Parameter Uncertainty Quantification in an Idealized GCM With a Seasonal Cycle

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Abstract Climate models are generally calibrated manually by comparing selected climate statistics, such as the global top-of-atmosphere energy balance, to observations. The manual tuning only targets a limited subset of observational data and parameters. Bayesian calibration can estimate climate model parameters and their uncertainty using a larger fraction of the available data and automatically exploring the parameter space more broadly. In Bayesian learning, it is natural to exploit the seasonal cycle, which has large amplitude compared with anthropogenic climate change in many climate statistics. In this study, we develop methods for the calibration and uncertainty quantification (UQ) of model parameters exploiting the seasonal cycle, and we demonstrate a proof-of-concept with an idealized general circulation model (GCM). UQ is performed using the calibrate-emulate-sample approach, which combines stochastic optimization and machine learning emulation to speed up Bayesian learning. The methods are demonstrated in a perfect-model setting through the calibration and UQ of a convective parameterization in an idealized GCM with a seasonal cycle. Calibration and UQ based on seasonally averaged climate statistics, compared to annually averaged, reduces the calibration error by up to an order of magnitude and narrows the spread of the non-Gaussian posterior distributions by factors between two and five, depending on the variables used for UQ. The reduction in the spread of the parameter posterior distribution leads to a reduction in the uncertainty of climate model predictions.

Plain Language Summary Climate models rely on empirical representations of physical processes that cannot be resolved with available computational resources. Empirical representations of physical processes, such as turbulence and cloud physics, reduce the computational cost of simulations, but introduce new unknown parameters into the climate model. The unknown parameters contribute to uncertainties associated with climate model predictions. Historically, fixed values of the model parameters have been hand-tuned using scientific intuition and a limited amount of available data. We develop methods for the computationally efficient estimation of the unknown climate model parameters and their uncertainty from data, by using optimization and machine learning. Many processes and observable statistics of Earth’s climate used to produce this data are influenced by seasonal variations. We demonstrate that the incorporation of seasonal information into these statistics significantly improves the resulting calibration of climate model parameters, in contrast to using annually averaged information alone. We show that including seasonal information also reduces the uncertainty associated with the model parameters, which consequently reduces the uncertainty of climate model predictions.

1. Introduction

The objective of quantifying uncertainty in computational models arises in a wide range of applications, including weather and climate modeling (Schneider, Lan, et al., 2017), fluid dynamics (Duraisamy et al., 2019), and energy systems (Constantinescu et al., 2010). Often, uncertainty associated with predictions from computational models is the result of processes that cannot be resolved, either due to computational complexity limitations (Meneveau & Katz, 2000) or due to uncertainty associated with the process itself (Schneider, Teixeira, et al., 2017). In general circulation models (GCMs), primary uncertainties arise from the representation of subgrid-scale turbulence, convection, and cloud physics, which have a significant impact on the evolution of climate under increasing concentrations of greenhouse gases (Bony & Dufresne, 2005; Brient & Schneider, 2016; Cess et al., 1989; Suzuki et al., 2013; Webb et al., 2013). While clouds are associated with turbulence and convective updrafts with scales of O(10 m), modern climate simulations have a typical horizontal resolution of O(10 km) to O(100 km) (Hoegh-Guldberg et al., 2018). Climate simulations rely on physically motivated parameterizations that model the effects of subgrid-scale processes such as clouds and turbulence on the resolved scales (Hourdin et al., 2013).
Such parameterizations come with parametric and structural uncertainties; quantifying these uncertainties and how they percolate into climate projections remains an outstanding challenge (Schneider, Teixeira, et al., 2017).

Physical parameterizations have historically been individually developed and calibrated using data from isolated experiments (Golaz et al., 2013; Hourdin et al., 2017; Jakob, 2010). They are further adjusted so that global models that incorporate them satisfy selected large-scale observational or physical constraints, such as a closed top-of-atmosphere energy balance or reproduction of the 20th-century global-mean temperature evolution (Hourdin et al., 2017; Mauritsen et al., 2012; Schmidt et al., 2017). Model calibration is usually done manually, focusing on a small subset of model parameters and exploiting only a fraction of the available observational data.

As a step toward automating and augmenting this process, here, we further develop algorithms for model calibration and uncertainty quantification (UQ) that in principle allow models to learn from large datasets and that scale to high-dimensional parameter spaces. In previous work, these algorithms have been demonstrated in simple conceptual models (Cleary et al., 2021) and in a statistically stationary idealized GCM (Dunbar et al., 2021). We take the next step and demonstrate how these algorithms can exploit seasonal variations, which for many climate statistics are large relative to the climate changes expected in the coming decades and contain exploitable information about the response of the climate system to perturbations (Knutti et al., 2006; Schneider et al., 2021).

Numerical weather prediction assimilates observations of atmospheric states that evolve in time to generate initial conditions for forecasts (e.g., Bauer et al., 2015; Kalnay, 2003). For parameter estimation in climate modeling, it is preferable to assimilate time-averaged climate statistics (Schneider, Lan, et al., 2017; Schirber et al., 2013). This focuses the learning problem on quantities of interest in climate predictions (i.e., climate statistics, including higher-order statistics such as precipitation extremes). Additionally, it avoids the need to estimate uncertain atmospheric initial conditions on which trajectories of states depend (Cleary et al., 2021), which reduces the complexity and computational cost of the data assimilation. Calibration and UQ of climate models on the basis of time-averaged statistics smooths the prediction-error based objective function and simplifies the use of data that have different resolution than the climate simulations (Dunbar et al., 2021; Schneider, Lan, et al., 2017). Given our dual desires to avoid having to estimate atmospheric initial conditions, which are forgotten over about 2 weeks (Zhang et al., 2019), and to exploit seasonal variations, it becomes natural to choose averaging timescales between around 30 and 90 days. Such averaging timescales are the focus of this study.

We consider the calibration and UQ of convective parameters in an idealized GCM with seasonally varying insolation (Bordoni & Schneider, 2008; Frierson et al., 2006; O’Gorman & Schneider, 2008). We develop an extension of the calibrate-emulate-sample (CES) Bayesian learning methodology (Cleary et al., 2021) to enable the use of statistics computed from a non-stationary statistical state. The qualitative and quantitative impacts of the time-averaging length of the climate statistics on the parameter calibration and UQ are assessed. GCMs are generally tuned in situations that have low parameter identifiability based on available climate data (Hourdin et al., 2017). We perform numerical experiments with observable climate statistics that are informative about the convective parameters we wish to calibrate, but are not in any simple and direct way related to them. We also perform UQ with less informative climate statistics. Compared to performing UQ with statistics which are highly informative for the convective parameters, UQ using less informative statistics highlights the benefits of incorporating seasonal variations in the climate statistics that are being used for UQ. When performing UQ with statistics which have limited relationships to model parameters, incorporating frequency content provides a mechanism to improve the parameter identifiability.

The remainder of this paper is organized as follows. In Section 2, the Bayesian learning methods for time-dependent problems are introduced. The numerical details of the seasonally forced GCM and UQ experiments are introduced in Section 3. The calibration and UQ results are shown in Section 4, and conclusions are provided in Section 5.

2. Uncertainty Quantification Methods

The goal of this study is to estimate the probability distribution associated with model parameters \( \theta \) that are used by a GCM and about which only imprecise prior information is known. We will consider a Bayesian approach to the estimation of the probability distribution of model parameters \( \theta \), where we seek \( P(\theta | y) \), the conditional probability distribution of parameters \( \theta \) given observed data \( y \). The GCM is a computationally expensive numerical
model that evolves weather states in time. From the GCM, we extract statistical information that is denoted by \( g(\theta) \). Here, \( g(\theta) \) includes the numerical integration of the GCM states and the aggregation of relevant climate statistics in time. Although the observed data \( y \) will, in general, not provide direct information about \( \theta \), we can estimate \( P(\theta|y) \) using Bayesian learning by comparing the GCM outputs \( g(\theta) \) with observed data \( y \).

