

Uncertainty decomposition in climate-change impact assessments: a Bayesian perspective

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Abstract

A climate-impact projection usually consists of several stages, and the uncertainty of the projection is known to be quite large. It is necessary to assess how much each stage contributed to the uncertainty. We call an uncertainty quantification method in which relative contribution of each stage can be evaluated as uncertainty decomposition. We propose a new Bayesian model for uncertainty decomposition in climate change impact assessments. The proposed Bayesian model can incorporate uncertainty of natural variability and utilize data in control period. We provide a simple and efficient Gibbs sampling algorithm using the auxiliary variable technique. We compare the proposed method with other existing uncertainty decomposition methods by analyzing streamflow data for Yongdam Dam basin located at Geum River in South Korea.

Keywords: climate change impact study, multimodel ensemble, Bayesian statistics, uncertainty decomposition

1. Introduction

The projected impacts of climate change involve large uncertainties. Considering these large uncertainties is critical to establish robust policies. For example, a water management plan which does not consider uncertainty causes large losses if extreme events occur that are not predicted by a projection of streamflow. Thus, quantifying uncertainties of a climate-impact projections has received significant attention.

A climate-impact projection consists of multiple stages. For example, a streamflow projection includes four stages: emission scenarios, global circulation models (GCMs), bias-correction techniques and hydrological models. Emission scenarios describe how future greenhouse gases emissions could evolve based on the set of assumptions about future energy use, technological change and population levels. GCMs are computational models simulating temperature and precipitation under the selected emission scenario. Bias correction techniques remove the systematic biases of the GCM outputs with respect to the observed data. Hydrological models produce projected values of the streamflow by using projected temperature and precipitation as an input. The projection of the streamflow in past can also be produced but in this case the projection is based on observed greenhouse gas emissions instead of the emission scenario. We use the terms *control period* and *scenario period* to distinguish the past and future periods of the projection. Usually the uncertainty of a climate-impact projection is measured by the projected data in the scenario period.

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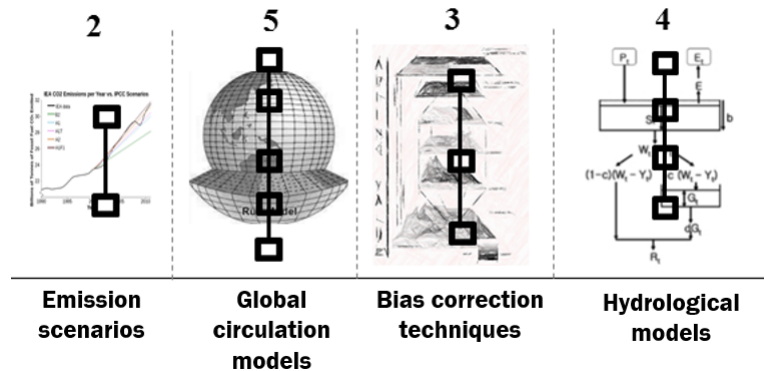


Figure 1: An example of the hydrological projection consisting of four stages.

A climate-impact projection usually gathers multiple projection results based on several scenarios/models/techniques in each stage. Figure 1 is an example of a streamflow projection generated by two emission scenarios, five GCMs, three bias correction techniques and four hydrological models, which results in $120 (= 2 \times 5 \times 3 \times 4)$ streamflow projection results.

The total of the uncertainties, is a mixture of uncertainties produced in each stage. However, quantifying the uncertainties of stages is not straightforward since we have only projected values obtained after the last stage. Various methods have been proposed in order to quantify uncertainties of each stage (e.g., Wilby and Harris, 2006; Bastola *et al.*, 2011; Najafi *et al.*, 2011; Dobler *et al.*, 2012; Lee *et al.*, 2017). Most of these studies quantified uncertainty of each source separately and compared the magnitudes of them. These studies, however, do not guarantee that the total sum of uncertainties of the multiple sources is equal to the total uncertainty, therefore they cannot assess the relative contribution of each source to the total uncertainty. It is also difficult to identify how uncertainty is propagated as uncertainty sources are added. For example, we would like to examine how uncertainty accumulates in a combination of emissions scenarios and GCMs compared to uncertainty with only emissions scenarios alone.

To address these issues, there is a need to quantify the uncertainty of each source such that it is possible to evaluate the relative contribution to the total uncertainty of each uncertainty source. We call such analysis *uncertainty decomposition*. One of the simplest approaches for uncertainty decomposition is to quantify the contribution of the each uncertainty source by partitioning total variance of the projected values into components attributable to the individual sources, as is done in the analysis of variance (ANOVA) framework. Based on the ANOVA approach, Yip *et al.* (2011) decomposed the total uncertainty of the surface air temperature projection into uncertainties of natural variability, emission scenarios, GCMs and interactions. Using the same method, Bosshard *et al.* (2013) decomposed the total uncertainty of the streamflow projection into uncertainties of GCMs, statistical post-processing, hydrological models and their interactions. On the other hand, Kim *et al.* (2019) proposed a cumulative uncertainty method that provides a unified view of how uncertainty is propagated as each source of the projection proceeds. Kim *et al.* (2019) considered a cascade of several stages (i.e., sources) for streamflow projection and quantified cumulative uncertainties as stages progress. Kim *et al.* (2019) defined uncertainties of individual stages as the successive differences of the cumulative uncertainties.

Even though several studies on uncertainty decomposition have been proposed, there are still many areas to be improved. The first problem of the existing methods is that they do not fully utilize the

information in projected values. Note that usually projected values are given at every month, but they use the average of all projected values instead of individual values. For example, the previous methods cannot be incorporated interannual variations. This is because the methods quantify uncertainty based on the averages of projected values over the future period (e.g., 50 years from 2030).

The second problem is that there is no room to use observed values in the analysis. Climate projection typically consists of two periods - control and scenario periods. In the control periods, we have projected values as well as observed values of a meteorological/hydrological variable of interest, while we have only projected values in the scenario period. The existing methods totally ignore the projected and observed values in the control period that may not yield a fully efficient result.

This study provides a Bayesian model for decomposing the total uncertainty into uncertainties of multiple sources. A Bayesian approach to analyze multimodel ensemble data has recently attracted much attention. Bayesian methods can express the complex relationships within the data, and between the data and prior knowledge in a transparent and clear structure. In addition, Bayesian methods provide a probabilistic representation of uncertainty. Tebaldi *et al.* (2005) is the initial work of the Bayesian uncertainty quantification. They obtained a posterior density of mean temperature change based on temporally averaged values of historic observations, GCM outputs for control period and GCM outputs for a scenario period. Tebaldi and Sansó (2009) introduced a Bayesian model that simultaneously analyzed temperature and precipitation. Smith *et al.* (2009) proposed a single Bayesian model that simultaneously deal with data for several regions known as the “multivariate” model in the sense that projected/observed values over different regions are treated as a multivariate response. Unlike the model of Tebaldi *et al.* (2005), which treated the data for the single region, the multivariate model of Smith *et al.* (2009) can distinguish biases of the simulators to the true climate values. Buser *et al.* (2009) analyzed annual data instead of temporally averaged values. The use of annual data makes it possible to take into account natural variability and linear trends in control and scenario periods.

Our proposed Bayesian model improves existing methods of decomposing the total uncertainty. First, it can model changes of mean as well as interannual variations, which makes it possible to incorporate the uncertainty of natural variability. As an illustration, we use a model in which the mean of projected values can change. In addition, the Bayesian model can utilize the projected and observed values in the observational period scientifically, which we believe leads to a better conclusion.

This article is organized as follows. Section 2 summarizes existing uncertainty decomposition methods. Section 3 provides a proposed Bayesian uncertainty decomposition method. Section 4 describes the study area and data involved and compares the proposed method with other uncertainty decomposition methods by analyzing the data. Finally, Section 5 provides the conclusion.

Terminology. Throughout this article, we refer to both emission scenarios, GCMs, bias correction techniques and hydrological models as “simulators”. There are two reasons: 1) to prevent confusion with statistical “models” for the simulator outputs and the observations; 2) to simplify sentences, e.g., we write “simulator combinations” instead of “combinations of model/scenario/technique”.

2. Review of uncertainty decomposition methods

In this section, we review existing methods for uncertainty decomposition. Suppose there are K many stages of a simulator cascade for projection. For example, a simulator cascade for streamflow projection generally consists of four stages: emission scenarios, GCMs, bias correction techniques and hydrological models. Suppose that for stage k , there are n_k simulators denoted by \mathcal{X}_k . Let Y_v be the projected value of a meteorological/hydrological variable of interest at the scenario period when

specific simulators in \mathbf{v} are used in projection, where $\mathbf{v} \in \mathcal{X} := \prod_{k=1}^K \mathcal{X}_k$. For example, if $\mathbf{v} = (2, 4, 3)$, then $Y_{\mathbf{v}}$ is the projected value when simulator 2 in stage 1, simulator 4 in stage 2 and simulator 3 in stage 3 are used.

