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Assessment of Air Pollutants of Dhanbad Using Machine Learning **Techniques**

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Abstract: A comprehensive assessment of air pollution in Dhanbad for the months of April 2019 through March 2023 was conducted using a support vector machines. random forest, xgboost, and decision tree methods in relation to seven main pollutants into the air (PM10, NO, NO2, NH3, SO2, CO, and O3). A randomly selected 30-day period was used to create line and bar graphs using accuracy & error matrices for 7 pollutants in various models. Our investigation found that the Random Forest model estimates Dhanbad air contaminants with the lowest MAE. Among the aforementioned models, the Random Forest one stands head and shoulders above the others.

Keywords: Machine learning models, air quality, air pollutants, CPCB, data preprocessing, performance criteria

1. Introduction

Air is one of the basic elements of PanchTatva for surviving on the surface of the earth because the life systems are dependent on smooth as well as harmonious functions of basic five elements of PanchTatva. It is therefore, necessary to maintain purity as well as quality of air for the existence of civilization. However, quality of air has been deteriorating continuously because of unplanned industrial as well as infrastructural development and deforestation in many countries including India. In particular, increase of PM₁₀, CO, SO₂, NO, NH₃ etc. have attained threatened level and hence, reduction of the quantity of air pollutants are utmost urgent for the betterment of the society and civilization. It is therefore, necessary to analyze the air quality of different places and necessary measures should be taken accordingly.

In this work, we consider sevenair pollutants such as $PM_{10}[1]$, NO[2]], $NO_2[3]$, $NH_3[4]$, $SO_2[5]$, CO[6], and O₃[7]for comparison of different models of Machine Learning algorithm. We have chosen five important models like Linear Regression, Support Vector Machine,

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Random Forest, XGBoostand Decision Tree foranalyzing the data available for estimation of air pollutants of Dhanbad.

Data on the seven pollutants have been gathered throughout the period of time from the Central Pollution Control Board (CPCB)[8]website from April 2019 to March 2023 in order to analyze the air pollutants of Dhanbad.

According to the current investigation, the Random Forest algorithm is the most accurate model for forecasting pollutant concentrations, with the exception of NO2, where XGBoost demonstrates the highest accuracy. Thus, based on data available throughout the aforementioned annular period, it can be said without reservation that the Random Forest technique provides highest accuracy with minimal mistakes for measuring the air contaminants of Dhanbad.

The plan for the paper is presented after this. In the second section, 5 machine learning algorithms are briefly explained. The third section covers the work's methodology, while the fourth section covers results and a summary. A conclusion is in the fifth section.

2. Machine Learning Models

Machine learning models show how computers can categorise, anticipate, and evaluate physical occurrences and their qualities by training them with complex data. This research builds a variety of models using various methods to predict efficient pattern recognition & model self-learning. Five algorithms' fundamentals presented here.

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2.1. Linear Regression (LR)

Linear regression is the first machine learning method most academics use [9]."Linear regression" is supervised machine learning that identifies the linear connection between a dependent variable & one or more variables that are independent. The approach finds the best linear equation to predict the dependent variable given the independent variables. The equation shows a straight line between independent and dependent variables.

2. 2. Support Vector Machine [SVM]

Support vector machines are supervised learning methods for regression, classification, and more [10]. The hyperplane divides data points into intervals from which the model output may be calculated. SVMs come in two varieties. Regression problems are different from classification problems. Most issues are nonlinear, therefore the kernel approach converts low-dimensional data to high-dimensional data multidimensional feature space. Kernels are used to separate indivisible input data. such as linear, Use polynomial, RBF, sigmoid, hyperbolic tangent, etc. Radial Basis Function Kernel is more accurate than other kernels, hence we used it [11].

2.3. The Random Forest (RF)

An other well-known machine learning approach is random forest[12]. This method employs supervised ensemble learning to finish problems related to regression or classification. Using a training set of data, it constructs many decision trees. The mean value of the decision tree set is then used to forecast the value for fresh input data.

