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Article

# MGGTSP-CAT: Integrating Temporal Patterns and LSTM for Multi-Scale Greenhouse Gas Time Series Prediction via Cross-Attention Mechanism

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Abstract: Accurate greenhouse gas (GHG) concentration forecasting is pivotal for evaluating climate policies, yet existing models face challenges in capturing multi-scale temporal patterns. To address this limitation, we propose MGGTSP-CAT, a novel multi-encoder framework that integrates daily and monthly data through an Input Attention encoder and an Autoformer encoder, enhanced by a Temporal Attention mechanism. Evaluated on NOAA datasets spanning five decades from Mauna Loa, Barrow, American Samoa, and Antarctica, MGGTSP-CAT outperforms 14 baseline models, achieving a Test\_R² of 0.9627 and Test\_MAPE of 1.47%. Crucially, the model demonstrates robustness in multi-step forecasting, with Test\_R² declining by only 3.3% over ten steps—significantly lower than Transformer-based counterparts. These results underscore the framework's capability to decouple short-term fluctuations from long-term trends, offering a reliable tool for cross-scale climate predictions and policy design.

**Keywords:** greenhouse gas forecasting; time series prediction; multi-scale modeling; cross-attention mechanism

# 1. Introduction

Climate change is a global challenge impacting all of humanity, and greenhouse gas emissions data plays a crucial role in assessing the effectiveness of environmental policies and understanding the dynamic changes in the global climate system [1–3]. As a key parameter for quantifying climate system states, greenhouse gas concentration reflects the imbalance in the Earth's carbon cycle, far surpassing the significance of regional or national emissions data. By integrating diverse data sources, including atmospheric monitoring, satellite remote sensing, and ground observations, greenhouse gas concentrations provide accurate insights into human activity, ecosystem carbon sink capacity, and the effectiveness of climate policies. This information serves as a foundation for predicting global climate trends, supporting international climate negotiations, and aiding governments in optimizing emission reduction strategies and identifying critical points in the carbon peak process, ultimately enhancing climate governance [4].

Accurately capturing the balance between carbon sources and sinks is essential for reliable greenhouse gas concentration predictions. To improve prediction accuracy, multi-dimensional modeling approaches are employed to assess natural systems, human activities, policy interventions, and technological innovations. Policy analysis plays a vital role in evaluating international agreements like the Paris Agreement, carbon pricing reforms, and clean energy promotion, all of which shape emission intensity and carbon removal efficiency [5,6]. Historical data analysis, including trend decomposition and anomaly detection, unveils the interaction mechanisms between human activities and natural systems, especially near carbon cycle critical thresholds [7,8].

However, policy analysis and scenario simulations are often fraught with uncertainty due to their dependence on political will, economic transitions, and unpredictable factors like technology

diffusion. Data analysis, in contrast, relies more heavily on observational data and scientific principles, offering a more stable and repeatable basis for prediction [9]. Despite the strong track record of traditional climate models, such as Global Circulation Models (GCMs) and Regional Climate Models (RCMs), these models struggle with accurately simulating short-term concentration fluctuations and fine-scale predictions. Smith et al. constructed a global climate model and made predictions for the next decade, significantly improving prediction accuracy [10]. However, these models still exhibit significant uncertainty, especially in characterizing short-term fluctuations and ecosystem feedback mechanisms, such as methane release due to permafrost thawing, which complicates their application in greenhouse gas forecasting.

With the growing demand for more precise climate predictions, data-driven approaches are becoming increasingly important. Machine learning models, such as Random Forests [11] and Gradient Boosting Decision Trees [12], have demonstrated their ability to handle complex, high-dimensional data, uncovering nonlinear relationships and identifying regional emission hotspots. For example, Zhang Jianxun et al. used the XGBoost algorithm to predict carbon emissions in urban expansion, highlighting the model's strength in predicting emissions for rapidly expanding megacities [13]. However, these models lack physical interpretability and are highly dependent on data quality, which can lead to biases in regions with sparse observation networks.

Recent breakthroughs in deep learning, particularly with architectures like Long Short-Term Memory (LSTM) [14] and Spatio-Temporal Graph Convolution Networks (STGCNs), have enhanced the prediction of greenhouse gas concentrations by capturing spatio-temporal correlations in atmospheric transport. Zhang Lei et al. demonstrated the use of CNN-LSTM for predicting soil organic carbon content, proving its effectiveness in regional carbon content forecasting [15]. Panja et al. developed the E-STGCN model for air quality forecasting in Delhi, India, achieving consistent performance across all seasons [16]. These models show promise but still face challenges, particularly in regions with limited data coverage and physical inconsistencies in extrapolated predictions. The Transformer architecture [17], with its self-attention mechanism, offers a new avenue for modeling global atmospheric transport, improving prediction accuracy for carbon emissions. Wu Xingping et al. implemented a temporal graph transformer-based attentional neural network to predict carbon emissions, achieving an 89.5% accuracy rate [18]. Despite its potential, the application of Transformers in climate prediction still requires further development, especially for extreme weather conditions.

