

**A Project Report**

on

**Heart Disease Detection Using Machine Learning**

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**May, 2023**

**(PCS23-15)**

**\_DECLARATION\_**

We affirm that this submission is entirely our original creation and, as far as we know and believe, it does not include any content that has been previously published or authored by someone else. Furthermore, it does not incorporate material that has been substantially accepted for the attainment of any other academic degree or diploma from the university or any other institution of higher education, except where proper credit has been acknowledged within the text.

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I hereby confirm that the Project Report titled "Utilizing Machine Learning for Heart Disease Detection," authored by Dhirendra Yadav, Utkarsh Singh, Devesh Chandra and Dharmendra Singh Yadav, is a genuine representation of their independent work towards fulfilling the requirements for the B. Tech. degree in the Department of Computer Science & Engineering at Dr. A.P.J. Abdul Kalam Technical University, Lucknow. I have personally supervised their progress throughout this endeavor. The contents of this report are authentic and have not been previously presented for the attainment of any other academic qualification

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**\_ACKNOWLEDGEMENT\_**

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**\_ABSTRACT\_**

Machine Learning has widespread applications worldwide, including in the healthcare sector. It can be instrumental in predicting various conditions such as locomotor disorders and heart diseases. These predictions, when made ahead of time, can offer valuable insights to doctors, enabling them to adjust their diagnoses and treatments on a per-patient basis. Our research focuses on utilizing Machine Learning algorithms to predict potential heart diseases in individuals. We conduct a comparative analysis of classifiers, including Decision trees, Logistic Regression, and K-Nearest Neighbor. Additionally, we propose an ensemble classifier that combines multiple approaches to achieve hybrid classification.

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**\_LIST OF ABBREVIATIONS\_**

**ML** : Machine Learning

**AI** : Artificial Intelligence

**NN** : Neural Network

**SVM** : Support Vector Machine

**XG** : Extreme Gradient

**EXANG** : Exercise induced angina

**CHAPTER 1**

**INTRODUCTION**

* 1. **Introduction**

The World Health Organization reports that heart disease is responsible for 12 million deaths annually on a global scale. It is a significant contributor to both illness and mortality worldwide. Analyzing data to predict cardiovascular disease is considered a crucial area of study. Over the past few years, the burden of cardiovascular disease has been steadily increasing worldwide. Numerous research studies have been conducted to identify the most influential factors contributing to heart disease and to develop accurate risk prediction models. Heart disease is often referred to as a silent killer, as it can lead to a person's demise without obvious symptoms. Timely detection of heart disease plays a critical role in making informed decisions regarding lifestyle modifications for individuals at high risk, ultimately reducing complications.

Machine learning has demonstrated its effectiveness in aiding decision-making and forecasting within the healthcare sector by processing the substantial volume of data generated. The objective of this project is to employ machine learning algorithms to analyze patient data, thereby predicting future occurrences of Heart Disease. Utilizing machine learning techniques can prove advantageous in this context. While heart disease may manifest in various forms, there exists a common set of fundamental risk factors that determine an individual's susceptibility to the condition. By gathering data from diverse sources, organizing it into relevant categories, and subsequently conducting thorough analysis, it can be inferred that this approach is well-suited for heart disease prediction.

* 1. **Motivation of Work**

The primary objective of this research is to develop a heart disease prediction model that can accurately predict the occurrence of heart disease. Additionally, the study aims to determine the most effective classification algorithm for identifying the likelihood of heart disease in patients. To justify this research, a comparative analysis was conducted using three commonly employed classification algorithms: Naïve Bayes, Decision Tree, and Random Forest. Given the critical nature of heart disease prediction and the need for utmost accuracy, these algorithms were extensively evaluated using various levels and types of assessment strategies. The findings from this research will offer valuable insights to researchers and medical professionals, enabling them to establish improved approaches in this field.

* 1. **Problem Statement**

Detecting heart disease poses a significant challenge due to various factors. Existing tools capable of predicting heart disease are either prohibitively expensive or lack the necessary efficiency to accurately assess the risk in individuals. Timely identification of cardiac ailments plays a crucial role in reducing mortality rates and minimizing complications. However, constant monitoring of patients is often unfeasible, and round-the-clock doctor consultation is limited by the need for extensive knowledge, time, and expertise. Fortunately, with the abundance of data available today, we can leverage diverse machine learning algorithms to scrutinize this data for concealed trends. These hidden patterns can be harnessed to facilitate health diagnosis in the field of medicine.

**CHAPTER 2**

**LITERATURE REVIEW**

With growing development in the field of medical science alongside machine learning various experiments and research has been carried out in these recent years releasing the relevant significant papers.

[1] A study conducted by Purushottam et al. introduced a research paper titled "Efficient Heart Disease Prediction System." The authors employed hill climbing and decision tree algorithms in their approach. The Cleveland dataset was utilized, and data preprocessing was performed before applying the classification algorithms. To handle missing values in the dataset, the authors employed Knowledge Extraction based on Evolutionary Learning (KEEL), an open-source data mining tool. The decision tree was constructed in a top-down manner, with the hill-climbing algorithm selecting nodes at each level based on a test. The parameters used in the study included confidence, with a minimum value of 0.25. The system achieved an accuracy rate of approximately 86.7%.

[2] In their study titled "Heart Disease Prediction Using Machine Learning Algorithms," Santhana Krishnan. J, et al. introduced a research paper that employed decision tree and Naive Bayes algorithms to predict heart disease. The decision tree algorithm constructs a tree structure based on specific conditions, enabling True or False decisions. Other algorithms such as SVM and KNN utilize vertical or horizontal split conditions depending on the dependent variables. The decision tree algorithm, characterized by its tree-like structure with root nodes, leaves, and branches, also aids in comprehending the significance of attributes within the dataset. The researchers utilized the Cleveland dataset, dividing it into a 70% training set and a 30% testing set using certain methods. This algorithm achieved a 91% accuracy rate. The second algorithm employed was Naive Bayes, which is well-suited for classification tasks and capable of handling complex, nonlinear, and dependent data. Considering the intricate, dependent, and nonlinear nature of the heart disease dataset, Naive Bayes exhibited an accuracy of 87%.

[3] Sonam Nikhar and colleagues presented a research paper titled "Utilizing Machine Learning Algorithms for Heart Disease Prediction." In their study, they provide a comprehensive explanation of the Naïve Bayes and Decision Tree classifiers, which are commonly employed in heart disease prediction. The researchers conducted an analysis to compare the performance of these predictive data mining techniques using the same dataset. The findings revealed that the Decision Tree classifier exhibited superior accuracy compared to the Bayesian classifier.

[4] In their publication titled "Forecasting Heart Disease through Machine Learning," Aditi Gavhane and colleagues introduced a novel approach to predicting heart disease. The study utilized the multi-layer perceptron neural network algorithm for training and testing the dataset. The algorithm incorporates an input layer, an output layer, and one or more hidden layers between them. Each input node is connected to the output layer through these hidden layers, with random weights assigned to the connections. Additionally, a bias input is included, and its weight can be adjusted according to the specific needs of the connection, allowing for both feedforward and feedback connections between the nodes.

[5] In their paper titled "Heart Disease Prediction Using Effective Machine Learning Techniques," Avinash Golande and colleagues introduced a novel approach to assist doctors in distinguishing heart diseases. The study employs various data mining techniques, including k-nearest neighbor, Decision tree, and Naïve Bayes algorithms. Additionally, distinctive characterization-based methods such as packing algorithm, Part thickness, consecutive minimal optimization, neural networks, linear Kernel self-organizing map, and Support Vector Machines (SVM) are utilized.

[6] In their paper titled "Heart Disease Prediction using Machine Learning Techniques," Lakshmana Rao and colleagues addressed the challenge of identifying heart disease due to the numerous contributing factors involved. To assess the severity of heart disease in individuals, various neural networks and data mining methods were employed.

