

Learning Across Scales: A Physics-Informed Approach to Climate Modeling

Arantxa Vicario

Universidad de Los Lagos

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Abstract

Climate modeling plays a pivotal role in understanding Earth’s complex systems, but traditional methods struggle with computational demands across spatial and temporal scales. Machine learning (ML) offers a promising alternative, yet purely data-driven approaches often lack physical consistency. To address this, we propose a physics-informed approach to learning across scales in climate modeling. Our framework integrates physics-informed neural networks (PINNs) with hierarchical representations to model multiscale processes efficiently. We demonstrate improved accuracy and efficiency on benchmark climate datasets, paving the way for more reliable predictions of complex climate phenomena. These findings underscore the potential of combining ML with domain knowledge to advance climate science.

1 Introduction

Climate modeling plays a pivotal role in understanding and predicting the Earth’s complex systems, from localized weather patterns to global climate trends. However, these systems are inherently multiscale, encompassing interactions across spatial scales (e.g., microscale turbulence to planetary-scale circulations) and temporal scales (e.g., seconds to centuries). Traditional numerical methods, while powerful, often struggle with the computational demands of resolving fine-scale details while maintaining accuracy at larger scales. This limitation has spurred growing interest in leveraging machine learning (ML) techniques to bridge these scales efficiently [14].

Recent advances in ML have demonstrated significant potential for integrating and accelerating multiscale processes. However, purely data-driven approaches often lack interpretability and fail to incorporate the rich physical principles that govern climate systems [1]. To address this gap, we propose a physics-informed approach to learning across scales in climate modeling. By combining the strengths of ML with domain-specific knowledge encoded in partial differential equations (PDEs), our method ensures both predictive accuracy and physical consistency [13].

In this work, we introduce a novel framework that integrates physics-informed neural networks (PINNs) with hierarchical representations to model climate processes across multiple scales. Our contributions include:

1. A scalable architecture that learns interactions between fine- and coarse-grained representations.
2. Demonstrations of improved accuracy and efficiency on benchmark climate datasets.
3. Insights into the interpretability and generalizability of our approach, paving the way for broader applications in Earth system science.

Through this work, we aim to advance the state-of-the-art in climate modeling by bridging the gap between data-driven and physics-based paradigms, enabling more accurate and efficient predictions of complex multiscale phenomena.

2 Related Work

2.1 Machine Learning for Climate Modeling

The application of machine learning to climate science has gained significant traction in recent years. For instance, Rasp et al. [14] introduced a deep learning framework for emulating atmospheric models, demonstrating the potential of neural networks to replace computationally expensive components of climate simulations. Similarly, Brenowitz and Bretherton [4] developed a convolutional neural network (CNN) to parameterize subgrid-scale processes in global climate models, achieving notable improvements in accuracy and speed. More recently, Chattopadhyay et al. [5] employed recurrent neural networks (RNNs) to predict chaotic weather patterns, highlighting the ability of ML to capture nonlinear dynamics in climate systems. While these approaches excel at capturing statistical patterns, they often lack explicit constraints from physical laws, which can lead to unphysical predictions [1].

2.2 Physics-Informed Machine Learning

To address this limitation, researchers have explored physics-informed machine learning, which incorporates physical principles into ML models. Raissi et al. [13] pioneered the use of physics-informed neural networks (PINNs) by embedding PDEs as constraints during training. PINNs have since been applied to various domains, including fluid dynamics [7] and material science [10]. In climate science, Beucler et al. [1] combined PINNs with thermodynamic constraints to model cloud processes, ensuring physically consistent predictions. These works underscore the importance of integrating domain knowledge into ML frameworks to enhance their reliability and interpretability.

2.3 Multiscale Modeling in Climate Science

Multiscale modeling remains a central challenge in climate science due to the wide range of interacting processes. Traditional approaches rely on scale separation or averaging techniques, such as Reynolds averaging in turbulence modeling [12]. However, these methods often oversimplify fine-scale interactions, leading to inaccuracies at larger scales. Recent efforts have focused on bridging scales using hierarchical models. For example, Grooms et al. [6] proposed a stochastic superparameterization framework to represent unresolved scales in ocean models. Similarly, Bolton and Zanna [2] used a multiscale CNN to learn coarse-grained representations of ocean currents, achieving better alignment with high-resolution simulations. These studies highlight the potential of hierarchical approaches to capture cross-scale interactions effectively.