2.4. XGBoost, or extreme gradient boosting

Another significant machine learning method is XGBoost[13]. The machine learning technique XGBoost combines a set of decision trees with gradient boosting to generate predictions. The XGBoost method uses the community-based weak learning technique to solve regression and classification problems. This method produces useful results since it relies predictions on parallel tree structures and takes hardware and software parameters into account while building them.

2.5.Decision Tree algorithm or DT

Decision tree machine learning is used for regression and classification[14]. This tree-like model has leaf nodes that represent class labels or numerical values, inner nodes that represent attributes or features, and branch nodes that provide decision rules depending on those features.

3. Methodology

Present section deals with the collection and Preprocessing of data, designing of models and criteria of performance for the current work.

3.1. Assemblage of Data

Datasets for the current investigation were gathered from the Central Pollution Control Board (CPCB), New Delhi, website as well asContinuous Ambient Air Quality Monitoring Station (CAAQMS),Dhanbad, Jharkhand. Particulate matter (PM₁₀), nitric oxide (NO), nitrogen dioxide (NO₂), ammonia (NH₃), sulphur dioxide (SO₂), carbon monoxide (CO), and ozone (O₃) are the seven features in the datasets, which also include meteorological parameters and pollutant concentrations.

3.2. Pre-processing of Data

High-quality data and effective data representations are key factors in a model's effectiveness. The following explanations pertain to crucial data pre-processing procedures, which include missing value incorporation, outlier elimination, feature scaling, and feature selection [15].

- (i) Incorporation of Missing Values: The nearest data points should be taken into consideration for analysis in lieu of the missing data, which should be replaced by linear interpolation estimation.
- (ii) Removal of Outliers: For CO, an unusual pattern was seen for the period from October to November in 2019. We have not considered values of CO for this period in the present work.
- (iii) **Features scaling:** One crucial stage in the preprocessing of data before creating a model is feature scaling. The range of characteristics is normalised or standardised using it. Min-max scaling has been used in this research to normalise the values between 0 and 1. The following is the normalisation formula: $Z' = \frac{Z Z_{min}}{Z_{max} Z_{min}}$

Z is the initial value, Z^{\prime} is the normalised value, $[\![Z]\!]$ _max is the maximum value that can be achieved, & Z_min is the lowest value that can be achieved for the feature.

(iv)Selection of features: A portion of the extensive feature space's features are selected in order to lower the dataset's dimensionality. This phase uses Pearson's Correlation coefficient to choose the characteristics.

3.3.Designing of Model

The model is created on a 64-bit machine running Python 3.7.6[16] on a 2.2GHz Intel Core i3 CPU with 4GB RAM. Using common charting tools like matplotlib[17],

seaborn[18], and sklearn[19], the model is trained and forecasted. A large portion of the work was implemented using Panda's dataframe. We were able to use the training set to train and evaluate the previously mentioned models by splitting the dataframe into a testing set and a training set with a ratio of 7:3.

3.4. Models'Performance Standards

Prediction models are examined for effectiveness using statistical matrices like Root Mean Squared Error or RMSE, Mean Absolute Error or MAE, Mean Squared Error or MSE, & R-Squared or R²[20].

R-Squared or R²: R-Squared, often known as R², is a statistical measure that illustrates how well the model and data fit in a regression model. A higher R2 value indicates a better model. The formula used to determine

$$R^2 = 1 - \frac{(x_i - x_i)^2}{(x_i - x_i)^2}$$

Mean Absolute Error /MAE is a statistical phrase that measures the average of the absolute differences between the observed values and the values predicted for the whole collection of data. The formula used to calculate the MAE is

$$MAE = \frac{1}{n} \sum_{i=1}^{n} |x_i - x_i|$$

Mean Squared Error/MSE total of the squared deviations for all values in the data set, both observed and anticipated.

$$MSE = {}_{n}^{1} \sum_{i=1}^{n} (x_{i} - \hat{x}_{i})^{2}$$

Squared Error of Root Mean: As one of the most frequently employed metrics for assessing the precision of predictions, RMSE (root mean squared error) is frequently utilised. Sqrt (mean squared error) It may be expressed as follows:

RMSE =
$$\sqrt{\frac{1}{n}} \sum_{i=1}^{n} (x_i - \hat{x}_i)^2$$

The expected x-value is represented by by $x^{\perp \wedge}$, the actual value of x is represented by x_i in all of the above formulas, and the number of mistakes is indicated by 'n'.