Emerging deep learning models like Temporal Convolutional Networks (TCN) [19], DA-RNN[20], Autoformer [21] have shown promise in modeling greenhouse gas concentrations by capturing seasonal periodicity and inter-annual variations in the carbon cycle. These architectures have yet to achieve breakthroughs in accuracy for emissions forecasting with a single time resolution.

To address the limitations of current models, this study introduces a multi-encoder fusion cross-scale forecasting framework, MGGTSP-CAT. By constructing a multi-time-resolution dataset, we integrate optimized Input-Attention and Autoformer Encoders, alongside LSTM, to extract features from both daily (short-term) and monthly (long-term) data. The multi-step forecasting is achieved through Temporal Attention, providing robust prediction accuracy and stability across different datasets. The proposed framework overcomes existing model limitations and sets a new standard for greenhouse gas concentration forecasting.

### 2. Materials and Methods

# 2.1. Data Curation

The dataset consists of daily carbon dioxide mole fraction data from continuous atmospheric measurements at stations in Mauna Loa (Hawaii), Barrow (Alaska), American Samoa, and Antarctica, collected by the National Oceanic and Atmospheric Administration (NOAA) [22]. The time span of the dataset is from July 24, 1973, to April 30, 2024, comprising 16,597 daily data points.

This dataset provides a benchmark platform for multi-scale prediction models, offering both high time resolution (daily) and long-term coverage (over half a century). The spatial distribution

characteristics of different stations allow for effective validation of the model's ability to simultaneously model regional specificity (e.g., the Arctic amplification effect) and global consistency (e.g., long-term trends). In particular, the multi-modal features contained in the daily data, such as short-term fluctuations (influenced by weather systems) and long-term trends (driven by human emissions), offer an ideal experimental scenario for testing the performance of the multi-time-resolution hybrid architecture in decoupling temporal scales.

#### 2.2. Methods

Overall, our model is divided into several parts: First, daily data is processed into monthly data for subsequent use (Input Phase). The encoding phase follows, where two different encoders are applied for different time resolution data: the Input-Attention Mechanism Encoder is used for daily data, and the Autoformer Encoder is used for monthly data. After all encoding is completed, the encoded daily and monthly data matrices are concatenated and fed into the decoder for decoding. The decoding phase utilizes a Temporal-Attention-based decoder to generate the output results (Output). The network diagram is shown in Figure 1.

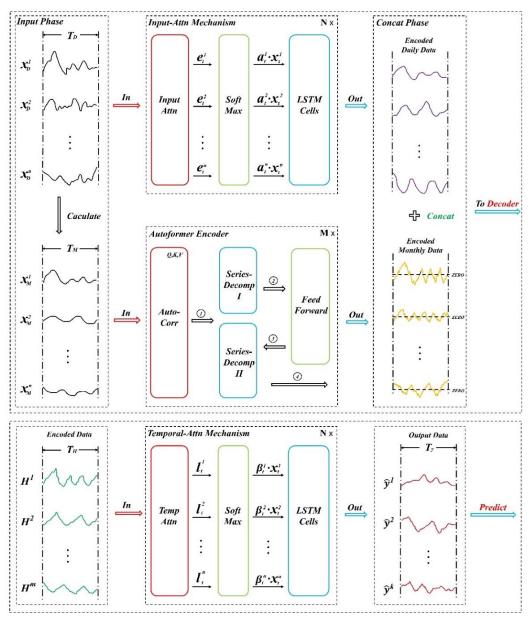


Figure 1. MGGTSP-CAT Encoding and Decoding Mechanism.

#### 2.2.1. Encoder

We first describe how to compute weekly and monthly data from daily data. For the daily input sequence  $X_D = (x_{D_1}, x_{D_2}, ..., x_{D_{T_d}})$ , monthly data  $X_M = (x_{M_1}, x_{M_2}, ..., x_{M_{T_m}})$  is computed by averaging every 30 data points:

$$X_{M} = \left(\frac{\sum_{i=1}^{30} x_{W_{i}}}{30}, \frac{\sum_{i=31}^{60} x_{W_{i}}}{30}, \dots, \frac{\sum_{i=T_{w}-29}^{T_{w}} x_{W_{i}}}{30}\right)$$
(1)

For daily data, we use an encoder based on Input Attention. The daily input sequence  $X_D = (x_{D_1}, x_{D_2}, ..., x_{D_{T_d}})$  is processed through the Input Attention Mechanism, which includes the Input Attention Unit and LSTM units. The encoder uses this input attention mechanism to select and weight the input sequence, and uses LSTM units to generate hidden state sequences.

