



Article

Weighted Averages and Polynomial Interpolation for PM2.5 Time Series Forecasting

Anibal Flores ^{1,*}, Hugo Tito-Chura ¹, Victor Yana-Mamani ¹, Charles Rosado-Chavez ¹ and Alejandro Ecos-Espino ²

- Departamento Académico de Ingeniería de Sistemas e Informática, Universidad Nacional de Moquegua, Urb. Ciudad Jardin-Pacocha-Ilo, Moquegua 18611, Peru
- Departamento de Ciencias Básicas, Universidad Nacional de Moquegua, Prolongación Calle Ancash S/N, Moquegua 18001, Peru
- * Correspondence: afloresg@unam.edu.pe

Abstract: This article describes a novel method for the multi-step forecasting of PM2.5 time series based on weighted averages and polynomial interpolation. Multi-step prediction models enable decision makers to build an understanding of longer future terms than the one-step-ahead prediction models, allowing for more timely decision-making. As the cases for this study, hourly data from three environmental monitoring stations from Ilo City in Southern Peru were selected. The results show average RMSEs of between 1.60 and 9.40 ug/m³ and average MAPEs of between 17.69% and 28.91%. Comparing the results with those derived using the presently implemented benchmark models (such as LSTM, BiLSTM, GRU, BiGRU, and LSTM-ATT) in different prediction horizons, in the majority of environmental monitoring stations, the proposed model outperformed them by between 2.40% and 17.49% in terms of the average MAPE derived. It is concluded that the proposed model constitutes a good alternative for multi-step PM2.5 time series forecasting, presenting similar and superior results to the benchmark models. Aside from the good results, one of the main advantages of the proposed model is that it requires fewer data in comparison with the benchmark models.

Keywords: PM2.5; multi-step forecasting; weighted averages; polynomial interpolation; deep learning



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1. Introduction

The global burden of disease associated with the exposure to air pollution causes millions of deaths and years of healthy life lost annually [1]. The burden of disease attributable to air pollution is estimated to be on par with other major global health risks, such as unhealthy diet and smoking [2,3]; air pollution is now recognized as the greatest environmental threat [4] to human health.

PM2.5 is one of the main pollutants [5–7], and it can accumulate in the respiratory system. It is associated with numerous negative effects on the health, such as increased respiratory diseases and decreased lung function. Thus, PM2.5 can cause premature death in people with heart [8] or lung disease [9,10] and can lead to non-fatal myocardial infarctions [11], irregular heartbeats [12], aggravated asthma [13,14], reduced lung function, and increased respiratory symptoms, such as irritation in the airways [15], cough, and difficulty breathing.

In this context, it is clear that there is a need to implement air quality monitoring stations and, from these, develop tools that allow for the analysis and prediction of the volumes of PM2.5 and other pollutants. According to the literature, numerous works have been found that address the topic of PM2.5 forecasting; however, most studies have addressed PM2.5 forecasting through the implementation of one-step-ahead forecasting. This means that the models predict only one value, which is not very useful if the prediction of more than one value is required; thus, in this context, multi-step forecasting models become important.

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The proposed model is a novel and simple multi-step model for PM2.5 forecasting. It is inspired by weighted moving averages [16,17] and polynomial interpolation [18]; it uses data that are closer to the predicted values because these are more relevant [19], assuming that the near future is close to the near past. In this sense, the proposed model uses data from just two days (48 h) to predict 24, 48, and 72 h in the future; thus, it requires a small amount of data to produce similar and superior results to those made by the state-of-the-art techniques such as deep learning-based approaches, including long short-term memory (LSTM), bidirectional LSTM (BiLSTM), gated recurrent unit (GRU), bidirectional GRU (BiGRU), and LSTM with attention mechanism (LSTM-ATT).

The main contributions of this work are as follows:

- A novel model for multi-step forecasting of PM2.5 time series;
- A comparative analysis between proposal results and benchmark models based on deep learning;
- A web application for PM2.5 time series forecasting.

Some of the limitations of this work lie in the type of forecasting model that is proposed here (a univariate model) and in the number of case studies which is limited to three monitoring stations located in Southern Peru.

2. Literature Review

2.1. Overview of Forecasting Models

Forecasting models for time series have evolved from statistical models to machine learning and deep learning models. Among the statistical models, one of the best-known models is ARIMA [20]. Machine learning-based models include the support vector machine (SVM) [21], nearest neighbors regression [22], and ensemble models such as random forests [23], gradient boosting [24], and Adaboost [25]. Deep learning models such as those based on recurrent neural networks include LSTM, GRU, BiLSTM, BiGRU, and RNNs with attention layers; finally, those based on Transformers [26,27] are gaining popularity.

2.2. One-Step-Related Works in PM2.5 Forecasting

Most of the related works have proposed and implemented LSTMs; among these, the research presented in [18,28–33] has included decomposition techniques, and others have combined LSTM with other techniques. In [30], the authors added a decomposition technique named Singular Spectrum Analysis (SSA). In [28,29], bidirectional LSTM (BiL-STM) and LSTM were used, respectively. In [31], LSTM was combined with convolutional neural networks (CNNs). In [33], a bidirectional LSTM model was combined with a CNN model. In [32], an LSTM was used with decomposition techniques such as CEEMDAN and FCM. Additionally, in [34,35], GRU was implemented with data augmentation and with Q-Learning, respectively.

Support vector regression (SVR) and random forest have been used in works such as [36–38]. The first uses SVR in a hybrid way with quantum particle swarm optimization (QPSO), while the second adds a decomposition technique named hybrid modified variational mode decomposition, and the last one uses a random forest model alone.