[7] In their work titled "Heart Attack Prediction Utilizing Deep Learning," Abhay Kishore and colleagues introduced a system that employs deep learning techniques to predict heart attacks and identify potential factors associated with heart-related infections in patients. The model utilizes Recurrent Neural Systems to achieve accurate results and minimize errors. This study serves as a valuable point of reference for future developments in heart attack prediction models.

[8] In their study titled "Enhancing Accuracy in Cardiovascular Disease Prediction through Hybrid Machine Learning Approaches," Senthil Kumar Mohan and colleagues aimed to improve the precision of cardiovascular issue diagnosis. The authors employed a combination of KNN, LR, SVM, and NN algorithms, resulting in an enhanced performance level. Their prediction model for heart disease, known as hybrid random forest with linear model (HRFLM), achieved a precision level of 88.7%.

[9] In their study, Anjan N. Repaka and colleagues presented a model that evaluates the predictive performance of two classification models. This model was analyzed and compared to prior research. The experimental findings demonstrate that our proposed method outperforms other models in accurately determining the percentage of risk prediction.

[10] In their study titled "Heart Disease Prediction through Evolutionary Rule Learning," Aakash Chauhan et al. introduced a method that utilizes electronic records to streamline data retrieval and reduce manual efforts. By employing frequent pattern growth association mining on patients' datasets, the researchers generated robust associations that significantly contributed to accurate heart disease prediction. This approach effectively decreased the number of services required for prediction, enhancing the overall efficiency of the process.

**CHAPTER 3**

**METHODOLOGY**

**3.1 Existing System**

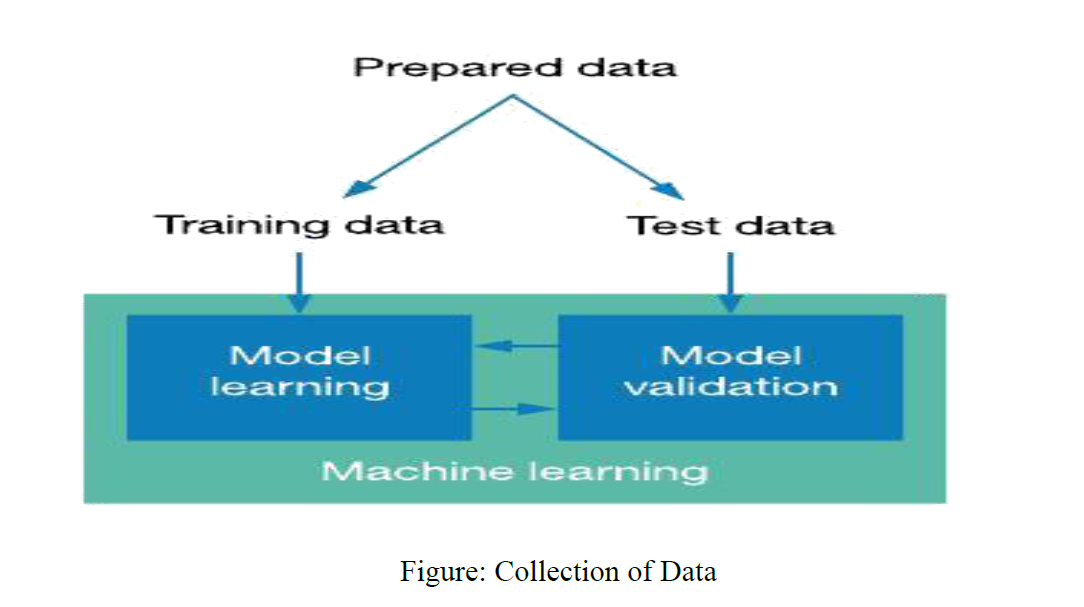
Heart disease is increasingly recognized as a silent killer, causing death without obvious symptoms. This characteristic of the disease has led to growing concerns and anxiety about its impact. Therefore, ongoing efforts are being made to forecast the likelihood of this fatal illness in advance. Various tools and techniques are continuously being explored to meet present-day healthcare needs, and machine learning methods hold great potential in this regard. Although heart disease can manifest in different forms, there exists a common set of core risk factors that determine an individual's susceptibility. By gathering data from diverse sources, organizing it into appropriate categories, and conducting thorough analysis, we can draw conclusions. This approach can effectively be employed for predicting heart disease. Following the well-known adage, "Prevention is better than cure," early prediction and control measures can significantly aid in reducing heart disease-related mortality rates.

**3.2 Proposed System**

The system begins by gathering data and identifying significant attributes. Subsequently, the necessary data undergoes preprocessing to achieve the desired format. The data is then split into two sets: training and testing data. Algorithms are employed to train the model using the training data. To evaluate the system's accuracy, it is tested with the testing data. This implementation incorporates the following modules.

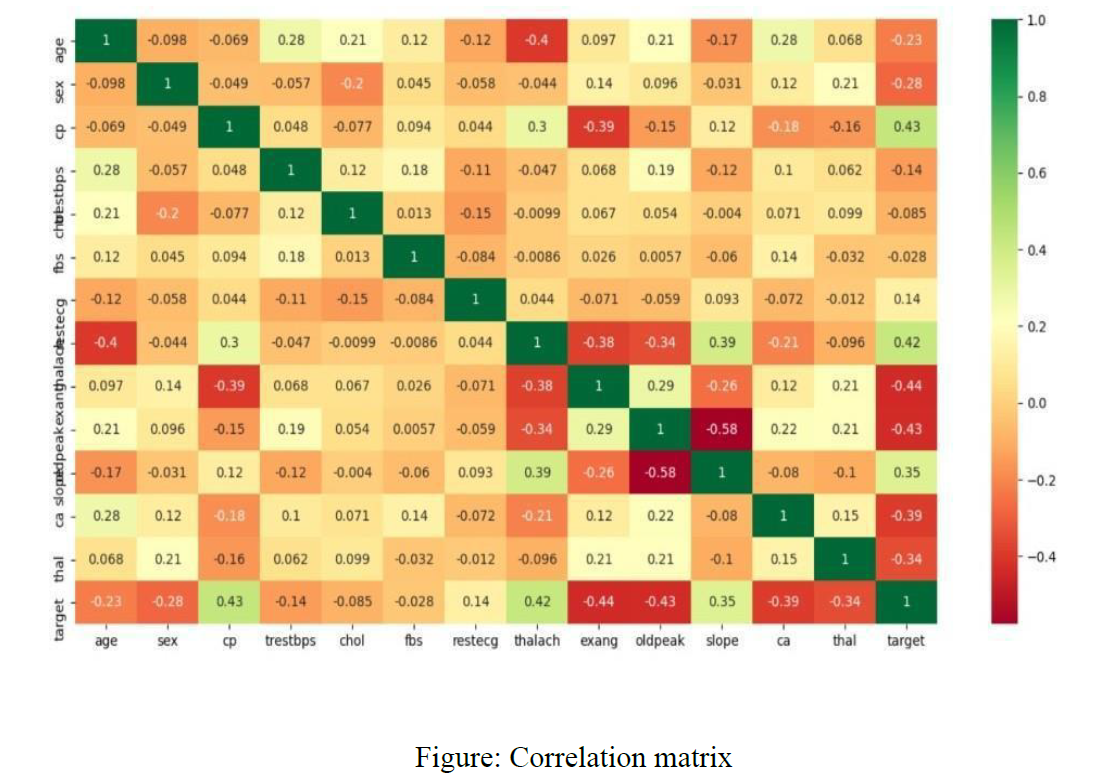
**3.2.1 Collection of Dataset**

In the beginning, we gather a dataset specifically for our heart disease prediction system. Following the dataset collection phase, we divide it into two subsets: training data and testing data. The training dataset is employed to train the prediction model, while the testing data is utilized to assess the performance of the model. In this project, we allocate 70% of the data for training and 30% for testing. We utilize the Heart Disease UCI dataset, which encompasses a total of 76 attributes. However, for our system, we only utilize 14 of these attributes.