Specifically, for each time step t, we first compute the basic hidden state  $h_t$  through the LSTM unit:

$$h_t = f_1(h_{t-1}, x_t) (2)$$

Next, we calculate the attention weights  $a_t^k$ , via the input attention mechanism and extract the new relevant driving sequence:

$$e_t^k = v_e^T \tanh(W_e[h_{t-1}; s_{t-1}] + U_e x_t^k)$$
 (3)

$$\alpha_t^k = \frac{\exp(e_t^k)}{\sum_{i=1}^n \exp(e_t^i)} \tag{4}$$

$$\tilde{\mathbf{x}}_t = (\alpha_t^1 \mathbf{x}_t^1, \alpha_t^2 \mathbf{x}_t^2, \dots, \alpha_t^n \mathbf{x}_t^n)^T \tag{5}$$

Using the new driving sequence, we update the hidden state:

$$h_t = f_1(h_{t-1}, \tilde{x}_t) \tag{6}$$

where  $f_1$  is the LSTM unit, and its internal computation is as follows:

$$\begin{cases} f_{t} = \sigma(W_{i}[h_{t-1}; x_{t}] + b_{f}) \\ i_{t} = \sigma(W_{i}[h_{t-1}; x_{t}] + b_{i}) \\ o_{t} = \sigma(W_{i}[h_{t-1}; x_{t}] + b_{o}) \end{cases}$$
(7)

$$s_t = f_t \odot s_{t-1} + i_t \odot \tanh(W_s[h_{t-1}; x_t] + b_s)$$
 (8)

$$h_t = o_t \odot tanh(s_t) \tag{9}$$

where  $\sigma$  represents the *Sigmoid* activation function, and  $\odot$  represents element-wise multiplication.  $W_f$ ,  $W_i$ ,  $W_o$ ,  $W_s \in \mathbb{R}^{m \times (m+n)}$  are weight matrices, and  $b_f$ ,  $b_i$ ,  $b_o$ ,  $b_s \in \mathbb{R}^m$  are bias vectors.

Thus, we obtain the hidden state sequence for daily data: $H_D = (h_{D_1}, h_{D_2}, ..., h_{D_{T_d}})$ 

For monthly data, we use the Autoformer Encoder. The monthly input sequence  $X_M = \left(x_{M_1}, x_{M_2}, ..., x_{M_{T_m}}\right)$  is processed Autoformer Encoder , which mainly consists of the series decomposition block (Series Decomp), auto-correlation mechanism (Auto-Correlation), and feedforward neural network (Feed Forward). The encoder uses these components to process the input sequence, extract long-term trend information, and model seasonal patterns, thus generating the hidden state sequence for monthly data.

Specifically, for each time step t, we first process the input sequence through the series decomposition block. Since directly decomposing future sequences is not realistic, Autoformer proposes the series decomposition block as an internal operation to progressively extract the long-term stable trend from the predicted intermediate hidden variables. The process smoothes periodic

fluctuations through moving averages (using AvgPool and Padding operations) to highlight long-term trends. For an input sequence  $X \in \mathbb{R}^{L \times d}$  of length L, the calculation is as follows:

$$X_t = AvgPool(Padding(x))$$
 (10)

$$X_{s} = x - x_{t} \tag{11}$$

where  $X_s$ ,  $X_t \in \mathbb{R}^{L \times d}$  represent the seasonal part and the extracted trend-periodic part. We summarize this process in the internal block  $X_s$ ,  $X_t = SeriesDecomp(X)$ .

Next, the output of the auto-correlation mechanism is calculated. The auto-correlation mechanism identifies cycle-based dependencies and aggregates similar sub-series through time-delay aggregation. First, we introduce two basic mechanisms: cycle-based dependencies and time-delay aggregation, then extend them to single-head and multi-head auto-correlation mechanism outputs.

Cycle-based Dependencies: For a real discrete-time process  $\{X_t\}$ , the auto-correlation is computed as follows:

$$R_{xx}(\tau) = \lim_{L \to \infty} \frac{1}{L} \sum_{t=1}^{L} X_t X_{t-\tau}$$
 (12)

(In practical computations, we select the most likely k periodicities  $\tau_1, ..., \tau_k$ , where  $k = \lfloor c \times log L \rfloor$ , with c being a hyperparameter) This autocorrelation reflects the temporal similarity between  $\{X_t\}$  and its  $\tau$ -lagged series  $\{X_{t-\tau}\}$ , thus capturing the periodic dependencies.

Time-delay Aggregation: Based on the selected time delays  $\tau_1, ..., \tau_k$ , the series is rolled using a *Roll* operation to align similar subsequences that share the same phase of the estimated period. These subsequences are then aggregated using a *Softmax*-normalized confidence score.

Thus, for the single-head case and a time series x of length L, after projecting the sequence to obtain the query Q, key K and value V matrices, the auto-correlation mechanism is computed as follows:

$$\tau_1, \dots, \tau_k = \operatorname{argTopk}\left(R_{Q,K}(\tau)\right), \tau \in \{1, \dots, L\}$$
(13)

$$\hat{R}_{Q,K}(\tau_1), \dots, \hat{R}_{Q,K}(\tau_k) = softmax\left(R_{Q,K}(\tau_1), \dots, R_{Q,K}(\tau_k)\right) \tag{14}$$

$$Auto-Correlation(Q,K,V) = \sum_{i=1}^{k} Roll(V,\tau_i) \, \hat{R}_{Q,K}(\tau_i)$$
 (15)

For multi-head mechanisms with hidden variable channels  $d_{model}$ , and number of heads h, the calculation process is:

$$head_{i} = Auto - Correlation(Q_{i}, K_{i}, V_{i}), Q_{i}, K_{i}, V_{i} \in \mathbb{R}^{L \times \frac{d_{model}}{h}}, i \in \{1, \dots, h\}$$
 (16)

$$MultiHead(Q, K, V) = W_0 * Concat(head_1, ..., head_h)$$
 (17)

Thus, the calculation of the auto-correlation mechanism output is completed.

