ChaosBench: A Multi-Channel, Physics-Based Benchmark for Subseasonal-to-Seasonal Climate Prediction

Supplementary Material

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A Accountability and Reproducibility Statement

ChaosBench is published under the open source GNU General Public License. Further development and potential updates discussed in the limitations section will take place on the ChaosBench page. Furthermore, we are committed to maintaining and preserving the ChaosBench benchmark. Ongoing maintenance also includes tracking and resolving issues identified by the broader community after release. User feedback will be closely monitored via the GitHub issue tracker. All assets are hosted on GitHub and HuggingFace, which guarantees reliable and stable storage.

Dataset: All our dataset, present and future (e.g., with more years, multi-resolution support, etc) are available at https://huggingface.co/datasets/LEAP/ChaosBench.

Model Checkpoints: All of our model checkpoints used for the purposes of ablation in this work are available at https://huggingface.co/datasets/LEAP/ChaosBench/tree/main/logs.

Code: Our code and its future extension based on community feedback is accessible at https://github.com/leap-stc/ChaosBench.

Documentation: Finally, our main webpage will keep track of all important updates and latest documentation, and is accessible at https://leap-stc.github.io/ChaosBench.

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B Getting Started

Here, we provide a detailed description on how to prepare the necessary data, perform training, and benchmark your own model. However, we refer users to our webpage https://leap-stc.github.io/ChaosBench for the most updated how-to guides.

The following sections assume successful cloning of our Github repository https://github.com/leap-stc/ChaosBench. If you find any problems, feel free to contact us or raise an issue.

B.1 Data Preparation

First, navigate to the repository directory and install the necessary dependencies.

```
$ cd ChaosBench
$ pip install -r requirements.txt
```

Second, download the dataset using the following commands.

```
$ cd data/
$ wget https://huggingface.co/datasets/LEAP/ChaosBench/resolve/main/
    process.sh
$ chmod +x process.sh
```

Third, process the following required and optional dataset.

```
# Required for inputs and climatology (e.g., normalization)
$ ./process.sh era5
$ ./process.sh lra5
$ ./process.sh oras5
$ ./process.sh climatology
# Optional: control (deterministic) forecasts
$ ./process.sh ukmo
$ ./process.sh ncep
$ ./process.sh cma
$ ./process.sh ecmwf
# Optional: perturbed (ensemble) forecasts
$ ./process.sh ukmo_ensemble
$ ./process.sh ncep_ensemble
$ ./process.sh cma_ensemble
$ ./process.sh ecmwf_ensemble
# Optional: SoTa (deterministic) forecasts
$ ./process.sh panguweather
$ ./process.sh graphcast
$ ./process.sh fourcastnetv2
```

B.2 Training

We will cover how training can generally be performed, followed by how one can switch between different training strategies by manipulating the config . yaml file.

First, define your model class.

```
# An example can be found for e.g. <YOUR_MODEL> == fno
$ touch chaosbench/models/<YOUR_MODEL>.py
```

Second, import and initialize your model in the main chaosbench/models/model.py file, given the pseudocode below.

```
# Examples for lagged_ae, fno, resnet, unet are provided
import lightning.pytorch as pl
from chaosbench.models import YOUR_MODEL

class S2SBenchmarkModel(pl.LightningModule):

    def __init__(
        self,
        ...
):
        super(S2SBenchmarkModel, self).__init__()

        # Initialize your model
        self.model = YOUR_MODEL.BEST_MODEL(...)

# The rest of model construction logic
```

Third, run the train.py script. We recommend using GPUs for training.

```
# The _s2s suffix identifies data-driven models

$ python train.py --config_filepath chaosbench/configs/<YOUR_MODEL>
    _s2s.yaml
```

Now you will notice that there is a .yaml file. We define the definition of each field, allowing for greater control over different training strategies.

Note,

- 1. If only_headline is set to True, then the model is optimized only for a subset of variables defined in config.HEADLINE_VARS (default: False).
- 2. If n_step is set to values greater than 1, the models will train over n-autoregressive steps (default: 1).
- 3. If lead_time is set to values greater than 1, the models will be able to forecast *n*-days ahead. For example, in our direct forecasts, if lead_time is set to 4, our model will predict the states 4 days into the future (default: 1).
- 4. If land_vars and/or ocean_vars are set with entries from the acronyms in Tables S3 and S2, these will be used as additional inputs and targets, on top of ERA5 variables (default: []).

B.3 Evaluation

Once training is done, we can perform evaluation depending on the use case. We recommend using GPUs for evaluation.

First, if we have an autoregressive model, we can simply run:

Second, if we have a collection of models trained specifically for unique lead_time, we can run:

```
# Evaluating direct model with the default sequence of
# lead_time = [1, 5, 10, 15, 20, 25, 30, 35, 40, 44] e.g.,
# --model_name 'unet_s2s'
# --eval_years 2022
# --version_nums 0 4 5 6 7 8 9 10 11 12
# --lra5 't2m' 'tp'  ## Additional LRA5 vars to be evaluated
# --oras5 'sosstsst'  ## Additional ORAS5 vars to be evaluated

$ python eval_direct.py --model_name <str> --eval_years <int> --
version_nums [...] --lra5 [...] --oras5 [...]
```

Third, if we have a probabilistic model that generates ensemble forecasts (e.g., one checkpoint represents one ensemble member) and are supposed to be evaluated with additional probabilistic metrics, we can run:

```
# Evaluating ensembles with additional probabilistic metrics e.g.,
# --model_name 'unet_ensemble_s2s'
# --eval_years 2022
# --version_nums 0 1 2 ## One ensemble member per version
# --lra5 't2m' 'tp' ## Additional LRA5 vars to be evaluated
# --oras5 'sosstsst' ## Additional ORAS5 vars to be evaluated

$ python eval_ensemble.py --model_name <str> --eval_years <int> --version_nums [...] --lra5 [...] --oras5 [...]
```

B.4 Optional: Processing Multi-Resolution Input

We open-source the data processing script to allow users to process the inputs given different resolution (highest is 0.25-degree):

```
# Process inputs with e.g., 0.25-degree resolution
$ python scripts/process_atmos.py --resolution 0.25 # ERA5
$ python scripts/process_ocean.py --resolution 0.25 # ORAS5
$ python scripts/process_land.py --resolution 0.25 # LRA5
```

C Related Work

Here we discuss the criteria used to compare different S2S benchmark. This list is by no means exhaustive and there exists many ways to interpret the different contribution, strength, and scope of each. We refer interested reader to the respective benchmark paper and website.

On Input Variables. The number of input channels indicates the number of *unique* variables used for training data-driven models. For instance, in the case of SubseasonalClimateUSA, these include tmin, tmax, tmean, precip_agg, precip_mean, SST, SIC, z-10, z100, z500, z850, u-250, u-925, v-250, v-925, surface_P, RH, SSP, precipitable water, PE, DEM, KG, MJO-phase, MJO-amp, ENSO-I, despite them having similar (25) variables across data sources.

On Target Variables and Agencies. Similarly, the number of target channels represent the variables these benchmarks are aiming for. This is closely related to the number of benchmark agencies, which refers to the number of physics-based simulations used as target, rather than inputs. In the case for SubseasonalClimateUSA, for instance, the number of target channels correspond to two: precipitation and surface temperature, while the number of benchmark agencies is also two: CFSv2 (NCEP) and IFS (ECMWF), despite them using multiple other simulations generated from agencies but as inputs; though evaluated on all in their follow-up work [50] despite not initially described in the dataset paper.

