Milestone 4: Model Building

Activity 1: Training the model in multiple algorithms

Now our data is cleaned and it's time to build the model. We can train our data on different algorithms. For this project we are applying four classification algorithms. The best model is saved based on its performance.

Activity 1.1: ANN Model

Building and training an Artificial Neural Network (ANN) using the Keras library with TensorFlow as the backend. The ANN is initialised as an instance of the Sequential class, which is a linear stack of layers. Then, the input layer and two hidden layers are added to the model using the Dense class, where the number of units and activation function are specified. The output layer is also added using the Dense class with a sigmoid activation function. The model is then compiled with the Adam optimizer, binary cross-entropy loss function, and accuracy metric. Finally, the model is fit to the training data with a batch size of 100, 20% validation split, and 100 epochs

```
# Importing the Keras libraries and packages
import tensorflow.keras.models import Sequential
from tensorflow.keras.layers import Dense

# Creating ANN skleton view

classification = Sequential()
classification.add(Dense(30,activation='relu'))
classification.add(Dense(128,activation='relu'))
classification.add(Dense(64,activation='relu'))
classification.add(Dense(32,activation='relu'))
classification.add(Dense(1,activation='sigmoid'))
```

```
[======0.05 4ms/step - loss: 0.1139 - accuracy: 0.9531 - val_loss: 0.2799 - val_accuracy: 0.8906 fg ...

[poch 99/100 26/26 [===========] - 05 3ms/step - loss: 0.1074 - accuracy: 0.9570 - val_loss: 0.2439 - val_accuracy: 0.9570 - val_loss: 0.2572 - val_accuracy: 0.9062 - accuracy: 0.9570 - val_loss: 0.2572 - val_accuracy: 0.9062 - accuracy: 0.9570 - val_loss: 0.2572 - val_accuracy: 0.9062
```

Activity 1.2: Random Forest model

A function named random Forest is created and train and test data are passed as the parameters. Inside the function, Random Forest Classifier algorithm is initialised and training data is passed to the model with .fit() function. Test data is predicted with. predict() function and saved in a new variable. For evaluating the model, a confusion matrix and classification report is done.

Activity 1.3: Decision tree model

A function named decision Tree is created and train and test data are passed as the parameters. Inside the function, Decision Tree Classifier algorithm is initialised and training data is passed to the model with the .fit() function. Test data is predicted with. predict() function and saved in a new variable. For evaluating the model, a confusion

matrix and classification report is done.

```
from sklearn.tree import DecisionTreeClassifier
dtc = DecisionTreeClassifier(max_depth~4,splitter~'best',criterion~'entropy')

dtc.fit(x_train,y_train)

DecisionTreeClassifier(criterion~'entropy', max_depth~4)

y_predict~ dtc.predict(x_test)
y_predict

array([0, 0, 0, 0, 1, 0, 0, 0, 1, 0, 0, 0, 1, 1, 0, 0, 0, 1, 1, 0, 1, 1, 0, 1, 0, 1, 0, 0, 1, 0, 0, 0, 1, 0, 0, 0, 1, 0, 0, 0, 1, 0, 0, 0, 0, 1, 0, 0, 0, 0, 1, 0, 0, 0, 0, 1, 0, 0, 0, 0, 1, 0, 0, 0, 0, 1, 0, 0, 0, 1, 0, 0, 0, 1, 0, 0, 0, 1, 0, 0, 0, 1, 0, 0, 0, 1, 0, 0, 0, 1, 0, 0, 0, 1, 0, 0, 0, 1, 0, 0, 0, 1, 0, 0, 0, 1, 0, 0, 0, 0, 1, 0])

y_predict_train = dtc.predict(x_train)
```

Activity 1.4: Logistic Regression

```
from sklearn.linear_model import LogisticRegression
lgr = LogisticRegression()
lgr.fit(x_train,y_train)

C:\Users\Saumya\Anaconda3\lib\site-packages\sklearn\utils\validation.py:72: DataConversionMart
Please change the shape of y to (n_samples, ), for example using ravel().
return f(**Nowargs)

LogisticRegression()

Predicting our output with the model which we build

from sklearn.metrics import accuracy_score,classification_report
y_predict = lgr.predict(x_test)
```

Activity 2: Testing the model

In ANN we first have to save the model to the test the inputs

This code defines a function named "predict_exit" which takes in a sample_value as an input. The function then converts the input sample_value from a list to a numpy array. It reshapes the sample_value array as it contains only one record. Then, it applies feature scaling to the reshaped sample_value array using a scaler object 'sc' that should have been previously defined and fitted. Finally, the function returns the prediction of the classifier on the scaled sample_value.

```
def predict_exit(sample_value):
    ## Convert list to numpy array
    | sample_value = np.array(sample_value)

## Reshape because sample_value contains only 1 record
    | sample_value = sample_value.reshape(1, -1)

## Feature Scaling
    | sample_value = sc.transform(sample_value)

return classifier.predict(sample_value)

test=classification.predict([[1,1,121.000000,36.0,0,0,1,0]]))
    if test==1:
        print('Prediction: High chance of CKD1')
else:
        print('Prediction: Low chance of CKD.')

**Prediction: Low chance of CKD. ***

**Prediction: Low chance of CKD. **

**Prediction: Low chance of CKD. ***

**Prediction: Low chance of CKD. **

**Prediction: Low chance of CKD. ***

**Prediction: Low chance of CKD. **

**Prediction: Lo
```