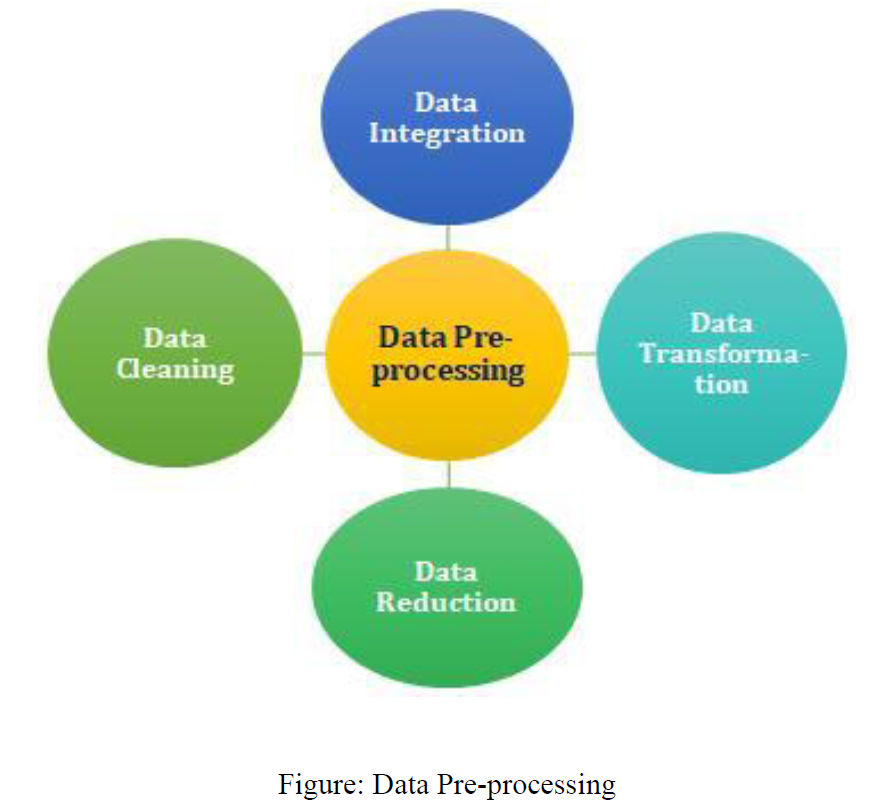
**3.2.2 Selection of Attributes**

The process of attribute or feature selection involves choosing suitable attributes for the prediction system in order to enhance its efficiency. This selection includes various patient attributes such as gender, chest pain type, fasting blood pressure, serum cholesterol, exang, and more, which are utilized for prediction purposes. The model employs a correlation matrix to aid in the attribute selection process.



**3.2.3 Pre-processing Data**

Data preprocessing plays a crucial role in developing a machine learning model. At the outset, the data may be unrefined or not in the desired format, leading to potentially misleading results. To address this, data preprocessing involves converting the data into the necessary format. Its purpose is to handle issues such as noise, duplicates, and missing values within the dataset. Activities involved in data preprocessing include importing datasets, splitting them, and scaling attributes, among others. Ultimately, data preprocessing is essential for enhancing the model's accuracy.

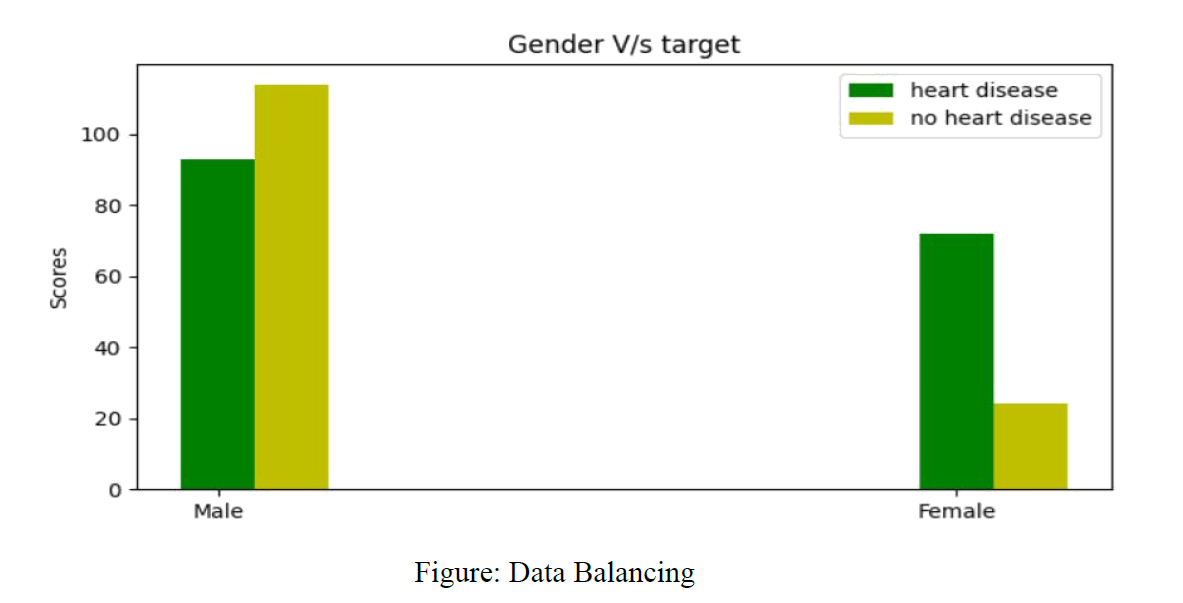


**3.2.4 Balancing of Data**

Two methods can be employed to address imbalanced datasets, namely Under Sampling and Over Sampling.

(a) Under Sampling involves reducing the size of the majority class in order to balance the dataset. This approach is used when there is a sufficient amount of data available.

(b) Over Sampling involves increasing the size of the minority class in order to achieve dataset balance. This approach is used when there is an insufficient amount of data available.



**3.2.5 Prediction of Disease**

Different classification algorithms, including SVM, Naive Bayes, Decision Tree, Random Tree, and Logistic Regression, are utilized in machine learning for predicting heart disease. A comparative evaluation is conducted among these algorithms to determine the one that yields the highest accuracy in heart disease prediction.

Diagram

Description automatically generated

**CHAPTER 4**

**WORKING OF SYSTEM**

**4.1 System Architecture**

The system architecture gives an overview of the working of the system.

The operational procedure of this system can be outlined as follows:

The initial step involves gathering a dataset comprising patient information. Next, a process is carried out to select the relevant attributes that are useful for predicting heart disease. Once the available data resources are identified, they undergo a selection and cleaning process to be transformed into the required format. Subsequently, various classification techniques mentioned earlier are applied to the preprocessed data in order to determine the accuracy of heart disease prediction. Finally, an accuracy measure is utilized to compare the performance of different classifiers.

**Diagram

Description automatically generated**

**4.2 Machine Learning**

In the field of machine learning, the task of classification involves making predictions about the class label associated with a given input data instance.

* **Supervised Learning:**

Supervised learning refers to a type of machine learning where machines are trained using labeled training data, enabling them to predict outputs based on this information. Labeled data signifies input data that has been associated with the correct output. In supervised learning, the provided training data acts as a supervisor, instructing the machines on how to accurately predict the output. This process parallels the way a student learns under the guidance of a teacher. The objective of a supervised learning algorithm is to discover a mapping function that establishes a relationship between the input variable (x) and the output variable (y), utilizing both input and correct output data provided to the machine learning model.

* + **Unsupervised learning:**

Unsupervised learning cannot be directly applied to regression or classification tasks because it lacks corresponding output data, unlike supervised learning. Instead, its objective is to uncover the inherent structure of a dataset, group similar data points, and represent the dataset in a condensed form.

i) Unsupervised learning is valuable for extracting meaningful insights from data.

ii) Unsupervised learning mimics the way humans learn through personal experiences, making it closer to true AI.

iii) Unsupervised learning operates on unlabeled and uncategorized data, emphasizing its significance.

iv) Unsupervised learning is necessary for situations where input data and corresponding outputs are unavailable in real-world scenarios.

