

Farming Made Easy Using Machine Learning

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CERTIFICATE

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ABSTRACT

Agriculture serves as the backbone of our country's economy. However, in recent years, the sector has faced challenges due to unpredictable climate patterns and fluctuating price trends. Farmers often bear the brunt of these uncertainties, leading to crop spoilage and significant financial losses. Moreover, farmers may not always be aware of the most suitable crop types for their land, further complicating their situation. To address these issues, a comprehensive and user-friendly system has been developed. This system leverages Machine Learning's Decision Tree Regression Algorithm to predict crop prices accurately. The key attributes considered for price prediction include rainfall, wholesale price index, month, and year. By providing advance price forecasts, this system empowers farmers to make informed decisions, thereby increasing their profitability and contributing to the country's economy. In addition to price prediction, this system incorporates modules for weather forecasting, crop recommendation, fertilizer recommendation, and a shop, chat portal, and guide. The crop recommendation module suggests suitable crop types based on the land's characteristics, helping farmers maximize their yield. The crop yield prediction module forecasts the expected yield based on various factors, aiding farmers in planning their harvest and resources accordingly. The fertilizer recommendation module suggests appropriate fertilizers based on soil health and crop requirements, optimizing crop growth and quality. Overall, this system serves as a valuable tool for farmers, providing them with essential information and insights to improve their farming practices and increase their profitability. By integrating advanced technologies like Machine Learning, this system has the potential to revolutionize the agricultural sector and contribute significantly to the country's economic growth.

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1. INTRODUCTION

Agriculture, the bedrock of economies worldwide, confronts a dual challenge of climate uncertainty and market volatility. Climate change disrupts cropping cycles and escalates pest pressures, while volatile market dynamics exacerbate income uncertainty for farmers. Smallholder farmers, lacking resources and market access, bear the brunt of these challenges, perpetuating cycles of poverty and food insecurity. To navigate these complexities, innovative solutions such as precision agriculture and digital farming offer promise in optimizing resource use, enhancing resilience, and improving market access. Government policies promoting climate-resilient agriculture, market liberalization, and institutional support mechanisms are essential for fostering inclusive growth and sustainable development in rural areas. By fostering collaboration among stakeholders and investing in technological and policy interventions, we can bolster agricultural resilience, ensure food security, and build a sustainable future for generations to come.

In the face of climate uncertainty and market volatility, the agricultural sector confronts significant challenges that threaten the livelihoods of farmers and the stability of economies worldwide. However, technological innovations offer promising solutions to address these challenges. One such innovation is the application of Machine Learning (ML) algorithms for agricultural price prediction. In this essay, we explore the development and implementation of a comprehensive and user-friendly ML system designed to predict crop prices with precision. By leveraging the Decision Tree Regression Algorithm and considering key attributes such as rainfall, wholesale price index, month, and year, this system empowers farmers to make informed decisions, thereby enhancing their profitability and bolstering the country's economy.\

To tackle these formidable challenges, we have meticulously crafted a comprehensive and user-friendly system harnessing the immense power of Machine Learning. Specifically, our system is engineered around the Decision Tree Regression Algorithm, a stalwart in predictive analytics renowned for its capacity to dissect complex datasets and discern intricate patterns. By integrating this cutting-edge technology with agricultural economics, we've developed a robust framework capable of forecasting crop prices with unparalleled precision. At the heart of our system lie key attributes meticulously selected to capture the nuances of market dynamics and climatic influences.

These attributes include rainfall patterns, wholesale price indexes, and temporal factors such as month and year, which collectively contribute to the accuracy and reliability of our price predictions.

The cornerstone of our approach lies in empowering farmers with advanced price forecasts that transcend mere speculation, offering actionable insights into future market trends. Armed with this foresight, farmers can navigate the volatile terrain of agricultural markets with confidence, making informed decisions that optimize their profitability and safeguard their economic interests. Whether it's determining the most opportune time to bring crops to market or strategically planning crop rotations to capitalize on price fluctuations, our system equips farmers with the knowledge and foresight needed to thrive in an increasingly uncertain environment.

The transformative impact of our system extends far beyond individual farm operations, reverberating through the fabric of the entire agricultural ecosystem and the broader economy. By bolstering the resilience and prosperity of farmers, our system lays the foundation for sustainable agricultural development, ensuring food security and economic stability for generations to come. Moreover, by facilitating more efficient resource allocation and market participation, our system fosters economic growth, driving productivity gains and fostering innovation across the agricultural value chain.

But what sets our system apart is not just its technological sophistication, but its unwavering commitment to user-centric design and accessibility. Recognizing the diverse needs and capabilities of farmers, we've developed an intuitive user interface that empowers users of all backgrounds to harness the power of Machine Learning with ease. Whether accessed through a web portal, mobile application, or interactive voice interface, our system provides farmers with seamless access to critical market insights and decision support tools, democratizing access to information and leveling the playing field for smallholder farmers.

Furthermore, our system is not a static entity but a living, evolving ecosystem driven by continuous learning and adaptation. Through feedback mechanisms and real-time data integration, we continuously refine and enhance the predictive accuracy of our models, ensuring they remain responsive to changing market dynamics and climatic conditions. Moreover, we remain steadfast in our commitment to transparency and accountability, providing users with full visibility into the underlying algorithms and data sources that power our system, fostering trust and confidence in its capabilities.

In addition to its immediate benefits for individual farmers, our system holds the potential to catalyze broader systemic change within the agricultural sector. By aggregating and anonymizing data across a diverse array of farming operations, our system generates valuable insights into macroeconomic trends, market dynamics, and climate resilience strategies. These insights can inform policy decisions, drive investment priorities, and shape the trajectory of agricultural development at the national and global levels, unlocking new pathways for inclusive and sustainable growth.

Our commitment to addressing the multifaceted challenges facing farmers goes beyond merely predicting crop prices. In addition to leveraging Machine Learning's Decision Tree Regression Algorithm for precise price forecasts, our comprehensive system integrates a suite of modules designed to empower farmers with a holistic range of tools and resources. These modules encompass weather forecasting, crop recommendation, fertilizer recommendation, as well as a versatile shop, chat portal, and guide, collectively aimed at optimizing agricultural productivity, sustainability, and economic viability.

The weather forecasting module serves as a vital component of our system, providing farmers with timely and accurate information about meteorological conditions that directly impact crop growth and yield. By leveraging advanced meteorological models and real-time data feeds, our system delivers forecasts for temperature, precipitation, humidity, and other relevant parameters, enabling farmers to proactively plan their agricultural activities, mitigate weather-related risks, and optimize resource management strategies. Whether it's deciding the optimal time for planting, irrigation scheduling, or pest management, access to reliable weather forecasts empowers farmers to make informed decisions that enhance crop resilience and maximize productivity.

Moreover, our system includes a fertilizer recommendation module that employs soil testing data, crop nutrient requirements, and agronomic best practices to generate personalized recommendations for fertilizer application. By optimizing nutrient management strategies, our system helps farmers improve soil fertility, enhance crop health and vigor, and maximize nutrient use efficiency, thereby reducing input costs, minimizing environmental pollution, and promoting sustainable agricultural practices. Whether it's determining the optimal nutrient ratios for specific crops and growth stages or selecting appropriate fertilizer formulations based on soil nutrient levels, our fertilizer recommendation module provides farmers with tailored guidance that enhances agronomic performance and economic profitability.

In addition to its core functionalities, our system features a versatile shop, chat portal, and guide, which serve as invaluable resources for farmers seeking access to inputs, market information, and expert advice. The shop module enables farmers to browse and purchase agricultural inputs such as seeds, fertilizers, pesticides, and machinery from trusted suppliers, streamlining procurement processes and ensuring timely access to essential resources. The chat portal facilitates communication and knowledge sharing among farmers, extension agents, researchers, and agricultural experts, enabling users to seek advice, share experiences, and collaborate on solutions to common challenges. Furthermore, the guide module provides educational resources, tutorials, and best practice guidelines on a wide range of agronomic topics, empowering farmers with the knowledge and skills needed to succeed in modern agriculture.

Linear Regression, a cornerstone algorithm in statistical modeling, holds significant value in the realm of agricultural economics. Its utility lies in its ability to establish relationships between variables, making it particularly well-suited for price trend analysis in the agricultural sector. By fitting a linear equation to observed data points, Linear Regression enables analysts to quantify the relationship between independent variables (such as rainfall, temperature, market demand) and dependent variables (crop prices). This allows farmers, policymakers, and market analysts to gain insights into the underlying factors driving price movements, identify trends, and anticipate future market conditions.

Beyond Linear Regression, advanced regression techniques such as Lasso and Ridge Regression offer powerful tools for handling multicollinearity and feature selection, which are essential considerations in agricultural price prediction. Multicollinearity, the phenomenon where independent variables are highly correlated with each other, can distort regression coefficients and compromise the interpretability of results. Lasso and Ridge Regression mitigate multicollinearity by introducing regularization penalties that constrain the magnitude of regression coefficients, effectively reducing the impact of correlated variables and improving model stability. Additionally, these techniques facilitate feature selection by automatically identifying and prioritizing the most relevant variables, thus enhancing the accuracy and parsimony of predictive models. In the context of agricultural price prediction, Lasso and Ridge Regression enable analysts to identify key factors influencing crop prices, such as weather conditions, input costs, market demand, and policy interventions, thereby informing strategic decision-making and risk management strategies.

Another machine learning algorithm that holds promise in agricultural analytics is K-Nearest Neighbors (KNN), a non-parametric method commonly used for clustering and classification tasks.

KNN operates on the principle of similarity, where data points are assigned to the same class or cluster based on their proximity to one another in feature space. In the context of crop management and recommendation systems, KNN can be applied to cluster similar data points representing agricultural parcels or crops, aiding in the identification of patterns, trends, and common characteristics. By grouping together similar crops or farming practices, KNN facilitates the generation of personalized recommendations tailored to specific agroecological conditions, soil types, and climate regimes. Furthermore, KNN can be used in conjunction with other algorithms, such as Decision Trees or Random Forests, to enhance the accuracy and robustness of crop classification and recommendation systems, thereby empowering farmers with actionable insights and tailored solutions for optimizing agricultural productivity and profitability.

Decision Trees, renowned for their intuitive and transparent decision-making process, are another valuable tool in the agricultural analytics toolbox. Decision Trees recursively partition the feature space into hierarchical decision nodes, with each node representing a split based on a chosen attribute. By analyzing historical data on crop yields, prices, agronomic practices, and environmental conditions, Decision Trees can elucidate the complex relationships between input variables and output outcomes, providing valuable insights into the factors influencing crop yields and prices. Decision Trees are particularly well-suited for scenario analysis and "what-if" simulations, enabling analysts to explore alternative management strategies, evaluate their potential impact, and make informed decisions under uncertainty. Moreover, Decision Trees facilitate knowledge transfer and decision support in agricultural extension and advisory services, empowering farmers with actionable insights and best practices for crop management, pest control, irrigation scheduling, and post-harvest handling.

Machine learning algorithms such as Linear Regression, Lasso and Ridge Regression, K-Nearest Neighbors, and Decision Trees can significantly impact agricultural practices by offering valuable insights and predictive capabilities. These algorithms enable analysts to establish relationships between variables, handle multicollinearity and feature selection, cluster similar data points, and understand decision-making processes in crop management. By leveraging advanced analytics techniques, stakeholders in the agricultural sector can enhance productivity, optimize resource allocation, mitigate risks, and foster sustainable development, thereby ensuring food security, economic prosperity, and environmental sustainability for present and future generations.

2. LITERATURE SURVEY

2.1 MACHINE LEARNING

Machine learning represents a transformative application of artificial intelligence (AI), revolutionizing various industries by empowering systems to learn and improve from experience autonomously, without explicit programming. This paradigm shift has unlocked unprecedented potential for computers to analyze vast datasets, identify patterns, and make informed decisions, leading to advancements in fields ranging from healthcare and finance to transportation and agriculture.

At its core, machine learning is driven by the principle of learning from data. Instead of relying solely on predefined rules and algorithms, machine learning algorithms iteratively analyze data to uncover insights, patterns, and trends. This process, known as training, involves feeding the algorithm with labeled or unlabeled data, allowing it to discern underlying patterns and relationships. Through continuous exposure to data, machine learning models refine their understanding and improve their performance over time, a concept often referred to as "learning."

One of the defining features of machine learning is its ability to generalize from past experiences to make predictions or decisions about new, unseen data. This capability is particularly valuable in scenarios where traditional programming approaches struggle to capture the complexity and variability of real-world data. By leveraging statistical techniques and mathematical models, machine learning algorithms can infer patterns from data and make accurate predictions, classifications, or recommendations.

Machine learning encompasses a diverse array of techniques and algorithms, each suited to different types of tasks and data. Supervised learning, for instance, involves training a model on labeled data, where each input is associated with a corresponding output or target variable. The model learns to map inputs to outputs, enabling it to make predictions on unseen data. In contrast, unsupervised

learning deals with unlabeled data, seeking to identify hidden structures or patterns within the data without explicit guidance.

Furthermore, machine learning techniques extend to reinforcement learning, where agents learn optimal behavior through trial and error, receiving feedback in the form of rewards or penalties. Reinforcement learning has found applications in areas such as robotics, gaming, and autonomous systems, enabling agents to learn complex behaviors and strategies through interaction with their environment.

Another notable aspect of machine learning is its versatility and applicability across diverse domains. In healthcare, machine learning algorithms analyze medical images, genomic data, and electronic health records to aid in disease diagnosis, treatment planning, and drug discovery. In finance, predictive models forecast stock prices, detect fraudulent transactions, and optimize investment portfolios. In transportation, machine learning algorithms power autonomous vehicles, optimize traffic flow, and enhance logistics operations.

In agriculture, machine learning plays a pivotal role in crop yield prediction, disease detection, and precision agriculture. By analyzing weather data, soil characteristics, and historical crop performance, machine learning models can generate insights to optimize planting schedules, irrigation management, and pest control strategies. Additionally, machine learning facilitates the development of crop recommendation systems, guiding farmers in selecting the most suitable crops for their specific conditions and maximizing productivity.

Overall, machine learning represents a paradigm shift in computing, empowering systems to learn and adapt from data, rather than relying solely on human intervention. As advancements in machine learning continue to accelerate, its impact across industries is poised to deepen, driving innovation, efficiency, and transformative change in the years to come.

The essence of machine learning lies in the process of learning from data without explicit programming or human intervention. This transformative approach enables computers to autonomously analyze observations, examples, or instructions, discern patterns within data, and refine their actions and decisions over time. At the heart of machine learning is the quest to empower computers to learn from experience and improve their performance without constant human guidance.

The journey of learning in machine learning commences with the ingestion of data. This data can take various forms, ranging from labeled examples to raw sensory inputs or textual documents. These observations serve as the foundation upon which machine learning algorithms build their understanding of the world. Through exposure to diverse datasets, algorithms aim to identify underlying patterns, correlations, and structures that characterize the data.