Furthermore, some related works, including [39–43], implemented different approaches to those cited before. In [39], a Hammerstein recurrent neural network was proposed; in [24], a decomposition ensemble learning model was used, based on the variation-mode decomposition approach and a whale-optimization algorithm (IWOA); in [42], an attention-based deep neural network was used; in [43], a multivariate deep belief network using PM2.5 and temperature data was implemented; in [40], a multiple model adaptive unscented Kalman filter was utilized.

2.3. Multi-Step-Model-Related Works in PM2.5 Forecasting

In [44,45], the authors proposed an extreme learning machine (ELM), which was used to obtain MAPEs between 5.11% and 37.23% for three-step forecasting and MAPEs between 5.12% and 22.2% for seven-step forecasting.

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In [46], a hybrid CNN-LSTM model was used for a ten-step forecasting, obtaining MAPEs between 22.09% and 25.94%.

In [47], AdaBoost was proposed, obtaining MAPEs between 15.32% and 23.75% for three-step forecasting.

In [48], a support vector machine (SVM) was used for four-step predictions, obtaining RMSEs between 8.16 and 39.04 ug/m³.

In [49], a self-organizing memory neural network was implemented for 1-, 4-, 8-, and 12-day forecasting, obtaining MAPEs between 6.66% and 14.08%.

In [50], a feature-selection approach based on a genetic algorithm was used to implement an LSTM model for one- to six-step forecasting, obtaining MAEs between 3.592 and 6.684.

In [51], a BiLSTM was used for three-step forecasting, obtaining MAPEs between 27.33% and 40.73%.

In [52,53], a CNN was proposed for ten-step predictions and four-step predictions, obtaining MAPEs between 28.51% and 33.29%.

Finally, in [54], a point system for three-step forecasting was implemented, obtaining MAPEs between 7.53% and 16.18%.

According to the literature review for single and multi-step PM2.5 forecasting, most researchers have focused on machine learning and deep learning techniques; some researchers considered hybrid models, decomposition techniques, and data augmentation. Most researchers addressed the problem of PM2.5 forecasting in a single step. According to [55], decisions about air quality often need to be made more than 10 h in advance; according to the literature, the maximum number of steps used for PM2.5 forecasting was 10 steps ahead (10 h). This finding served as a motivation for this work, prompting the authors to propose a different multi-step forecasting proposal with a greater number of steps than those reported in the literature. It is important to highlight that, according to the literature, the term "steps" refers to the number of forecasted points.

The main differences between the related literature and the present study are summarized in Table 1.

Table 1. Differences between related research and the present study.

Rel	Related Research		This Work	
-	Most existing studies are based on machine learning and deep learning.	-	The present study is based on weighted averages and polynomial interpolation.	
-	Most existing studies were implemented for one-step forecasting.	-	The present study was implemented for multi-step forecasting.	
-	The models proposed in existing studies require a lot of training data.	-	The model proposed in the present paper requires very few training data.	
-	The models proposed in existing studies are based on different and more complex processes.	-	The model proposed in the present paper is based on the simple correlation of days.	

3. Materials and Methods

3.1. Data Collection

The hourly data used for this work were downloaded from OEFA's server (located at https://pifa.oefa.gob.pe/VigilanciaAmbiental/, accessed on 2 May 2023), and they span the period between 1 August 2020 and 30 April 2023. They correspond to three environmental monitoring stations (Pacocha, Bolognesi, and Pardo) located at Ilo city, Moquegua region, Southern Peru.

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The downloaded data had several missing values, so the days that presented missing values were discarded. The available records for each station remaining after this process are shown in Table 2. Likewise, since the present study addresses a regression problem, according to the literature for statistical models, the dataset was split into the following two sets: the training set (80%) and the test (20%) set.

Table 2. Data of environmental monitoring stations.

Station	Total Hours	Train 80%	Test 20%	
Pardo	21,960	17,568	4392	
Bolognesi	18,120	14,496	3624	
Pacocha	16,344	13,080	3265	

3.2. Selection of Days

The proposed model does not work with all the training data; it only works with the two historical days. For this selection, first, the correlations of the hours between the last 24 days of the training data of the three stations were analyzed separately. For this, the data were organized in a matrix manner, and the correlation matrix that is commonly used for feature selection [56] in classification tasks using machine learning and deep learning models was utilized. The respective correlation matrices for each monitoring station are shown in Figures 1–3.

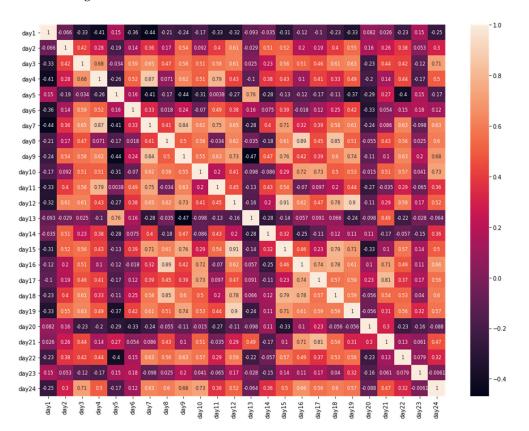


Figure 1. The 20-day correlation of Pacocha station.