* **Reinforcement Learning:**

Reinforcement learning, a branch of Machine Learning, focuses on selecting optimal actions to maximize rewards within a given context. This approach is utilized by different software and machines to determine the most favorable behavior or path in specific situations. Unlike supervised learning, where the training data provides explicit answers, reinforcement learning relies on the reinforcement agent's decision-making to accomplish a given task without predetermined solutions. As a result, in the absence of training datasets, the agent learns through its own experiences.

**4.3 Algorithms**

An AI system relies on a machine learning algorithm to perform its task, typically making predictions based on input data. Classification and regression are the primary techniques employed by machine learning algorithms.

* **Decision Tree Algorithm:**

The technique utilized by an AI system to perform its task is known as a machine learning algorithm. Typically, these algorithms predict output values based on provided input data. Machine learning algorithms primarily encompass two fundamental procedures: classification and regression.

Decision nodes are utilized to make determinations and exhibit numerous branches, while Leaf nodes are the resulting outcomes of those determinations and do not possess additional branches. The determinations or tests are executed depending on characteristics found in the provided dataset. This form of graphical representation offers a comprehensive array of potential solutions for a problem or decision based on given conditions. It acquires the name "Decision Tree" due to its resemblance to a tree structure, commencing with a root node that extends into further branches. The construction of the tree involves the implementation of the CART algorithm, which stands for Classification and Regression Tree algorithm. Essentially, a Decision Tree poses a question and proceeds to split the tree into subtrees based on the response (Yes/No).

The Decision Tree Algorithm is a type of supervised machine learning algorithm that can be utilized for both classification and regression tasks.

The objective of this algorithm is to develop a predictive model for determining the value of a target variable. To solve the problem, a decision tree is employed, utilizing a tree structure where class labels are assigned to leaf nodes, and attributes are represented within the internal nodes of the tree.

The goal of this algorithm is to create a model that predicts the value of a target variable, for which the decision tree uses the tree representation to solve the problem in which the leaf node corresponds to a class label and attributes are represented on the internal node of the tree.

* Decision Trees usually mimic human thinking ability while deciding, so it is easy to understand.
* The logic behind the decision tree can be easily understood because it shows a tree-like structure.

The primary difficulty in Decision Tree lies in determining the attribute for the root node at each level, which is referred to as attribute selection. There are two well-known measures for selecting attributes.

* **Information Gain:**

The utilization of a node in a Decision Tree for dividing the training instances into smaller subsets results in a modification of entropy. Information gain quantifies the extent of this entropy alteration. Entropy represents the level of uncertainty associated with a random variable and defines the impurity of a given set of examples. A greater entropy value indicates a higher amount of information contained within.

* **Gini Index:**

The Gini Index is a measurement used to assess the likelihood of misidentifying a randomly selected element. It indicates that attributes with lower Gini Index values are more desirable. Sklearn includes support for the "Gini" criteria in calculating the Gini Index, with the default value being set to "Gini".

The most notable types of Decision Tree algorithms are: -

* **IDichotomiser 3 (ID3):**

This algorithm employs the concept of Information Gain to determine the most appropriate attribute for classifying the present subset of data. At each level of the tree, the information gain is computed iteratively for the remaining data.

* **C4.5:**

The algorithm in question is a follow-up to the ID3 algorithm, and it employs either Information gain or Gain ratio to determine the attribute for classification. It represents a significant advancement over the ID3 algorithm, as it has the capability to manage continuous and missing attribute values.

* **Classification and Regression Tree (CART):**

It is a dynamic learning algorithm which can produce a regression tree as well as a classification tree depending upon the dependent variable.

* **Working:**

When utilizing a Decision Tree to predict the class of a given dataset, the process begins at the root node of the tree. This involves comparing the attribute values of the root with the corresponding attributes of the dataset, and subsequently proceeding to the appropriate branch and advancing to the next node based on the outcome of the comparison.

This procedure is then repeated for each subsequent node, with the algorithm comparing attribute values to determine the path to follow. This iterative process continues until the algorithm reaches a leaf node in the tree. To gain a clearer understanding of the entire process, refer to the algorithm provided below.

Step 1: Start constructing the tree by establishing a root node called S, which encompasses the entire dataset.

Step 2: Determine the most suitable attribute from the dataset using an Attribute Selection Measure (ASM).

Step 3: Partition S into subsets based on the potential values associated with the selected attribute.

Step 4: Formulate a Decision Tree node that incorporates the chosen attribute.

Step 5: Repeat the procedure recursively for each subset generated in step 3, generating new decision trees. Continue this iterative process until reaching a stage where further classification of the nodes is not possible, designating the final node as a leaf node.

* **Logistic Regression Algorithm:**

Logistic regression is a widely employed algorithm in Machine Learning that falls under the Supervised Learning category. Its purpose is to forecast the outcome of a categorical dependent variable by utilizing a specified group of independent variables. By employing logistic regression, the algorithm produces probabilistic values ranging between 0 and 1 instead of precise 0s and 1s, allowing for predictions of categorical outcomes such as Yes or No, 0 or 1, true or false, and similar.

Logistic regression is similar to linear regression, but their applications differ. While linear regression is employed to address regression problems, logistic regression is utilized for classification problems. In logistic regression, we don't fit a regression line; instead, we fit a sigmoid or "S" shaped logistic function that predicts either 0 or 1 as the maximum values.

The logistic function's curve portrays the probability of various outcomes, such as determining if cells are cancerous or not, if a mouse is obese or not based on its weight, and so on. Logistic Regression holds immense importance in the field of machine learning as it enables the calculation of probabilities and facilitates the classification of new data using both continuous and discrete datasets.

* **Advantages:**

Logistic Regression stands out as a straightforward machine learning technique that is both simple to implement and capable of delivering efficient training in certain scenarios. Moreover, the algorithm's effectiveness eliminates the need for substantial computational resources when training a model.

The inferred parameters (trained weights) provide insight into the significance of each feature and also indicate the direction of the association, whether positive or negative. Therefore, Logistic Regression can be employed to determine the connection between the features.

The ability of this algorithm to incorporate new data and update models sets it apart from Decision Tree or Support Vector Machine, with stochastic gradient descent being employed for the updating process.

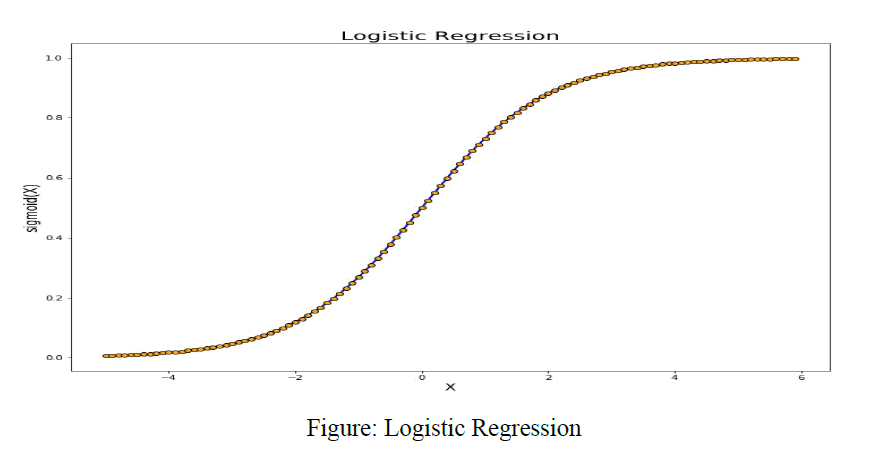
Logistic Regression offers the benefit of providing both well-calibrated probabilities and classification outcomes. This distinguishes it from models that solely produce final classifications. By considering the probabilities assigned to each class, we can determine the relative accuracy of training examples for the specific problem at hand. For instance, if one training example has a 95% probability for a certain class, while another has a 55% probability for the same class, we can make an inference about their respective accuracy levels.