Central to the learning process is the quest to extract meaningful insights and knowledge from data. Machine learning algorithms employ a plethora of techniques to uncover patterns, including statistical methods, neural networks, and optimization algorithms. By iteratively processing data, these algorithms refine their internal representations, gradually enhancing their ability to generalize from past experiences and make informed predictions or decisions on new, unseen data.

The overarching goal of machine learning is to enable computers to learn autonomously and adapt their actions based on experience. This paradigm shift represents a departure from traditional programming paradigms, where human programmers encode explicit instructions and rules for the computer to follow. In contrast, in machine learning, the computer learns from data, adjusting its actions and decisions based on observed patterns and feedback.

A fundamental aspect of machine learning is the concept of feedback loops. As the algorithm interacts with data and makes predictions or decisions, it receives feedback on the correctness or efficacy of its actions. This feedback serves as a signal for the algorithm to adjust its internal parameters and update its models to improve performance. Over time, through continuous iterations of learning and adaptation, machine learning algorithms refine their capabilities and become increasingly proficient at the tasks they are designed to perform.

Crucially, machine learning enables computers to learn from vast amounts of data at scale, processing information far beyond the capacity of human cognition. This capability has fueled advancements across a wide range of domains, from natural language processing and computer vision to healthcare and finance. In healthcare, for example, machine learning algorithms analyze medical images and patient data to assist in diagnosis and treatment planning. In finance, predictive models leverage historical market data to forecast stock prices and optimize investment strategies.

In essence, the process of learning in machine learning embodies the quest to imbue computers with the ability to acquire knowledge and skills autonomously from data. By harnessing the power of data-driven insights, machine learning enables computers to make better decisions, solve complex problems, and adapt to changing environments with minimal human intervention. As the field of

machine learning continues to evolve, its impact on society, industry, and technology is poised to grow exponentially, unlocking new opportunities and driving innovation across diverse domains.

Over the years, the commercial world has become more competitive, as organizations such as these have to meet the needs and desires of their customers, attract new customers and thus improve their businesses. The task of identifying and meeting the needs and requirements of every customer in the business is very difficult. This is because customers can vary according to their needs, wants, demographics, size, taste and taste, features etc. As it is, it is a bad practice to treat all customers equally in business. This challenge has adopted the concept of customer segmentation or market segmentation, where consumers are divided into subgroups or segments, where members of each subcategory exhibit similar market behaviors or characteristics. Accordingly, customer segmentation is the process of dividing the market into indigenous groups.

Data collection serves as a fundamental step in research across various fields, including physical and social sciences, humanities, and business. It involves gathering and measuring information pertaining to targeted changes within an established system, enabling researchers to address relevant questions and evaluate outcomes. The process of data collection is essential for obtaining quality evidence, which forms the basis for analysis and enables researchers to draw meaningful conclusions.

In essence, data collection serves as the bedrock upon which research endeavors are built. Whether conducting experiments, surveys, observations, or interviews, researchers rely on collected data to inform their investigations and generate insights. By systematically gathering information, researchers can examine phenomena, test hypotheses, and uncover patterns or relationships within the data.

The purpose of data collection extends beyond mere accumulation of information; it aims to provide quality evidence that guides analysis and facilitates the construction of concrete answers to research questions. Quality evidence is characterized by its relevance, accuracy, reliability, and validity. Researchers strive to collect data that is pertinent to their research objectives, obtained through reliable methods, and accurately represents the phenomenon under study.

Data collection methodologies vary depending on the nature of the research and the objectives of the study. In experimental research, data may be collected through controlled experiments, where

variables are manipulated and outcomes are measured. Surveys and questionnaires are commonly used in social sciences and business research to gather information from individuals or groups. Observational studies involve systematically observing and recording behavior or phenomena in natural settings. In qualitative research, methods such as interviews and focus groups are utilized to explore subjective experiences and perspectives.

The UCI Machine Learning Repository serves as a valuable resource for researchers seeking to access and utilize datasets for analysis and experimentation. Hosted by the University of California, Irvine, the repository provides a diverse collection of datasets across various domains, ranging from healthcare and finance to environmental science and engineering. These datasets are curated and made available to researchers for academic and research purposes, enabling them to conduct experiments, develop models, and explore data-driven insights.

Utilizing datasets from repositories such as the UCI Machine Learning Repository offers several advantages for researchers. It provides access to a wide range of datasets encompassing diverse topics and domains, saving researchers the time and effort required to collect data independently. Additionally, datasets from reputable repositories are often well-documented and standardized, ensuring consistency and facilitating reproducibility of research findings.

In conclusion, data collection is a crucial aspect of research in all fields of study, serving as the foundation for analysis and interpretation. The process of collecting and measuring information enables researchers to address pertinent questions, evaluate outcomes, and construct meaningful answers. Leveraging resources such as the UCI Machine Learning Repository enhances the accessibility and quality of data available to researchers, fostering innovation and advancement across diverse domains of knowledge.

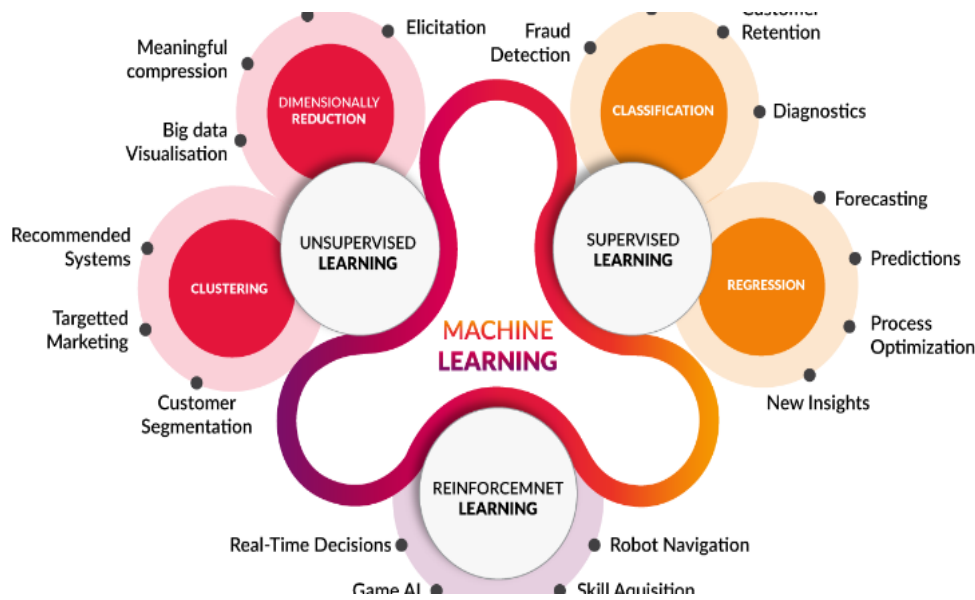


Fig 2.1: Machine Learning

2.1 Some machine learning methods

Machine learning algorithms are often categorized as supervised and unsupervised.

- **Supervised machine learning algorithm:** This algorithm utilize labeled data to predict outcomes for new, unseen instances. Through analysis of training data, algorithms discern patterns to produce a function mapping input features to output labels. Following training, the model can accurately predict outcomes for new data inputs by applying the inferred function. By comparing predictions with actual outcomes, the algorithm refines its parameters, enhancing accuracy. This iterative process enables continuous improvement, making supervised learning valuable for tasks like classification and regression. Its applicability spans various domains, including finance, healthcare, marketing, and natural language processing, where it aids in tasks such as stock price prediction, disease diagnosis, and sentiment analysis. Supervised learning's ability to generalize from known examples to predict outcomes for unseen instances underscores its significance in enabling computers to learn from data and make informed decisions, driving advancements across diverse fields.

Unsupervised machine learning algorithms operate without the need for labeled data, making them invaluable for uncovering hidden structures and patterns within datasets. Unlike supervised learning, where algorithms rely on labeled examples to predict outputs, unsupervised learning autonomously explores data to infer underlying relationships and groupings. By analyzing patterns, correlations, and similarities among data points, unsupervised algorithms can cluster similar data points together

or identify anomalies and outliers. This process enables the algorithm to gain insights into the inherent structure of the data, facilitating decision-making and knowledge discovery. Applications of unsupervised learning span various domains, including data mining, anomaly detection, and dimensionality reduction. In data mining, unsupervised algorithms reveal hidden patterns or associations, while in anomaly detection, they identify unexpected deviations from the norm. Additionally, unsupervised techniques aid in dimensionality reduction by extracting essential features while preserving critical information. Overall, unsupervised machine learning plays a pivotal role in exploring unlabeled data, enabling organizations to uncover insights, detect anomalies, and make informed decisions across diverse domains.

Reinforcement learning is a learning paradigm where an agent interacts with its environment, taking actions and receiving feedback in the form of rewards or penalties. Key characteristics of reinforcement learning include search and delayed reward. In this approach, the agent seeks to discover optimal behaviors through trial and error, aiming to maximize cumulative rewards over time.

Unlike supervised learning, where labeled examples guide the learning process, reinforcement learning relies on simple reward feedback to learn which actions lead to desirable outcomes. The reinforcement signal serves as a guide for the agent to adjust its behavior, reinforcing actions that lead to positive outcomes and discouraging those that lead to negative ones.

Reinforcement learning finds applications in various domains, including robotics, gaming, and autonomous systems. For example, in robotics, reinforcement learning algorithms enable robots to learn complex tasks such as grasping objects or navigating environments through interaction with their surroundings. In gaming, reinforcement learning powers intelligent agents capable of learning optimal strategies in games like chess or Go.

Overall, reinforcement learning empowers machines and software agents to autonomously determine optimal behaviors within specific contexts, leading to enhanced performance and adaptability in dynamic environments. By leveraging simple reward feedback, reinforcement learning enables agents to learn and improve their decision-making processes over time, making it a powerful tool in the realm of artificial intelligence.

2.2 APPLICATIONS OF MACHINE LEARNING

1. Virtual Personal Assistants
2. Predictions while Commuting
3. Videos Surveillance
4. Social Media Services
5. Email Spam and Malware Filtering
6. Online Customer Support
7. Search Engine Result Refining
8. Product Recommendations
9. Online Fraud Detection

2.3 IMPORTANT ALGORITMS IN MACHINE LEARNING

Linear Regression is a fundamental algorithm used in machine learning for predicting a continuous value based on one or more input features. It establishes a linear relationship between the input variables and the target variable, making it suitable for tasks like predicting house prices or stock prices. Logistic Regression, on the other hand, is specifically designed for binary classification problems. It estimates probabilities using a logistic or sigmoid function to predict the likelihood of a particular outcome, such as whether an email is spam or not. Decision Trees are versatile algorithms used for both classification and regression tasks. They partition the data into subsets based on features and predict the target variable in each subset, making them useful for tasks like customer segmentation or predicting loan default. Random Forest is an ensemble method that uses multiple decision trees to improve classification or regression performance. It combines the predictions of multiple trees to reduce overfitting and is commonly used in tasks like predicting customer churn or diagnosing diseases. K-Nearest Neighbors (KNN) is a simple yet intuitive algorithm for classification and regression. It predicts the label of a data point by majority voting of its k nearest neighbors, making it useful for tasks like image recognition or recommendation.

systems. Each of these algorithms has its strengths and weaknesses, and the choice of algorithm depends on the specific problem and the nature of the data.

2.4 TYPES OF FEATURE SELECTION (FS) TECHNIQUES

When building a machine learning model in real-life, it's almost rare that all the variables in the dataset are useful to build a model. Adding redundant variables reduces the generalization capability of the model and may also reduce the overall accuracy of a classifier. Furthermore adding more and more variables to a model increases the overall complexity of the model. As per the Law of Parsimony of 'Occam's Razor', the best explanation to a problem is that which involves the fewest possible assumptions. Thus, feature selection becomes an indispensable part of building machine learning models.

The goal of feature selection in machine learning is to find the best set of features that allows one to build useful models of studied phenomena.

The techniques for feature selection in machine learning can be broadly classified into the following categories:

Supervised Techniques: These techniques can be used for labeled data, and are used to identify the relevant features for increasing the efficiency of supervised models like classification and regression.

Unsupervised Techniques: These techniques can be used for unlabeled data.

From a taxonomic point of view, these techniques are classified as under:

- A. Filter methods
- B. Wrapper methods
- C. Embedded methods
- D. Hybrid methods

A. Filter methods

Filter methods pick up the intrinsic properties of the features measured via univariate statistics instead of cross-validation performance. These methods are faster and less computationally expensive than wrapper methods. When dealing with high-dimensional data, it is computationally cheaper to use filter methods.

Let's, discuss some of these techniques:

Information Gain

Information gain calculates the reduction in entropy from the transformation of a dataset. It can be used for feature selection by evaluating the Information gain of each variable in the context of the target

Chi-square Test

The Chi-square test is used for categorical features in a dataset. We calculate Chi-square between each feature and the target and select the desired number of features with the best Chi-square scores. In order to correctly apply the chi-squared in order to test the relation between various features in the dataset and the target variable, the following conditions have to be met: the variables have to be categorical, sampled independently and values should have an expected frequency greater than 5.

Fisher's Score

Fisher score is one of the most widely used supervised feature selection methods. The algorithm which we will use returns the ranks of the variables based on the fisher's score in descending order. We can then select the variables as per the case.

Correlation Coefficient

Correlation is a measure of the linear relationship of 2 or more variables. Through correlation, we can predict one variable from the other. The logic behind using correlation for feature selection is that the good variables are highly correlated with the target. Furthermore, variables should be correlated with the target but should be uncorrelated among themselves.

If two variables are correlated, we can predict one from the other. Therefore, if two features are correlated, the model only really needs one of them, as the second one does not add additional information

B. Wrapper Methods:

Wrappers require some method to search the space of all possible subsets of features, assessing their quality by learning and evaluating a classifier with that feature subset. The feature selection process is based on a specific machine learning algorithm that we are trying to fit on a given dataset. It follows a greedy search approach by evaluating all the possible combinations of features against the evaluation criterion. The wrapper methods usually result in better predictive accuracy than filter methods.

Let's, discuss some of these techniques:

Forward Feature Selection

This is an iterative method wherein we start with the best performing variable against the target. Next, we select another variable that gives the best performance in combination with the first selected variable. This process continues until the preset criterion is achieved.

Backward Feature Elimination

This method works exactly opposite to the Forward Feature Selection method. Here, we start with all the features available and build a model. Next, we remove the variable from the model which gives the best evaluation measure value. This process is continued until the preset criterion is achieved.

Exhaustive Feature Selection

This is the most robust feature selection method covered so far. This is a brute-force evaluation of each feature subset. This means that it tries every possible combination of the variables and returns the best performing subset.