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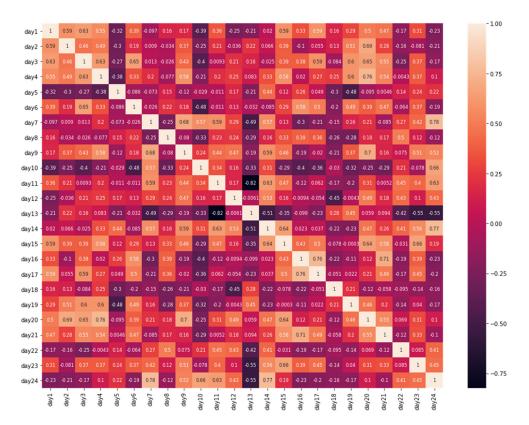


Figure 2. The 20-day correlation of Bolognesi station.

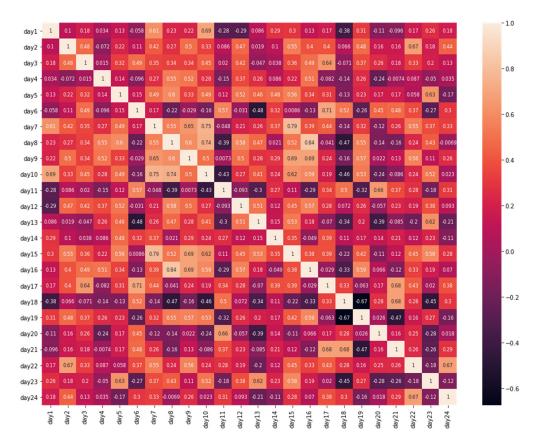


Figure 3. The 20-day correlation of Pardo station.

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In this study, the correlation between a given day and the subsequent days is of interest; this correlation can be determined by looking at the values above the main diagonal of the respective correlation matrix, C. If the day to be analyzed is located on the main diagonal in row i, then the prior days to i will begin in row i-1, column j=i; thus, the average correlation is given by Equation (1). For the following previous day, it also starts in row i-1 but in column j=i+1 (2). The following previous day starts in row i-1, column j=i+2 (3), and so on.

$$\overline{C}_1 = \frac{\sum_{i=1}^{n-1} C_{i-1,i}}{n-1} \tag{1}$$

$$\overline{C}_2 = \frac{\sum_{i=1}^{n-2} C_{i-1,i+1}}{n-2} \tag{2}$$

$$\overline{C}_3 = \frac{\sum_{i=1}^{n-3} C_{i-1,i+2}}{n-3} \tag{3}$$

Equations (1)–(3) can be summarized in (4) as follows:

$$\overline{C_j} = \frac{\sum_{i=1}^{n-j} C_{i-1,i-1+j}}{n-j} \tag{4}$$

where $\overline{C_j}$ is the average correlation of day j. Using Equation (4), Table 3 was formulated.

Table 3. Average correlations between days.

Day	Pacocha	Bolognesi	Pardo	Avg
1	0.2637	0.1952	0.1450	0.2013 ± 0.06
2	0.1903	0.1301	0.1378	0.1527 ± 0.03
3	0.3079	0.1488	0.2632	0.2400 ± 0.08
4	0.2539	0.0417	0.2411	0.1789 ± 0.12
5	0.2413	0.1026	0.1729	0.1723 ± 0.07
6	0.2404	0.0639	0.2595	0.1879 ± 0.11
7	0.3393	0.0650	0.2604	0.2216 ± 0.14
8	0.3589	0.0984	0.2152	0.2242 ± 0.13
9	0.1939	0.0779	0.2572	0.1763 ± 0.09
10	0.2409	0.1284	0.1451	0.1715 ± 0.06
11	0.1685	0.0992	0.1410	0.1362 ± 0.03
12	0.2378	0.1378	0.1965	0.1907 ± 0.05
13	0.2683	0.2056	0.1785	0.2175 ± 0.05
14	0.1554	0.2822	0.2514	0.2297 ± 0.07
15	0.2515	0.2429	0.2888	0.2611 ± 0.02
16	0.2057	0.2890	0.1607	0.2185 ± 0.07
17	0.0924	0.4475	0.0662	0.2020 ± 0.21
18	0.1612	0.2635	0.2781	0.2343 ± 0.06
19	0.0845	0.2231	0.0313	0.1130 ± 0.10
20	0.1950	0.1938	0.2017	0.1968 ± 0.00
21	0.1806	-0.1380	0.1579	0.0668 ± 0.18
22	0.2255	0.0505	0.3517	0.2092 ± 0.15
23	-0.2536	-0.2328	0.1823	-0.1014 ± 0.25

According to Table 3, on day i, there is an average correlation of 0.2013 with the previous day, i - 1, an average correlation of 0.1527 with day i - 2, an average correlation of 0.2400 with day i - 3, and so on.

It is important to highlight that, according to Figure 4, the different numbers of days were used for estimating average correlations. Thus, for day i - 1, there are n - 1 days,

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for day i - 2, there are n - 2 days, and so on; for the last day, or the one furthest from the main diagonal, there is only one day.

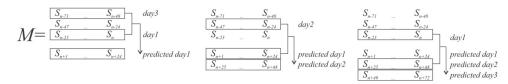


Figure 4. How matrix *M* is used to make predictions.

Just two days were chosen for the implementation of the proposed model. A greater number of days can be chosen, but the results do not improve significantly. According to Table 1, for the implementation of the proposed model, days 1 and 3 were chosen. Day 1 was chosen because—despite not having the highest average correlation—it showed a good correlation with the highest number of days (0.2013), with little deviation. It can be seen in all stations that there are correlations greater than 0.14. Similarly, day 3 was selected because it has an average correlation of 0.2400, which is higher than those of day 1, day 2, day 4, day 5, day 6, and others. Days 7 and 8 present average correlations higher than 0.22 and are higher than day 1, but they were discarded due to their deviation. For the Pacocha and Pardo stations, their correlations are higher than 0.30 and 0.20, respectively; for the Bolognesi station, the correlations are less than 0.10.