On Physics Metrics. The flag for physics-based metrics indicates whether these benchmarks incorporate not just physical explanation, but also formulate them as scalar and differentiable metrics for future optimization problem.

On Probabilistic Metrics. The flag for probabilistic metrics indicates whether these benchmarks incorporate probabilistic (e.g., CRPS, CRPSS, Spread, SSR), in addition to deterministic metrics.

On Spatial Extent. Furthermore, the spatial extent indicates the extent of the target benchmark, rather than of the input dataset. This is because some of the more challenging S2S forecasting task is to get the correct global space-time correlation, and having a full global coverage provides a more complete evaluation.

D ChaosBench

D.1 Observations from Reanalysis Products

D.1.1 ERA5

The following table indicates the 48 variables that are inferred by physics-based models. Note that the Input ERA5 observations contains **ALL** fields, including the unchecked boxes:

Table S1: List of ERA5 reanalysis variables

Parameters/Levels (hPa)	1000	925	850	700	500	300	200	100	50	10
Geopotential height, z (gpm)		√	√	✓	√	✓	√	√	\	√
Specific humidity, q $(kg kg^{-1})$	✓	✓	\checkmark	✓	✓	✓	✓			
Temperature, t (K)	✓	✓	\checkmark	✓	✓	✓	✓	✓	✓	✓
U component of wind, u (ms^{-1})	✓	✓	\checkmark	✓	✓	✓	✓	✓	✓	✓
V component of wind, v (ms^{-1})	✓	✓	\checkmark	✓	✓	✓	✓	✓	✓	✓
Vertical velocity, w (Pas^{-1})					\checkmark					

D.1.2 ORAS5

The variables for ORAS5 consist of the following as described in Table S2.

Table S2: List of ORAS5 reanalysis variables

Acronyms	Long Name	Units						
iicethic	sea ice thickness	m						
iicevelu	sea ice zonal velocity	ms^{-1}						
iicevelv	sea ice meridional velocity	ms^{-1}						
ileadfra	sea ice concentration	(0-1)						
so14chgt	depth of 14° isotherm	m						
so17chgt	depth of 17° isotherm	m						
so20chgt	depth of 20° isotherm	m						
so26chgt	depth of 26° isotherm	m						
so28chgt	depth of 28° isotherm	m						
sohefldo	net downward heat flux	Wm^{-2}						
sohtc300	heat content at upper 300m	Jm^{-2}						
sohtc700	heat content at upper 700m	Jm^{-2}						
sohtcbtm	heat content for total water column	Jm^{-2}						
sometauy	meridonial wind stress	Nm^{-2}						
somx1010	mixed layer depth 0.01	m						
somx1030	mixed layer depth 0.03	m						
sosaline	salinity	PSU						
sossheig	sea surface height	m						
sosstsst	sea surface temperature	$^{\circ}C$						
sowaflup	net upward water flux	$kg/m^2/s$						
sozotaux	zonal wind stress	Nm^{-2}						

D.1.3 LRA5

The variables for LRA5 consist of the following as described in Table S3.

Table S3: List of LRA5 reanalysis variables

	Table S3: List of LRA5 reanalysis variables							
Acronyms	Long Name	Units						
asn	snow albedo	(0 - 1)						
d2m	2-meter dewpoint temperature	K						
е	total evaporation	m of water equivalent						
es	snow evaporation	m of water equivalent						
evabs	evaporation from bare soil	m of water equivalent						
evaow	evaporation from open water	m of water equivalent						
evatc	evaporation from top of canopy	m of water equivalent						
evavt	evaporation from vegetation transpiration	m of water equivalent						
fal	forecaste albedo	(0 - 1)						
lai_hv	leaf area index, high vegetation	m^2m^{-2}						
lai_lv	leaf area index, low vegetation	m^2m^{-2}						
pev	potential evaporation	m						
ro	runoff	m						
rsn	snow density	kgm^{-3}						
sd	snow depth	m of water equivalent						
sde	snow depth water equivalent	m ¹						
sf	snowfall	m of water equivalent						
skt	skin temperature	K						
${ t slhf}$	surface latent heat flux	Jm^{-2}						
smlt	snowmelt	m of water equivalent						
snowc	snowcover	%						
sp	surface pressure	Pa						
src	skin reservoir content	m of water equivalent						
sro	surface runoff	m ¯						
sshf	surface sensible heat flux	Jm^{-2}						
ssr	net solar radiation	Jm^{-2}						
ssrd	download solar radiation	Jm^{-2}						
ssro	sub-surface runoff	m						
stl1	soil temperature level 1	K						
stl2	soil temperature level 2	K						
stl3	soil temperature level 3	K						
stl4	soil temperature level 4	K						
str	net thermal radiation	Jm^{-2}						
strd	downward thermal radiation	Jm^{-2}						
swvl1	volumetric soil water layer 1	m^3m^{-3}						
swvl2	volumetric soil water layer 2	m^3m^{-3}						
swv13	volumetric soil water layer 3	m^3m^{-3}						
swvl4	volumetric soil water layer 4	m^3m^{-3}						
t2m	2-meter temperature	K						
tp	total precipitation	m						
tsn	temperature of snow layer	K						
u10	10-meter u-wind	ms^{-1}						
v10	10-meter v-wind	ms^{-1}						

D.2 Physics-Based Simulations

In this section, we describe in detail the physics-based models used as baselines in ChaosBench. Wherever possible, we discuss specific strategies regarding coupling to the ocean, sea ice, wave, land, initialization and perturbation strategies, specifications of initial/boundary conditions, as well as other numerical considerations to generate forecast.

D.2.1 The UK Meteorological Office (UKMO) [33]

- **Initialization and Ensemble**. The UKMO model employs the lagged initialization strategy to generate an ensemble of forecasts (4 in this case) at different initialization time to improve prediction stability.
- Coupling with ocean is performed with the Global Ocean 6.0 model [51], based on NEMO3.6 [52] with 0.25 degree horizontal resolution and 75 vertical pressure levels. The ocean model is initialized and calibrated using Nonlinear Evolutionary Model VARiation (NEMOVAR) [53], a specific data assimilation strategy that uses temperature, salinity profiles, altimeter-derived sea level anomalies to calibrate forecasts. Frequency of coupling is 1-hourly.
- Coupling with sea ice is performed with the Global Sea Ice 8.1 (CICE5.1.2) model [54], and again initialized from NEMOVAR.
- Coupling with wave model is not yet operational.
- Coupling with land surface is performed with the Joint UK Land Environment Simulator (JULES) [55]. Soil moisture, soil temperature, and snow are initialized using JULES and forced using the the Japanese 55-year Reanalysis (JRA-55) data [56]. The land surface model is paramaterized by land cover type from a combination of satellite (e.g., MODIS LAI [57]) and radiometer data (e.g., AVHRR [58]). In addition, another parameterization in the form of soil characteristics is derived from the Harmonized World Soil Database [59].
- **Model grid** uses the Arakawa C-grid [60] to solve partial differential equations on a spherical surface. In particular, the velocity components (such as zonal and meridional wind) are defined at the center of each face of the grid cells (in the case of a rectilinear grid) or along cell edges (in the case of a curvilinear grid). The scalar quantities such as pressure or temperature are computed at the corners of the grid cells.
- Large-scale dynamics uses the Semi-Lagrangian approach. It does not strictly follow fluid parcels (i.e., Lagrangian), but it does calculate the value of a field, such as temperature (i.e., Eulerian) by tracing back along the trajectory that a fluid parcel would have taken to reach a specific point at the current time step. This backward trajectory is used to find the origin of the fluid parcel and determine its properties, which are then used to update the model fields. This hybrid approach is therefore termed Semi-Lagrangian.