Recursive Feature Elimination

‘Given an external estimator that assigns weights to features (e.g., the coefficients of a linear model), the goal of recursive feature elimination (RFE) is to select features by recursively considering smaller and smaller sets of features. First, the estimator is trained on the initial set of features and the importance of each feature is obtained either through a `coef_` attribute or through a `feature_importances_` attribute.

Then, the least important features are pruned from the current set of features. That procedure is recursively repeated on the pruned set until the desired number of features to select is eventually reached.

C. Embedded Methods:

These methods encompass the benefits of both the wrapper and filter methods, by including inter-actions of features but also maintaining reasonable computational cost. Embedded methods are iterative in the sense that takes care of each iteration of the model training process and carefully extracts those features which contribute the most to the training for a particular iteration.

Let’s, discuss some of these techniques [click here](#):

LASSO Regularization (L1)

Regularization consists of adding a penalty to the different parameters of the machine learning model to reduce the freedom of the model, i.e. to avoid over-fitting. In linear model regularization, the penalty is applied over the coefficients that multiply each of the predictors. From the different types of regularization, Lasso or L1 has the property that is able to shrink some of the coefficients to zero. Therefore, that feature can be removed from the model.

Random Forest Importance

Random Forests is a kind of a Bagging Algorithm that aggregates a specified number of decision trees. The tree-based strategies used by random forests naturally rank by how well they improve the purity of the node, or in other words a decrease in the impurity (Gini impurity) over all trees. Nodes with the greatest decrease in impurity happen at the start of the trees, while nodes with the least decrease in impurity occur at the end of trees. Thus, by pruning trees below a particular node, we can create a subset of the most important features.

2.5 MACHINE LEARNING ALGORITHMS IN THE FIELD OF FEATURE SELECTION (FS)

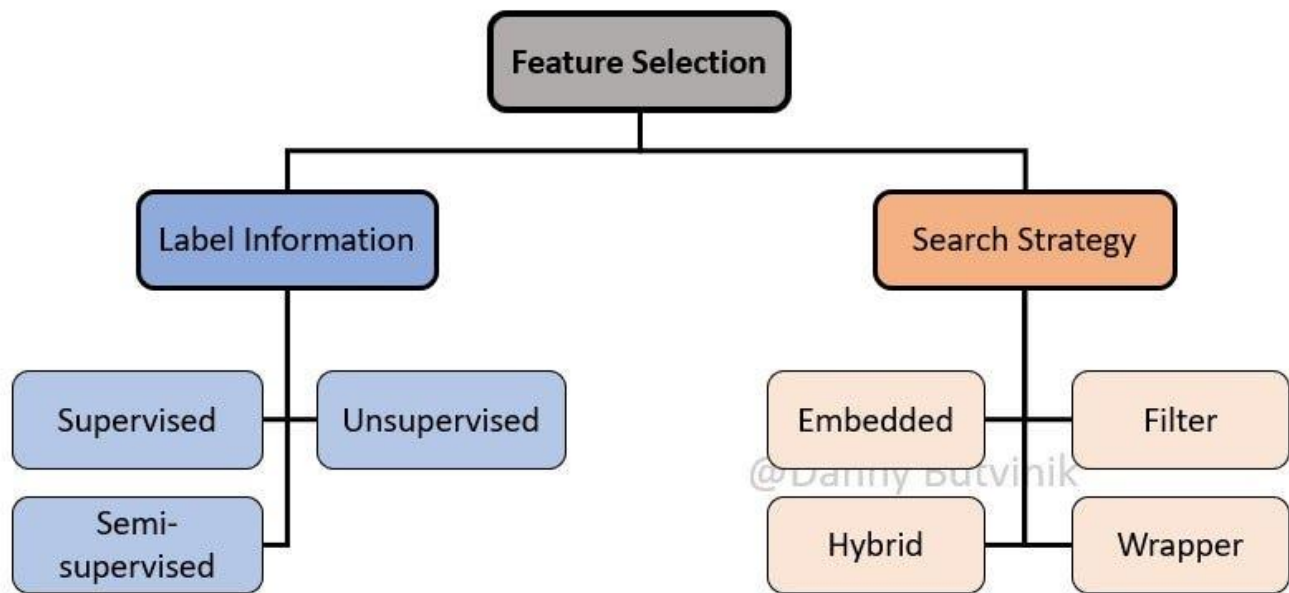


Fig 2.5.1 Classification of machine learning algorithms

Label Information:

In the context of machine learning, the term "label information" typically refers to the known outcomes or target variable in a dataset that the model aims to predict. This information is crucial for supervised learning algorithms, where the model learns from labeled training data to make predictions on new, unlabeled data.

1. **Supervised Learning:** In supervised learning, the dataset consists of input-output pairs, where the output (or label) is known for each input. The goal is to learn a mapping from inputs to outputs, so the model can predict the output for new, unseen inputs. Common algorithms include linear regression, logistic regression, decision trees, and neural networks.
2. **Unsupervised Learning:** In unsupervised learning, the dataset consists only of input data, without any corresponding output labels. The goal is to explore the structure of the data, find patterns, and group similar data points together. Clustering algorithms like K-means and hierarchical clustering are common in unsupervised learning.

3. **Semi-Supervised Learning:** Semi-supervised learning is a combination of supervised and unsupervised learning. It uses a small amount of labeled data along with a large amount of unlabeled data to improve the learning process. This can be useful when labeling data is expensive or time-consuming. Semi-supervised learning algorithms include self-training, co-training, and multi-view learning.

In feature selection, the label information is particularly important in supervised learning, as it guides the selection of features that are most relevant for predicting the target variable. Unsupervised learning techniques, on the other hand, do not rely on label information and instead focus on finding patterns and relationships in the data based solely on the input features. Semi-supervised learning leverages both labeled and unlabeled data to improve the performance of the model.\

Search Strategy:

When deciding on a feature selection strategy for a machine learning project, it's crucial to consider the nature of the dataset, the computational resources available, and the specific goals of the analysis.

1. **Filter methods** are efficient for large datasets as they evaluate features based on statistical properties like correlation or information gain, making them suitable for preprocessing before more complex algorithms.
2. **Wrapper methods**, on the other hand, are computationally intensive, evaluating subsets of features using a machine learning model, making them ideal for optimizing model performance but less practical for large datasets due to their computational cost.
3. **Embedded methods** offer a balance, performing feature selection as part of model training, making them efficient and suitable for high-dimensional data. They're particularly useful for interpreting feature importance.
4. **Hybrid methods** combine different techniques to leverage their strengths, offering a compromise between accuracy and computational cost.

Ultimately, the choice of feature selection strategy depends on the specific characteristics of the dataset and the goals of the analysis. Experimentation and comparison of different strategies are often necessary to determine the most effective approach for a given task

3. EXISTING SYSTEM

The system offers comprehensive support for farmers through its crop recommendation, crop yield prediction, and crop fertilizer prediction features. By utilizing data on soil type, climate conditions, and historical crop performance, the system suggests the most suitable crops for a given piece of land, optimizing yield and profit potential. Moreover, by analyzing historical data, weather forecasts, and soil conditions, the system predicts crop yields, aiding farmers in planning their harvest and resource allocation. Additionally, by evaluating soil health, crop requirements, and nutrient levels, the system recommends the appropriate fertilizers for optimal crop growth, ensuring that crops receive the necessary nutrients for improved yields and quality. These features collectively empower farmers with, enabling them to make informed decisions and enhance their agricultural practices.

3.1 FEATURE SELECTION TECHNIQUES

There are mainly two types of Feature Selection techniques, which are:

- Supervised Feature Selection technique**

Supervised Feature selection techniques consider the target variable and can be used for the labelled dataset.

- Unsupervised Feature Selection technique**

Unsupervised Feature selection techniques ignore the target variable and can be used for the unlabelled dataset.

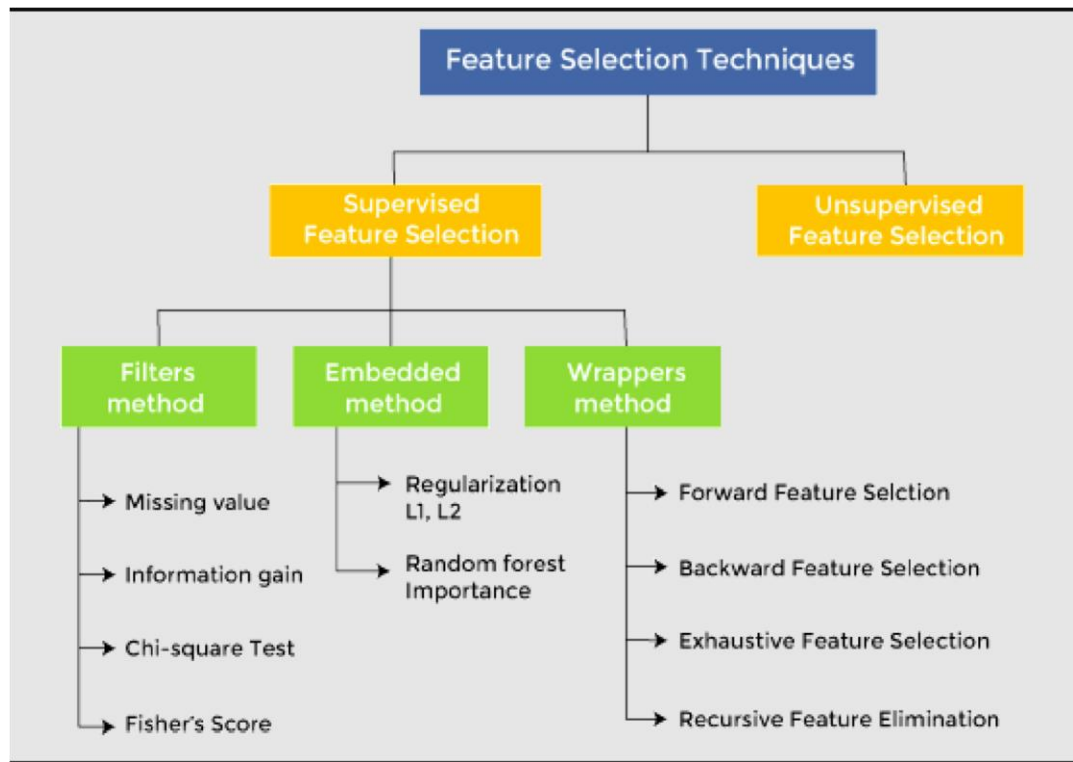


Fig 3.1.1 Feature Selection methods

3.2 WRAPPER METHODS

Forward selection-

Forward selection is an iterative process, which begins with an empty set of features. After each iteration, it keeps adding on a feature and evaluates the performance to check whether it is improving the performance or not. The process continues until the addition of a new variable/feature does not improve the performance of the model.

Backward elimination-

Backward elimination is also an iterative approach, but it is the opposite of forward selection. This technique begins the process by considering all the features and removes the least

significant feature. This elimination process continues until removing the features does not improve the performance of the model.

Exhaustive Feature Selection-

Exhaustive feature selection is one of the best feature selection methods, which evaluates each feature set as brute-force. It means this method tries & make each possible combination of features and return the best performing feature set.

Recursive Feature Elimination-

Recursive feature elimination is a recursive greedy optimization approach, where features are selected by recursively taking a smaller and smaller subset of features. Now, an estimator is trained with each set of features, and the importance of each feature is determined using `coef_attribute` or through a `feature_importances_attribute`.

3.3 FILTER METHODS

In Filter Method, features are selected on the basis of statistics measures. This method does not depend on the learning algorithm and chooses the features as a pre-processing step. The filter method filters out the irrelevant feature and redundant columns from the model by using different metrics through ranking. The advantage of using filter methods is that it needs low computational time and does not overfit the data.

Information Gain:

Information gain determines the reduction in entropy while transforming the dataset. It can be used as a feature selection technique by calculating the information gain of each variable with respect to the target variable.

Chi-square Test:

Chi-square test is a technique to determine the relationship between the categorical variables. The chi-square value is calculated between each feature and the target variable, and the desired number of features with the best chi-square value is selected.

Fisher's Score:

Fisher's score is one of the popular supervised technique of features selection. It returns the rank of the variable on the fisher's criteria in descending order. Then we can select the variables with a large fisher's score.

Missing Value Ratio:

The value of the missing value ratio can be used for evaluating the feature set against the threshold value. The formula for obtaining the missing value ratio is the number of missing values in each column divided by the total number of observations. The variable is having more than the threshold value can be dropped.

3.4 EMBEDDED METHODS

Embedded methods combined the advantages of both filter and wrapper methods by considering the interaction of features along with low computational cost. These are fast processing methods similar to the filter method but more accurate than the filter method.

These methods are also iterative, which evaluates each iteration, and optimally finds the most important features that contribute the most to training in a particular iteration. Some techniques of embedded methods are:

Regularization-Regularization adds a penalty term to different parameters of the machine learning model for avoiding overfitting in the model. This penalty term is added to the coefficients; hence it shrinks some coefficients to zero. Those features with zero coefficients can be removed from the dataset. The types of regularization techniques are L1 Regularization (Lasso Regularization) or Elastic Nets (L1 and L2 regularization).

Random Forest Importance- Different tree-based methods of feature selection help us with feature importance to provide a way of selecting features. Here, feature importance specifies which feature has more importance in model building or has a great impact on the target variable.

Random Forest is such a tree-based method, which is a type of bagging algorithm that aggregates a different number of decision trees. It automatically ranks the nodes by their performance or decrease in the impurity (Gini impurity) over all the trees. Nodes are arranged as per the impurity values, and thus it allows to pruning of trees below a specific node.

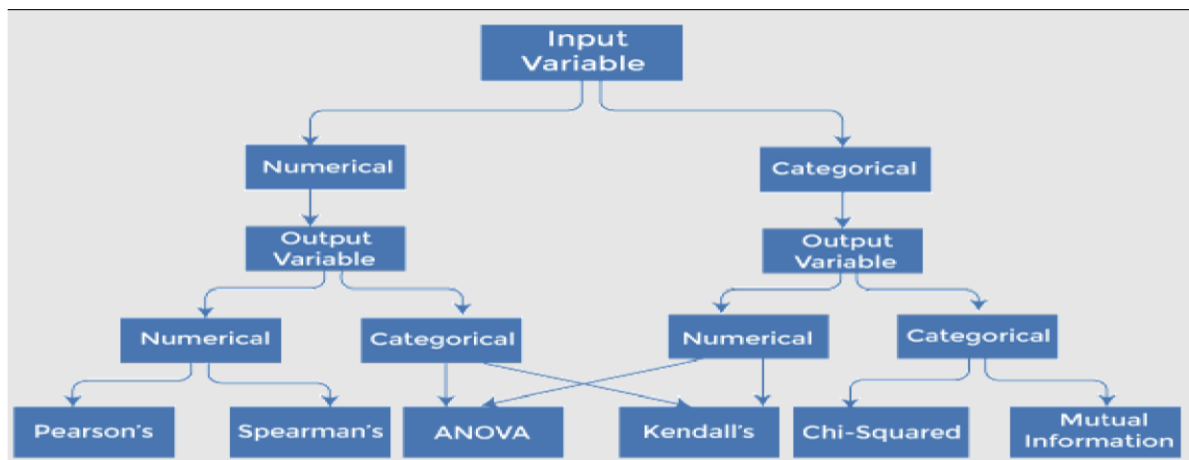


Fig 3.4.1 Tree Structure of Feature Selection

3.5 ALGORITHMS

3.5.1 Random Forest Algorithm

Machine learning, a fascinating blend of computer science and statistics, has witnessed incredible progress, with one standout algorithm being the Random Forest. Random forests or Random Decision Trees is a collaborative team of decision trees that work together to provide a single output. Originating in 2001 through Leo Breiman, Random Forest has become a cornerstone

for machine learning enthusiasts. In this article, we will explore the fundamentals and implementation RandomForestAlgorithm.

