This work presents a hybrid modeling approach to data-driven learning and representation of unknown physical processes and closure parameterizations. These hybrid models are suitable for situations where the mechanistic description of dynamics of some variables is unknown, but reasonably accurate observational data can be obtained for the evolution of the state of the system. In this work, we propose machine learning to account for missing physics and then data assimilation to correct the prediction. In particular, we devise an effective methodology based on a recurrent neural network to model the unknown dynamics. A long short-term memory (LSTM) based correction term is added to the predictive model in order to take into account hidden physics. Since LSTM introduces a black-box approach for the unknown part of the model, we investigate whether the proposed hybrid neural-physical model can be further corrected through a sequential data assimilation step. We apply this framework to the weakly nonlinear Lorenz model that displays quasiperiodic oscillations, the highly nonlinear chaotic Lorenz model, and two-scale Lorenz model. The hybrid neural-physics model yields accurate results for the weakly nonlinear Lorenz model with the predicted state close to the true Lorenz model trajectory. For the highly nonlinear Lorenz model and the two-scale Lorenz model, the hybrid neural-physics model deviates from the true state due to the accumulation of prediction error from one time step to the next time step. The ensemble Kalman filter approach takes into account the prediction error and updates the diverged prediction using available observations in order to provide a more accurate state estimate for the highly nonlinear chaotic Lorenz model and two-scale Lorenz system. The successful synergistic integration of neural network and data assimilation for low-dimensional system shows the potential benefits of the proposed hybrid-neural physics model for complex dynamical systems.

**Keywords** Chaotic systems · Deep learning · Neural networks · Data assimilation · Hybrid modeling
1 Introduction

The advancement in high-performance computing has made numerical simulation of atmosphere and climate models on computational grid consisting of $O(10^7)$ cells, spaced $O(10 \text{ km}) - O(100 \text{ km})$ possible [1]. However, the geophysical flows are governed by scales smaller than the mesh size of the climate models, and resolving these scales is essential for accurate prediction. For example, the dynamical scales of boundary layer clouds are $O(10 \text{ m})$ [2], and the submesoscale dynamics of oceans have length scales of $O(100 \text{ m})$ [3], and will remain unresolvable in the near future [4]. Therefore, the physics in climate models is essentially incomplete and the dynamics of smaller scales in the atmosphere and oceans must be represented through some parameterization schemes. Another example of an incomplete dynamical system is a hybrid coupled model where the physical equations are replaced by empirical relations to achieve a large computational speedup and greater stability. One example of a hybrid coupled model is the coupling of the statistical atmospheric model with a fully nonlinear ocean general circulation model, where the wind stress is empirically estimated from the ocean model variables [5].

The dynamics of the missing physics in geophysical flows is inherently nonlinear and replacing it with linear statistical methods will not deliver accurate prediction for longer periods. The ‘universal approximator theorem’ [6] of the neural network makes them an attractive choice for learning nonlinear dynamics in incomplete dynamical systems. Some of the early use of the neural network for dynamical systems are turbulence control for drag reduction [7], prediction of wind stress field from the ocean state in a hybrid coupled atmosphere-ocean model [8], modeling missing dynamics of the Lorenz system [9,10], and modeling parameterization in climate models [11]. All these early works have utilized a simple feed-forward neural network with a single hidden layer. In the past decade, ‘deep learning’ has made rapid progress in diverse areas such as speech translation [12], classification of images [13], and playing the game of Go [14]. These applications can require a neural network with a large number of layers, special types of convolutional layers, and specialized architectures, such as generative adversarial networks [15] along with the vast amount of data for the training. The machine learning (ML) methods can also be used to construct data-driven models of geophysical flows from the extensive data collected from satellites, in-situ scientific measurements, and futuristic internet-of-things (IoT) devices. Indeed, many researchers have explored the use of ML methods for earth science [1,16–18].

One of the main challenges with ML includes the difficulty of incorporating existing domain knowledge and handling of uncertainty [16,19,20]. The geophysical community utilizes the data assimilation (DA) framework to make use of physical laws and has robust ways of handling uncertainty in all parts of the problem [21]. The DA framework is also capable of handling sparse and irregularly distributed observations gathered through satellite measurements [22]. On the other hand, DA has mainly been employed for state estimation, such as the initial state of the system for the weather forecast. The forecast model can generally assumed to be perfect. However, the weather and climate models work on grid size around 10 km upwards, but the processes such as turbulence, radiation, and precipitation can take place at a scale around 1 km or less. Hence, the parameterization scheme for these unresolved processes should take into account the average impact at the model grid scale. The ML can be applied to discover accurate parameterization schemes using the data from high resolution models [23,25]. The ML has also been utilized to replace the complete physical model of chaotic systems [26,27]. These data-driven models are computationally cheap and can be integrated within the DA framework for state estimation and forecasting [28].

The DA and ML share a lot of similarities as both approaches use the data to learn about the system using inverse methods [29,31] and the synergistic integration of the two is essential for learning improved models of the earth system [32,36]. One way to integrate the DA and ML is to build a hybrid neural-dynamical model via variational DA [9]. This approach involves replacing the missing dynamics of the dynamical system with the neural network, and the parameters of the neural network were determined by the 4DVAR assimilation approach. The variational DA methods are commonly used to estimate the initial condition or model parameters, and are also implemented on operational weather and climate models [37,40]. Particularly, the backpropagation algorithm used to train the neural network has a strong mathematical equivalence with the adjoint method for calculating gradients in variational DA [29], and this parallelism was utilized to determine parameters of the neural network to build a hybrid neural-dynamical model [9]. Recently, the sequential data assimilation technique was applied to estimate the full state of the system using the truncated model (i.e., without accounting for unresolved processes), and then ML was utilized to learn the effect of unresolved part [41]. The work by Brajard et al. [41] goes beyond high-resolution simulations and they develop ML-based parameterization using direct data in the realistic scenario of sparse and noisy observations. The combination of ML-based parameterization to the physical core produces a hybrid model [42]. The ML can be applied to learn the closure term accounting for the effect of subgrid-scale processes and the hybrid model can be corrected using sequential data assimilation in the online deployment phase [43,45]. Also, DA and ML can be applied iteratively to build a fully data-driven model from sparse and noisy observations [46]. The fully data-driven model can then be used for state estimation and forecasting.
In this work, a hybrid modeling approach is presented with a data-driven representation for unknown physical processes and sequential data assimilation to correct the modeling error to achieve accurate analysis for chaotic dynamical systems. The paper is structured as follows. In Section 2, the baseline Lorenz model, the Lorenz model with incomplete dynamics, and the two-scale Lorenz model with data-driven parameterization is presented along with the working of long short-term memory (LSTM) neural network. Section 3 details the formulation of sequential data assimilation problem and lists the deterministic ensemble Kalman filter (DEnKF) algorithm. Section 4 discusses the results of numerical experiments. In particular, to test the proposed approach, we consider the following joint experiments: (i) incomplete physics-based model with LSTM for missing physics, (ii) incomplete physics-based model with LSTM for missing physics and DEnKF for data assimilation, and (iii) incomplete physics-based model and DEnKF for data assimilation. Finally, the summary of this work and concluding remarks are provided in Section 5. Python codes for this work are available on GitHub [47].

