ClimateBench v1.0: A benchmark for data-driven climate projections

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Key Points:

- We introduce the first benchmark for emulation of key spatially resolved climate variables derived from a full complexity Earth System Model
- Three baseline emulators are presented which are able to predict regional temperature and precipitation with varying skill
- Evaluation metrics and areas for future research are presented to encourage further development of trustworthy data-driven climate emulators

Abstract

Many different emission pathways exist that are compatible with the Paris climate agreement, and many more are possible that miss that target. While some of the most complex Earth System Models have simulated a small selection of Shared Socioeconomic Pathways, it is impractical to use these expensive models to fully explore the space of possibilities. Such explorations therefore mostly rely on one-dimensional impulse response models, or simple pattern scaling approaches to approximate the physical climate response to a given scenario. Here we present ClimateBench - the first benchmarking framework based on a suite of CMIP, AerChemMIP and DAMIP simulations performed by a full complexity Earth System Model, and a set of baseline machine learning models that emulate its response to a variety of forcers. These emulators can predict annual mean global distributions of temperature, diurnal temperature range and precipitation (including extreme precipitation) given a wide range of emissions and concentrations of carbon dioxide, methane and aerosols, allowing them to efficiently probe previously unexplored scenarios. We discuss the accuracy and interpretability of these emulators and consider their robustness to physical constraints such as total energy conservation. Future opportunities incorporating such physical constraints directly in the machine learning models and using the emulators for detection and attribution studies are also discussed. This opens a wide range of opportunities to improve prediction, robustness and mathematical tractability. We hope that by laying out the principles of climate model emulation with clear examples and metrics we encourage engagement from statisticians and machine learning specialists keen to tackle this important and demanding challenge.

Plain Language Summary

Many different emission pathways exist that are compatible with the Paris climate agreement, and many more are possible that miss that target. While some of the most complex Earth System Models have simulated a small selection of possible futures, it is impractical to use these expensive models to fully explore the space of possibilities. Such explorations therefore mostly rely on simple approximations of the global mean temperature response to a given scenario. Here we present ClimateBench - the first benchmarking framework based on a suite of state-of-the-art simulations performed by a full complexity Earth System Model, and a set of baseline machine learning models that emulate its response to a variety of forcers. These emulators can predict annual mean global distributions of temperature, diurnal temperature range and precipitation (including extreme precipitation) given a wide range of emissions and concentrations of carbon dioxide, methane and aerosols, allowing them to efficiently probe previously unexplored scenarios. We also describe a set of evaluation metrics which we hope will entice statisticians and machine learning experts to tackle this important and demanding challenge.

1. Introduction

Many different emission pathways exist that are compatible with the Paris Agreement of limiting global mean temperatures to "well below 2°C above pre-industrial levels and pursuing efforts to limit the temperature increase to 1.5 °C", and many more are possible that miss that target. Sampling possible emissions scenarios is therefore crucial for policy makers to weigh the economic cost and societal impact of different mitigation and adaptation strategies. While many of the most complex Earth System Models (ESMs) have simulated a small selection of 'Shared Socioeconomic Pathways' (SSPs; self-consistent emissions scenarios based on assumptions about future socio-economic changes and imperatives) it is impractical to use these expensive models to fully explore the space of possibilities (O'Neill et al., 2016). Therefore, such explorations mostly rely on one-dimensional impulse response models, or simple pattern scaling approaches to approximate the physical climate response to a given scenario (e.g., Millar et al., 2017).

Impulse response models (Smith et al., 2018; Meinshausen et al., 2011; Nicholls et al., 2020) are physically interpretable and can capture the general non-linear behaviour of the system, but are inherently unable to model regional climate changes, while pattern scaling approaches rely on a simple scaling of spatial distributions of temperature (e.g., Tebaldi et al., 2014) by global mean temperature changes. This approach breaks down when considering precipitation, however, because of the strong non-linearities in its response to temperature (e.g., Cabré et al., 2010). Statistical emulators of the regional climate have been developed although these have been quite bespoke (Castruccio et al., 2014) or focus on the relatively simple problem of emulating temperature (Holden and Edwards, 2010). These approaches also do not account for the influence of aerosol, which can be important for both regional temperature and precipitation (e.g. Kasoar et al. 2018, Wilcox et al. 2020). As has been noted recently (Watson-Parris, 2021), approaches including non-linear pattern scaling (Beusch et al., 2020) and Gaussian process (GP) regression of long-term climate responses (Mansfield et al., 2020) suggest the possibility of using modern machine learning (ML) tools to produce robust and general emulators of future scenarios. However, comparing and contrasting these approaches is currently hindered by the lack of a consistent benchmark.

ClimateBench defines a set of criteria and metrics for objectively evaluating such climate model emulation; aims to demonstrate the feasibility of such emulators; and provides a curated dataset that will allow, and hopefully encourage, broader engagement with this challenge in the same way WeatherBench (Rasp et al., 2020) has achieved for weather modelling. The target is to predict annual mean global distributions of temperature (T), diurnal temperature range (DTR), precipitation (PR) and the 90^{th} percentile of precipitation (PR90). These variables are chosen to represent a range of important climate variables which respond differently to each forcing and include extreme changes (PR90) that might not be expected to scale in the same way as the mean. For example, while T has been shown to scale roughly linearly with global mean temperature changes (Castruccio et al., 2014), PR responds non-linearly, and DTR is more sensitive to aerosol perturbations than global mean temperature changes (Hansen et al., 1995). Four of the main anthropogenic forcing agents are provided as emulator inputs (predictors): carbon dioxide (CO2), sulphur dioxide (SO2; a precursor to sulphate aerosol), black carbon (BC) and methane (CH4). To enable spatially accurate emulators ClimateBench includes (annual mean): spatial

distributions of emissions for the short-lived aerosol species (SO2 and BC), globally averaged emissions of CH4, and global cumulative emissions of CO2.

The training data which is provided in order to support such predictions is generated from the simulations performed by the second (and latest) version of the Norwegian Earth System Model (NorESM2; Seland et al., 2020) as part of the sixth coupled model intercomparison project (CMIP6; Eyring et al., 2016). The provided inputs are constructed from the same input data that is used to drive the original simulations. While we could have included simulations from multiple different models, only one model submitted all of the DECK (Diagnostic, Evaluation, and Characterization of Klima), historical, AerChemMIP (Collins et al., 2017) and ScenarioMIP (O'Neill et al., 2016) experiments required for our purposes, making it impossible to provide a harmonised dataset. Further, there is no agreed way of robustly combining multiple models, and while statistically combining multiple different models can lead to improved skill (Pincus et al., 2008) the resulting variance is not reliable since the models are not truly independent (Knutti et al., 2013). Nevertheless, this single model dataset still allows us to explore both scenario uncertainty and internal variability. Further, since even very simple models are able to capture a variety of forcing responses (Smith et al. 2021), there is reason to believe that the response of the models to a given forcing is more consistent than the range of responses (e.g., Richardson et al. 2019). We thus suggest that an emulator that works best for NorESM2 will also have the tendency to perform better in emulating other CMIP models, mainly because the data characteristics are by design similar (CMIP models represent the same physical system). In contrast, variations in the structure of learning algorithms vary more significantly and follow entirely different ways of building a regression model.

As a demonstration of the variety of possible approaches to tackle this benchmark we also introduce three distinct baseline emulators trained and evaluated against ClimateBench. These constitute the first data driven models for the projection of multiple climatic variables and show promising skill in both the global-mean and spatial responses. We discuss the merits and challenges in using each class of (regression) model and hope these provide a useful starting point for researchers wishing to develop more advanced emulators.

The remainder of this paper describes the development of the dataset including the underlying ESM and all post-processing (Section 2), the evaluation metrics used to rank ClimateBench submissions (Section 3), the baseline emulators (Section 4), a discussion of such approaches and future opportunities for diverse approaches (Section 5) before providing a few concluding remarks in Section 6.

2. Data set description and preparation

The data provided as part of ClimateBench is a heavily curated version of that publicly available in the CMIP6 data archive. Here we describe the data extraction and processing steps, but the scripts used to perform this are also freely available (as described in the data availability statement).

We use a selection of complementary simulations in order to provide as large a training dataset as possible while attempting to avoid unnecessary redundancy. Table 1 details the full list of simulations included, the period they cover and a brief description of their purpose in this context. Given that the primary purpose of ClimateBench is to train emulators over different emission scenarios, ScenarioMIP simulations are a key component of the dataset. ScenarioMIP prescribes a limited set of possible future emissions pathways exploring different socio-economic scenarios representing plausible narratives. These scenarios are designed to span a range of mitigation scenarios (denoted by the first number in each scenario) and end-of-century forcing possibilities (denoted by the last two numbers in each scenario). We include all available simulations, including the AerChemMIP *ssp370-lowNTCF* variation of *ssp370* which includes lower emissions of near-term climate forcers (NTCFs) such as aerosol (but not methane). We choose *ssp245* as our test dataset against which all ClimateBench emulators are to be evaluated. This scenario represents a medium mitigation and medium forcing scenario, ensuring trained emulators are able to interpolate a solution rather than extrapolate (as discussed further in Section 5). The CMIP6 *historical* experiment is also included since it provides useful training data at low emissions values.

Table 1: Details of post-processed simulations provided as part of the ClimateBench dataset. Experiments denoted (*) are ancillary data that, while potentially useful, are not used in training the baseline emulators presented here.

Protocol	Experiment	Period	Notes
ScenarioMIP (O'Neill et al., 2016)	ssp126	2015 - 2100	A high ambition scenario designed to produce significantly less than 2 degrees warming by 2100.
	ssp245	2015 - 2100	Designed to represent a medium forcing future scenario. This is the test scenario to be held back for evaluation.
	ssp370	2015 - 2100	A medium-high forcing scenario with high emissions of near-term climate forcers (NTCF) such as methane and aerosol.
	ssp370-lowNTCF	2015 - 2054	Variation of SSP370 with lower emissions of aerosol and their precursors
	ssp585	2015 - 2100	This scenario represents the high end of the range of future pathways in the IAM literature and leads to a very large forcing of 8.5 Wm^{-2} in 2100.
CMIP6 (Eyring et al., 2016)	historical	1850 – 2014	A simulation using historical emissions of all forcing agents designed to recreate the historically observed climate.
	abrupt-4xCO2*	500 years	Idealised simulation in which CO2 is abruptly quadrupled. Other forcing agents remain unchanged.
	1pctCO2*	150 years	Idealised simulation in which CO2 is gradually increased by 1% / year. Other forcing agents remain unchanged.
	piControl*	500 years	Baseline simulation in which all forcing agents remain unchanged.