The result of this stage is a matrix, M, of the order 3×4 , with the data of the last 72 h, according to Equation (5). The first row corresponds to day 3, the second to day 2, and the third to day 1. New rows (4th, 5th, 6th, . . .) are added according to the performed predictions.

$$M = \begin{bmatrix} S_{n-71} & \dots & S_{n-48} \\ S_{n-47} & \dots & S_{n-24} \\ S_{n-23} & \dots & S_n \end{bmatrix}$$
 (5)

Here, S is the training data vector of an environmental monitoring station and n is the vector length.

The first row contains the 24 h of day 3, the second contains the 24 h of day 2, and the third contains the 24 h of day 1.

According to Figure 4, in the first step, *M* was used to obtain the predicted day 1 (24 h ahead) using day 1 and day 3. Then, the predicted day 2 was obtained using day 2 and the predicted day 1. Finally, the predicted day 3 was obtained using day 1 and the predicted day 2. More details about the estimation of predicted days are presented in the next section.

3.3. Estimating Weighted Averages (WAs)

The hours of the two selected days in the previous stage were averaged; however, this was not performed in the traditional way, considering the same weight for each day. Instead, the values from Table 2 were used to create the matrix *Corrs* (6). The weights were important since the days did not have the same correlation.

$$Corrs = \begin{bmatrix} 0.2637 & 0.3079 \\ 0.1952 & 0.1488 \\ 0.1450 & 0.2632 \end{bmatrix}$$
 (6)

The first row of the matrix represents the weights of days 1 and 3 for the Pacocha station, the second represents the weights of days 1 and 3 for the Bolognesi station, and the third represents the weights of days 1 and 3 for the Pardo station.

From (5), the weights for each station were obtained according to (7) as follows:

$$w_j = Corrs_{i,j} / (Corrs_{i,1} + Corrs_{i,2})$$
 (7)

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where i is the environmental monitoring station. According to (5), j is the day, 1 is the first day, and 2 is the third day. w_1 is the weight for the first day and w_3 is the weight for the third day.

According to Figure 4, in the predicted day 1, S_{n-71} and S_{n-23} were used for the first hour, S_{n-70} and S_{n-22} were used for the second hour, S_{n-69} and S_{n-21} were used for the third hour, and so on.

Then, the weighted average (WA) Equations (8)–(10) for different predicted days were implemented.

$$E_1 = M_{1,i} * w_1 + M_{3,i} * w_3 \tag{8}$$

$$E_2 = M_{2,j} * w_1 + M_{4,j} * w_3 (9)$$

$$E_3 = M_{3,j} * w_1 + M_{5,j} * w_3 (10)$$

 E_1 , E_2 , and E_3 are the vectors of predicted days 1, 2, and 3, respectively. Equations (8)–(10) can be summarized in (11) as follows:

$$E_i = M_{i,j} * w_1 + M_{i+2,j} * w_3 (11)$$

where i is the predicted day, j is the j-th hour of the day, w_1 is the weight for day 1, and w_3 is the weight for day 3.

An example of how WA (11) predicts 24 h for Pardo station can be seen in Figure 5.

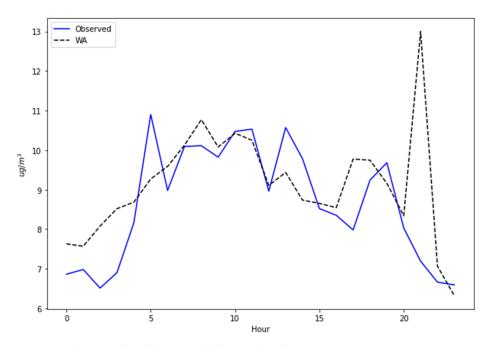


Figure 5. The 24 predicted hours with the weighted average equation.

3.4. Applying Polynomial Interpolation (PI)

To reduce the error of the weighted average (WA) predictions, in this stage, polynomial interpolation (PI) was used to smooth the curve of the predicted values, bringing them closer to the observed ones.

For this, from the 24 predicted points in each iteration of WA, a representative number was selected; in this case, the optimal number was seven, being the seven positions, 0, 3, 7, 11, 15, 19, and 23. Then, NumPy's polyfit function was used with the optimal degree of five. Finally, the obtained function was used to predict the 24 values. The corresponding polynomial function is expressed in Equation (12) as follows:

$$pi = a_0 + a_1 x + a_2 x^2 + a_3 x^3 + a_4 x^4 + a_5 x^5$$
 (12)

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where a_0 , a_1 , a_2 , a_3 , a_4 , and a_5 are the coefficients of the polynomial function.

Figure 6 shows a comparison between the observed values, the WA, and the WA + PI predicted values for 24 h for the Pacocha station.

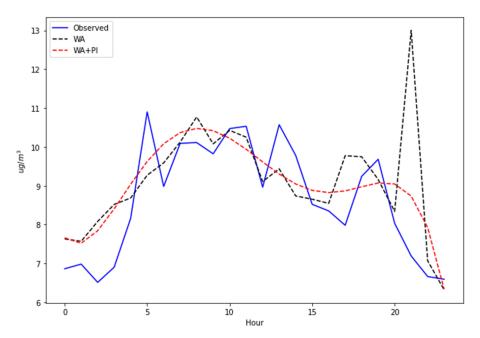


Figure 6. Multi-step predictions of WA and WA + PI.

