INTERDISCIPLINARY PROJECT REPORT

at

Sathyabama Institute of Science and Technology (Deemed to be University)

Submitted in partial fulfillment of the requirements for the award of Bachelor of Engineering Degree in Computer Science and Engineering

Ву

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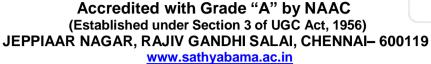
APRIL 2023



SATHYABAMA

INSTITUTE OF SCIENCE AND TECHNOLOGY

(DEEMED TO BE UNIVERSITY)





DEPARTMENT OF COMPUTER SCIENCE AND ENGINEERING

BONAFIDE CERTIFICATE

This is to certify that this Project Report is the bonafide work of **LEVAKU VENKATA KOWSHIK REDDY (Reg No : 40731054)** who carried out the project entitled " **Early risk prediction of forest fires based on Machine Learning Algorithm**" under my supervision from Feb 2023 to April 2023.

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Submitted for Viva voce Examination held on	

Internal Examiner

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DECLARATION

I, LEVAKU VENKATA KOWSHIK REDDY(40731054) hereby declare that the Project Report

entitled " Early risk prediction of forest fires based on Machine Learning Algorithm"

done by me under the guidance of Ms.G.Anbu Selvi, M.Tech.,(PhD) is submitted in partial

fulfillment of the requirements for the award of Bachelor of Engineering degree in

Computer Science and Engineering.

DATE:

PLACE:

SIGNATURE OF THE CANDIDATE

ACKNOWLEDGEMENT

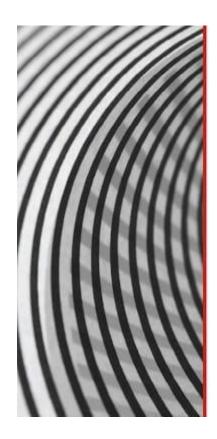
I am pleased to acknowledge my sincere thanks to **Board of Management** of **SATHYABAMA** for their kind encouragement in doing this project and for completing it successfully. I am grateful to them.

I convey my thanks to **Dr.T.Sasikala M.E.,Ph.D**, **Dean**, School of Computing, **Dr.S.Vigneshwari M.E.**, **Ph.D.**, Head of the Department of Computer Science and Engineering for providing me necessary support and details at the right time during the progressive reviews.

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I wish to express my thanks to all Teaching and Non-teaching staff members of the **Department of Computer Science and Engineering** who were helpful in many ways for the completion of the project.

TRAINING CERTIFICATE





Certificate of Training

is hereby granted to

LEVAKU VENKATA KOWSHIK REDDY (40731054)

from Sathyabama Institute of Science and Technology for successfully completing the 45 hours professional training program on Machine Learning conducted by Cognibot between 10th Feb, 2023 and 16th Apr, 2023

Manhoosechon

16th April, 2023

DATE

ABSTRACT

Algerian forest fires have become a major environmental and socioeconomic concern due to their devastating impacts on forests, ecosystems, and human lives. Accurately predicting forest fires is critical for timely detection, management, and prevention of these disasters. Machine learning algorithms have shown promising results in predicting forest fires, and their potential for use in Algeria is currently being investigated. This study aims to explore the use of machine learning algorithms for predicting forest fires in Algeria by analyzing historical data on weather, vegetation, and forest fires. The research will focus on building predictive models using machine learning algorithms such as Random Forest, Support Vector Machines, and Artificial Neural Networks. The models will be trained and tested on a dataset of forest fire incidents in Algeria to evaluate their accuracy and performance. The results will provide insights into the effectiveness of machine learning algorithms in predicting forest fires in Algeria and their potential use in the country's fire management strategies. This study will contribute to the development of early warning systems for forest fires in Algeria, which could help minimize the devastating impacts of these disasters on the environment and society.

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CHAPTER-1 INTRODUCTION

1.1 ESTIMATING FOREST FIRES

Estimating forest fires is a crucial aspect of fire management, as it helps identify the size, intensity, and spread of the fire. One of the primary methods used for estimating forest fires is remote sensing, which involves using satellite data to detect heat signatures and smoke plumes. This data is processed to create maps and images that provide insights into the location and intensity of the fire. Other methods used for estimating forest fires include ground-based observations, aerial surveys, and fire behavior modeling. These methods provide additional information on fire behavior, fuel types, and weather conditions that can help in predicting the fire's future behavior and potential impact.

Accurate estimation of forest fires is essential for timely and effective fire management. Early detection of forest fires and accurate estimation of their size and intensity can help fire managers make informed decisions about resource allocation and prioritization of firefighting efforts. It can also help in the evacuation of affected communities and the implementation of preventative measures to minimize the fire's impact. Therefore, ongoing research on improving estimation methods and developing more advanced technologies, such as machine learning algorithms, is critical for enhancing forest fire management and reducing the negative impacts of these disasters.