4. Result and Summary

The accuracy, mean squared error, root mean square error, & mean absolute error are now shown for each of the five machine learning models, which cover seven distinct contaminants, that we previously examined. To make things simple, line and bargraphs covering a 30day period selected at random have been made. The analytical assessment of the air quality in Dhanbad is shown in the diagrams.

4.1 PM₁₀

Estimation of accuracy and error metrics for PM₁₀ in different models

+
Coefficient of determination R^2 on Test Data
0.713524
0.56395
0.738381
0.657383
0.71769

Table 1 (a): Estimation of accuracy for PM₁₀ in different models

+		+	+
Models	MAE	MSE	
	0.0476522	0.00580864	0.0762145
Decision Tree		0.00904255	
Random Forest	0.0445463	0.00529851	
Support Vector Machine	0.0595149	0.00694697	0.0833485
XGBoost	0.0478898	0.00572418	0.0756583

Table 1 (b): Estimation of error metrics for PM₁₀ in different models

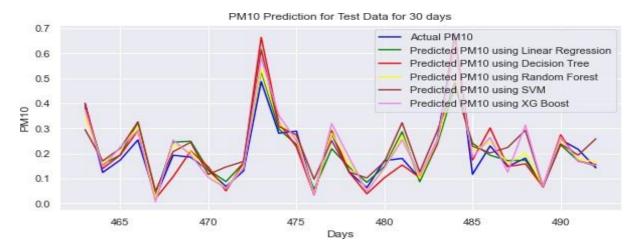


Fig 1 (a): Line graph Comparing PM_{10} levels between the models' actual and predicted levels

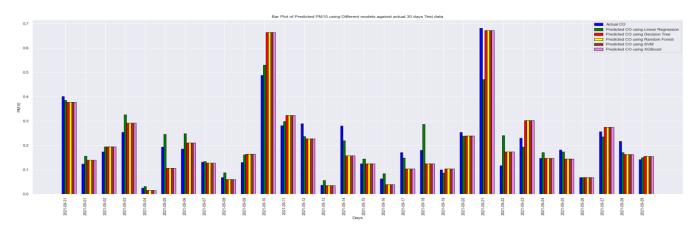


Fig 1 (b): Bar graph comparing PM₁₀ levels between models' actual and anticipated values

4.2. NO₂

Estimation of accuracy and error metrics for NO2 in different models

Models	Coefficient of determination R^2 on Test Data
Linear Regression	0.886757
Decision Tree	0.940375
Random Forest	0.961305
Support Vector Machine	0.808075
XGBoost	0.966593

Table 2 (a): Estimation of accuracy for NO2 in different models

Models	MAE		
	0.0205374	0.00186783	0.0432185
	0.00847022	0.000983462	0.0313602
	0.00894517	0.000638241	0.0252634
Support Vector Machine	0.0465254	0.00316562	0.0562639
XGBoost	0.00917657	0.000551015	0.0234737

Table 2 (b): Estimation of error metrics for NO₂ in different models

NO2 Prediction for Test Data for 30 days

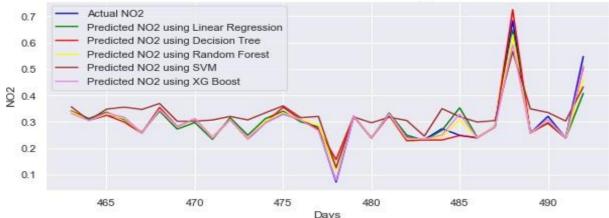


Fig 2 (a): A line graph that compares the models' real and anticipated NO₂ levels