* **Disadvantages:**

Logistic Regression is a statistical model used for analyzing data and making predictions based on independent variables. However, when dealing with datasets that have a high number of dimensions, there is a risk of overfitting the model to the training data. Overfitting occurs when the model becomes too specialized to the training set, resulting in inflated accuracy during prediction. Consequently, the model may struggle to provide accurate results when applied to the test set. This issue commonly arises when training data is limited, but there are numerous features. To address overfitting in high-dimensional datasets, it is advisable to employ regularization techniques. However, it is worth noting that regularization can increase the complexity of the model. Excessive regularization factors may even lead to the model underfitting the training data.

Logistic regression is limited in its ability to solve non-linear problems due to its reliance on a linear decision surface. In practical situations, it is uncommon to encounter data that is easily separable in a linear manner. To address this, transforming non-linear features becomes necessary. One approach is to increase the number of features, enabling the data to become linearly separable in higher dimensions.

Non-Linearly Separable Data: Logistic regression struggles to capture intricate relationships, making it challenging. However, more advanced and intricate algorithms like Neural Networks can surpass the performance of this method effortlessly.



* **KNN (K-Nearest Neighbor):**

The K-Nearest Neighbor (K-NN) algorithm is a straightforward Supervised Learning technique widely used in Machine Learning. It operates under the assumption of similarity between new and existing data points, assigning the new case to the category that closely matches the available categories. By retaining all available data, the K-NN algorithm can effectively classify new data points into appropriate categories.

K-NN is a versatile algorithm applicable to both Regression and Classification tasks, although it is primarily employed for the latter. Being a non-parametric algorithm, K-NN does not rely on assumptions about the underlying data distribution. This characteristic has earned it the nickname of a "lazy learner" since it postpones learning from the training set. Instead, the algorithm stores the dataset and takes action when classifying new data points.

During the training phase, the K-NN algorithm simply stores the dataset. When encountering new data, it classifies it into the category that closely resembles the new data, based on similarity.

**Example:** Consider the following scenario: We possess an image of a creature bearing resemblance to both a cat and a dog, and our objective is to determine whether it belongs to the cat or dog category. In this situation, the KNN algorithm can be employed for identification purposes, as it operates on a measure of similarity. By utilizing the KNN model, we can discover the shared characteristics between the newly acquired dataset and the images of cats and dogs. Subsequently, based on the most analogous features, the algorithm will assign the creature to either the cat or dog classification.

Graphical user interface, application

Description automatically generated

**Working:**

The K-NN working can be explained on the basis of the below algorithm:

**Step-1:** Select the number K of the neighbors.

**Step-2:** Calculate the Euclidean distance of **K number of neighbors**

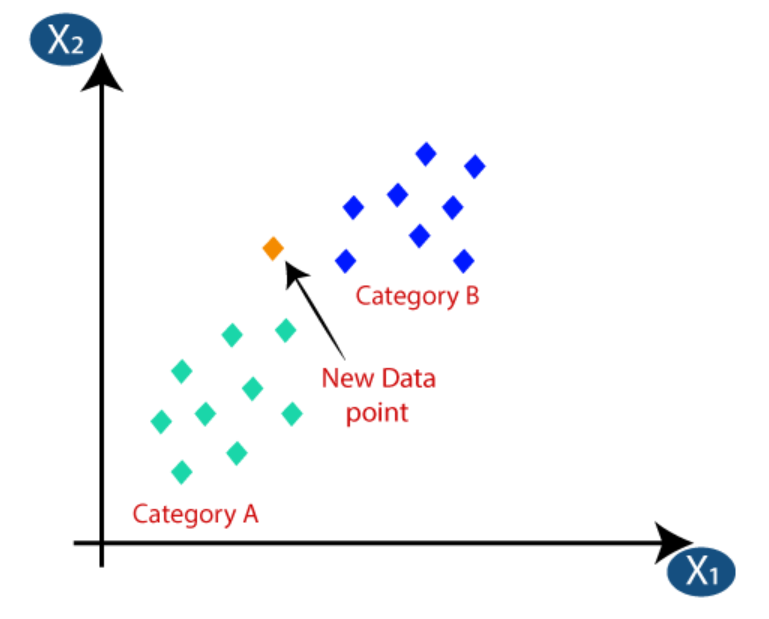
**Step-3:** Take the K nearest neighbors as per the calculated Euclidean distance.

**Step-4:** Among these k neighbors, count the number of the data points in each category.

**Step-5:** Assign the new data points to that category for which the number of the neighbor is maximum.

**Step-6:** Our model is ready.

Suppose we have a new data point, and we need to put it in the required category. Consider the below image:



* Firstly, we will choose the number of neighbors, so we will choose the k=5.
* Next, we will calculate the **Euclidean distance** between the data points. The Euclidean distance is the distance between two points, which we have already studied in geometry. It can be calculated as:

Chart, line chart

Description automatically generated

* By calculating the Euclidean distance we got the nearest neighbors, as three nearest neighbors in category A and two nearest neighbors in category B. Consider the below image:

Chart, scatter chart

Description automatically generated

* As we can see the 3 nearest neighbors are from category A, hence this new data point must belong to category A.

## **Advantages of KNN Algorithm:**

* It is simple to implement.
* It is robust to the noisy training data.
* It can be more effective if the training data is large.

## **Disadvantages of KNN Algorithm:**

* Always needs to determine the value of K which may be complex some time.
* The computation cost is high because of calculating the distance between the data points for all the training samples.

**4.4 Code Implementation:**

* **Import Libraries**

import pandas as pd

import numpy as np

import matplotlib.pyplot as plt

import seaborn as sns

import missingno as msn

* **Read CSV file**

df=pd.read\_csv("heart.csv");

df.head()

df.tail()

df.shape

print('Sex', df['Sex'].unique())

print('ChestPainType', df['ChestPainType'].unique())

print('RestingECG', df['RestingECG'].unique())

print('ExerciseAngina', df['ExerciseAngina'].unique())

print('ST\_Slope', df['ST\_Slope'].unique())

sex\_map = {'M':1, 'F':2}

chest\_pain\_type\_map = {'ATA':1, 'NAP':2, 'ASY':3, 'TA':4 }

resting\_ecg\_map = {'Normal':1, 'ST':2, 'LVH':3}

excercise\_angina\_map = {'N':1, 'Y':2}

st\_slop\_map = {'Up':1, 'Flat':2, 'Down':3 }

df['Sex']=df['Sex'].map(sex\_map)

df['ChestPainType']=df['ChestPainType'].map(chest\_pain\_type\_map)

df['RestingECG']=df['RestingECG'].map(resting\_ecg\_map)

df['ExerciseAngina']=df['ExerciseAngina'].map(excercise\_angina\_map)

df['ST\_Slope']=df['ST\_Slope'].map(st\_slop\_map)

df.head()

sns.pairplot(df);

fig,ax=plt.subplots(figsize=(8,4));

sns.lineplot(data=df['Cholesterol']);

fig,ax=plt.subplots(figsize=(8,4));

sns.lineplot(data=df['Age']);

df.columns

X=df[['Age', 'Sex', 'ChestPainType', 'RestingBP', 'Cholesterol', 'FastingBS',

'RestingECG', 'MaxHR', 'ExerciseAngina', 'Oldpeak', 'ST\_Slope']]

X.head()

X = X.values

y=df['HeartDisease']

y.head()

y = y.values

* **Splitting Dataset into training and testing by train\_test\_split**

from sklearn.model\_selection import train\_test\_split

X\_train,X\_test,y\_train,y\_test=train\_test\_split(X,y,test\_size=0.3,random\_state=22)

1. **Logistic Regression Model**

from sklearn.linear\_model import LogisticRegression

model\_lgr=LogisticRegression()

model\_lgr.fit(X\_train,y\_train)

y\_predict\_lgr=model\_lgr.predict(X\_test)