First we note that the total uncertainty can be defined as

$$U_{\text{tot}} = U(\{Y_{\mathbf{v}} : \mathbf{v} \in \mathcal{X}\}),$$

where U denotes an uncertainty measure such as range, variance or standard deviation. The objective of the uncertainty decomposition is to quantify the uncertainty U_k of stage k for $k = 1, \dots, K$, which satisfies $\sum_{k=1}^K U_k = U_{\text{tot}}$.

The ANOVA approach (Yip *et al.*, 2011; Bosshard *et al.*, 2013) is based on the ANOVA decomposition. In the terminology of ANOVA, the stages are ‘‘effects’’, which could have an influence on the variability of the variable. The ANOVA approach quantifies the influence of the each effect by partitioning the total sum of the squares into sums of squares due to individual effects and the sum of squares due to errors. The uncertainty of stage k defined by the ANOVA approach is then

$$U_k^{\text{anova}} = \frac{1}{n_k} \sum_{i \in \mathcal{X}_k} (\bar{Y}_{k,i} - \bar{Y})^2,$$

which is equivalent to the sum of squares due to stage k divided by the number of projected values $|\mathcal{X}|$. Here \bar{Y} denotes the average of all projection values and $\bar{Y}_{k,i}$ denotes the average of projection values where simulator i in stage k is involved, that is, $\bar{Y} = \sum_{\mathbf{v} \in \mathcal{X}} Y_{\mathbf{v}} / |\mathcal{X}|$ and $\bar{Y}_{k,i} = \sum_{\mathbf{v}: v_k=i} Y_{\mathbf{v}} / \prod_{l=1, l \neq k}^K n_l$. We have relationship such that

$$U_{\text{tot}} = \sum_{k=1}^K U_k^{\text{anova}} + U_{\text{err}},$$

where U_{tot} is equal to $(1/|\mathcal{X}|) \sum_{\mathbf{v} \in \mathcal{X}} (Y_{\mathbf{v}} - \bar{Y})^2$ and U_{err} represents the variability not explained by the effects. Since we can write $U_k^{\text{anova}} = U^{\text{var}}(\{\bar{Y}_{k,i} : i \in \mathcal{X}_k\})$ with U^{var} being the variance measure, the ANOVA approach can be generalized to any other uncertainty measure U by replacing U^{var} with U . This generalization is equivalent to the ‘‘standard method’’ referred in Kim *et al.* (2019).

The disadvantage of the ANOVA approach is that it ignores relationships or interactions between stages. To clarify the problem, consider a simple example where there are two stages having two simulators each. Suppose that four projection values are given as $Y_{(1,1)} = 0, Y_{(1,2)} = 0, Y_{(2,1)} = 10$ and $Y_{(2,2)} = -10$. Even though the final projection values significantly change with respect to the choice of the simulator in the first stage, the uncertainty of the first stage quantified by using the ANOVA approach is 0 since $\bar{Y}_{1,1} = \bar{Y}_{1,2} = 0$.

To overcome this problem, Lee *et al.* (2017) proposed an uncertainty quantification method considering the interactions between simulators from different stages. Lee *et al.* (2017) defined the uncertainty of stage k by

$$U_k^{\text{lee}} = \frac{1}{n_k} \sum_{i \in \mathcal{X}_k} U(\{Y_{\mathbf{v}} : v_k = i\}).$$

For the previously described example, the uncertainty of the first stage is 50 when the variance measure is used. In this case, Lee’s approach quantifies the uncertainties of individual stages more reasonably than the ANOVA approach. But Lee’s approach does not decompose uncertainty properly because $\sum_{k=1}^K U_k^{\text{lee}} \neq U_{\text{tot}}$.

Kim *et al.* (2019) proposed the cumulative uncertainty approach that is the modified version of Lee's approach. The cumulative uncertainty approach captures the interaction effect reasonably and decomposes the total uncertainty properly. Kim *et al.* (2019) defines cumulative uncertainty upto stage k by

$$\text{CU}_k = \frac{1}{n_{k+1} \cdots n_K} \sum_{i_k \in \mathcal{X}_K} \cdots \sum_{i_{k+1} \in \mathcal{X}_{k+1}} U(\{Y_{\mathbf{v}} : v_{k+1} = i_{k+1}, \dots, v_K = i_K\}).$$

The cumulative uncertainty can be viewed as the average of conditional uncertainties of the stages up to k given the simulators in stages after k . Then they defines uncertainty of stage k by

$$U_k^{\text{cu}} = \text{CU}_k - \text{CU}_{k-1}.$$

with $\text{CU}_0 = 0$. Kim *et al.* (2019) showed that U_k^{cu} 's are nonnegative for most reasonable uncertainty measures including range, variance and standard deviation. The cumulative uncertainty up to the last stage $\text{CU}_K = U(\{Y_{\mathbf{v}} : \mathbf{v} \in \mathcal{X}\})$ is equal to the total uncertainty; therefore, by definition, the total uncertainty is equal to the sum of the uncertainties of stages.

3. Bayesian model for uncertainty decomposition

3.1. Model

For the control period, we let $X_{t,\mathbf{w}}$ be the projected value of a meteorological/hydrological variable of interest at time $t \in \mathcal{T}_c$ when specific simulators in \mathbf{w} are used in projection, where $\mathbf{w} \in \prod_{k=2}^K \mathcal{X}_k$. Here \mathcal{T}_c denotes the set of time points in the control period. For the control period, emission scenarios are not used for simulator combinations, thus the projection is generated by the simulators of stage 2 to stage K . We let $X_{t,o}$ be the observed value at time $t \in \mathcal{T}_c$. Similarly, we let $Y_{t,\mathbf{v}}$ be the projected value of time $t \in \mathcal{T}_s$ at the scenario period when $\mathbf{v} \in \mathcal{X}$ simulator combination is used, where \mathcal{T}_s denotes the set of time points in the scenario period. In this paper, we consider annual data, that is, elements in \mathcal{T}_c and \mathcal{T}_s are years in the control and scenario periods, respectively. Thus we can ignore seasonal variation. Note that we do not observe the "true" value $Y_{t,o}$ in the scenario period.

For the observed values in the control period, our Bayesian model assumes

$$X_{t,o} \stackrel{\text{ind}}{\sim} N(\mu_o + \gamma_c t, \sigma_o^2), \quad t \in \mathcal{T}_c,$$

where γ_c represents a linear time trend, σ_o^2 represents an interannual variation. Here we centered the time variable t around the median year of the control period so that μ_o can be interpreted as the (true) mean value of the meteorological/hydrological variable at the median year.

For the description of the distributions of the projected values in the control period, we let $\Delta\mu_{c,\mathbf{w}}$ denote the bias of the mean of $X_{t,\mathbf{w}}$ relative to μ_o and $\sigma_{c,\mathbf{w}}^2$ denote the interannual variation of $X_{t,\mathbf{w}}$. We permit the linear time trends in the projected values to be the same as the trend in the observed values. Our Bayesian model assumes that the projected values in the control period are distributed as

$$X_{t,\mathbf{w}} \stackrel{\text{ind}}{\sim} N(\mu_o + \Delta\mu_{c,\mathbf{w}} + \gamma_c t, \sigma_{c,\mathbf{w}}^2), \quad t \in \mathcal{T}_c, \mathbf{w} \in \prod_{k=2}^K \mathcal{X}_k.$$

We also centered the time variable t around the median year of the scenario period. We let μ_s be the mean of the all projected values at the median year in the scenario period and $\Delta\mu_{s,\mathbf{v}}$ and $\sigma_{s,\mathbf{v}}^2$

be the bias of the mean of $Y_{t,\mathbf{v}}$ relative to μ_s and the interannual variation of $Y_{t,\mathbf{v}}$, respectively. We let the linear time trends in the projected values be the same for all the simulator combinations. Our Bayesian model assumes that the projected values in the scenario period are distributed as

$$Y_{t,\mathbf{v}} \stackrel{\text{ind}}{\sim} N\left(\mu_s + \Delta\mu_{s,\mathbf{v}} + \gamma_s t, \sigma_{s,\mathbf{v}}^2\right), \quad t \in \mathcal{T}_s, \mathbf{v} \in \prod_{k=1}^K \mathcal{X}_k.$$

We impose a certain structural assumption on $\Delta\mu_{c,\mathbf{w}}$, $\Delta\mu_{s,\mathbf{v}}$, $\sigma_{\mathbf{w}}^2$, and $\sigma_{\mathbf{v}}^2$ for uncertainty decomposition. We assume that

$$\begin{aligned} \Delta\mu_{c,\mathbf{w}} &= \sum_{k=2}^K \Delta\mu_{c,k,w_k}, & \sigma_{c,\mathbf{w}}^2 &= \sigma_o^2 + \sum_{k=2}^K \sigma_{k,w_k}^2 \\ \Delta\mu_{s,\mathbf{v}} &= \sum_{k=1}^K \Delta\mu_{s,k,v_k}, & \sigma_{s,\mathbf{v}}^2 &= \sigma_{1,v_1}^2 + \nu \sum_{k=2}^K \sigma_{k,v_k}^2 \end{aligned}$$

where $\sigma_{k,i}^2$ represents the interannual variation of simulator i in stage k , $\Delta\mu_{c,k,i}$'s and $\Delta\mu_{s,k,i}$'s represent the biases of simulator i in stage k in the control and scenario periods respectively, and ν is the multiplicative change in simulator interannual variations between the control and scenario periods. That is, we assume that the bias (or the interannual variation) of each simulator combination is the sum of the biases (or the interannual variation) of the individual simulators that compose the simulator combination.