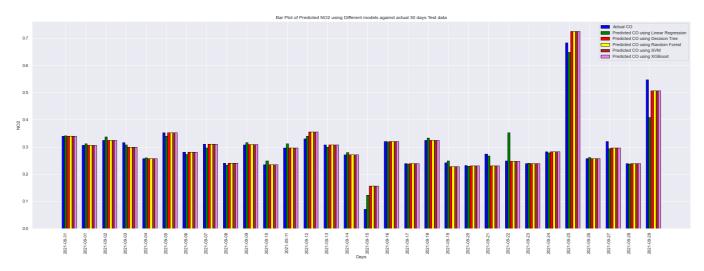


Fig 2 (b): Bar graph showing how the models' actual and anticipated NO₂ levels compare

4.3.NO

Estimation of accuracy and error metrics for NO in different models

Models	Coefficient of determination R^2 on Test Data
Linear Regression	0.846018
Decision Tree	0.746021
Random Forest	0.850919
Support Vector Machine	0.799639
XGBoost	0.822513

Table 3 (a): Estimation of accuracy for NO in different models

+	+	+	
Models	MAE	MSE	RMSE
Linear Regression	0.0241867	0.00255732	
Decision Tree	0.0269907	0.00375439	
Random Forest	0.0222771	0.00254159	0.0504142
Support Vector Machine	0.0452136	0.00332758	0.0576851
XGBoost	0.0245447	0.00294768	0.0542926

Table 3 (b): Estimation of error metrics for NO in different models

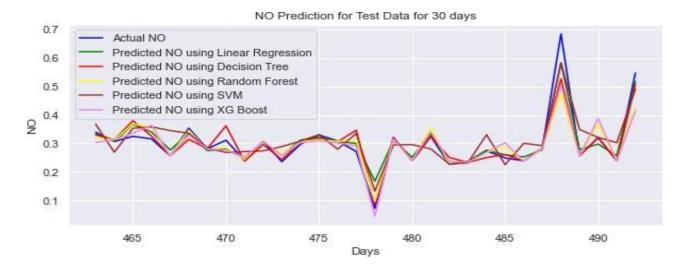


Fig 3 (a): A line graph comparing the models' actual and anticipated levels of NO

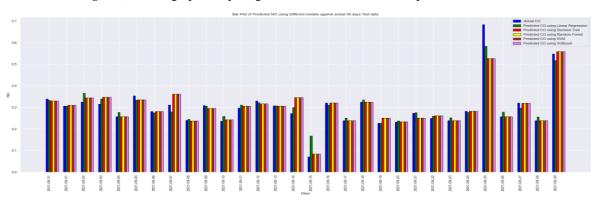


Fig 3 (b): Bar graph comparing the models' actual and anticipated levels of NO

4.4. NH₃

Estimation of accuracy and error metrics for NH₃ in different models

Models	Coefficient of determination R^2 on Test Data
Linear Regression	0.82798
Decision Tree	0.834401
Random Forest	0.863692
Support Vector Machine	0.768598
XGBoost	0.849097

Table 4 (a): Estimation of accuracy for NH₃ in different models

Models	MAE	MSE	RMSE
		0.00296068	
Decision Tree		0.00292416	0.0540755
Random Forest		0.00236551	0.0486365
Support Vector Machine	0.0461031		0.0631087
XGBoost	0.019611	0.00259721	0.0509629