* **Model Evaluation**

from sklearn.metrics import accuracy\_score,confusion\_matrix,classification\_report

print("Accuracy of the model by Logistic Regression is :",accuracy\_score(y\_test,y\_predict\_lgr)\*100)

print(confusion\_matrix(y\_test,y\_predict\_lgr))

print(classification\_report(y\_test,y\_predict\_lgr))

1. **KNN**

from sklearn.neighbors import KNeighborsClassifier

model\_knn=KNeighborsClassifier()

model\_knn.fit(X\_train,y\_train)

y\_predict\_knn=model\_knn.predict(X\_test)

* **Model Evaluation**

from sklearn.metrics import accuracy\_score,confusion\_matrix,classification\_report

print("Accuracy of the model by KNN is :",accuracy\_score(y\_test,y\_predict\_knn)\*100)

print(confusion\_matrix(y\_test,y\_predict\_knn))

print(classification\_report(y\_test,y\_predict\_knn))

1. **Decision Tree**

from sklearn import tree

model\_tree = tree.DecisionTreeClassifier() # empty model of the decision tree

model\_tree.fit(X\_train,y\_train)

y\_predict\_tree=model\_tree.predict(X\_test)

* **Model Evaluation**

from sklearn.metrics import accuracy\_score,confusion\_matrix,classification\_report

print("Accuracy of the model by Decision Tree is : ",accuracy\_score(y\_test,y\_predict\_tree)\*100)

print(confusion\_matrix(y\_test,y\_predict\_tree))

print(classification\_report(y\_test,y\_predict\_tree))

* **Implementing Decision Tree, KNN & Logistic Regression**

import ipywidgets as widgets

import ipywidgets

from ipywidgets import interact

selected\_algo = None

def onSelectAlgoritham(x):

global selected\_algo

selected\_algo = x

selected\_sex = None

def onSelectSex(x):

global selected\_sex

selected\_sex = x

selected\_chest\_pain\_type = None

def onSelectChestPainType(x):

global selected\_chest\_pain\_type

selected\_chest\_pain\_type = x

selected\_resting\_ecg = None

def onSelectRestingECG(x):

global selected\_resting\_ecg

selected\_resting\_ecg = x

selected\_exercise\_angina = None

def onSelectExerciseAngina(x):

global selected\_exercise\_angina

selected\_exercise\_angina = x

selected\_st\_slope = None

def onSelectST\_Slope(x):

global selected\_st\_slope

selected\_st\_slope = x

interact(onSelectAlgoritham, x=ipywidgets.Combobox(options=["LogisticRegression", "KNeighborsClassifier", "DecisionTreeClassifier"]));

input\_age=float(input("Enter age :-"))

interact(onSelectSex, x=ipywidgets.Combobox(options=['Male', 'Female']));

interact(onSelectChestPainType, x=ipywidgets.Combobox(options=['ATA', 'NAP', 'ASY', 'TA']));

input\_resting\_bp=float(input("Enter RestingBP : "))

input\_cholesterol=float(input("Enter Cholesterol : "))

input\_fasting\_bs=float(input("Enter FastingBS : "))

interact(onSelectRestingECG, x=ipywidgets.Combobox(options=['Normal', 'ST', 'LVH']));

input\_max\_hr=float(input("Enter Max HR : "))

interact(onSelectExerciseAngina, x=ipywidgets.Combobox(options=['Yes', 'No']));

input\_oldpeak = float(input("Enter Oldpeak : "))

interact(onSelectST\_Slope, x=ipywidgets.Combobox(options=['Up', 'Flat', 'Down']));

print(selected\_algo)

print(input\_age)

print(selected\_sex)

print(selected\_chest\_pain\_type)

print(input\_resting\_bp)

print(input\_cholesterol)

print(input\_fasting\_bs)

print(selected\_resting\_ecg)

print(input\_max\_hr)

print(selected\_exercise\_angina)

print(input\_oldpeak)

print(selected\_st\_slope)

sex\_int = sex\_map[selected\_sex[0]]

sex\_int = float(sex\_int)

chest\_pain\_type\_int = chest\_pain\_type\_map[selected\_chest\_pain\_type]

chest\_pain\_type\_int = float(chest\_pain\_type\_int)

resting\_ecg\_int = resting\_ecg\_map[selected\_resting\_ecg]

resting\_ecg\_int = float(resting\_ecg\_int)

selected\_exercise\_int = excercise\_angina\_map[selected\_exercise\_angina[0]]

selected\_exercise\_int = float(selected\_exercise\_int)

selected\_st\_slope\_int = st\_slop\_map[selected\_st\_slope]

selected\_st\_slope\_int = float(selected\_st\_slope\_int)

if selected\_algo == "LogisticRegression":

result = model\_lgr.predict([[input\_age, sex\_int, chest\_pain\_type\_int, input\_resting\_bp

, input\_cholesterol, input\_fasting\_bs, resting\_ecg\_int, input\_max\_hr,

selected\_exercise\_int, input\_oldpeak, selected\_st\_slope\_int]])

print("LogisticRegression:")

print(result)

if result[0]==0:

print("Heart disease not deteacted!")

elif result[0]==1:

print("Heart disease deteacted!")

elif selected\_algo == "KNeighborsClassifier":

result = model\_knn.predict([[input\_age, sex\_int, chest\_pain\_type\_int, input\_resting\_bp

, input\_cholesterol, input\_fasting\_bs, resting\_ecg\_int, input\_max\_hr,

selected\_exercise\_int, input\_oldpeak, selected\_st\_slope\_int]])

print("KNeighborsClassifier:")

print(result)

if result[0]==0:

print("Heart disease not deteacted!")

elif result[0]==1:

print("Heart disease deteacted!")

elif selected\_algo == "DecisionTreeClassifier":

result = model\_tree.predict([[input\_age, sex\_int, chest\_pain\_type\_int, input\_resting\_bp

, input\_cholesterol, input\_fasting\_bs, resting\_ecg\_int, input\_max\_hr,

selected\_exercise\_int, input\_oldpeak, selected\_st\_slope\_int]])

print("DecisionTreeClassifier:")

print(result)

if result[0]==0:

print("Heart disease not deteacted!")

elif result[0]==1:

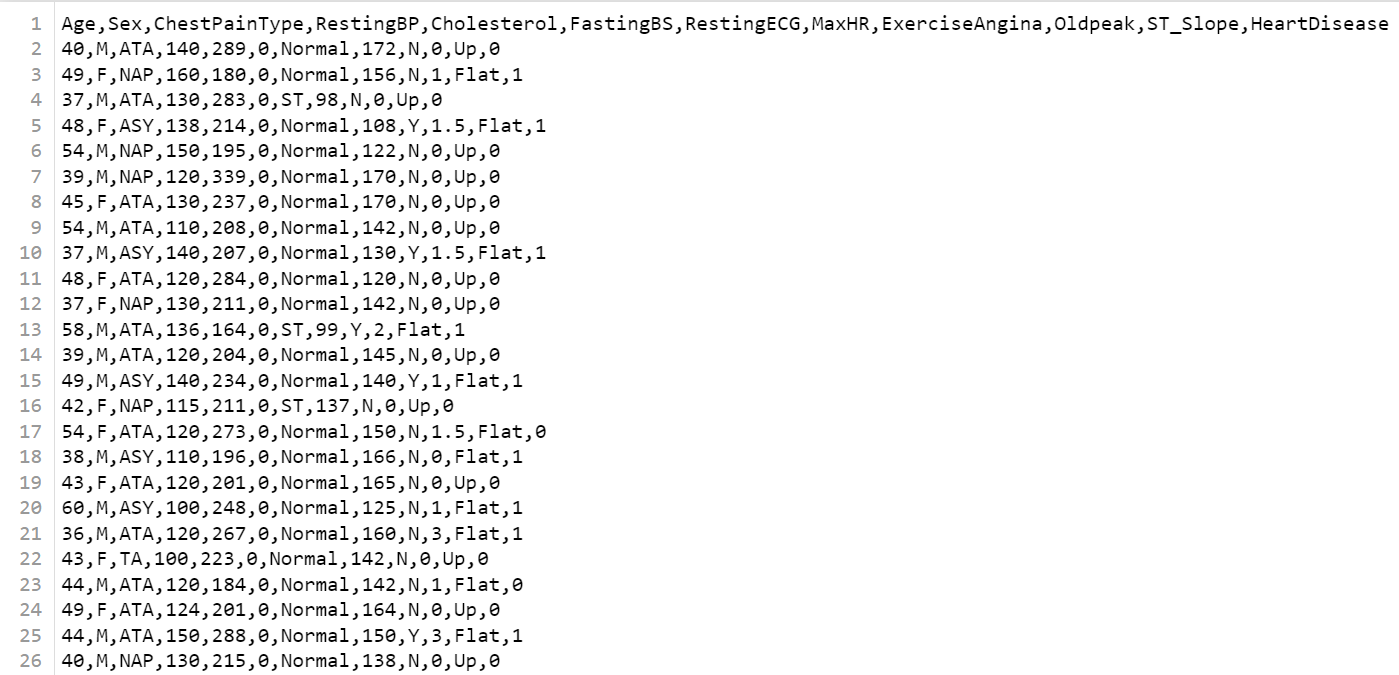
print("Heart disease deteacted!")