D.2.2 National Centers for Environmental Prediction (NCEP) [61]

- **Initialization and Ensemble**. The NCEP model adds small perturbation to the atmospheric, oceanic and land analysis at each cycle across 4 ensemble to reduce sensitivity to initial conditions.
- Coupling with ocean is performed with the GFDL Modular Ocean Model version 4 (MOM4) model that has a spatial resolution of 0.5-degree and 0.25-degree in the longitude-latitude directions [62]. There are 40 vertical pressure levels.

- Coupling with sea ice is also performed with the GFDL Sea Ice Simulator (SIS), which models the thermodynamics and overall dynamics of sea ice [62].
- Coupling with wave model is not yet operational.
- Coupling with land surface is performed with 4-layer Noah Land surface model 2.7.1 [63]. Soil moisture, soil temperature, and snow are initialized using Noah and forced using the Climate Forecast System [61] and the Global Land Data Assimilation System [64] reanalysis data. The land surface model is parameterized by land cover type AVHRR data. In addition, another paramaterization in the form of soil characteristics is derived from the world soil climate database [65].
- **Model grid** uses the Gaussian grid [66], where the longitude (x-axis) are evenly spaced while the latitudes (y-axis) are not. Instead, they are determined by the roots of the associated Legendre polynomials, which correspond to the Gaussian quadrature points for the sphere. This ensures that the actual area represented by each grid cell is more uniform.
- Large-scale dynamics uses the Spectral approach. It solves partial differential equations by transforming them from the physical space into the spectral domain. In the latter case, the equations are transformed into a series of coefficients that represent the amplitude of waves across scales. The transformations are usually done using Fourier series for periodic domains or spherical harmonics when dealing with the whole Earth's surface [66]. This method is especially beneficial for smooth functions and for representing large-scale wave phenomena, such as the Rossby waves, which are important for understanding weather and climate.

D.2.3 China Meteorological Administration (CMA) [35]

- **Initialization and Ensemble**. The CMA model uses the lagged average forecasting (LAF) method across 4 ensemble members to ensure that the mean forecast is less sensitive to initial conditions.
- **Coupling with ocean** is performed with the GFDL MOM4 model, which has 40 vertical pressure levels [62]. Frequency of coupling is 2-hourly.
- Coupling with sea ice is performed with the GFDL Sea Ice Simulator (SIS), similar to that used by NCEP [62].
- Coupling with wave model is not yet operational.
- Coupling with land surface is performed with the Atmosphere-Vegetation Interaction Model version 2 (AVIM2) model [67] and the NCAR NCAR Community Land Model version 3.0 (CLMv3) [68]. Soil moisture, soil temperature, and snow are not initialized directly using reanalysis data, as used by other land surface models. Rather, air-sea-land-ice coupled model is forced by near-surface atmospheric and ocean reanalysis in a long-term integration, and the land initial conditions are produced as a by-product. As a result, the parameterization of land cover type is done by this process, while soil characteristics is derived from the Harmonized World Soil Database [59].
- Model grid uses the Gaussian grid [66], similar to that used by the NCEP.
- Large-scale dynamics uses a mixture of Spectral approach for the vorticity, temperature, and surface pressure, as well as Semi-Lagrangian for specific humidity and cloud waters other tracers.

D.2.4 European Center for Medium-Range Weather Forecasts (ECMWF) [36]

- **Initialization and Ensemble**. The operational IFS forecast is generated through Singular Vectors (SV) method: it creates a variety of initial conditions by adjusting certain parameters slightly, thus generating different starting points.
- **Coupling with ocean** is performed with NEMO3.4.1 with 1-degree resolution and 42 vertical pressure levels. Frequency of coupling is 3-hourly.
- Coupling with sea ice is not operational for this model's version (but it is in the newer generation, though the forecast start-date is much later than 2016). As a result, sea ice initial conditions are persisted up to day 15 and then relaxed to climatology up to day 45.
- Coupling with wave model is performed with ECMWF wave model with 0.5-degree resolution [69].
- Coupling with land surface is relatively more complex than the rest, and we refer readers to their documentation. Regardless, it is based on Land Data Assimilation System (LDAS) that combines heterogenous high-quality dataset from satellite to ground sensors, and integrated with the operational IFS model. The parameterization for land cover type is primarily based on MODIS collection 5 [57] and soil characteristics from the FAO dominant soil texture class [70].
- **Model grid** uses the Cubic Octohedral grid [71], where the Earth's surface is projected onto a cube. Then, the cube is further subdivided to form an octahedron, where the faces represent finer grid cells. This multi-scale gridding scheme allows for parallelization where processes at different scales could be solved simultaneously.
- Large-scale dynamics uses a mixture of Spectral and Semi-Lagrangian approach, similar to that used by CMA.

E Data-Driven Baseline Models

In this section, we describe in detail implementation and hyperparemeter selections of our data-driven models used as baselines to ChaosBench. Most of the choices are based on the original works that are adapted to weather and climate applications using similar input dataset. All training are performed using 2x NVIDIA A100 GPUs.

E.1 Lagged Autoencoder (AE)

We implement lagged AE from [72] with 5 encoder blocks and 5 decoder block, with detailed specification in Table S4. Each encoder block is comprised of MaxPool2D \circ (Conv2D \rightarrow BatchNorm2D \rightarrow ReLU \rightarrow Conv2D \rightarrow BatchNorm2D \rightarrow ReLU). Similarly, the decoder block is comprised of ConvTranspose2D \rightarrow BactNorm2D \rightarrow ReLU) \bigoplus (ConvTranspose2D \rightarrow BactNorm2D \rightarrow Sigmoid) \circ (Conv2D).

Table S4:	Hyperparameters	s for Lagged AE

Hyperparameters	Values			
Channels	[64, 128, 256, 512, 1024]			
Encoder Kernel	3×3			
Decoder Kernel	2×2			
Max Pooling Window	2×2			
Batch Normalization	True			
Optimizer	ADAMW [73]			
Learning Rate	CosineAnnealing $(10^{-2} \rightarrow 10^{-3})$			
Batch Size	32			
Epochs	500			
Tmax	500			

E.2 ResNet

We adapt ResNet implementation from [42] using ResNet-50 as feature extractor and 5 decoder blocks, following specification in Table S5. Each decoder block is composed of ConvTranspose2D → Bactnorm2D → Leakyrelu.

Table S5: Hyperparameters for ResNet

Hyperparameters	Values				
Backbone	RESNET-50				
Decoder Channels	[1024, 512, 256, 128, 64]				
Decoder Activation	LEAKYRELU(0.15)				
Optimizer	AdamW				
Learning Rate	CosineAnnealing $(10^{-2} \rightarrow 10^{-3})$				
Batch Size	32				
Epochs	500				
Tmax	500				

E.3 UNet

We adapt UNet implementation from [22] using 5 *encoder* and 5 *decoder* blocks, with skip connections, following specification in Table S6. The composition of the encoder and decoder components are similar to those described for Lagged Autoencoder, with the addition of SKIP connection between each corresponding contracting-expansive path.