Table 4 (b): Estimation of error metrics for NH₃ in different models

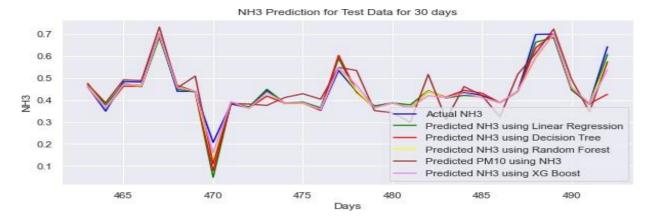


Fig 4 (a): Line graph comparing the models' actual and expected values for NH₃

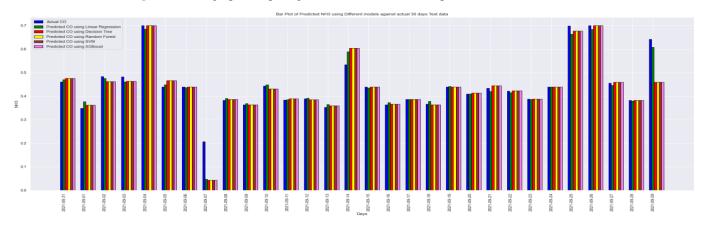


Fig 4 (b): Bar graph comparing the models' actual and expected values for NH₃

4.5. SO₂

Estimation of accuracy and error metrics for SO₂ in different models

Models	Coefficient of determination R^2 on Test Data
Linear Regression	0.83404
Decision Tree	0.770791
Random Forest	0.885418
Support Vector Machine	0.813137
XGBoost	0.883537

Table 5 (a): Estimation of accuracy for SO₂ in different models

Models	MAE	MSE	
Linear Regression	0.0400133	0.00484021	0.0695716
Decision Tree	0.0431597	0.0057056	0.0755355
Random Forest	0.0337864	0.0033607	0.0579716
Support Vector Machine			
XGBoost		0.00339664	0.0582807

Table 5 (b): Estimation of error metrics for SO₂ in different models

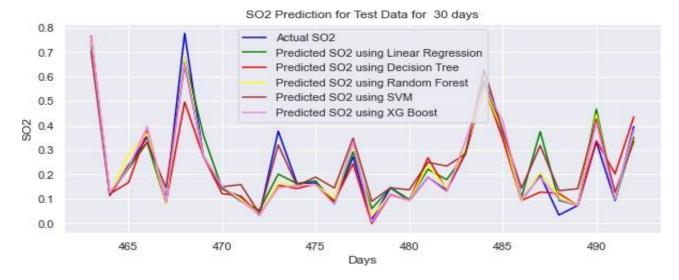


Fig 5 (a): A line graph comparing the models' estimated and real SO₂ levels

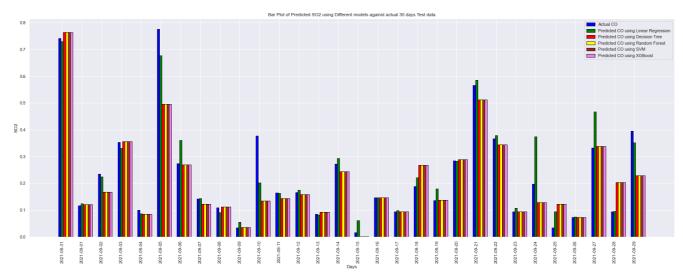


Fig 5 (b): Bar graph showing how the models' anticipated and real SO₂ levels compare

4.6. CO

Assessment of precision and inaccuracy measures for CO in various models

Models	Coefficient of determination R^2 on Test Data
Linear Regression	0.921752
Decision Tree	0.848762
Random Forest	0.928037
Support Vector Machine	0.872661
XGBoost	0.907438
Support Vector Machine	

Table 6 (a): Accuracy estimation for CO in various models

Models	+			
Decision Tree	Models	MAE	MSE	RMSE
Random Forest	Linear Regression	0.0254029	0.00170739	0.0413206
Support Vector Machine 0.0405885 0.00277858 0.0527122	Decision Tree	0.0274254	0.00309524	0.0556349
†	Random Forest	0.0187556	0.00157835	0.0397284
XGBoost	Support Vector Machine	0.0405885	0.00277858	0.0527122
	XGBoost	0.0206101	0.00201973	0.0449414

