

Uncertainty component estimates in transient climate projections.

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▶ To cite this version:

Benoit Hingray, Juliette Blanchet, Guillaume Evin, Jean-Philippe Vidal. Uncertainty component estimates in transient climate projections.: Precision of estimators in a single time or time series approach. Climate Dynamics, 2019, 53 (5-6), pp.2501-2516. 10.1007/s00382-019-04635-1. hal-02327452

HAL Id: hal-02327452 https://hal.science/hal-02327452

Submitted on 26 Nov 2020

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- ¹ Uncertainty component estimates in transient ² climate projections
- ³ Precision of estimators in a single time or time series
- 4 approach
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8 Received: date / Accepted: date

Abstract Quantifying model uncertainty and internal variability components in climate projections has been paid a great attention in the recent years. For 10 multiple synthetic ensembles of climate projections, we compare the precision 11 of uncertainty component estimates obtained respectively with the two Analy-12 sis of Variance (ANOVA) approaches mostly used in recent works: the popular 13 Single Time approach (STANOVA), based on the data available for the con-14 sidered projection lead time and a time series based approach (QEANOVA), 15 which assumes quasi-ergodicity of climate outputs over the available simula-16 tion period. 17 We show that the precision of all uncertainty estimates is higher when more 18 members are used, when internal variability is smaller and/or the response-to-19 uncertainty ratio is higher. QEANOVA estimates are much more precise than 20 STANOVA ones: QEANOVA simulated confidence intervals are roughly 3 to 21

- ²² 5 times smaller than STANOVA ones. Except for STANOVA when less than
- $_{\rm 23}$ $\,$ 3 members is available, the precision is rather high for total uncertainty and
- ²⁴ moderate for internal variability estimates. For model uncertainty or response-
- ²⁵ to-uncertainty ratio estimates, the precision is low for QEANOVA to very low
- ²⁶ for STANOVA. In the most unfavorable configurations (small number of mem-

²⁷ bers, large internal variability), large over- or underestimation of uncertainty

²⁸ components is thus very likely. In a number of cases, the uncertainty analysis
²⁹ should thus be preferentially carried out with a time series approach or with a

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local-time series approach, applied to all predictions available in the temporal
 neighborhood of the target prediction lead time.

32 Keywords Uncertainty Sources · Climate Projections · ANOVA · Internal

³³ Variability · Model Uncertainty · Scenario Uncertainty · Precision of Estimates

34 1 Introduction

A critical issue in climate change studies is the estimation of uncertainties in projections along with the contribution of the different uncertainty sources, including scenario uncertainty, the different components of model uncertainty,

³⁸ and internal variability (e.g. Hawkins and Sutton, 2009).

Scenario uncertainty is related to the poorly known future of greenhouse 39 gas emissions. Model uncertainty corresponds to the dispersion between the 40 41 different climate responses obtained with different models for the same forcing configuration. Model uncertainty concerns Global Climate Models (GCMs) but 42 also all subsequent models of the climate change impact modeling chain, such 43 as Regional Downscaling Models (regional climate models and/or statistical 44 downscaling methods) and impact models like Hydrological Models (HMs). 45 Internal variability first originates from the chaotic variability of the climate 46 at the global scale (e.g. Räisänen, 2001; Deser et al, 2012). At the local scale, it 47 also results from the fact that very similar large scale atmospheric circulation 48 configurations can lead to very different meteorological observations (Braun 49 et al, 2012; Lafaysse et al, 2014). 50 Estimating and partitioning uncertainty in future climate projections is 51

first intended to help evaluating the significance of estimated changes for adaptation purposes. Besides, this is intended to highlight the most important uncertainty sources. This thus allows estimating the fraction of total uncertainty that could be narrowed via scenario refinement and model improvement, and its irreducible fraction pertaining to natural variability (e.g. Hawkins and Sutton, 2011; Lafaysse et al, 2014).