Table S6: Hyperparameters for UNet

Hyperparameters	Values			
Channels	[64, 128, 256, 512, 1024]			
Activation	LEAKYRELU(0.15)			
Encoder Kernel	3×3			
Decoder Kernel	2×2			
Max Pooling Window	2×2			
Optimizer	ADAMW			
Learning Rate	CosineAnnealing $(10^{-2} \rightarrow 10^{-3})$			
Batch Size	32			
Epochs	500			
Tmax	500			

E.4 Fourier Neural Operator (FNO)

We adapt FNO implementation from [43], following specification in Table S7 and illustrated in S1. We implement the encoder-decoder structure, where we (1) first transform our input X_t by convolutional layers both in the Fourier (applying fast fourier transform; FFT) and physical domains, before we concatenate both (applying inverse FFT for the former convolved features), and apply non-linear GELU activation function [74]. We select only the first 4 main Fourier modes to make the number of trainable parameters comparable with the other data-driven baseline models. The (2) decoder block then applies deconvolutional operation to the latent features to generate output Y_t .

Table S7: Hyperparameters for FNO

Hyperparameters	Values
Non-Spectral Channels	[64, 128, 256, 512, 1024]
Spectral Channel	[64, 128, 256, 512, 1024]
Activation	GELU
Fourier Modes	(4,4)
Optimizer	ADAMW
Learning Rate	CosineAnnealing $(10^{-2} \rightarrow 10^{-3})$
Batch Size	32
Epochs	500
Tmax	500

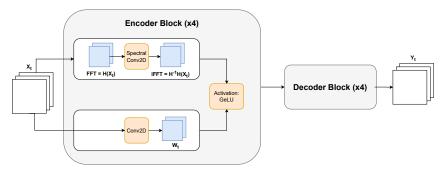


Figure S1: FNO architecture: (1) in the encoder block, we transform our input X_t by convolutional layers both in the Fourier and physical domains, before we concatenate and apply non-linear GELU activation function. The (2) decoder block is then applying deconvolutional operation to the latent features to generate forecast Y_t .

E.5 ClimaX

ClimaX is based on the ViT model [75] with variational positional embedding in variable-time space. We use ClimaX model as is described and implemented in the original paper and is pre-trained using CMIP6 [17]. We fine-tune the original pre-trained model given our training setup.

E.6 PanguWeather, FourCastNetV2, GraphCast

We perform inference using their latest checkpoints using the API provided here: https://github.com/ecmwf-lab/ai-models.

For this work, we process the forecasts at biweekly temporal resolution. In the codebase, we provide the script for further flexibility, for instance:

```
# Process biweekly, 1.5-degree forecasts for the year 2022

## Panguweather
$ python scripts/process_sota.py --model_name panguweather --years 2022

## Graphcast
$ python scripts/process_sota.py --model_name graphcast --years 2022

## FourCastNetV2
$ python scripts/process_sota.py --model_name fourcastnetv2 --years 2022
```

F Evaluation Metrics

F.1 Deterministic Metrics

We describe in detail the four primary vision-based metrics used for this benchmark, including RMSE, Bias, ACC, and MS-SSIM.

F.1.1 Root Mean-Squared Error (RMSE)

As described in the main text, we apply latitude-adjustment to RMSE computation.

$$\mathcal{M}_{RMSE} = \sqrt{\frac{1}{|\boldsymbol{\theta}||\boldsymbol{\gamma}|} \sum_{i=1}^{|\boldsymbol{\theta}|} \sum_{j=1}^{|\boldsymbol{\gamma}|} w(\theta_i) (\hat{\mathbf{Y}}_{i,j} - \mathbf{Y}_{i,j})^2}$$
(S1)

F.1.2 Bias

Similarly, we apply latitude-adjustment to Bias computation.

$$\mathcal{M}_{Bias} = \frac{1}{|\boldsymbol{\theta}||\boldsymbol{\gamma}|} \sum_{i=1}^{|\boldsymbol{\theta}|} \sum_{j=1}^{|\boldsymbol{\gamma}|} w(\theta_i) (\hat{\mathbf{Y}}_{i,j} - \mathbf{Y}_{i,j})$$
 (S2)

F.1.3 Anomaly Correlation Coefficient (ACC)

We remove the indexing for a more compact representation where the summation is performed over each grid cell (i,j). The predicted and observed anomalies at each grid-cell are denoted by $A_{\hat{\mathbf{Y}}_{i,j}} = \hat{\mathbf{Y}}_{i,j} - C$ and $A_{\mathbf{Y}_{i,j}} = \mathbf{Y}_{i,j} - C$, where C is the observational climatology. We apply latitude-adjustment to ACC computation.

$$\mathcal{M}_{ACC} = \frac{\sum w(\theta) [A_{\hat{\mathbf{Y}}} \cdot A_{\mathbf{Y}}]}{\sqrt{\sum w(\theta) A_{\hat{\mathbf{Y}}}^2 \sum w(\theta) A_{\mathbf{Y}}^2}}$$
(S3)

F.1.4 Multi-scale Structural Similarity Index Measure (MS-SSIM)

Let \mathbf{Y} and $\hat{\mathbf{Y}}$ be two images to be compared, and let $\mu_{\mathbf{Y}}$, $\sigma_{\mathbf{Y}}^2$ and $\sigma_{\mathbf{Y}\hat{\mathbf{Y}}}$ be the mean of \mathbf{Y} , the variance of \mathbf{Y} , and the covariance of \mathbf{Y} and $\hat{\mathbf{Y}}$, respectively. The luminance, contrast and structure comparison measures are defined as follows:

$$l(\mathbf{Y}, \hat{\mathbf{Y}}) = \frac{2\mu_{\mathbf{Y}}\mu_{\hat{\mathbf{Y}}} + C_1}{\mu_{\mathbf{Y}}^2 + \mu_{\hat{\mathbf{Y}}}^2 + C_1},$$
 (S4)

$$c(\mathbf{Y}, \hat{\mathbf{Y}}) = \frac{2\sigma_{\mathbf{Y}}\sigma_{\hat{\mathbf{Y}}} + C_2}{\sigma_{\mathbf{Y}}^2 + \sigma_{\hat{\mathbf{Y}}}^2 + C_2},$$
 (S5)

$$s(\mathbf{Y}, \hat{\mathbf{Y}}) = \frac{\sigma_{\mathbf{Y}\hat{\mathbf{Y}}} + C_3}{\sigma_{\mathbf{Y}}\sigma_{\hat{\mathbf{Y}}} + C_3},$$
(S6)

where C_1 , C_2 and C_3 are constants given by

$$C_1 = (K_1 L)^2, C_2 = (K_2 L)^2, \text{ and } C_3 = C_2/2.$$
 (S7)

L=255 is the dynamic range of the gray scale images, and $K_1\ll 1$ and $K_2\ll 1$ are two small constants. To compute the MS-SSIM metric across multiple scales, the images are successively low-pass filtered and down-sampled by a factor of 2. We index the original image as scale 1, and the desired highest scale as scale M. At each scale, the contrast comparison and structure comparison

are computed and denoted as $c_j(\mathbf{Y}, \hat{\mathbf{Y}})$ and $s_j(\mathbf{Y}, \hat{\mathbf{Y}})$ respectively. The luminance comparison is only calculated at the last scale M, denoted by $l_M(\mathbf{Y}, \hat{\mathbf{Y}})$. Then, the MS-SSIM metric is defined by

$$\mathcal{M}_{MS-SSIM} = [l_M(\mathbf{Y}, \hat{\mathbf{Y}})]^{\alpha_M} \cdot \prod_{j=1}^{M} [c_j(\mathbf{Y}, \hat{\mathbf{Y}})]^{\beta_j} [s_j(\mathbf{Y}, \hat{\mathbf{Y}})]^{\gamma_j}$$
(S8)

where α_M , β_j and γ_j are parameters. We use the same set of parameters as in [37]: $K_1=0.01$, $K_2=0.03$, M=5, $\alpha_5=\beta_5=\gamma_5=0.1333$, $\beta_4=\gamma_4=0.2363$, $\beta_3=\gamma_3=0.3001$, $\beta_2=\gamma_2=0.2856$, $\beta_1=\gamma=0.0448$. The predicted and ground truth images of physical variables are re-scaled to 0-255 prior to the calculation of their MS-SSIM values.