1.1.1 TYPES OF FOREST FIRES

There are three basic types of wildfires:

- Crown Fires
- Surface Fires
- Ground Fires

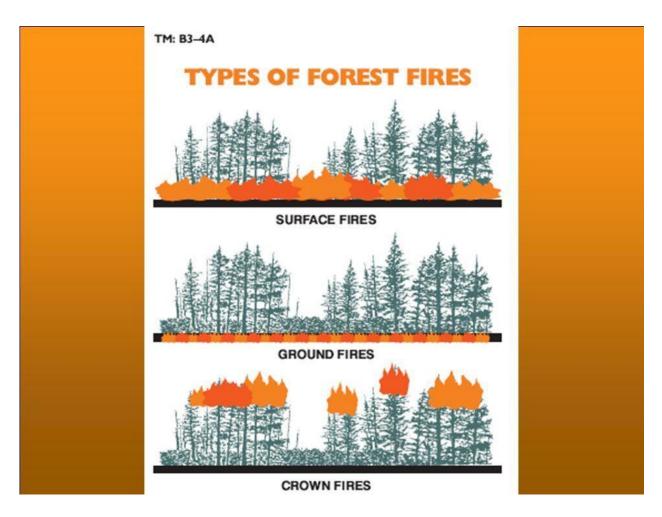


Fig: 1.1.1 Types of Forest Fires

1.2 ESTIMATING FOREST FIRES DATASET INFORMATION

There are 14 attributes in total related to the habits of forest fires and values that are likely to determine forest fires such as day, month, year, temperature, relative humidity, wind speed, rain, fine fuel, druff moisture code, drought code, initial spread index, buildupindex, fire weather index, output

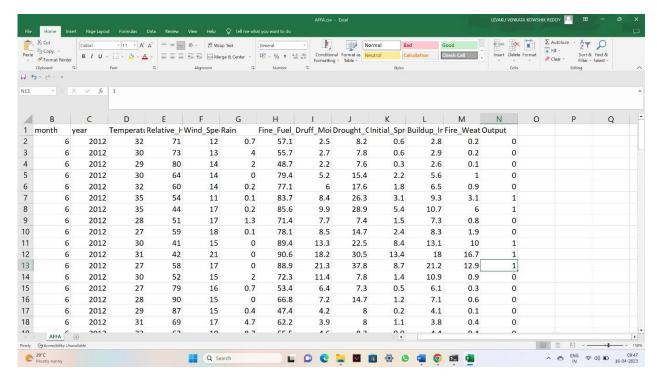


Fig 1.2- Excel sheet of the given estimated forest fires predictions dataset

1.3 COMMON MACHINE LEARNING ALGORITHMS AND GOALS

There are three types of Machine learning algorithms which are been widely used. They are:

- Supervised Learning
- Reinforcement Learning
- Unsupervised Learning

•

Supervised Learning:

Supervised Learning is a machine learning paradigm for problems where the available data consists of labelled examples, meaning that each data point contains features and an associated label. The model is used widely to predict the label of new observations using the features. Depending on the characteristics of the target variable i.e., it can be either classification(discrete variable) or regression(continuous variable).

Unsupervised Learning:

Unsupervised learning is a Machine learning paradigm for problems where the available data consists of un labelled examples, meaning that each data point contains features only, without an associated label. The goal of unsupervised learning algorithms is learning useful patterns or structural properties of the data. It finds the structures in un labelled data.

Reinforcement Learning:

Reinforcement learning is an area of machine learning concerned with how intelligent agents ought to take actions in an environment in order to maximize the notion of cumulative reward. It works on the action-reward principle. An agent learned to reach the goal by continuously calculating the rewards that it gained from the actions.

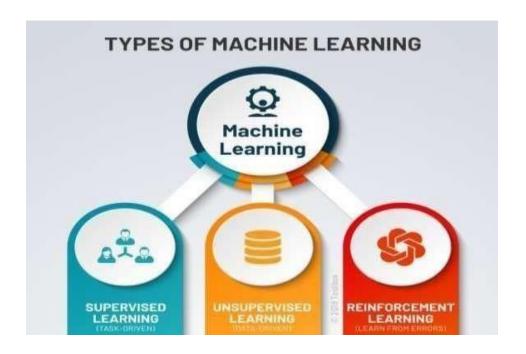


Fig 1.3.1 -Types of machine learning with field of use

ALGORITHMS

1. Linear Regression:

Linear regression is the type of supervised machine learning algorithm where the anticipated output is continuous and features a constant slope. It predicts the values withinendless range, instead of trying to classify them into categories. It is used to predict the worth of a variable supported the worth of another variable. For choosing this algorithm, there needs be a linear relation between independent and target variable. As scatter plot shows the positive correlation between an independent variable(x-axis) and dependent variable (y-axis).

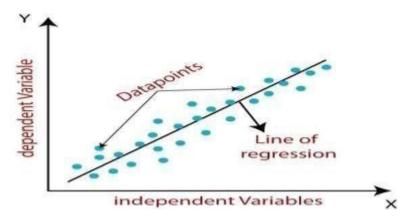


Fig 1.3.2-Linear Regression

2. Naive Bayes classifiers:

Naive Bayes classifiers is a supervised machine learning model for constructing classifiers or models that assign class labels to problem instances, represented as vectors of feature values, where the class labels are drawn from some finite set. It assumes that the factors are independent of each other and there is no correlation between features. As assumption of features are being uncorrelated, it gets its name as Naive Bayes.

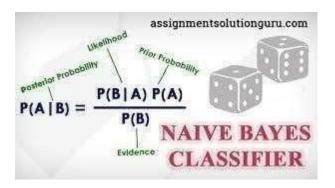


Fig 1.3.3-Naive Bayes Classifier

Where,

p(A|B): Probability of event A given event B has already

occurredp(B|A): Probability of event B given event A has

already occurredp(A): Probability of event A

p(B): Probability of event B

3. Logistic Regression:

Logistic Regression is a supervised learning algorithm which is mostly used in binary classification problems. Even when regression contradicts with classification, here the spotis for logistic that refers to logistic function which does the classification task. It is simple but effective classification algorithms most commonly used for binary classification problems. It is also called sigmoid function.

Sigmoid Function:
$$y = \frac{1}{1 + e^{-x}}$$

Logistic regression takes linear equation as input and uses sigmoid function and logs oddsto performs a binary problem. As a result 's' or sigmoid curve will be obtained as the output.