DAMIP (Gillett et al., 2016)	hist-GHG	1850 – 2014	A historical simulation with varying concentrations for CO2 and other long-lived greenhouse-gases (only).
	hist-aer	1850 – 2014	A historical simulation only forced by changes in anthropogenic aerosol.
	ssp245-aer	2015 - 2100	A medium forcing scenario with only changes in anthropogenic aerosol, which provides an alternative test scenario for emulator evaluation.

ClimateBench also includes a selection of more idealised simulations which are intended to provide training data at the 'corners' of the four-dimensional input space, again helping reduce the chances of extrapolation in the resulting emulators (as demonstrated in Figure A1). Two simulations that are commonly used to diagnose the equilibrium and transient climate sensitivity are *abrupt-4xCO2* and *1pctCO2,* respectively. As the name suggests, the *abrupt-4xCO2* includes an abrupt quadrupling of CO2 over the pre-industrial concentrations while all other forcing agents remain unchanged. This level of concentration represents the high end of future scenarios, broadly in line with *ssp585* but with no contribution from the other forcers, simplifying its interpretation. The abrupt nature of the forcing also allows the timescale of the responses to be determined which can be useful for emulators which account for this. The *1pctCO2* simulation gradually increases the atmospheric concentration of CO2 by 1% per year, again with other forcing agents unchanged. While potentially very useful, they are not used in the training of the emulators presented in this work. Two other idealised simulations performed as part of the Detection-Attribution Model Intercomparison Project (DAMIP; Gillett et al., 2016) represent the historical period forced by only CO2 and other long-lived greenhouse gases (*hist-GHG*), or only anthropogenic aerosol (*hist-aer*). These provide opportunities to train emulators in regions of the input (emissions) space that are at the limits of plausible future scenarios and were used in training the emulators described in Section 4.

Finally, the *piControl* simulation provides a baseline simulation with all forcings remaining unchanged from their pre-industrial values. All target variables are calculated as a change against this climatology to simplify the training and interpretation of the results. This long (500 year) simulation also enables a robust estimation of internal variability of the climate system for those emulators which are able to represent it in future work, as discussed further in Section 5.1.

2.1 Input variables

The input data for these simulations is prescribed by the experimental protocol and provided by the input4MIPS project (https://esgf-node.llnl.gov/search/input4mips/), which we collate and pre-process for ease of use. Specifically, we extract the provided global mean emissions of CO2 and CH4 for each of the realistic (historical, ScenarioMIP and DAMIP) experiments from the checksum files provided by the Community Emissions Data System (CEDS) dataset (Hoesly et al., 2018). We sum over each sector and each month in order to derive annual total emissions and convert from Kg to Gt of CO2. Some historical and future periods are only provided in 5-yearly increments, so we linearly interpolate to yearly values for consistency. The CO2 emissions are summed cumulatively since, for realistic scenarios, a compensation between forcing

efficiency and ocean uptake means the temperature response to CO2 is approximately linear in the cumulative emissions (Matthews and Caldeira 2008; Allen et al. 2009). Figure 1 shows the global mean emissions of each of the forcing agents under different future emissions scenarios, showing a wide range of possible pathways.

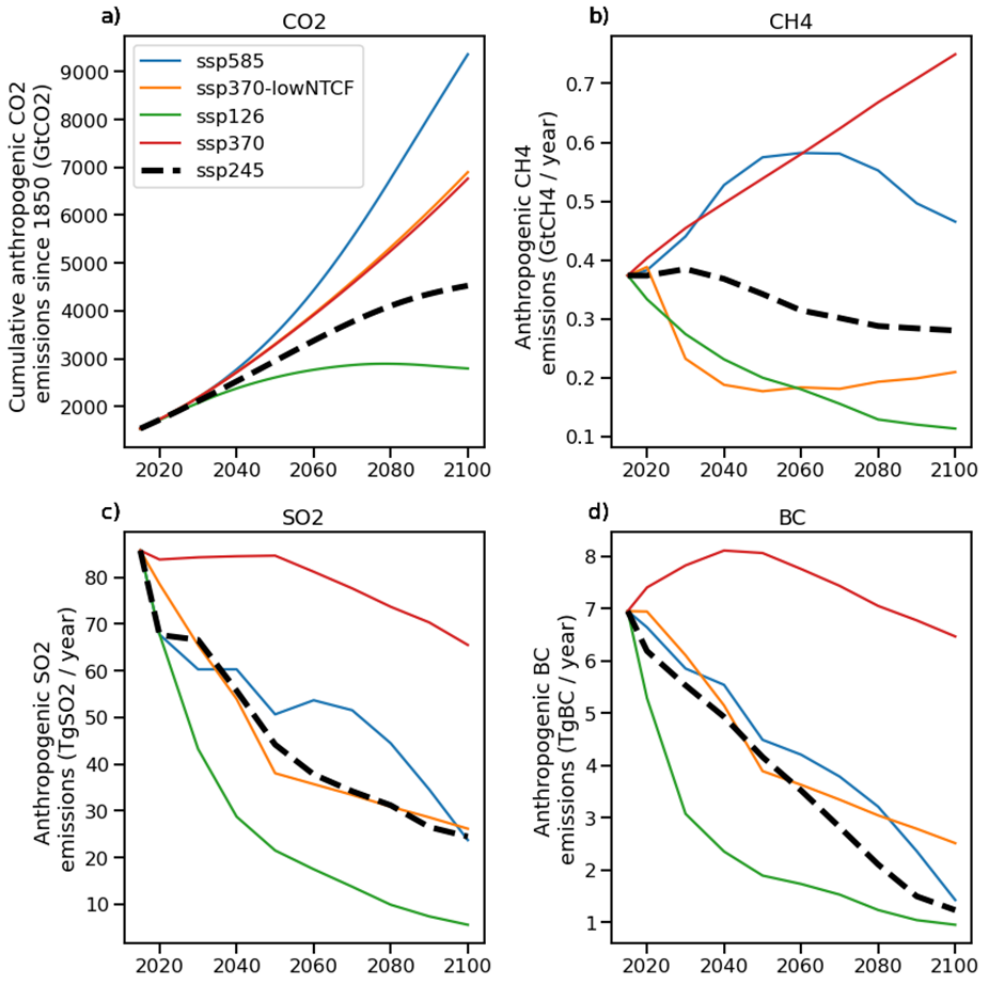

Figure 1: Time series of cumulative anthropogenic CO2 emissions since 1850 (a); emissions of CH4 (b); global mean emissions of SO2 (c) and black carbon (BC; d) derived from NorESM2 ScenarioMIP simulations available within ClimateBench, including the SSP245 test scenario (shown in black).

The aerosol (precursor) emissions are derived from the latest version of the spatially resolved CEDS dataset and again summed over sectors and months to produce maps of annual total emissions, as shown in Figure 2 for SO2 in different years. While the spatial distribution clearly evolves over the historical period and into the future scenarios, the emissions are fairly

localised around industrialised regions and dimensionality reduction can be used to reduce the size of these input features (as discussed for the baseline emulators in Section 4). An area preserving interpolation is performed so that the emission data are provided on the same spatial grid as the NorESM2 output fields to simplify its use in ML workflows. Again, as used for NorESM2 the 5-yearly data is interpolated to a yearly frequency for consistency.

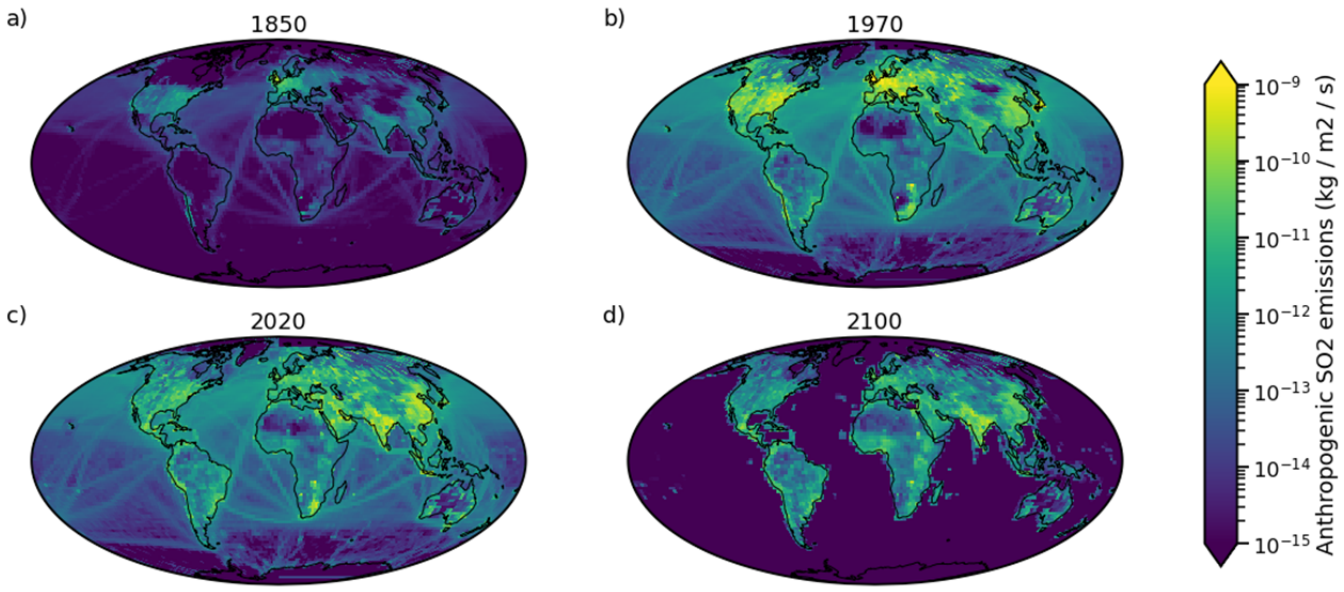

Figure 2: Maps showing the evolution of the spatial distribution of anthropogenic SO2 emissions in the pre-industrial era represented by 1850 (a); the peak emissions era of 1970 (b); current emissions (c); and future emissions under SSP 245 (d).