Standard methods for Bayesian learning include Markov Chain Monte Carlo (MCMC; Brooks et al., 2011), which typically requires \( \mathcal{O}(10^5) \) forward model evaluations of \( g(\theta) \) to sample the posterior distribution (Geyer, 2011). Instead, we perform calibration and UQ using the recently developed CES methodology (Cleary et al., 2021), which consists of three steps: (a) Ensemble Kalman processes (Garbuno-Inigo et al., 2020; Schillings & Stuart, 2017) are used to calibrate parameters and to generate input-output pairs of the mapping \( \theta \rightarrow g(\theta) \); (b) Gaussian process (GP) regression, or other machine learning tools, are used to train an emulator \( g_{\text{GP}}(\theta) \) of the mapping \( \theta \rightarrow g(\theta) \) using the training points generated in the calibration step; and (c) MCMC sampling with the computationally efficient emulator \( g_{\text{GP}}(\theta) \) rather than the expensive forward model \( g(\theta) \) is used to estimate the posterior distribution \( P(\theta|y) \). The CES methodology has previously been used for calibration and UQ of parameters in simple model problems such as Darcy flow and Lorenz systems (Cleary et al., 2021) and for convective parameters in a statistically stationary GCM (Dunbar et al., 2021). More recent methodological developments by Lan et al. (2021) enabled the CES framework to perform simultaneous UQ on \( \mathcal{O}(1000) \) parameters using deep neural network-based emulation and MCMC suited to high-dimensional spaces (Beskos et al., 2008, 2011). A schematic of the CES methodology is shown in Figure 1.

The Bayesian learning methodology used in this study is described in the following sections. In Section 2.1, the inverse problem of estimating \( P(\theta|y) \) in a setting with a periodic cycle is introduced. In Section 2.2, the ensemble Kalman process calibration method is outlined. Section 2.3 introduces the GP emulation in an uncorrelated transformed space, obtained by a singular value decomposition (SVD; principal component analysis) of the noise covariance matrix. Section 2.4 describes how the Bayesian posterior distribution is approximated using the GP emulator.

### 2.1. Seasonal GCM Inverse Problem

With fixed insolation (Frierson et al., 2006; O’Gorman & Schneider, 2008), statistics from the idealized GCM are stationary and ergodic. With the seasonal cycle incorporated, the insolation varies as a function of the ordinal day in the simulation. The resulting GCM states are statistically cyclostationary, with a dependence on the ordinal day. The GCM outputs we use are constructed accounting for the seasonally varying boundary conditions, as the time-averaged GCM statistics

\[
g_T(\theta, \xi, \tau) = \frac{1}{T} \int_{\tilde{t} = \tau}^{\tau+T} H(t' ; \theta, \xi) dt'. \tag{1}
\]
Here, $H(t; \theta, \xi) \in \mathbb{R}^N$, represents the forward operator of the GCM that outputs time-dependent states, depending on time $t$, given parameters $\theta$ and the initial state $\xi$. The number of states being measured is $N_y$. The integration length is specified as $T$. The time-averaged statistics are $G_T(\theta, \xi) \in \mathbb{R}^{N_y}$.

With seasonally varying insolation, the time-averaged statistics in Equation 1 depend on the ordinal day ($t$) in the GCM. With a specified integration length $T$, there are a corresponding set of time-averaged statistics

$$G_T(j, \xi) = \frac{1}{T} \int_{t_{0}+j(T-1)}^{t_{0}+jT} H(t'; \theta, \xi)dt',$$

where, $j = 1$ through $N_j$ is an index representing the time of year in the GCM simulation, starting from $j = 1$ at vernal equinox (time $t_0$). The integration windows are non-overlapping. The length of the year in the GCM is $T_y = 360$ days, and the resulting number of non-overlapping time-averaged statistics are $N_j = T_y/T$. In this framework, the number of batches of statistics $N_b$ is a design parameter set by the integration timescale $T$. For $T = 90$ days, the GCM outputs are aggregated seasonally, with $N_j = 4$. For $T = 360$ days, the statistics are averaged over the full year in the GCM, corresponding to annually averaged climate statistics, and $N_j = 1$.

The GCM outputs are the concatenation of the time-averaged batches over collated ordinal days

$$\mathcal{G}(\theta, \xi) = [G_{T,1}(\theta, \xi_1), G_{T,2}(\theta, \xi_2), \ldots, G_{T,N_j}(\theta, \xi_{N_j})].$$

In this study, we will consider perfect-model numerical experiments, with $y$ being constructed using the same GCM with the parameters set to the target parameters $\theta = \theta'$. The synthetic data are given by

$$y = [G_{T,1}(\theta', \xi_1), G_{T,2}(\theta', \xi_2), \ldots, G_{T,N_j}(\theta', \xi_{N_j})].$$

The sizes of $y$ and $\mathcal{G}$ are $\mathbb{R}^N$, where $N = N_j \cdot N_y$. Note that the total data size $N$ is of length $N_j \cdot N_y$, where, $N_j$ is the number of integration windows per year and $N_y$ is the number of states we are tracking. The resulting relationship between parameters and data is

$$y = \mathcal{G}(\theta, \xi) + \eta,$$

where, $\eta \sim N(0, \Delta)$ is a realization of normal measurement error with zero mean and covariance matrix, $\Delta$. We generate synthetic data and forward model output starting from the same initial conditions $\xi$ in Equation 5. Since, the states of the GCM depend on the ordinal day, the averaging operation in Equation 2 must be consistently aligned in ordinal day between the synthetic data and GCM outputs.

The states of the GCM depend on the boundary conditions, specific to the ordinal day, and the initial conditions $\xi$. The ensemble average of independent realizations of $\mathcal{G}(\theta, \xi_i)$ over different initial conditions $\xi_i$ is

$$\mathcal{G}_{av}(\theta) = \lim_{M \to \infty} \frac{1}{M} \sum_{i=1}^{M} \mathcal{G}(\theta, \xi_i).$$

In the limit of infinite realizations of the climate statistics, the dependence of the ensemble average $\mathcal{G}_{av}(\theta)$ on the initial conditions trends to zero by the central limit theorem. The central limit theorem applies since $\mathcal{G}(\theta, \xi_i)$ contains a full year of GCM output by construction (Equation 3) and is therefore a statistically stationary object with respect to the cycle. But due to computational limitations, only a finite number of ensemble realizations is available in practice. For the synthetic data, we can similarly define $y_{av}$; however, $y_{av}$ is generally not accessible from observations.

Our focus is on the UQ of the model parameters. We aim to avoid the estimation of the atmospheric initial condition. We therefore reformulate the inverse problem such that $\mathcal{G}$ and $y$ can have different initial conditions, and such that we do not require the estimation of the initial conditions. To do so, we assume the forward model output $\mathcal{G}(\theta, \xi)$, is a noisy, finite average approximation of the infinite ensemble average of the climate statistics, that is, $\mathcal{G}(\theta, \xi) = \mathcal{G}_{av}(\theta) + N(0, \Sigma)$, where $\Sigma$ is the internal variability covariance matrix of the GCM (Cleary et al., 2021). The realizations $\mathcal{G}(\theta, \xi)$ and $y$ are both subject to the internal variability of the climate system. In this study, given the definition of $y$ and $\mathcal{G}(\theta, \xi)$, the internal variability is the interannual variability of the GCM. As a result, the relationship between parameters and data is
\[ y = G_m(\theta) + \gamma, \]

where, \( \gamma \sim N(0, \Delta + \Sigma) \). In Equation 7, the initial conditions are removed from the inference problem. The synthetic data are collected into the matrix \( Y \in \mathbb{R}^{N \times N_d} \), where \( N \) is the dimension of the data space and \( N_d \) is the number of data samples, or years in this study. The internal variability covariance matrix \( \Sigma \) is computed from the synthetic data matrix \( Y \) such that \( \Sigma = \text{cov}(Y) \). We assume that \( \Sigma \) does not vary as a function \( \theta \) such that \( \Sigma(\theta^*) \approx \Sigma(\theta) \) and that \( \Sigma \) does not depend on initial conditions \( \xi \). Invoking Gaussian error assumptions based on the central limit theorem, the corresponding negative log-likelihood objective function then is

\[ \Phi(\theta, y) = \frac{1}{2} \| y - G_m(\theta) \|_{\Delta + \Sigma}^2 \]

where, \( \| \cdot \|_k = \| A^{-1/2} \cdot \|_2 \). The likelihood is (Kaipio & Somersalo, 2006)

\[ P(\theta|y) \propto \exp(-\Phi(\theta, y)). \]