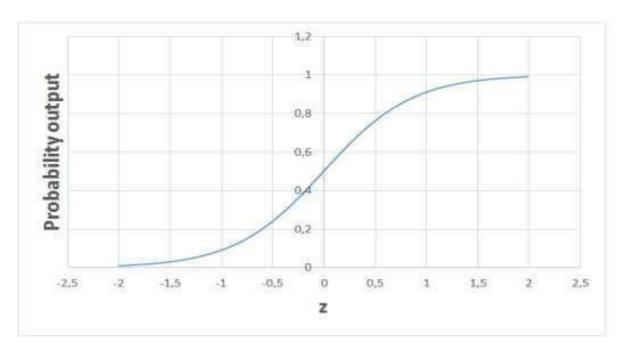


Fig 1.3.4 – Logistic Regression

4. Decision Trees:

Decision trees are a non parametric supervised learning method used for classification andregression. The goal is to create a model that predicts the value of a target variable by learning simple decision rules inferred from the data features. A tree can be seen as a piecewise constant approximation. Though it achieves high accuracy with training set but poorly on new. The depth of the tree is controlled by max_depth parameter for decision treealgorithm in scikit-learn

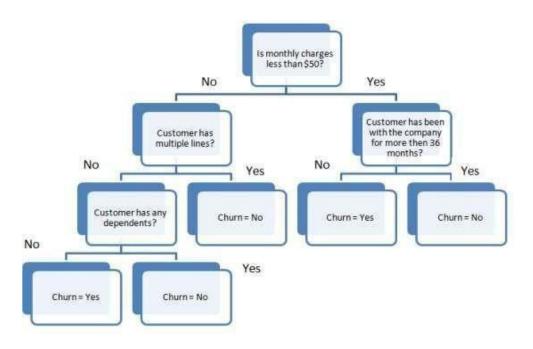


Fig 1.3.5-Decision Tree

5. Random Forest:

Random forests or Random decision forests is an ensemble learning method for classification, regression and other tasks that operates by constructing a multitude of decisiontrees at training time. For classification tasks, the output of the random forests is the class selected by the most trees.

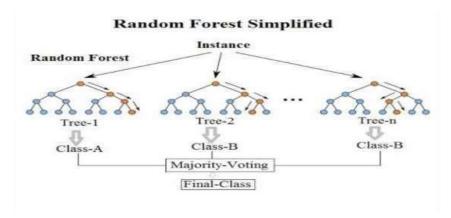


Fig1.3.6-Random Forest

6. K Means Clustering:

K-means clustering is a method of vector quantization, originally from signal processing, that aims to partition n observations into k clusters in which each observation belongs to the cluster with the nearest mean, serving as a prototype of the cluster. It was a way to group ofset of data points are together. Thus, they took look for dissimilarities or similarities among data points. It is an unsupervised learning so there is no label associated with data points. They try to find the underlying structures of the data. Clustering is not classification.

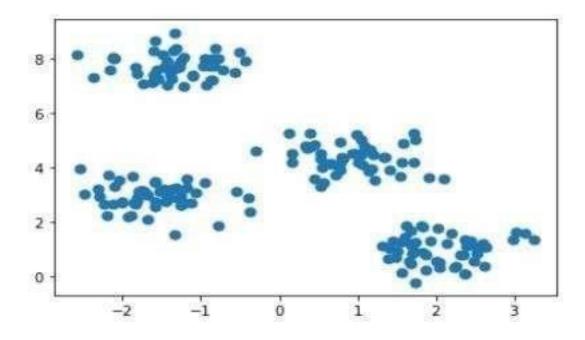


Fig 1.3.7-K Means Clustering

CHAPTER-2

AIM AND SCOPE OF PRESENT INVESTIGATION

2.1 AIM:

To predict the forest fires from the given forest fires dataset.

2.2 SCOPE:

There is a significant scope for predicting forest fires using advanced technologies such as satellite imagery, artificial intelligence (AI), and machine learning (ML) algorithms. These tools can help monitor and analyze various environmental factors that contribute to the risk of forest fires, such as temperature, humidity, wind speed and direction, vegetation moisture content, and topography. Satellite imagery, in particular, can provide valuable information about the extent and severity of forest fires, as well as the location and distribution of smoke and other airborne pollutants. AI and ML algorithms can also be used to analyze historical data and real-time environmental data to identify patterns and predict the likelihood of a forest fire in a particular region.

In the given dataset,

- Data set characteristics are Multivariate
- Attribute characteristics are categorical, integer.
- Associated task Classification
- no.of instances 122
- no.of attributes 14

2.3 DATA PREPARATION:

In this project, Python serves as the key tool which carries out important machine learningalgorithms. With the help of Anaconda Navigator, a desktop Graphical User Interface(GUI), a web based interactive application called Jupyter Notebook that allows editing and running notebook documents via web browser. It is an incredibly powerful

tool for interactively developing and presenting Machine learning and Data Science projects. After importing the required libraries, the dataset will be read in the notebook with the help of data frame (two-dimensional labeled data structure with columns of potentially different types) and read_csv(desired file type).