Random Forest algorithm is a powerful tree learning technique in Machine Learning. It works by creating a number of Decision Trees during the training phase. Each tree is constructed using a random subset of the data set to measure a random subset of features in each partition. This randomness introduces variability among individual trees, reducing the risk of overfitting and improving overall prediction performance. In prediction, the algorithm aggregates the results of all trees, either by voting (for classification tasks) or by averaging (for regression tasks) This collaborative decision-making process, supported by multiple trees with their insights, provides an example stable and precise results. Random forests are widely used for classification and regression functions, which are known for their ability to handle complex data, reduce overfitting, and provide reliable forecasts in different environments.

Random Forest

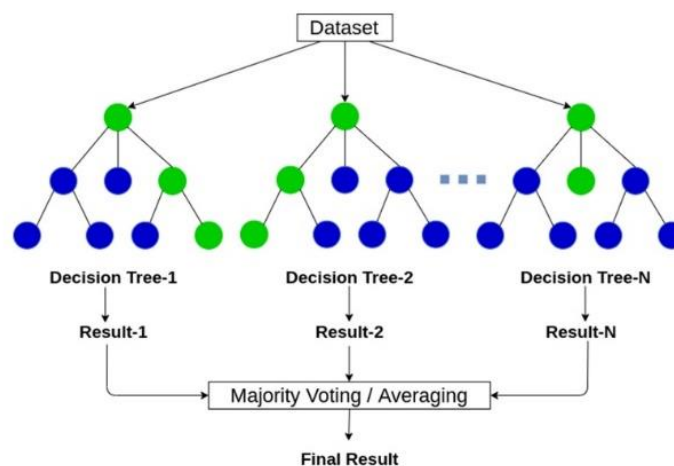


Fig 3.5.1.1 Random Forest tree Algorithm

3.5.2 K-Nearest Neighbor(KNN) Algorithm

The **K-Nearest Neighbors (KNN) algorithm** is a supervised machine learning method employed to tackle classification and regression problems. Evelyn Fix and Joseph Hodges developed this algorithm in 1951, which was subsequently expanded by Thomas Cover. The article explores the fundamentals, workings, and implementation of the KNN algorithm.

What is the K-Nearest Neighbors Algorithm?

KNN is one of the most basic yet essential classification algorithms in machine learning. It belongs to the supervised learning domain and finds intense application in pattern recognition, data mining, and intrusion detection.

It is widely disposable in real-life scenarios since it is non-parametric, meaning it does not make any underlying assumptions about the distribution of data (as opposed to other algorithms such as GMM, which assume a Gaussian distribution of the given data). We are given some prior data (also called training data), which classifies coordinates into groups identified by an attribute.

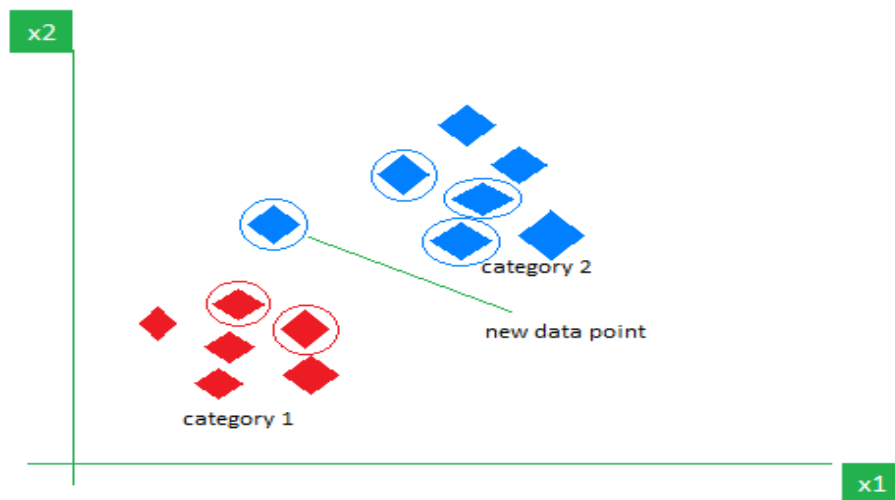


Fig 3.5.2.1 KNN Algorithm working visualization

3.5.3 Naive Bayes Classifiers

A Naive Bayes classifiers, a family of algorithms based on Bayes' Theorem. Despite the “naive” assumption of feature independence, these classifiers are widely utilized for their simplicity and efficiency in machine learning. The article delves into theory, implementation, and applications, shedding light on their practical utility despite oversimplified assumptions.

What is Naive Bayes Classifiers?

Naive Bayes classifiers are a collection of classification algorithms based on Bayes' Theorem. It is not a single algorithm but a family of algorithms where all of them share a common principle, i.e. every pair of features being classified is independent of each other. To start with, let us consider a dataset.

One of the most simple and effective classification algorithms, the Naïve Bayes classifier aids in the rapid development of machine learning models with rapid prediction capabilities.

Naïve Bayes algorithm is used for classification problems. It is highly used in text classification. In text classification tasks, data contains high dimension (as each word represent one feature in the data). It is used in spam filtering, sentiment detection, rating classification etc. The advantage of using naïve Bayes is its speed. It is fast and making prediction is easy with high dimension of data.

3.5.4 Linear Regression in Machine learning

Machine Learning is a branch of Artificial intelligence that focuses on the development of algorithms and statistical models that can learn from and make predictions on data. **Linear regression** is also a type of machine-learning algorithm more specifically a **supervised machine-learning algorithm** that learns from the labelled datasets and maps the data points to the most optimized linear functions. which can be used for prediction on new datasets.

First of we should know what supervised machine learning algorithms is. It is a type of machine learning where the algorithm learns from labelled data. Labeled data means the dataset whose respective target value is already known. Supervised learning has two types:

- **Classification:** It predicts the class of the dataset based on the independent input variable. Class is the categorical or discrete values. like the image of an animal is a cat or dog?
- **Regression:** It predicts the continuous output variables based on the independent input variable. like the prediction of house prices based on different parameters like house age, distance from the main road, location, area, etc.
-

4. PROPOSED SYSTEM

Our proposed system focuses on developing a robust crop price prediction model through the integration of machine learning algorithms and historical data. By analyzing factors such as rainfall patterns, wholesale cost index fluctuations, and temporal variables like month and year, our framework aims to generate accurate forecasts of crop prices. These predictions will empower farmers by providing valuable insights for crop selection, optimal harvest timing, and strategic pricing decisions. By increasing farmers' profitability and facilitating informed decision-making, our system seeks to enhance the stability and sustainability of the agricultural sector. Through the utilization of advanced technology and data-driven approaches, we aim to address the challenges faced by farmers and contribute to the overall growth and resilience of agricultural communities.

5. SYSTEM REQUIREMENTS

5.1 Hardware Requirements:

- SystemType : Intel Core i3 or above
- Cachememory : 4MB(Megabyte)
- RAM : 8 gigabyte (GB)
- BusSpeed : 5 GT/s DBI2
- Number of cores : 2
- Number of threads : 4

5.2 Software Requirements:

- Operating System : Windows 10 Home, 64 bit Operating System
- Coding Language : Python
- Python distribution: Anaconda

6. SYSTEM ANALYSIS

6.1 SCOPE OF PROJECT

The scope of the project encompasses the development of a comprehensive agricultural system integrating Decision Tree Regression for crop price prediction and various modules including weather forecasts, crop recommendations, and fertilizer suggestions. This system aims to empower farmers by providing advance price forecasts, enhancing decision-making, maximizing crop yield, and optimizing resource allocation. Additionally, it includes features such as a shop, chat portal, and guide functionalities to facilitate user interaction and support. By leveraging advanced technologies like Machine Learning, the project seeks to transform farming practices, increase profitability, and contribute to the country's economic growth and agricultural sustainability.

6.2 ANALYSIS

The provided system, leveraging Machine Learning's Decision Tree Regression Algorithm, addresses critical challenges faced by farmers in the agricultural sector. By accurately predicting crop prices based on key attributes such as rainfall, wholesale price index, month, and year, the system enables farmers to make informed decisions, mitigating financial risks and increasing profitability. This aspect of price prediction is crucial, as it empowers farmers to anticipate market trends and plan their crop sales accordingly, reducing the impact of fluctuating prices and minimizing losses due to spoilage.

Moreover, the system goes beyond price prediction by incorporating additional modules for weather forecasting, crop recommendation, and fertilizer recommendation. These modules provide farmers with comprehensive insights and recommendations tailored to their specific needs and circumstances. For instance, the crop recommendation module assists farmers in selecting suitable crop types based on their land's characteristics, optimizing yield potential and resource utilization.

Similarly, the fertilizer recommendation module suggests appropriate fertilizers based on soil health and crop requirements, further enhancing crop growth and quality.

Additionally, the inclusion of a shop, chat portal, and guide enhances the system's usability and accessibility, allowing farmers to access essential information and support services conveniently. By integrating advanced technologies like Machine Learning, the system not only addresses immediate challenges but also has the potential to revolutionize the agricultural sector by fostering innovation, efficiency, and sustainability. Overall, this comprehensive system serves as a valuable tool for farmers, empowering them to improve their farming practices, increase productivity, and contribute to the country's economic growth.

6.3 DATA PREPROCESSING & FEATURE SCALING

Agriculture, as the backbone of many economies, faces significant challenges in recent years, primarily due to unpredictable climate patterns and fluctuating market trends. These uncertainties pose substantial risks to farmers, often resulting in crop spoilage and financial losses. Moreover, farmers may struggle to identify the most suitable crops for their land, exacerbating their predicament. To address these issues, a comprehensive agricultural decision support system has been developed, leveraging advanced machine learning (ML) techniques. This system aims to empower farmers with accurate predictions, personalized recommendations, and valuable insights to enhance their decision-making processes, ultimately increasing profitability and contributing to economic growth.

At the heart of the system lies Decision Tree Regression, a powerful ML algorithm capable of analyzing diverse datasets to predict crop prices accurately. By considering factors such as rainfall patterns, wholesale price indices, and temporal variations, Decision Tree Regression offers a flexible and robust framework for price forecasting. Its ability to handle both numerical and categorical data makes it particularly well-suited for the agricultural context, where variables can exhibit varying levels of complexity and interdependence.

In addition to price prediction, the system incorporates ML algorithms into several key modules to address different aspects of agricultural decision-making. The crop recommendation module utilizes collaborative filtering or content-based filtering techniques to match land characteristics with suitable crop types. Drawing insights from historical data and expert knowledge, this module assists farmers in

selecting crops that are well-suited to their specific agricultural conditions, thereby maximizing yield potential.

Similarly, the crop yield prediction module employs regression models such as Random Forest Regression or Long Short-Term Memory networks to forecast crop yields accurately. By integrating various factors such as weather conditions, soil health, and agricultural practices, this module provides farmers with valuable insights into expected yields, enabling them to plan their harvest and allocate resources more effectively.

The fertilizer recommendation module complements these efforts by suggesting appropriate fertilizers based on soil nutrient levels and crop requirements. By leveraging rule-based systems or collaborative filtering techniques, this module ensures that farmers apply fertilizers optimally, thereby promoting crop growth and quality while minimizing environmental impact.

Furthermore, the system seamlessly integrates advanced technologies such as data preprocessing, feature engineering, and model training to ensure the accuracy and reliability of its predictions and recommendations. Through a user-friendly interface encompassing weather forecasts, crop and fertilizer suggestions, a shop, chat portal, and guide features, the system provides farmers with easy access to essential information and tools to support their decision-making processes.

Data integration from multiple sources, including weather stations, agricultural databases, and market prices, further enhances the system's effectiveness. By aggregating and analyzing data from these diverse sources, the system can provide farmers with comprehensive insights into market trends, weather patterns, and soil health, enabling them to make informed decisions.

Scalability considerations are also paramount, given the potential for the system to serve a large and diverse user base. By leveraging distributed computing and cloud infrastructure, the system can handle large volumes of data and accommodate a growing number of users without sacrificing performance or reliability.