For the idealised CMIP simulations (*abrupt-4xCO2* and *1pctCO2*) no emissions files are used and so the cumulative anthropogenic CO2 emissions are calculated from the difference in the diagnosed CO2 atmospheric mass concentrations in these and the *piControl* experiment. Emissions of all other species are also provided but set to zero (as they represent no change since the pre-industrial).

2.2 Target ESM

We use the output from simulations performed by the NorESM2 model in its low atmosphere-medium ocean resolution (LM) configuration (Seland et al., 2020). This model consists of a fully coupled earth system with online atmosphere, land, ocean, ice and biogeochemistry components. It shares many components with the Community Earth System Model Version 2 (Danabasoglu et al., 2020) but has a replaced aerosol and atmospheric chemistry scheme (including their interactions with clouds) and a different ocean model. It has a relatively low equilibrium climate sensitivity (ECS; equilibrium global mean

temperature after a doubling of CO2) of 2.5 K, particularly compared to the 5.3K of CESM2 (Gettelman et al., 2019), which has been attributed to ocean heat uptake and convective mixing in the Southern Ocean (Gjermundsen et al., 2021). This is, nevertheless, well within the assessed likely range of ECS (90% probability between 2–5°C; Forster et al., 2021) and makes the emulation task harder than it might be for other CMIP6 models since the warming signal is weaker at the end of SSP245 (by which point CO2 is the dominant forcing), as shown in Fig. A4. The combination of a weak ECS with a relatively strong aerosol forcing (−1.36 W m^{-2} for 1850 to 2014), likely accounts for the somewhat anomalous cooling between 1950-1980 in the historical simulations (Seland et al., 2020), although it has been noted that the combined anthropogenic response in NorESM is realistic (Gillett et al. 2021).

2.3 Output variables

The output of these simulations are aggregated to annual mean values but kept at their native spatial resolution (approximately 2°). The temperature (T) and precipitation (P) are exactly equivalent to the archived surface air temperature (tas) and total precipitation (pr) output variables respectively. The DTR is calculated as the annual mean of the difference in daily maximum and minimum surface air temperatures: $|tasmax - tasmin|_{ann}$. The PR90 is calculated as the 90th percentile of the daily precipitation in each year. The annual mean baseline values (from the full *piControl* simulation) for each variable are then subtracted from each experiment so that they represent a difference from pre-industrial. Temperature changes under anthropogenic climate change are routinely reported in this way, and it also makes the downstream emulation task somewhat easier as it removes an offset. The values are not scaled to have unit variance, but users of the dataset might choose to do this with certain emulators. Many of the NorESM2 simulations include three ensemble members sampling internal variability by choosing different initial model states from the start of the piControl simulation at intervals of 30 model years apart. These are included to allow (optional) emulation of internal variability.

Samples of these output fields from the target ssp245 dataset are shown in Figure 3. The relative increase in warming in the northern polar regions (known as Arctic amplification) is clearly seen in Fig. 3a, as well as the north Atlantic warming hole (Woollings et al., 2012; Drijfhout et al., 2012; Manabe and Stouffer, 1993), the emergence of which is also affected by aerosol radiative forcing (Dagan et al., 2020). Figure 3b shows the strong land/sea contrast in DTR, since most of the change is confined to land, and largely caused by changes in aerosol (particularly sulfate) forcing. Most of the precipitation response shown in Figure 3c-d is due to the shift in the inter-tropical convergence zone (ITCZ) which results from a shift in the cross-equatorial energy balance under increased warming (Schneider et al., 2014), but some features, particularly in South-East Asia might be due to local aerosol effects (particularly due to BC; e.g., Bollasina et al. 2014, Wilcox et al. 2020, Mansfield et al. 2020).

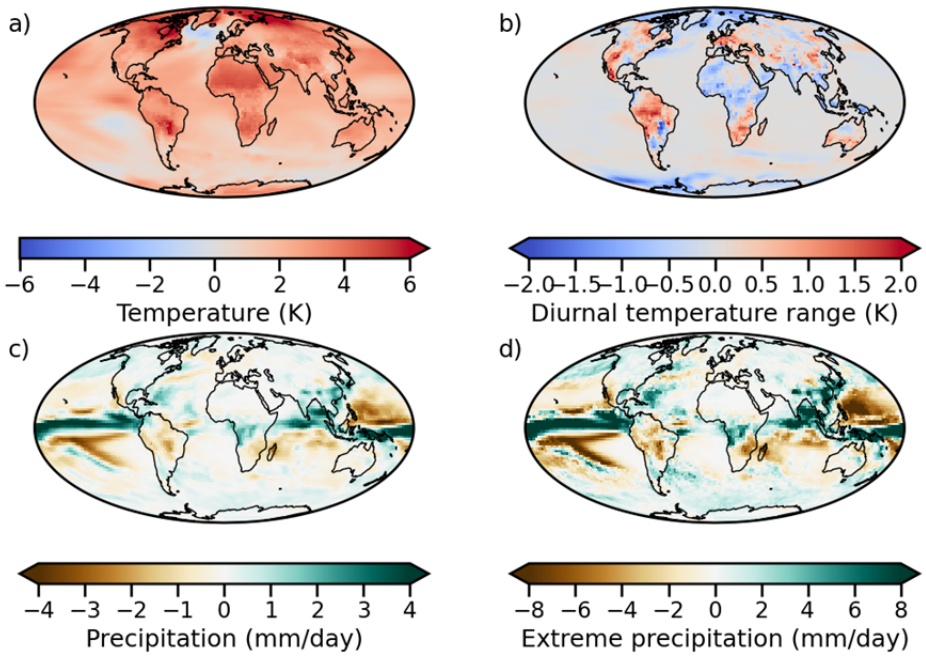

Figure 3: Maps of target outputs from the SSP245 held-back test scenario at 2100 (as an anomaly to the pre-industrial control run) performed by NorESM2: (a) Annual mean surface temperature; (b) annual mean diurnal surface temperature range; (c) annual mean precipitation; and (d) 90th percentile of the daily precipitation.

Also included in the dataset are the top-of-atmosphere Effective Radiative Forcings (ERFs) for this model for each forcing agent over the historical period. These are based on diagnostics of the fixed sea-surface temperature experiments of the Radiative Forcing Model Intercomparison Project (RFMIP; Pincus et al., 2016; Smith et al., 2021) and provide a more direct estimate of the radiative climate effect of each forcer over this period than simply emissions. It also allows an estimate of the efficacy of each forcer in this model (the temperature response per unit of forcing). This might be useful for normalising the inputs by their efficacy or developing more physically interpretable emulators that derive the climate response via the forcing, but these are not used in the present study.

3. Benchmark task

The task defined by ClimateBench is the prediction of the output variables described in Section 2.3 using only the inputs available from Section 2.1 under the chosen test scenario - ssp245. Emulators may choose to use as much or as little of the data presented in Table 1 in order to train their models as appropriate for a given approach. They may also choose to predict the contemporaneous response to emissions (as used in our RF and GP baseline emulators), account for a lagged response (as in our baseline NN emulator), or even predict the full time-series simultaneously.

3.1. Evaluation metrics

The evaluation criteria are a crucial aspect to any benchmark dataset and need to be concretely defined and accurately reflect the objectives of the machine learning task. Ideally, the criteria are also simple to implement such that they can be used as a target in any loss function that might be used to train emulators. The global mean changes in temperature and precipitation are key climatic variables but the spatial characteristics of the outputs in this task also need to be considered if the emulators are to be used for regional projections. As a primary metric we choose to combine the normalised, global mean root-mean square error ($NRMSE_s$) and the NRMSE in the global mean ($NRMSE_g$), calculated following:

$$NRMSE_s = \sqrt{\langle\left(\left|x_{i,j,t}\right|_t - \left|y_{i,j,t,n}\right|_{n,t}\right)^2\rangle} \Big/ \left|\langle y_{i,j}\rangle\right|_{t,n} \tag{1}$$

$$NRMSE_g = \sqrt{\left|\left(\langle x_{i,j,t}\rangle - \langle\left|y_{i,j,t,n}\right|_n\rangle\right)^2\right|_t} \Big/ \left|\langle y_{i,j}\rangle\right|_{t,n} \tag{2}$$

$$NRMSE_t = NRMSE_s + \alpha * NRMSE_g, \tag{3}$$

where the global mean denoted $\langle x_{i,j}\rangle$ includes a weighting function that accounts for the decreasing grid-cell area towards the poles and is defined as: $\langle x_{i,j}\rangle = \frac{1}{N_{lat}N_{lon}} \sum_i^{N_{lat}} \sum_j^{N_{lon}} cos\big(lat(i)\big)x_{i,j}$, and α is a coefficient empirically chosen to be 5 so that each component provides roughly equal weight.

Combining these commonly used metrics in this way provides a single number summarising the mismatch between the predictions (x) and the target variables (y). By squaring the difference, the RMSE also weighs large discrepancies more heavily, penalising larger errors. We average the target variables over the three available ensemble members (n) and a relatively long period of the target scenario (2080 – 2100) in order to minimise the contribution of internal variability. We choose the final years of the century since the start of the *ssp245* is quite similar to some of the training scenarios. We normalise the RMSEs so that the metrics are broadly comparable across the target variables.