Over the recent years, uncertainty has been mostly explored and parti-58 tioned based on Multiscenarios Multimodel Multimember Ensembles (MMEs) 59 of transient climate projections. Various methods have been proposed for this, 60 most of them based on an Analysis of Variance (ANOVA) of projections avail-61 able for the specific projection lead time considered (Hingray et al, 2007; Yip 62 et al, 2011; Bosshard et al, 2013; Giuntoli et al, 2015; van Pelt et al, 2015; 63 Paeth et al, 2017). In this single time approach - called STANOVA in the 64 following -, and provided multiple members are available for each modeling 65 chain, the model uncertainty components are estimated from the dispersion 66 between the climate responses of the different modeling chains, obtained for 67 each chain from the multimember mean of the projections. Similarly, the in-68 ternal variability component for each modeling chain is estimated from the 69 inter-member variance of the projections. 70

The different uncertainty components can also be estimated from a time reserves approach, benefiting from the long time series now available for most

climate experiments (e.g. Johns et al, 2011; Jacob et al, 2014). A quasi-ergodic 73 assumption for climate simulations in transient climate is often used in this 74 context (Hingray and Saïd, 2014). It considers that the temporal variations 75 of the climate response of a particular simulation chain are necessarily grad-76 ual and smooth, the higher frequency variations of the time series being due 77 to internal variability alone. It considers next that the internal variability 78 can be assumed to remain roughly constant over the considered period or to 79 vary as a gradual and smooth function of the climate response of the chain. 80 These assumptions were used for instance by Räisänen (2001); Hawkins and 81 Sutton (2009, 2011); Charlton-Perez et al (2010); Lafaysse et al (2014); Brace-82 girdle et al (2014); Reintges et al (2017). In this time series approach called 83 QEANOVA in the following, the climate change response of each simulation 84 chain and its possible evolution with time is obtained from the long term trend 85 estimated from the time series of the chain. The variance over time of the de-86 viation from the climate response defines the internal variability of each chain. 87 The climate responses of the different chains can be used to estimate the com-88 ponents of model uncertainty for any projection lead time, thanks again to a 89 usual ANOVA. 90 In both approaches, a potentially critical step is the possibility to identify 91 a precise estimate of the climate response of each modeling chain. A number of 92 previous works advocated the use of multiple runs to improve this identification 93

(e.g. Kendon et al, 2008; Deser et al, 2012; Kew et al, 2011). Unfortunately 94 a limited number of members have until now been usually available for most 95 modeling chains, logically deriving from the small number of GCM runs (e.g. 96 Johns et al, 2011; Jacob et al, 2014; Paeth et al, 2017; Reintges et al, 2017). In 97 a number of analyses, the discrimination between internal climate variability 98 and the climate response of each modeling chain is thus expected to be quite 99 inaccurate owing to the limited ensemble size especially when the internal 100 variability is non-negligible compared to the chain's climate responses (e.g. 101 Deser et al, 2012; van Pelt et al, 2015). 102

These inaccuracies translate to inaccuracies in estimates of uncertainty components and of derived characteristics (e.g. response to uncertainty ratio, significance of projected changes, time of emergence of climate change, etc.). In specific analysis configurations, the precision of uncertainty component estimates is potentially low to very low, but to our knowledge this issue has never been considered.

In the following, we characterize the precision of uncertainty components 109 estimates when obtained either with the single time or the time series ap-110 proaches introduced previously. The precision is obviously expected to depend 111 on the multimodel multimember ensemble of projections (MME) available for 112 the analysis and on the climate variable considered for the analysis. In the 113 following, both ANOVA approaches are applied to multiple synthetic MMEs 114 obtained via Monte Carlo simulations. We consider a simplified configuration 115 where MME account for two uncertainty sources: model uncertainty and in-116 ternal variability. The analysis framework we follow here is rather crude but it 117 allows comparing both ANOVA approaches for a large number of MME config-118

urations with different characteristics. We especially discuss how the precision 119 depends on the number of members used in the analysis, and we characterize 120 the gain in precision obtained with the time series based approach, owing to 121 the larger size of the data set accounted for. We also discuss the results for 122 different levels of internal climate variability contribution to total uncertainty 123 variance. We first focus on the precision of estimates obtained for model un-124 certainty and internal variability. We next consider the precision of estimates 125 obtained on the one hand for the mean climate response from the model en-126 semble and on the other hand for the response to uncertainty ratio frequently 127 used to estimate the significance of estimated changes. 128

All recent analyses of uncertainty sources in climate projections make use of the moment estimators of variance components. The present study also makes use of such estimators in order to put the precision of uncertainty estimates derived from these analyses into perspective.

The statistical bases for the estimation of uncertainty components in a given MME are summarized in section 2 for both the single lead time ANOVA and the quasi-ergodic ANOVA approaches. The Monte Carlo experiment for the simulation of multiple synthetic MMEs is presented in section 3. The precision of uncertainty estimates is presented in section 4, and discussed in section 5. Section 6 concludes.