2 Hybrid neural-physics modeling

2.1 Lorenz model

We consider the following Lorenz system which is used as a simplified model for atmospheric convection [48]. The nonlinear system of three ordinary differential equations is

\[
\begin{align*}
\frac{dX}{dt} &= \sigma(Y - X), \\
\frac{dY}{dt} &= X(\rho - Z) - Y, \\
\frac{dZ}{dt} &= XY - \beta Z,
\end{align*}
\]

where \(X\), \(Y\), and \(Z\) are the state of the Lorenz system, and \(\sigma, \rho, \beta\) are the system’s parameter. In the above equations \(X\) is proportional to the intensity of convection motion, \(Y\) is proportional to the temperature difference in the horizontal direction, and \(Z\) is proportional to the distortion of the temperature profile in the vertical direction. In the vector notation, the state of the Lorenz system can be defined as

\[
\mathbf{x} = [X, Y, Z]
\]

and the Equations (1-3) can be written as

\[
\dot{\mathbf{x}} = \mathbf{f}(t, \mathbf{x}).
\]

The Lorenz system is integrated using the third-order Adams-Bashforth method [49] as follows

\[
\mathbf{x}^{(k+1)} = \mathbf{x}^{(k)} + \frac{\Delta t}{12}[23\mathbf{f}^{(k-1)} - 16\mathbf{f}^{(k-2)} + 5\mathbf{f}^{(k-3)}],
\]

where \(\mathbf{f} = [\sigma(Y - X), X(\rho - Z) - Y, XY - \beta Z]\), and \(\Delta t\) is the time step size. The Lorenz system is very sensitive to the initial condition as well as the choice of parameters. Based on the selection of the parameters, the Lorenz system can exhibit either chaotic behavior or the transient chaotic behavior [50]. Following the previous study [9], we investigate two cases here to evaluate the performance of the neural network in learning the chaotic dynamics. The first case is called the weakly nonlinear case, with the parameters \(\sigma, \rho, \beta\) set to 10, 28, and 8/3, respectively. The initial condition for the weakly nonlinear case is \([-9.42, -9.43, 28.3]\) for \([X(0), Y(0), Z(0)]\) [51]. The Lorenz system displays near-regular oscillations with a gradually increasing amplitude in the devised integration period for the weakly nonlinear case. The initial condition for the highly nonlinear case is set to be \([22.8, 35.7, 114.9]\), and the parameters \(\sigma, \rho, \beta\) are set to 16.0, 120.1, and 4.0, respectively [52]. The dataset for training the neural network is generated with these reference solutions and is regarded as the ’true’ solution of the Lorenz equations. The Lorenz attractor for the weakly and highly nonlinear case is shown in Figure 1.

Next, we assume that the complete dynamics of the Lorenz model is unknown. Specifically, we assume that the evolution equation for the distortion of the temperature profile in the vertical direction, i.e., \(Z\) is not available and must be approximated. The hybrid neural-physics Lorenz model composed of the physics-based equations for the known dynamics and the neural network to approximate the unknown dynamics can be written as

\[
\begin{align*}
\frac{dX}{dt} &= \sigma(Y - X), \\
\frac{dY}{dt} &= X(\rho - Z) - Y, \\
\frac{dZ}{dt} &= N(x(t), \cdots, x(t - l\Delta t)),
\end{align*}
\]
where $l$ is the number of previous time steps used for the forecast of the variable $Z$ at next time step, i.e., $t + \Delta t$.

In this study, we train the neural network using the teacher forcing method [53], which consists of using the ground truth data as the label for each training example. We assume that the data for the unknown state of the system (i.e., $Z$ variable in the Lorenz system) is either coming from measurements or from high-fidelity numerical simulations. The state derivative for the $Z$ variable at the $k$th time step is computed using the forward-difference scheme as follows

$$\frac{dZ}{dt} = \frac{Z^{(k+1)} - Z^{(k)}}{\Delta t}. \quad (9)$$

Any type of machine learning model can be utilized to learn the relationship between the known and unknown dynamics. Tang et al. [9] applied the feedforward neural network with a single hidden layer and five neurons to learn the missing physics in the Lorenz model. They observed that the neural network can approximate the unknown dynamics reasonably well for the weakly nonlinear case, but it fails for the highly nonlinear case. They proposed a better approach that makes use of the variational data assimilation where the dynamical constraints can be imposed for learning the parameters (i.e., weights and biases) of the neural network. Their proposed hybrid neural–dynamical variational data assimilation procedure leads to a very good prediction with almost the same forecast skill as the original Lorenz model. However, for the highly nonlinear case, the proposed method proposed a reasonable forecast only for few numerical experiments. This might have been due to the difficulty with the deterministic optimization algorithms within the variational assimilation in finding the global minima [54-55]. Even if the global optimum is found, the data in the assimilation window might be from one wing of the butterfly-shaped attractor, whereas the data in the forecast period lies on the second wing, thereby resulting in the poor forecast. This can be mitigated by applying stochastic data assimilation approaches like ensemble Kalman filter [55-57]. In this work, we revisit the problem of recovering missing physics in a dynamical system using the same examples as investigated by Tang et al. [9] with the application of recurrent neural network (RNN) integrated within the deterministic ensemble Kalman filter algorithm [58].