F.2 Physics-Based Metrics

In this section, we describe in detail the definition and implementation of our physics-based metrics, including PYTORCH psuedocode implementation.

Let Y be a 2D image of size $h \times w$ for a physical variables at a specific time, variable, and level. Let f(x,y) be the intensity of the pixel at position (x,y). First, we compute the 2D Fourier transform of the image by

$$F(k_x, k_y) = \sum_{x=0}^{w-1} \sum_{y=0}^{h-1} f(x, y) \cdot e^{-2\pi i (k_x x/w + k_y y/h)},$$
 (S9)

where k_x and k_y correspond to the wavenumber components in the horizontal and vertical directions, respectively, and i is the imaginary unit. The power at each wavenumber component (k_x, k_y) is given by the square of the magnitude spectrum of $F(k_x, k_y)$, that is,

$$S(k_x, k_y) = |F(k_x, k_y)|^2 = \operatorname{Re}[F(k_x, k_y)]^2 + \operatorname{Im}[F(k_x, k_y)]^2.$$
 (S10)

The scalar wavenumber is defined as:

$$k = \sqrt{k_x^2 + k_y^2},\tag{S11}$$

which represents the magnitude of the spatial frequency vector, indicating how rapidly features change spatially regardless of direction. Then, the energy distribution at a spatial frequency corresponding to k is defined as

$$S(k) = \sum_{(k_x, k_y): \sqrt{k_x^2 + k_y^2} = k} S(k_x, k_y).$$
 (S12)

Given the spatial energy frequency distribution for observations E(k) and predictions $\hat{S}(k)$, we perform normalization for each over \mathbf{K}_q , the set of wavenumbers corresponding to high-frequency components of energy distribution, as defined in Equation S13. This is to ensure that the sum of the component sums up to 1 which exhibits pdf-like property.

$$S'(k) = \frac{S(k)}{\sum_{k \in \mathbf{K}_q} S(k)}, \quad \hat{S}'(k) = \frac{\hat{S}(k)}{\sum_{k \in \mathbf{K}_q} \hat{S}(k)}, \quad k \in \mathbf{K}_q$$
 (S13)

```
import torch
import torch.nn as nn
class SpectralDiv(nn.Module):
    Compute Spectral divergence given the top-k percentile wavenumber
       (higher k means higher frequency)
   def
        __init__(
        self,
        percentile=0.9,
       input_shape=(121,240)
        super(SpectralDiv, self).__init__()
        self.percentile = percentile
        # Compute the discrete Fourier Transform sample frequencies
           for a signal of size
        nx, ny = input_shape
        kx = torch.fft.fftfreq(nx) * nx
        ky = torch.fft.fftfreq(ny) * ny
        kx, ky = torch.meshgrid(kx, ky)
        # Construct discretized k-bins
        self.k = specify_k_bins(...)
        # Get k-percentile index
        self.k_percentile_idx = int(len(self.k) * self.percentile)
   def forward(self, predictions, targets):
        # Preprocess data, including handling of missing values, etc
        predictions = preprocess_data(...)
        targets = preprocess_data(...)
        # Compute along mini-batch
        predictions, targets = torch.nanmean(predictions, dim=0),
           torch.nanmean(targets, dim=0)
        # Transform prediction and targets onto the Fourier space and
           compute the power
        predictions_power = torch.fft.fft2(predictions)
        predictions_power = torch.abs(predictions_power)**2
        targets_power = torch.fft.fft2(targets)
        targets_power = torch.abs(targets_power)**2
        # Normalize as pdf
        predictions_Sk = predictions_power / torch.nansum(
           predictions_power)
        targets_Sk = targets_power / torch.nansum(targets_power)
        # Compute spectral Sk divergence
        div = torch.nansum(targets_Sk * torch.log(torch.clamp(
           targets_Sk / predictions_Sk, min=1e-9)))
        return div
```

Listing S1: Psuedocode for computing SpecDiv using PyTorch

```
import torch
import torch.nn as nn
class SpectralRes(nn.Module):
    Compute Spectral residual given the top-k percentile wavenumber (
       higher k means higher frequency)
   def
        __init__(
        self,
        percentile=0.9,
       input_shape=(121,240)
        super(SpectralRes, self).__init__()
        self.percentile = percentile
        # Compute the discrete Fourier Transform sample frequencies
           for a signal of size
        nx, ny = input_shape
        kx = torch.fft.fftfreq(nx) * nx
        ky = torch.fft.fftfreq(ny) * ny
        kx, ky = torch.meshgrid(kx, ky)
        # Construct discretized k-bins
        self.k = specify_k_bins(...)
        # Get k-percentile index
        self.k_percentile_idx = int(len(self.k) * self.percentile)
   def forward(self, predictions, targets):
        # Preprocess data, including handling of missing values, etc
        predictions = preprocess_data(...)
        targets = preprocess_data(...)
        # Compute along mini-batch
        predictions, targets = torch.nanmean(predictions, dim=0),
           torch.nanmean(targets, dim=0)
        # Transform prediction and targets onto the Fourier space and
           compute the power
        predictions_power = torch.fft.fft2(predictions)
        predictions_power = torch.abs(predictions_power)**2
        targets_power = torch.fft.fft2(targets)
        targets_power = torch.abs(targets_power)**2
        # Normalize as pdf
        predictions_Sk = predictions_power / torch.nansum(
           predictions_power)
        targets_Sk = targets_power / torch.nansum(targets_power)
        # Compute spectral Sk residual
        res = torch.sqrt(torch.nanmean(torch.square(predictions_Sk -
           targets_Sk)))
        return res
```

Listing S2: Psuedocode for computing SpecRes using PyTorch

F.3 Probabilistic Metrics

Here, we broadly define $n \in N$ as an ensemble member, and $N \in \mathbb{R}$ the total number of ensemble members.