2.4 READING THE DATASET:

Fig 2.4.1 -syntax for reading the dataset

The data set provided cannot always be fully valued set, in that case we need to prepare the data in such way the machine understands what is the value that has been entered.

	day	month	year	Temperature	Relative_Humidity	Wind_Speed	Rain	Fine_Fuel_Moisture_Code	Druff_Moisture_Code	Drought_Code	Initial_Spread_Index
0	1	6	2012	32	71	12	0.7	57.1	2.5	8.2	0.6
1	2	6	2012	30	73	13	4.0	55.7	2.7	7.8	0.6
2	3	6	2012	29	80	14	2.0	48.7	2.2	7.6	0.3
3	4	6	2012	30	64	14	0.0	79.4	5.2	15.4	2.2
4	5	6	2012	32	60	14	0.2	77.1	6.0	17.6	1.8
											1

Fig 2.4.2- imported estimated forest fires dataset in jupyter notebook

day	0
month	0
Temperature	0
Relative_Humidity	0
Wind_Speed	0
Rain	0
Fine_Fuel_Moisture_Code	0
Druff_Moisture_Code	0
Drought_Code	0
<pre>Initial_Spread_Index</pre>	0
Buildup_Index	0
Fire_Weather_Index	0
Output	0
dtype: int64	

Fig 2.4.3 -using isnull()function checking for null values

CHAPTER 3 EXPERIMENTAL OR MATERIAL AND METHODS ALGORITHMS USED

The given data set is in the form of classification algorithm. So we used classification types to predict the accuracy

3.1 TYPES OF CLASSIFICATIONS ALGORITHMS USED:

- 1. Decision Tree Algorithm
- 2. Random Forests Algorithm

3.2 DECISION TREE ALGORITHM:

Decision Tree is a decision support tool that uses a tree like model of decisions and theirpossible consequences, including chance event outcomes, resource costs, and utility. It is one way to display an algorithm that only contains conditional control statements.

The goal is to create a model that predicts the value of a target variable by learning simpledecision rules inferred from the data features. A tree can be seen as a piecewise constantapproximation.

For instance, decision trees learn from data to approximate a sine curve with a set of if- then-else decision rules. The deeper the tree, the more complex the decision rules and the fitter the model.

Some of the advantages of decision trees are being simple to understand and to interpretand it requires very little data preparation whereas other methods require data normalization, dummy variables etc.

Decision Tree classifier is a class capable of performing multi-class classification on a dataset. As with other classifiers, Decision Tree Classifier takes as input two arrays: an array X, sparse or dense, of shape holding the training samples, and an array Y of integervalues, shape holding the class labels for the training samples.

Why use Decision Trees?

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

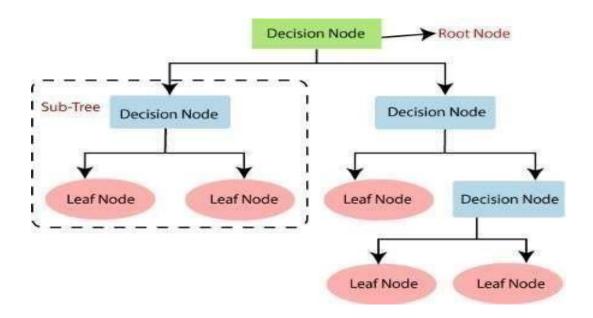


Fig 3.2.1-Decision Tree Algorithm

Decision Tree Terminologies

Root Node: Root node is from where the decision tree starts. It represents the entire dataset, which further gets divided into two or more homogeneous sets.

Leaf Node: Leaf nodes are the final output node, and the tree cannot be segregated further after getting a leaf node.

Splitting: Splitting is the process of dividing the decision node/root node into subnodes according to the given conditions

Parent/Child node: The root node of the tree is called the parent node, and other nodes are called the child nodes.

How does the Decision Tree algorithm Work?

In a decision tree, for predicting the class of the given dataset, the algorithm starts from the root node of the tree. This algorithm compares the values of root attribute with the record (real dataset) attribute and, based on the comparison, follows the branch and jumps to the next node. For the next node, the algorithm again compares the attribute value with the other sub-nodes and move further. It continues the process until it reachesthe leaf node of the tree. The complete process can be better understood using the below algorithm:

- **Step-1:** Begin the tree with the root node, says S, which contains the complete dataset.
- **Step-2**: Find the best attribute in the dataset using Attribute Selection Measure (ASM).
- Step-3: Divide the S into subsets that contains possible values for the best attributes.
- **Step-4:** Generate the decision tree node, which contains the best attribute.
- **Step-5**: Recursively make new decision trees using the subsets of the dataset created in step -3. Continue this process until a stage is reached where you cannotfurther classify the nodes and called the final node as a leaf node.

Example: Suppose there is a candidate who has a job offer and wants to decide whetherhe should accept the offer or not. So, to solve this problem, the decision tree starts with the root node (Salary attribute by ASM). The root node splits further into the next decision node (distance from the office) and one leaf node based on the corresponding labels. The next decision node further gets split into one decision node (Cab facility) and one leaf node.

Finally, the decision node splits into two leaf nodes (Accepted offers and Declined offer).

3.3 RANDOM FORESTS:

Random forests or Random decision forests is an ensemble learning method for classification, regression and other tasks that operates by constructing a multitude of decision trees at training time. For classification tasks, the output of the random forest is the class selected by most trees. It is a supervised learning machine learning algorithm made up of decision trees which are used in classification and regression. It plays a majorrole in our model. It is called a "forest" because it grows a forest of decision trees. The data from these trees are then merged together to ensure the most accurate predictions. While a solo decision tree has one outcome and a narrow range of groups, the forest assures a more accurate result with a bigger number of groups and decisions. It has the added benefit of adding randomness to the model by finding the best feature among a random subset of features. Overall, these benefits create a model that has wide diversity that many data scientists favor.