### 2.2 Two-scale Lorenz model

The two-scale Lorenz model is given by a following set of ordinary differential equations (ODEs):

$$\frac{dX_n}{dt} = -X_{n-1}(X_{n-2} - X_{n+1}) - X_n + F - \frac{hc}{b} \sum_{m=M(n-1)+1}^{nM} Y_m, \quad (10)$$

$$\frac{dY_m}{dt} = -c b Y_{m+1}(Y_{m+2} - Y_{m-1}) - c Y_m + \frac{hc}{b} X_{\lfloor (m-1)/M \rfloor + 1}, \quad (11)$$
where \( n = 1, \ldots, N, m = 1, \ldots, MN \), and \([\cdot]\) is the modulo operation. The \( X \) and \( Y \) variables are periodic, i.e., \( X_1 = X_N \) and \( Y_1 = Y_M N \). In the present study, the number of \( X \) variables is set at \( N = 8 \) and the number of \( Y \) variables per \( X \) variables is set to \( M = 32 \). We also set the coupling constant between \( X \) and \( Y \) variables to \( b = 1 \), the spatial-scale ratio to \( c = 10 \), and the temporal-scale ratio to \( c = 10 \). The forcing term is set at a large value of \( F = 20 \) to ensure chaotic behavior. These parameters are such that one model time unit (MTU) is approximately equivalent to five atmospheric days [59]. Recently, this model with the same parameter settings was applied to test the feasibility of generative adversarial networks to learn stochastic parameterization in multiscale systems [25].

The two-scale Lorenz model given by Equation [10] and Equation [11] is regarded as the truth model that must be simulated. In weather and climate modeling, the governing equations of motion of the system are known. However, it is not possible to resolve every small scale due to limited computational resources. Therefore, the effect of small scale variables are typically parameterized as a function of resolved variables. A forecast model for the two-scale Lorenz equations with the parameterization of the small \( Y \) scale variables on the resolved \( X \) variables can be written as

\[
\frac{d\tilde{X}_n}{dt} = -\tilde{X}_{n-1}(\tilde{X}_{n-2} - \tilde{X}_{n+1}) - \tilde{X}_n + F - G_n, \tag{12}
\]

where \( \tilde{X}_n \) is the forecast estimate of \( X_n \), and \( G_t \) represent the effect of unresolved variables. The parameterization is usually a function of the resolved variables, i.e., \( \tilde{X} \in \mathbb{R}^n \), and can be written mathematically as

\[
\frac{hc}{b} \sum_{m=N(n-1)+1}^{nM} Y_m \approx G_n = \mathcal{M}(\tilde{X}), \tag{13}
\]

where \( \mathcal{M}(\cdot) \) is the nonlinear mapping for representing the effect of unresolved scales on resolved scale variables. This mapping can be based on physical arguments or some prior information about subgrid scale processes. Alternatively, data-driven models can be used to exploit the data generated from experimental measurements or high-fidelity numerical simulations to learn the relation between resolved and unresolved scales. Indeed, in recent years, data-driven models have been successfully applied for parameterizing subgrid scale processes in complex geophysical flows [23, 25, 44, 60].

In the present study, we utilize the LSTM neural network to learn the subgrid scale parameterization of unresolved scales. Once we have the ground truth data available for large scale \( X \) variables, we can compute the \textit{true} subgrid scale parameterization as follows

\[
G^{(k)}_n = -X_{n-1}^{(k)}(X_{n-2}^{(k)} - X_{n+1}^{(k)}) - X_n^{(k)} + F - \left( \frac{X_n^{(k+1)} - X_n^{(k)}}{\Delta t} \right), \tag{14}
\]

where \( \Delta t \) is the time step of integration of the forecast model. The ground truth data for training the neural network is generated by temporally integrating the two-scale Lorenz system using the fourth-order Runge-Kutta (RK4) time stepping scheme.

### 2.3 Long short-term memory neural network

In this work, we apply the long short-term memory (LSTM) neural network to predict the missing physics (i.e., incomplete dynamics) from the resolved physics of the dynamical system. The motivation behind applying LSTM is due to its success in modeling high-dimensional spatio-temporal chaotic time series of physical systems [26, 27, 61, 62]. LSTM is a type of recurrent neural network (RNN) that can capture the long-term dependencies in the evolution of time series data [63]. RNNs contain loops that allow them to persist information from one time step to another and can be expressed as

\[
\begin{align*}
\mathbf{h}^{(t)} &= f_{h\rightarrow h}(\mathbf{o}^{(t)}, \mathbf{h}^{(t-1)}), \\
\tilde{\mathbf{o}}^{(t+1)} &= f_{h\rightarrow o}(\mathbf{h}^{(t)}),
\end{align*}
\tag{15}
\]

\[
\begin{align*}
\mathbf{h}^{(t)} &= f_{h\rightarrow h}(\mathbf{o}^{(t)}, \mathbf{h}^{(t-1)}), \\
\tilde{\mathbf{o}}^{(t+1)} &= f_{h\rightarrow o}(\mathbf{h}^{(t)}),
\end{align*}
\tag{16}
\]

where \( \mathbf{h}^{(t)} \in \mathbb{R}^{dh} \) is the hidden state at time \( t \), \( \mathbf{o}^{(t)} \in \mathbb{R}^{dh} \) is the input vector at time \( t \), \( f_{h\rightarrow h} \) is the hidden to hidden mapping, and \( f_{h\rightarrow o} \) is the hidden to output mapping. The output of the model is the forecast \( \tilde{\mathbf{o}}^{(t+1)} \) at time step \( t + 1 \).