Note that we assume that $\{\sigma_{k,i}^2 : i \in \mathcal{X}_k, k = 2, \dots, K\}$ are shared by the control and scenario periods to borrow information in the control period for the inference of the scenario period.

Our Bayesian model is similar to the model of Buser *et al.* (2009). We simplify the model by removing unidentifiable parameters such as additive changes in simulator biases between the control and scenario periods. We complete such a modification since our focus is to decompose the total uncertainty rather than provide physical meanings of the model.

3.2. Uncertainty decomposition

The total uncertainty can be understood as the variance of all detrended projected values $\mathcal{Y}_s = \{Y_{t,\mathbf{v}}^{\text{det}} : t \in \mathcal{T}_s, \mathbf{v} \in \mathcal{X}\}$, where $Y_{t,\mathbf{v}}^{\text{det}} = Y_{t,\mathbf{v}} - \gamma t$ and $\mathcal{X} = \prod_{k=1}^K \mathcal{X}_k$. Note that the distribution of a randomly selected value from \mathcal{Y}_s is a mixture of Gaussians given as

$$f(y) = \frac{1}{|\mathcal{X}|} \sum_{\mathbf{v}} N\left(\mu_s + \Delta\mu_{s,\mathbf{v}}, \sigma_{s,\mathbf{v}}^2\right).$$

We define the total uncertainty τ_{tot} as the variance of $f(y)$. By the variance formula of mixture distributions, τ_{tot} is given as

$$\tau_{\text{tot}} = \frac{1}{|\mathcal{X}|} \left[\sum_{\mathbf{v}} \sigma_{s,\mathbf{v}}^2 + \sum_{\mathbf{v}} (\Delta\mu_{s,\mathbf{v}} - \overline{\Delta\mu})^2 \right],$$

where $\overline{\Delta\mu} = \sum_{\mathbf{v}} \Delta\mu_{s,\mathbf{v}} / |\mathcal{X}|$.

Under the assumption in Subsection 3.1, τ_{tot} can be rewritten as

$$\tau_{\text{tot}} = \sum_{k=1}^K \tau_k,$$

where

$$\tau_k = \frac{1}{n_k} \sum_{i=1}^{n_k} \left[\nu_k \sigma_{k,i}^2 + \left(\Delta\mu_{k,i} - \overline{\Delta\mu_k} \right)^2 \right],$$

where $\nu_k = 1$ for $k = 1$, $\nu_k = \nu$ for $k \geq 2$ and $\overline{\Delta\mu_k} = \sum_{i=1}^{n_k} \Delta\mu_{s,k,i} / n_k$. We can interpret τ_k as the uncertainty raised by stage k . The total uncertainty is then the sum of uncertainties of each stage, which is a property required by Jones (2000). Furthermore, we define the uncertainty of each simulator by

$$\tau_{k,i} = \frac{1}{n_k} \left(\nu_k \sigma_{k,i}^2 + \left(\Delta\mu_{k,i} - \overline{\Delta\mu_k} \right)^2 \right)$$

which satisfies a natural requirement that $\tau_k = \sum_{i=1}^{n_k} \tau_{k,i}$.

Lee *et al.* (2017) also quantified the uncertainty of each simulator, but they considered only the dispersion of the projected values which use the corresponding simulator, not the bias to the total mean. However, our method quantifies the simulator uncertainty more reasonably by considering both bias and dispersion.

The above definition of the total uncertainty, uncertainties of each stage and each simulator are given as conditional on the parameters. Our Bayesian method estimates the posterior distribution of the parameters as well as the uncertainties based on the data. Considering the posterior distribution, we report the posterior expectations of the uncertainties as our estimated uncertainties. Let Z be the set of all projected and observed data and $E(\cdot | Z)$ denote the posterior expectation given the data Z . Then we define the total uncertainty, uncertainties of each stage and each simulator as

$$E(\tau_{\text{tot}} | Z) = \frac{1}{|X|} E \left[\sum_{\nu} \sigma_{s,\nu}^2 + \sum_{\nu} \left(\Delta\mu_{s,\nu} - \overline{\Delta\mu} \right)^2 | Z \right], \quad (3.1)$$

$$E(\tau_k | Z) = \frac{1}{n_k} \sum_{i=1}^{n_k} E \left[\nu_k \sigma_{k,i}^2 + \left(\Delta\mu_{k,i} - \overline{\Delta\mu_k} \right)^2 | Z \right], \quad (3.2)$$

$$E(\tau_{k,i} | Z) = \frac{1}{n_k} E \left[\nu_k \sigma_{k,i}^2 + \left(\Delta\mu_{k,i} - \overline{\Delta\mu_k} \right)^2 | Z \right],$$

respectively.

3.3. Prior and Markov chain Monte Carlo algorithm

We let the prior of our model be as

$$\begin{aligned} \mu_o, \mu_s &\sim N(0, q_\mu^2), \\ \gamma_c, \gamma_s &\sim N(0, q_\gamma^2), \\ \Delta\mu_{c,k,i}, \Delta\mu_{s,k,i} | \phi &\sim N(0, \phi^2), \\ (\phi^2)^{-1} &\sim \text{Gamma}(a_\phi, b_\phi), \\ (\sigma_o^2)^{-1}, (\sigma_{k,i}^2)^{-1} &\sim \text{Gamma}(a_\sigma, b_\sigma), \\ \nu^{-1} | \alpha &\sim \text{Gamma}(\alpha, \alpha), \\ \alpha &\sim \text{Gamma}(a_\alpha, b_\alpha). \end{aligned}$$

We take a hierarchical prior for the simulator biases to prevent a nonidentifiable problem. Note that if any constant C is added to all the $\Delta\mu_{s,k,i}$'s and subtracted from all the $\Delta\mu_{s,k',i}$'s for $k' \neq k$, the distributions of the data remain unchanged since they are determined only by the sums of the simulator biases $\Delta\mu_{s,v} = \sum_{k=1}^K \Delta\mu_{s,k,v_k}$. The identifiability can be improved by the prior $N(0, \phi^2)$ with $(\phi^2)^{-1} \sim \text{Gamma}(a_\phi, b_\phi)$, which implies the constraint on the $\Delta\mu_{s,k,i}$'s such that $\sum_{i=1}^{n_k} \Delta\mu_{s,k,i} = 0$ (Jackman, 2009).

The hyperparameters $a_\phi, b_\phi, a_\sigma, b_\sigma, a_\alpha, b_\alpha, q_\mu^2$, and q_γ^2 are selected so that the corresponding priors are diffused. In our case studies, we use $a_\phi = b_\phi = a_\sigma = a_\alpha = b_\alpha = 0.001$ and $q_\mu^2 = q_\gamma^2 = 10,000$.