The corresponding algorithm for the WA + PI proposal is shown in Figure 7.

```
procedure WA PI(RW,w1,w2)
     begin
         n=length(RW[0])
 4
         preds=[]
 5
         days=3
 6
         i=0
 7
         points=[0,3,7,11,15,19,23]
 8
         x=points
 9
         while i<days
         begin
11
             res=[]
             for j = 0 \rightarrow n
13
             begin
14
                  p=RW[i][j]*w1+RW[i+2][j]*w2
15
                  res.append(p)
16
             end
17
             y=[]
18
             for k = 0 -> length (points)
19
                  y.append(res[points[k]])
             poly = numpy.poly1d(numpy.polyfit(x, y, 5))
21
             res1=[]
22
             for k = 0 \rightarrow n
                  res1.append(poly(k))
24
             preds.append(res1)
25
             RW.append(res1)
26
             i+=1
27
         end
    end
28
```

Figure 7. The WA + PI algorithm.

According to Figure 7, the procedure receives RW, w_1 , and w_2 as parameters. RW is a matrix with the data from the last three days of the training data. The algorithm was set to predict 3 days (72 h); for this, an iterative process (while) was performed, going from 0 to 2, predicting 24 h in each iteration. Between lines 12 and 16, the WA predictions were made. Between lines 17 and 20, from the WA predictions, the polynomial function was obtained using the array points. Between lines 21 and 25, the polynomial function was used to make predictions, and the RW matrix was updated. The process continued, with the next day being used to predict the following 24 h. The loop ended when 72 h had been predicted.

The corresponding obtained results are described in the Results and Discussion section.

3.5. Evaluation

The evaluation of the results was carried out through the root mean squared error (RMSE), which allows one to measure the error in terms of ug/m^3 , mean absolute percentage error (MAPE), which allows one to measure the error of the predictions in percentage terms, and finally the correlation coefficient R^2 , which allows one to measure the correlation between the predicted and observed values. These metrics are implemented through Equations (13)–(15) as follows:

$$RMSE = \sqrt{\frac{\sum_{i=1}^{n} (Pi - Oi)^2}{n}}$$
 (13)

$$MAPE = \frac{1}{n} \sum_{i=1}^{n} \left| \frac{(O_i - P_i)}{O_i} * 100 \right|$$
 (14)

$$R^{2} = \frac{\sum_{i=1}^{n} \left(O_{i} - \overline{O}\right) \left(P_{i} - \overline{P}\right)}{\sum_{i=1}^{n} \left(O_{i} - \overline{O}\right)^{2} \sum_{i=1}^{n} \left(P_{i} - \overline{P}\right)^{2}}$$
(15)

where P is the predicted value vector, O is the observed value vector, \overline{O} is the mean of observed values, and \overline{P} is the mean of the predicted values.

4. Results and Discussion

In this section, the achieved results are described.

4.1. Results

4.1.1. Weighted Averages and Polynomial Interpolation (WA + PI)

Tables 4–6 show comparisons between WA and WA + PI for each environmental monitoring station.

In Tables 4–6, it can be seen that, in terms of RMSE (66.67%) and MAPE (50.00%), in most cases, polynomial interpolation (PI) allows one to improve the results derived using the weighted average (WA), improving the WA between 0.0582% and 1.6433%; hence, one can see the importance of PI in the implementation of the model proposed in this work.

Table 4.	Results	of WA and	l WA +	PI (Pacocha).
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	Tashmiana	P	Predicted Hours	
	Technique ——	24	48	72
RMSE				
	WA	4.8293	10.9212	12.9891
	WA + PI	4.0954	11.1827	12.9212
MAPE				
	WA	37.8821	31.0839	30.6611
	WA + PI	27.2208	29.95	29.5541
\mathbb{R}^2				
	WA	0.7781	0.1642	0.1293
	WA + PI	0.5412	0.1175	0.1209

Table 5.	Results of	WA a	nd WA +	PΙ	(Bolognesi).
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	Technique ——	P	redicted Hours	
	recinique ——	24	48	72
RMSE				
	WA	2.9791	6.5504	5.7880
	WA + PI	3.2104	6.6086	5.7738
MAPE				
	WA	19.9318	25.1731	26.1533
	WA + PI	21.5751	25.2608	25.3131
\mathbb{R}^2				
	WA	0.0111	0.0025	0.0031
	WA + PI	0.0086	0.0000	0.0054

Table 6. Results of WA and WA + PI (Pardo).

	Tashmiana	P	Predicted Hours	
	Technique ——	24	48	72
RMSE				
	WA	1.7193	2.1458	2.1535
	WA + PI	1.2283	1.7856	1.8002
MAPE				
	WA	13.3863	18.0499	20.9966
	WA + PI	13.5280	19.1461	20.3845
\mathbb{R}^2				
	WA	0.2852	0.1614	0.0000
	WA + PI	0.6414	0.2315	0.0014

4.1.2. A Web Application Based on WA + PI

The constructed model was used for the implementation of a web application. It has a simple architecture, which includes the front-end and the back-end. HTML5, Javascript, Bootstrap, and Chart.js were used to implement the front-end. The back-end was Python, and the Flask framework was used for the development of the web application. Figure 8 shows the main view of the web application.