139 2 ANOVA methods

¹⁴⁰ 2.1 Transient ensembles of climate experiments

Let us consider a hypothetical ensemble of climate experiments where M dif-141 ferent members are available for each of G different climate modeling chains. 142 For the sake of simplicity, we assume that the same number of members is 143 available for each chain. A chain refers for instance to a given GCM and the 144 members to the different runs available for each GCM. A chain could also re-145 fer to a given GCM/RDM combination or to a given GCM/RDM/HM model 146 combination. Members would respectively refer to the potentially multiple re-147 alizations obtained with the different runs for each GCM/RDM combination 148 or for each GCM/RDM/HM combination (e.g. Lafaysse et al, 2014; Vidal et al, 149 2016). 150

Note Y(g, m, t) the climate projection obtained for the m^{th} member of chain g for time t. Y(g, m, t) can be expressed as:

$$Y(g,m,t) = \lambda(g,t) + \nu(g,m,t), \tag{1}$$

where $\lambda(g, t)$ is the climate response of model g for year t and $\nu(g, m, t)$ is the deviation from the climate response obtained with the member m for this year as a result of internal variability (IV).

As in most climate impact studies, the uncertainty analysis can be carried out for the change variable X. In the following, we consider absolute changes,

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¹⁵⁸ but a similar analysis could be obtained for relative changes. Absolute changes
 ¹⁵⁹ are expressed as:

$$X(g, m, t) = Y(g, m, t) - Y(g, m, t_C),$$
(2)

where X(g,m,t) are the experiment outputs of the change variable for a given

¹⁶¹ member m of each chain g, each year t and where t_C refers to a given reference

year (or *n*-yr reference period centered on year t_C). These change outputs can

¹⁶³ be written as:

$$X(g,m,t) = \varphi(g,t) + \eta(g,m,t), \tag{3}$$

where $\varphi(g, t)$ is the climate change response of chain g for t and where $\eta(g, m, t)$ are the residuals obtained for the m^{th} member of chain g as a result of IV.

¹⁶⁶ In the following, we assume that the climate change response can be further

¹⁶⁷ decomposed as:

$$\varphi(g,t) = \mu(t) + \alpha(g,t), \tag{4}$$

where $\mu(t)$ is the mean climate change response from all available chains at projection lead time t, $\alpha(g, t)$ is the deviation from $\mu(t)$ for the climate change response of chain g.

By assumption, no correlation is expected between the change responses 171 and the residuals $\eta's$. If the residuals are additionally independent and iden-172 tically distributed (i.i.d.), the total uncertainty of the change variable at t is 173 $\sigma_X^2(t) = s_\alpha^2(t) + \sigma_\eta^2(t)$ where $s_\alpha^2(t)$ is the sample variance of the α 's in equa-174 tion (4) and $\sigma_n^2(t)$ the variance of the $\eta's$ in equation (3). They respectively 175 correspond to the model uncertainty and IV components of total uncertainty 176 variance for X. Note that, for any given chain g, no correlation is expected 177 between the values of the residuals $\nu(q, m, t)$ obtained from the raw variable 178 Y in the reference and in a future climate. The IV of the change variable X179 is thus the sum of the IV estimates for Y obtained for the reference and the 180 future periods respectively. 181

In the following, we detail the two ANOVA approaches used to estimate (1)182 the mean climate change response $\mu(t)$, (2) the deviations $\alpha(q, t)$ for the differ-183 ent models, (3) the total uncertainty variance $\sigma_X^2(t)$, (4) its model uncertainty 184 $s_{\alpha}^{2}(t)$, and (5) its IV component $\sigma_{\eta}^{2}(t)$. We will also consider the Response-to-185 Uncertainty ratio R2U(t) defined as R2U(t) = $\mu(t)/\sigma_X(t)$, the fraction $F_{\alpha}(t)$ 186 of total variance explained by model uncertainty and the fraction $F_{\eta}(t)$ of to-187 tal variance explained by internal variability. Moment estimators of variance 188 components are used here for both ANOVA approaches. 189

¹⁹⁰ 2.2 Single Time ANOVA method

The Single Time ANOVA method (STANOVA in the following) is by far the most frequently used approach for the last years. It consists in estimating an ANOVA model on the projected changes X(g, m, t) for each time step t in turn. For any given t, climate projections data available for the time steps that precede and follow t are disregarded.