The joint distribution of the data and parameters are proportional to

$$\begin{aligned} & \left(\frac{1}{\sigma_o^2} \right)^{\frac{|\mathcal{T}_c|}{2}} \exp \left(-\frac{1}{2\sigma_o^2} \sum_{t \in \mathcal{T}_c} (X_{t,o} - \mu_o - \gamma_c t)^2 \right) \\ & \times \left(\prod_{\mathbf{w}} \left(\frac{1}{\sigma_{c,w}^2} \right)^{\frac{|\mathcal{T}_c|}{2}} \right) \exp \left(-\sum_{\mathbf{w}} \frac{1}{2\sigma_{c,w}^2} \sum_{t \in \mathcal{T}_c} (X_{t,w} - \mu_o - \Delta\mu_{c,w} - \gamma_c t)^2 \right) \\ & \times \left(\prod_{\mathbf{v}} \left(\frac{1}{\sigma_{s,v}^2} \right)^{\frac{|\mathcal{T}_s|}{2}} \right) \exp \left(-\sum_{\mathbf{v}} \frac{1}{2\sigma_{s,v}^2} \sum_{t \in \mathcal{T}_s} (Y_{t,v} - \mu_s - \Delta\mu_{s,v} - \gamma_s t)^2 \right) \\ & \times \exp \left(-\frac{1}{2\phi^2} \left(\sum_{k=2}^K \sum_{i=1}^{n_k} \Delta\mu_{c,k,i}^2 + \sum_{k=1}^K \sum_{i=1}^{n_k} \Delta\mu_{s,k,i}^2 \right) \right) \\ & \times \exp \left(-\frac{1}{2q_\mu^2} (\mu_o^2 + \mu_s^2) \right) \times \exp \left(-\frac{1}{2q_\gamma^2} (\gamma_c^2 + \gamma_s^2) \right) \\ & \times \left[(\sigma_o^2)^{-a_\sigma+1} e^{-b_\sigma(\sigma_o^2)^{-1}} \right] \times \left[\prod_{k=1}^K \prod_{i=1}^{n_k} (\sigma_{k,i}^2)^{-a_\sigma+1} e^{-b_\sigma(\sigma_{k,i}^2)^{-1}} \right] \\ & \times [(\gamma)^{-a_\alpha+1} e^{-a_\alpha(\gamma^{-1})}] \times [\alpha^{a_\alpha-1} e^{-b_\alpha\alpha}] \times [(\phi^2)^{-a_\phi+1} e^{-b_\phi(\phi^2)^{-1}}]. \end{aligned}$$

The conditional distributions of the parameters (except the interannual variation parameters) for the Gibbs sampling algorithm are given as:

$$\mu_o | - \sim N \left(\tilde{\mu}_o, \frac{1}{|\mathcal{T}_c| \left((\sigma_o^2)^{-1} + \sum_{\mathbf{w}} (\sigma_{c,w}^2)^{-1} \right) + (q_\mu^2)^{-1}} \right), \quad (3.3)$$

$$\mu_s | - \sim N \left(\tilde{\mu}_s, \frac{1}{|\mathcal{T}_s| \sum_{\mathbf{v}} (\sigma_{s,v}^2)^{-1} + (q_\mu^2)^{-1}} \right), \quad (3.4)$$

$$\gamma_c | - \sim N \left(\tilde{\gamma}_c, \frac{1}{\left((\sigma_o^2)^{-1} + \sum_{\mathbf{w}} (\sigma_{c,w}^2)^{-1} \right) \sum_{t \in \mathcal{T}_c} t^2 + (q_\gamma^2)^{-1}} \right), \quad (3.5)$$

$$\gamma_s | - \sim N \left(\tilde{\gamma}_s, \frac{1}{\left(\sum_{\mathbf{v}} (\sigma_{s,v}^2)^{-1} \right) \sum_{t \in \mathcal{T}_s} t^2 + (q_\gamma^2)^{-1}} \right), \quad (3.6)$$

$$(\phi^2)^{-1}|-\sim \text{Gamma}\left(a_\phi + \frac{1}{2}n_1 + \sum_{k=2}^K n_k, b_\phi + \frac{1}{2}\left(\sum_{k=2}^K \sum_{i=1}^{n_k} \Delta\mu_{c,k,i}^2 + \sum_{k=1}^K \sum_{i=1}^{n_k} \Delta\mu_{s,k,i}^2\right)\right), \quad (3.7)$$

$$\Delta\mu_{c,k,i}|-\sim N\left(\widetilde{\Delta\mu}_{c,k,i}, \frac{1}{|\mathcal{T}_c| \sum_{\mathbf{w}:w_k=i} (\sigma_{c,\mathbf{w}}^2)^{-1} + (\phi^2)^{-1}}\right), \quad (3.8)$$

$$\Delta\mu_{s,k,i}|-\sim N\left(\widetilde{\Delta\mu}_{s,k,i}, \frac{1}{|\mathcal{T}_s| \sum_{\mathbf{v}:v_k=i} (\sigma_{s,\mathbf{v}}^2)^{-1} + (\phi^2)^{-1}}\right), \quad (3.9)$$

where

$$\begin{aligned} \widetilde{\mu}_o &= \frac{(\sigma_o^2)^{-1} \sum_{t \in \mathcal{T}_c} (X_{t,o} - \gamma_c t) + \sum_{\mathbf{w}} \left\{ (\sigma_{c,\mathbf{w}}^2)^{-1} \sum_{t \in \mathcal{T}_c} (X_{t,\mathbf{w}} - \Delta\mu_{c,\mathbf{w}} - \gamma_c t) \right\}}{|\mathcal{T}_c| (\sigma_o^2)^{-1} + |\mathcal{T}_c| \sum_{\mathbf{w}} (\sigma_{c,\mathbf{w}}^2)^{-1}}, \\ \widetilde{\mu}_s &= \frac{\sum_{\mathbf{v}} \left\{ (\sigma_{s,\mathbf{v}}^2)^{-1} \sum_{t \in \mathcal{T}_s} (Y_{t,\mathbf{v}} - \Delta\mu_{s,\mathbf{v}} - \gamma_s t) \right\}}{|\mathcal{T}_s| \sum_{\mathbf{v}} (\sigma_{s,\mathbf{v}}^2)^{-1}}, \\ \widetilde{\gamma}_c &= \frac{(\sigma_o^2)^{-1} \sum_{t \in \mathcal{T}_c} t(X_{t,o} - \mu_o) + \sum_{\mathbf{w}} (\sigma_{c,\mathbf{w}}^2)^{-1} \sum_{t \in \mathcal{T}_c} t(X_{t,\mathbf{w}} - \mu_o - \Delta\mu_{c,\mathbf{w}})}{\left((\sigma_o^2)^{-1} + \sum_{\mathbf{w}} (\sigma_{c,\mathbf{w}}^2)^{-1} \right) \sum_{t \in \mathcal{T}_c} t^2 + (q_y^2)^{-1}}, \\ \widetilde{\gamma}_s &= \frac{\sum_{\mathbf{v}} (\sigma_{s,\mathbf{v}}^2)^{-1} \sum_{t \in \mathcal{T}_s} t(Y_{t,\mathbf{v}} - \mu_s - \Delta\mu_{s,\mathbf{v}})}{\left(\sum_{\mathbf{v}} (\sigma_{s,\mathbf{v}}^2)^{-1} \right) \sum_{t \in \mathcal{T}_s} t^2 + (q_y^2)^{-1}}, \\ \widetilde{\Delta\mu}_{c,k,i} &= \frac{\sum_{\mathbf{w}:w_k=i} (\sigma_{c,\mathbf{w}}^2)^{-1} \sum_{t \in \mathcal{T}_c} (X_{t,\mathbf{w}} - \mu_o - \sum_{h \in \{2, \dots, K\} \setminus \{k\}} \Delta\mu_{c,h,w_h} - \gamma_c t)}{|\mathcal{T}_c| \sum_{\mathbf{w}:w_k=i} (\sigma_{c,\mathbf{w}}^2)^{-1} + (\phi^2)^{-1}}, \\ \widetilde{\Delta\mu}_{s,k,i} &= \frac{\sum_{\mathbf{v}:v_k=i} (\sigma_{s,\mathbf{v}}^2)^{-1} \sum_{t \in \mathcal{T}_s} (Y_{t,\mathbf{v}} - \mu_s - \sum_{h \in \{1, \dots, K\} \setminus \{k\}} \Delta\mu_{s,h,v_h} - \gamma_s t)}{|\mathcal{T}_s| \sum_{\mathbf{v}:v_k=i} (\sigma_{s,\mathbf{v}}^2)^{-1} + (\phi^2)^{-1}}. \end{aligned}$$