One of the limitations of RNNs is vanishing (or exploding) gradient to capture the long-term dependencies. This stems from the fact that the gradient is multiplied with the weight matrix repetitively during backpropagation through time (BPTT). The LSTM mitigates the issue with vanishing (or exploding) gradient by employing the gating mechanism that allows information to be forgotten. The equations that implicitly define the mapping from hidden state from the previous time step (i.e., \( \mathbf{h}^{(t-1)} \)) and input vector at the current time step (i.e., \( \mathbf{o}^{(t)} \)) to the forecast hidden state (i.e., \( \mathbf{h}^{(t)} \)) take the following form.
Mathematically, the LSTM mapping for learning parameterizations in a two-scale Lorenz model can be represented as

\[ \mathbf{G}^{(l)} \in \mathbb{R}^{l \times N} \rightarrow \{ \mathbf{G}^{(k)} \} \in \mathbb{R}^N. \]  

where \( \mathbf{X} \) is the resolved flow variables and \( \mathbf{G} \) is true parameterization.
3 Sequential data assimilation

Data assimilation (DA) is a technique to incorporate sparse and irregularly distributed noisy measurements with the physical model of the system to achieve an accurate prediction of the state trajectory. DA is generally classified into two categories: variational and sequential DA. In variational DA, all observations over a particular time window are utilized to minimize the cost function with certain constraints to compute the model trajectory that best fits the observations. The variational DA is particularly well suited for re-analyses problems to obtain the best possible state of the system at some time \( t \) using observations before and after this time [71]. In sequential DA, the system’s state is evolved in time using background information (i.e., the physical model) until observations become available. At this instant, an update (correction) to the system’s state is determined and the solver is re-initialized with this analyzed state. This procedure is continued until new observations get available, and so on. There is a rich literature on DA [22–74] and here we discuss only sequential DA problem and then outline the algorithm procedure for the deterministic ensemble Kalman filter (DEnKF).

We consider the dynamical system whose evolution can be represented as

\[
\mathbf{x}^{(k+1)} = \mathbf{M}_{t_k \rightarrow t_{k+1}}(\mathbf{x}^{(k)}) + \mathbf{w}^{(k+1)},
\]

where \( \mathbf{x}^{(k)} \in \mathbb{R}^n \) is the state of the system at discrete time \( t_k \), and \( \mathbf{M} : \mathbb{R}^n \rightarrow \mathbb{R}^n \) is the nonlinear model operator that defines the evolution of the system. The term \( \mathbf{w}^{(k+1)} \) denotes the model error that takes into account any type of uncertainty in the model that can be attributed to boundary conditions, imperfect models, etc. Let \( \mathbf{z}^{(k)} \in \mathbb{R}^m \) be observations of the state vector obtained from sparse, noisy measurements and can be written as

\[
\mathbf{z}^{(k+1)} = \mathbf{q}(\mathbf{x}^{(k+1)}) + \mathbf{v}^{(k+1)},
\]

where \( \mathbf{q}() \) is a nonlinear function that maps \( \mathbb{R}^n \rightarrow \mathbb{R}^m \), and \( \mathbf{v}^{(k+1)} \in \mathbb{R}^m \) is the measurement noise. We assume that the measurement noise is a white Gaussian noise with zero mean and the covariance matrix \( \mathbf{R}^{(k+1)} \), i.e., \( \mathbf{v}^{(k+1)} \sim \mathcal{N}(0, \mathbf{R}^{(k+1)}) \). Additionally, the noise vectors \( \mathbf{w}^{(k+1)} \) and \( \mathbf{v}^{(k+1)} \) are assumed to be uncorrelated to each other at all time steps. The sequential DA can be considered as a problem of estimating the state \( \mathbf{x}^{(k+1)} \) of the system given the observations up to time \( t_{k+1} \), i.e., \( \mathbf{z}^{(1)}, \ldots, \mathbf{z}^{(k+1)} \). When we utilize observations to estimate the state of the system, we say that the data are assimilated into the model, and use the notation \( \hat{\mathbf{x}}^{(k+1)} \) to denote an analyzed state estimate of the system at time \( t_{k+1} \). When all the observations before (but not including) time \( t_{k+1} \) are applied for estimating the state of the system, then we call it the forecast estimate and denote it as \( \mathbf{x}^{(k+1)} \).

The ensemble Kalman filter (EnKF) [75] follows the Monte Carlo estimation method to approximate the covariance matrix in the Kalman filter equations [76]. Inflation and covariance localization approaches have been often used in the EnKF framework to mitigate small number of ensembles [77–80]. Instead of modeling the exact evolution of a probability density function under nonlinear dynamics, ensemble methods maintain an empirical approximation to the target distribution in the form of a set of ensemble members \( \mathbf{x}^{(k)}(i) \) for \( i = 1 \ldots N \). We begin by initializing the state of the system for different ensemble members \( \hat{\mathbf{x}}^{(0)}(i) \) drawn from the distribution \( \mathcal{N}(\mathbf{x}^{(0)}, \mathbf{P}^{(0)}) \), where \( \mathbf{x}^{(0)} \) represents the best-known state estimate at time \( t_0 \), and \( \mathbf{P}^{(0)} \) is the initial covariance error matrix.

The propagation of the state for each ensemble member over the time interval \( [t_k, t_{k+1}] \) can be written as

\[
\mathbf{x}^{(k+1)}(i) = \mathbf{M}_{t_k \rightarrow t_{k+1}}(\hat{\mathbf{x}}^{(k)}(i)) + \mathbf{w}^{(k+1)},
\]

where \( \mathbf{w}^{(k+1)} \) accounting for model imperfections is usually assumed to be Gaussian noise. The prior state and the covariance matrix are approximated using the sample mean and error covariance matrix \( \mathbf{P}^{(k+1)} \) as follows

\[
\mathbf{x}^{(k+1)} = \frac{1}{N} \sum_{i=1}^{N} \mathbf{x}^{(k+1)}(i),
\]

\[
\mathbf{A}^{(k+1)}(i) = \mathbf{x}^{(k+1)}(i) - \mathbf{x}^{(k+1)},
\]

\[
\mathbf{P}^{(k+1)} = \frac{1}{N-1} \sum_{i=1}^{N} \mathbf{A}^{(k+1)}(i)(\mathbf{A}^{(k+1)}(i))^T,
\]

where the superscript \( T \) denotes the transpose, and \( \mathbf{A}^{(k+1)}(i) \) is the anomalies between the forecast estimate for the \( i \)th ensemble and the sample mean. Once the observations gets available at time \( t_{k+1} \), the forecast state estimate is assimilated using the Kalman filter analysis equation as follows

\[
\hat{\mathbf{x}}^{(k+1)} = \mathbf{x}^{(k+1)} + \mathbf{K}^{(k+1)}(\mathbf{z}^{(k+1)} - \mathbf{q}(\mathbf{x}^{(k+1)})),
\]

where

\[
\mathbf{A}^{(k+1)}(i) = \mathbf{x}^{(k+1)}(i) - \mathbf{x}^{(k+1)}.
\]
In contrast to the ensemble Kalman filter algorithm, the DEnKF does not employ any perturbed observations. The Kalman gain matrix is computed using its square root version (without storing or computing $P^{(k+1)}$ explicitly) as follows

$$K^{(k+1)} = \frac{A_f^{(k+1)}(Q^{(k+1)}A_f^{(k+1)})}{N-1} \left[ \frac{(Q^{(k+1)}A_f^{(k+1)})(Q^{(k+1)}A_f^{(k+1)})}{N-1} + R^{(k+1)} \right]^{-1}, \quad (31)$$

where $Q \in \mathbb{R}^{m \times n}$ is the Jacobian of the observation operator (i.e., $Q_{kl} = \frac{\partial q}{\partial x_l}$), and the matrix $A_f^{(k+1)} \in \mathbb{R}^{n \times N}$ is concatenated as follows

$$A_f^{(k+1)} = [A_f^{(k+1)}(1), A_f^{(k+1)}(2), \ldots, A_f^{(k+1)}(N)]. \quad (32)$$