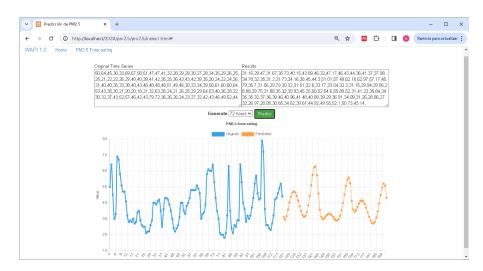


Figure 8. Web application for PM2.5 forecasting.

4.2. Discussions

4.2.1. Comparison with Benchmark Models

To evaluate how the proposed model contributes to advancing the state of knowledge in this field, well-known, multi-step, deep learning-based, state-of-the-art models—such as LSTM, BiLSTM, GRU, BiGRU, and LSTM-ATT—were implemented.

LSTM is a type of recurrent neural network (RNN) that is used to process and predict data sequences. An LSTM cell is composed of an input gate, an output gate, and a forget gate, which enables useful information to be remembered.

The BiLSTM is based on LSTMs, with the main difference being that it is bidirectional. In LSTMs, information flows in a single direction from left to right; meanwhile, with BiLSTM, information flows in both directions, from left to right and from right to left. BiLSTMs can learn more complex relationships, but the computational cost is higher.

GRU is a type of RNN; GRU has fewer gates (only two: reset and update) and fewer parameters than LSTMs. LSTMs are more complex and effective in modeling complex patterns; meanwhile, GRU is more efficient and easier to train. The choice between an LSTM and a GRU depends on the specific problem and the available resources.

BiGRU is similar to BiLSTM, with the main difference being that the BiGRU is based on GRU.

LSTM-ATT is a variant of LSTMs that incorporates an attention mechanism to improve the ability to model data sequences. The attention mechanism allows the model to focus on certain parts of the input sequence that are more relevant to a specific task. This is useful when working with long sequences, where only some parts of the sequence are important for predictions.

A graphical view of the benchmark model architectures can be seen in Figure 9. The hyperparameters of the benchmark models are listed in Table 7.

Table 7. Hyperparameters of the ben	chmark models.
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Model	Hyperparameters
LSTM	$[30,30,30,n *]$, $lr = 0.001$, $drop_rate = [",0.1,0.1]$
BiLSTM	[30,30,30,n *], lr = 0.001, drop_rate = [",0.1,0.1]
GRU	[30,30,30,n *], lr = 0.001, drop_rate = [",0.1,0.1]
BiGRU	$[30,30,30,n *]$ lr = 0.001, drop_rate = [",0.1,0.1]
LSTM-ATT	[100,1100,n *], lr = 0.001

^{*} n is the number of hours to be predicted (24, 48 or 72).

The lock-back for each model was set to 24 h. RNNs, with no attention mechanism, include three layers of 30 neurons each, while RELU is the activation function, and drop_rates of 0.1 were used after each layer to avoid overfitting.

An LSTM with an attention mechanism includes one LSTM layer with 100 neurons, followed by a dense layer with one neuron and tanh as activation, an activation layer (softmax), a repeatVector(100), a permute layer, a multiply layer, a lambda, and a final dense layer with n neurons, depending on the number of hours to be predicted.

All models were trained with 100 epochs, using adam as the optimizer and mse as the loss function. The 100 epochs used produced the best results for each model.

The obtained results are summarized in Tables 8–10.

According to Table 8, for the Pacocha station, in terms of RMSE, the proposed model outperformed the benchmark models in most cases (66.67%) for 24 and 72 predicted hours, obtaining RMSEs of 4.0954 and 12.9212, respectively. For 48 predicted hours, the proposed model was only surpassed by the BiLSTM and LSTM. In terms of MAPE, the proposed model outperformed all the benchmark models (100%). In terms of R², it outperformed all benchmark models for 24 predicted hours; it was outperformed by BiLSTM and BiGRU for 48 predicted hours; and for 72 predicted hours, it was outperformed by BiLSTM only.

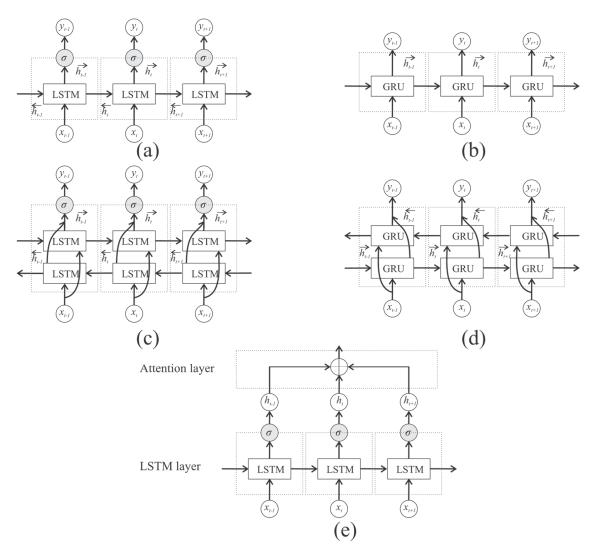


Figure 9. Architectures of the benchmark models: (a) LSTM, (b) GRU, (c) BiLSTM, (d) BiGRU, and (e) LSTM-ATT.

Table 8. The benchmark models vs. the proposed model in Pacocha station.