The conditional distributions of the interannual variation parameters are awkward forms; therefore, we use an auxiliary variable (Damlén *et al.*, 1999) instead of a straightforward Gibbs sampling algorithm technique. Note that the distributions of $X_{t,\mathbf{w}}$ and $Y_{t,\mathbf{v}}$ can be written as

$$\begin{aligned} X_{t,\mathbf{w}} &\stackrel{d}{=} \mu_o + \Delta\mu_{\mathbf{w}} + \gamma_c t + \epsilon_{t,\mathbf{w},o}^c + \sum_{k=2}^K \epsilon_{t,\mathbf{w},k}^c, \\ Y_{t,\mathbf{v}} &\stackrel{d}{=} \mu_s + \Delta\mu_{\mathbf{v}} + \gamma_s t + \sum_{k=1}^K \epsilon_{t,\mathbf{v},k}^s, \end{aligned}$$

where $\epsilon_{t,\mathbf{w},o}^c \sim N(0, \sigma_o^2)$, $\epsilon_{t,\mathbf{w},k}^c \sim N(0, \sigma_{k,w_k}^2)$, $\epsilon_{t,\mathbf{v},1}^s \sim N(0, \sigma_{1,v_1}^2)$ and $\epsilon_{t,\mathbf{v},k}^s \sim N(0, \nu\sigma_{k,v_k}^2)$ for $k \geq 2$, and they are independent. The joint distribution of $(\epsilon_{t,\mathbf{w},o}^c, \epsilon_{t,\mathbf{w},2}^c, \dots, \epsilon_{t,\mathbf{w},K-1}^c, \epsilon_{t,\mathbf{w},o}^c + \sum_{k=2}^K \epsilon_{t,\mathbf{w},k}^c)$ is the

normal distribution with mean $\mathbf{0}$ and variance

$$\begin{pmatrix} \sigma_o^2 & 0 & \cdots & 0 & \sigma_o^2 \\ 0 & \sigma_{2,w_2}^2 & \cdots & 0 & \sigma_{2,w_2}^2 \\ \vdots & \vdots & \ddots & \vdots & \vdots \\ 0 & 0 & \cdots & \sigma_{K-1,w_{K-1}}^2 & \sigma_{K-1,w_{K-1}}^2 \\ \sigma_o^2 & \sigma_{2,w_2}^2 & \cdots & \sigma_{K-1,w_{K-1}}^2 & \sigma_o^2 + \sum_{k=2}^K \sigma_{k,w_k}^2 \end{pmatrix}.$$

Note that if a multivariate normal random vector \mathbf{x} is partitioned into $\mathbf{x} = (\mathbf{x}_1, \mathbf{x}_2)^\top$, then the distribution of \mathbf{x}_1 conditional on $\mathbf{x}_2 = \mathbf{a}$ is $\mathbf{x}_1 | \mathbf{x}_2 = \mathbf{a} \sim N(\boldsymbol{\mu}_1 + \boldsymbol{\Sigma}_{22}^{-1}(\mathbf{a} - \boldsymbol{\mu}_2), \boldsymbol{\Sigma}_{11} - \boldsymbol{\Sigma}_{12}\boldsymbol{\Sigma}_{22}^{-1}\boldsymbol{\Sigma}_{12}^\top)$ where $\boldsymbol{\mu}_i = \mathbf{E}\mathbf{x}_i$, $\boldsymbol{\Sigma}_{ii} = \text{var}(\mathbf{x}_i)$ for $i = 1, 2$ and $\boldsymbol{\Sigma}_{12} = \text{cov}(\mathbf{x}_1, \mathbf{x}_2)$. Thus the distribution of $(\epsilon_{t,w,o}^c, \epsilon_{t,w,2}^c, \dots, \epsilon_{t,w,K-1}^c)^\top$ conditional on $\epsilon_{t,w,o}^c + \sum_{k=2}^K \epsilon_{t,w,k}^c = X_{t,w} - \mu_o - \Delta\mu_w - \gamma_c t$ is the normal distribution with mean

$$\mathbf{m}_{\epsilon^c, \mathbf{w}} := \frac{X_{t,w} - \mu_o - \Delta\mu_w - \gamma_c t}{\sigma_o^2 + \sum_{k=2}^K \sigma_{k,w_k}^2} \boldsymbol{\Sigma}_{\mathbf{w}}$$

and variance

$$\mathbf{S}_{\epsilon^c, \mathbf{w}} := \text{diag}(\sigma_o^2, \sigma_{2,w_2}^2, \dots, \sigma_{K-1,w_{K-1}}^2) - \frac{1}{\sigma_o^2 + \sum_{k=2}^K \sigma_{k,w_k}^2} \boldsymbol{\Sigma}_{\mathbf{w}} \boldsymbol{\Sigma}_{\mathbf{w}}^\top,$$

where $\boldsymbol{\Sigma}_{\mathbf{w}} = (\sigma_o^2, \sigma_{2,w_2}^2, \dots, \sigma_{K-1,w_{K-1}}^2)^\top$. We generate

$$(\epsilon_{t,w,o}^c, \epsilon_{t,w,2}^c, \dots, \epsilon_{t,w,K-1}^c)^\top \sim N(\mathbf{m}_{\epsilon^c, \mathbf{w}}, \mathbf{S}_{\epsilon^c, \mathbf{w}}) \quad (3.10)$$

and calculate

$$\epsilon_{t,w,K}^c = (X_{t,w} - \mu_o - \Delta\mu_w - \gamma_c t) - \left(\epsilon_{t,w,o}^c + \sum_{k=2}^{K-1} \epsilon_{t,w,k}^c \right). \quad (3.11)$$

Similarly we generate

$$(\epsilon_{t,v,1}^s, \dots, \epsilon_{t,v,K-1}^s)^\top \sim N(\mathbf{m}_{\epsilon^s, \mathbf{v}}, \mathbf{S}_{\epsilon^s, \mathbf{v}}), \quad (3.12)$$

where we define

$$\mathbf{m}_{\epsilon^s, \mathbf{v}} := (Y_{t,v} - \mu_s - \Delta\mu_v - \gamma_s t) \left(\sigma_{1,v_1}^2 + v \sum_{k=2}^K \sigma_{k,v_k}^2 \right)^{-1} \boldsymbol{\Sigma}_{\mathbf{v}}$$

and

$$\mathbf{S}_{\epsilon^s, \mathbf{v}} := \text{diag}(\sigma_{1,v_1}^2, v\sigma_{2,v_2}^2, \dots, v\sigma_{K-1,w_{K-1}}^2) - \left(\sigma_{1,v_1}^2 + v \sum_{k=2}^K \sigma_{k,v_k}^2 \right)^{-1} \boldsymbol{\Sigma}_{\mathbf{v}} \boldsymbol{\Sigma}_{\mathbf{v}}^\top,$$

where $\boldsymbol{\Sigma}_{\mathbf{v}} = (\sigma_{1,v_1}^2, v\sigma_{2,v_2}^2, \dots, v\sigma_{K-1,w_{K-1}}^2)^\top$. Then we calculate

$$\epsilon_{t,v,K}^s = (Y_{t,v} - \mu_s - \Delta\mu_v - \gamma_s t) - \left(\epsilon_{t,v,o}^s + \sum_{k=1}^{K-1} \epsilon_{t,v,k}^s \right). \quad (3.13)$$

Then we can sample the interannual variation parameters from the following conditional distributions

$$(\sigma_o^2)^{-1} | - \sim \text{Gamma} \left(a_\sigma + \frac{|\mathcal{T}_c|}{2} \prod_{k=2}^K n_k, \tilde{b}_{\sigma_o} \right), \quad (3.14)$$

$$(\sigma_{k,i}^2)^{-1} | - \sim \text{Gamma} \left(a_\sigma + \frac{1}{2} \left(\frac{\prod_{h=2}^K n_h}{n_k} |\mathcal{T}_c| \mathbb{I}_{\{k \neq 1\}} + \frac{|\mathcal{X}|}{n_k} |\mathcal{T}_s| \right), \tilde{b}_{\sigma_{k,i}} \right), \quad (3.15)$$

$$(\nu)^{-1} | - \sim \text{Gamma} \left(\alpha + \frac{K-1}{2} |\mathcal{X}| |\mathcal{T}_s|, \tilde{b}_\nu \right), \quad (3.16)$$

where

$$\begin{aligned} \tilde{b}_{\sigma_o} &= b_\sigma + \frac{1}{2} \left\{ \sum_{t \in \mathcal{T}_c} (X_{t,o} - \mu_o - \gamma_c t)^2 + \sum_{\mathbf{w}} \sum_{t \in \mathcal{T}_c} (\epsilon_{t,\mathbf{w},o}^c)^2 \right\}, \\ \tilde{b}_{\sigma_{k,i}} &= b_\sigma + \frac{1}{2} \left\{ \sum_{\mathbf{w}: w_k=i} \sum_{t \in \mathcal{T}_c} (\epsilon_{t,\mathbf{w},k}^c)^2 \mathbb{I}_{\{k \neq 1\}} + \sum_{\mathbf{v}: v_k=i} \sum_{t \in \mathcal{T}_s} (\epsilon_{t,\mathbf{v},k}^s)^2 \right\}, \\ \tilde{b}_\nu &= \alpha + \frac{1}{2} \sum_{\mathbf{v}} \sum_{k=2}^K \left((\sigma_{k,v_k}^2)^{-1} \sum_{t \in \mathcal{T}_s} (\epsilon_{t,\mathbf{v},k}^s)^2 \right). \end{aligned}$$

Lastly, the Metropolis-Hasting algorithm is used for the updating of α . Following Smith *et al.* (2009), we propose $\alpha^* = \alpha e^{U-1/2}$ with U being a random number generated from the uniform distribution on $(0, 1)$ and calculate

$$\begin{aligned} l(\alpha) &= \alpha \log(\alpha) - \log(\Gamma(\alpha)) + (\alpha - 1) \log(\nu) - \alpha \nu - a_\alpha \log(\alpha) - b_\alpha \alpha, \\ l(\alpha^*) &= \alpha^* \log(\alpha^*) - \log(\Gamma(\alpha^*)) + (\alpha^* - 1) \log(\nu) - \alpha^* \nu - a_\alpha \log(\alpha^*) - b_\alpha \alpha^*. \end{aligned}$$

Then set

$$\alpha = \begin{cases} \alpha^*, & \text{with probability } \min(e^{l(\alpha^*)-l(\alpha)}, 1), \\ \alpha, & \text{otherwise.} \end{cases} \quad (3.17)$$

One iteration of the Gibbs sampler is summarized in Algorithm 1.