The anomalies for all ensemble members are then updated separately with half the Kalman gain as shown below

$$\tilde{A}^{(k+1)}(i) = A_f^{(k+1)}(i) - \frac{1}{2} K^{(k+1)} Q^{(k+1)} A_f^{(k+1)}(i). \quad (33)$$

The state for all ensemble members is updated by adding ensemble anomalies to analysis state estimate and can be written as

$$\hat{X}^{(k+1)}(i) = \tilde{X}^{(k+1)} + \lambda \cdot \tilde{A}^{(k+1)}(i), \quad (34)$$

where $\lambda$ is the inflation factor to account for modeling errors. The above ensembles are used as initial ensembles for the next assimilation cycle and the procedure is continued.

4 Results and discussions

The general structure of the proposed hybrid neural-physics modeling approach is illustrated in Figure 3. We stress here that for many physical systems the complete description for dynamics of the system is not available and the missing physics for such systems can be modeled using machine learning tools. This provides us with the hybrid neural-physics model. The forecast of the hybrid neural-physics model can be further corrected using the sparse and noisy observations through data assimilation (specifically, we utilize the DEnKF algorithm). We consider a series of numerical experiments: (i) incomplete physics-based model with LSTM for missing physics (without data assimilation), (ii) incomplete physics-based model with LSTM for missing physics and DEnKF for data assimilation, and (iii) incomplete physics-based model (without LSTM closure) and DEnKF for data assimilation. We emphasize that we apply machine learning to improve the model of the system (e.g., see Figure 3). Data assimilation is then used in the analysis step to improve the forecast of the system.

4.1 Lorenz model

4.1.1 Hybrid neural-physics model

The Lorenz model is integrated using the third-order Adams-Bashforth method with the time step $\Delta t = 1 \times 10^{-3}$ for 6000 time steps, i.e., $t = 6$. The LSTM network is trained using the data from time $t = 0$ to $t = 3$. For training the LSTM network, we apply the lookback $l = 6$ meaning that the state of the system for six previous time steps is used to predict the unknown dynamics (i.e., $dZ/dt$ in our case) at the next time step. We use the full state of the system $[X, Y, Z]$ as the input features to the LSTM network. During the deployment of the trained LSTM network, the variable $Z$ at the next time step is calculated using Equation 6 with the predicted $dZ/dt$. Since the LSTM network is deployed in an auto-regressive method, there is an accumulation of errors from one time step to another. The chaotic systems are very sensitive to the initial condition and this makes the accurate forecast of chaotic systems difficult over a longer period. Figure 4 displays the prediction of the Lorenz system using the true model and the hybrid neural-physics model for the weakly and highly nonlinear case. The hybrid neural-physics model can predict the correct dynamics of the $Z$ variable and the predicted state trajectory is very close to the true state of the system. However, for the highly nonlinear Lorenz system, the predicted state trajectory starts deviating from the true state trajectory at around $t \approx 1.5$. The largest Lyapunov exponent of the highly nonlinear Lorenz system is 2.33. Therefore, the time $t \approx 1.5$ corresponds to 3.5 Lyapunov time ($L_1^{-1}$, where $L_1$ is the largest Lyapunov exponent of the Lorenz system). Even though the hybrid neural-physics model can predict the switching between lobes, the predicted state trajectory is considerably different from the true state trajectory for the highly nonlinear case. Therefore, the online data should be integrated using sequential DA to achieve an accurate forecast by limiting the accumulation of prediction errors.
Figure 3: The general structure of the proposed hybrid neural-physics modeling approach. Machine learning is used to learn the missing dynamics or parameterization of unresolved processes and this provides us with the hybrid neural-physics model. The hybrid neural-physical model can be further corrected through the data assimilation step making use of sparse and noisy observations in the online deployment phase.

Figure 4: The trajectory of the Lorenz system computed using the true physics-based model and the hybrid model where the $dZ/dt$ is approximated using the neural network for the weakly nonlinear case (left) and the highly nonlinear case (right). Here, LSTM refers to forecast with the hybrid model.

4.1.2 Hybrid neural-physics model coupled with DA

Next, we show the performance of the DEnKF algorithm that uses the hybrid neural-physics model as the forward model and compare it with only using the hybrid neural-physics model. Figure 5 shows the trajectory of the weakly nonlinear Lorenz system where the observations are assimilated after every 50 time steps. The observations are obtained by adding the Gaussian noise with the variance $\sigma_b^2 = 1.0$ and we assume that the full state is observable. We initialize 10 different ensemble members by adding the Gaussian noise with the variance $\sigma_0^2 = 1.0$ to the true state of the system at time $t = 0$. From Figure 5, we can notice that the prediction with the hybrid neural-physics model integrated within
Figure 5: The trajectory of the weakly nonlinear Lorenz system computed using the true physics-based model, the hybrid model where the $dZ/dt$ is approximated using the neural network, and the assimilated results with the hybrid model. The observation noise is drawn from $\mathcal{N}(0, 1)$, and are collected after every 50 time steps, i.e., the time interval between two observations is 0.05. Here, LSTM refers to forecast with the hybrid model and LSTM-DEnKF refers to an analysis state of the system.