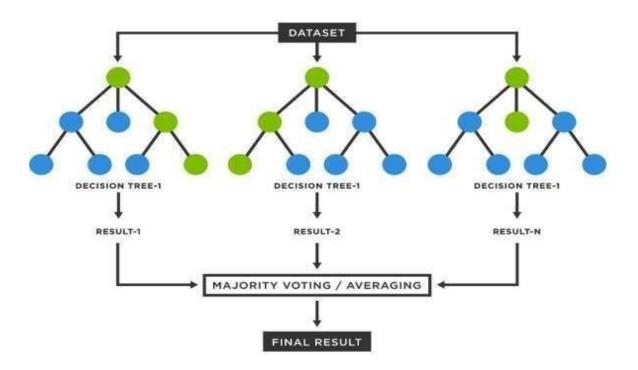


Fig 3.3.1 -Random Forest Algorithm

Assumptions for Random Forest

Since the random forest combines multiple trees to predict the class of the dataset, it is possible that some decision trees may predict the correct output, while others may not. But together, all the trees predict the correct output. Therefore, below are two assumptions for a better Random Forest classifier:

There should be some actual values in the feature variable of the dataset so that the classifier can predict accurate results rather than a guessed result.

The predictions from each tree must have very low correlations.

Why use Random Forest?

- Below are some points that explain why we should use the Random Forestalgorithm. It takes less training time as compared to other algorithms.
- It predicts output with high accuracy, even for the large dataset it runs efficiently.
- It can also maintain accuracy when a large proportion of data is missing.

How does Random Forest algorithm work?

Random Forest works in two-phase first is to create the random forest by combining N decision tree, and second is to make predictions for each tree created in the first phase. The Working process can be explained in the below steps and diagram:

- Step-1: Select random K data points from the training set.
- Step-2: Build the decision trees associated with the selected data points 13
- Step-3: Choose the number N for decision trees that you want to build.
- Step-4: Repeat Step 1 & 2.
- Step-5: For new data points, find the predictions of each decision tree, and assign the new data points to the category that wins the majority votes.

Applications of Random Forest:

There are mainly four sectors where Random Forest mostly used:

- 1. Banking: Banking sector mostly uses this algorithm for the identification of loan risk.
- 2. Medicine: With the help of this algorithm, disease trends and risks of the disease canbe identified.
- 3. Land Use: We can identify the areas of similar land use by this algorithm.
- 4. Marketing: Marketing trends can be identified using this algorithm.

Advantages of Random Forest:

- Random Forest is capable of performing both Classification and Regression tasks.
- It is capable of handling large datasets with high dimensionality.
- It enhances the accuracy of the model and prevents the overfitting issue.

Disadvantages of Random Forest:

Although random forest can be used for both classification and regression tasks, it is not more suitable for Regression tasks.

3.4 MACHINE LEARNING LIBRARIES:

Libraries are collections of prewritten code that users can use to optimize tasks. In projetas python is used for implementation tool, it has the most libraries as compared to otherprogramming languages. More than of 60% machine learning developers use and goes for python as it is easy to learn. As python has comparatively large collection of librarieslets look at the libraries that came handy for our dataset.

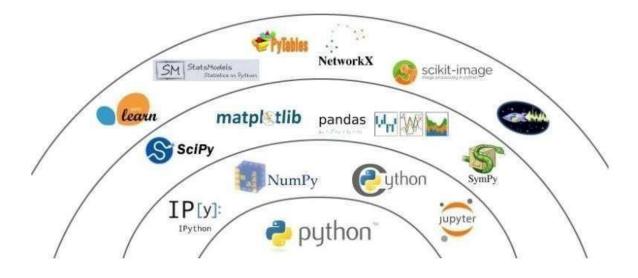


Fig 3.4.1-Various python libraries for machine learning

1. Sklearn:

Sklearn stands for Scikit-learning, a machine learning library. It is imported for various classification, regression and clustering algorithms including k-means, random forest, support vector machines, gradient boosting and DBSCAN. It is designed using libraries Numpy and Scipy. From the sklearn library and from the tree inside the library DecisionTreeClassifier.It is a class capable of performing multi-class classifier on a dataset. When compared with other classifiers, DecisionTreeClassifier takes input as two arrays:an array X, aparse or dense,of shape(n_samples,n_features) holding training samples and an array Y of integer values, shape holding class labels for training sample. From sklearn another one called model selection for training and testing the model imports train_test_split. It is a method setting a blueprint to analyze data and using it to measure new data. Selecting a proper model allows to generate accurate results while making prediction. For proceeding, we need to train the model by using a specific dataset and test the model by using a specific dataset and test the model against another dataset. By default, sklearntrain test split will make random partitions for two subsets. We can also specify a random state for the operation. First, we need to split the dataset and then allocate the size for train and test.

2. Math:

Math is a built-in module that you can use for mathematical tasks. It has set of methods and constants. It is a standard module in python and is always available

3. Seaborn:

Seaborn is a library built on top of matplotlib. It is used for data visualization and exploratory data analysis. They work easily with dataframes and pandas library. The graphs created can also be customized easily. It provide default styles and color palettesto make statistical plots more attractive. Also closely integrated to the data structures from pandas

4. Matplotlib.pyplot:

Matplotlib.pyplot is a state-based interface to matplotlib. It provides a MATLAB-like way of plotting. It make changes to figures.

3.5 IMPORTED LIBRARIES

1.Pandas:

Pandas is a widely used data analysis and manipulation library for python. It provides a lot of functions and methods that expedite the data analysis and preprocessing steps. It also provides fast, flexible and expressive data structures working with relational or labeled or both easy and intuitive. Considered as fundamental high-level building block in performing practical, real world data analysis in python. It has Data Frame and series for analyzing.