4. Case studies

4.1. Study area

The target watershed is Yongdam Dam located at the upper basin of the Geum River in South Korea (Figure 2). The Yongdam Dam basin is located at latitude $35^\circ 35' - 36^\circ 00'$ and longitude $127^\circ 20' - 127^\circ 45'$, and it lies on the borders of Chungcheongnam-do, Jeollabuk-do and Gyeongsangnam-do. The basin area is 930 km^2 accounting for about 9.5% of the Geum River basin area. The climate of the Geum River region dominated by the East Asian monsoon, which is divided into a warm and wet summer monsoon and a cold and dry winter monsoon. The monsoon drives an intense precipitation during the summer season, but sometimes it leads to drought. Therefore the projections for precipitation and streamflow are associated with a large uncertainty. Various water-related issues such as frequent extreme droughts and increased water demands have recently occurred in the Geum River region; therefore, uncertainty analysis is becoming increasingly important for water resource planning.

Algorithm 1: One iteration of the Gibbs sampler for Bayesian uncertainty decomposition

```

1 for  $\mathbf{w} \in \prod_{k=2}^K \mathcal{X}_k$  do
2   for  $t \in \mathcal{T}_c$  do
3     Sample  $\epsilon_{t,\mathbf{w},o}^c, \epsilon_{t,\mathbf{w},1}^c, \dots, \epsilon_{t,\mathbf{v},K-1}^c$  according to Eq. (3.10) and compute  $\epsilon_{t,\mathbf{w},K}^s$  according to
       Eq. (3.11)
4   end
5 end
6 for  $\mathbf{v} \in \prod_{k=1}^K \mathcal{X}_k$  do
7   for  $t \in \mathcal{T}_s$  do
8     Sample  $\epsilon_{t,\mathbf{v},1}^s, \dots, \epsilon_{t,\mathbf{v},K-1}^s$  according to Eq. (3.12) and compute  $\epsilon_{t,\mathbf{v},K}^s$  according to Eq.
       (3.13)
9   end
10 end
11 Sample  $\mu_0$  and  $\mu_s$  according to Eq. (3.3) and Eq. (3.4).
12 Sample  $\gamma_c$  and  $\gamma_s$  according to Eq. (3.5) and Eq. (3.6).
13 Sample  $(\phi^2)^{-1}$  according to Eq. (3.7)
14 Sample  $\nu$  according to Eq. (3.16)
15 Sample  $\alpha$  according to Eq. (3.17)
16 Sample  $(\sigma_o^2)^{-1}$  according to Eq. (3.14)
17 for  $k = 1, \dots, K$  do
18   for  $i = 1, \dots, n_k$  do
19     Sample  $\Delta\mu_{c,k,i}$  according to Eq. (3.8) (not for  $k = 1$ )
20     Sample  $\Delta\mu_{s,k,i}$  according to Eq. (3.9)
21     Sample  $(\sigma_{k,i}^2)^{-1}$  according to Eq. (3.15)
22   end
23 end

```

4.2. Data

Daily streamflow data for the Yongdam Dam basin were obtained from the Water Resources Management Information System webpage (<http://www.wamis.go.kr>). For the projection data, we used the emission scenarios, GCMs, bias correction techniques, and hydrological models presented in Table 1 or emission scenarios. We also adopted two representative concentration pathways (RCPs) scenarios, RCP4.5 and RCP8.5. Out of 27 GCMs contained in the Coupled Model Intercomparison Project Phase 5 (CMIP5), we selected the following four GCMs: CESM1-BGC (Moore *et al.*, 2013), HadGEM2-ES (Collins *et al.*, 2011), MIROC-ESM-CHEM (Watanabe *et al.*, 2011), MRI-CGCM3 (Yukimoto *et al.*, 2012). The Katsavounidis-Kuo-Zhang (KKZ) algorithm (Katsavounidis *et al.*, 1994) was used for GCM selection. For bias correction techniques, bias correction spatial disaggregation (BCSD) (Wood *et al.*, 2004) and bias correction constructed analogs (BCCA) (Hidalgo *et al.*, 2008) were used to remove the systematic biases of GCM outputs with respect to the observations. Lastly, we used two hydrological models GR4J (Perrin *et al.*, 2003) and the Tank model (Sugawara 1979). These two hydrological models have been widely applied to hydrological studies including flood control and groundwater modeling (Amireche *et al.*, 2017).

Daily projection and observation data are aggregated to the annual data for two seasons, DJF (December, January and February) and JJA (June, July and August). Our model was applied indepen-

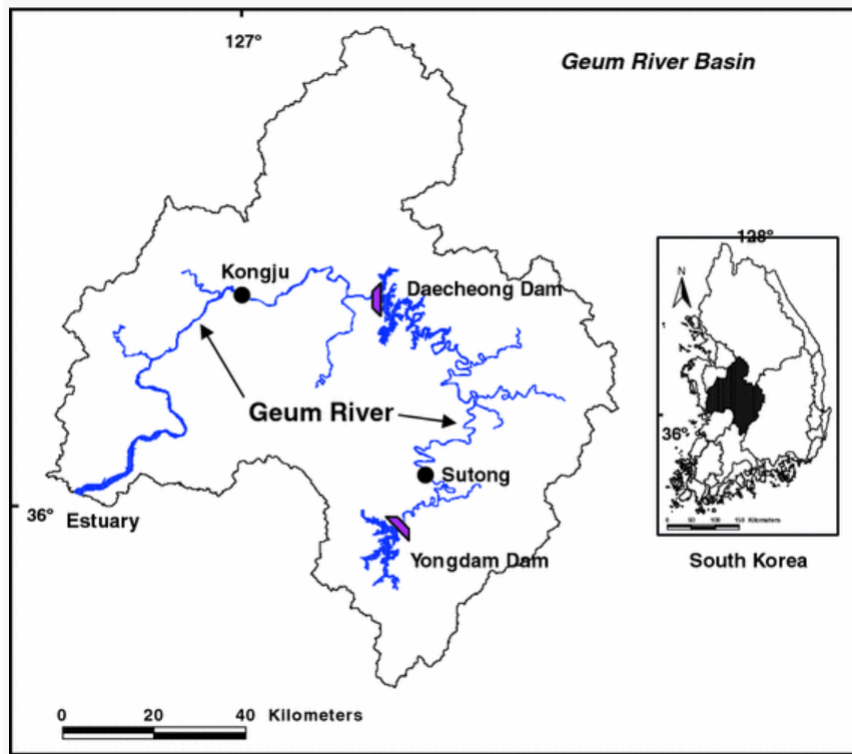


Figure 2: Location of Yongdam Dam basin. This figure is from Lee et al. (2013).

Table 1: Emission scenarios (ESs), global circulation models (GCMs), bias correction techniques (BCs), and hydrological models (HMs) used for streamflow projections

Stage	Simulator
ESs	RCP 4.5, RCP 8.5
GCMs	CESM1-BGC, HadGEM2-ES, MIROC-ESM-CHEM, MRI-CGCM3
BCs	BCCA, BCSD
HMs	GR4J, Tank

dently for both seasons. The control period is from 1976 to 2005 and the scenario period is from 2016 to 2045.

Figures 3 and 4 draw the averaged projection values over the scenario period according to the simulators in each stage. For example, the dots above ‘GR4J’ in the right lower panel of Figure 3 are 16 projection values where hydrological model GR4J was used and the diamond shape dot is the average of the 16 projection values. The dots connected by lines are projection values using the same combination of simulators except the simulator at the x-axis. For the JJA season the hydrological model stage seems to be the most influential to the projection. The two hydrological models yield fairly different projection values. In contrast, the choice of bias correction technique does not appear to have a significant impact on projection values. The two projection values from the same combination of models, except the bias correction technique, are almost similar. For the DJF season, emission scenario and GCM seem to be more influential to the projection than the other stages.