Estimates of this internal variability can be very valuable for climate projections however and since ClimateBench includes three ensemble members for each training dataset emulators are encouraged to include estimates of it if they are able. A natural extension of the RMSE for probabilistic estimates commonly used in weather forecasting is the Continuous Ranked Probability Score (CRPS):

$$CRPS = \int_{x=-\infty}^{x=\infty} \big(\langle F_{i,j,t}(x)\rangle - \langle F_{i,j,t}(y)\rangle\big)^2 dx, \tag{4}$$

where $F(x)$ and $F(y)$ are the cumulative distribution functions (CDFs) over the predicted and target ensembles respectively (Gneiting et al. 2005). This measures the area between the two CDFs so that smaller values are better and has the benefit of retaining a well-defined interpretation in the case of only a single target observation (whose CDF would be the Heaviside function). The CDFs can be approximated over finite ensembles using quadrature, or direct integration if the PDFs can be assumed to be Gaussian. It should be noted that the relatively low number of ensemble members available in ClimateBench

will likely underestimate full internal variability and a larger ensemble (e.g., 100 members in Rogers et al., 2021) should be used for robust estimation. Indeed, the formulation above only includes variability in the global mean since such small ensembles are unlikely to capture regional variability. Methods to calculate both metrics based on the *climpred* (Brady and Spring, 2021) package are provided in the example notebooks included with the dataset. While this metric is not included in the headline ranking of ClimateBench approaches, we include an example approach using GPs which is discussed in more detail in Section 4.1.

3.2. Baseline evaluation

Before evaluating some baseline statistical emulators, it is useful to consider two cases with which we hope to place the skill of the data-driven approaches in a broader context. The first is the internal variability of the NorESM2 target ensemble which provides an upper bound on the predictability of the scenario in the presence of the natural variability of the Earth system. This is estimated as the standard deviation across the three NorESM2 ensemble members in *ssp245*. In practice, the emulators can (and do) outperform this baseline because they target the mean over all three ensemble members, reducing the effect of internal variability. The second is a comparison against the inter-model spread encountered within CMIP6 for the variables of interest which, despite (as discussed above) not providing a robust model uncertainty, represents a lower bound on the accuracy we would like our emulators to achieve.

Additionally, we introduce a linear pattern scaling model which uses independent linear regressions of each of the output variables at each model grid cell given the global mean temperature response to the emissions (e.g., Tebaldi et al., 2014). This approach is somewhat simpler than the other data driven models since it assumes access to an accurate impulse response (or box) model to determine the global mean temperature but provides a useful baseline. We train the regression models using the same training output data as the other emulators (described in the next section) but the only input is the global mean temperature. We assume this is available at prediction time as well so that this constitutes a 'perfect' pattern scaling approach.

4. Baseline emulators

Three baseline emulators are developed to demonstrate various potential approaches to tackling the machine learning problem this dataset provides. These are performed using the Earth System Emulator (ESEm; Watson-Parris et al., 2021) to provide a simple interface for non-ML experts and permit sampling the emulators for potential use in detection and attribution workflows (as discussed in the Section 5). All three emulators are trained using all the available training data: the historical data; ssp126; ssp370; ssp585; and the historical data with aerosol (hist-aer) and greenhouse gas (hist-GHG) forcings only, leading to 754 training / validation points (which are nevertheless not fully independent). More details on emulator specific data pre-processing, training procedure and results are described in each of the following subsections.

The emulators all perform skilfully, as summarised in Table 2 and Figure 4. The emulators also show broadly similar biases, particularly for precipitation where they all slightly underestimate increases (decreases) in tropical (subtropical) rainfall in the western Pacific. They also tend to overpredict northern-hemisphere warming while underpredicting warming elsewhere. This might suggest that these particular changes are driven by different climate forcers or longer time-scale changes than modelled in this study. A direct comparison of the emulator predictions and NorESM is shown in Figure A2. Overall, the neural network emulator performs the best in predicting temperature and precipitation changes, while the Gaussian process emulator performs best at predicting changes in the diurnal temperature range.

Table 2: The spatial, global and total NRMSE of the different baseline emulators for the years 2080-2100 against the ClimateBench task of estimating key climate variables under future scenario SSP245. The best (lowest) emulator scores for each task are highlighted in bold. The normalised standard deviation in each variable over 22 different CMIP6 models and across the NorESM ensemble members are also included as indications of inter-model and internal variability, respectively.

	NRMSE surface air temperature (1)			**NRMSE diurnal temperature range (1)**			**NRMSE precipitation (1)**			**NRMSE 90th percentile precipitation (1)**		
	Spatial	**Global**	**Total**	**Spatial**	**Global**	**Total**	**Spatial**	**Global**	**Total**	**Spatial**	**Global**	**Total**
Gaussian Process	0.109	0.074	0.478	9.207	2.675	22.582	2.341	0.341	4.048	**2.556**	0.429	4.702
Neural Network	**0.107**	**0.044**	**0.327**	9.917	**1.372**	**16.778**	**2.128**	**0.209**	**3.175**	2.610	**0.346**	**4.339**
Random Forest	0.108	0.058	0.400	**9.195**	2.652	22.457	2.524	0.502	5.035	2.682	0.543	5.399
Pattern Scaling	0.080	0.048	0.320	8.083	2.327	19.719	2.006	0.331	3.662	2.400	0.412	4.461
Variability	0.052	0.072	0.414	2.513	1.492	9.973	1.350	0.268	2.691	1.757	0.457	4.043
CMIP6	-	0.206	-	-	0.815	-	-	0.417	-	-	-	-

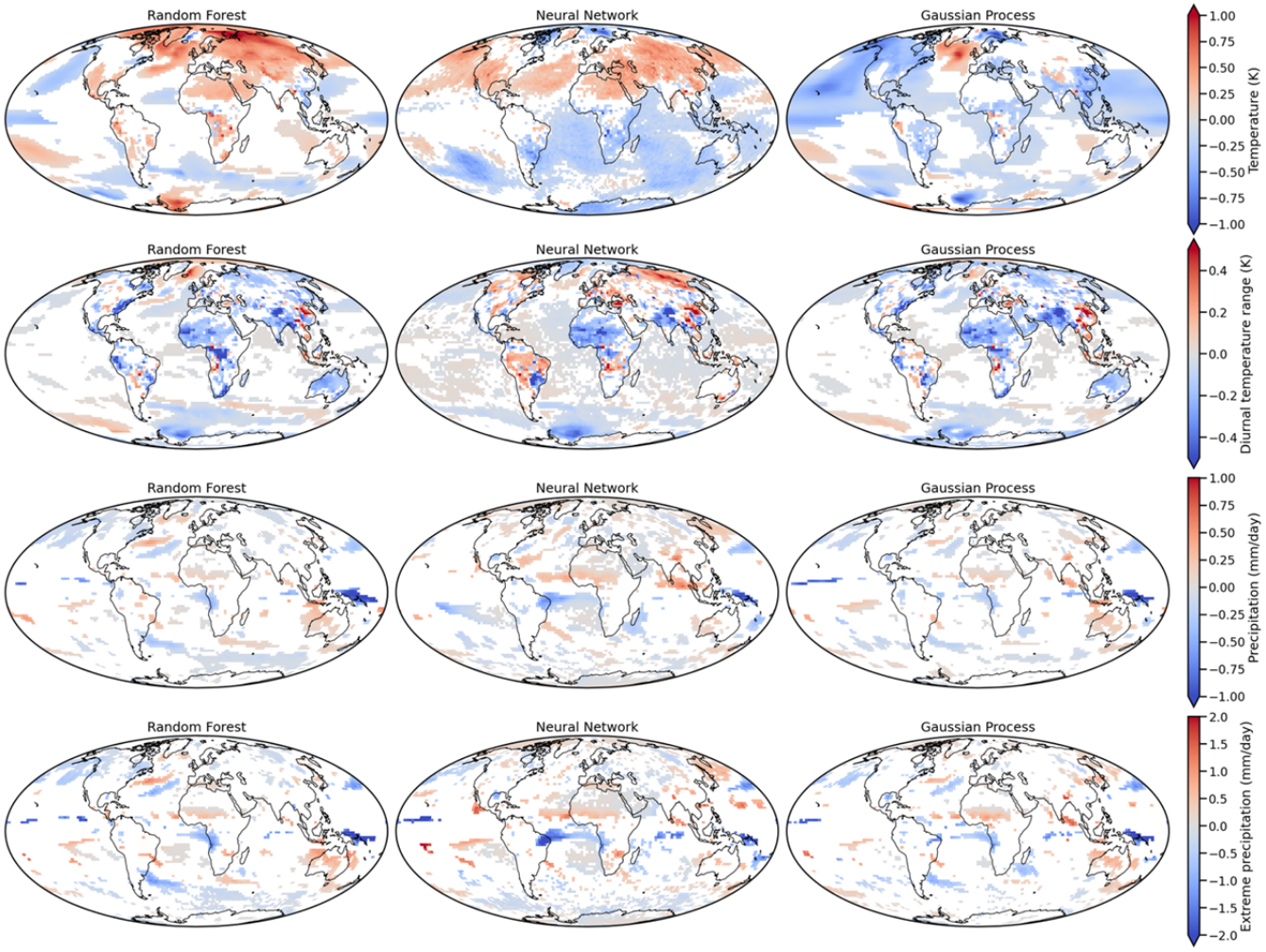

Figure 4: Maps of the mean difference in the ClimateBench target variables for each baseline emulator against the target NorESM values under the test ssp245 scenario averaged between 2080-2100. Differences insignificant at the p<5% level are masked from the plots.

4.1. Gaussian process regression

Gaussian processes (GPs) (Rasmussen and Williams, 2005) are probabilistic models which assume predictions can be modelled jointly as normally distributed. GPs have been widely used for nonlinear and nonparametric regression problems in the geosciences (Camps-Valls et al., 2016). A GP is fully determined by the expectation of individual predictions – referred to as the mean – and the covariance between pairs of predictions. Such covariance is typically user-specified as a bivariate function of the input data called the kernel function. The choice of the kernel function allows to restrict the functional class the GP belongs to, offering, for example, control over functional smoothness. GPs for regression solve a supervised problem

where the observed input-output sample pairs are used to: (1) infer the emulator parameters (typically only the noise variance and the kernel parameters) by maximising the log-likelihood of the observations under the evidence; and then (2) allow to obtain its posterior probability distribution that is used to make predictions over unseen inputs.

To prepare the input samples, the dimensionality of the SO2 and BC emission maps are reduced with principal component analysis, and we only use the 5 first principal components of each as inputs, corresponding to 96% and 98% of the explained variance, respectively. All input covariates and target outputs are standardised using training data mean and standard deviation.