In this perfect-model setting, where the synthetic data and forward model are obtained from the same GCM, there is no direct measurement error (\( \eta \)) and no structural model error. To emulate measurement error, we add Gaussian noise to the GCM output, with zero mean and covariance matrix

\[ \Delta = \text{diag}(\delta_k^2), \]

where, \( k \) is an index from 1 through \( N \). We define the noise standard deviation to keep the noise-perturbed data within physical boundaries of the data \( \partial \Omega \) with high probability (e.g., to keep relative humidities between 0 and 1). Here, the noise standard deviation is defined such that

\[ \delta_k = \min \left( C \min \left[ \text{dist}(y_k + 2\sqrt{\Sigma_{kk}}, \partial \Omega_k), \text{dist}(y_k - 2\sqrt{\Sigma_{kk}}, \partial \Omega_k) \right], C_m \cdot y_k \right). \]

The first term takes a fraction \( C \) of the minimum distance from the bounds of the 95% confidence interval around \( y_k \) to the boundary \( \partial \Omega_k \). The second term caps the maximum measurement error standard deviation. We select \( C_m = 0.1 \) to cap the maximum measurement error standard deviation to 10% of the mean data values and \( C = 0.2 \) to control how close the noise-perturbed data can come within physical boundaries \( \partial \Omega \).

### 2.2. Calibrate: Ensemble Kalman Inversion

The first stage of the CES methodology is to calibrate the parameters \( \theta \) of the model based on \( y \). We perform calibration with independent realizations of \( G(\theta, \xi) \), viewed as noisy approximations of \( G_m(\theta) \). Calibration is performed using ensemble Kalman methods, which demonstrate theoretical success, in idealized problems, and empirical success, in complex problems, to optimize parameters under such noise (Duncan et al., 2021).

The utility of the calibration stage is two-fold: (a) optimize parameters to minimize the mismatch between model output and data; and (b) provide good parameter–model output pairs \((\theta, G(\theta, \xi))\) for training an emulator of the parameter-to-data map, with a higher density of training points near the optimal parameters. The ensemble Kalman filter (EnKF) is a Kalman filter implementation in which the covariances are approximated using Monte Carlo sampling (Evensen, 2003). The EnKF has been used widely for derivative-free state estimation in numerical weather prediction (e.g., Houtekamer & Zhang, 2016) and model-based control (e.g., Howland, Ghate, et al., 2020). Ensemble Kalman methods for Bayesian inversion were introduced by Chen and Oliver (2012) and Emerick and Reynolds (2013). These methods provably draw samples from the posterior of linear inverse problems subject to additive Gaussian noise; however, they fail to do so more in more general, nonlinear problems. Recognizing this, the offline ensemble Kalman inversion (EKI, Iglesias et al., 2013) algorithm was introduced for classical, optimization-based inversion. EKI generally drives the ensemble members toward consensus near the optimal solution of the inverse problem (Schillings & Stuart, 2017). The parameter update of ensemble member \( m \) at iteration step \( n \) is

\[ \theta_m^{n+1} = \theta_m^n + C_{m}^{n} (\Sigma + \Delta + C_{m}^{n})^{-1} (y - G(\theta_m^n, \xi_m)), \]

where, \( C_{m}^{n} \) is the ensemble covariance matrix. The update computes the optimal solution of the inverse problem at iteration step \( n \) for training an emulator of \( G_m(\theta) \) to cap the maximum measurement error standard deviation to 10% of the mean data values and

\[ \text{dist}(y_k + 2\sqrt{\Sigma_{kk}}, \partial \Omega_k), \text{dist}(y_k - 2\sqrt{\Sigma_{kk}}, \partial \Omega_k) \]

such that

\[ \delta_k = \min \left( C \min \left[ \text{dist}(y_k + 2\sqrt{\Sigma_{kk}}, \partial \Omega_k), \text{dist}(y_k - 2\sqrt{\Sigma_{kk}}, \partial \Omega_k) \right], C_m \cdot y_k \right). \]

The first term takes a fraction \( C \) of the minimum distance from the bounds of the 95% confidence interval around \( y_k \) to the boundary \( \partial \Omega_k \). The second term caps the maximum measurement error standard deviation. We select \( C_m = 0.1 \) to cap the maximum measurement error standard deviation to 10% of the mean data values and \( C = 0.2 \) to control how close the noise-perturbed data can come within physical boundaries \( \partial \Omega \).
where, $C_{\text{GC}}$ is the empirical cross-covariance between the parameters and the model outputs, and $C_{\text{GG}}$ is the empirical covariance of the model outputs.

EKI is guaranteed to find the optimal parameters in linear problems (Schillings & Stuart, 2017) and has empirical success in nonlinear problems (Dunbar et al., 2021). Several other approaches exist for estimation in nonlinear problems. Li and Reynolds (2009) and Sakov et al. (2012) developed iterative EnKF methods which improve state estimation in strongly nonlinear problems. However, the distribution of the ensemble does not converge to the posterior distribution in the limit of infinite members for nonlinear problems (Annan & Hargreaves, 2007; Garbuno-Inigo et al., 2020; Le Gland et al., 2009; Zhou et al., 2006), necessitating the emulation and sampling described in the following sections. The ensemble Kalman sampler (Garbuno-Inigo et al., 2020) could be used in place of EKI to prevent ensemble collapse. We performed UQ using CES with the ensemble Kalman sampler instead of EKI; the results were not significantly different, but the ensemble Kalman sampler took longer to converge.

The number of ensemble members and EKI iterations are hyperparameters; they are set to standard values of $N_{\text{ens}} = 100$ and $N_i = 5$ in this study (Cleary et al., 2021; Dunbar et al., 2021; Schillings & Stuart, 2017). In CES, the primary purpose of the calibration stage is to provide training samples for the emulator. Increasing $N_{\text{ens}}$, for example, will increasingly collapse the ensemble and limit the utility of the ensembles from the ensuing iterations for emulator training. For our application, $N_{\text{ens}} = 100$ and $N_i = 5$ empirically demonstrate a good balance in hyperparameter tuning experiments.

Each ensemble member is run through the GCM, initialized from the same initial conditions. A spin-up period of one year (360 days) is run before the statistics are computed using Equation 2, to ensure the forward model realizations are subject to differing instantiations of internal variability in the chaotic GCM system. We select a spin-up of one year such that it is sufficiently larger than the Lyapunov timescale of atmospheric motions (about 2 weeks) and so that the statistics are averaged starting from the same ordinal day (see Section 2.1).

### 2.3. Emulate: Gaussian Process Emulation

We emulate the mapping from parameters to model output using a machine learning method that enables the rapid execution of the mapping, compared to the computationally expensive forward model. The calibration stage (Section 2.2) results in $N_i = N_{\text{ens}}$. $N_i$ input-output pairs $(\theta_i, G(\theta, \xi_i))_{i=1}^{N_i}$ of model parameter to model output. Harnessing the input-output pairs as training points, GP regression is used (Rasmussen, 2003) to create an emulator, composed of a mean function and covariance function pair, where $G_{\text{GP}}(\theta) \approx N(\tilde{G}_{\text{GP}}(\theta), \tilde{\Sigma}_{\text{GP}}(\theta))$ and $\Sigma_{\text{GP}} \approx \Sigma$. Since, the input-output pairs are subject to different realizations of the chaotic system, the emulator mean approximates $G_{\text{ens}}(\theta)$ rather than $G(\theta, \xi)$ (Cleary et al., 2021; Dunbar et al., 2021).