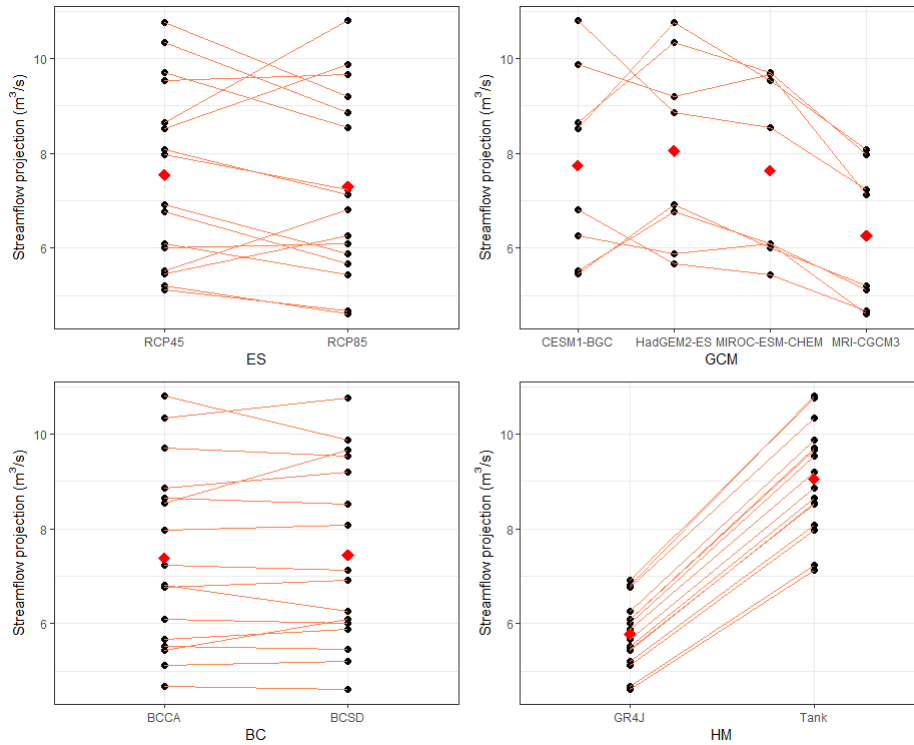


Figure 3: Averaged projection values over the scenario period for the JJA season. We draw averaged projection values according to simulators in each stage. Each diamond shape dot is the average of the circle shaped dots on the same vertical line. The dots connected by lines are projection values using the same combination of simulators except the simulator at the x-axis. ES = emission scenarios; GCM = global circulation model; BC = bias correction technique; HM = hydrological model.

4.3. Estimated posterior distributions

We ran a single Markov chain with length 60,000. We saved every 50th sample after a burn-in period of 10,000 samples and obtained 1,000 samples used for further analysis.

Figures 5 and 6 present the estimated posterior densities of some easily interpretable parameters for the JJA and DJF seasons, respectively. For the JJA season, there is an apparent mean shift between the control and scenario periods but not for the DJF season. For both seasons, the dispersion of μ_s is larger than that of μ_o , which is due to the absence of the true values $Y_{t,o}$. On the other hand, the posterior densities of the natural variability parameters are almost similar for both the control and scenario periods in both seasons, which implies that the natural system is expected not to change much.

There is a reduction in the linear trend for the JJA season. It shows that the degree of increase in streamflow gradually decreases over time. For the DJF season, the linear trend is negative in the control period, but positive in the scenario period.

Tables 2 and 3 present the posterior expectations of bias, interannual variation and uncertainty of each simulator for the JJA and DJF seasons, respectively. For both seasons, the interannual variation of RCP4.5 is greater than that of RCP8.5 and the interannual variation of Tank hydrological model is

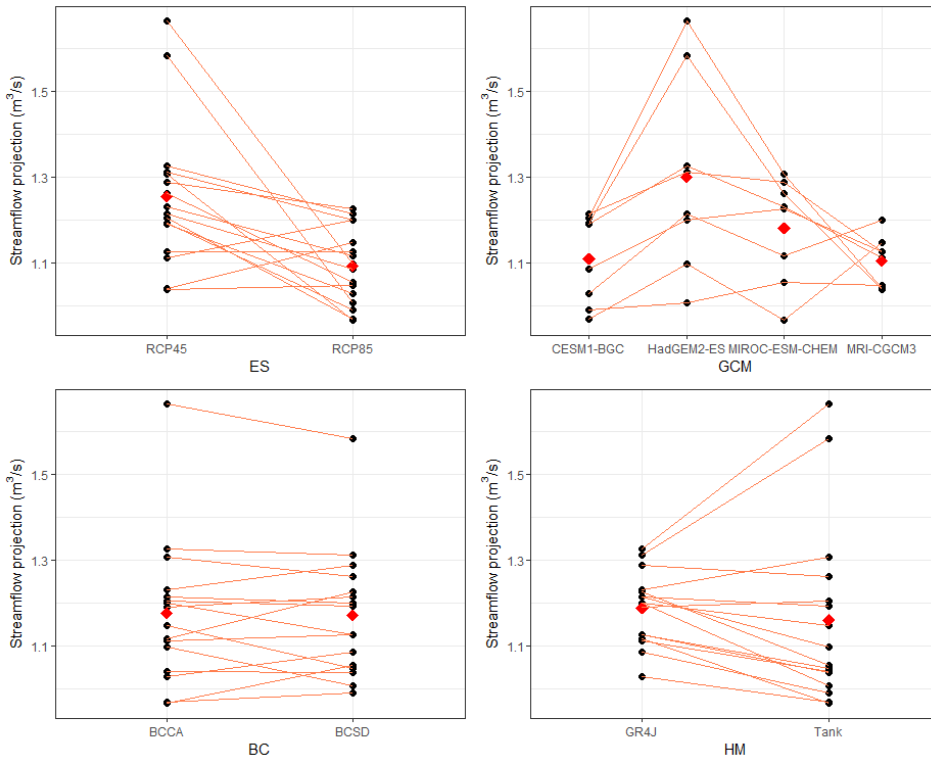


Figure 4: Same as Figure 3, but for the DJF season. ES = emission scenarios; GCM = global circulation model; BC = bias correction technique; HM = hydrological model.

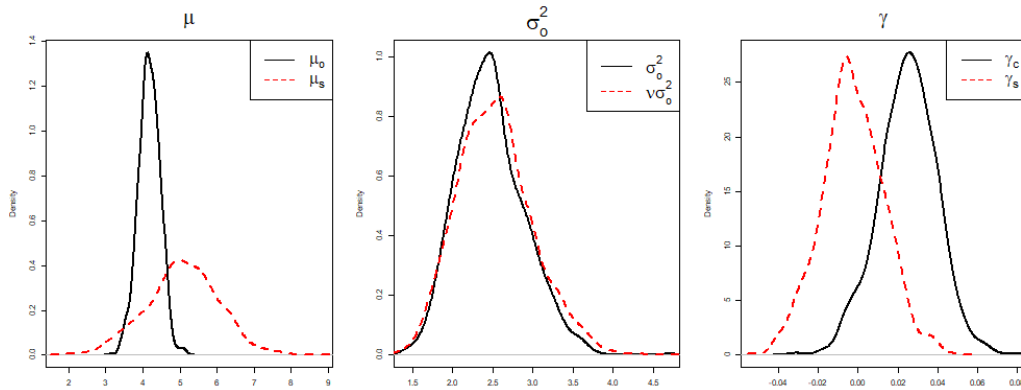


Figure 5: Posterior densities for μ_o and μ_s (left); σ_o^2 and σ_s^2 (middle); γ_c and γ_s (right) for the JJA season. The solid black lines and the red dashed lines indicate the parameters in the control and scenario periods, respectively.

greater than that of GR4J hydrological model. This implies that RCP4.5 and Tank lead to the larger interannual variations. On the other hand, the bias of Tank hydrological model is higher than the bias of the GR4J hydrological model for the JJA season, but the relationship is reverse for the DJF

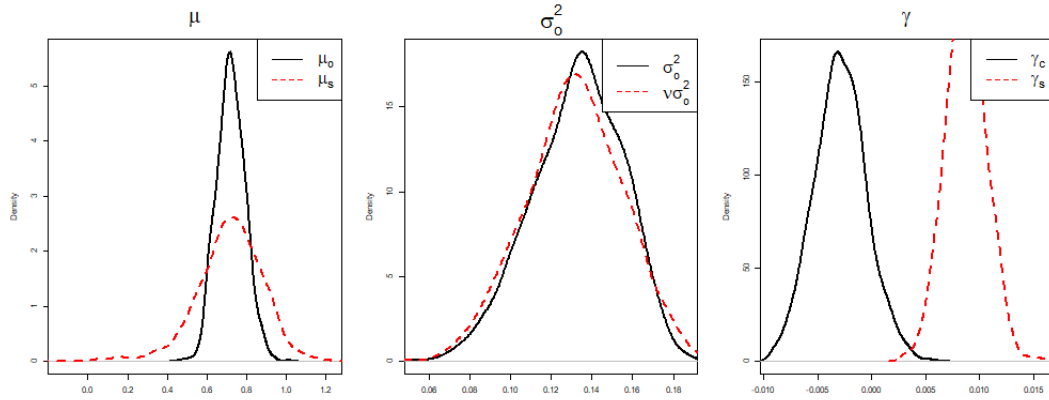


Figure 6: Same as Figure 5, but for the DJF season.