The GP is set with a constant mean prior and separate kernels are devised for each species. Automatic relevance determination (ARD) kernels are used for SO2 and BC, allowing each principal component to be treated independently with its own lengthscale parameter. The GP covariance function is obtained by summing all kernels together, thus accounting for multiscale feature relations (see Camps-Valls et al., 2016 for several composite kernel constructions in remote sensing and geoscience problems). To account for internal variability between ensemble members, we consider an additional white noise term with constant variance over the output targets, which is also inferred from the training phase.

We use Matérn-1.5 kernels for each input. This guarantees the GP is a continuous, once differentiable function; details are provided in Section A1. The mean value, kernels parameters and internal variability variance are jointly tuned against the training data by marginal likelihood maximisation with the limited memory Broyden–Fletcher–Goldfarb–Shanno (L-BFGS) optimisation algorithm. The emulators used have 18 parameters in total: 5 lengthscale and one variance parameter for each aerosol kernel; one lengthscale and one variance for each of the GHG kernels; one mean and one likelihood variance.

As reported in Table 2, the total NRMSE of the mean predictions with the GPs are competitive with the neural network for all the variables. This is remarkable given the limited number of parameters that are learnt. It suggests the GP prior is an adequate choice for the purposes of emulation. Study of the inferred kernel variance (not shown) suggests that cumulative CO2 emissions generally influence all predictions, and unequivocally dominate the predictions for surface air temperature and diurnal temperature range. CH4 and BC emissions on the other hand appear to have negligible influence on the predictions. Since the GP also provides posterior estimates of the variance (which will incorporate an estimate of internal variability) we also calculate the CPRS for this emulator (see Table A5). While we are unable to compare these scores with the other baseline methods the similarity to the global NRMSE indicates that the GP is also predicting the internal variability accurately (otherwise it would be penalised in the CPRS relative to the NRMSE).

4.2. Random Forests

Random forests aggregate predictions of multiple decision trees (Ho, 1995; Breiman, 2001). These trees repeatedly split data into subsets according to its features such that in-subset variance is low and between-subset variance is high. This makes decision trees good at modelling non-linear functions, in particular interactions between different variables. However, they

are prone to overfitting (Ho, 1995). This problem is alleviated by ensemble methods which train a large number of different trees. Weak learners are combined to give strong learners. Bagging, used in Random Forests, describes training different trees on different subsets of the data or holding back some of the data dimensions for each individual tree. The Forest makes a prediction by averaging over the predictions of all individual trees.

Two main arguments support an ensemble method approach to climate model emulation: These methods are skillful at interpolation tasks, but by construction are unable to extrapolate (Breiman, 2001). However, for applications of climate model emulation, interesting predictions will likely lie inside the hypercube delimited by historical data, low-emissions (*ssp126*) and business-as-usual (*ssp585*) scenarios. A major advantage of ensemble methods over more complex ML methods such as neural networks (and even ESMs) is their interpretability. This is important as ultimately predictions should inform decision-making. Being able to provide explanations why a given input led to a prediction helps to understand the consequences of decisions about emission pathways.

Analogously to the GP emulator, the dimensionality of aerosol emission maps is reduced with principal component analysis. The first five principal components of SO2 and BC together with the global emission maps of CO2 and CH4 form the input features of the model. Separate random forest emulators are trained for the four target variables. The following hyperparameters are tuned using random search of the training data without replacement: number of trees, tree depth, number of samples required to split a node and to be at each leaf node. The hyperparameters used for each emulator are indicated in Section A2.

As shown in Table 2, the spatial NRMSE scores of the random forest regressors are comparable to the performance of the other emulators for all variables but the global NRMSE is significantly worse for temperature and precipitation (as can also be seen in Figure 6). Discontinuities in the predicted global mean temperature change time series over this period (not shown) perhaps indicate a deeper tree structure is required. To assess the impact of the four input features on the prediction, we calculate the permutation feature importance. It is defined as the decrease in a model score when a single feature value is randomly shuffled (Breiman, 2001). Figure 5 shows that CO2 concentrations dominate the predictions. For temperature predictions the other features are negligible. SO2 and BC aerosol emissions have a small impact on the global mean temperature and precipitation predictions. This is in line with the physical understanding that while anthropogenic aerosol can influence precipitation rates (both radiatively and through aerosol-cloud interactions), aerosol contributions play a negligible role at the end of the century in the *ssp245* test scenario. The regional influences may be more significant however and this will be explored separately.

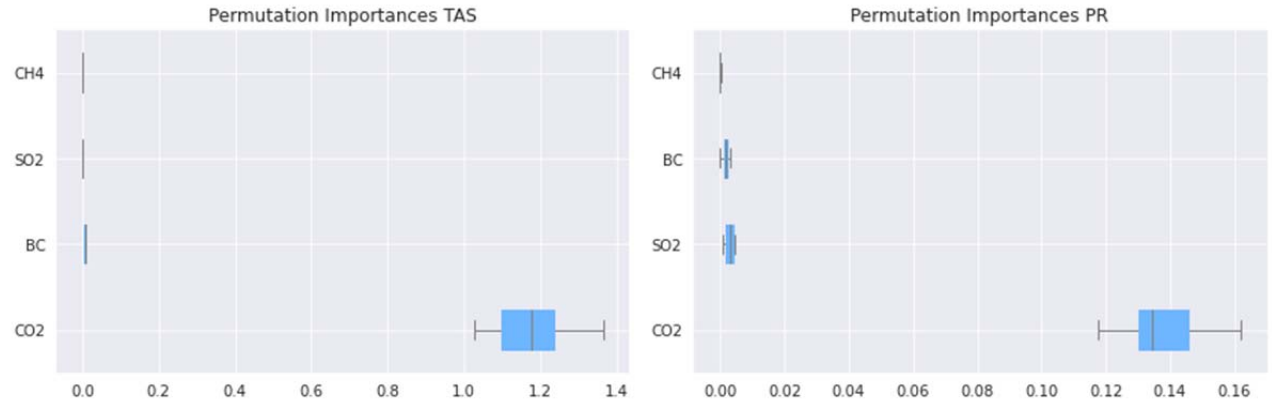

Figure 5: Permutation importances for the most important component of each variable in predicting global mean temperature (TAS) and precipitation (PR). Each emulator input variable is shuffled in turn to determine the relative contribution to prediction skill. Note that these average estimates do not account for potential regional contributions which may be particularly relevant for aerosol.

4.3. Neural Networks

Artificial Neural Networks (ANNs) are algorithms inspired by the biological neural networks of human brains that have shown outstanding success in areas like Computer Vision and Natural Language Processing. Two major ANN architectures are Convolutional Neural Networks (CNNs) (LeCun et al., 1990), that are able to model spatial dependencies, and Recurrent Neural Networks (RNNs), that are able to process time series and sequential data. ANNs have recently been employed to tackle a variety of problems in earth system science (Camp-Valls et al., 2021). CNNs are helpful for modeling climate data with a spatial structure, for instance, precipitation patterns or satellite imagery, and are frequently applied in climate science and weather forecasting (Trebing et al., 2020, Harder et al., 2020). Long short-term memory (LSTM) networks (Hochreiter et al., 1997), an advanced type of RNNs, have proven skillful for modeling climate time series, for example, for the prediction El Niño-Southern Oscillation (Broni-Bedaiko et al., 2019).

For time series of spatial variables, as in the ClimateBench dataset, we can use the two types of networks in sequence to model both spatial and temporal dependencies. The chosen architecture consists of a CNN followed by an LSTM built with the Keras library. The CNN includes one convolutional layer with 20 filters, a filter size of 3, and a ReLU activation function. The 3x3 pixel filters scan the input images to detect spatial patterns and feed these patterns to the next layer. These next layers are average pooling layers that reduce the spatial dimensionality ahead of the LSTM layer. The LSTM uses 25 units (i.e., the output dimension of each LSTM cell) and a ReLU activation function. The LSTM is followed by a dense layer and reshaping layer to (96, 144), i.e., the (latitude, longitude) dimension of the output variables.

The training data time-series is segmented into 10-year chunks, using a moving-time window in one-year increments, leading to 754 training samples of shape (10, 96, 144, 4) corresponding to the number of years, latitude, longitude and then number of variables. We trained four different emulators for the four different output variables. Each emulator is trained for 30 epochs, using a batch size of 16. For this baseline approach, we chose not to do any hyperparameter optimization, and all the parameters were chosen manually.

RMSE scores obtained with the CNN-LSTM architecture are somewhat better than those achieved with the other methods, particularly in the global-mean. This might be because the LSTM is able to better capture the temporal autocorrelation than the other emulators which treat the prediction instantaneously. The CNN-LSTM architecture also captures spatial changes in temperature well (e.g., the Arctic amplification), even though warming at the poles is somewhat underestimated. In general, warming in the Northern hemisphere is overestimated while it is underestimated in the Southern Hemisphere. Given the overestimated temperature response in the *ssp245-aer* simulations shown in Figure A3, this may be due to an overestimation of the effect of aerosol on the temperature by this emulator. The diurnal temperature range is well predicted, with a lower performance over land. The CNN-LSTM also captures spatio-temporal changes in precipitation (e.g., the ICTZ shift) quite well.

5. Discussion

5.1. Climate-specific challenges

The emulation of future climate states presents particular challenges for machine learning and other statistical approaches. Chiefly among those is the limited amount of training data that is typically available; current ML approaches are not prepared to learn such complex scenarios in small data regimes under a covariate shift. As pointed out, the complex ESMs that are trusted to model the future climate are extremely computationally expensive to run and the observational record cannot inform us about unseen future scenarios. By harnessing a large selection of simulations performed as part of CMIP6, ClimateBench attempts to alleviate this difficulty, but nevertheless only around 500 training points (years) represent realistic climate states, many of which are not independent (as shown in Fig. A1). This presents a challenge for deep learning approaches which typically require tens of thousands of training samples to avoid over-fitting. The inclusion of longer idealised simulations does provide opportunities for pre-training however, particularly the 500-year long *piControl* simulations which could be used with contrastive learning to reduce the training samples required for neural network architectures.