The variables of interest in the synthetic data $Y \in \mathbb{R}^{N \times N_d}$ are correlated. The output variables from the GCM forward model $G(\theta, \xi)$ are similarly correlated. The correlation between the GCM statistics results in nonzero off-diagonal covariance matrix elements. In order to maintain a diagonal covariance in the GP emulator $\Sigma_{\text{GP}}$, we transform the GCM statistics into a decorrelated space using principal component analysis on $\Sigma \in \mathbb{R}^{N \times N}$ (Cleary et al., 2021). That is, we decompose the covariance matrix using the SVD

$$\Sigma = V D^2 V^T,$$

with a matrix of principal component vectors (or singular vectors), $V$ and the diagonal matrix, $D$ containing the square roots of the singular values $\sigma_i$. A similar SVD-based variable decorrelation transformation is used in the Ensemble Adjustment Kalman filter (Anderson, 2001). Often, in practical data assimilation problems, $N > N_d$ (Houtekamer & Zhang, 2016). In this case, the covariance matrix is rank deficient, with $\text{rank}(\Sigma) \leq \min(N, N_d)$, with singular values $\sigma_i^2 = 0$ for $i > \text{rank}(\Sigma)$. Methods to decorrelate the data and model output with rank deficient covariance matrices are discussed in Appendix A.

The GP is trained using input-output pairs in the decorrelated space $\{\theta_i, \tilde{G}_{\text{GP}}(\theta, \xi_i)\}_{i=1}^{N_i}$ with the decorrelated space denoted by tildes (\~{}). The resulting input-output mapping is approximated as

$$\tilde{G}_{\text{ens}}(\theta) \approx N(\tilde{G}_{\text{GP}}(\theta), \tilde{\Sigma}_{\text{GP}}(\theta)).$$
The GP kernels used are a white-noise kernel added to an Automatic Relevance Determination radial basis function kernel; further details provided in Dunbar et al. (2021); Cleary et al. (2021). The GP hyperparameters are trained using the input-output pairs. The training process results in a GP regression function \( \hat{G}_{GP}(\theta) \) which takes \( \theta \) as input and emulates \( G_e(\theta) \) in a computationally efficient fashion.

### 2.4. Sample: Posterior Sampling Using MCMC

The Bayesian posterior distribution is approximated through MCMC sampling with the trained GP emulator \( \hat{G}_{GP}(\theta) \). The data \( y \) is normalized and transformed into the decorrelated space \( \hat{y} \), as in Section 2.3. The Bayesian posterior distribution is approximated as (Stuart, 2010).

\[
P(\theta|\hat{y}) \propto P(\hat{y}|\theta)P(\theta)
\]

\[
P(\theta|\hat{y}) \propto \exp\left( -\frac{1}{2}||\hat{y} - \hat{G}_{GP}(\theta)||_2^2 - \frac{1}{2} \log\det (\Sigma_{GP}(\theta) + \Delta) \right) P(\theta).
\]

With a normal prior distribution governed by mean \( \bar{\theta} \) and covariance \( \Gamma_y \), the resulting MCMC objective function is

\[
\Phi_{MCMC}(\theta, \hat{y}) = \exp\left( -\frac{1}{2}||\hat{y} - \hat{G}_{GP}(\theta)||_2^2 \right.
\]

\[
- \frac{1}{2} \log\det (\Sigma_{GP}(\theta) + \Delta) - \frac{1}{2}||\theta - \bar{\theta}||_2^2.
\]

We use a random walk Metropolis MCMC algorithm. The number of MCMC samples is set to \( N_{MCMC} = 200,000 \) with a burn-in of 10,000.

### 3. Seasonal GCM Uncertainty Quantification

We perform numerical experiments with an idealized GCM with a seasonal cycle. The GCM simulation setup is presented in Section 3.1. Various climate statistics that we extract from the GCM and use in the numerical experiments are discussed in Section 3.2 and shown in Section 3.3.

#### 3.1. Seasonal Simulation Setup

The GCM used in this study is based on the Geophysical Fluid Dynamics Laboratory’s Flexible Modeling System (Frierson et al., 2006). The GCM simulates an idealized aquaplanet with a homogeneous mixed-layer slab ocean bottom boundary condition with a depth of 1 m. The GCM has been used previously for simulations of the hydrological cycle over a range of climates (O’Gorman & Schneider, 2008) and to characterize seasonal variability in the tropics (Bischoff & Schneider, 2014, 2016; Bordoni & Schneider, 2008; Kaspi & Schneider, 2011; Merlis et al., 2013a, 2013b; Wei & Bordoni, 2018). The GCM is axisymmetric and statistically cyclostationary. The spectral transform method is used in the horizontal directions, and finite differencing in sigma coordinates is used in the vertical direction. The horizontal resolution used is T21 with \( N_{\phi} = 32 \) discrete latitude points on the transform grid, and 20 sigma levels (\( \sigma_p = p/p_s \) where \( p \) is the pressure and \( p_s \) is the local surface pressure).

A two-stream gray radiation scheme is used. The top-of-atmosphere insolation is prescribed and varies according to a seasonal cycle (Bordoni & Schneider, 2008; Wei & Bordoni, 2018). The diurnal cycle insolation variations are neglected, with a daily average insolation applied. The longwave and shortwave optical thicknesses depend on the latitude and pressure. The radiative effects of variations of atmospheric water vapor or clouds are neglected, and therefore, water vapor and cloud feedbacks are not included in the GCM.

The convection is parameterized using a simplified quasi-equilibrium Betts-Miller scheme (Betts, 1986; Betts & Miller, 1986; Frierson, 2007). Vertical profiles of temperature and humidity are used to calculate precipitation and associated temperature and humidity changes through a relaxation to moist adiabatic reference profiles (Frierson, 2007). The relaxation is included as a forcing to the temperature and humidity balances.

\[
\frac{dT}{dt} + \cdots = S_T - \frac{T - T_{ref}}{\tau}
\]

\[ 10.1029/2021MS002735 \]
\[
\frac{\partial q}{\partial t} + \cdots = S_q - f_q f_T \frac{q - q_{ref}}{\tau},
\]

where, the dots represent the dynamical terms in the equations, and \(S_q\) and \(S_T\) represent the temperature and specific humidity forcings aside from the convection scheme. The term \(f_q\) governs the spatiotemporal activation of the convection scheme, depends on the thermodynamic state, and is a function of \(z\). The term \(f_T\) modifies the specific humidity relaxation (O’Gorman & Schneider, 2008). The reference temperature \(T_{ref}\) is a moist adiabat, chosen so that the convection scheme conserves enthalpy integrated over vertical columns (Frierson, 2007; O’Gorman & Schneider, 2008). The reference specific humidity \(q_{ref}\) is that which corresponds to a prescribed relative humidity with respect to the moist adiabat \(T_{ref}\). For the UQ experiments, our foci are two parameters: the prescribed reference relative humidity parameter \(\theta_{RH}^{ref}\) and the relaxation timescale parameter \(\theta_f\).

The GCM simulation starts from vernal equinox, and the year length is \(T_c = 360\) days. All climate statistics used for UQ are zonally averaged. We use the Betts-Miller convection scheme with standard reference values of the parameters of \(\theta = (\theta_{RH}, \theta_f) = (0.7, 7200\) s) (Frierson, 2007; O’Gorman & Schneider, 2008). UQ of \(\theta\) relies on a prior knowledge about the convective parameters. We use wide prior distributions of the parameters, which enforce physical constraints, such as \(0 < \theta_{RH} \leq 1\) and \(\theta_f > 0\), but are otherwise uninformative (Dunbar et al., 2021). The selected priors are \(\theta_{RH} \sim \text{Logit}(N[0, 1])\) and \(\theta_f \sim \text{Log}(N[12\) hr, \((12\) hr\(^2\))], that is, normal distributions of logit- and log-transformed parameters. Given the true parameter value of \(\theta_f^* = 2\) h, the prior is wide and relatively uninformative. The parameter priors are independent, although joint prior distributions could be used in future work.

### 3.2. Climate Statistics Used for UQ

We calibrate and perform UQ on the parameters of the convection scheme using more and less informative time-averaged observations from GCM simulations. The informative statistics, such as the mid-tropospheric relative humidity, are chosen because they are strongly affected by the choice of the convective parameters (e.g., O’Gorman & Schneider, 2008; O’Gorman et al., 2011; Schneider & O’Gorman, 2008). For comparison, we also choose less informative statistics that are affected less by the choice of convective parameters, such as surface wind speeds. Numerical UQ experiments using the differing degrees of information in the climate statistics will illustrate the impact of incorporating higher frequencies in the climate statistics used for parameter estimation.