**CHAPTER 5**

**EXPERIMENTAL ANALYSIS**

|  |  |  |
| --- | --- | --- |
| **5.1 SYSTEM CONFIGURATION**  **5.1.1 Hardware requirements:** | | |
| Processer | : | Any Update Processer |
| Ram | : | Min 4GB |
| Hard Disk | : | Min 100GB |
| **5.1.2 Software requirements:** | | |
| Operating System | : | Windows family |
| Technology | : | Python3.7 |
| IDE | : | Jupiter notebook |

**5.2 Dataset Details:**

****

**Text

Description automatically generated with medium confidence**

**Input Dataset Attributes:**

* + Age in year
  + Sex: value (M: Male, F: Female)
  + Chest Pain Type
  + Resting BP: Resting Blood Pressure
  + Cholesterol
  + FastingBS: Fasting Blood Sugar value (1: >120md/dl, 0: <120mg/dl)
  + RestingECG: Resting electrocardiograph can detect coronary heart diseases.
  + MaxHR: Maximum Heart Rate
  + Exercise Angina
  + OldPeak
  + ST\_Slope: The ST segment shift relative to exercise-induced increments in heart rate, the ST/heart rate slope, has been proposed as a more accurate ECG criterion for diagnosing significant coronary artery disease (CAD).

|  |  |  |  |
| --- | --- | --- | --- |
| S. No. | Attributes | Description | Type |
| 1. | Age | Patient’s age (29 to 77) | Numerical |
| 2. | Sex | Gender of patient | Nominal |
| 3. | Chest Pain Type | Chest pain type | Nominal |
| 4. | RestingBP | Resting blood pressure | Numerical |
| 5. | Cholesterol | Serum cholesterol in mg/dl | Numerical |
| 6. | FastingBS | Fasting blood sugar | Numerical |
| 7. | RestingECG | Resting electrocardiograph | Nominal |
| 8. | MaxHR | Maximum Heart Rate | Numerical |
| 9. | Exercise Angina |  | Nominal |
| 10. | OldPeak |  | Numerical |
| 11. | ST\_Slope | ST segment shift | Nominal |

**Attributes of the Dataset**

**5.3 Performance Analysis**

In this study, heart disease prediction is conducted using different machine learning algorithms including Decision Tree, Logistic Regression, and KNN. The Heart Disease UCI dataset contains a total of 76 attributes, but only 11 attributes are taken into account for predicting heart disease. Patient attributes such as gender, chest pain type, fasting blood pressure, serum cholesterol, and exang are among the factors considered. Each algorithm's accuracy is calculated, and the algorithm with the highest accuracy is chosen for heart disease prediction. To evaluate the experiment, several metrics including accuracy, confusion matrix, precision, recall, and f1-score are utilized.

Accuracy is defined as the proportion of correct predictions to the total number of inputs in the dataset, and it can be calculated as follows:

Accuracy = (TP + TN) / (TP + FP + FN + TN)

**Confusion Matrix**: - It gives us a matrix as output and gives the total performance of the system.

Chart

Description automatically generated

where

TP: True positive

FP: False Positive

FN: False Negative

TN: True Negative

**Correlation Matrix**: - The correlation matrix in machine learning is used for feature selection. It represents dependency between various attributes.

**Precision: -** It is the ratio of correct positive results to the total number of positive results predicted by the system.

It is expressed as:

* Recall-It is the ratio of correct positive results to the total number of positive results.
* predicted by the system.
* It is expressed as:
* F1 Score-It is the harmonic mean of Precision and Recall. It measures the test accuracy. The range of this metric is 0 to 1.

**5.4 Performance Measures**

|  |  |  |  |  |
| --- | --- | --- | --- | --- |
|  | **Precision** | **recall** | **F1-Score** | **Support** |
| **0** | 0.80 | 0.84 | 0.82 | 122 |
| **1** | 0.87 | 0.84 | 0.85 | 154 |
| **accuracy** |  |  | 0.84 | 276 |
| **Micro avg.** | 0.83 | 0.84 | 0.84 | 276 |
| **Weighted avg** | 0.84 | 0.84 | 0.84 | 276 |

**Logistic Regression Model Evaluation**

|  |  |  |  |  |
| --- | --- | --- | --- | --- |
|  | **Precision** | **recall** | **F1-Score** | **Support** |
| **0** | 0.65 | 0.57 | 0.61 | 122 |
| **1** | 0.69 | 0.76 | 0.72 | 154 |
| **accuracy** |  |  | 0.68 | 276 |
| **Micro avg.** | 0.67 | 0.67 | 0.67 | 276 |
| **Weighted avg** | 0.68 | 0.68 | 0.67 | 276 |

**KNN Model Evaluation**

|  |  |
| --- | --- |
| 97 | 25 |
| 34 | 120 |

**Decision Tree Confusion Matrix**

|  |  |  |  |  |
| --- | --- | --- | --- | --- |
|  | **Precision** | **recall** | **F1-Score** | **Support** |
| **0** | 0.74 | 0.80 | 0.77 | 122 |
| **1** | 0.83 | 0.78 | 0.80 | 154 |
| **accuracy** |  |  | 0.79 | 276 |
| **Micro avg.** | 0.78 | 0.79 | 0.78 | 276 |
| **Weighted avg** | 0.79 | 0.79 | 0.79 | 276 |

**Decision Tree Model Evaluation**

* **The highest accuracy is given by Logistic Regression Model.**

**5.5 Result**

In comparison to other algorithms, we discover that the accuracy of the Logistic Regression is superior after applying the machine learning approach for training and testing. With the help of the confusion matrices for each algorithm, accuracy is calculated. Here, the number of TP, TN, FP, and FN is given. Using the equation for accuracy, value has been calculated. It is concluded that extreme gradient boosting is best with 84% accuracy.

TABLE: Accuracy comparison of algorithms Algorithm Accuracy

|  |  |
| --- | --- |
| **Algorithm** | **Accuracy** |
| Logistic Regression | 83.5% |
| KNN | 67.5% |
| Decision Tree | 78.5% |

**Accuracy Table**

**CHAPTER 6**

**CONCLUSION AND FUTURE SCOPE**

The utilization of promising technology, such as machine learning, in the initial prediction of heart problems could have a noteworthy societal impact, considering that heart diseases are a prominent cause of death both in India and worldwide. The early identification of cardiac ailments has the potential to assist high-risk individuals in making informed choices regarding lifestyle adjustments, representing a significant advancement in the field of medicine. Each year, an increasing number of individuals receive a diagnosis of cardiac illnesses, necessitating the need for early detection and appropriate intervention. The application of suitable technological support in this domain has the potential to greatly benefit both the medical community and patients. To assess performance, this study employed a range of machine learning algorithms, including SVM, Decision Tree, Random Forest, Naive Bayes, Logistic Regression, Adaptive Boosting, and Extreme Gradient Boosting.