The *piControl* simulation could also be used to inform emulators more explicitly about the internal variability of climate (as produced by NorESM2). The signal, particularly for the precipitation target variables, can be small compared to this variability and this proves challenging for some emulators to reproduce. An explicit model of the internal variability (Castruccio et al., 2019) could help to alleviate this.

Another challenge in applying statistical learning approaches to this dataset is the relatively high dimensional inputs and outputs (96 x 144). Most approaches to emulating the regional temperature response to a CO2 forcing have been carried out at, at most, dozens of locations, but accounting for the spatial correlations is something which CNNs can excel at and have recently been shown to produce accurate emulations of temperature across similar dimensionality (Beusch et al., 2020). Such approaches typically assume a regular spacing, however, and neglect the reducing area of each grid-cell towards the poles.

While more traditional approaches of dimensionality reduction can also be used, such as (weighted) empirical orthogonal functions (EOFs), these may not be appropriate for the non-linear precipitation fields which might require kernel-based approximations (e.g., Bueso et al., 2019).

For practical purposes, an estimate of the uncertainty in any prediction would be extremely valuable. This uncertainty should encompass that due to the internal variability and the emulator approximation (and ideally that of the underlying physical model). In the ML community, these are known as the epistemic and the model uncertainties, and are being studied intensively (Kendall et al., 2017). Quantifying these two uncertainties would allow increased trust (a concept explored in the next section) in the prediction as well as quantitative comparison to other predictions. We encourage the estimation of uncertainty wherever possible, using the provided CRPS metric to evaluate such probabilistic projections. The ability to sample from such distributions would also permit the generation of so-called 'superensembles' which can provide very large ensembles of multiple models under given scenarios (Beusch et al., 2020).

As previously discussed, and shown in Figure A4, there is large inter-model variability in the projected climate variables in CMIP6, even across a single scenario. Future work should explore the ability of a given emulator to robustly recreate each of these model responses, and could allow a deeper understanding of their discrepancies.

5.2. Emulator trustworthiness

For climate model emulators to be useful for policy decisions they must be trusted by their users. The trustworthiness of any model is a subjective concept that broadly represents one's belief that the model faithfully represents some underlying 'truth'. Model verification attempts to objectively assert this view (indeed the word derives from the Latin, *verus,* meaning true) but is formally impossible for an open system like the Earth (see e.g., Oreskes et al., 1994). While weather models can be regularly validated against observations, in the climate sciences we often instead resort to necessarily incomplete model evaluation and rely on underlying physical principles to provide reassurances of broader validity. The ClimateBench emulators side-step this issue by aiming only to accurately reproduce an existing physical model which is assumed to already be well evaluated, and therefore attain trustworthiness through proxy. It would nevertheless be reassuring if the emulators could be demonstrated to respect some of the same physical constraints.

In this spirit, Figure 6 shows the relative change in global mean precipitation as a function of global mean temperature change (the hydrological sensitivity) of the baseline emulators and NorESM2. While locally precipitation can change in accordance with the Clausius-Clapeyron relationship (6-7% / K), energy conservation requires that the global changes in precipitation are balanced by radiative cooling and limited to 2-3% / K (Allen and Ingram, 2002; Pendergrass and Hartmann, 2013; Jeevanjee and Romps, 2018; Dagan et al., 2019). While the RF emulator underestimates the hydrological sensitivity of NorESM, it is clear that the emulators learn the physical relationship from the underlying model. Since the emulators were trained on the precipitation and temperature this is to be expected to some degree, but this demonstrates the principle that emulators trained correctly can retain the physical laws of the underlying models over the range of their training data. Future

efforts to introduce these invariances directly have the potential to significantly ease the training and improve the inference of climate model emulators (Beucler et al., 2021), ultimately improving their trustworthiness.

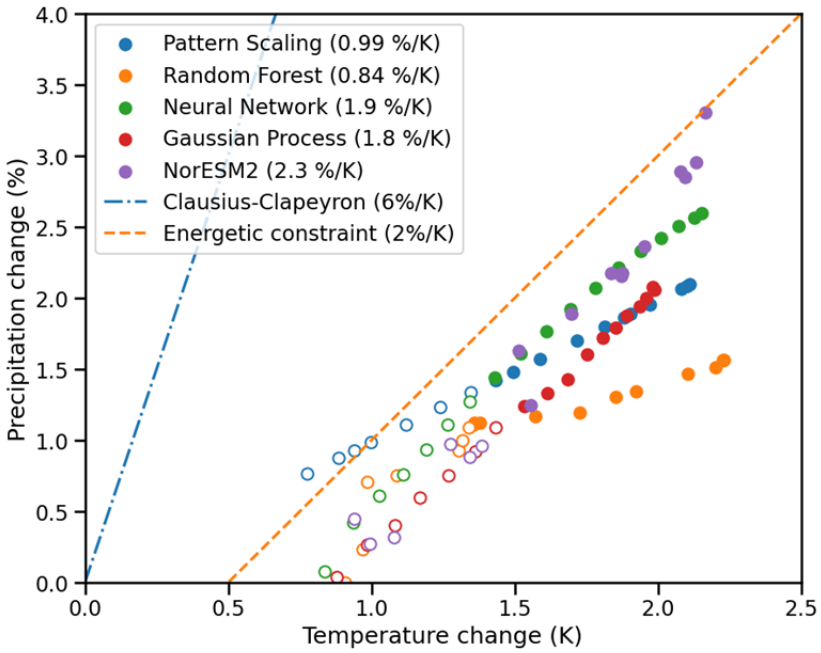

Figure 6: The relative change in global mean precipitation as a function of global mean temperature change in the baseline emulators and NorESM2 averaged in 5-year increments to reduce internal-variability. Hollow and solid points indicate years before and after 2050 respectively. The change predicted by the Clausius-Clapeyron relationship and energy conservation considerations are shown as dashed lines.

There has been much attention recently given to 'interpretable' and 'explainable' machine learning models, the former of which are said to behave in a-priori understandable ways (Barnes et al., 2020), while the latter provide mechanisms to determine post-hoc understanding (McGovern et al., 2019). While not as robust as physical laws, these techniques provide useful indications that such models are getting the right answer for the right reasons. Indeed, the physical ESMs currently considered the 'gold standard' of climate modelling are often only interpretable or explainable by expert practitioners and it is hoped that (interpretable) ClimateBench emulators will be useful in analysing and understanding the response of the underlying physical models themselves.

5.3. Research opportunities

While the challenges outlined above are mostly surmountable with modern architectures and carefully chosen workflows, there are also several broader opportunities ClimateBench presents to develop the state-of-the-art in climate model emulation.

As already mentioned, one area of particular interest is the use of hybrid modelling whereby statistical or ML based emulators embed physical equations, constraints or symmetries in order to improve accuracy, robustness and generalisability (Camps-Valls et al., 2021; Reichstein et al., 2019; Karpatne et al., 2017). One obvious way in which to apply such approaches to ClimateBench is to marry the simple impulse response models discussed in Section 1 with more complex methods to predict the spatial response. Such an approach has recently been demonstrated for temperature (Beusch et al., 2021) but could conceivably be extended to modelling each of the fields targeted in ClimateBench. A more unified, and ambitious, approach would be to model the ordinary differential equations of the response to a forcing directly in the statistical emulator using either numerical GPs (Raissi et al., 2018) or Fourier neural operators (Li et al., 2020).

Another important open question when using data-driven approaches to emulate the climate is how to ensure predictions are performed at locations within the distribution of the training data. In other words, how to ensure the emulator is being used to interpolate existing model simulations rather than extrapolating to completely unseen regions of input space. This can be easy to test for in low dimensions, but it becomes increasingly difficult in higher dimensions and while the training and test data in ClimateBench have been chosen to minimise the risk of extrapolation broader use could be hindered by the risk of inadvertently asking for an out-of-distribution prediction. While the predictive variance of GPs provide such indications (out of the sample range the GP mean returns to the prior and the covariance is maximised), it is not so easy for other techniques and the use of modern techniques to detect such occurrences (e.g., Lee et al., 2018; Rabanser et al., 2018) could be of great value to minimise this risk.

5.4. Application to detection and attribution

The use of an efficient and accurate way of estimating the climate impacts of different emission scenarios is not limited to exploring future pathways. We may also ask: ‘What observed climate states and events can be attributed to anthropogenic emissions?’. A whole field, which started with the seminal work of Hasselmann (1993) has developed rapidly in the last decade (Stott et al., 2016; Barnett et al., 2005; Stott et al., 2010; Shindell et al., 2009; Otto et al., 2016) attempting to answer this question. A common approach is to use climate model (or ESM) simulations to determine optimal ‘fingerprints’ with which to test observations as well as the power of such a fingerprint under internal variability. These typically have to make fairly strong assumptions about the form of the climate response however (often relying on multiple linear regression) and can incorporate observations of only a few dimensions.

One possible application of the efficient emulators trained using ClimateBench could then be to allow the inference of higher dimensional attribution problems, incorporating more information (such as the DTR and PR) and potentially providing more confident assessments. It would be straightforward to implement such an approach using the ESEm package which provides a convenient interface for such inferences using e.g., approximate Bayesian computation, variational inference or Markov Chain Monte-Carlo sampling. Future work will investigate these possibilities.

As a simple demonstration of potential of such an approach we have included a prediction by the emulators compared to the original NorESM2 simulations of the *ssp245-aer* DAMIP experiment in which only the aerosol species are emitted, shown in Figure A3. This is a more challenging scenario than the *ssp245* test case due to the much smaller total forcing and the emulators do not perform as well (see NRMSE in Table A4). It is interesting to note that the emulators particularly struggle with temperature changes in the North Atlantic where slow ocean circulation changes (e.g., Dagan et al. 2020) may not be fully captured. They nevertheless capture the main features of the response and show promise for future work disentangling the forcings and feedbacks in NorESM2, other ESMs and ultimately observations.

6. Conclusions

The application of machine learning to the prediction of future climate states has, perhaps justifiably due to the challenges laid out above, been cautious to date. Particular applications however, with carefully chosen training data and objectives, can provide fruitful avenues for research and open exciting opportunities for improvement over the current state-of-the-art. This paper introduces the ClimateBench dataset in order to galvanise existing research in this area, provide a standard objective with which to compare approaches and also introduce new researchers to the challenge of climate emulation. It provides a diverse set of training data with clear objectives and challenging target variables, some of which have been extensively studied (surface air temperature) and some which have been somewhat neglected (diurnal temperature range and precipitation).