The output of the auto-correlation mechanism is further transformed through a feed-forward neural network. The computation of the feed-forward network is similar to the Transformer Encoder (including fully connected layers with an activation function, such as ReLU)

$$FeedForward(x) = ReLU(xW_1 + b_1)W_2 + b_2$$
(18)

Let the output of the auto-correlation mechanism Auto-Correlation(Q,K,V) or MultiHead(Q,K,V) be  $h_t^{attn}$ , then:

$$h_t^{ffn} = FeedForward(h_t^{attn}) \tag{19}$$

Finally, the output hidden state is obtained through residual connections and layer normalization:

$$h_t = LayerNorm(h_t^{attn} + h_t^{ffn})$$
 (20)

Thus, we obtain the hidden state sequence for monthly data  $H_M = \left(h_{M_1}, h_{M_2}, \dots, h_{M_{T_m}}\right)$ 

#### 2.2.2. Decoder

After encoding, the hidden state sequences  $H_D$ ,  $H_M$  from the two encoders are concatenated to form the final hidden sequence  $H = (h_1, h_2, ..., h_T)$ :

$$H = Concat(H_D, H_M) (21)$$

This concatenated sequence is then passed into the Temporal Mechanism to decode the final prediction sequence. The Temporal Mechanism includes the Temporal Attention Unit and LSTM unit, where the Temporal Attention Unit is used to select the hidden states of the encoders, and the LSTM unit updates the decoder's hidden state and outputs the final prediction  $\hat{y}_T$ .

Specifically, for each time step t, we first compute the time attention scores:

$$l_t^i = v_d^T \tanh(W_d[d_{t-1}; s'_{t-1}] + U_d h_i)$$
(22)

Then, we calculate the attention weights:

$$\beta_t^i = \frac{\exp(l_t^i)}{\sum_{j=1}^T \exp(l_t^j)}$$
 (23)

Next, we compute the context vector:

$$c_t = \sum_{i=1}^T \beta_t^i h_i \tag{24}$$

The decoder input is combined with the context vector to generate the new input:

$$\tilde{y}_{t-1} = \tilde{w}^T[y_{t-1}; c_{t-1}] + \tilde{b}$$
 (25)

The decoder's hidden state is updated:

$$d_t = f_2(d_{t-1}, \tilde{\gamma}_{t-1}) \tag{26}$$

Finally, we output the prediction:

$$\hat{\mathbf{y}}_T = v_v^T (W_v[d_t; c_t] + b_w) + b_v \tag{27}$$

where  $f_2$  is the LSTM unit, and its internal computation is:

$$\begin{cases} f'_{t} = \sigma(W'_{f}[d_{t-1}; \tilde{y}_{t-1}] + b'_{f}) \\ i'_{t} = \sigma(W'_{i}[d_{t-1}; \tilde{y}_{t-1}] + b'_{i}) \\ o'_{t} = \sigma(W'_{o}[d_{t-1}; \tilde{y}_{t-1}] + b'_{o}) \end{cases}$$
(28)

$$s'_{t} = f'_{t} \odot s'_{t-1} + i'_{t} \odot tanh (W'_{s}[d_{t-1}; \tilde{y}_{t-1}] + b'_{s})$$
 (29)

$$d_t = o'_t \odot \tanh(s'_t) \tag{30}$$

where  $\sigma$  represents the *Sigmoid* activation function,  $\odot$  denotes element-wise multiplication,  $W_y \in \mathbb{R}^{p \times (p+m)}$  and  $v_y \in \mathbb{R}^p$  are the weight matrices and weight vectors, while  $b_w \in \mathbb{R}^p$  and  $b_v \in \mathbb{R}$  are the bias vectors.

#### 2.2.3. Metric

The following are the formulas for calculating the goodness-of-fit of the model:  $R^2$ , Coefficient of Determinatio:

$$R^{2} = 1 - \frac{\sum_{i=1}^{n} (y_{i} - \hat{y}_{i})^{2}}{\sum_{i=1}^{n} (y_{i} - \bar{y})^{2}}$$
(31)

where  $y_i$  represents the true value,  $\hat{y}_i$  represents the predicted value,  $\bar{y}$  represents the mean of the true values, n represents the sample size.

MSE, Mean Squared Error:

$$MSE = \frac{1}{n} \sum_{i=1}^{n} (y_i - \hat{y}_i)^2$$
 (32)

where  $y_i$  represents the true value,  $\hat{y}_i$  represents the predicted value, n represents the sample size. MAE, Mean Absolute Error:

$$MAE = \frac{1}{n} \sum_{i=1}^{n} |y_i - \hat{y}_i|$$
 (33)

where  $y_i$  represents the true value,  $\hat{y}_i$  represents the predicted value, n represents the sample size. MAPE, Mean Absolute Percentage Error:

$$MAPE = \frac{100\%}{n} \sum_{i=1}^{n} \left| \frac{y_i - \hat{y}_i}{y_i} \right|$$
 (34)

where  $y_i$  represents the true value,  $\hat{y}_i$  represents the predicted value, n represents the sample size.