F.3.1 Deterministic Extension

This includes the ensemble version of deterministic and physics-based metrics, including RMSE, Bias, ACC, MS-SSIM, SpecDiv, and SpecRes.

$$\mathcal{M}_{RMSE}^{ens} = \frac{1}{N} \sum_{n=1}^{N} \mathcal{M}_{RMSE}^{n}$$
 (S14)

$$\mathcal{M}_{Bias}^{ens} = \frac{1}{N} \sum_{n=1}^{N} \mathcal{M}_{Bias}^{n}$$
 (S15)

$$\mathcal{M}_{ACC}^{ens} = \frac{1}{N} \sum_{n=1}^{N} \mathcal{M}_{ACC}^{n}$$
 (S16)

$$\mathcal{M}_{MS-SSIM}^{ens} = \frac{1}{N} \sum_{n=1}^{N} \mathcal{M}_{MS-SSIM}^{n}$$
 (S17)

$$\mathcal{M}_{SpecDiv}^{ens} = \frac{1}{N} \sum_{n=1}^{N} \mathcal{M}_{SpecDiv}^{n}$$
 (S18)

$$\mathcal{M}_{SpecRes}^{ens} = \frac{1}{N} \sum_{r=1}^{N} \mathcal{M}_{SpecRes}^{n}$$
 (S19)

F.3.2 CRPS

CRPS measures the accuracy of probabilistic forecasts by integrating the square of the difference between the cumulative distribution function (CDF) of the forecast and the CDF of the observed data over all possible outcomes. It can be thought of as probabilistic MAE, where a smaller value is desirable and a deterministic forecast reduces to MAE. We first apply latitude-adjustments for the forecasts and target fields.

$$\mathcal{M}_{CRPS}(F, x) = \int_{-\infty}^{\infty} (F(y) - H(y - x))^2 dy$$
 (S20)

where F(y) is the CDF of the forecast, H(y-x) is the Heaviside step function at the observed value x, and y ranges over all possible outcomes.

F.3.3 CRPSS

CRPSS measures the skillfulness of an ensemble forecasts, with positive being skillful, zero unskilled, and negative being worse than baseline climatology.

$$\mathcal{M}_{CRPSS} = 1 - \frac{CRPS_{\text{forecast}}}{CRPS_{\text{climatology}}}$$
 (S21)

F.3.4 Spread

We apply latitude-adjusted spread of the ensemble members, and std is the standard deviation operator.

$$\mathcal{M}_{Spread} = \frac{1}{|\boldsymbol{\theta}||\boldsymbol{\gamma}|} \sum_{i=1}^{|\boldsymbol{\theta}|} \sum_{j=1}^{|\boldsymbol{\gamma}|} \operatorname{std}\left(\{w(\boldsymbol{\theta}_i) \hat{\mathbf{Y}}_{i,j}^n\}_{n=1}^N\right) \tag{S22}$$

F.3.5 Spread/Skill Ratio (SSR)

We use ensemble RMSE as the skill in the SSR computation.

$$\mathcal{M}_{SSR} = \frac{\mathcal{M}_{Spread}}{\mathcal{M}_{RMSE}^{ens}} \tag{S23}$$

G Extended Results

We provide extended results accompanying the main text.

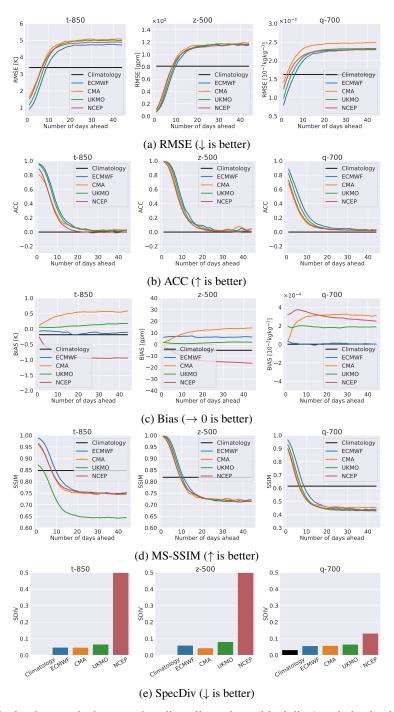


Figure S2: Evaluation results between baseline climatology (black line) and physics-based control (deterministic) forecasts. At longer forecasting horizon, most physics-based deterministic forecasts perform worse than climatology while maintaining structures as evidenced from their low SpecDiv (barring NCEP).

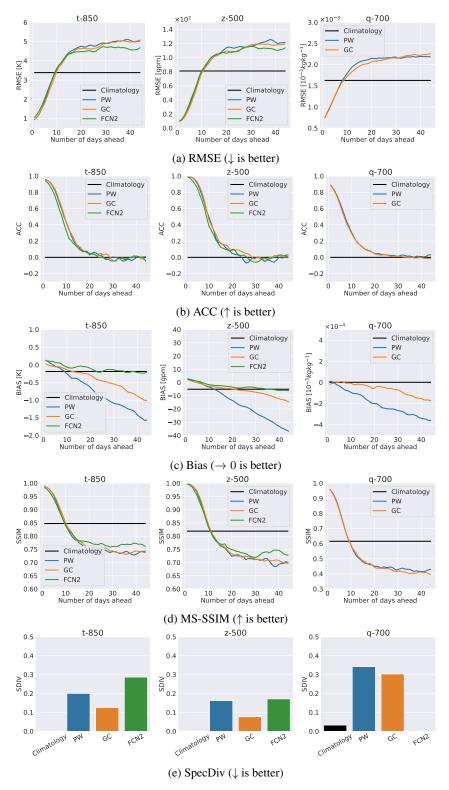


Figure S3: Evaluation results between baseline climatology (black line) and data-driven models including PanguWeather (PW), FourCastNetV2 (FCN2), and GraphCast (GC). Overall, we observe that data-driven models perform significantly worse than climatology on S2S timescale. They also perform poorly on physics-based metrics indicating the lack of predictive power on multi-scale structures. Note: FCN2 lacks q-700 and climatology naturally has low SpecDiv (direct observations).

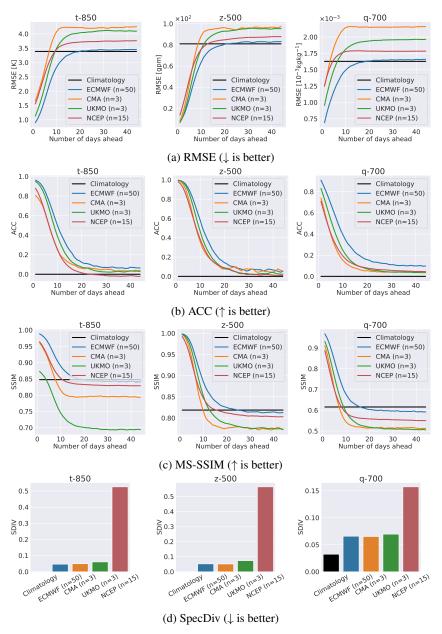


Figure S4: Evaluation results between baseline climatology (black line) and physics-based ensembles from ECMWF, CMA, UKMO, NCEP. Overall, we observe that ensemble forecasts perform better than their deterministic counterparts.

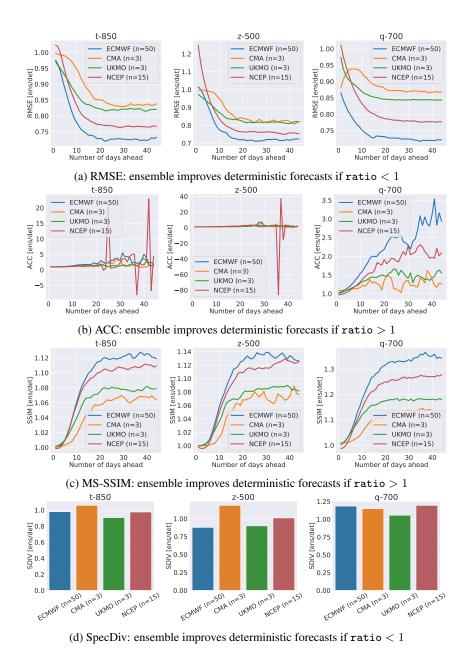


Figure S5: Metrics ratio e.g., ${\rm RMSE}_{ens}/{\rm RMSE}_{det}$ between ensemble and deterministic forecasts, where the former improves the latter by accounting for IC uncertainty that can lead to long-range instability and trajectory divergences. Note: n represents the number of ensemble members. The ratio for ACC fluctuates as the scalar value approaches 0.