We also introduce three quite distinct approaches for undertaking this challenge: a random forest; a Gaussian process; and a neural network model. These different models are based on different principles, have distinct assumptions and rely on quite different learning paradigms. Each has their strengths and weaknesses but all perform well in the evaluation metrics and generally reproduce the NorESM2 temperature and precipitation response well in a realistic (but unseen) future scenario, especially compared to CMIP6 inter-model diversity. The neural network model performs best overall and shows good skill both in the global mean and spatially. All the models perform less well in the aerosol only test, suggesting that they have not fully learnt the distinct response due to each forcer and future emulators should aim to rectify this.

Current impact assessments are often based on simple emulators, which are then scaled to match modelled patterns, but which are unable to predict non-linear responses in e.g., precipitation. A robust, trustworthy emulator which is able to provide such predictions could be immensely valuable in quantifying and understanding the changes and associated risks of different socio-economic pathways. Given the importance of faithfully and accurately reproducing the response of ESMs, we hope the challenge will also spur innovation in nascent physically informed ML techniques.

In order to meet these objectives, we have provided open, easy to access datasets and training notebooks which reproduce the results shown in this manuscript and demonstrate the use of the different baseline emulators. All software is open-source and readily available using commonly used package managers. We hope this dataset will provide a focus for climate and ML

555 researchers to advance the field of climate model emulation and provide policy makers with the tools they require to make
556 well informed decisions.

557

5. Data and code availability

The baseline code is available on GitHub (https://github.com/duncanwp/ClimateBench) and a DOI for the specific version, including that used to generate the plots in this paper, will be made available on acceptance.

The benchmark data is available here: https://doi.org/10.5281/zenodo.5196512. The raw CMIP6 data used here are available through the Earth System Grid Federation and can be accessed through different international nodes e.g.: https://esgf-index1.ceda.ac.uk/search/cmip6-ceda/.

6. Author Contributions

DWP conceptualised ClimateBench and performed the data-processing and initial analysis. YD and PN contributed to the definition and setup of the framework. DO and ØS performed the original NorESM2 simulations used for training. SB, MD, EF, PH, KJ, JL, PM, MN, LR, CR and JV developed the baseline emulators. DWP prepared the manuscript with contributions from all co-authors.

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9. Appendix 1

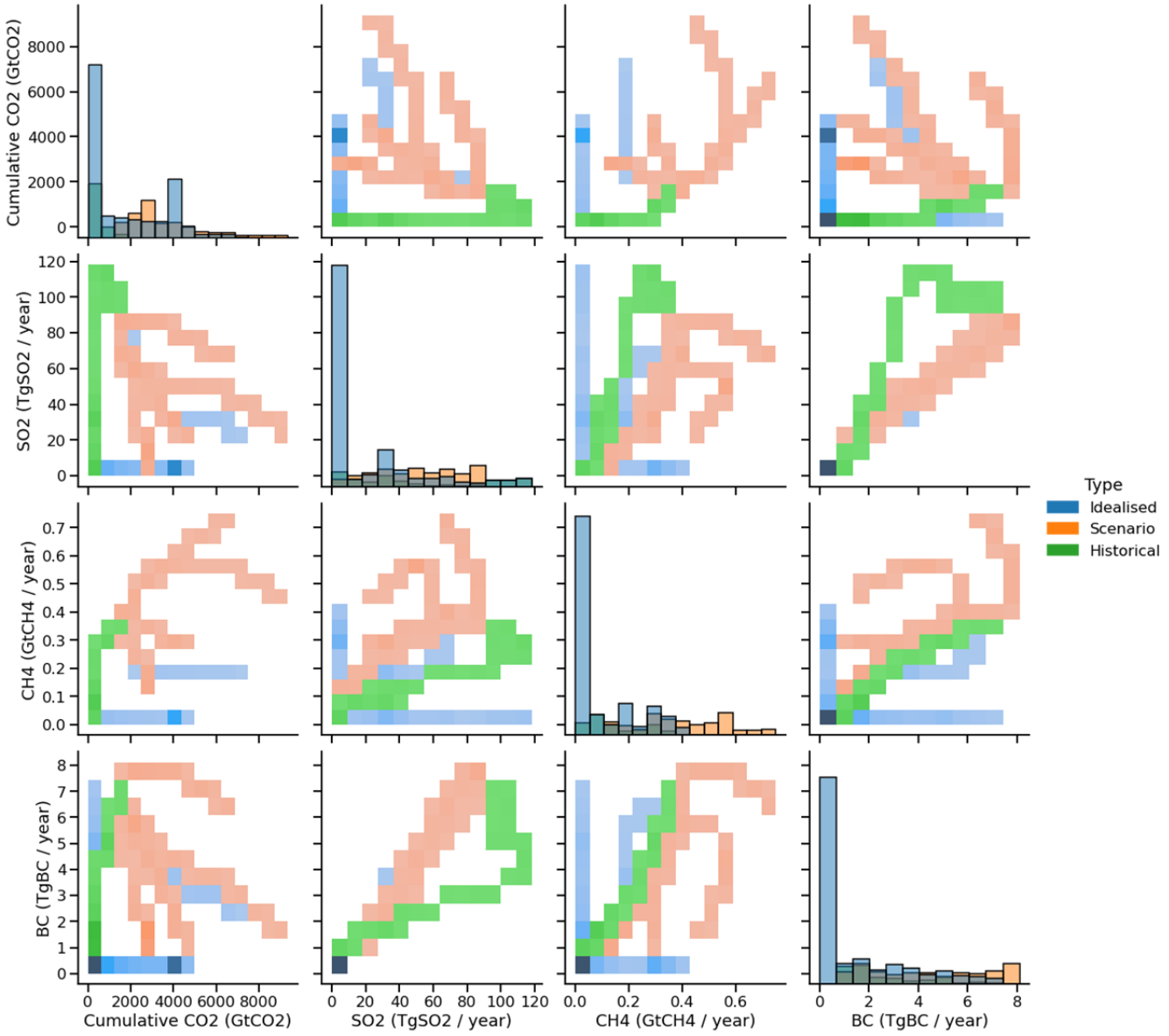

Figure A1: Joint and marginal distributions of annual global mean emissions and concentrations across the ClimateBench training dataset. Input datasets are classified as Idealised (such as *1pctCO2* and *abrupt4xCO2,* and including *ssp370-lowNTCF*), Historical and Scenario to demonstrate the contribution of each to sampling the full input space.

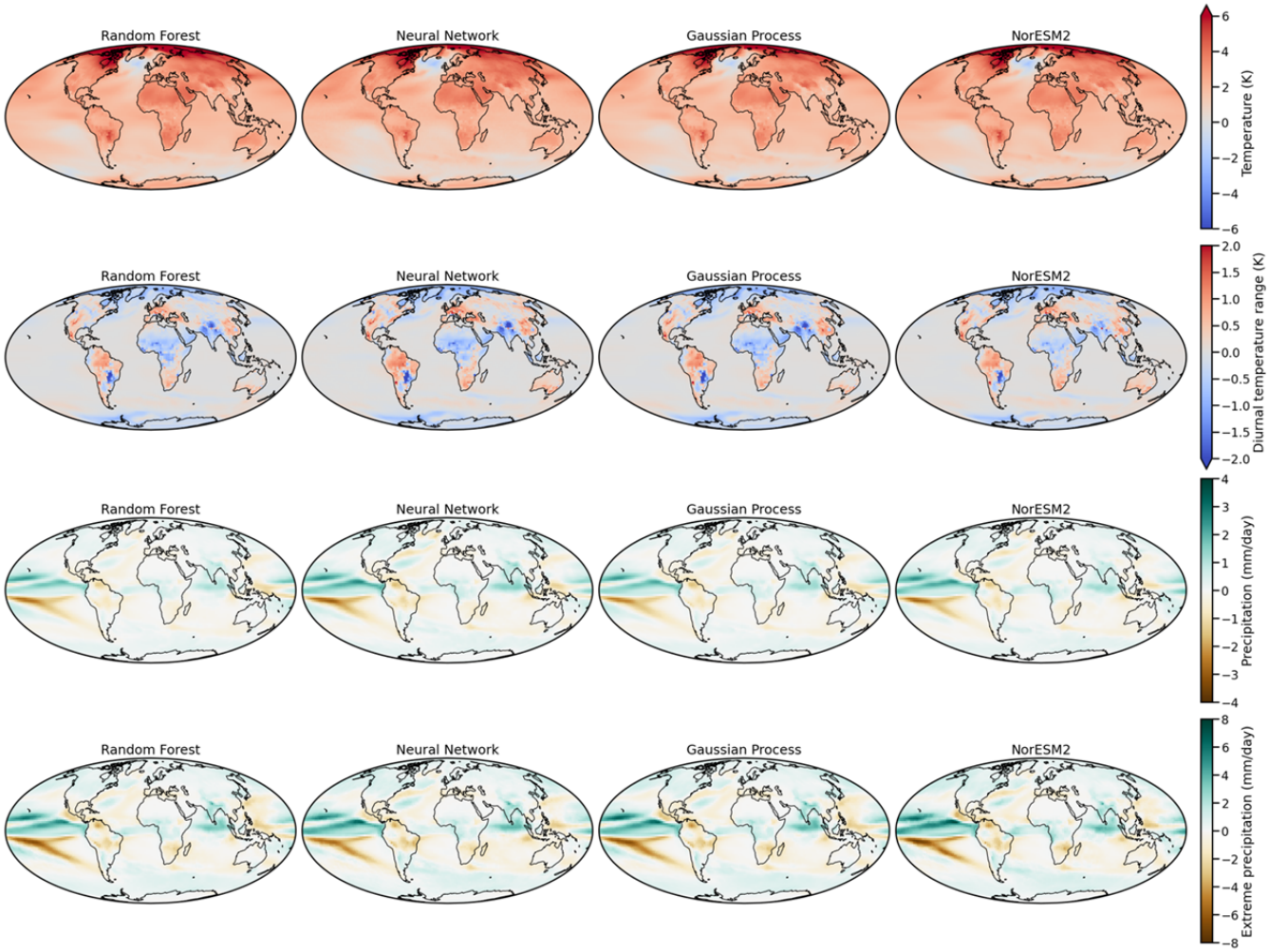

Figure A2: Maps of ClimateBench target variables for each baseline model and the target NorESM values under the test *ssp245* scenario averaged between 2080-2100.