In conclusion, the agricultural decision support system represents a significant step forward in leveraging ML and advanced technologies to address the challenges faced by farmers. By providing

accurate predictions, personalized recommendations, and valuable insights, the system empowers farmers to make informed decisions that enhance productivity, profitability, and sustainability. As agriculture continues to evolve in the face of changing climate patterns and market dynamics, systems like these will play an increasingly vital role in driving economic growth and ensuring food security for future generations

7. DESIGN ANALYSIS

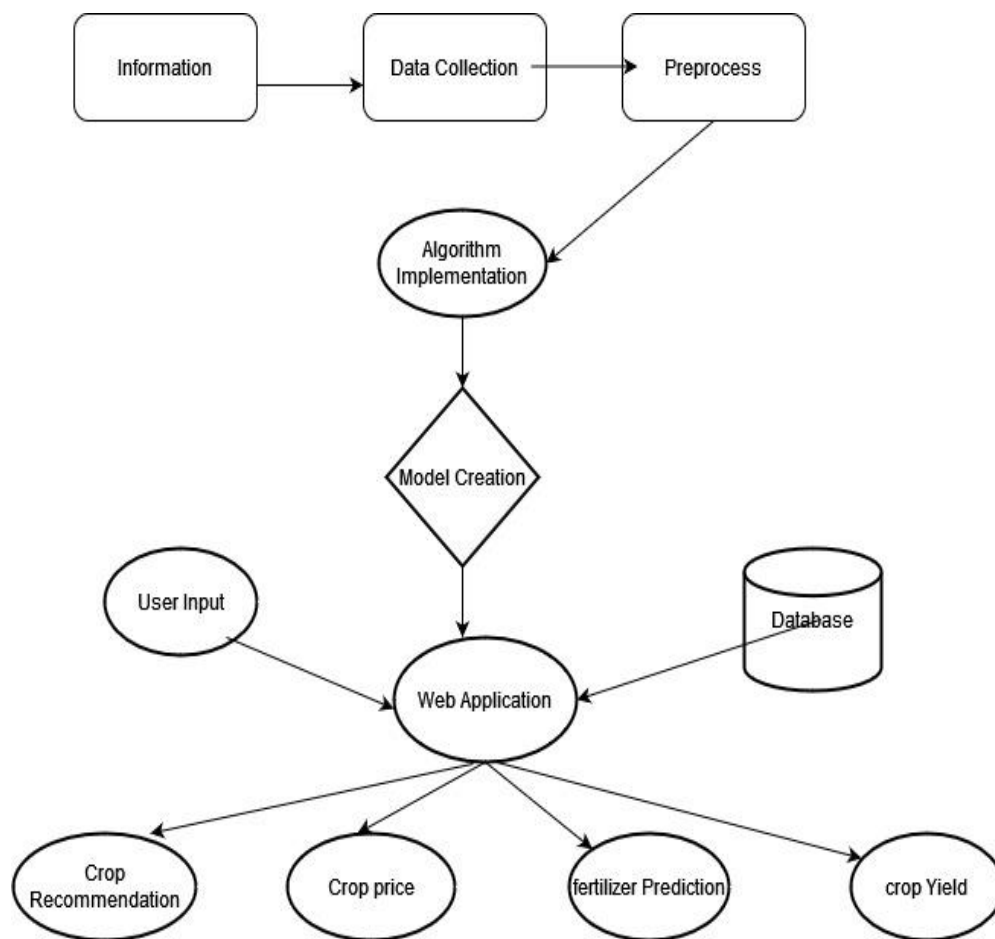


Fig 7.1 Design of Binary Dragonfly Algorithm

A. Research Objective: By integrating these elements into your research, you can offer a more holistic analysis of the factors influencing crop production and pricing, helping farmers make informed decisions. The crop recommendation model will assist farmers in selecting optimal crops, considering soil quality and land suitability, climate, and historical data. This can lead to improved crop yields and reduced risk of crop failure.[4]

The crop fertilizer prediction model will assist farmers decide on the type and amount of fertilizer needed for optimal crop growth. By considering soil nutrient levels, crop type, and other relevant factors, this model can ensure that farmers apply the right amount of fertilizer, reducing waste and environmental impact. The crop yield prediction model will enable farmers to anticipate their crop yields based on elements like weather patterns, soil health, and crop type.

This information can help farmers plan their harvest and allocate resources more effectively, leading to increased efficiency and profitability. B. Data Description: Getting the data is a crucial step in any machine learning endeavor, as the quality of the data affects how well the model works. In this study, we obtained our data from Kaggle, a popular website where data scientists share datasets. After collecting the data, which comprised over 7418 records and 9 different pieces of information[8], we uploaded it to Google Colab, an online platform for analyzing data and performing machine learning tasks. We utilized four datasets from Kaggle, each containing a different number of records

- Fertilizer Dataset: This dataset consists of 100 records and 9 columns, providing information on the types and quantities of fertilizers suitable for different crops.[12]
- Crop Price Dataset: This dataset consists of 7418 records and 9 columns, providing information on the types and quantities of Crop Price suitable for different crops.[9]
- Yield Prediction Dataset: With 120 records and 6 columns, this dataset contains information Utilized for forecasting crop yield considering weather, soil, and crop type factors.[8].
- Crop Recommendation Dataset: This dataset comprises 2201 records and 7 columns, offering

recommendations for suitable crops based on factors like soil quality, climate, and historical data [9].

C. Pre-processing of Dataset

Before we could use the data, we had to clean it up by getting rid of any missing information and unusual values. This makes the data ready for training and testing our models.

D. Data Visualization

We used graphs and charts to look at the data and see if there were any patterns. In below figure (1) we divided the dataset by states on one side and rainfall in different states. it was use analyzing average in every state. [21]

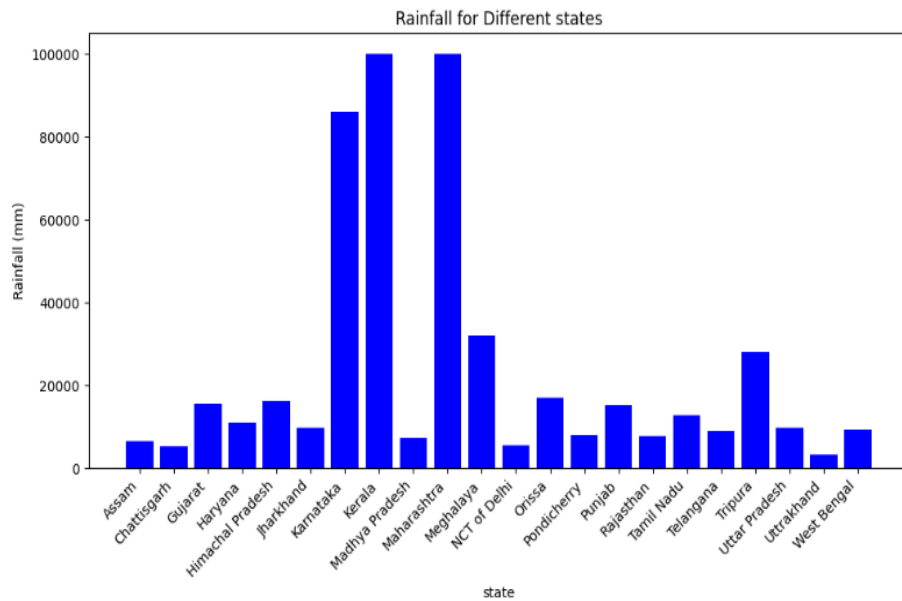


Fig (1): Rain for different states

We used graphs and charts to look at the data and see if there were any patterns. In below figure (2) we divided the dataset by varieties on one side and rainfall in different states. It was use analyzing average in every districts [12].

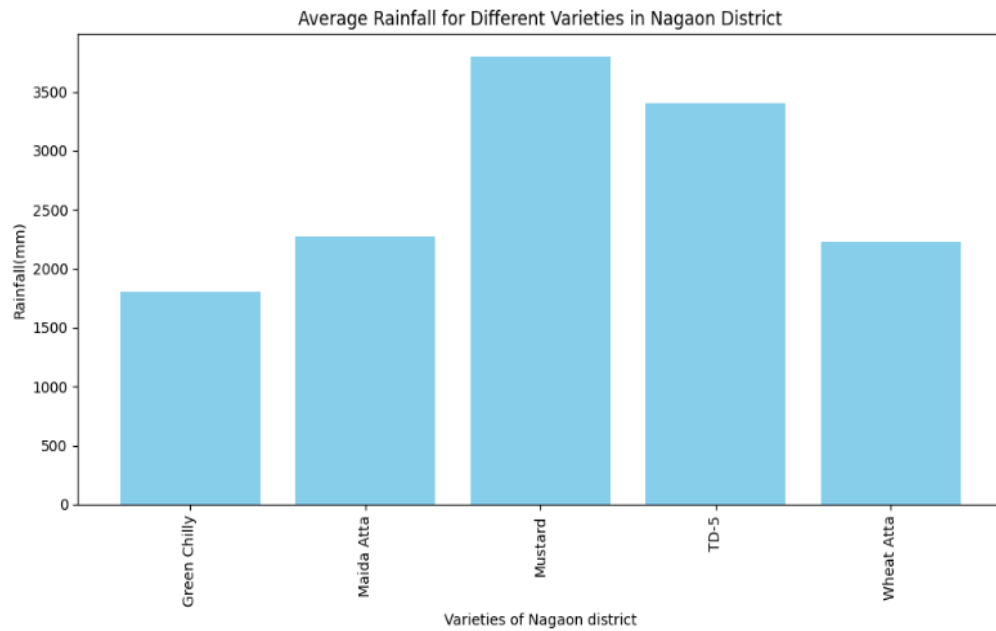


Fig (2): Average rain fall for different varieties in particular district

We used graphs and charts to look at the data and see if there were any patterns. In below figure (3) we divided the dataset by varieties on one side and Croptype in different crops . It was use analyzing average phosphorous on soi [12]l.

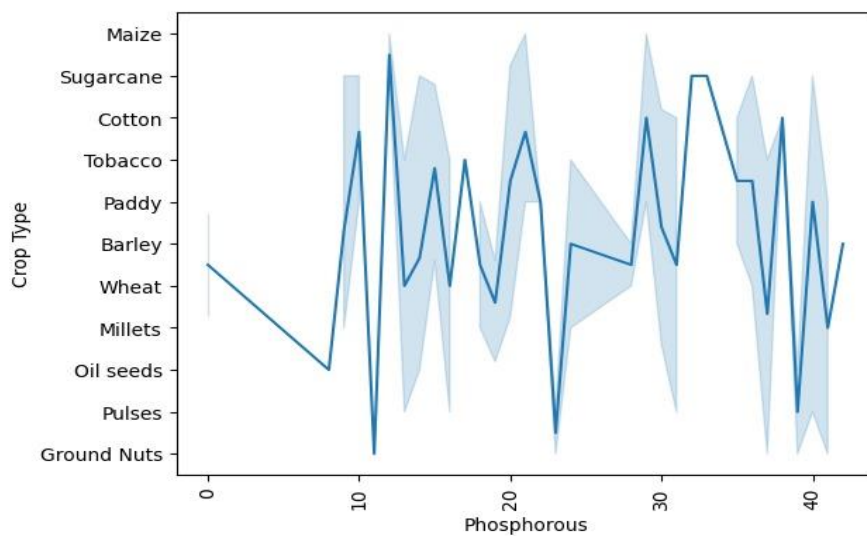


Fig (3): Phosphorous levels in different crops

We used graphs and charts to look at the data and see if there were any patterns. In below figure (4) we divided the dataset by varieties on one side and Crop type in different crops . It was use analyzing average Nitrogen on soil [8].

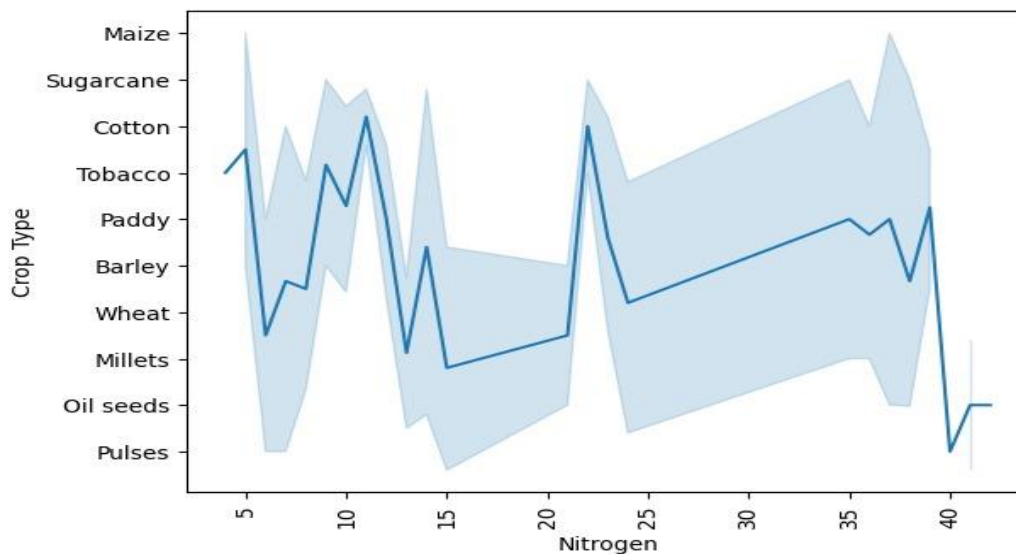


Fig (4): Nitrogen levels in different crops

We used graphs and charts to look at the data and see if there were any patterns. In below figure (5) we divided the dataset by varieties on one side and Crop Yield in different crops. It was use analyzing average Cost on every Crop [8].

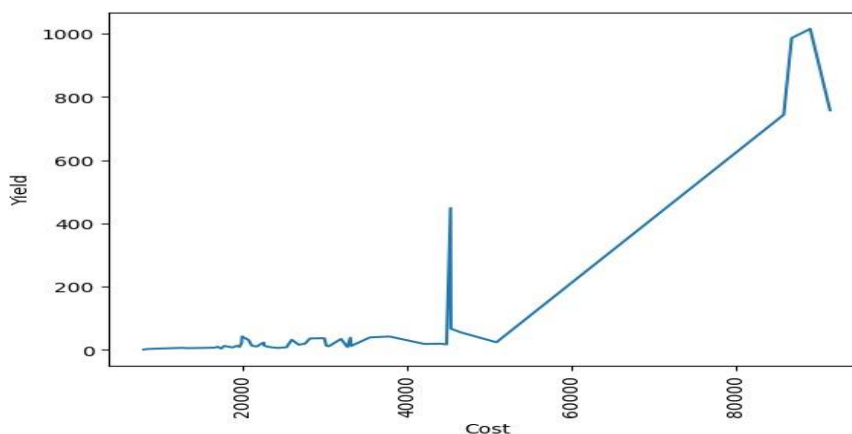


Fig (5): Analyzing yield and cost

We used graphs and charts to look at the data and see if there were any patterns. In below figure (6) we divided the dataset by varieties on one side and Crop Yield in different crops. It was use analyzing Different States on every Crop [9].

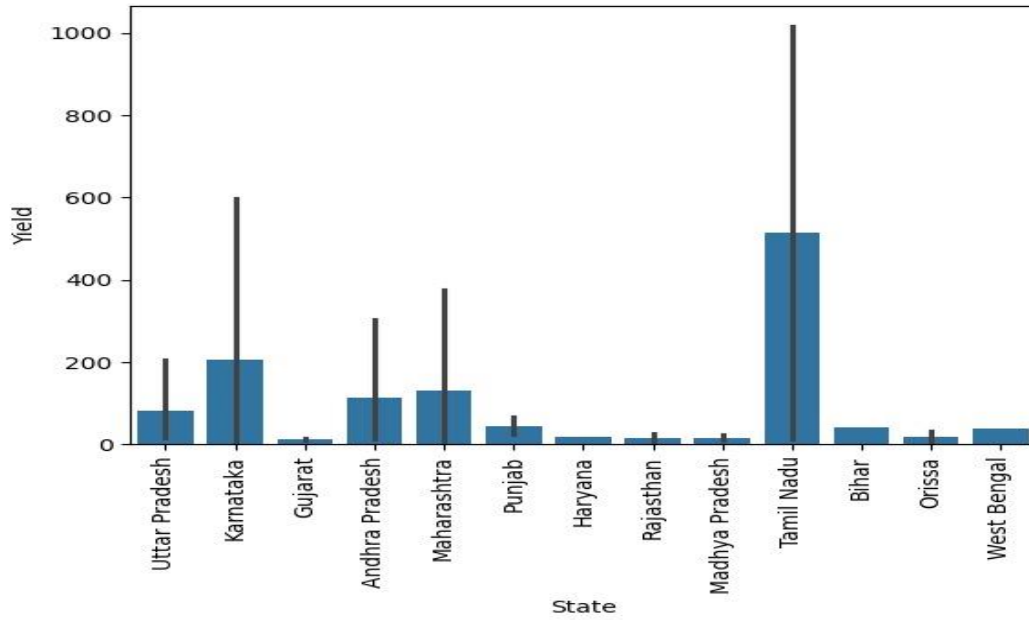


Fig (6): yield production in different states

E. Separating and Splitting the Data

Separating the input features and target data. Then, we split the data into two parts: one part, which was 80% of the data, was used for training our models, and the other part, which was 20% of the data, was used to see how well the models worked.

