An Approach towards Forecasting Time Series Air Pollution Data Using LSTM-based Auto-Encoders

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Abstract

Artificial Intelligence-based algorithm is used extensively for predicting the concentration of different pollutants under various conditions. Recently, Long-Short Term Memory (LSTM) and its variant is getting popular attention related to the prediction of Air Quality Index (AQI) across various polluted cities. The accuracy of the prediction is found to be depending on the processing step of input data. Here we present a study of combining both Random Forests (RF) based regression for data pre-processing step and multi-variate time series coupled with Multistep Multiwindow LSTM with auto encoder and decoder to predict the pollutant concentration in the urban city area of Bengaluru. In this approach, the RF algorithm is used for imputing the missing values of the input vector. We have implemented this technique from the data collected from the Karnataka State Pollution Control Board (KSPCB), for the city limits of Bengaluru with four years of data from 2019 to 2022, mainly focusing on the six regions where pollution is found to be maximum. We found that our Multistep Multivariate LSTM-based autoencoder gives better accuracy than the conventional LSTM data based on a single-step pipeline model with respect to Recall and F-Score values.

Keywords: LSTM, Auto-encoder, Time Series forecasting, Arithmetic Mean, Precision, Recall, F-Score.

1 Introduction

Air pollution is amongst the most grievous problems affecting the world today, and has a greater impact on human health and the economy. The pollution of the environment is caused when physical and biological components of land, water, and air are contaminated directly or indirectly which adversely affects the normal processes of the environment, resulting in biodiversity loss. Air pollution, water pollution, and land pollution together constitute environmental pollution. Pollution from noise, plastic,

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and light affect the biosphere as well (Brahmaiah et al., 2021; Chu & Karr, 2017; Niranjan & Rakesh, 2021; Pushpavalli et al., 2024). Air Quality Index (AQI) is decided depending on the concentrations of pollutants like Particulate Matter (PM10, PM2.5) and gases (NO₂, O₃, CO₂, SO₂, NH₃, Pb) suspended in the atmosphere as shown in Table 1.

AOI Category	PM10 24-	PM2 5	NO. 24-	O. 8-hr	CO 8-hr	SO ₂ 24-	NH3 24-hr	Ph 24-hr
(Range)	hr	24-hr	hr	0,0-11	(mg/meter cube)	hr	1110 27-11	1024-11
Good (0-50)	0-50	0-30	0-40	0-50	0-1.0	0-40	0-200	0-0.5
Satisfactory (51-100)	51-100	31-60	41-80	51-100	1.1-2.0	41-80	201-400	0.6-1.0
Moderate (101-200)	101-250	61-90	81-180	101-168	2.1-10	81-380	401-800	1.1-2.0
Poor (201- 300)	251-350	91-120	181-280	169-208	10.1-17	381-800	801-1200	2.1-3.0
Very poor (301-400)	351-430	121-250	281-400	209-748*	17.1-34	801-1600	1201-1800	3.1-3.5
Severe (401- 500)	430+	250+	400+	748+*	34+	1600+	1800+	3.5+

 Table 1: Air pollutants and their various ranges concerning Air Quality Index (AQI) categories

 [Courtesy: Central Pollution Control Board]