# 3. Results

#### 3.1. Overall Performances

The proposed MGGTSP-CAT framework demonstrates superior predictive capabilities across all evaluation metrics compared to 14 baseline models. As summarized in Table 1, MGGTSP-CAT achieves the highest Test\_R<sup>2</sup> of 0.9627, significantly outperforming DARNN (0.9519) and Autoformer (0.9480). This improvement is attributed to the model's ability to effectively integrate multi-scale temporal patterns, capturing both short-term fluctuations (e.g., daily weather variations) and long-term trends (e.g., seasonal and policy-driven changes). The model also exhibits the lowest error metrics, with Test\_MSE = 0.000288 (22.7% reduction vs. DARNN) and Test\_MAE = 0.0116 (11.3% improvement vs. Autoformer). Notably, the Test\_MAPE of 1.47% confirms the model's robustness in minimizing percentage-based deviations, which is critical for practical applications where relative error margins are more meaningful than absolute values. Figure 2 visually illustrates the alignment between predicted and true values over ten forecasting steps.

Table 1. All test set data of 15 models.

Model	Test_MSE	Test_MAE	Test_R <sup>2</sup>	Test_MAPE
MGGTSP	0.0002878	0.0116377	0.9627339	0.0147310
DARNN	0.0003716	0.0131211	0.9518857	0.0167605
Autoformer	0.0004013	0.0136400	0.9480379	0.0174719
TCN	0.0004571	0.0146228	0.9408147	0.0187982
BiTransfomer_LSTM	0.0005962	0.0184426	0.9227868	0.0231243
Informer	0.0006874	0.0182407	0.9109833	0.0236269
LSTM	0.0008636	0.0243913	0.8881612	0.0295327
Bi_GRU	0.0012059	0.0257306	0.8438375	0.0333236
Bi_LSTM	0.0012679	0.0260454	0.8358032	0.0338971
GRU	0.0012723	0.0292467	0.8352157	0.0367302
RNN	0.0013303	0.0306885	0.8276975	0.0373433
CNN1D	0.0018290	0.0351233	0.7631485	0.0422021
Bi_RNN	0.0025301	0.0413129	0.6723005	0.0485256
CNN1D_LSTM	0.0028931	0.0447957	0.6253179	0.0527006
ANN	0.0043728	0.0612601	0.4336226	0.0744804

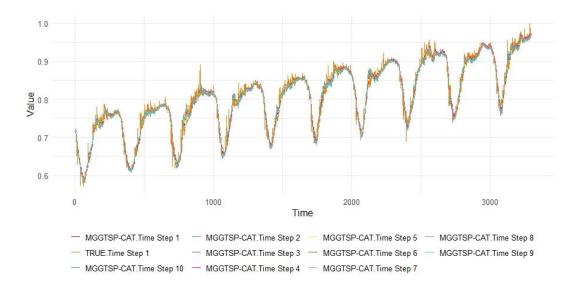


Figure 2. Ten Steps MGGTSP-CAT Predicted Values and True Values on dataset.

# 3.2. Single Step

The model's stability under extended forecasting horizons is rigorously validated through multi-step experiments (Tables 2–5). From Step 1 to Step 10, MGGTSP-CAT maintains a gradual decline in Test\_R² (0.9788  $\rightarrow$  0.9470,  $\Delta$  = 3.3%), outperforming DARNN ( $\Delta$  = 3.9%) and Autoformer ( $\Delta$  = 4.2%). This gradual decline indicates that the model's Temporal Attention mechanism effectively mitigates error accumulation, a common challenge in long-term forecasting. Critically, the rate of error accumulation remains constrained: Test\_MSE increases from 0.000163 (Step 1) to 0.000410 (Step 10), representing a 2.5× rise, compared to Autoformer's 2.3× increase (0.000236  $\rightarrow$  0.000554). This advantage is particularly evident in Steps 5–10, where traditional LSTM/GRU models exhibit significant performance degradation due to their inability to model long-term dependencies effectively. Figure 3 further demonstrates the model's ability to preserve prediction accuracy across diverse baseline comparisons, particularly in Steps 5–10, where traditional LSTM/GRU models exhibit significant performance degradation. The model's ability to maintain high accuracy over extended forecasting horizons is a significant improvement over existing methods, which often struggle with error propagation in multi-step predictions.