the DEnKF algorithm matches almost exactly with the true trajectory of the weakly nonlinear Lorenz system. This is to be expected as even the hybrid neural-physics model was able to predict the state trajectory close to the true trajectory of the weakly nonlinear Lorenz system and ensemble Kalman filter approaches are robust for sparse and noisy observations in improving the state estimation. In Figure 6, we compare the power spectral density for all three states of the true model, hybrid neural-physics model, and the hybrid neural-physics model coupled with DEnKF. The power spectral density (PSD) is calculated as the square of the absolute value of the discrete Fourier transform of the time series of interest and can be written as

$$\text{PSD} = |\hat{u}(\omega)|^2,$$

where $\hat{u}(\omega)$ is the Fourier transform of a time series. We observe some noise in the PSD for the $Y$ variable, especially at higher frequencies which can be attributed to the Gaussian noise of observations that are assimilated for the state estimation in the DEnKF algorithm. We further increase the time interval at which the observations are assimilated to 100 time steps. As shown in Figure 7, the hybrid neural-physics model integrated into the DEnKF algorithm is able to accurately estimate the full state of the weakly nonlinear Lorenz system from time $t = 0$ to $t = 6$ even with the large time interval between two observations. The largest Lyapunov exponent for the weakly nonlinear Lorenz system is 0.86 and the system eventually becomes chaotic if we integrate it for a long time. To demonstrate the capability of the proposed framework, we integrate the weakly nonlinear Lorenz system for a long lead time up to $t = 30$. The LSTM network is trained using the true trajectory up to $t = 15$. From Figure 8 we can see that the LSTM model diverges from the true trajectory after approximately 15 Lyapunov times. If we assimilate the observations through the DEnKF algorithm, then the analysis state is close to the true state trajectory.

After a successful demonstration of the DEnKF algorithm with the hybrid neural-physics model for the weakly nonlinear Lorenz system, we discuss the results for the highly nonlinear Lorenz system. Figure 9 depicts the state trajectory of the highly nonlinear Lorenz system for the true model, hybrid neural-physics model, and the hybrid neural-physics model coupled with the DEnKF. The observations for the highly nonlinear Lorenz system are contaminated using the Gaussian noise with variance $\sigma^2 = 5$ and are assimilated after every 50 time steps. Even though the LSTM is trained for time $t = 0$ to $t = 3$ the state predicted by the hybrid neural-physics model starts deviating from the true model trajectory at around $t \approx 1.5$. This is due to the accumulation of prediction error in the deployment of the trained LSTM network and high sensitivity to the initial condition in the strongly nonlinear Lorenz system. When the hybrid neural-physics model
Figure 6: Power spectra analysis of the weakly nonlinear Lorenz system computed using the true physics-based model, the hybrid model where the $\frac{dZ}{dt}$ is approximated using the neural network, and the assimilated results with the hybrid model. The observation noise is drawn from $\mathcal{N}(0, 1)$, and are collected after every 50 time steps, i.e., the time interval between two observations is 0.05. Here, LSTM refers to forecast with the hybrid model and LSTM-DEnKF refers to an analysis state of the system.

is coupled with the DEnKF algorithm, the predicted state is accurate beyond the training period up to time $t = 6$. This shows the success of the LSTM network to learn the mapping between resolved dynamics and the missing physics of the Lorenz system and also testifies the functioning of sequential DA. Figure 10 shows the PSD for all three states of the highly nonlinear Lorenz system for the true model, hybrid neural-physics model, and hybrid neural-physics model coupled with DEnKF. We can observe that there is a mismatch in the distribution of PSD at lower frequencies for the state predicted by the hybrid neural-physics model compared to the true model state trajectory. The PSD for the state trajectory predicted with the hybrid neural-physics model coupled with DEnKF agrees very well with the PSD of the true model state trajectory.

In the next set of numerical experiments, we study the effect of observation noise and the frequency of assimilation on the prediction of the state of the highly nonlinear Lorenz system. Figure 11 shows the prediction of the state trajectory of the highly nonlinear Lorenz system with observations assimilated after every 100 time steps. The observation noise is taken to be from the Gaussian distribution with variance $\sigma_b^2 = 5$ and the number of ensembles is set at 10. We get a very good prediction between the true state of the Lorenz system and the state predicted by the hybrid neural-physics model integrated with the DEnKF. We further increase the observation noise to the Gaussian distribution with the variance of $\sigma_b^2 = 10$ and assimilate observations every 50 time steps. We note here that the number of ensembles is kept fixed at 10. As shown in Figure 12, there is a discrepancy between the analyzed state estimate of the Lorenz system and the true state for the highly nonlinear case. This discrepancy can be due to the small number of ensembles in the DEnKF algorithm. In the DEnKF algorithm, the error covariance matrix is approximated by computing the statistics of all ensemble samples and ensemble size of 10 might not be enough for the large measurement noise. We repeat the numerical experiment with the same set of parameters (i.e., the observation noise and the frequency of observations) and increase the number of ensembles to 20. Figure 13 displays the analyzed state estimate using an ensemble of 20 members for the hybrid neural-physics model coupled with DEnKF. The analyzed state estimate is improved and there is a very good agreement between the true state of the highly nonlinear Lorenz system and the analyzed state estimate by the hybrid neural physics model coupled with DEnKF.

These findings illustrate the robustness of sequential DA in handling sparse and noisy observations to improve the state estimate of the hybrid neural-physics model. The underestimation of the error covariance matrix can also be
Figure 7: The trajectory of the weakly nonlinear Lorenz system computed using the true physics-based model, the hybrid model where the $dZ/dt$ is approximated using the neural network, and the assimilated results with the hybrid model. The observation noise is drawn from $N(0, 1)$, and are collected after every 100 time steps, i.e., the time interval between two observations is 0.1. Here, LSTM refers to forecast with the hybrid model and LSTM-DEnKF refers to an analysis state of the system.

handled using practical DA techniques like covariance inflation [81, 82] and covariance localization [83, 84]. These techniques also prevent the filter divergence when the number of ensembles is less [85, 86]. Indeed for high-dimensional geophysical systems, increasing the number of ensembles will be computationally expensive and adaptive inflation and localization techniques can be adopted [87–89]. The quantitative performance of all numerical experiments with sequential data assimilation for the hybrid neural-physics Lorenz model is evaluated using the root mean squared error (RMSE) defined as

$$RMSE = \sqrt{\frac{1}{n T} \sum_{i=1}^{n} \sum_{t=1}^{T} (x_T^i(t) - x_P^i(t))^2},$$

where $x_T^i$ is the true state of the system and $x_P^i$ is the predicted state of the system. Table 1 reports the RMSE for all numerical experiments carried out with a hybrid neural-physics model coupled with DA. We can notice that an increase in the time between two observations leads to poor prediction for both weak and strongly nonlinear Lorenz systems. We also observe that for the strongly nonlinear Lorenz system, the increase in observation noise deteriorates the prediction for an ensemble of 10 members. When we increase the number of ensemble members to 20, the prediction is improved and the RMSE is decreased.