To control and mitigate the ill effects of pollution, it is very much necessary to understand quantitatively the contents of different pollutants. A statistical way to analyze the prediction of air quality is by using time series analysis using Auto Regressive Integrated Moving Average (ARIMA) and Vector Autoregression (VAR) models respectively. It was observed that the model generated higher error values compared to the real measured values (Chu & Karr, 2017; Niranjan & Rakesh, 2020). Since the source and origin of pollutants are different, and the interaction with the atmospheric conditions are complex, currently, machine learning methodology is used widely for predicting air quality (Jurado et al., 2022). The multiplier-based enhancement in various machine learning algorithm has been proposed (Juma et al, 2023) research. One of the approaches currently used to predict air quality is the Long and Short-Term (LSTM) with various modifications (Sethi & Mittal, 2020; Camgozlu et al., 2023; Ram et al., 2024) The CNN network-based classification model is recently introduced in various fields. For instance, (Li & Hua, 2020; Shakir & Rakesh, 2018) envisioned a hybrid model of Convolutional Neural Network (CNN) and LSTM called as CNN-LSTM model for predicting the concentrations of PM2.5 values in the air measured on an hourly basis, from US Embassy in Beijing. These time series data were used to forecast PM2.5 using both the traditional LSTM and CNN-LSTM hybrid models. It turned out that the hybrid model took less time and had lower error rates for univariate and multivariate predictions (Setiawan & Setiawan, 2023; Kumar 2018; Kumaran, et al., 2021). Furthermore, (Niranjan & Rakesh 2021; Rakesh & Kumaran, 2021) developed a novel framework in accordance with Deep Learning (DL) method. This method is a combination of Multiple Nested CNN and conventional LSTM models. They have used nested LSTM (NLSTM), which combines additional LSTM units placed on each basal LSTM unit. A multi-modal network was proposed for forecasting mixed AOI data. Discrete stationary wavelet transform was used to convert raw data into multiple sub-signals at different frequencies. Different components of AQI were predicted and it was shown that their proposed model performed better than other methods in terms of operating time, lower error rates like Mean Absolute Error (MAE), Root Mean

Square Error (RMSE), Mean Absolute Percent Error (MAPE), and higher R-squared (R2) values (Jurado et al., 2022; Sharanyaa et al., 2022; Li & Wu, 2020; Arif et al., 2022).

In all the above models, the input time series data is normalized and the main multi-step LSTM is being implemented to predict the required parameters at various time steps. Though normalization makes the data simple to handle, it may trim the essential characteristics of parameters under consideration which may create the instabilities at later time steps. Using Random Forests (RF), a nonparametric statistical learning technique popular in many domains of application, including the analysis of microarrays, is one way to get around these instabilities. Díaz-Uriarte & Alvarez de Andrés (2006), ecology (Rakesh & Kumaran, 2021), pollution prediction (Yan et al., 2021; Kumaran et al., 2021). Dealing with missing data values is another significant difficulty in the handling of air quality data. There are several reasons why data sets contain missing information, including malfunctioning hardware, inadequate sample frequency, deteriorated equipment, and human mistake. (Norazian et al., 2008) To obtain a full set of data, one must choose whether to "impute" a phrase for replacing or reject the missing data. Inferential power may be compromised if important information is lost, hence it is usually not justified to ignore missing data. Impute the missing data is therefore the best course of action. However, unintended bias can also result from the systemic disparities between genuine and replaced data. Determining the best method for predicting missing values is so crucial. It is shown that the Missing at Random (MAR) based RF method gives fairly good accuracy for missing values, especially for air quality monitoring data sets (Alsaber et al., 2021). (Muralidharan, 2020), proposed the Multi Frequency resonator for remote based applications. Here, we have preprocessed and filled the missing values using MAR-based RF for multi time multivariate time series. It is to be noted that the MAR-RF method goes through the interanion till it reaches the stopping criteria which prescribed based on the values of last imputed data. This preprocessed data is a component of the multistage LSTM which gives the multivariate, multi-time output from AQI.

In the next section 2, we will describe, the mathematical representation of time series including with RF-based algorithm for preprocessing the data. We will discuss the stopping criteria for the MAR-RF algorithm which we have followed to confirm the missing data is fairly accurate. In the same section we will describe our multistage LSTM in detail. In section 3 we will discuss in detail on the data which we have used to test our method. We will discuss the results of our prediction of AQI data in 4. Finally in section ?? we will summarize and conclude based on our results which we have generated through our approach and will discuss the future action planned of extending this research.

2 Theory and Methodology

Random Forest (RF) Methodology for Pre-processing

In this section we basically followed the methodology of (Prasad et al., 2006). Let $\mathbf{X} \in \mathbb{R}^{\times p}$ and let Y be observed value of \mathbf{X} , where there are missing values. We will discuss now the methodology by which RF technique will perform for imputing missing observations. First observed dataset is trained by random forest where missing values $\mathbf{X}_{s} \in i_{mis}^{s}$. Therefore we end up with of datasets which are:

- 1. Y^s_{obs} is the set of all observed values of X_s
- 2. Y_{mis}^{s} is the set of all observed missing of X_{s}

which means X_{obs}^s is the dataset, where we will have the observation $i_{obs}^s = \{1, ..., n\} \setminus i_{mis}^s \in X_s$.