**Table 2.** 10 step Test\_R<sup>2</sup> data of 15 models participating in the experiment.

Model	STEP1	STEP2	STEP3	STEP4	STEP5	STEP6	STEP7	STEP8	STEP9	STEP10
MGGTSP	0.97878	0.97443	0.97074	0.96684	0.96564	0.96125	0.95805	0.95458	0.95005	0.94698
DARNN	0.97102	0.96757	0.96023	0.95383	0.95822	0.94851	0.94829	0.94355	0.93422	0.93341
Autoformer	0.96930	0.96559	0.95608	0.94813	0.95578	0.94324	0.94548	0.93990	0.92843	0.92844
TCN	0.96514	0.96146	0.94898	0.93800	0.95049	0.93385	0.93957	0.93361	0.91780	0.91926
BiTransfomer	0.94296	0.92876	0.93205	0.92333	0.91395	0.94672	0.90693	0.91436	0.91156	0.90723
_LSTM	0.71270	0.72070	0.70200	0.72000	0.71070	0.71072	0.70070	0.71100	0.71100	0.50725
Informer	0.94596	0.94410	0.92072	0.89425	0.92664	0.89500	0.91480	0.90963	0.87846	0.88027
LSTM	0.91190	0.90502	0.90029	0.88391	0.89488	0.90956	0.88113	0.85724	0.86588	0.87179
Bi_GRU	0.88297	0.83717	0.90740	0.88788	0.84569	0.76714	0.91339	0.78506	0.74577	0.86590
Bi_LSTM	0.85418	0.92533	0.88067	0.78546	0.82317	0.80695	0.76901	0.85771	0.80601	0.84954
GRU	0.85939	0.84737	0.86361	0.88095	0.79092	0.82703	0.83580	0.79108	0.85289	0.80312
RNN	0.85730	0.84250	0.84315	0.82455	0.82357	0.82105	0.80826	0.81793	0.82042	0.81823
CNN1D	0.79477	0.79974	0.86495	0.76666	0.71630	0.78686	0.73372	0.79676	0.71642	0.65530
Bi_RNN	0.72101	0.70851	0.72330	0.66788	0.61899	0.65601	0.62097	0.69373	0.66151	0.65108
CNN1D_LSTM	0.70959	0.63854	0.65612	0.62404	0.70758	0.59694	0.67335	0.53487	0.60110	0.51105
ANN	0.51734	0.44780	0.37191	0.43568	0.46613	0.43010	0.44904	0.45915	0.47785	0.28123

**Table 3.** 10 step Test\_MSE data of 15 models participating in the experiment.

Model	STEP1	STEP2	STEP3	STEP4	STEP5	STEP6	STEP7	STEP8	STEP9	STEP10
MGGTSP	0.000163	0.000197	0.000226	0.000256	0.000265	0.000299	0.000324	0.000351	0.000386	0.000410
DARNN	0.000223	0.000250	0.000307	0.000356	0.000322	0.000398	0.000400	0.000436	0.000509	0.000515
Autoformer	0.000236	0.000265	0.000339	0.000400	0.000341	0.000438	0.000421	0.000465	0.000553	0.000554
TCN	0.000268	0.000297	0.000393	0.000478	0.000382	0.000511	0.000467	0.000513	0.000636	0.000625
BiTransfomer_LSTM	1 0.000439	0.000549	0.000524	0.000591	0.000664	0.000411	0.000719	0.000662	0.000684	0.000718
Informer	0.000416	0.000431	0.000611	0.000816	0.000566	0.000811	0.000658	0.000699	0.000940	0.000926
LSTM	0.000678	0.000732	0.000769	0.000896	0.000811	0.000698	0.000919	0.001104	0.001037	0.000992
Bi_GRU	0.000901	0.001255	0.000714	0.000865	0.001191	0.001798	0.000669	0.001662	0.001966	0.001037
Bi_LSTM	0.001123	0.000575	0.000920	0.001655	0.001365	0.001491	0.001785	0.001100	0.001500	0.001164
GRU	0.001083	0.001176	0.001052	0.000918	0.001614	0.001336	0.001269	0.001615	0.001138	0.001523
RNN	0.001099	0.001214	0.001209	0.001354	0.001362	0.001382	0.001482	0.001407	0.001389	0.001406
CNN1D	0.001581	0.001543	0.001041	0.001800	0.002190	0.001646	0.002058	0.001571	0.002193	0.002667
Bi_RNN	0.002149	0.002246	0.002133	0.002562	0.002941	0.002657	0.002929	0.002368	0.002618	0.002699
CNN1D_LSTM	0.002237	0.002785	0.002651	0.002900	0.002257	0.003113	0.002524	0.003596	0.003085	0.003783
ANN	0.003717	0.004255	0.004843	0.004354	0.004121	0.004401	0.004257	0.004181	0.004038	0.005561