For the informative climate statistics, we choose three variables: the mid-tropospheric \((\sigma_p = 0.5)\) relative humidity, the total precipitation rate, and, as a measure of precipitation intensity, the probability that a daily precipitation amount exceeds the latitude-dependent ninetieth precipitation percentile threshold from a long control simulation with the true parameters (Dunbar et al., 2021). Since intense precipitation events are influenced by the convection scheme (O’Gorman & Schneider, 2009a, 2009b), exceedances of precipitation over a high threshold are anticipated to be informative about the convection scheme parameters. The three statistics, are evaluated at each of the \(N_p = 32\) discrete latitudes, giving 96 total quantities of interest. For less informative statistics, we use the zonally averaged precipitation rate and surface wind speed. As with the informative statistics, \(N_p = 32\) discrete latitudes are considered, giving 64 total quantities of interest (see Table 1 for a summary).

To ensure the different GCM statistics are weighted equally in the UQ, we nondimensionalize the GCM statistics (Equation A1) with the median \(y^*_p\), taken over latitude and time, of each specified data type. There is one characteristic value for each data type, for example, one characteristic value \(y^*_p\) for precipitation and a separate characteristic value for relative humidity. The nondimensionalization is described in more detail in Appendix A.

### 3.3. Synthetic Data Used for UQ

As synthetic data, we used GCM output with different initial conditions, and perturbed with measurement error with covariance \(\Delta\). The synthetic states are constructed by running the seasonal GCM for 150 years at the true parameters \(\theta^*\). For each case of \(T = 90\) days and \(T = 360\) days, the GCM states are aggregated according to the method introduced in Section 2.1. The informative synthetic statistics for \(T = 360\) days are shown in Figures 2a–2c, which compares the \(N_p = 150\) years ensemble average of the statistics with a 1 year sample subject to internal variability and perturbed with measurement noise \(N(0, \Delta)\). The synthetic data for \(T = 90\) days are shown in Figures 2d–2f. For the data averaged over 90 days, a distinct seasonal cycle emerges with, for example, relatively
low precipitation and relative humidity in the northern hemisphere winter (days 271–360). The less informative statistics of precipitation rate and surface wind speed are shown for $T = 360$ days and $T = 90$ days in Figure 3.

4. GCM Calibration and UQ Results

4.1. Informative Statistics

The convective parameters $\theta_{RH}$ and $\theta_{\tau}$ are calibrated using EKI with the informative statistics with $T = 90$ days and $T = 360$ days (see Table 1). The EKI calibration is performed with a synthetic data sample $y$ (Equation 7). Since the synthetic data is subject to internal variability, the corresponding calibration is influenced by the synthetic sample values. The synthetic data sample is randomly selected from the 150 years of historical observed data constructed with the true convective parameters $\theta^t$ (see Section 3.2). Calibration is performed with 10

| Table 1 | Summary of the Climate Observed Data Used |
|---------|------------------------------------------|
| **Variables** | Informative | Less informative |
| 1. Relative humidity ($\sigma_p = 0.5$) | | |
| 2. Precipitation rate (mm/day) | | |
| 3. Probability of 90th percentile precipitation | | |
| Latitudes, $N_\phi$ | 32 | 32 |
| $N_\psi$ | 96 | 64 |
| $N$, $T = 360$ days (annual) | 96 | 64 |
| $N$, $T = 90$ days (seasonal) | 384 | 256 |

*Note. Calibration and uncertainty quantification (UQ) is performed using observed data that are more or less informative about the convective scheme. The total size of the observed data used for UQ is $N$."

Figure 2. Informative synthetic data with (a)–(c) $T = 360$ days and (d)–(f) $T = 90$ days. The light solid lines correspond to $y^\infty$. The shaded regions correspond to a 95% confidence interval around $y^\infty$ with covariance $\Sigma + \Delta$. The dark solid lines with circle markers correspond to a randomly selected sample of the synthetic data $y$ that has been subjected to measurement noise $N(0, \Delta)$. Day 0 corresponds to vernal equinox, and the northern hemisphere winter and fall seasons are provided. Northern spring and summer are not shown since the general circulation model is symmetric about the equator.
independent synthetic samples, and the results in this section are presented as means and standard deviations over the 10 independent realizations. The cases with \( T = 90 \) days and \( T = 360 \) days are run with the same 10 synthetic samples.

The modes of the EKI ensemble distributions are shown in Figure 4. The modes demonstrate a reduction in estimation error of the true convective parameters when the statistics are aggregated seasonally with \( T = 90 \) days. The calibrated convective parameters using annually averaged GCM statistics have higher mean bias and variability between the instantiations with different synthetic data samples. The percent error of the average mode at the last EKI iteration for \( \theta_{RH} \) is 0.3% and 0.8% for \( T = 90 \) days and \( T = 360 \) days, respectively. For \( \theta_{\tau} \), the percent errors are 1.2% and 11.5% for \( T = 90 \) days and \( T = 360 \) days, respectively. For both integration timescales, the percent errors are 1.2% and 11.5%

![Figure 3](image)

![Figure 4](image)
error associated with the \( \theta_r \) estimation was larger than the estimation error for \( \theta_{RH} \). The percent error metric accounts for differences in the order of magnitude of the different parameters of interest. This indicates that the integration of the GCM states over an annual cycle filters information from the resulting GCM statistics that is informative, especially about the relaxation timescale.

The normalized mean square error (MSE) for \( \theta_{RH} \) at EKI iteration \((n)\) is

\[
\epsilon_{RH}^{(n)} = \frac{1}{\theta_{RH}} \sqrt{\frac{1}{N_{ens}} \sum_{i=1}^{N_{ens}} (\theta_{RH,i} - \theta_{RH}^{\dagger})^2},
\]

(20)

The error is computed similarly for \( \theta_r \). For both the relative humidity and relaxation timescale parameters, \( T = 90 \) days reduces the parameter estimation MSE relative to \( T = 360 \) days by about a factor 2–3: The MSE for \( \theta_{RH} \) is 0.006 and 0.015 for \( T = 90 \) days and \( T = 360 \) days, respectively; for \( \theta_r \), the MSE is 0.06 and 0.23 for \( T = 90 \) days and \( T = 360 \) days, respectively. The seasonally aggregated data also have a smaller spread among the realizations, as visualized by the uncertainty bands around the mean values. The reductions in the MSE standard deviations were 43% and 66% for \( \theta_{RH} \) and \( \theta_r \), respectively.

The input-output pairs generated during EKI parameter calibration are used for training the GP emulator. These pairs of parameters and model evaluations are taken from four 100-member EKI iterations as well the initial 100-member ensemble drawn from the prior, giving a total of 500 input-output pairs for GP training. The results of this study were not materially different with more training points. The GCM outputs are mapped into a normalized and decorrelated space according to the SVD of the internal variability covariance matrix \( \Sigma \), truncated to contain 95% of the energy in the singular values. For \( T = 90 \) days, the truncation contains \( k = 47 \) singular values, while \( T = 360 \) days retains \( k = 23 \) (see Appendix B). A scalar-valued GP is trained for each of the \( k \) outputs. For comparison, we also ran the emulation and sampling steps for \( T = 360 \) days without SVD truncation.

The posterior distributions are approximated using MCMC sampling with the trained GP emulators. The posterior distribution for the two convective parameters for a randomly selected synthetic data sample are shown in Figure 5. In Figure 5a, the posterior distribution was computed with normalized input data, aggregated annually with \( T = 360 \) days, but the SVD was not truncated before emulation and sampling. The posterior resulting from the same initial ensemble and synthetic data sample but with SVD truncation is shown in Figure 5b. The truncation of the SVD smooths and inflates the posterior distribution. While the posteriors are qualitatively similar, they differ quantitatively. The true parameters are outside the region containing 99% of the posterior mass for the untruncated case but are within the region containing 75% of the posterior mass for the truncated case. Further investigation of the influence of SVD regularization on the posteriors is in Appendix B. Hereafter, all posterior results will focus on the posterior distributions estimated using emulation and sampling in the truncated SVD space (see Appendix A). The convective parameter posterior distribution using seasonally aggregated data with \( T = 90 \) days with the SVD truncated at 95% of the total energy is shown in Figure 5c. As in the calibration stage, the posterior mode for \( T = 90 \) days is closer to the true parameters than for \( T = 360 \) days. Compared to the posterior with \( T = 360 \) days, the \( T = 90 \) days posterior is also more compact.