Table 2: Posterior expectations of bias, variance, and uncertainty of each simulator for the JJA season

Stage	Simulator	$\Delta\mu_{k,i}$	$\sigma_{k,i}^2$	$\tau_{k,i}$
ES	RCP 4.5	0.0358	5.5000	2.8207
	RCP 8.5	-0.1292	2.7917	1.4372
GCM	CESM1-BGC	0.6574	1.1452	0.3080
	HadGEM2-ES	0.7098	6.6361	1.7193
	MIROC-ESM-CHEM	0.6176	4.0201	1.0402
	MRI-CGCM3	-0.0782	0.6669	0.2542
BC	BCCA	0.9462	1.2556	0.6480
	BCSD	0.9441	2.1061	1.0825
HM	GR4J	-0.5993	0.8461	1.5742
	Tank	2.4136	14.5822	8.5870

ES = emission scenarios; GCM = global circulation model; BC = bias correction technique; HM = hydrological model.

Table 3: Posterior expectations of bias, variance, and uncertainty of each simulator for the DJF season

Stage	Simulator	$\Delta\mu_{k,i}$	$\sigma_{k,i}^2$	$\tau_{k,i}$
ES	RCP 4.5	0.0697	0.0889	0.0463
	RCP 8.5	-0.0644	0.0049	0.0048
GCM	CESM1-BGC	0.0591	0.0132	0.0037
	HadGEM2-ES	0.1579	0.0477	0.0129
	MIROC-ESM-CHEM	0.1054	0.0360	0.0090
	MRI-CGCM3	0.0619	0.0107	0.0031
BC	BCCA	0.1701	0.0072	0.0036
	BCSD	0.1679	0.0167	0.0084
HM	GR4J	0.2019	0.0111	0.0058
	Tank	0.1599	0.1726	0.0857

ES = emission scenarios; GCM = global circulation model; BC = bias correction technique; HM = hydrological model.

season. It implies that Tank hydrological model projects more extreme streamflow compared to the GR4J hydrological model.

4.4. Uncertainty decomposition

We compared the proposed Bayesian uncertainty decomposition method with the AVOVA approach (Yip *et al.*, 2011; Bosshard *et al.*, 2013) and the cumulative uncertainty approach (Kim *et al.*, 2019).

Table 4: Results of decomposition of the total uncertainty into the sources including NV, ESs, GCMs, BCs, and HMs for the JJA season

Source	ANOVA	Cumulative	Bayesian (whole)	Bayesian (scenario)
NV	-	-	2.5331 (11.51%)	2.2324 (9.63%)
ES	0.0150 (0.43%)	0.3027 (8.63%)	4.2579 (19.35%)	1.1619 (5.01%)
GCM	0.4751 (13.55%)	0.5276 (15.05%)	3.3217 (15.1%)	6.8689 (29.63%)
BC	0.0010 (0.03%)	0.0011 (0.03%)	1.7305 (7.86%)	1.2469 (5.38%)
HM	2.6745 (76.29%)	2.6745 (76.29%)	10.1611 (46.18%)	11.6702 (50.34%)
RSS	0.3403 (9.71%)	-	-	-

ANOVA denotes the ANOVA approach and Cumulative denotes the cumulative uncertainty approach, and Bayesian (whole) and Bayesian (scenario) denote the proposed Bayesian methods using the whole data and using only the data in the scenario period, respectively. RSS stands for residual sum of squares for the ANOVA approach. NV = natural variability; ES = emission scenarios; GCM = global circulation model; BC = bias correction technique; HM = hydrological model.

Table 5: Results of decomposition of the total uncertainty into the sources including NV, ESs, GCMs, BCs, and HMs for the DJF season

Source	ANOVA	Cumulative	Bayesian (whole)	Bayesian (scenario)
NV	-	-	0.1314 (41.75%)	0.0310 (9.61%)
ES	0.0066 (28.26%)	0.0158 (67.47%)	0.0511 (16.24%)	0.0859 (26.61%)
GCM	0.0058 (24.93%)	0.0073 (30.99%)	0.0287 (9.13%)	0.0528 (16.36%)
BC	4×10^{-6} (0.02%)	1.48×10^{-4} (0.63%)	0.0120 (3.81%)	0.0349 (10.82%)
HM	2.12×10^{-4} (0.91%)	2.12×10^{-4} (0.91%)	0.0915 (29.07%)	0.1181 (36.60%)
RSS	0.0108 (45.88%)	-	-	-

ANOVA denotes the ANOVA approach and Cumulative denotes the cumulative uncertainty approach, and Bayesian (whole) and Bayesian (scenario) denote the proposed Bayesian methods using whole data and only the data in the scenario period. RSS stands for residual sum of squares for the ANOVA approach. NV = natural variability; ES = emission scenarios; GCM = global circulation model; BC = bias correction technique; HM = hydrological model.

Both the ANOVA and cumulative uncertainty approach used only the average of all projected values over the scenario period.

Our Bayesian model can be applied when there is only data in the scenario period. In this case, we exclude the unnecessary parameters μ_o and ν . The total uncertainty and the uncertainty of stage k are then computed by using Equation (3.1) and Equation (3.2) with $\nu = 1$, respectively. The purpose of this analysis is to examine the impact of the data in the control period on uncertainty decomposition.

Tables 4 and 5 summarize the results of the uncertainty decomposition using the ANOVA method, the cumulative uncertainty method and the proposed Bayesian method using whole data and the Bayesian method using only the data in the scenario period for the JJA and DJF seasons. The percentage in the bracket indicates the proportion of the uncertainty of each stage that contributed to the total uncertainty. For the uncertainty measure in the cumulative uncertainty approach, we used the variance to quantify uncertainties for fair comparison. For the Bayesian model, the presented uncertainties are posterior expectation values and the proportions are computed as the ratio of each posterior expectation to the posterior expectation of the total uncertainty.

For the JJA season, the bias correction technique is the smallest contributor to the total uncertainty and the hydrological model is the largest contributor for all four methods. The uncertainty of GCMs is larger than that of emission scenarios for the three other methods except for the Bayesian method using whole data. For the DJF season, the hydrological model is the largest contributor among the four simulator stages for the two Bayesian methods but the emission scenario is the largest contributor for the other two methods. This is due to the large interannual variation of the Tank simulator, which

is quantified by Bayesian methods.

As mentioned in Section 1, unlike the other methods, the proposed Bayesian method can incorporate the uncertainty of natural variability. When we use the whole data in the control and scenario periods, the estimated proportion of uncertainty of natural variability is the fourth largest for the JJA season and the largest for the DJF season. However, the uncertainty of natural variability when using only the data in the scenario period is similar to the JJA season, but different for the DJF season. The inconsistency of the results is due to the absence of historical observations.

The uncertainty decomposition results between the Bayesian model using whole data and the model using only the data in the scenario period, are quite different. This implies that the information from the data in the control period depends significantly on the uncertainty decomposition results.

5. Concluding remarks

We proposed a Bayesian method for decomposing the total uncertainty into uncertainty of each source with the following advantages: (1) it can borrow the information from projected and observed data in the control period; and (2) it can quantify the uncertainty of natural variability. We also provide a simple and efficient Gibbs sampling algorithm using auxiliary variables.

Future research will apply the proposed Bayesian method to meteorological/hydrological variables that are not expected to follow a normal distribution, e.g., maximum precipitation which may be assumed to follow a generalized extreme value distribution as in Jo *et al.* (2016). Another potential extension of the proposed Bayesian method is to treat datasets for multiple regions within a single Bayesian model. This multivariate Bayesian model could ease the distributional assumption due to the plenty of the data. For example, the multivariate Bayesian model can introduce changes in parameters between the control and scenario periods or correlations between the uncertainty sources.

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