**Table 4.** 10 step Test\_MAE data of 15 models participating in the experiment.

Model	STEP1	STEP2	STEP3	STEP4	STEP5	STEP6	STEP7	STEP8	STEP9	STEP10
MGGTSP	0.008435	0.009374	0.010210	0.010994	0.011186	0.012021	0.012577	0.013206	0.013918	0.014457
DARNN	0.010061	0.010630	0.011907	0.012869	0.012302	0.013680	0.013829	0.014469	0.015657	0.015807
Autoformer	0.010274	0.010891	0.012558	0.013691	0.012639	0.014409	0.014199	0.014928	0.016397	0.016415
TCN	0.010958	0.011553	0.013643	0.015087	0.013372	0.015661	0.014940	0.015786	0.017707	0.017521
Informer	0.013944	0.014224	0.017385	0.020133	0.016555	0.020105	0.018042	0.018612	0.021806	0.021601
BiTransfomer _LSTM	0.016098	0.018339	0.017876	0.018638	0.020110	0.013880	0.020849	0.019472	0.019489	0.019675
LSTM	0.021328	0.022347	0.022861	0.024917	0.023618	0.021431	0.025366	0.028250	0.027284	0.026510
Bi_GRU	0.022423	0.026983	0.021235	0.022717	0.026333	0.032263	0.019261	0.030710	0.032259	0.023122
Bi_LSTM	0.024449	0.017924	0.022438	0.029404	0.027164	0.028833	0.031595	0.024125	0.027837	0.026685
GRU	0.027698	0.027475	0.026573	0.025132	0.033664	0.030206	0.029865	0.032262	0.028023	0.031569
RNN	0.027605	0.029178	0.029061	0.030809	0.031127	0.031362	0.032624	0.031747	0.031580	0.031791
CNN1D	0.032972	0.032569	0.026367	0.035236	0.038554	0.033392	0.037443	0.033172	0.038725	0.042803
Bi_RNN	0.037570	0.038244	0.037586	0.041467	0.044881	0.042510	0.044928	0.040119	0.042524	0.043299
CNN1D_LSTM	0.038234	0.043631	0.042643	0.044788	0.038907	0.046889	0.041832	0.051089	0.047142	0.052802
ANN	0.057103	0.060965	0.065537	0.061544	0.059535	0.061573	0.060271	0.059627	0.057873	0.068573

**Table 5.** 10 step Test\_MAPE data of 15 models participating in the experiment.

Model	STEP1	STEP2	STEP3	STEP4	STEP5	STEP6	STEP7	STEP8	STEP9	STEP10
MGGTSP	0.010664	0.011856	0.012908	0.013898	0.014153	0.015210	0.015928	0.016728	0.017639	0.018326
DARNN	0.012742	0.013492	0.015195	0.016444	0.015651	0.017497	0.017647	0.018523	0.020128	0.020286
Autoformer	0.013061	0.013875	0.016077	0.017557	0.016118	0.018489	0.018148	0.019151	0.021127	0.021117
TCN	0.013990	0.014770	0.017531	0.019441	0.017128	0.020177	0.019164	0.020293	0.022875	0.022614
BiTransfomer _LSTM	0.019778	0.022580	0.022051	0.023331	0.025092	0.017770	0.026130	0.024649	0.024767	0.025097
Informer	0.017983	0.018345	0.022511	0.026143	0.021399	0.026094	0.023321	0.024085	0.028329	0.028060
LSTM	0.025587	0.026832	0.027559	0.030017	0.028598	0.026056	0.030931	0.034351	0.033116	0.032278
Bi_GRU	0.028922	0.034812	0.027270	0.028817	0.033980	0.042075	0.024814	0.040054	0.042387	0.030106
Bi_LSTM	0.032075	0.022765	0.029086	0.038758	0.035504	0.037625	0.041396	0.031411	0.036513	0.033838
GRU	0.034505	0.034609	0.033364	0.031164	0.042297	0.037924	0.037472	0.041016	0.035105	0.039846
RNN	0.033020	0.035437	0.035304	0.037536	0.037904	0.038326	0.039871	0.038640	0.038453	0.038941
CNN1D	0.039474	0.038810	0.032060	0.042228	0.046183	0.040666	0.045155	0.039777	0.046488	0.051180
Bi_RNN	0.043991	0.044804	0.044306	0.048538	0.052553	0.049851	0.052712	0.047334	0.050156	0.051010
CNN1D_LSTM	0.044628	0.051063	0.050001	0.052544	0.045748	0.055173	0.049362	0.060271	0.055763	0.062453
ANN	0.069636	0.074062	0.079720	0.074864	0.072489	0.074870	0.073280	0.072609	0.070289	0.082984

#### 3.3. Robustness Against Baseline Models

MGGTSP-CAT consistently outperforms both classical and state-of-the-art architectures. For instance, the Test\_R² gap between MGGTSP-CAT and the second-best model (DARNN) widens from 0.7% at Step 1 to 1.4% at Step 10 (Table 2). Similarly, error metrics such as Test\_MAE (0.01446 at Step 10) remain 8.5% lower than DARNN (0.01581), underscoring the efficacy of the multi-scale fusion mechanism in mitigating error propagation. The model's advantage is particularly pronounced in high-variability scenarios, as evidenced by its stable MAPE trajectory (1.07%  $\rightarrow$  1.42% over ten steps), contrasting sharply with the erratic performance of Transformer-based models like BiTransformer\_LSTM. This robustness is critical for real-world applications, where models must perform reliably under diverse and often unpredictable conditions. Additionally, MGGTSP-CAT's computational efficiency is a notable advantage, as it achieves these results with lower computational overhead compared to Transformer-based models, making it more feasible for deployment in resource-constrained environments. The predicted values and true values of the nine best models participating in the experiment for Step 1, Step 2, Step 5, and Step 10 are shown in Figure 3.

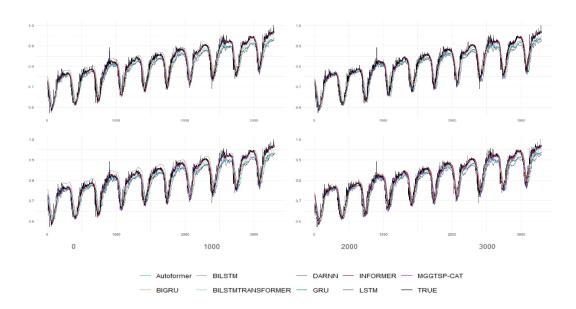


Figure 3. The predicted values and true values of the nine best models.