Fig 3.5.1- Pandas library is used to read the dataset

2.Numpy:

Numpy stands for Numerical python, is a library consisting of multidimensional array objects and a collection of countless of routines for processing those arrays. Using this mathematical and logical operations on arrays can be performed. The difference in Numpy from Pandas is, it works on numerical data whereas pandas on tabular data.

```
import numpy as np
import pandas as pd
import seaborn as sns
import matplotlib.pyplot as plt
from matplotlib import pyplot as plt
import sklearn
from sklearn import tree
from sklearn.model_selection import train_test_split
from sklearn.tree import DecisionTreeClassifier
from sklearn.metrics import confusion_matrix
from sklearn.metrics import accuracy_score
from sklearn.metrics import classification_report
```

Fig 3.5.2 – Imported libraries in jupyter notebook

3.6 IMPLEMENTATION OF DECISION TREE ALGORITHM

Decision trees are a non parametric supervised learning method used for classification and regression. The goal is to create a model that predicts the value of a target variable by learning simple decision rules inferred from the data features. A tree can be seen as a piecewise constant approximation. Though it achieves high accuracy with training set but poorly on new. The depth of the tree is controlled by max_depth parameter for decision tree algorithm in scikit-learn. For instance, decision trees learn from data to approximate a sine curve with a set of if-then-else decision rules. The deeper the tree, the more complex the decision rules and the fitter the model. Some of the advantages ofdecision trees are being simple to understand and to interpret and it requires very little data preparation whereas other methods require data normalization, dummy

variables etc.

- 1. Import the packages and classes you need.
- 2. Provide data to work with and eventually do appropriate transformations.
- 3. Create a classification model and fit it with existing data.
- 4. Check the results of model fitting to know whether the model is satisfactory.
- 5. Apply the model of predictions.

	day	month	year	Temperature	Relative_Humidity	Wind_Speed	Rain	Fine_Fuel_Moisture_Code	Druff_Moisture_Code	Drought_Code	Initial_Spread_Inde
0	1	6	2012	32	71	12	0.7	57.1	2.5	8.2	0.
1	2	6	2012	30	73	13	4.0	55.7	2.7	7.8	0.
2	3	6	2012	29	80	14	2.0	48.7	2.2	7.6	0.
3	4	6	2012	30	64	14	0.0	79.4	5.2	15.4	2.
4	5	6	2012	32	60	14	0.2	77.1	6.0	17.6	1.
17	26	9	2012	30	65	14	0.0	85.4	16.0	44.5	4.
18	27	9	2012	28	87	15	4.4	41.1	6.5	8.0	0.
19	28	9	2012	27	87	29	0.5	45.9	3.5	7.9	0.
20	29	9	2012	24	54	18	0.1	79.7	4.3	15.2	1.
21	30	9	2012	24	64	15	0.2	67.3	3.8	16.5	1.2

Fig 3.6.1-Given dataset

```
X = np.asarray(forest_data[['month', 'Temperature', 'Relative_Humidity', 'Wind_Speed', 'Fine_Fuel_Moisture_Code', 'Druff_Moisture
Y = np.asarray(forest_data['Output'])
X_train, X_test, y_train, y_test = train_test_split(X, Y, test_size = 3, random_state = 100)

gini_classifier = DecisionTreeClassifier(criterion = "gini", random_state = 100, max_depth=None, min_samples_leaf=5)
gini_classifier.fit(X_train, y_train)
DecisionTreeClassifier(min_samples_leaf=5, random_state=100)
```

Fig 3.6.2-Applying decision tree algorithm

3.7 IMPLEMENTATION OF RANDOM FORESTS ALGORITHM:

Random forests or Random decision forests is an ensemble learning method for classification, regression and other tasks that operates by constructing a multitude of decision trees at training time. For classification tasks, the output of the random forest is the class selected by most trees. It is a supervised learning machine learning algorithm made up of decision trees which are used in classification and regression. It plays a majorrole in our model. As we know that a forest is made up of trees and more trees means more robust forest. Similarly, random forest algorithm creates decision trees on data samples and then gets the prediction from each of them and finally selects the best solution by means of voting. It is an ensemble method which is better than a single decision tree because it reduces the over-fitting by averaging the result.

random_classifier.fit(X_train, y_train)

Fig 3.7.1 Applying random forest algorithm

CHAPTER 4

RESULTS AND DISCUSSION, PERFORMANCE ANALYSIS

4.1 MODEL ANALYSIS

The above algorithms are written in python with the help of numpy and executed using jupyter notebook. The accuracy rate of decision tree algorithm when executed for the given dataset i.e Forest Fires prediction Dataset is 100.00 and for random forest is 100.00

The dataset contains 122 training samples with 14 features (day, month, year, temperature, relative humidity, wind speed, rain, fine fuel, druff moisture code, drought code, initial spread index, buildup index, fire weather index, output) our task is to determine the accuracy of working of decision tree and random forest algorithm when trained with the provided dataset

ACCURACY, PRECISION, RECALL, F1-SCORE

```
y_pred = gini_classifier.predict(X_test)
print("Predicted Values using Gini: ")
print(y_pred)
print()
print("Results using Gini")
print("Confusion Matrix: ", confusion_matrix(y_test, y_pred))
print ("Accuracy : ", accuracy_score(y_test,y_pred)*100)
print("Report : ", classification_report(y_test, y_pred))
Predicted Values using Gini:
[1 1 0]
Results using Gini
Confusion Matrix: [[1 0]
 [0 2]]
Accuracy : 100.0
Report :
                          precision recall f1-score support
             0 1.00 1.00
1 1.00 1.00
                                    1.00
                                              1.00
                                              1.00
                                                                3
    accuracy
accuracy 1.00 macro avg 1.00 1.00 1.00 weighted avg 1.00 1.00 1.00
```

Fig 4.1.1 Accuracy, precision, recall, f1-score for decision tree algorithm

```
print("ACCURACY OF THE MODEL: ", metrics.accuracy_score(y_test, y_pred))
```

ACCURACY OF THE MODEL: 1.0

Fig 4.1.2 Accuracy of Random forest algorithm

CHAPTER 5 5.1 SUMMARY AND CONCLUSIONS

Our goal was to predict diabetes risk accurately using Random forests and Decision Tree algorithms. Here, we will train both the models and test the accuracy of them. There was a significant improvement in the accuracy levels upto 100 in decision tree classification and 100 in random forest classifications respectively. we got a significant accuracy while training the model using random forest classifier. While the accuracy is high, the model is not substantially overfitted, as the cross-validation scores for each of the ensemble methods all differed from the model accuracy by less than 1%. Also, we concluded that f1_score, precision, recall by using these random forest and decision tree classifier.