The dataset contains 76 features that pertain to various characteristics associated with heart disease in individuals. To evaluate the system, the author selects 14 significant features from this dataset. However, the author notices a decrease in system efficiency when considering all the features. In order to enhance efficiency, attribute selection is implemented. This involves choosing a specific number of characteristics (referred to as 'n') that yield higher accuracy for evaluating the model. Some features in the dataset exhibit nearly identical correlations, leading to their elimination. It is observed that system efficiency declines significantly when all attributes in the dataset are considered.

The accuracies of seven different machine learning methods were compared to determine the best prediction model. The objective was to employ various evaluation metrics, such as the confusion matrix, accuracy, precision, recall, and f1-score, to efficiently predict the disease. Among the seven methods, the extreme gradient boosting classifier yielded the highest accuracy of 84%.

**APPENDIX**

**Python**

Python, which was developed by Guido Van Rossum and made available in 1991, is a general-purpose programming language known for its interpretive nature and high-level capabilities. The language places a strong emphasis on code readability by utilizing significant whitespace. With its object-oriented methodology and various language constructs, Python aids programmers in producing coherent and well-structured code for projects of all sizes. Notably, Python is dynamically typed and features automatic garbage collection. It further facilitates diverse programming paradigms such as procedural, object-oriented, and functional programming.

**Sklearn**

Sklearn, also known as Scikit-learn, is an exceptionally valuable and reliable Python library extensively employed for machine learning purposes. It offers a comprehensive range of efficient resources for tasks such as classification, regression, clustering, and dimensionality reduction, all accessible through a unified Python interface. Primarily coded in Python, this library relies on the foundational frameworks of NumPy, SciPy, and Matplotlib.

**Numpy**

NumPy is a Python library that enhances the capabilities of the Python programming language by introducing extensive support for large, multi-dimensional arrays and matrices. It also provides a vast collection of high-level mathematical functions that enable efficient operations on these arrays. The precursor to NumPy, known as Numeric, was initially developed by Jim in collaboration with multiple other programmers. In 2005, Travis took the initiative to merge the features of the rival Num array with Numeric, resulting in the creation of NumPy with significant enhancements. Being open-source software, NumPy benefits from numerous contributors who actively contribute to its development and improvement.

**Librosa**

Librosa serves as a Python toolkit designed for music and audio analysis purposes. It proves particularly useful in various applications involving audio data, such as music generation with LSTMs and Automatic Speech Recognition. This package offers essential components for constructing music information retrieval systems. By employing diverse signal processing techniques, Librosa facilitates audio signal visualization and enables feature extraction.

**Matplotlib**

Matplotlib serves as a plotting library designed for Python and its numerical mathematics extension, NumPy. It offers an object-oriented interface that allows developers to incorporate plots into applications using various GUI toolkits such as Tkinter, wxPython, Qt, or GTK. Additionally, there exists a procedural "pylab" interface, which operates on a state machine similar to OpenGL, aiming to mimic MATLAB's interface, although its usage is not recommended.

**Seaborn**

Seaborn, a Python library for data visualization, is built upon matplotlib and closely integrated with pandas' data structures. It offers a user-friendly interface for creating visually appealing and informative statistical graphics. Seaborn primarily focuses on visualization, facilitating the exploration and comprehension of data.

**SciPy**

SciPy encompasses a range of modules catering to various scientific and engineering tasks, such as optimization, linear algebra, integration, interpolation, special functions, FFT, signal and image processing, ODE solvers, and more. In addition, there are dedicated conferences organized for both users and developers of these tools: SciPy (held in the United States), Euro SciPy (hosted in Europe), and SciPy.in (conducted in India). The original SciPy conference was initiated by Enthought in the United States and remains involved in sponsoring international conferences while also serving as the host for the SciPy website. Functioning as a scientific computation library, SciPy is built upon the foundation of NumPy and offers an expanded set of utility functions encompassing optimization, statistics, and signal processing.

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**Research Paper**

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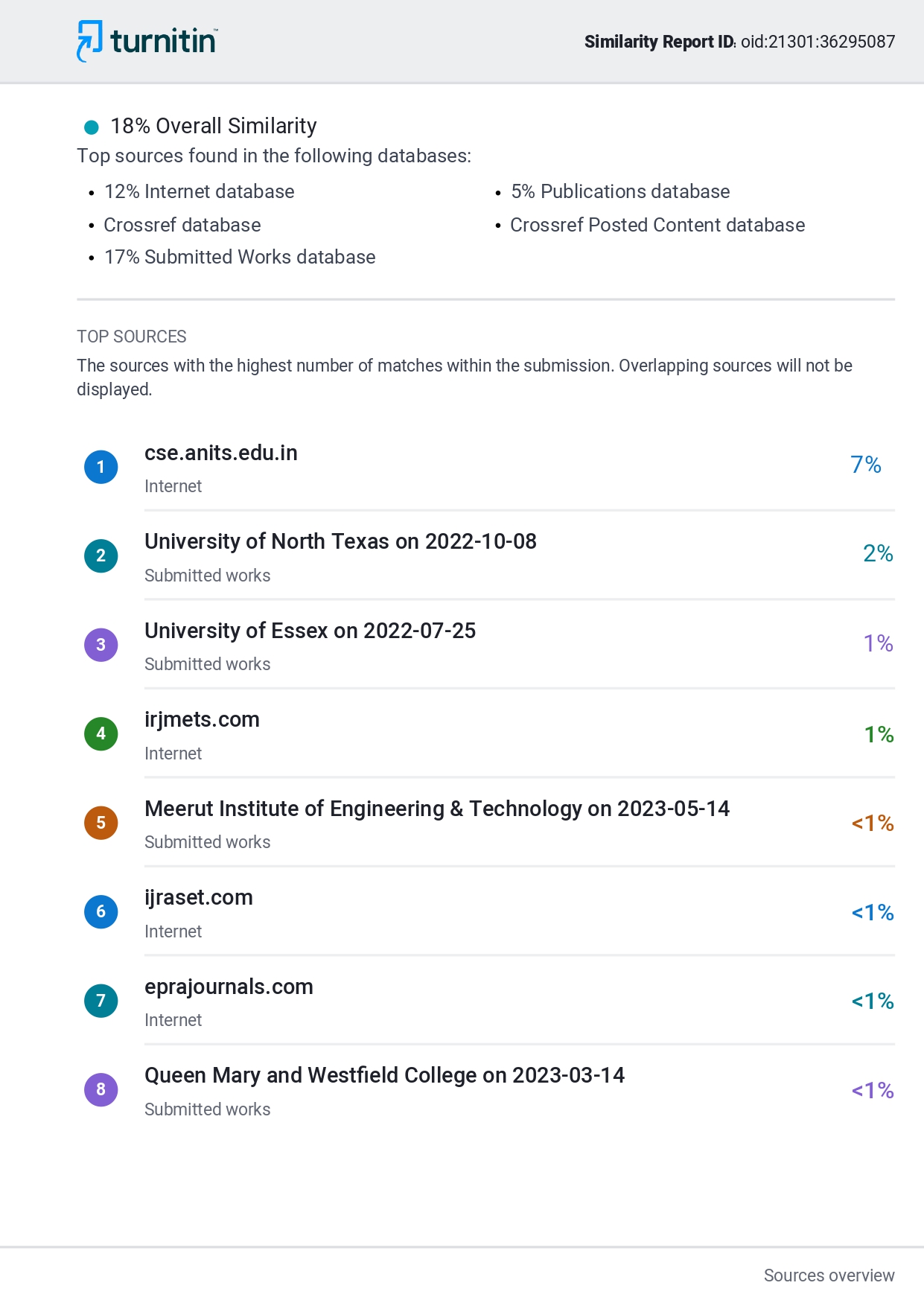
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