sigmoid({a}x)')
    plt.plot(x, tanh_fn(a * x), '--', label=f'tanh({a}x)', alpha=0.6)

plt.title('Bipolar Sigmoid funtion vs Tanh function for different values of a')
plt.xlabel('x')
plt.ylabel('f(x)')
plt.legend()
plt.grid(True)
plt.show()
```



2.4 PART 4

```
[22]: def linear_range(a,x):
    y = bipolar_sigmoid(a * x)

# Define the threshold for linearity
    threshold = 0.05
```

```
# Finding derivative of y
    dy_dx = np.gradient(y, x)
    # Find the index where x is closest to 0 as the function is clearly linear.
 \Rightarrownear x = 0
    x0 idx = np.abs(x).argmin()
    # Use the derivative at x = 0 as the reference
    dy_dx_ref = dy_dx[x0_idx]
    # Check for linearity based on deviations from the derivative at x = 0
    linear_idx = np.where(np.abs(dy_dx - dy_dx_ref) < threshold)[0]</pre>
    return linear_idx
x = np.linspace(-100, 100, 100000)
for a in a_values:
    linear_idx = linear_range(a, x)
    if len(linear_idx) > 0:
        print(f"Linear range for a = {a}: {x[linear_idx[0]]} to__
 \hookrightarrow \{x[linear idx[-1]]\}")
```

```
Linear range for a = -5: -0.05500055000550219 to 0.05500055000548798

Linear range for a = -1: -0.6530065300653121 to 0.6530065300652979

Linear range for a = -0.1: -100.0 to 100.0

Linear range for a = 0.01: -100.0 to 100.0

Linear range for a = 0.01: -100.0 to 100.0

Linear range for a = 0.01: -100.0 to 100.0

Linear range for a = 0.1: -100.0 to 100.0

Linear range for a = 1: -0.6530065300653121 to 0.6530065300652979

Linear range for a = 5: -0.05500055000550219 to 0.05500055000548798
```

Clearly, we can see that for small 'a' values, the linearity range is high. Here I have considered x values to range from -100 to 100 so that it avoids arithmetic overflow.