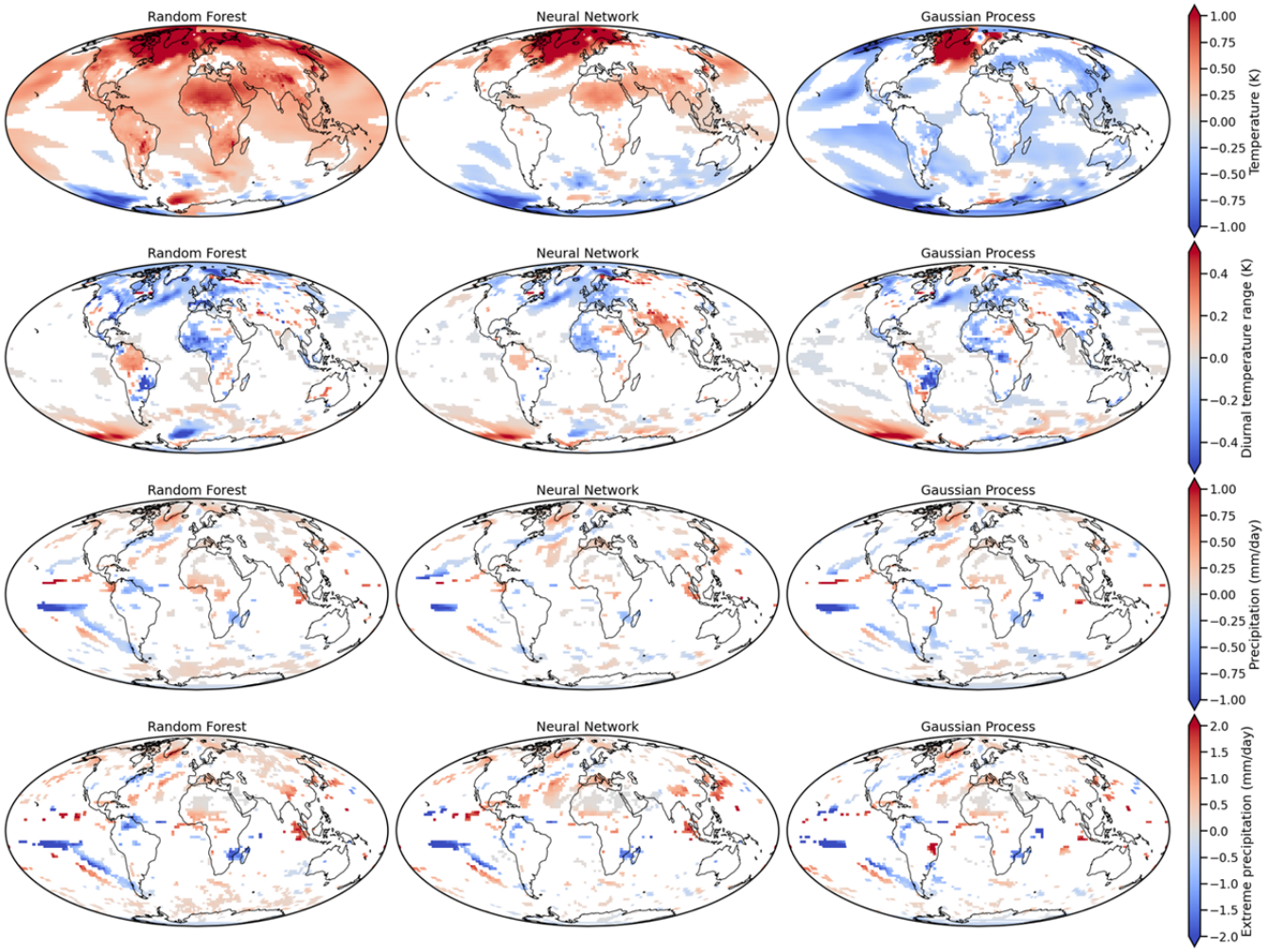

Figure A3: Maps of the mean difference in the ClimateBench target variables for each baseline emulator against the target NorESM values under the test ssp245-aer scenario averaged between 2080-2100. Differences insignificant at the p<5% level are masked from the plots.

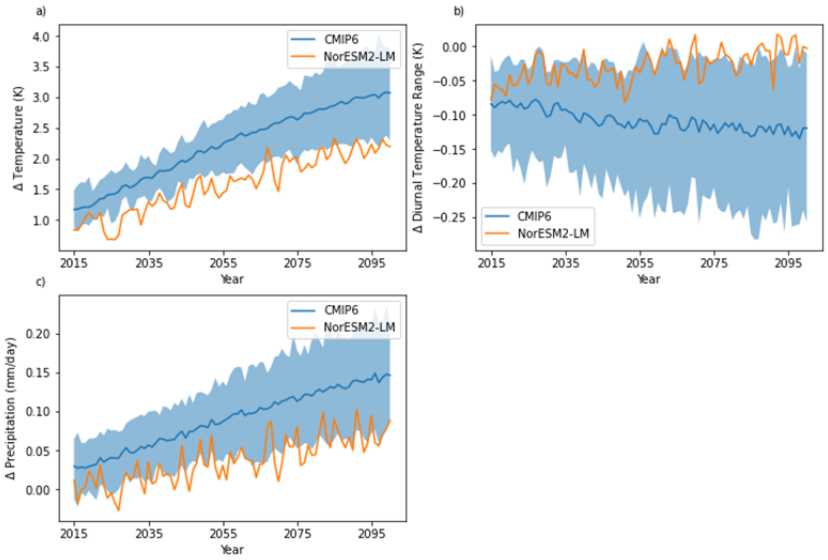

Figure A4: Global mean NorESM-LM projections under *ssp-245* as compared to all other available CMIP6 models for three of the target variables

Table A4: The spatial, global and total NRMSE of the different baseline emulators for the years 2080-2100 against the ClimateBench task of estimating key climate variables under the idealised future scenario SSP245-aer.

	NRMSE surface air temperature (1)			NRMSE diurnal temperature range (1)			NRMSE precipitation (1)			NRMSE 90th percentile precipitation (1)		
	Spatial	Global	Total	Spatial	Global	Total	Spatial	Global	Total	Spatial	Global	Total
Gaussian Process	2.138	1.165	7.963	14.298	2.868	28.636	12.100	**0.933**	16.767	13.486	**1.353**	20.252
Neural Network	**2.116**	**1.011**	**7.173**	**12.387**	**2.200**	**23.386**	**10.316**	0.977	**15.199**	**12.224**	1.438	**19.414**
Random Forest	2.977	2.041	13.182	16.222	3.284	32.642	11.562	1.291	18.017	12.302	1.616	20.382

Table A5: The CRPS for the Gaussian process emulator for the years 2080-2100 against the ClimateBench task of estimating key climate variables under future scenario SSP245.

	CRPS surface air temperature (K)	CRPS diurnal temperature range (K)	CRPS precipitation (mm / day)	CRPS 90th percentile precipitation (mm / day)
Gaussian Process	0.4765	0.3601	1.0753	1.0029

1. Gaussian process model specifications

The GP models kernel *k* have the same form for all four climate response variables

$$k = k_{CO2} + k_{CH4} + k_{BC} + k_{SO2}$$

where k_{CO2} and k_{CH4} are kernels that respectively take as inputs CO2 and CH4 emissions. k_{BC} and k_{SO2} are kernels that take as inputs the 5 principal components of BC and SO2 emission maps respectively, each principal component being rescaled by an independent length scale term. We choose the Matérn-1.5 class of kernel,

$$k_X(x, x') = \left(1 + \sqrt{3}d(x, x')\right) exp\left(-\sqrt{3}d(x, x')\right),$$

where X is a general notation for CO2, CH4, BC or SO2, and $d(x, x')$ is a distance between inputs typically given by

$$d(x, x') = \sum_i \quad |x_i - x_i'|/l_i.$$

l_i is a length scale associated to the i^{th} coordinate x_i. Global CO2 and CH4 emissions are scalar inputs, hence the corresponding distances only involve one length scale parameter. The principal components decompositions of BC and SO2 emission maps both have 5 coordinates, hence we set each principal component to be a different coordinate with its own length scale parameter. The Matérn-1.5 kernel guarantees that the corresponding GP lies in a space of continuous functions, hence providing regularity to the climate response predictions. We refer the reader to Rasmussen and Williams, 2005, Chapter 4 for more details on the Matérn kernel. Each kernel is multiplied by a variance term ${\sigma_X}^2$, which rescales the kernel in the above sum and allows to balance relative features importance. Variances and length scales are tuned during the optimization step.

2. Random forest model specification

Hyperparameter	number of trees	min samples split	min samples leaf	maxdepth
Surface air temperature	250	5	7	5
Diurnal temperature range	150	15	8	40
Precipitation	250	15	12	25
90^{th} percentile of precipitation	300	10	12	20

3. Neural Network model specification

The parameters are the same for all four models.

Model architecture

Layer	Hyperparameter value (if not specified, the default parameters are used)	Output Shape	Param #
Time distributed Conv2D	Number of filters: 20 Filter size: 3 Activation function: ReLu	(None, 10, 96, 144, 20)	740
Time distributed AveragePooling2D	Pool size: 2	(None, 10, 48, 72, 20)	0
Time distributed GlobalAveragePooling2D		(None, 10, 20)	0
LSTM	Number of units: 25 Activation function: ReLu	(None, 25)	4600
Dense	Units: 96*144	(None, 13824)	359424
Activation	Activation function: linear	(None, 13824)	0
Reshape		(None, 1, 96, 144)	0

Model training

Hyperparameter	Value

Batch size	16
Epochs	30
Optimizer	Rmsprop
Metric	MSE