2.4 Graph Neural Networks and Attention Mechanisms

Graph neural networks (GNNs) and attention mechanisms have emerged as promising tools for modeling structured and multiscale data. Li et al. [8] introduced GNNs for simulating molecular dynamics, demonstrating their ability to encode relationships between particles at different scales. Building on this, Pfaff et al. [11] applied GNNs to fluid dynamics, achieving

state-of-the-art performance in predicting flow fields. Attention mechanisms, on the other hand, have been used to identify critical interactions in complex systems. For instance, Vaswani et al. [15] showed how self-attention can be used to weigh the importance of different inputs in sequence modeling tasks. These innovations provide a foundation for developing flexible and interpretable models for multiscale climate processes. The GNNs needs to be optimized using large scale optimization models like [9].

3 Methodology

Our framework integrates **physics-informed neural networks (PINNs)** with **hierarchical representations** to model multiscale climate processes efficiently. The key components of our approach include a hierarchical graph structure, a physics-informed loss function, and an attention mechanism. Below, we detail each component mathematically and explain their roles in learning across scales.

3.1 Hierarchical Representation

To capture interactions across multiple scales, we represent the system as a hierarchical graph $G = (V, E)$, where:

- V denotes the set of nodes representing physical entities (e.g., grid points or particles).
- E denotes the set of edges encoding interactions between nodes.

Each node $v_i \in V$ is associated with a feature vector $\mathbf{x}_i \in \mathbb{R}^d$, which encodes information at a specific scale. For example, in climate modeling, \mathbf{x}_i could represent atmospheric variables such as temperature, pressure, or velocity at a given spatial resolution.

The hierarchical structure is defined by partitioning the graph into multiple levels L_1, L_2, \dots, L_k , where each level corresponds to a different scale. For instance:

- L_1 : Fine-scale representation (e.g., microscale turbulence).
- L_k : Coarse-scale representation (e.g., planetary-scale circulation).

Nodes at adjacent levels are connected via aggregation functions. Let $\mathcal{N}(v_i)$ denote the neighborhood of node v_i . The feature update rule for node v_i at level l is given by:

$$\mathbf{h}_i^{(l)} = \sigma \left(W^{(l)} \cdot \text{AGG} \left(\{\mathbf{h}_j^{(l-1)}\}_{j \in \mathcal{N}(v_i)} \right) + b^{(l)} \right),$$

where:

- $\mathbf{h}_i^{(l)}$ is the updated feature vector of node v_i at level l .
- $\text{AGG}(\cdot)$ is an aggregation function (e.g., mean or max pooling) that combines features from neighboring nodes.
- $W^{(l)}$ and $b^{(l)}$ are learnable parameters at level l .
- $\sigma(\cdot)$ is a nonlinear activation function (e.g., ReLU).

This hierarchical representation allows us to encode multiscale interactions explicitly while maintaining computational efficiency.

3.2 Physics-Informed Loss Function

To ensure physical consistency, we incorporate governing partial differential equations (PDEs) into the training process. Let $\mathcal{L}_{\text{data}}$ denote the data-driven loss term, which measures the

discrepancy between predicted outputs $\hat{\mathbf{y}}$ and ground truth labels \mathbf{y} :

$$\mathcal{L}_{\text{data}} = \frac{1}{N} \sum_{i=1}^N \|\hat{\mathbf{y}}_i - \mathbf{y}_i\|^2,$$

where N is the number of samples in the dataset.

Additionally, let $\mathcal{L}_{\text{physics}}$ denote the physics-based loss term, which enforces constraints derived from PDEs. For example, consider the Navier-Stokes equation for fluid dynamics:

$$\frac{\partial \mathbf{u}}{\partial t} + (\mathbf{u} \cdot \nabla) \mathbf{u} = -\frac{1}{\rho} \nabla p + \nu \nabla^2 \mathbf{u} + \mathbf{f},$$

where:

- \mathbf{u} is the velocity field.
- p is the pressure field.
- ρ is the density.
- ν is the kinematic viscosity.
- \mathbf{f} represents external forces.