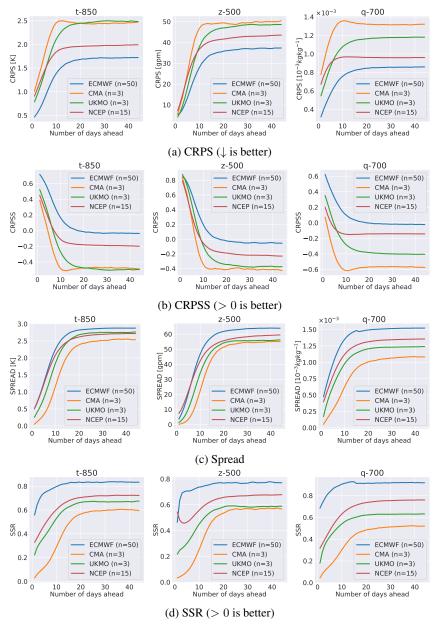


Figure S6: Probabilistic evaluation on ensemble forecasts indicating current skill limits of 15 days. Note: *n* represents the number of ensemble members.

G.1 Effects of Different Autoregressive Training Steps; lead_time

We showcased more results for autoregressive training strategy. In this case, we performed autoregressive training using either 1 or 5 iterative steps (n_step; s). As illustrated in Figure S7, we observe that incorporating temporal information improve the vision-based metrics even at longer forecasting timesteps, with lower RMSE, higher MS-SSIM. However, the converse trend is true incorporating temporal context makes S2S forecast worse off in some physics-based scores. The modified loss function for training a model with multiple autoregressive steps is:

$$\mathcal{L} = \frac{1}{|S|} \sum_{i=1}^{s} \mathcal{L}(\hat{\mathbf{Y}}_{\mathbf{t}+\mathbf{s_i}} \mathbf{Y}_{\mathbf{t}+\mathbf{s_i}}), \forall s_i \in S$$
 (S24)

Here $S = \{1, \dots, s\}$ and $s \in \mathbb{N}^+$ is the autoregressive steps. For this work, we set s = 5.

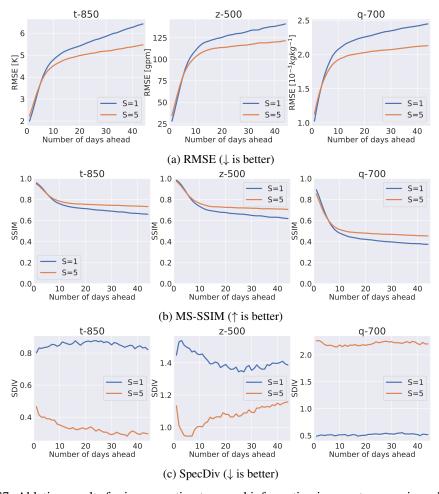


Figure S7: Ablation results for incorporating temporal information in an autoregressive scheme for long-range forecast using UNet models. The x-axis represents the number of forecasting days for t-850, z-500, q-700 representative tasks. Blue and orange lines illustrate autoregressive scheme with s=1 and s=5 respectively. Overall we observe that incorporating temporal information improve the vision-based metrics even at longer forecasting timesteps. However, the converse trend is true where incorporating temporal context makes S2S forecast worse off in some physics-based scores.

G.2 Effects of Subset Optimization; headline_vars

In many cases, we seek to train data-driven models so that they are able to perform well on *all* states by optimizing for the full state of the next forecasting timestep t+1, that is,

$$\phi^* = \operatorname*{argmin}_{\phi} \mathcal{L}(\hat{\mathbf{Y}}_{t+1}, \mathbf{Y}_{t+1})$$

where \mathcal{L} is any loss function. This task is especially useful for building emulators that act as surrogates for the more expensive physics-based NWP models [17].

Although the first task is useful for learning the full complex interaction between variables, it is relatively difficult due to the intrinsic high-dimensionality of the data. As a result, we introduce a second task that allows for the optimization on a subset of variables of interest $(Y' \in Y)$:

$$\phi^* = \operatorname*{argmin}_{\phi} \mathcal{L}(\hat{\mathbf{Y}}'_{t+1}, \mathbf{Y'}_{t+1})$$

Here $\mathbf{Y}'_{t+1} = \{\text{t-850, z-500, q-700}\}$, and we train them using 5 autoregressive steps i.e., $\mathbf{n_step} = 5$.

Table S8: Long-range forecasting ($\Delta t = 44$) results on select metrics and target variables between physics-based and data-driven models. Results are for Task 1 (full) and Task 2 (sparse). (*) Baseline model that uses privileged information (observations) to make prediction.

	RMSE ↓		MS-SSIM ↑			SpecDiv ↓			
Models	T850 (K)	Z500 (gpm)	Q700 $(\times 10^{-3})$	T850	Z500	Q700	T850	Z500	Q700
Climatology*	3.39	81.0	1.62	0.85	0.82	0.62	0.01	0.01	0.03
Persistence*	5.88	127.8	2.47	0.71	0.69	0.41	0.02	0.03	0.05
UKMO	5.00	116.2	2.32	0.64	0.71	0.43	0.06	0.09	0.07
NCEP	4.90	116.7	2.30	0.75	0.71	0.43	0.53	0.55	0.10
CMA	5.08	118.7	2.49	0.75	0.72	0.45	0.05	0.04	0.06
ECMWF	4.72	115.1	2.30	0.75	0.72	0.44	0.06	0.07	0.06
	Task 1: Full Dynamics Prediction								
Lagged AE	5.55	122.4	2.03	0.74	0.71	0.47	0.18	2.44	0.21
ResNet	5.67	125.3	2.07	0.73	0.70	0.47	0.21	0.37	0.26
UNet	5.47	121.5	2.13	0.73	0.71	0.45	0.30	1.16	2.20
FNO	5.06	112.5	1.95	0.75	0.73	0.51	0.18	0.11	0.10
	Task 2: Sparse Dynamics Prediction								
Lagged AE	5.39	119.0	2.12	0.75	0.73	0.48	0.52	1.41	0.29
ResNet	5.80	124.1	2.18	0.74	0.72	0.46	0.33	1.22	0.09
UNet	5.57	120.2	2.18	0.74	0.71	0.45	1.20	1.08	0.07
FNO	4.73	101.8	1.91	0.79	0.76	0.52	0.18	0.23	0.21

Overall, we find models that attempt to preserve spectral structures (e.g., FNO) perform better on all metrics, deterministic and physics-based. Also, Task 2 (sparse) appears to be easier than Task 1 (full). Nonetheless, they are still performing worse than climatology.

G.3 Effects of Ensemble Forecasts

This section provides additional results for data-driven ensemble approach, and follow similar evaluation process as the physics-based counterpart.