Tarkeriana	Predicted Hou	ırs	
Technique	24	48	72
RMSE			
LSTM	6.2482	11.1098	12.9222
BiLSTM	5.0760	11.0872	12.8388
GRU	5.9334	11.4898	13.2358
BiGRU	5.2568	11.7640	13.5366
LSTM-ATT	5.7577	12.7558	15.3852
WA + PI	4.0954	11.1827	12.9212
MAPE			
LSTM	49.9085	45.6897	43.9890
BiLSTM	36.7623	38.0884	38.0875
GRU	49.1408	45.6031	44.4407
BiGRU	32.0877	36.7171	37.1657
LSTM-ATT	30.5190	36.9551	40.6333
WA + PI	27.2208	29.9500	29.5541

 Table 8. Cont.

Tashniaus	Predicted Hou	urs	
Technique	24	48	72
R ²			
LSTM	0.1101	0.1057	0.0336
BiLSTM	0.4451	0.3815	0.1410
GRU	0.0069	0.0068	0.0062
BiGRU	0.1653	0.1986	0.0377
LSTM-ATT	0.0227	0.0738	0.0725
WA + PI	0.5412	0.1175	0.1209

Table 9. The benchmark models vs. the proposed model in Bolognesi station.

Talkaiana	Predicted Hours			
Technique	24	48	72	
RMSE				
LSTM	2.6880	5.6406	5.1198	
BiLSTM	2.7096	6.0241	5.3871	
GRU	2.5525	5.7999	5.2771	
BiGRU	2.5477	6.2492	5.5734	
LSTM-ATT	2.2942	5.7173	5.0636	
WA + PI	3.2104	6.6086	5.7738	
MAPE				
LSTM	20.3258	19.8867	24.0068	
BiLSTM	20.0818	21.6240	24.8745	
GRU	18.9890	19.5996	24.2701	
BiGRU	17.4580	21.2114	24.5218	
LSTM-ATT	17.2521	18.6097	20.9124	
WA + PI	21.5751	25.2608	25.3131	
R ²				
LSTM	0.0543	0.0455	0.2520	
BiLSTM	0.0569	0.0002	0.0890	
GRU	0.0248	0.0203	0.1359	
BiGRU	0.1528	0.1011	0.0001	
LSTM-ATT	0.1875	0.0003	0.2360	
WA + PI	0.0086	0.0000	0.0005	

Table 10. The benchmark models vs. the proposed model in Pardo station.

Tarkeriana	Predicted Hours			
Technique	24	48	72	
RMSE				
LSTM	1.5667	2.2181	2.5044	
BiLSTM	1.4403	2.0565	2.3202	
GRU	1.5048	2.0773	2.3194	
BiGRU	1.5725	1.8373	1.8743	
LSTM-ATT	1.4485	1.8306	2.4208	
WA + PI	1.2283	1.7856	1.8002	

Table 10. Cont.

Tashmiana	Predicted Hours			
Technique	24	48	72	
MAPE				
LSTM	17.0770	25.9310	30.9449	
BiLSTM	15.4638	23.7230	28.3142	
GRU	16.0796	23.7334	28.2635	
BiGRU	17.0590	21.0457	22.1567	
LSTM-ATT	15.7723	19.3737	28.1934	
WA + PI	13.5280	19.1461	20.3845	
R ²				
LSTM	0.6403	0.3008	0.1022	
BiLSTM	0.6472	0.3075	0.1133	
GRU	0.4997	0.2422	0.0007	
BiGRU	0.0876	0.0505	0.1475	
LSTM-ATT	0.6343	0.1627	0.0170	
WA + PI	0.6414	0.2315	0.0014	

According to Table 9, for the Bolognesi station, the proposed model obtained the worst results. In terms of RMSE, for the 24 predicted hours, the proposed model was the worst, with 3.2104 $\rm ug/m^3$; for the 48 predicted hours, the proposed model was outperformed by LSTM-ATT and LSTM. For the 72 predicted hours, the proposed model was the worst model, with 7.7738 $\rm ug/m^3$. In terms of MAPE, the proposed model was outperformed by all the benchmark models. In terms of $\rm R^2$, the proposed model was outperformed by all models for 24 and 48 predicted hours, and for 72 predicted hours, it outperformed only BiGRU. The best model for Bolognesi station was LSTM-ATT.

According to Table 10, for the Pardo station, the proposed model obtained the best results. In terms of RMSE and MAPE, the proposed model outperformed all the benchmark models. In terms of R², for 24 predicted hours, the proposed model (0.6414) was outperformed by BiLSTM (0.6472); for 48 predicted hours, the proposed model (0.2315) was outperformed by BiLSTM (0.3075), LSTM (0.3008), and GRU (0.2422); for 72 predicted hours, the proposed model was outperformed by all the benchmark models except GRU.

Graphical comparisons can be viewed in Figures 10–12.

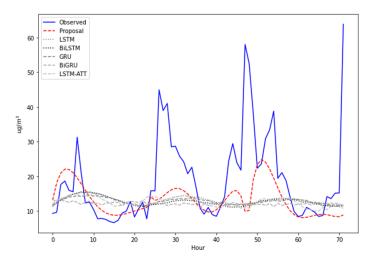


Figure 10. The 72 predicted hours for Pacocha station.

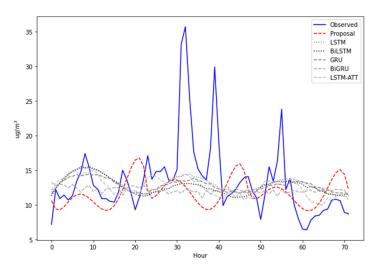


Figure 11. The 72 predicted hours for Bolognesi station.