F. Model Assessment and Selection

By using K-Fold Validation we tested four various machine learning algorithms —Ridge, Lasso, Decision Tree, Random forest, linear regression, and KNeighbors Regressor regression—to see how well they could predict maximum crop price. After conducting testing, the models will undergo evaluation based on their scores obtained through performing 10 iterations of K-fold validation. The model exhibiting the highest mean score in predicting the target variable is chosen as the optimal model. Additionally, prediction Performance metrics including Mean Square Error (MSE), R-Squared (R²), and Mean Absolute Error (MAE) are utilized for additional evaluation of model performance. The model exhibiting the lowest prediction errors will be deemed the most suitable choice.

H. Findings and Discussion

In the results section, we will analyze the accuracy of models trained on both past and current data, examine various types of errors encountered, present bar charts illustrating the accuracy of different algorithms, and provide visual representations of evaluation metrics for each algorithm. Additionally, we will showcase the predicted results from our web Interface and compare our findings to recent studies in the field.

7.1 Model Accuracies of Different Models

Our assessment involved rigorous validation of our models utilizing K-fold cross-validation with 10 iterations, a widely recognized technique in machine learning evaluations. This method ensures robust validation, fostering dependable performance assessment. The accuracy and reliability of our models are pivotal, offering users valuable insights to inform decisions in various agricultural domains.

In interfaces such as crop price prediction, crop recommendation, fertilizer prediction, and crop yield prediction, users benefit immensely from accurate predictions. These predictions empower users to make informed decisions regarding crop cultivation, pricing strategies, and market trends. Presenting this information in an intuitive and user-friendly interface enhances usability, enabling individuals to leverage predictions according to their specific needs effectively.

Throughout our evaluation, all models consistently demonstrated high accuracies, affirming their effectiveness. Notably, the Decision Tree model exhibited an impressive average accuracy of 95.76%, while the Random Forest model surpassed expectations with an average accuracy of 98%. These exceptional results underscore the reliability and efficacy of these models in predicting maximum crop prices.

The remarkable accuracies observed render these models highly suitable for integration into web applications, ensuring dependable and effective predictions. The bar graph below illustrates the average accuracies of the regression algorithms, providing a visual depiction of their performance.

In conclusion, our thorough assessment utilizing K-fold cross-validation with 10 iterations reaffirms the robustness and reliability of our models. Their high accuracies make them invaluable tools for agricultural decision-making, offering users actionable insights to optimize crop cultivation and pricing strategies. By presenting predictions through user-friendly interfaces, we aim to empower individuals in the agricultural sector to make informed choices, ultimately enhancing productivity and profitability.

Table -1 Accuracy of different algorithms

Algorithms	Accuracy
Random Forest	0.98
Decision Tree	0.95
Lasso	0.57
Ridge	0.57
Linear Regression	0.57
KNN	0.75

The performance of various algorithms was evaluated based on their accuracy scores for a specific task. Random Forest proved to be the most effective, attaining an accuracy of 0.98. Following closely was the Decision Tree algorithm with an accuracy of 0.95, indicating its suitability for the task. In contrast, the linear models including Lasso, Ridge, and Linear Regression all displayed similar accuracy scores of 0.57, which, although lower than the tree-based models, can still be valuable in certain contexts. K-Nearest Neighbours (KNN) achieved an accuracy of 0.75, respectable but falling behind the ensemble methods. These results emphasize the effectiveness of ensemble methods for the task, particularly Random Forest and Decision Tree, while also highlighting the importance of considering other factors such as interpretability, computational efficiency, and scalability when choosing an algorithm for a specific application. Random Forest, with its impressive accuracy of 0.98, stands out as the top performer in his evaluation.

8. IMPLEMENTATION

```
import numpy as np

import matplotlib.pyplot as plt

import pandas as pd

data= pd.read_csv('CropPrice.csv')

print(data)

data.dtypes

data.isnull()

import pandas as pd

data = pd.read_csv('CropPrice.csv')

data.drop(8, inplace=True)

data.head(10)

plt.figure(figsize=(10, 6))

plt.bar(data['state'], data['Rainfall'], color='blue')

plt.xlabel('state')

plt.ylabel('Rainfall (mm)')

plt.title('Rainfall for Different states')

plt.xticks(rotation=45, ha='right')

plt.tight_layout()

plt.show()

Gujarat_data = data[data['state'] == 'Gujarat']

districts = Gujarat_data['district'].tolist()

rainfall = Gujarat_data['Rainfall'].tolist()

plt.figure(figsize=(10, 6))
```

```
plt.bar(districts, rainfall, color='skyblue')

plt.xlabel('Districts')

plt.ylabel('Rainfall (in mm)')

plt.title('Rainfall in Different Districts of Gujarat')

plt.xticks(rotation=45, ha='right')

plt.tight_layout()
```

```
plt.show()

Assam_data = data[data['state'] == 'Assam']
```

```
districts = Assam_data['district'].tolist()

variety = Assam_data['variety'].tolist()
```

```
plt.figure(figsize=(10, 10))

plt.bar(districts, variety, color='skyblue')

plt.xlabel('Districts')

plt.ylabel('Variety')

plt.title('Varieties in different districts of Assam state')

plt.xticks(rotation=45, ha='right')

plt.tight_layout()
```

```
plt.show()

import numpy as np

import pandas as pd

import matplotlib.pyplot as plt

from sklearn.preprocessing import PolynomialFeatures

from sklearn.linear_model import LinearRegression, Lasso, Ridge
```



```

from sklearn.neighbors import KNeighborsRegressor

from sklearn.tree import DecisionTreeRegressor

from sklearn.ensemble import RandomForestRegressor, GradientBoostingRegressor


from sklearn.metrics import mean_squared_error, mean_absolute_error, r2_score

# Load the data

df = pd.read_csv('CropPrice.csv')

from sklearn.preprocessing import LabelEncoder

from sklearn.model_selection import train_test_split

# Handling missing values

crop_data = crop_data.dropna()

# Converting categorical data to numerical form using LabelEncoder

le = LabelEncoder()

categorical_columns = ['state', 'district', 'market', 'commodity', 'variety']

for col in categorical_columns:

    crop_data[col] = le.fit_transform(crop_data[col])

# Convert 'arrival_date' to a numerical value (e.g., timestamp)

crop_data['arrival_date'] = pd.to_datetime(crop_data['arrival_date']).values.astype(np.int64) // 10**9

# Splitting the dataset into features (X) and the target variable (y)

X = crop_data.drop(['max_price'], axis=1)

y = crop_data['max_price']

# Splitting the dataset into training and testing sets

X_train, X_test, y_train, y_test = train_test_split(X, y, test_size=0.2, random_state=42)

print(X_train.iloc[0])

# Define the regression models

models = {

    'Linear Regression': LinearRegression(),

```

```

'Polynomial Regression' : LinearRegression(),

'Lasso': Lasso(),

'Ridge': Ridge(),

'K-Nearest Neighbors': KNeighborsRegressor(n_neighbors=5),

'Decision Tree': DecisionTreeRegressor(),

'Random Forest': RandomForestRegressor(),

}

# Evaluate each model and calculate feature importance

results = []

for model_name, model in models.items():

    model.fit(X_train, y_train)  # Fit the model on the training data

    # Make predictions on the testing data

    y_pred = model.predict(X_test)

    # Calculate evaluation metrics

    mse = mean_squared_error(y_test, y_pred)

    mae = mean_absolute_error(y_test, y_pred)

    r2 = r2_score(y_test, y_pred)

    # Append results to a DataFrame

    results.append({'model_name': model_name, 'mse': mse, 'mae': mae, 'r2': r2})

# Create a Pandas DataFrame

df = pd.DataFrame(results)

# Print the DataFrame

print(df.to_string())

import matplotlib.pyplot as plt

import pandas as pd

```

```

# Assuming you have a DataFrame named 'df_results' with the evaluation results

# Replace 'df_results' with the name of your DataFrame

# df_results should have columns: 'model_name', 'mse', 'mae', 'r2'

df_results=df

# Plotting the evaluation metrics for each model

fig, axes = plt.subplots(3, 1, figsize=(12, 18))

metrics = ['mse', 'mae', 'r2']

titles = ['Mean Squared Error (MSE)', 'Mean Absolute Error (MAE)', 'R-squared (R2)']

for ax, metric, title in zip(axes, metrics, titles):

    df_results.plot(x='model_name', y=metric, kind='bar', ax=ax, legend=False, color='skyblue',
edgecolor='black')

    ax.set_ylabel(metric.upper(), fontsize=12)

    ax.set_xlabel('Model Name', fontsize=12)

    ax.set_title(title, fontsize=14)

    ax.tick_params(axis='x', rotation=45)

    ax.grid(True)

# Annotate each bar with the value

for p in ax.patches:

    ax.annotate(f"{p.get_height():.2f}",

        (p.get_x() + p.get_width() / 2., p.get_height()),

            ha='center', va='center',

            xytext=(0, 10),

            textcoords='offset points')

plt.tight_layout()

plt.show()

import pandas as pd

data_set= pd.read_csv('CropPrice.csv')

```

```

print(data_set)

#Extracting Independent and dependent Variable

x= data_set.drop(['max_price'],axis=1)

x=x[['state','district','market','variety','Rainfall']]

y= data_set['max_price']

#Catgorical data

from sklearn.preprocessing import LabelEncoder

x_encoded = x.copy() # Create a copy of x to store the encoded values


label_encoder_x = LabelEncoder()

cols=x.columns

cols.drop('Rainfall')


# Iterate over each column in x and encode it

for column in x.columns:

    x_encoded[column] = label_encoder_x.fit_transform(x[column])

#Catgorical data

print(x_encoded.head())

x=x_encoded

from sklearn.model_selection import train_test_split

x_train, x_test, y_train, y_test= train_test_split(x, y, test_size= 0.25, random_state=0)

print(x.iloc[0])

RM=RandomForestRegressor(n_estimators=10,random_state=0,oob_score=True)


RM.fit(x_train, y_train)

#Prediction of Test and Training set result

y_pred= RM.predict(x_test)

```

```

x_pred= RM.predict(x_train)

print('Random Forest')

print('Train Score: ', RM.score(x_train, y_train))

print('Test Score: ', RM.score(x_test, y_test))

new_data = np.array([x_test.iloc[0]])

print(new_data)

predictions = RM.predict(new_data)

print(predictions)

RI = Ridge(random_state=0)

RI.fit(x_train, y_train)

y_pred = RI.predict(x_test)

x_pred = RI.predict(x_train)


print('Ridge')

print('Train Score: ', RI.score(x_train, y_train))

print('Test Score: ', RI.score(x_test, y_test))

from sklearn.tree import DecisionTreeRegressor


DT = DecisionTreeRegressor()

DT.fit(x_train, y_train)


# Prediction of Test and Training set result

y_pred = DT.predict(x_test)

y_pred_train = DT.predict(x_train)


print('Decision Tree')

print('Train Score:', DT.score(x_train, y_train))

```

```

print('Test Score:', DT.score(x_test, y_test))

from sklearn.preprocessing import PolynomialFeatures
from sklearn.linear_model import LinearRegression
from sklearn.metrics import r2_score

# Create polynomial features
poly = PolynomialFeatures(degree=2)
x_train_poly = poly.fit_transform(x_train)
x_test_poly = poly.transform(x_test)

# Fit linear regression model on polynomial features
poly_reg = LinearRegression()
poly_reg.fit(x_train_poly, y_train)

# Predict on both training and test set
y_pred_train_poly = poly_reg.predict(x_train_poly)
y_pred_test_poly = poly_reg.predict(x_test_poly)

# Calculate R^2 score
train_score_poly = r2_score(y_train, y_pred_train_poly)
test_score_poly = r2_score(y_test, y_pred_test_poly)

print('Polynomial Regression')
print('Train Score:', train_score_poly)
print('Test Score:', test_score_poly)
import matplotlib.pyplot as plt

```

```
models = ['KNN', 'Lasso', 'Ridge', 'Random Forest', 'Decision Tree', 'Linear Regression', 'Polynomial Regression']
```

```
accuracies = [0.75, 0.57, 0.57, 0.98, 0.95, 0.57, 0.54]
```

```
colors = ['blue', 'orange', 'green', 'red', 'purple', 'brown', 'pink']
```

```
plt.figure(figsize=(14, 6))
```

```
bars = plt.bar(models, accuracies, color=colors)
```

```
plt.xlabel('Models')
```

```
plt.ylabel('Accuracy')
```

```
plt.title('Accuracy of Different Models')
```

```
plt.ylim(0.5, 1.0)
```

```
plt.show()
```

```
import joblib
```

```
joblib.dump(RM, "crop_price.pkl")
```

#Crop Yield Prediction:

```
import numpy as np
```

```
import pandas as pd
```

```
import matplotlib.pyplot as plt
```

```
import seaborn as sns
```

```
# loading the dataset
```

```
crop_data=pd.read_csv("dataset.csv")
```

```
crop_data
```

```

# dataset columns

crop_data.columns

# statistical inference of the dataset


crop_data.describe()

# viewing unique crops available in the dataset


print(crop_data['Crop'].unique())

# Dropping missing values
crop_data = crop_data.dropna()

crop_data

import seaborn as sns

import matplotlib.pyplot as plt


# Assuming crop_data is a pandas DataFrame containing Crop_Year and Production columns


sns.barplot(x=crop_data["Crop"], y=crop_data["Yield"])

plt.xticks(rotation=90)

plt.show()

from sklearn import preprocessing

lb = preprocessing.LabelEncoder()

crop_data['State'] = lb.fit_transform(crop_data['State'])

crop_data['Crop'] = lb.fit_transform(crop_data['Crop'])

print(crop_data)

from sklearn.model_selection import train_test_split


x = crop_data.drop(["Yield", "Cost"], axis=1)