We quantify the size of the posterior distributions for the two integration timescales by computing the two-dimensional area of the convex hull containing 75% of the posterior mass (middle contour level in Figure 5). The area containing 75% of the posterior mass for each value of \( T \) is normalized by the area containing 75% of the prior mass for reference. The percentages of the prior area taken up by the posteriors for \( T = 90 \) days and \( T = 360 \) days, averaged over the 10 CES instantiations, are shown in Table 2. The areas are computed in the \( \logit(\theta_{RH}) - \logit(\theta / 1 \text{s}) \) transformed space, in which the sampling is performed. The area of the posterior distribution resulting from the seasonal integration is approximately a factor of two smaller than the posterior from the annually integrated statistics. However, compared to the wide prior, both posteriors are tight, with the posterior areas occupying around 1% of the prior area.

### 4.2. Less Informative Statistics

In many UQ applications, intuition about which quantities of interest will lead to improved parameter estimation is not available before experimentation. While the idealized GCM and convective parameterization used in this study have a rich set of investigations to provide prior knowledge about the relationship between \( \theta \) and \( G(\theta) \), we...
also consider UQ of the parameters using less informative GCM statistics. To do so, we use the precipitation rate and surface wind speed, zonally averaged and integrated with $T = 360$ days and $T = 90$ days (Figure 3). Four instantiations of CES with differing synthetic data samples are run for each integration length of $T = 90$ days and $T = 360$ days.

The MSE of the EKI parameter calibration with the less informative GCM statistics are shown in Figure 6 for $T = 90$ days and $T = 360$ days. With precipitation rate and surface wind speed as the basis for UQ, the annually averaged data result in significantly higher parameter estimation error than seasonally aggregated data. The $T = 360$ days MSEs are 0.26 and 1.56 for $\theta_{RH}$ and $\theta_{\tau}$, respectively. The $T = 90$ days MSEs are 0.02 and 0.28 for $\theta_{RH}$ and $\theta_{\tau}$, respectively. The ensemble means and modes of the two parameters are shown in Figure 7. In Appendix C, the calibration is performed by considering the seasonal data sequentially, rather than collectively, to restrict the data size for both $T = 90$ days and $T = 360$ days to $N = 64$. The results demonstrate that the reductions in MSE are the result of the incorporation of seasonal information in the data, rather than the dimensionality of the data space.

The posterior distributions computed using Markov Chain Monte Carlo using a Gaussian process emulator that was trained using observed data and general circulation model outputs aggregated with an integration timescale of $T = 360$ days and $T = 90$ days. The contours correspond to 50%, 75%, and 99% of the posterior distribution, the star is the truth, and the circle is the average of the ensemble members after the last ensemble Kalman inversion iteration. Posterior distributions are shown for (a) $T = 360$ days, decorrelated with full singular value decomposition (SVD), and (b) a 95%-energy truncation of the SVD. (c) Posterior for $T = 90$ days, decorrelated with 95%-energy truncated SVD. Panels (d)–(f) same as (a)–(c) with the posterior distributions shown in the physical parameter space.

![Figure 5](image-url)

**Figure 5.** Convective parameter posterior distributions computed using Markov Chain Monte Carlo using a Gaussian process emulator that was trained using observed data and general circulation model outputs aggregated with an integration timescale of $T = 360$ days and $T = 90$ days. The contours correspond to 50%, 75%, and 99% of the posterior distribution, the star is the truth, and the circle is the average of the ensemble members after the last ensemble Kalman inversion iteration. Posterior distributions are shown for (a) $T = 360$ days, decorrelated with full singular value decomposition (SVD), and (b) a 95%-energy truncation of the SVD. (c) Posterior for $T = 90$ days, decorrelated with 95%-energy truncated SVD. Panels (d)–(f) same as (a)–(c) with the posterior distributions shown in the physical parameter space.

### Table 2

| Statistics for UQ | $T = 90$ days | $T = 360$ days |
|-------------------|---------------|---------------|
| Informative statistics | $0.5\% \pm 0.1\%$ | $1.3\% \pm 0.3\%$ |
| Less informative statistics | $5.5\% \pm 0.9\%$ | $26.0\% \pm 7.0\%$ |

Note. The area of the convex hull containing 75% of the posterior mass for each value of $T$ and data type is normalized by the area containing 75% of the prior mass. The means and standard deviations over independent synthetic data realizations are provided. Ten and four realizations are used for UQ with the informative and less informative statistics, respectively.
compared to the true parameters. For $T = 90$ days, the uncertainty for $\theta_{RH}$ estimated with the less informative statistics has collapsed to a similar order as the uncertainty estimated with informative statistics. While the uncertainty is larger for $\theta_\tau$ when estimated using the less informative statistics, the overall posterior area is reasonably collapsed, as it is of similar order to the $T = 360$ days posterior with informative statistics. These results suggest that seasonally averaged precipitation rate and surface wind speed are sufficient statistics to estimate convective parameters while annually averaged precipitation and wind speed introduce higher estimation error and parameter uncertainty.

As with the informative GCM statistics, we compare the sizes of the posterior distributions by computing the two-dimensional area of the convex hull containing 75% of the posterior mass. The posterior areas, normalized by the prior area, averaged over the four CES instantiations, are shown in Table 2. The posterior area for $T = 360$ days is approximately 5 times larger than the area produced with $T = 90$ days. The implications of the reduction in the size of the posterior distribution on the parametric uncertainty in the GCM are tested in Section 4.2.1.

Figure 6. Parameter calibration performed with ensemble Kalman inversion (EKI) using less informative statistics. Mean square error of (a) the estimated relative humidity parameter $\theta_{RH}$ compared to truth value $\theta_{RH}^{\dagger}$ and of (b) the estimated relaxation time scale $\theta_\tau$ compared to truth value $\theta_\tau^{\dagger}$. Solid lines are $T = 360$ days and dashed lines are $T = 90$ days.

Figure 7. Parameter calibration performed with ensemble Kalman inversion (EKI) using less informative statistics. Mean of (a) the estimated relative humidity parameter $\theta_{RH}$ compared to truth value $\theta_{RH}^{\dagger}$ and of (b) the estimated relaxation time scale parameter $\theta_\tau$ compared to truth value $\theta_\tau^{\dagger}$. (c), (d) Same as (a), (b) for the modes of the ensemble members.
4.2.1. Prediction Experiments

To demonstrate the effect of parametric uncertainty in the GCM on climate predictions, we draw samples of 100 parameter pairs from the posterior distribution resulting from the less informative statistics of precipitation rate and surface wind speed (Figure 8). The samples of parameter pairs using seasonal integration ($T = 90$ days) and annual integration ($T = 360$ days) in the UQ are shown in Figure 9. With each sample of parameters, we produce climate quantities of interest, averaged over a 20 year period. We compute ensemble statistics over the outputs from the samples (prediction uncertainty) and compare these with a 20 year simulation with the true parameters $\theta^\dagger$.

The relative humidity at $\sigma_p = 0.5$, the precipitation rate, and the intense precipitation probability (the probability of exceeding a 90th percentile threshold averaged over the 20 year simulation at $\theta^\dagger$) are shown in Figure 10. The posterior distribution was estimated using (seasonally averaged) precipitation rate data directly, but neither relative humidity nor intense precipitation probability were used. The mean absolute percent errors for the mean prediction values, averaged across latitudes, for the $T = 90$ days posterior are 1.2%, 0.6%, and 1.4%, for the relative humidity, precipitation rate, and intense precipitation probability, respectively. For the predictions with the $T = 360$ days posterior, the mean absolute percent errors are 4.7%, 1.8%, and 21.5%.