We approximate the residuals of the PDE using automatic differentiation. The physics-based loss term is then defined as:

$$\mathcal{L}_{\text{physics}} = \frac{1}{M} \sum_{j=1}^M \|\mathcal{R}(\hat{\mathbf{u}}_j, \hat{p}_j)\|^2,$$

where $\mathcal{R}(\cdot)$ represents the residual of the PDE, and M is the number of collocation points used to enforce the physics constraints.

The total loss function is a weighted combination of the data-driven and physics-based terms:

$$\mathcal{L}_{\text{total}} = \alpha \mathcal{L}_{\text{data}} + \beta \mathcal{L}_{\text{physics}},$$

where α and β are hyperparameters controlling the trade-off between the two terms.

3.3 Attention Mechanism

To prioritize critical interactions at each scale, we introduce an attention mechanism. Specifically, we compute attention weights a_{ij} for each edge $(v_i, v_j) \in E$ using a softmax function:

$$a_{ij} = \frac{\exp\left(\text{LeakyReLU}\left(\mathbf{w}^\top [\mathbf{h}_i^{(l)} \parallel \mathbf{h}_j^{(l)}]\right)\right)}{\sum_{k \in \mathcal{N}(v_i)} \exp\left(\text{LeakyReLU}\left(\mathbf{w}^\top [\mathbf{h}_i^{(l)} \parallel \mathbf{h}_k^{(l)}]\right)\right)},$$

where:

- \mathbf{w} is a learnable weight vector.
- $[\mathbf{h}_i^{(l)} \parallel \mathbf{h}_j^{(l)}]$ denotes the concatenation of feature vectors.

The aggregated feature vector for node v_i is then computed as:

$$\mathbf{z}_i^{(l)} = \sum_{j \in \mathcal{N}(v_i)} a_{ij} \mathbf{h}_j^{(l)}.$$

This attention mechanism enables the model to focus on the most relevant interactions at each scale, improving both accuracy and interpretability.

3.4 Training Process

The training process alternates between optimizing the data-driven loss $\mathcal{L}_{\text{data}}$ and enforcing the physics-based constraints $\mathcal{L}_{\text{physics}}$. At each iteration, we update the model parameters θ using gradient descent:

$$\theta \leftarrow \theta - \eta \nabla_{\theta} \mathcal{L}_{\text{total}},$$

where η is the learning rate.

By combining hierarchical representations, physics-informed constraints, and attention mechanisms, our framework achieves both scalability and interpretability, making it well-suited for multiscale climate modeling. Such parameter optimization has been proved to be useful in [3].

4 Experiments and Results

We evaluate our framework on two benchmark tasks:

1. Predicting turbulent flows in atmospheric models.
2. Simulating ocean currents at multiple resolutions.

Table 1 and Table 2 summarize the performance of our method compared to state-of-the-art baselines. Metrics include mean absolute error (MAE), root mean squared error (RMSE), and computational runtime.

Table 1: Performance on Turbulent Flow Prediction

| Method | MAE ($\times 10^{-3}$) | RMSE ($\times 10^{-3}$) | Runtime (s) |
|---------------|--------------------------|---------------------------|-------------|
| Baseline CNN | 5.2 | 7.8 | 120 |
| Baseline PINN | 4.1 | 6.5 | 200 |
| Our Method | 2.8 | 4.9 | 90 |

Table 2: Performance on Ocean Current Simulation

| Method | MAE ($\times 10^{-3}$) | RMSE ($\times 10^{-3}$) | Runtime (s) |
|---------------|--------------------------|---------------------------|-------------|
| Baseline CNN | 6.7 | 9.4 | 150 |
| Baseline PINN | 5.3 | 7.8 | 250 |
| Our Method | 3.4 | 5.6 | 110 |

Our results demonstrate superior performance compared to state-of-the-art baselines, with up to 30% improvement in accuracy while reducing computational costs by 50%.

5 Discussion

Our work highlights the potential of physics-informed learning to advance climate modeling. By integrating hierarchical representations and attention mechanisms, we achieve both scalability and interpretability. Future work will explore applications to other Earth system processes, such as land-atmosphere interactions and biogeochemical cycles.

6 Conclusion

We presented a novel framework for learning across scales in climate modeling, combining physics-informed neural networks with hierarchical representations. Our approach achieves state-of-the-art performance on benchmark tasks, offering a promising direction for advancing multiscale climate science.

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