4.2 Two-scale Lorenz 96 model

The ground truth data for learning the subgrid scale parameterization in a two-scale Lorenz system is generated by temporally integrating the model using the fourth-order Runge-Kutta (RK4) time-stepping scheme with the time step $\Delta t = 0.001$ MTU. The equilibrium initial condition for the slow variables is set as $X_n = 0$ for $n \in 2, \ldots, N$ and $X_1 = 1$. In a similar manner, the fast variables are assigned $Y_1 = 1.0$ and $Y_m = 0$ for $m \in 2, \ldots, MN$ as an initial condition. The initial transient period of 5 MTU is disregarded. The periodic boundary condition is applied for both slow and fast variables in the two-scale Lorenz model. The output for the first 10 MTU is utilized for training and the performance of the trained network is evaluated for the next 10 MTU. Since the training and forecast period are different, there is no overlap between the train and test data. The trained LSTM network predicts the subgrid-scale parameterization at $k$th time step, i.e., $G^{(k)}$ using the temporal history of resolved flow variables for six consecutive
Figure 8: The trajectory of the weakly nonlinear Lorenz system for a long time computed using the true physics-based model, the hybrid model where the \( dZ/dt \) is approximated using the neural network, and the assimilated results with the hybrid model. The observation noise is drawn from \( N(0, 1) \), and are collected after every 100 time steps, i.e., the time interval between two observations is 0.1. Here, LSTM refers to forecast with the hybrid model and LSTM-DEnKF refers to an analysis state of the system.

Table 1: Quantitative assessment of different numerical experiments carried out with hybrid neural-physics model coupled with DA. The root mean square error (RMSE) is computed using Equation (36). Here, \( \sigma^2_b \) is the observation noise, \( N_s \) is the number of time steps between two observations and \( N \) is the number of ensembles used in DEnKF algorithm.

| Type | \( \sigma^2_b \) | \( N_s \) | \( N \) | RMSE  |
|------|-----------------|---------|------|-------|
| Weak | 1.0             | 50      | 10   | 0.169 |
| Weak | 1.0             | 100     | 10   | 0.233 |
| Strong | 5.0           | 50      | 10   | 2.417 |
| Strong | 5.0           | 100     | 10   | 4.804 |
| Strong | 10.0          | 50      | 10   | 15.385|
| Strong | 10.0          | 50      | 20   | 6.889 |

During the online deployment, we also use the DEnKF algorithm to improve the state prediction. The observation data for the DEnKF algorithm are generated by adding the Gaussian noise with the variance \( \sigma^2_b = 1.0 \) and the observations are assimilated every 10 time steps, i.e., 0.01 MTU. We assume that 50% of the full state is observable and the observation points correspond to \( X_1, X_3, X_5, \) and \( X_7 \). We initialize 10 different ensemble members by adding the Gaussian noise with the variance \( \sigma^2_0 = 1 \times 10^{-2} \) to the true initial state of the system. We also apply an inflation factor \( \lambda = 1.02 \) to account for modeling errors. Figure 14 displays the full-state trajectory of the slow variables in the two-scale Lorenz system from 10 MTU to 20 MTU (forecast period) along with the difference between the true and predicted state. The true state is calculated by solving both the evolution of slow and fast variables. We see that the LSTM based parameterization provides accurate prediction only for approximately 1 MTU and then the prediction has diverged from the true state of the two-scale Lorenz system. We also notice that the LSTM based subgrid scale parameterization model coupled with DEnKF renders accurate state prediction for the entire period of the forecast. In past time steps. We highlight here that we utilize the same time step for the forecast as the ground truth data generation. However, one can also use the different time step size for the forecast period by using the forecast time step size in Equation (14) to compute the parameterization term.
Figure 9: The trajectory of the strongly nonlinear Lorenz system computed using the true physics-based model, the hybrid model where the $dZ/dt$ is approximated using the neural network, and the assimilated results with the hybrid model. The observation noise is drawn from $\mathcal{N}(0, 5)$, and are collected after every 50 time steps, i.e., the time interval between two observations is 0.05. Here, LSTM refers to forecast with the hybrid model and LSTM-DEnKF refers to an analysis state of the system.

Figure 15, the trajectory of each slow variable is shown and we see that the LSTM-DEnKF framework can estimate the state for all slow variables with very high accuracy. The temporal variation of the RMSE also suggests that the sequential data assimilation aids in considerable improvement in the state prediction compared to employing only the trained LSTM network.

We highlight here that it may be argued that the forecast of a free run and an analysis (corrected forecast from DEnKF) of the system do not have access to the same information. Specifically, the LSTM-DEnKF model utilizes sparse and noisy observations that are not available to the LSTM model. For a fair comparison, we repeat the numerical experiments with the incomplete physics-based model (IPM) coupled with the DEnKF model. Figure 16 depicts the full state trajectory of the slow variables in the two-scale Lorenz system from 10 MTU to 20 MTU along with the difference between the true and predicted state. We can see that the analysis with the LSTM model is significantly more accurate than the analysis with the IPM model. This result demonstrates that machine learning offers possibilities of learning missing physics or model imperfections and the data assimilation prediction with such hybrid models is superior compared to truncated models.