```



```

y = crop_data["Yield"]

print(x)

x_train, x_test, y_train, y_test = train_test_split(x, y, test_size=0.25, random_state=5)

x_train

from sklearn.ensemble import RandomForestRegressor

model = RandomForestRegressor()

model.fit(x_train, y_train)

import joblib

joblib.dump(model, 'mymodel_for_cropprice.pkl')

#crop Fertilizer Prediction:

crop_data=pd.read_csv("Fertilizer Prediction.csv")

crop_data

import joblib

joblib.dump(model, 'mymodel_for_cropprice.pkl')

crop_data.columns

import seaborn as sns

import matplotlib.pyplot as plt

# Assuming crop_data is a pandas DataFrame containing State_Name and Production columns

sns.barplot(x=crop_data["Fertilizer Name"], y=crop_data["Soil Type"])

plt.xticks(rotation=0)

plt.show()

import seaborn as sns

import matplotlib.pyplot as plt

# Assuming crop_data is a pandas DataFrame containing Season and Yield columns

```

```

sns.lineplot(x=crop_data["Nitrogen"], y=crop_data["Crop Type"])
plt.xticks(rotation=90)
plt.show()

import seaborn as sns
import matplotlib.pyplot as plt

# Assuming crop_data is a pandas DataFrame containing Season and Yield columns

sns.lineplot(x=crop_data["Phosphorous"], y=crop_data["Crop Type"])
plt.xticks(rotation=90)
plt.show()

import seaborn as sns
import matplotlib.pyplot as plt

# Assuming crop_data is a pandas DataFrame containing Season and Yield columns

sns.lineplot(x=crop_data["Potassium"], y=crop_data["Soil Type"])
plt.xticks(rotation=90)
plt.show()

from sklearn import preprocessing
lb = preprocessing.LabelEncoder()
crop_data['Soil Type'] = lb.fit_transform(crop_data['Soil Type'])
crop_data['Crop Type'] = lb.fit_transform(crop_data['Crop Type'])
print(crop_data)

from sklearn.model_selection import train_test_split
from sklearn.ensemble import RandomForestClassifier

```

```

from sklearn.metrics import accuracy_score

# Assuming crop_data contains your dataset

# Splitting the data into features (x) and target variable (y)
x = crop_data.drop(["Fertilizer Name"], axis=1)
print(x)
y = crop_data["Fertilizer Name"]

# Splitting the data into training and testing sets
x_train, x_test, y_train, y_test = train_test_split(x, y, test_size=0.25, random_state=5)

# Initializing the Random Forest classifier
rf_classifier = RandomForestClassifier(n_estimators=100, random_state=42)

# Training the classifier on the training data
rf_classifier.fit(x_train, y_train)

# Making predictions on the testing data
y_pred = rf_classifier.predict(x_test)

# Calculating the accuracy of the classifier
accuracy = accuracy_score(y_test, y_pred)
print("Accuracy:", accuracy)

```

#Flask Code:

```

from flask import Flask, render_template, request
import joblib

```

```

import json

import numpy as np

import pandas as pd

import pickle

from sklearn.preprocessing import StandardScaler

from sklearn.model_selection import train_test_split

from sklearn.naive_bayes import GaussianNB

from sklearn.ensemble import RandomForestClassifier


app = Flask(__name__)


with open('df_dict.json', 'r') as json_file:

    state_mapping = json.load(json_file)


with open('Crop.json','r') as crop_json:

    crop_mapping=json.load(crop_json)


with open('Fertilizer.json','r') as fertilizer_json:

    fertilizer_mapping=json.load(fertilizer_json)


def predict_price(state, district, market, variety, rainfall):

    # Process the input data and make prediction

    input_data = [[state, district, market, variety, rainfall]] # Input data for the model

    print(input_data)

    prediction = model.predict(input_data)[0] # Use the model to make prediction

```

```

    return prediction

@app.route('/')
def navbar():
    return render_template('navbar.html')

@app.route('/crop')
def home():
    return render_template('index.html')

@app.route('/predict', methods=['GET', 'POST'])
def predict():
    if request.method == 'POST':
        state = request.form['state']
        district = request.form['district']
        market = request.form['market']
        variety = request.form['variety']
        rainfall = request.form['Rainfall']

        # Convert rainfall to float if it's not None
        rainfall = float(rainfall) if rainfall is not None and rainfall != " " else 0.0

        state_number = state_mapping['state'].get(state, None)
        district_number = state_mapping['district'].get(district, None)
        market_number = state_mapping['market'].get(market, None)
        variety_number = state_mapping['variety'].get(variety, None)

```

```
prediction = predict_price(state_number,district_number,market_number,variety_number,
rainfall)
```

```
return render_template('result.html', prediction=prediction)
```

```
# Handle the case where the method is not POST
```

```
return render_template('result.html', prediction=None)
```

```
@app.route('/rem')
```

```
def rem():
```

```
    return render_template('recommend.html')
```

```
@app.route('/rempredict',methods=['POST','GET'])
```

```
def rempredict():
```

```
    N=float(request.form.get('N'))
```

```
    P = float(request.form.get('P'))
```

```
    K = float(request.form.get('K'))
```

```
    temperature = float(request.form.get('temperature'))
```

```
    humidity = float(request.form.get('humidity'))
```

```
    ph= float(request.form.get('ph'))
```

```
    rainfall = float(request.form.get('rainfall'))
```

```
    data = pd.read_csv('Crop_recommendation.csv')
```

```
    data.dropna()
```

```
    ss = StandardScaler()
```

```
    y = data['label']
```

```
    x = data.drop('label', axis=1)
```

```
    ss.fit_transform(x)
```

```
    x_train, x_test, y_train, y_test = train_test_split(x, y, train_size=0.75)
```

```

model = GaussianNB()
model.fit(x_train, y_train)
a=[N, P, K, temperature, humidity, ph, rainfall]
result=model.predict([a])
data=data[data.label !=result[0]]
y = data['label']
x = data.drop('label', axis=1)
ss.fit_transform(x)
x_train, x_test, y_train, y_test = train_test_split(x, y, train_size=0.75)
model.fit(x_train, y_train)
result1 = model.predict([a])
data = data[data.label != result1[0]]
y = data['label']
x = data.drop('label', axis=1)
ss.fit_transform(x)
x_train, x_test, y_train, y_test = train_test_split(x, y, train_size=0.75)
model.fit(x_train, y_train)
result2 = model.predict([a])

return render_template('result1.html',**locals())

@app.route('/pre')
def home3():
    return render_template('yield.html')

@app.route('/yieldpredict', methods=['POST','GET'])
def yieldpredict():

```

```

# Get the input fields from the form

crop = request.form['crop']

state = request.form['state']

cost_of_cultivation = float(request.form['cost_of_cultivation'])

cost_of_production = float(request.form['cost_of_production'])


crop_number = crop_mapping['state'].get(crop, None)

state_number = crop_mapping['district'].get(state, None)

# Use the loaded model to make predictions

predicted_yield = model1.predict([[crop_number, state_number, cost_of_cultivation,
cost_of_production]])


return render_template('result3.html', crop=crop, state=state, predicted_yield=predicted_yield[0])

@app.route('/ferti')

def home4():

    return render_template('fertilizer.html')


@app.route('/ferpredict', methods=['POST','GET'])

def ferpredict():

    Temperature= float(request.form['Temparature'])

    Humidity= float(request.form['Humidity'])

    Moisture= float(request.form['Moisture'])

    Soil_Type = request.form['Soil Type']

    Crop_Type= request.form['Crop Type']

    Nitrogen = float(request.form['Nitrogen'])

    Potassium = float(request.form['Potassium'])

```



```

Phosphorous= float(request.form['Phosphorous'])

soil_number = fertilizer_mapping['Soil_type'].get(Soil_Type, None)
crop_number = fertilizer_mapping['Crop_type'].get(Crop_Type, None)
# Use the loaded model to make predictions
Fertilizer = model2.predict([[soil_number, crop_number,
Temperature, Humidity, Moisture, Nitrogen, Potassium, Phosphorous]])

return render_template('result4.html', Soil_Type=Soil_Type, Crop_Type= Crop_Type,
Fertilizer=Fertilizer[0])
if __name__ == '__main__':
    model=joblib.load("crop_price.pkl")
    model1=joblib.load("mymodel_for_cropprice.pkl")
    model2=joblib.load("Fertilizer.pkl")
    app.run(host='0.0.0.0', port=5000, debug=True)

```

9. RESULT ANALYSIS

9.1 Crop price:

Crop cost prediction is a multifaceted process influenced by various factors such as geographical location, state policies, district-level conditions, crop varieties, and market dynamics. The cost prediction of crops varies significantly from one state to another due to differences in climate, soil quality, and agricultural practices. Similarly, within a state, different districts may have distinct agricultural landscapes, infrastructure, and resource availability, leading to varying cost predictions.

Moreover, the prediction of crop costs is contingent upon the specific crop being cultivated. Different crops have unique growth requirements, market demand, and price fluctuations, all of which impact cost projections. Additionally, market trends and demand-supply dynamics play a crucial role in determining the profitability of cultivating certain crops.

One of the critical factors influencing crop cost prediction is rainfall patterns. Rainfall directly affects crop yields and quality, thereby influencing input costs such as irrigation, fertilizers, and pest control measures. Regions experiencing inadequate rainfall may incur higher expenses to ensure crop survival, while those with ample rainfall may have lower input costs.

Overall, accurate crop cost prediction demands a comprehensive understanding of regional variations, crop-specific factors, market dynamics, and environmental conditions such as rainfall patterns. By considering these diverse elements, policymakers, farmers, and stakeholders can make informed decisions to optimize agricultural productivity and economic outcomes.

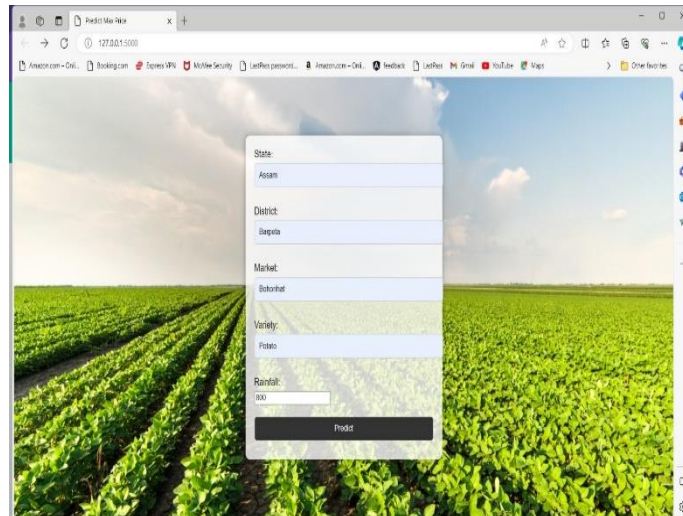


Fig 9.1 Crop price

Now we can see the how the crop prediction is working the major states of the crops can be seen by using this technique of cost prediction. For example we have taken the random state Assam in which the cost prediction and the yield of the crop is different. And his price varies from one state to another state. Here we have also tried crop prediction on different states like as Andhra Pradesh, Kerala, Tamil Nadu and etc.

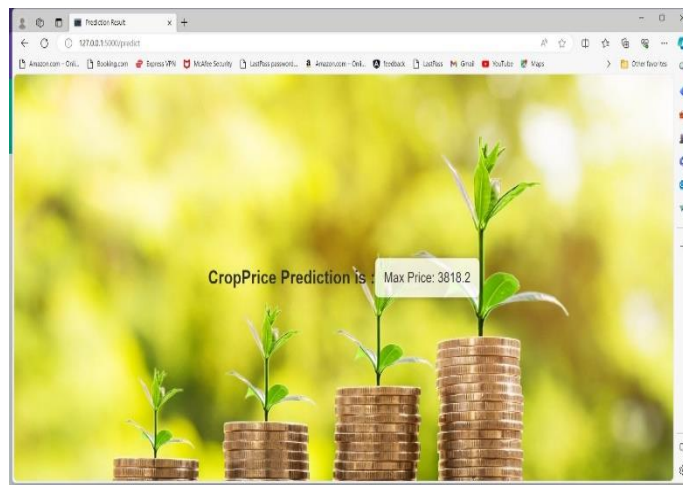


Fig9.1 Crop price result

The cost prediction of crops, exemplified by the case of potato cultivation in Assam, underscores the nuanced variations in agricultural economics across different states. In Assam, the maximum market price for potatoes is estimated at approximately 3818.2, reflecting the culmination of various factors

specific to the region. However, it's crucial to recognize that this price point is not universal and cannot be replicated in other states.

Each state possesses unique agro-climatic conditions, soil types, irrigation facilities, and market dynamics, all of which contribute to distinct costs of crop cultivation and market prices. Additionally, the cultivation process itself varies from crop to crop and from region to region, influencing input costs, labor requirements, and yield potentials.

Therefore, the cost prediction of crops encapsulates a complex interplay of factors, including geographical location, state-specific policies, infrastructure, technology adoption, and market demand. While Assam may fetch a certain price for potatoes, the same crop cultivated elsewhere may command a significantly different value due to diverse contextual factors.

Understanding and adapting to these variations is essential for stakeholders involved in agriculture, enabling informed decision-making, resource allocation, and sustainable production practices tailored to the specific requirements and potentials of each region.

9.2Crop Recommendation

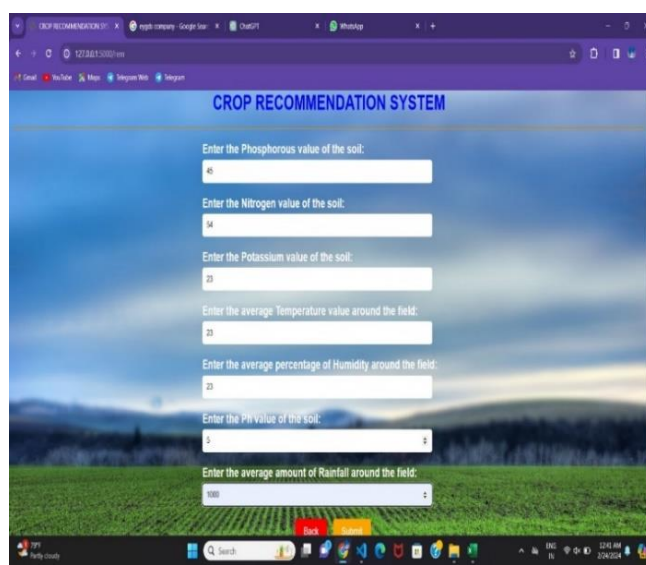
The image shows a web browser window displaying a 'CROP RECOMMENDATION SYSTEM' interface. The page has a purple header with the title 'CROP RECOMMENDATION SYSTEM' in white. Below the header, there is a form with seven input fields, each with a label and a value: 'Enter the Phosphorous value of the soil:' (45), 'Enter the Nitrogen value of the soil:' (34), 'Enter the Potassium value of the soil:' (23), 'Enter the average Temperature value around the field:' (23), 'Enter the average percentage of Humidity around the field:' (23), 'Enter the PH Value of the Soil:' (5), and 'Enter the average amount of Rainfall around the field:' (100). At the bottom of the form, there are two buttons: 'Back' and 'Submit'. The background of the page is a landscape image of a green field under a blue sky with clouds. The browser's address bar shows the URL '127.0.0.1:5000/index.html'.