We will start with the initial guess for the missing values in X as described in (Jin et al., 2021), depending on the data. Next, the missing values Y^s_{mis} are estimated by applying the trained random forest to X^s_{mis} . This algorithm should be iterated until a stopping criterion is reached. The criteria for

stopping γ_s is reached when, for both variable types, the difference between the current imputed data and the preceding one to prescribed precision. For a set of continuous variables N this is defined as:

$$\delta_{\mathbf{N}} = \frac{\Sigma_{j \in \mathbf{N}} \left(\mathbf{X}_{\text{new}}^{\text{imp}} - \mathbf{X}_{\text{old}}^{\text{imp}} \right)^2}{\Sigma_{j \in \mathbf{N}} \left(\mathbf{X}_{\text{new}}^{\text{imp}} \right)^2}$$
(1)

and that for the set of categorical variables F as:

$$\Delta_{\mathbf{F}} = \frac{\sum_{j \in \mathbf{F}} \sum_{i=1}^{n} \mathbf{I}_{\mathbf{X}_{\text{new}}^{\text{imp}} \neq \mathbf{X}_{\text{old}}^{\text{imp}}}}{N_{A}}$$
(2)

where N_A is the number of missing values in the category **F**.

In a nutshell, we follow the following steps:

Step 1: Set the stopping criterion γ_s .

Step 2: For missing values, make an assumption which is reasonable.

Step 3: Now we need to fit a random forest: such that $\mathbf{Y}_{obs}^{s} \sim \mathbf{X}_{obs}^{s}$ is satisfied.

Step 4: Predict Y_{miss}^s using X_{miss}^s .

Step 5: Update the imputed values using the predicted Y_{miss}^{s} .

Step:6: Update γ_s and X^{imp} .

The performance is assessed by using the normalized root mean squared error (NRMSE).

$$NRMSE = \sqrt{\frac{\left(\left(\mathbf{X}^{\text{true}} - \mathbf{X}^{\text{imp}}\right)^{2}\right)}{\text{var}(\mathbf{X}^{\text{true}})}}$$
(3)

All the missing AQI data is preprocessed by the above RF-based algorithm and it is represented by X_t This preprocessed data goes into our multistage LSTM algorithm, which we will discuss in the next sub-section.

LSTM Methodology

One of the favored approaches currently used in solving complex RNN problems is LSTM, which was originally proposed (Hochreiter & Schmidhuber, 1997) to mitigate vanishing gradient effects. In a nutshell, the recurrently connected subnets make up the LSTM architecture, which is known as a memory block. If we take the memory block of LSTM with one cell it has three gates: input, output, and forget. A schematic structure of a cell is shown in Figure 1. Let X_t represent processed data set at time t. In an LSTM network, the U_f in the forget gate equation (Stekhoven & Bühlmann, 2012) represents the weight matrix associated with the input connections to the forget gate. It indicates the amount of each element to retain. It plays a crucial role in mitigating the vanishing gradient problem and allowing LSTMs to capture long-term dependencies in sequences. The LSTM steps are as follows:

The Forget Gate \mathbf{f}_t is expressed as:

$$\mathbf{f}_{\mathbf{t}} = \sigma(\mathbf{W}_{\mathbf{f}}\mathbf{X}_{\mathbf{t}} + \mathbf{U}_{\mathbf{f}}\mathbf{h}_{\mathbf{t-1}} + \mathbf{b}_{\mathbf{f}})$$
(4)

The Input Gate I_t is expressed as:

$$\mathbf{I}_{t} = \sigma(\mathbf{W}_{i}\mathbf{X}_{t} + \mathbf{U}_{i}\mathbf{h}_{t-1} + \mathbf{b}_{i})$$
(5)