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REFERENCES

[1] Refferal code for forest fires

https://www.kaggle.com/datasets/elikplim/forest-fires-data-set/code

[2] Dataset

https://drive.google.com/drive/folders/1xf6uwCPMXsHxHhpax8agm4dkJQWNavDE

[3] Dataset Information

https://archive.ics.uci.edu/ml/datasets/Algerian+Forest+Fires+Dataset++

WORKING ENVIRONMENT

ANACONDA NAVIGATOR is desktop GUI used to launch applications and also manage packages in one place.

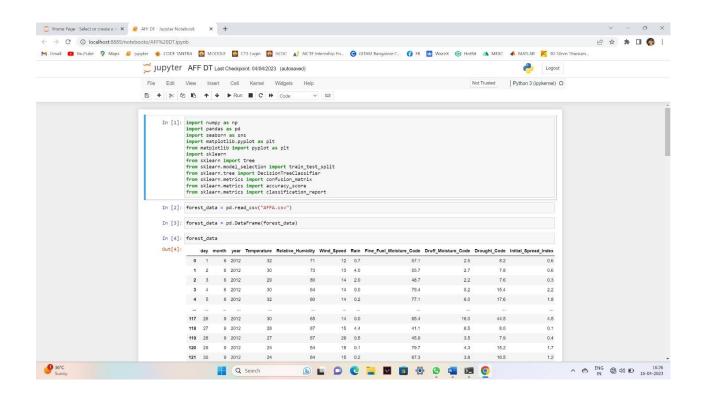


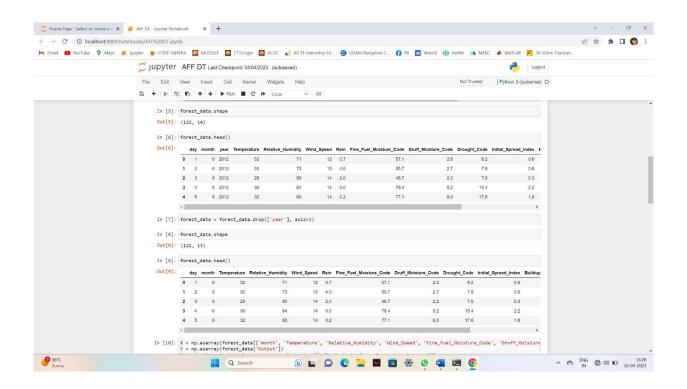
CODING ENVIRONMENT

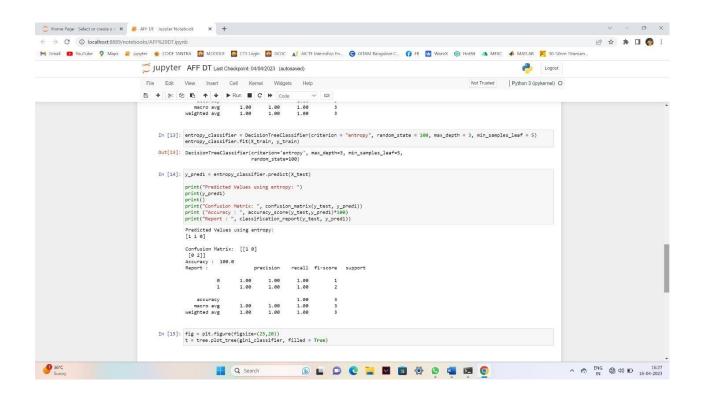
Jupyter notebook from the anaconda navigator is launched along with all the preinstalled packages for python.