As discussed in Section 2.2, the tendency (i.e., the time derivative) is approximated by a finite-difference method in the computation of subgrid scale parameterization. One of the important questions is the effect of this approximation on the generalizability of the trained neural network. We investigate this effect by deploying the trained neural network in a two-scale Lorenz 96 system with a different time step used during the deployment. A similar sensitivity study was also conducted in Brajard et al. [41], where they analyzed the effect of time step on the linear superposition assumption between the resolved and unresolved part of the model. Figure 17 shows the evolution of the RMSE for the two-scale Lorenz 96 model with a different time step used in the prediction and for different observation noise. In Figure 17, the $\Delta t_T$ denotes the time step used for generating the training data, and $\Delta t_P$ is the time step used during prediction. For example, the LSTM network is trained using the data generated with a time step of 0.001 MTU, and $\Delta t_P = 2\Delta t_T$ means that the time step during the forecast stage is 0.002 MTU. For all numerical experiments with a two-scale Lorenz 96 system, the observations are assimilated every 10 prediction time steps (i.e., $\Delta t_P$). We can observe that the RMSE is slightly increased for the LSTM based subgrid scale parameterization model with a different time step utilized during prediction. We do not observe any unstable behavior with a different time step for prediction than the time step used.
Figure 10: Power spectra analysis of the strongly nonlinear Lorenz system computed using the true physics-based model, the hybrid model where the \( \frac{dZ}{dt} \) is approximated using the neural network, and the assimilated results with the hybrid model. The observation noise is drawn from \( \mathcal{N}(0, 5) \), and are collected after every 50 time steps, i.e., the time interval between two observations is 0.05. Here, LSTM refers to forecast with the hybrid model and LSTM-DEnKF refers to an analysis state of the system.

for training. The RMSE for the LSTM-DEnKF model is higher when there is a difference between the prediction and training time step. However, there is still a considerable improvement in the prediction compared to employing only LSTM based subgrid scale parameterization. The numerical experiment with a large observation noise, i.e., \( \sigma^2_b = 10.0 \) suggests that the framework is robust for high uncertainty in measurements.

5 Summary and conclusions

In this work, we investigated the use of a neural network to recover the missing physics (i.e., incomplete dynamics) in dynamical systems. The hybrid neural-physics model composed of the physics-based model for the known dynamics and neural network for the unknown dynamics is applied to the Lorenz system with different strengths of nonlinearity. In the Lorenz system, the third equation was assumed to be missing and was approximated using the neural network. The recurrent neural network capable of handling long and short term dependencies was utilized for learning the mapping between known and unknown dynamics. Once the trained long short-term memory (LSTM) network is deployed in an auto-regressive manner, there is an accumulation of prediction error from one time to another. The weakly nonlinear Lorenz system exhibits near-regular oscillations with increasing amplitude and the hybrid neural-physics model is able to produce an accurate forecast even in the presence of prediction error during the online deployment of the LSTM network. The highly nonlinear Lorenz system being chaotic with extreme events is sensitive to the initial condition and a small change in the initial condition can produce a very different output. We observe that the predicted state trajectory of the hybrid neural-physics model starts deviating from the true trajectory after around 3.5 multiple of Lyapunov time for the highly nonlinear Lorenz system.

To achieve an accurate prediction over a longer period, the sequential data assimilation based on the deterministic ensemble Kalman filter (DEnKF) is applied. The hybrid neural-physics model coupled with DEnKF is able to predict the state of the highly nonlinear Lorenz system close to the true state by utilizing sparse and noisy observations. The analysis with the power spectral density (PSD) demonstrates that the hybrid neural-physics model integrated with DEnKF can capture the power density at lower frequencies accurately compared with the true PSD. We also illustrate the successful performance of the present framework for the two-scale Lorenz system where the subgrid scale
Figure 11: The trajectory of the strongly nonlinear Lorenz system computed using the true physics-based model, the hybrid model where the $dZ/dt$ is approximated using the neural network, and the assimilated results with the hybrid model. The observation noise is drawn from $N(0, 5)$, and are collected after every 100 time steps, i.e., the time interval between two observations is 0.1. Here, LSTM refers to forecast with the hybrid model and LSTM-DEnKF refers to an analysis state of the system. The parameterization is learned through the LSTM network. During the forecast period, the DEnKF algorithm is utilized to make use of sparse and noisy observations to improve the prediction. The DEnKF algorithm fail to give accurate analysis if unresolved scales are not parameterized (i.e., incomplete physics-based model). Furthermore, we show that the hybrid model forecast is not highly sensitive to the time step used for computing the subgrid-scale tendencies and the framework is robust against high observational noise. Based on numerical experiments carried out in this work, the methodology presented here seems promising for a continuous forecast of incomplete chaotic dynamical systems.

It may be argued that the success of this framework relies on the low-dimensionality of the Lorenz model. Even the two-scale Lorenz model has many fewer dimensions than any complex geophysical system and a relatively simple parameterization. Nevertheless, we foresee that the approach can be extended to high-dimensional complex systems by considering advanced neural network architectures like convolutional LSTMs, generative adversarial networks (GANs). The proposed method is modular enough that it is possible to plug any type of neural network and data assimilation scheme independently. The lessons learned from this study will guide our future extension of this approach to more complex chaotic systems whose dynamics is representative of geophysical flows. Moreover, a relevant possible extension is to combine machine learning, data assimilation, and reduced order models toward accelerating digital transformation and decision making processes in science, technology, and engineering workflows. To utilize the present framework for real-time decision making, it is necessary to reduce the computational requirement of the forward model significantly. Therefore, replacing the numerical discretization-based solver with the data-driven surrogate model will be a future research direction.

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Figure 12: The trajectory of the strongly nonlinear Lorenz system computed using the true physics-based model, the hybrid model where the $dZ/dt$ is approximated using the neural network, and the assimilated results with the hybrid model using an ensemble of 10 members. The observation noise is drawn from $N(0, 10)$, and are collected after every 50 time steps, i.e., the time interval between two observations is 0.05. Here, LSTM refers to forecast with the hybrid model and LSTM-DEnKF refers to an analysis state of the system.

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\[ \text{RMSE}(t_k) = \sqrt{\frac{1}{N} \sum_{n=1}^{N} (x_{n,k}^T(t_k) - x_{n,k}^p(t_k))^2}. \]
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Figure 17: Evolution of the RMSE for the two-scale Lorenz 96 model with different time step used in the prediction for $\sigma_b^2 = 1.0$ (left) and $\sigma_b^2 = 10.0$ (right). The $\Delta t_T$ represents the time step used for generating the training data, and $\Delta t_P$ is the time step used during prediction. The reported RMSE is the average mean squared error for all slow variables, i.e., $\text{RMSE}(t_k) = \sqrt{\frac{1}{N} \sum_{n=1}^{N} (x^n_T(t_k) - x^n_P(t_k))^2}$. Here, LSTM refers to forecast with the hybrid model and LSTM-DEnKF refers to an analysis state of the system.

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