The prediction uncertainty in each climate quantity of interest is indicated by a 95% confidence interval, estimated as the values of the 2.5th and 97.5th percentiles of the variables of interest at each latitude over the convective parameter pairs (Figure 9). For both $T = 90$ days and $T = 360$ days, there is limited prediction uncertainty in the mean precipitation rate. The widths of the confidence intervals are 6% and 4% of the mean precipitation rate, averaged across latitudes, for $T = 360$ days and $T = 90$ days, respectively. The prediction uncertainties in relative humidity and intense precipitation probability are significantly higher for UQ performed with annually averaged data ($T = 360$ days) compared to UQ with seasonally averaged data ($T = 90$ days). The $T = 90$ days posterior reduces the size of the 95% confidence interval for the relative humidity by 70%, averaged across latitudes, compared to $T = 360$ days. The reduction in the size of the 95% confidence interval for the intense precipitation probability is 9%.

We also performed idealized global-warming prediction experiments. To represent global warming in the idealized GCM, the longwave opacity in the atmosphere is increased by 50%, as in O’Gorman and Schneider (2008) which results in a global-mean surface air temperature increase from 284 K in the control climate to 292 K in the warm climate. We accumulated GCM statistics over a 20-year window after a spinup of one year, running simulations with the true GCM parameters $\theta^\dagger$ with increased longwave opacity for comparison. We use the same 100 parameter pairs drawn from the posterior distributions estimated from the less informative climate statistics.
The idealized global warming results are shown in Figure 11. As expected, the warming climate predictions with the $T = 90$ days posterior distribution are more accurate than with the $T = 360$ days posterior, especially for the relative humidity and intense precipitation probability. The mean absolute percent error for the mean prediction values, averaged across latitudes, for the $T = 90$ days posterior are 1.7%, 1.1%, and 2.2%, for the relative humidity, precipitation rate, and intense precipitation probability, respectively. For the predictions with the $T = 360$ days posterior, the mean absolute percent errors are 4.7%, 2.2%, and 10.4%. The $T = 90$ days posterior reduces the size of the 95% confidence interval for the relative humidity in a warmed climate by 75% compared to $T = 360$ days, averaged across latitudes. The reduction in the size of the 95% confidence interval for the intense precipitation probability in a warmed climate is 35%. These results demonstrate that the incorporation of the seasonal cycle in the UQ reduces the parametric uncertainty in climate predictions in this model.

5. Conclusions

We performed calibration and UQ of the parameters of a convective parameterization in a seasonally forced idealized GCM. While GCMs are typically tuned using annually and globally averaged climate statistics (Hourdin et al., 2017), our results demonstrate, in an idealized setting, the qualitative, quantitative and systematic refinement of the parameter distribution through the incorporation of seasonal information. Performing parameter calibration with seasonally averaged data significantly reduced the error associated with the estimated parameters compared to calibration using annually averaged data. The posterior distribution resulting from the Bayesian UQ with seasonally averaged data were reduced in size. One measure of success for Bayesian UQ, is to capture the true parameters within a high mass region of the posterior distribution (demonstrated in our results). We are also interested in the shape and size of the posterior distribution, as this provides valuable information about

Figure 10. Prediction experiments for the case with no imposed warming for (a) relative humidity at $\sigma_p = 0.5$, (b) precipitation rate, and (c) intense precipitation probability. The climate statistics are averaged over 20 years. The posteriors from the less informative statistics (surface wind speed and precipitation rate) are used. The lines correspond to the mean of the predictions and the shaded regions, correspond to 95% confidence intervals. The climate statistics at the true parameters $\theta^*$ are shown with a dashed line.

Figure 11. Prediction experiments for the case with imposed warming (50% increase in longwave opacity) for (a) relative humidity at $\sigma_p = 0.5$, (b) precipitation rate, and (c) intense precipitation probability. The climate statistics are averaged over 20 years. The posteriors from the less informative statistics (surface wind speed and precipitation rate) are used. The lines correspond to the mean of the predictions, and the shaded regions correspond to 95% confidence intervals. The climate statistics at the true parameters $\theta^*$ are shown with a dashed line.
parameter correlations and uncertainty, with respect to observed data. We demonstrate that choosing suitable data for UQ which reduce the size of the posterior distribution can lead to significant reductions of climate prediction uncertainty.

The impact of incorporating additional frequency content is pronounced when the climate statistics used for UQ are less informative about the parameterizations. Such situations often occur in the climate modeling setting where a number of parameterizations may simultaneously and nonlinearly influence a quantity of interest and where it is not always clear which climate statistics should be used to calibrate an unknown GCM parameter.

To enable UQ in the climate model setting, we used the CES methodology (Cleary et al., 2021), which enables the efficient and accurate estimation of Bayesian posterior distributions of parameters from noisy climate data. CES uses gradient-free optimization to calibrate parameters and generate parameter-data pairs, GP regression to emulate the parameter-to-data mapping, and MCMC to sample from the posterior distribution. The emulation and sampling is performed in a decorrelated dataspace through a transformation based on the principal component analysis of the noise covariance matrix. In this study, we modified the original CES methodology by first normalizing the data to ensure all statistics are weighted equally in the UQ and second by regularizing the covariance matrix to enable the UQ of applications with ill-conditioned or rank deficient covariance matrices. We quantified the impact of the normalization and regularization, with the regularization smoothing the posterior and slightly increasing its size.

Beyond convection, Earth system models rely on a number of parameterizations of cyclostationary processes, including models of carbon accumulation and storage (Bloom et al., 2016) and of atmospheric boundary layer (ABL) turbulence which parameterize unclosed subgrid processes (Stull, 2012). Carbon storage is inherently seasonal (Rowland et al., 2014). The ABL has distinct seasonal and diurnal variations (Wyngaard, 2010; Howland, González, et al., 2020). We expect our results to be relevant to calibration and UQ for statistical variations on different timescales, beyond the seasonal cycle. Higher frequency content can be incorporated in the framework proposed in this paper by including the amplitude and phase of the Fourier components associated with the cycle of interest. For example, based on our findings presented here, we anticipate that incorporating the diurnal variation of ABL turbulence statistics (Howland et al., 2021) may improve the UQ of subgrid scale turbulence models in GCMs.

The selection of the optimal aggregation timescale for climate statistics and choice of objective function remain open questions. In the context of parameter estimation, the integration timescale used to generate time-averaged statistics is a hyperparameter that dictates the trade-off between the frequency content and signal-to-noise ratio of the climate statistics. In the limit case of \( T \to \tau_y \), where \( \tau_y \) is the Lyapunov timescale, higher frequencies are incorporated in the statistics, but the signal-to-noise ratio is reduced. Conversely, as \( T \to \infty \), all frequency content is removed from the statistics, which, as this study demonstrates, can adversely impact parameter estimation. We anticipate that the selection of the filter timescale \( T \) may be problem and parameterization specific, as it relates to the question of parameter identifiability. In this study, we selected seasonal averages because of the large amplitude of the seasonal cycle in many climate statistics and indications that seasonal variations are informative about the climate change response in many climate variables (Schneider et al., 2021). Future work should develop a more generalized approach for the selection of \( T \) based on the frequency content of the quantities of interest.

**Appendix A: Data and Model Output Decorrelation for Rank Deficient Problems**

To facilitate the transformation into the uncorrelated space with rank deficient covariance matrices, the singular value decomposition/principal component analysis is truncated. Since the statistical quantities of interest have a range of magnitudes, the data are first normalized, and then the singular value decomposition is truncated as a form of regularization (Hansen, 1987). A detailed investigation of the influence of the normalization and truncation on the posteriors resulting from calibrate-emulate-sample (CES) is performed in Appendix B.