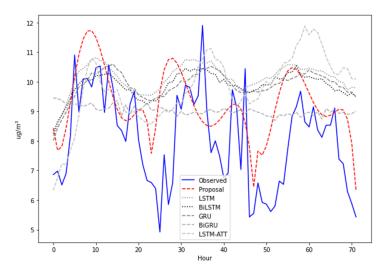


Figure 12. The 72 predicted hours for Pardo station.

In this work, three different prediction horizons were experimented with for multistep forecasting—24, 48, and 72 h. The proposed model produced similar and superior results to the benchmark models, surpassing them in most cases in terms of RMSE and MAPE. However, it is important to explain why the proposed model produced these results for the different environmental monitoring stations; for this purpose, correlations were determined between the first day of the test data and the last training days. Table 10 shows these correlations.

According to Table 11, it can be seen that in the case of the Bolognesi station—which is where the proposed model presented the worst results—the last training data that correspond to day 1 present a very low correlation (0.0115). Likewise, the data for the other day used, which is day 3, also present a very low correlation (-0.009).

However, for the Pacocha and Pardo stations, for the first day and third day, there are better correlations of between 0.293 and 0.6018, and this is reflected in the results of the proposed model.

The above means that the average correlation used for the selection of days may not be good in some cases; because it is an average, it may only be good for some cases. For future work, different day selections could be made for each station and not just the ones used in this work.

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Day	Pacocha	Bolognesi	Pardo
1	0.293	0.0115	0.5334
2	0.0054	0.5976	-0.2219
3	0.4315	-0.009	0.6018
4	0.3573	0.5646	0.1011
5	-0.3451	0.6147	0.2458
6	0.3808	0.1252	0.1812
7	0.323	0.0238	0.2447
8	0.2735	0.5718	0.3889
9	0.3125	0.3951	0.1212
10	0.275	0.7076	0.1224
11	-0.2403	0.2895	-0.0402

Table 11. Correlations between the first day of test data and the last training days.

4.2.2. Comparison with Related Works

Next, the proposed model results were compared with the results of multi-step related works.

Finally, the results of the proposed model were compared with those reported in related works; according to Table 12, in terms of the five studies addressing MAPE, the research presented in [44,45,47,49,54] obtained MAPEs of 5.11%, 5.12%, 15.32%, 14.08%, and 7.53%. These studies reported MAPEs higher than those obtained by the proposed model (17.69%); three of them used three-step forecasting, and two of them used twelve-and seven-step forecasting. Therefore, there is an enormous difference between the number of steps in the related works and the proposed model. This is a factor that is in favor of the results presented in these other studies; this is because, as the number of steps increases, the accuracy of the results worsens.

Tab!	e 12.	Related	works	s vs.	pro	posal.
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Work	Technique	Freq.	Steps	Metric	Value
[45]	ELM	Hourly	3	MAPE	[5.11–37.23]
[46]	ELM	Hourly	7	MAPE	[5.12-22.2]
[47]	CNN-LSTM	Hourly	10	MAPE	[22.09-25.94]
[48]	Adaboost	Hourly	3	MAPE	[15.32-23.75]
[49]	SVM	Hourly	4	RMSE	[8.16-39.04]
[50]	LSTM	Daily	1, 12	MAPE	[6.66–14.08]
[34]	LSTM	Hourly	1, 6	MAE	[4.395–7.246]
[51]	BiLSTM	Hourly	3	MAPE	[27.33-40.73]
[52]	CNN-BP	Hourly	10	RMSE	[5.42–10.21]
[53]	ResCNN	Hourly	4	MAPE	[28.51-33.29]
[54]	TMSG	Hourly	3	MAPE	[7.53–16.18]
Proposal	WA + PI	Hourly	24, 48, 72	MAPE RMSE	[17.69–28.91] [1.23–5.77]

On the other hand, the results for the proposed model surpass the results presented in [46,51,53]; the authors of these works implemented techniques similar to the benchmark models implemented in the present paper.

In terms of RMSE, the proposed model considerably exceeds the RMSEs reported in [48,52].

5. Conclusions and Future Work

According to the obtained results, it can be concluded that the proposed model based on weighted averages (WAs) and polynomial interpolation (PI) constitutes a good alternative for the multi-step forecasting of PM2.5 time series. The proposed model outperformed the other models for most of the analyzed environmental monitoring stations. The well-known, state-of-the-art models that were used for these comparisons included LSTM,

BiLSTM, GRU, BiGRU, and LSTM-ATT. The proposed model outperformed these models in terms of the average MAPE obtained (between 2.4010% and 17.6208%). One of the main advantages of the proposed model is that it requires few data compared to these benchmark models, which are based on deep learning.

Although the results are interesting, and the implemented benchmark models had been outperformed in most cases, some aspects still need improvement. The achieved MAPEs were between 17.69% and 28.91%. According to [57], if the MAPE is less than 10%, then the model is highly accurate; if the MAPE is between 10% and 20%, then the model is good; if the MAPE is between 20% and 50%, then the model is reasonable. Thus, the proposed model can be considered as being between good and reasonable; it should be highly accurate. In pursuit of a highly accurate rating, in future work, instead of working with only two days, it could be adjusted to function with three or more days, depending on the level of the correlations. Similarly, instead of polynomial interpolation, other types of interpolation, such as linear, Stineman, inverse distance weighting, or kriging, could be used. On the other hand, ensemble and hybrid models could be created with other forecasting techniques, such as autoregressive integrated moving average (ARIMA).

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Conflicts of Interest: The authors declare no conflict of interest.

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