Fig9.2 Crop recommendation

Crop recommendation is a highly intricate process influenced by a myriad of factors, as depicted in the provided image, particularly focusing on nitrogen and phosphorus levels in the soil. The essence lies in understanding that crop recommendations vary significantly across different soil types and pH levels. Soil type serves as a fundamental determinant, as each type harbors distinct nutrient compositions, drainage properties, and fertility levels. Similarly, pH levels profoundly impact nutrient availability and plant uptake, thereby necessitating tailored recommendations for optimal crop growth.

Moreover, crop recommendations hinge on a multitude of considerations, encompassing crop varieties, market dynamics, and economic viability. Different crops exhibit varying nutrient requirements, growth patterns, and tolerance levels, necessitating customized recommendations to maximize yields and profitability.

The significance of soil type cannot be overstated in crop recommendation algorithms. Soil-specific data guides precision agriculture practices, ensuring efficient resource utilization and minimizing environmental impact. By factoring in soil type alongside other variables, such as nutrient levels and pH, agricultural experts can deliver targeted recommendations tailored to the unique needs and constraints of each agricultural parcel.

In essence, the efficacy of crop recommendation systems relies on a holistic integration of soil science, agronomy, market analysis, and technological advancements. By leveraging comprehensive datasets and advanced analytics, stakeholders can optimize crop selection and management practices, fostering sustainable agricultural production while enhancing farm profitability and resilience.

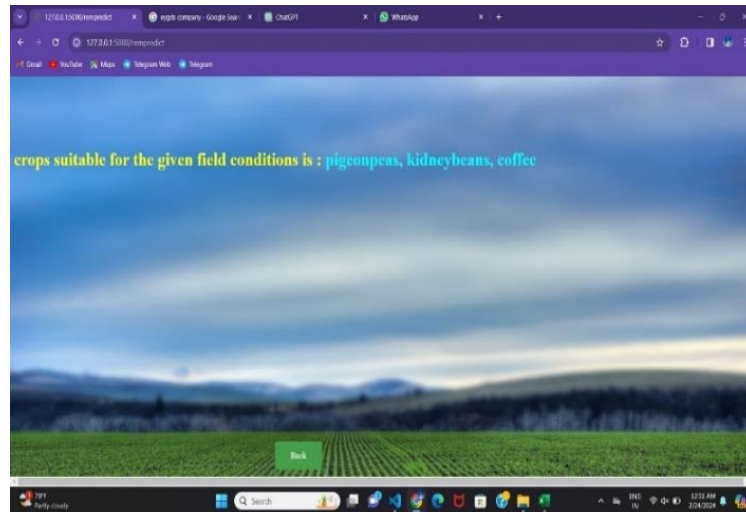


Fig 9.2 Crop recommendation result

Crop recommendation systems utilize diverse soil and environmental factors, including potassium and nitrogen levels in the soil, pH value, phosphorus value, temperature, and rainfall amount, to advise farmers on the most suitable crops for cultivation in a specific area. By analyzing these factors, the system can make informed predictions, optimizing agricultural productivity and sustainability. For instance, based on the input parameters provided, such a system might recommend coffee and kidney beans as the ideal crops for a given location. This approach not only aids in maximizing yield but also supports sustainable farming practices by aligning crop selection with the prevailing environmental conditions.

9.3 Crop Fertilizer Prediction:

The image showcases the results of a sophisticated crop fertilizer prediction model, which takes into account four key parameters: humidity level, soil type, moisture content, and crop type. This model represents a significant advancement in precision agriculture, offering farmers and agricultural planners invaluable insights into optimizing fertilizer usage.

Humidity level serves as an indicator of environmental conditions, influencing crop growth and nutrient uptake. Soil type, with its varying nutrient compositions and water-holding capacities, profoundly

impacts fertilizer requirements. Moisture content in the soil directly affects nutrient availability to plants, influencing the efficacy of fertilization practices.

Additionally, crop type plays a pivotal role in determining fertilizer needs, as different crops have distinct nutrient requirements at various growth stages. By analyzing these four attributes in conjunction, the prediction model can accurately estimate the optimal fertilizer dosage for a specific crop under prevailing environmental conditions.

This predictive capability empowers farmers and agricultural planners to fine-tune their resource management strategies, ensuring precise application of fertilizers to maximize crop yields while minimizing environmental impact and input costs. By tailoring fertilizer application to the specific needs of each crop and field, farmers can enhance efficiency, sustainability, and ultimately, profitability.

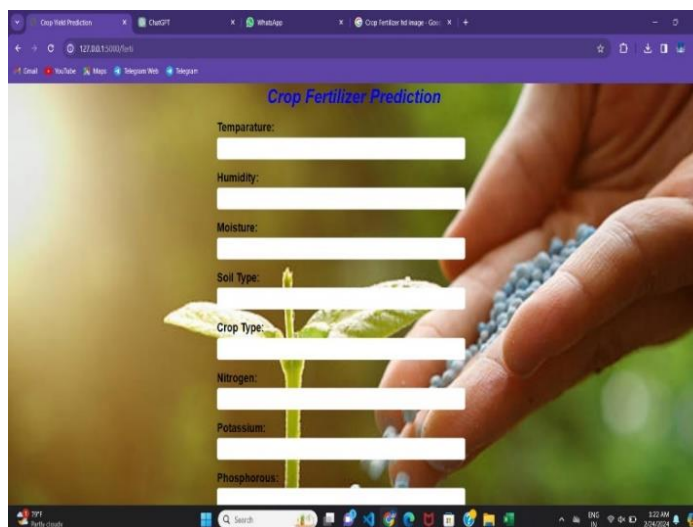
The image shows a web browser window with multiple tabs. The active tab is titled 'Crop Fertilizer Prediction'. The webpage has a purple header with the title 'Crop Fertilizer Prediction' in white. The background of the page is a blurred image of a hand holding blue fertilizer granules over a green plant. On the left side, there is a vertical list of input fields with labels: 'Temperature:', 'Humidity:', 'Moisture:', 'Soil Type:', 'Crop Type:', 'Nitrogen:', 'Potassium:', and 'Phosphorous:'. Each label is followed by a white rectangular input box. The browser's address bar shows the URL '127.0.0.1:5000/Veri'. The Windows taskbar is visible at the bottom, showing the time as 12:24 PM on 2/26/2024.

Fig9.3 Crop fertilizer prediction

The image displays the outcome of a crop fertilizer prediction model, highlighting the recommended fertilizer for a specific crop based on its soil type. This result provides critical information, including the name of the crop, the type of soil it thrives in, and the optimal fertilizer to use. Such detailed insights empower farmers to select the most suitable fertilizer for their crops, ensuring healthier growth and potentially higher yields by aligning agricultural practices with precise soil and crop needs.

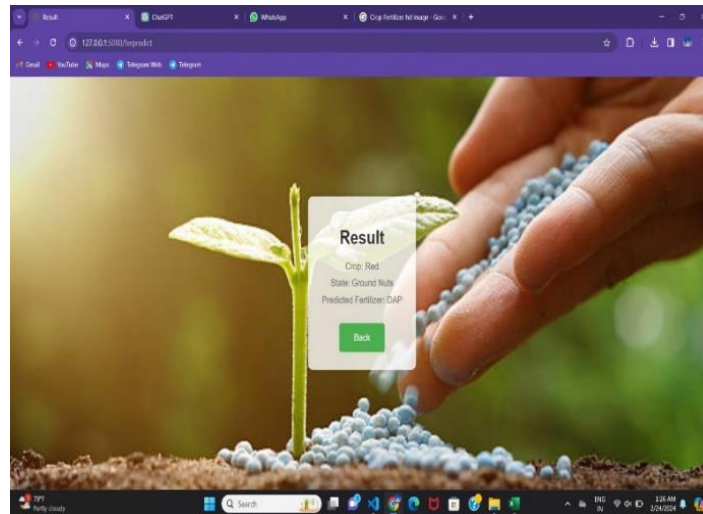


Fig 9.3 Crop fertilizer result

9.4 Crop Yield Prediction:

The image depicts a crop yield prediction model that integrates four crucial attributes: crop name, cultivation state, cultivation and production expenses. This model represents a pivotal advancement in agricultural analytics, offering farmers and planners actionable insights to optimize resource allocation and enhance profitability.

Crop name serves as a fundamental input, as different crops exhibit varying growth patterns, nutrient requirements, and susceptibility to environmental factors. The cultivation state provides context-specific information, considering variations in climate, soil quality, and agricultural practices. Expenses related to cultivation and production encompass a wide range of factors, including inputs like seeds, fertilizers, pesticides, labor costs, and equipment usage.

By analyzing these attributes collectively, the model generates accurate predictions of crop yield, enabling farmers and agricultural planners to make informed decisions. These insights facilitate optimal resource allocation, ensuring that inputs are applied judiciously to maximize productivity while minimizing costs and environmental impact.

Ultimately, the crop yield prediction model empowers stakeholders to fine-tune their strategies and mitigate risks associated with agricultural production. By leveraging data-driven insights, farmers can enhance efficiency, resilience, and profitability, contributing to sustainable agriculture practices and food security.

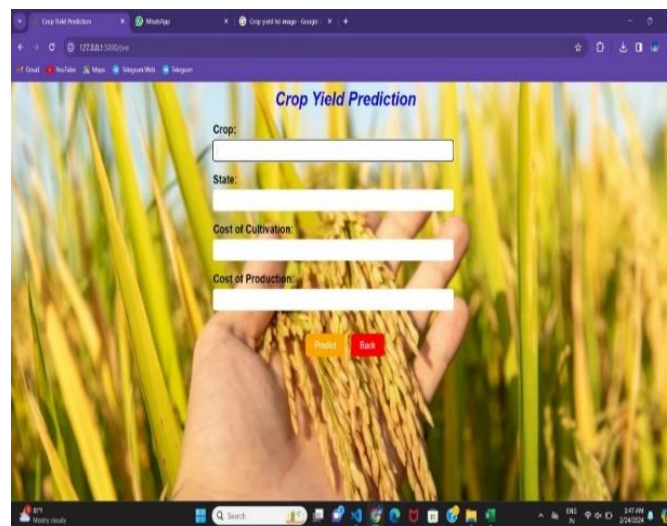


Fig 9.4 Crop yield prediction

The image showcases the results of a crop yield prediction, displaying critical information such as the crop name, the state where it was cultivated, and the total yield production measured in hectares. This output provides essential insights, enabling farmers and agricultural stakeholders to understand the potential productivity of different crops across various states, thereby facilitating informed decision-making for future cultivation strategies and resource allocation.

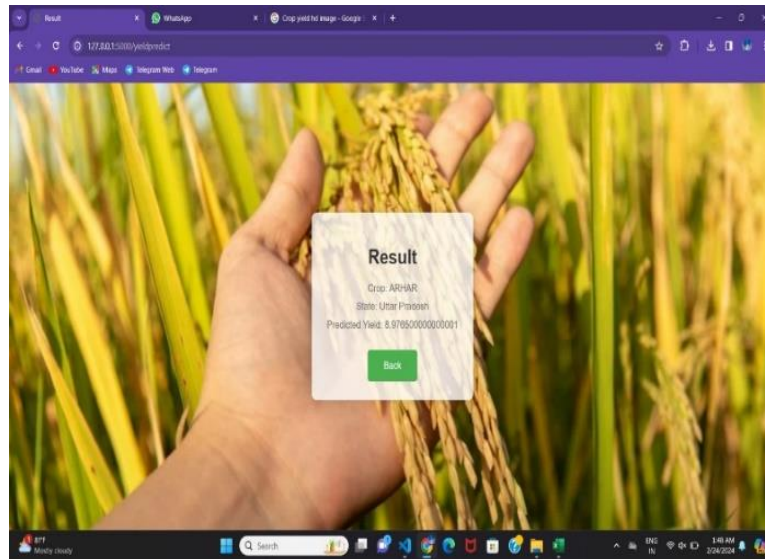


Fig9.4 Crop yield result

10. FUTURE SCOPE

Throughout this extensive project, the utilization of machine learning algorithms, particularly Random Forest, has revolutionized agricultural practices. The project's core focus was to evaluate the performance of various algorithms—KNN, Random Forest, Linear Regression, Ridge, Lasso, and Decision Tree—for predicting crop yields. Among the algorithms utilized, Random Forest proved to be the most effective, achieving an impressive accuracy rate of 98%. This high level of accuracy is particularly beneficial for predicting yields of extreme crop types, providing valuable insights into previously unexplored cultivation opportunities. Beyond yield prediction, the project explored the versatility of machine learning models in predicting crop prices, recommending suitable crops for cultivation, and determining optimal fertilizer requirements. These applications are instrumental in streamlining farming processes, empowering farmers with tailored insights for maximizing crop yield and profitability. By leveraging machine learning algorithms, farmers can make informed decisions that enhance agricultural productivity and sustainability. A key highlight of the project was the development of a userfriendly web application, designed to provide seamless access to machine learning predictions. With a testing accuracy exceeding 98%,

the application demonstrates the reliability and effectiveness of machine learning in predicting crop yields. This high level of accuracy instills confidence in farmers, encouraging the adoption of technology-driven practices for improved agricultural outcomes. The implications of this project are profound, offering a glimpse into the future of agriculture. By harnessing the predictive power of machine learning, farmers can optimize their crop selection, resource allocation, and cultivation practices. This not only leads to higher yields and profitability but also promotes sustainable agricultural practices. To conclude, this project underscores the transformative possibilities of machine learning within agriculture. By providing accurate predictions for crop yields, prices, and fertilizer requirements, machine learning empowers farmers to make informed decisions that drive agricultural innovation and durability. As the agricultural industry advances, the adoption of machine learning technologies evolves alongside it promises to revolutionize farming practices, ensuring a future for agriculture that is stronger and more productive..

11. CONCLUSION

Throughout this extensive project, the utilization of machine learning algorithms, particularly Random Forest, has revolutionized agricultural practices. The project's core focus was to evaluate the performance of various algorithms—KNN, Random Forest, Linear Regression, Ridge, Lasso, and Decision Tree—for predicting crop yields. Among the algorithms utilized, Random Forest proved to be the most effective, achieving an impressive accuracy rate of 98%. This high level of accuracy is particularly beneficial for predicting yields of extreme crop types, providing valuable insights into previously unexplored cultivation opportunities.

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