Table 6 (b): Estimation of error metrics for CO in different models

CO Prediction for Test Data for 30 days

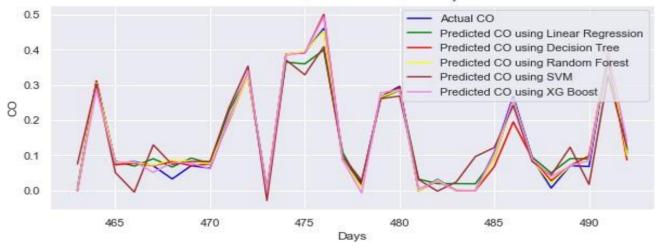


Fig 6 (a): A line graph that compares the models' actual and expected CO levels

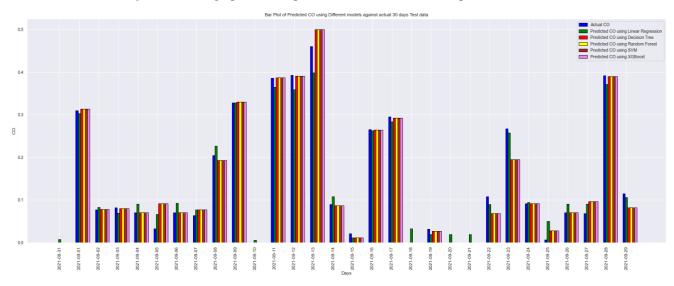


Fig 6 (b): Bar graph for comparing the models' actual and expected CO level

4.7. O₃

Estimation of accuracy and error metrics for O₃ in different models

Models	+
Linear Regression	0.471247
Decision Tree	0.151841
Random Forest	0.488067
Support Vector Machine	0.41508
XGBoost	0.458913

Table 7 (a): Estimation of accuracy for O₃ in different models

Models	MAE	MSE	
Linear Regression	0.0592842	0.00998107	0.0999053
	0.0717411	0.0162444	0.127454
Random Forest	0.0543182	0.00952098	0.0975755
Support Vector Machine	0.0705496	0.0110413	0.105078
XGBoost		0.0102139	

Table 7 (b): Estimation of error metrics for O₃ in different models

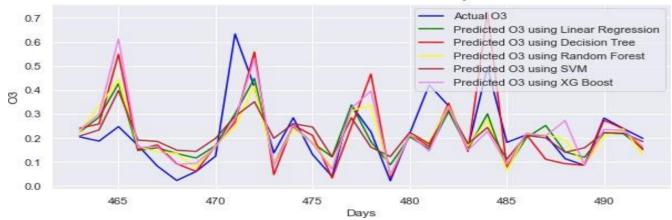


Fig 7 (a): A line graph that compares the models' actual and anticipated levels of O₃

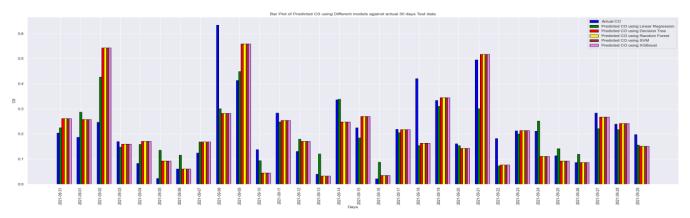


Fig 7 (b): Bar graph comparing the models' actual and anticipated levels of O₃

Examining the data and graphs for the various models stated above, it is evident that Random Forest has the highest accuracy for all pollutants, with the exception of NO₂, for which XGBoost is shown to have the highest accuracy. The results indicate that Random Forest produces the least amount of error for each of the seven contaminants in terms of Mean Absolute Error, or MAE.Once again, in terms of MSE and RMSE, the Random Forest model performs better for six pollutants (NO2 excluded). Regarding the results, XGBoost outperforms the Random Forest model in terms of NO2.

5. Conclusions

We have assessed Dhanbad's air quality for the period of April 2019 to March 2023, utilising five key machine learning-based prediction models to analyse pollutants such as PM₁₀, NO, NO₂, NH₃ SO₂, CO, and O₃. After carefully analysing Dhanbad's air quality rating, our purpose was to compare and contrast five different machine learning methods. Our research leads us to the conclusion that the Random Forest model is most suited for assessing the air pollutants in Dhanbad, namely PM₁₀, NO, SO₂, NH₃, CO, and O₃. On the other hand, XGBoost yields the most accurate findings for NO₂. The most inaccurate model is the Random Forest. Thus,

Dhanbad's air quality is best assessed using the Random Forest model..

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