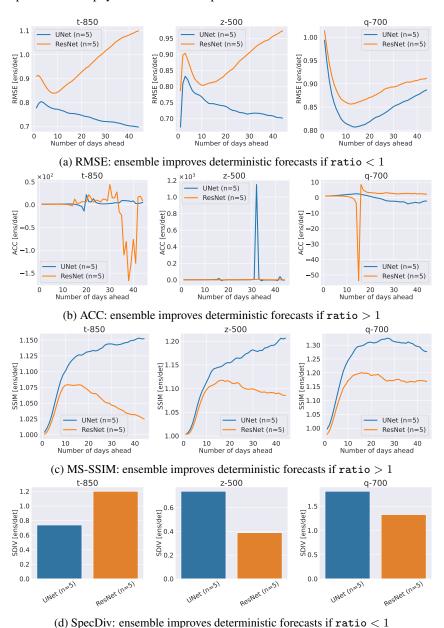


Figure S8: Metrics ratio e.g., $RMSE_{ens}/RMSE_{det}$ between ensemble and deterministic forecasts, where the former improves the latter by accounting for IC uncertainty that can lead to long-range instability and trajectory divergences. Note: n represents the number of ensemble members.

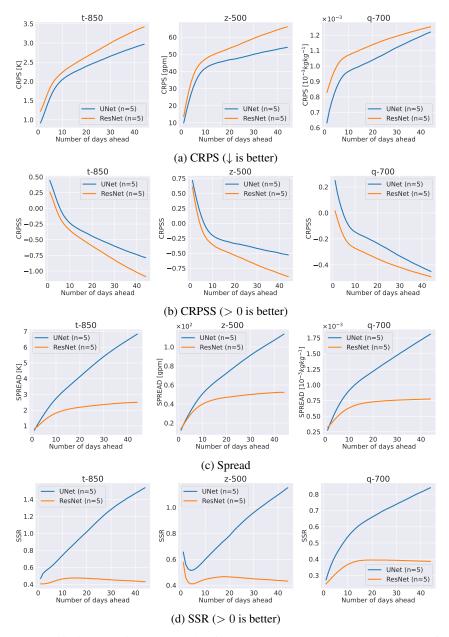


Figure S9: Probabilistic evaluation on ensemble forecasts. Note: n represents the number of ensemble members.

G.4 Power Spectra

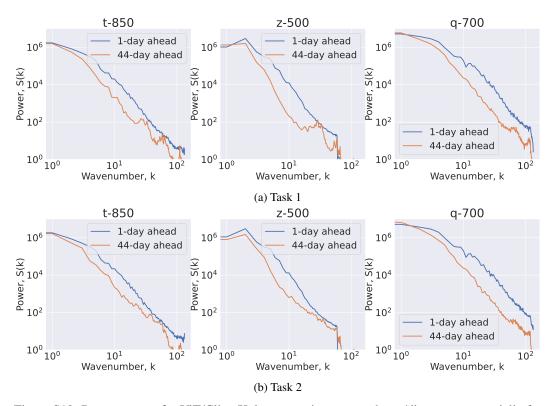


Figure S10: Power spectra for ViT/ClimaX demonstrating energy decay/divergence especially for high k as lead time grows.

G.5 Qualitative Evaluation

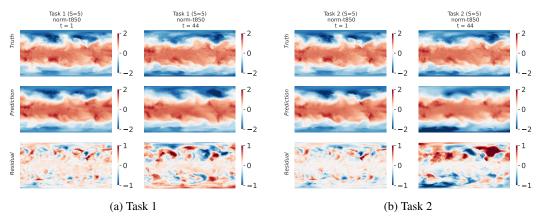


Figure S11: Normalized t@850-hpa qualitative results for UNet-autoregressive (S=5).

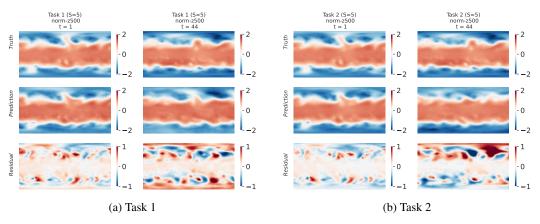


Figure S12: Normalized z@500-hpa qualitative results for UNet-autoregressive (S=5).

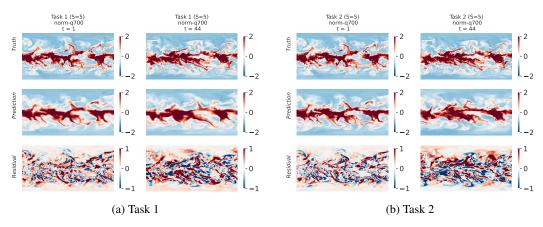


Figure S13: Normalized q@700-hpa qualitative results for UNet-autoregressive (S=5).

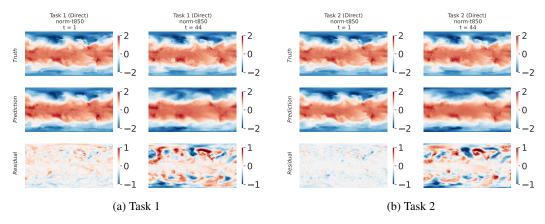


Figure S14: Normalized t@850-hpa qualitative results for UNet-direct.

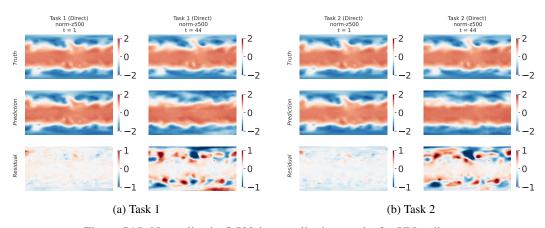


Figure S15: Normalized z@500-hpa qualitative results for UNet-direct.

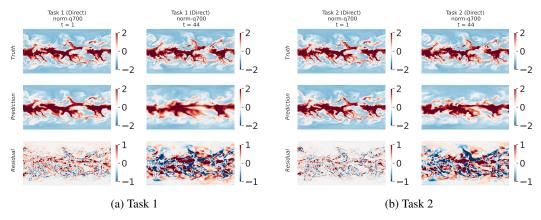


Figure S16: Normalized q@700-hpa qualitative results for UNet-direct.

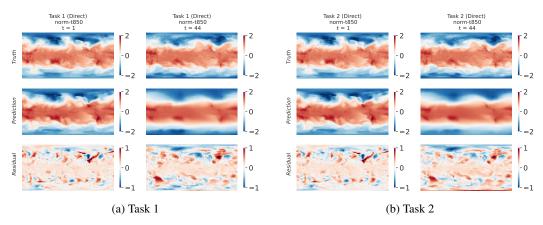


Figure S17: Normalized t@850-hpa qualitative results for ClimaX-direct.

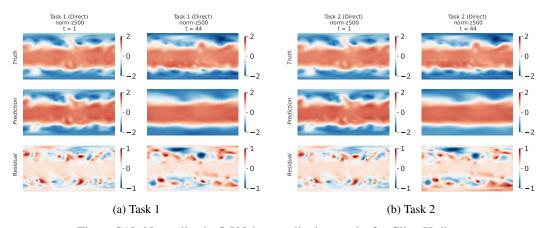


Figure S18: Normalized z@500-hpa qualitative results for ClimaX-direct.

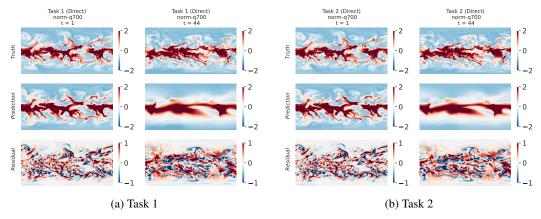


Figure S19: Normalized q@700-hpa qualitative results for ClimaX-direct.