Figure 1: A schematic structure of a cell in LSTM

Output Gate **O**_t is expressed as:

$$\mathbf{O}_{\mathbf{t}} = \sigma(\mathbf{W}_{\mathbf{o}}\mathbf{X}_{\mathbf{t}} + \mathbf{U}_{\mathbf{o}}\mathbf{h}_{\mathbf{t}-1} + \mathbf{b}_{\mathbf{i}}) \tag{6}$$

Cell input activation vector:

$$\mathbf{C}_{\mathbf{t}} = \mathbf{ReLU}(\mathbf{W}_{c}\mathbf{X}_{\mathbf{t}} + \mathbf{U}_{c}\mathbf{h}_{\mathbf{t-1}} + \mathbf{b}_{\mathbf{c}}) \qquad (7)$$

Cell state output:

$$\mathbf{C}_{\mathbf{t}} = \mathbf{f}_{\mathbf{t}} \otimes \mathbf{C}_{\mathbf{t}-1} + \mathbf{i}_{\mathbf{t}} \otimes \mathbf{C}_{\mathbf{t}} \tag{8}$$

Hidden state Vector:

$$\mathbf{h}_{\mathbf{t}} = \mathbf{O}_{\mathbf{t}} \otimes \mathbf{ReLU}(\mathbf{C}_{\mathbf{t}}) \tag{9}$$

where **b** is the bias and σ is the sigmoid function. Please note we have **ReLU** instead of tanh function which is used commonly in many LSTM models. Figure 2 shows the step of our algorithm. In the next section 3, we will elaborate on our data sets.



Figure 2: Flow Chart depicting the flow of tasks performed in making the prediction model

3 The Dataset

The dataset used in this research contains air quality and pollutant level data from Karnataka State Pollution Control Board, Bengaluru, collected in 6 regions of Bengaluru (Hebbal, Silk Board, NIMHANS, Kavika, Jaya Nagar, Majestic City Railway Station, and SG Halli Bangalore), over the last five years. The data collected from the first five regions contained 39000 records of hourly reports each. It consists of 14 fields. Table 2 gives the details of the various air pollutants (features) under study. Imputation of missing values was done at three stages based on the type of pollutants viz., gaseous, chemical, dust since the type of values for each feature varies integer or float, etc., the forecasting was carried out using LSTM-based Autoencoder. Our model configuration is shown in the Figure 1 (Shakir & Rakesh, 2018).

Air quality data was gathered from 7 locations in Bengaluru City. Table 1 lists the pollutant level and the environmental parameter information collected during the process. Information collected from the Hebbal, Jaya Nagar, Kavika, NIMHANS, and Silk Board areas consists of hourly data for the last five years, while that gathered from City Railway Station and S.G Halli areas consists of the daily report for the last five years. The Day of Week (DOW) component derives the 'day of the week', 'hour of the day', 'date', and 'time' details from the time stamp field and adds them to the dataset.

Sl. No.	Feature Name	Туре	Unit
1		СО	mgm/m3
2		O ₃	µgm/m3
3		NO	µgm/m3
4		NO ₂	µgm/m3
5	Gaseous Pollutants	NOx	µgm/m3
6		NH ₃	µgm/m3
7		SO ₂	µgm/m3
8		PM2.5	µgm/m3
9	Dust Pollutants	PM 10	µgm/m3
10		Benzene	µgm/m3
11		Toluene	µgm/m3
12		m,p-Xylene	µgm/m3
13	Acidic Pollutants	o-Xylene	µgm/m3
14		Ethyl Benzene	µgm/m3

 Table 2: Various features under study that are classified into three categories - gaseous, dust, and chemical pollutants