The data used for uncertainty quantification are provided concurrently to the CES pipeline in the concatenated vector \( y \). In this framework, the statistical quantities in the data \( y \) may have a range of magnitudes, and normalization is required before regularization (e.g., Tibshirani. 1996). The data are normalized by a characteristic value associated with each individual data type \( y_i \in \mathbb{R}^N \). The normalized data are

\[
y'_i = y_i \cdot y^{-1}_c, \quad (A1)
\]
where, (*) denotes normalized data. The characteristic values in \( y \) are described in the application of CES to the seasonally forced general circulation model in Section 3. Each column of \( Y \) is normalized using Equation A1 to yield the normalized data matrix \( Y^* \). The normalized covariance is \( \Sigma^* = \text{cov}(Y^*) \). The SVD transformation is performed using covariance matrix \( \Sigma^* \) and data \( y^* \).

The singular value decomposition (SVD) is truncated in order to account for rank deficient or ill-conditioned covariance matrices (Hansen, 1987). The truncated SVD is defined for covariance \( \Sigma^* \) as

\[
\Sigma^*_k \approx \Sigma^*_k = V_k D_k^2 V_k^\top \quad D_k^2 = \text{diag}(\sigma_1^*, \ldots, \sigma_k^*),
\]

truncated at the \( k \leq \text{rank}(\Sigma^*) \) singular value. The truncated singular vector matrix is

\[
A_k = [v_1, \ldots, v_k]
\]

where, \( v_i \) is the singular vector corresponding to the singular value \( \sigma_i^2 \). The truncated SVD space is given by

\[
\tilde{\Sigma}^*_k = D_k^{-1} V_k^\top \Sigma^*_k V_k D_k^{-1}
\]

\[
\tilde{y}_k = D_k^{-1} V_k^\top y^*.
\]

In this study, the truncation location is selected as the lowest value of \( k \) such that

\[
\sum_{i=1}^k \sigma_i^2 \geq 0.95 \sum_{i=1}^N \sigma_i^2,
\]

so that 95% of the variance is retained. The Gaussian process (GP) output can be mapped to the original normalized space using

\[
G^*_{\text{GP}} = V_k D_k \tilde{G}_{\text{GP}}
\]

\[
\Sigma^*_{\text{GP}} = V_k D_k \Sigma_{\text{GP}} D_k V_k^\top.
\]

Finally, the GP output can be transformed into the dimensional space through

\[
G_{\text{GP}} = G^*_{\text{GP}} \otimes y_c \quad \text{and} \quad \Sigma_{\text{GP}} = \Sigma^*_{\text{GP}} \otimes y_c y_c^\top,
\]

where \( \otimes \) denotes pointwise multiplication.

### Appendix B: Sensitivity of the Posterior to Covariance Regularization

In Section 2.3, the singular value decomposition (SVD) is truncated as a form of regularization. Here, we detail the sensitivity of the calibrate-emulate-sample posterior distributions to the SVD regularization using informative general circulation model statistics (see Table 1) and annually averaged data with \( T = 360 \) days, for which the data covariance matrix is full rank. The data are normalized, as discussed in Appendix A. The cumulative sum of the singular values is shown in Figure B1. Truncation at 95% of the singular value energy corresponds to \( k = 23 \). Regularization can also be performed using Tikhonov regularization, which inflates the diagonal elements (Hansen, 1987; Schneider, 2001). Following Hansen (1987), regularization in the form of diagonal inflation \( \Lambda = \text{diag}(\lambda^2) \) is added to \( \Sigma \) where, \( \lambda = (\sigma_1^2 / \sigma_{k+1}^2)^{1/4} \). The SVD is performed such that

\[
\Sigma + \Lambda = V D^2 V^\top,
\]

where, \( D^2 \) is a diagonal matrix of the singular values \( \sigma_1^2 + \lambda^2 \). The orthonormal eigenvectors are in \( V \). The influence of the truncation and Tikhonov regularizations on the posterior areas for \( T = 360 \) days are provided in Table B1. The \( T = 360 \) days posterior distributions are shown in Figure B2. The regularization slightly increases the posterior area and smooths the posterior roughness.

**Figure B1.** Cumulative sum of the singular values from the normalized \( T = 360 \) days synthetic data.
Parameter space
Informative statistics

Note. The area of the convex hull containing 75% of the posterior mass for each regularization method is normalized by the area containing 75% of the prior mass.

Table B1
The Ratio (%) of the Area Occupied by the Posterior to the Area Occupied by the Prior

| Parameter space    | Original | Normalized | Truncated | Tikhonov |
|--------------------|----------|------------|-----------|----------|
| Informative statistics | 0.5%     | 0.5%       | 0.8%      | 0.9%     |

Figure B2. Posterior distributions using the informative statistics with $T = 360$ days. Comparison of the posteriors estimated with (a) original, (b) normalized, (c) normalized and regularized (truncation), and (d) normalized and regularized (Tikhonov diagonal inflation) covariance matrix $\Sigma$.

Appendix C: Parameter Estimation With Statistics Batching (Filtering and Smoothing)

It is often of practical interest to estimate parameters in problems that are not statistically stationary, such as the seasonally forced general circulation model (GCM) with time-dependent boundary conditions. In the uncertainty quantification (UQ) methodology proposed in Section 2, the statistics for all seasons must be available to perform UQ. We refer to this approach as smoothing. In problems which are not statistically stationary (e.g., with time-evolving boundary conditions), all requisite data for parameter UQ may not be available at the initialization of the estimation. This arises in the ensemble Kalman filter setting for state estimation (e.g., Houtekamer & Zhang, 2016), and we refer to this as filtering.

We can instead pose ensemble Kalman filter (EKI) as a filtering approach for parameter estimation where the synthetic data are collected sequentially in a time-dependent problem. This differs from the definition of the synthetic data in Section 2.1, where the data from each season are collected together into $y$ (termed smoothing). For the filtering approach, the data is observed and averaged over time $t_j \rightarrow t_j + T$ in $y_{T,j}$, where $j$ indicates the...
The parameter calibration results comparing filtering and smoothing are shown in Figure C1. For $T = 360$ days, filtering and smoothing are identical. For $T = 90$ days, the filtering approach leverages the data from one season at a time to update the parameters $\theta$. The final parameter errors for the $T = 90$ days filtering and smoothing approaches are similar, and are both significantly less than the error for the $T = 360$ days case. This result also demonstrates that the improvement in parameter estimates with $T = 90$ days, compared to $T = 360$ days, is a result of increased temporal information, rather than the dimensionality of the data space alone. Depending on the application, the filtering approach may be beneficial for computational efficiency, since it reduces the length of the forward model simulations, instead of requiring temporal integration over the timescale corresponding to an impactful low-frequency cycle.

Appendix D: Gaussian Process Regression Validation and CES Computational Cost

We validate the Gaussian process (GP) regression by comparing the predictions of the GP at the true parameters $G_{GP}(\theta^\dagger)$ to the general circulation model (GCM) output at the true parameters $G(\theta^\dagger)$. The GP is not trained with an input-output pair at the true parameters $\theta^\dagger$. Therefore, $G(\theta^\dagger)$ is an out-of-sample prediction. The GCM output and the GP regression prediction for precipitation at $\theta^\dagger$ is shown in Figure D1. There is a mean absolute error of 0.012 mm/day in the GP prediction.

Given the SVD transformation into a decorrelated space (see Section 2.3 and Appendix A), the GP training and evaluations are parallelizable. The execution time for the GP training for the present application is $\mathcal{O}(\text{seconds})$. The execution time for GP evaluations (forward pass) is $\mathcal{O}(\text{milliseconds})$, substantially less than the GCM execution time we anticipate for our target application of climate model uncertainty quantification. However, the cost of calibrate-emulate-sample (CES) is serial, stemming from MCMC, which is challenging to parallelize. The cost of CES will scale linearly with the cost of GP evaluations ($\mathcal{O}(M \cdot \text{milliseconds})$ for the present application, where $M$ is the number of Markov Chain Monte Carlo steps).
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Figure D1. Annually averaged precipitation (T = 360) from the general circulation model at the true parameters \( \theta^* \) (\( G(\theta^*) \)) and predicted by the Gaussian process regression \( Q_{GP}(\theta^*) \).

Data Availability Statement
All computer code used in this paper is open source. The code is available at (http://doi.org/10.5281/zenodo.5138467). An open-source Julia implementation of CES is accessible at (https://github.com/ClimA/CalibrateEmulateSample.jl).
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