#### 4. Discussions

The MGGTSP-CAT framework introduces a structured approach to multi-scale greenhouse gas (GHG) concentration forecasting by integrating daily and monthly temporal patterns through a dual-encoder architecture. This design addresses a critical challenge in existing models, which often struggle to reconcile short-term variability with long-term trends. By employing an Input Attention encoder for high-frequency daily data and an Autoformer encoder for low-frequency monthly data, the model effectively captures both transient weather-driven fluctuations (e.g., synoptic-scale atmospheric changes) and persistent seasonal or policy-driven trends (e.g., annual emission cycles). The experimental results validate this approach, with MGGTSP-CAT achieving a Test\_ $R^2$  of 0.9627, outperforming baseline models such as DARNN (Test\_ $R^2$  = 0.9519) and Autoformer (Test\_ $R^2$  = 0.9480).

A notable strength of the model lies in its stability during multi-step forecasting. Over a ten-step horizon, the Test\_R² declines by only 3.3% (from 0.9788 to 0.9470), compared to sharper deteriorations in Transformer-based models like BiTransformer\_LSTM ( $\Delta$  = 5.7%) and traditional sequential architectures like LSTM ( $\Delta$  = 12.7%). This robustness is attributed to the Temporal Attention mechanism, which dynamically weights historical states to minimize error propagation—a common issue in long-term prediction tasks. For instance, the Test\_MSE increases by a factor of 2.5 from Step

1 to Step 10 (0.000163  $\rightarrow$  0.000410), a slower escalation than Autoformer's 2.3-fold rise (0.000236  $\rightarrow$  0.000554), despite the latter's reliance on self-attention for global dependencies.

The framework's computational efficiency further enhances its practicality. By decoupling daily and monthly processing, MGGTSP-CAT reduces computational overhead compared to monolithic Transformer architectures, which require quadratic memory for self-attention over long sequences. This efficiency is critical for real-world deployment, particularly in regions with limited computational resources. Additionally, the model's ability to handle multi-modal temporal features—such as isolating short-term noise from long-term signals—proves advantageous in complex scenarios like Arctic monitoring, where rapid temperature shifts and multi-decadal warming trends coexist.

While the current evaluation focuses on  $CO_2$  data from well-instrumented stations, the model's architecture is extensible to other GHGs (e.g.,  $CH_4$ ,  $N_2O$ ) and heterogeneous data sources, such as satellite observations or urban sensor networks. Future work could explore integrating exogenous variables (e.g., economic activity indices, land-use changes) to enhance scenario-based forecasting for policymakers. Nevertheless, the framework's reliance on continuous high-resolution input data highlights the importance of maintaining robust observational infrastructure, particularly in underrepresented regions.

# 5. Conclusions

This study presents an innovative multi-attention encoder model based on cross-time resolution, which combines daily and monthly data at different time scales and processes them through specific encoder architectures. Specifically, the model employs the Input Attention Mechanism to process daily data and the Autoformer Encoder to handle monthly data. Subsequently, the features extracted from these different time resolutions are concatenated and decoded through the Temporal Mechanism to generate the final prediction results. This approach provides a more accurate and robust solution for cross-scale forecasting of greenhouse gas concentrations and offers significant support for more complex climate models.

The contributions of this study are as follows:

- 1) A greenhouse gas dataset covering multiple global major climate monitoring stations was created. The data is sourced from long-term monitoring stations in Mauna Loa (Hawaii), Barrow (Alaska), American Samoa, and Antarctica, spanning over half a century [22]. This dataset not only offers high temporal resolution (daily data) but also effectively reflects global climate change trends, providing a solid experimental foundation for cross-scale climate forecasting.
- 2) An innovative fusion of the model's time resolution with the characteristics of greenhouse gas data. This study designs a multi-encoder fusion architecture that integrates the characteristics of different time resolution data. The Input Attention Mechanism and Autoformer Encoder are used to extract features from daily and monthly data, respectively, while the Temporal Attention Mechanism further enhances the model's ability to integrate information across different time scales. This method effectively captures the multi-scale features of greenhouse gas concentration changes, improving the model's adaptability and accuracy in both short-term and long-term climate predictions.
- 3) Exceptional accuracy and stability of the model. The experimental results show that the proposed model exhibits outstanding performance in prediction tasks across multiple climate monitoring stations, especially in high-variability daily data. The model can accurately capture key change patterns and provide stable and reliable predictions. For both short-term and long-term predictions at different time scales, the model performs excellently, showing less fluctuation in prediction results and significantly improving stability compared to other existing methods.
- 4) Future research could further extend the application of the model by considering more types of greenhouse gas monitoring data and a wider range of geographical areas. The effectiveness and generalization ability of the model under different climate conditions could be explored. Additionally, the model could be applied to the prediction of various climate events, further enhancing its predictive accuracy and practical application value.

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