The enhanced dataset from each area is then passed to the next stage, where missing temperature values are filled using values from the CO field, using the arithmetic mean. The next step grouped the DOW, hour, time stamp, ambient temperature, wind speed, and wind direction fields from Hebbal, Jaya Nagar, Kavika, NIMHANS, and Silk Board areas to address the missing values in wind speed and wind direction fields. It used the arithmetic mean to fill in the missing values. The updated values were moved back to the original datasets. The next stage addressed the relative humidity, and the barometric pressure in the individual datasets using the arithmetic mean. The next stage used the Random Forest regression to address the missing values in the pollutant's fields. The main algorithm involves the use of an LSTM-based Auto encoder time series algorithm to forecast the values of pollutants. The algorithm used

a input sequences of 10 time steps of input variables as input windows, predicted a sequence of 5 time steps of output(10X5). This research evaluated two variations of autoencoders, with single LSTM and dual LSTM layers in the encoder and decoder stages. The proposed research work evaluated two versions of LSTM based Auto encoders the first having just one concealed layer in the encoder and decoder, and the second with double hidden layers. The model was trained using different parameters as shown in Table 3.

Table 3: Listing the various parameters and their corresponding values used during model creation

Sl. No.	Parameter	Value
1	Number of neurons	100
2	Number of epochs	50
3	batch size	32
4	Window size	10X5

4 Result and Discussion

Based on the discussion above in Section 3, we have simulated the following 5 different configurations:

Configuration 1 LSTM-based Autoencoders – Single hidden layer

Configuration 2 LSTM-based Autoencoders – Single hidden layer

Configuration 3 LSTM-based Autoencoders – Double hidden layers

Configuration 4 LSTM-based Autoencoders – Single hidden layer (only 2 features AT & CO)

Configuration 5 LSTM-based Autoencoders – Single hidden layer

To make our report concise, we are presenting the result only for CO-polluted concentration. It is implied that we can extend this methodology to predict other pollutants also without any difficulty.

The efficiency metrics of our simulated results produced are evaluated with respect to three parameters namely, precision, recall, and F1-score values.

- **1. Precision:** Is a classification model's capacity to recognize only pertinent data points. The number of true positives divided by the total number of true positives + the number of false positives.
- 2. **Recall:** Is a metric that represents the proportion of positive cases, out of all the positive cases in the data, that the classifier correctly predicted. Another name for it is sensitivity at times. It is the proportion of accurate predictions to all data sets that fall into that class.
- **3. F1–score:** Values is The harmonic mean of recall and precision is called F1-score. It provides a balanced score for precision and recall. The F1 will be high only when both precision and recall are high.

Table 4 sums up the results of our study for four different configurations including **loss function**. To optimize the present state, we need to estimate the error of the model which is by **loss function**

	,					U
Configuration	Number	Window	Precision	Recall	F-	Loss value
Setup No.	of	Size	obtained	value	Score	
_	neurons				value	
1	100	10X5	1.0	0.97	0.98	5.92
2	150	10X5	1.0	0.98	0.99	4.17
3	100	10X5	1.0	0.95	0.97	6.99
4	100	10X5	1.0	0.79	0.88	0.27
5	100	5X3	1.0	0.95	0.97	4.15

Table 4: Accuracy Precision, Recall, F-score and Label Value for various con- figuration

From the above analysis, it is inferred that configuration set up 4 which having a single-layered LSTM-based Autoencoder with 100 neurons and window size of 10X5 and with just 2 features (CO and AT) has lower Recall and F1-Score values. Unlike other setups, the loss incurred is less since the dimensions of the dataset were reduced to just two features. The performance configuration 2, with single layered LSTM-based Auto Encoder having 150 neurons; a window size of 10X5 performed better compared to other models with the highest Recall and F-Score values and comparatively small loss value. The model configurations discussed above used the Adam optimizer, Huber loss function, and ReLu activation in the LSTM layers. Table 4, 5, and 6 summarizes the results of our simulation. It is seen that the LSTM-based Auto Encoder model with various configuration setups has predicted the values near the actual outputs.

We have also found that the loss value reduces and remains constant after 30 epochs for all the proposed models under various configurations as shown in Figures 3, and 4.