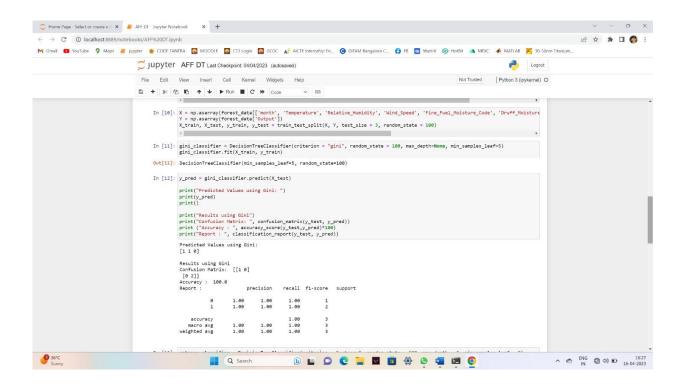


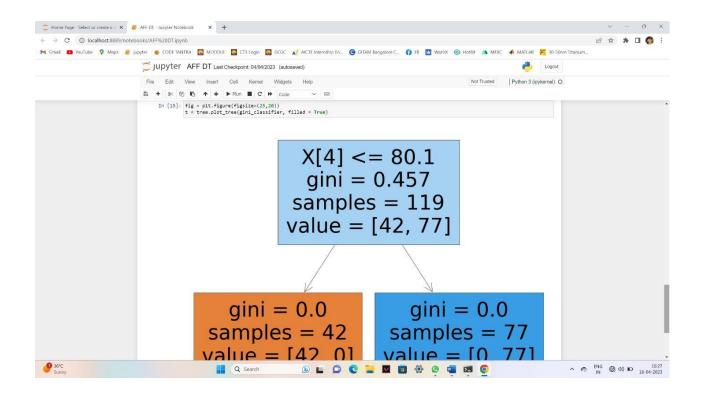
SCREENSHOTS AND OUTPUTS

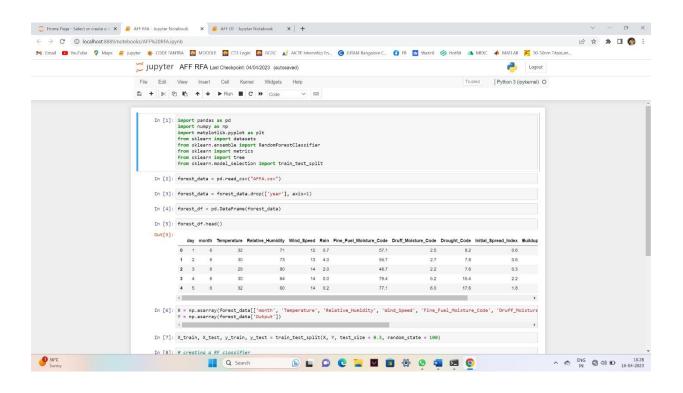


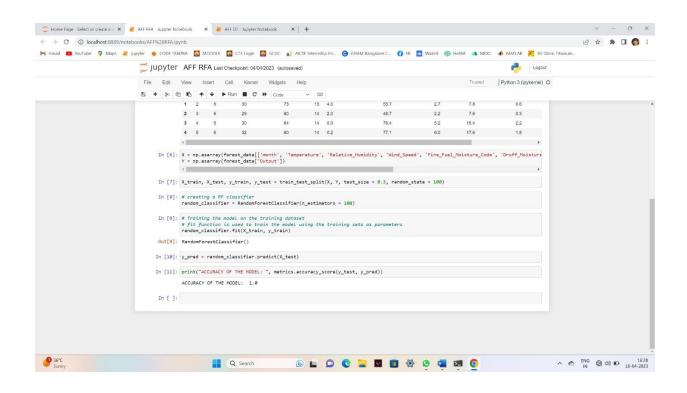












SOURCE CODE

```
import numpy as np
import pandas as pd
import seaborn as sns
import matplotlib.pyplot as plt
from matplotlib import pyplot as plt
import sklearn
from sklearn import tree
from sklearn.model_selection import train_test_split
from sklearn.tree import DecisionTreeClassifier
from sklearn.metrics import confusion_matrix
from sklearn.metrics import accuracy_score
from sklearn.metrics import classification_report
from sklearn.ensemble import RandomForestClassifier
from sklearn import metrics
from sklearn.model selection import train test split
forest_data = pd.read_csv("AFFA.csv")
forest data = pd.DataFrame(forest data)
forest data
forest_data.shape
forest_data.head()
forest_data = forest_data.drop(['year'], axis=1)
forest data.shape
forest_data.head()
X = np.asarray(forest_data[['month', 'Temperature', 'Relative_Humidity', 'Wind_Speed',
'Fine_Fuel_Moisture_Code', 'Druff_Moisture_Code', 'Drought_Code',
'Initial_Spread_Index', 'Buildup_Index', 'Fire_Weather_Index']])
Y = np.asarray(forest_data['Output'])
X_train, X_test, y_train, y_test = train_test_split(X, Y, test_size = 3, random_state = 100)
gini_classifier = DecisionTreeClassifier(criterion = "gini", random_state = 100,
max_depth=None, min_samples_leaf=5)
gini_classifier.fit(X_train, y_train)
y_pred = gini_classifier.predict(X_test)
```

```
print("Predicted Values using Gini: ")
print(y_pred)
print()
print("Results using Gini")
print("Confusion Matrix: ", confusion_matrix(y_test, y_pred))
print ("Accuracy : ", accuracy_score(y_test,y_pred)*100)
print("Report : ", classification_report(y_test, y_pred))
entropy_classifier = DecisionTreeClassifier(criterion = "entropy", random_state = 100,
max_depth = 3, min_samples_leaf = 5)
entropy_classifier.fit(X_train, y_train)
y_pred1 = entropy_classifier.predict(X_test)
print("Predicted Values using entropy: ")
print(y_pred1)
print()
print("Confusion Matrix: ", confusion_matrix(y_test, y_pred1))
print ("Accuracy : ", accuracy_score(y_test,y_pred1)*100)
print("Report : ", classification_report(y_test, y_pred1))
fig = plt.figure(figsize=(25,20))
t = tree.plot_tree(gini_classifier, filled = True)
X = np.asarray(forest_data[['month', 'Temperature', 'Relative_Humidity', 'Wind_Speed',
'Fine_Fuel_Moisture_Code', 'Druff_Moisture_Code', 'Drought_Code',
'Initial_Spread_Index', 'Buildup_Index', 'Fire_Weather_Index']])
Y = np.asarray(forest_data['Output'])
X_train, X_test, y_train, y_test = train_test_split(X, Y, test_size = 0.3, random_state =
100)
random_classifier = RandomForestClassifier(n_estimators = 100)
random_classifier.fit(X_train, y_train)
y_pred = random_classifier.predict(X_test)
print("ACCURACY OF THE MODEL: ", metrics.accuracy_score(y_test, y_pred))
                                     ----THE END-----
```