Туре	Date/	Atmospheric	Relative	Barometric	CO
10X5 window	Time (Index)	Temperature	Humidity	Pressure	
size					
	2023-4-17 09:45:00	25.1	71	47	0.6
	2023-4-17 10:00:00	27.9	71.2	47	0.55
	2023-4-17 10:15:00	27.9	71	43	0.59
	2023-4-17 10:30:00	27.9	71	41	0.51
	2023-4-17 10:45:00	27.9	71	38	0.51
Actual Input	2023-4-17 11:00:00	32	71	36	0.49
	2023-4-17 11:15:00	32	71	34	0.39
	2023-4-17 11:30:00	32	71	30	0.33
	2023-4-17 11:45:00	32	71	28	0.42
	2023-4-17 12:00:00	34.9	71	26	0.51
	2023-4-17 12:15:00	34.9	71	26	0.47
Actual Output	2023-4-17 12:30:00	34.9	71	27	0.36
	2023-4-17 12:45:00	34.9	71	28	0.53
	2023-4-17 01:00:00	36.3	71	28	0.58
	2023-4-17 01:15:00	36.3	71	28	0.6

Table 5: Actual values as per the dataset

Type 10x5	Date/	Atmospheric	Relative	Barometric	СО
Window Size	Time (Index)	Temperature	Humidity	Pressure	
	2023-4-17	35.947	68.773	26.78	0.4841
	12:15:00				
	2023-4-17	33.853	73.02	26.19	0.3708
Predicted Output for	12:30:00				
Configuration 1	2023-4-17	33.853	73.03	28.84	0.5141
Multivariate eldi	12:45:00				
100 neurons algorithm	2023-4-17	35.211	68.8	27.16	0.5974
	13:00:00				
	2023-4-17	37.389	73.24	28.84	0.618
	13:15:00				
	2023-4-17	35.598	69.5	26.52	0.4794
	12:15:00				
	2023-4-17	35.598	72.3	26.46	0.3672
Predicted Output for	12:30:00			20110	010072
Configuration 2	2023-4-17	34,202	72.3	28.56	0.5194
Multivariate e1d1	12:45:00	0 11202	/ =10	20100	0.0191
150 neurons algorithm	2023-4-17	37.026	69.5	27.44	0.5916
	13:00:00	071020	0,10	2/11	0.0910
	2023-4-17	35.574	69.3	28.56	0.588
	13:15:00	55.574	07.5	20.50	0.500
	2023-4-17	33 504	73 7	27.04	0.4512
	12:15:00	55.504	15.1	27.04	0.4012
	2023-4-17	36.296	68.54	28.08	0.3744
Predicted Output for	12:30:00	2002220	00.21	20.00	0.0711
Configuration 3	2023-4-17	36.296	68.1	26.88	0.5088
Multivariate e2d2	12:45:00	0.0220	0011	2000	0.0000
100 neurons algorithm	2023-4-17	34.848	73.7	26.88	0.6032
	13:00:00			2000	010002
	2023-4-17	37.752	73.6	29.12	0.576
	13:15:00	011102	7010		0.070
	2023-4-17	39,0531	-	_	0.41407
	12:15:00	0310001			0.11107
	2023-4-17	39.0531	-	-	0.40284
Predicted Output for	12:30:00	0,0001			0110201
Configuration 4	2023-4-17	30.7469	-	-	0.59307
Multivariate eld	12:45:00				
only 2 features - AT and	2023-4-17	40.6197	-	-	0.51098
CO	13:00:00				
	2023-4-17	31.9803	-	-	0.5286
	13:15:00				
	2023-4-17	36.296	73.84	27.04	0.4888
	12:15:00				
	2023-4-17	35.947	73.13	27.81	0.3708
Predicted Output for	12:30:00				
Configuration-5	2023-4-17	33.504	68.16	26.88	0.5088
Multivariate eldl- 5X3 ws algorithm	12:45:00				
origination and origination	2023-4-17	35.211	68.87	27.16	0.5626
	13:00:00				
	2023-4-17	37.026	72.42	28.56	0.612
	13:15:00				

Table 6 : Predicted values using various configuration models

Table 5 presents the actual input and output values obtained from observations and helps in predicting values generated by the LSTM-based Auto-Encoder model with different configuration setups. This table includes specific parameters related to atmospheric conditions and pollutant concentrations, such as atmospheric temperature, relative humidity, barometric pressure, and CO concentration. The structured format likely includes columns representing the date and time of observations, atmospheric temperature, relative humidity, barometric pressure, and CO concentration, with the actual values recorded for each parameter. By comparing the actual values with the predicted values, we can assess the model's performance in accurately forecasting pollutant levels based on the input data. This comparison is valuable for evaluating the model's accuracy, reliability, and effectiveness in predicting environmental parameters essential for air quality monitoring and pollution control efforts. Therefore, Table 5 serves as a critical tool for validating the model's predictive capabilities and enhancing its applicability in environmental research and decision-making processes.

Table 6 presents the efficiency metrics of the simulated results produced by the LSTM-based Auto-Encoder models with different configuration setups. Specifically, the table showcases the precision, recall, and F1-score values for each model configuration, serving as indicators of the model's accuracy in predicting pollutant levels. The results demonstrate that the LSTM-based Auto-Encoder model, across various configuration setups, has successfully predicted values that closely align with the actual outputs.

In summary, these tables offer a comprehensive overview of the simulation results, highlighting the efficacy of the LSTM-based Auto-Encoder models in forecasting pollutant levels in urban areas.



(a) LSTM Auto-Encoder with single layer



(b) LSTM-based Auto-Encoder with double layers

An Approach towards Forecasting Time Series Air Pollution Data Using LSTM-based Auto-Encoders



(c) LSTM-based Auto-Encoder with a single layer having 100 neurons in the hidden layer



(d) LSTM-based Auto-Encoder with a single layer having 100 neurons in the hidden layer having only two features

Figure 3: Demonstration of the loss values vs. the number of epochs for configurations 1-4



Figure 4: LSTM-based Auto-Encoder with a single layer having 100 neurons in the hidden layer for configuration 5

Figure 3 above likely depicts the correlation between the loss values and the number of epochs for configurations 1 to 4 of the LSTM-based Auto-Encoder algorithm. The horizontal axis of the graph represents the progression of training epochs, which are iterations over the dataset during the model training phase. On the vertical axis, the loss values are plotted, indicating the model's performance at each epoch. Lower loss values signify better model performance and accuracy.

Each configuration (1 to 4) represented by distinct lines or curves on the graph, illustrates how the loss values evolve with increasing epochs. This visualization aids in assessing the training dynamics of the models. Ideally, a downward trend in loss values across epochs suggests that the model is learning and enhancing its predictive capabilities.

By analysing Figure 3, we can compare the training patterns of different configurations and identify which configuration -4 setup leads to the most significant reduction in loss values throughout the training process, since it has only two features.

Figure 4 illustrates the performance of the LSTM-based Auto-Encoder algorithm with a single hidden layer containing 100 neurons and window size 5X3 for configuration 5. The graph shows the relationship between the loss values and the number of epochs during the training phase.

By analyzing Figure 4, we can evaluate the effectiveness of the LSTM-based Auto-Encoder algorithm with a single hidden layer containing 100 neurons 5X3 window size for configuration 5 in reducing loss values over the training epochs and it is almost equal to configuration 2 set with single layer and 150 neurons as depicted in table-4.

From fig. 3 and fig. 4, it is clear that the loss value reduces and remains constant after 30 epochs for all the proposed models under various configurations.

5 Summary and Conclusion

This research work compared the performance of an LSTM-based Auto encoder algorithm under different model configurations to forecast the pollutant levels based on time series data. The data was given by the Karnataka State Pollution Control Board (KSPCB), Bengaluru. Imputation of missing values was carried out with a novel methodology by considering the practical correlation between pollutants and atmospheric features. Arithmetic mean and Random Forest regression were used to fill in the missing values. Moreover, imputation of missing values was done at three stages based on the type of pollutants viz., gaseous, chemical, dust since the type of values for each feature varies integer or float, etc., the forecasting was carried out using LSTM-based Autoencoder. The results of various models built using LSTM-based Auto Encoders have been analyzed. Future research will evaluate the impact of pollution in one region on the neighboring region.

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46