The STANOVA method uses an ANOVA model with fixed effects of the 196 form of equations (3) and (4). The residuals are assumed to be i.i.d. In this 197 configuration, an unbiased estimator of the residual's variance σ_{η}^2 at time t is 198 the multimodel mean of the inter-member variance at t, expressed as (Mont-199 gomery, 2012): 200

$$\hat{\sigma}_{\eta,ST}^{2}(t) = \frac{1}{G(M-1)} \sum_{g=1}^{G} \sum_{m=1}^{M} \left\{ X(g,m,t) - \hat{\varphi}_{ST}(g,t) \right\}^{2}$$
(5)

where $\hat{\varphi}_{ST}(g,t)$ is an estimate of the mean climate change response for model 201 202

g at *t* estimated over all *M* members as $\widehat{\varphi}_{ST}(g,t) = \frac{1}{M} \sum_{m=1}^{M} X(g,m,t)$. The model uncertainty variance is the inter-model sample variance of the deviations $\alpha(g,t)$ for all *g*, namely $s_{\alpha}^{2}(t) = \frac{1}{G-1} \sum_{g=1}^{G} \{\alpha(g,t)\}^{2}$. An unbiased estimator of $s_{\alpha}^{2}(t)$ under constraint $\sum_{g=1}^{G} \widehat{\alpha}_{ST}(g,t) = 0$ is (Montgomery, 2012): 203 204 205 206

$$\hat{s}_{\alpha,ST}^{2}(t) = \frac{1}{G-1} \sum_{g=1}^{G} \left\{ \hat{\alpha}_{ST}(g,t) \right\}^{2} - \frac{1}{M} \hat{\sigma}_{\eta,ST}^{2}(t)$$
(6)

where $\widehat{\alpha}_{ST}(g,t)$ is an estimate of the deviation $\alpha(g,t)$ for model g, namely 207 $\hat{\alpha}_{ST}(g,t) = \hat{\varphi}_{ST}(g,t) - \hat{\mu}_{ST}(t) \text{ with } \hat{\mu}_{ST}(t) \text{ the overall mean climate change response estimate at } t, \text{ namely } \hat{\mu}_{ST}(t) = \frac{1}{G} \sum_{g=1}^{G} \hat{\varphi}_{ST}(g,t).$ An unbiased estimator of the total variance for X(g,m,t) finally reads: 208 209

210 $\widehat{\sigma}_{X,ST}^2(t) = \widehat{s}_{\alpha,ST}^2(t) + \widehat{\sigma}_{n,ST}^2(t).$ 211

2.3 Quasi-Ergodic ANOVA method 212

A time series ANOVA approach consists first in estimating for each chain g the 213 climate response from the raw projections Y(q, m, t). This is typically achieved 214 by fitting a trend model to one of the members available for each chain (e.g. 215 Hawkins and Sutton, 2009; Reintges et al, 2017). In the Quasi-Ergodic ANOVA 216 method (QEANOVA in the following) considered here, the trend model is 217 not obtained from one given member but from all the M members Y(q, m, t)218 available for the chain (Hingray and Saïd, 2014). The trend model is assumed 219 to give an unbiased estimator of the climate response function $\lambda(q, t)$ of the 220 chain (the limitations linked to the choice of the trend model are discussed in 221 section 5.2). In this case, the trend model for Y gives an unbiased estimator 222 of the climate change response function of the chain $\varphi(q,t)$, as: 223

$$\widehat{\varphi}_{QE}(g,t) = \widehat{\lambda}_{QE}(g,t) - \widehat{\lambda}_{QE}(g,t_C) \tag{7}$$

where $\widehat{\lambda}_{QE}(g,t)$ and $\widehat{\lambda}_{QE}(g,t_C)$ are the trend estimates of the raw projections 224 Y for the future and reference periods respectively. 225

Under the quasi-ergodic assumption, the IV of chain q for the raw data Y is 226 estimated as the variance over time of the residuals from the climate response 227 estimated for the chain, namely $\{Y(g, m, t) - \lambda_{QE}(g, t)\}$. Note that all residuals 228

from all members are used here. If the experiment period covers T time steps, the IV of chain g is thus estimated from a sample of size $T \times M$. Assuming no correlation between residuals obtained for the reference and future periods and assuming that the IV variance for Y is constant over the entire experiment period, the IV of chain g for the change variable X is twice that of the raw variable Y (Hingray and Saïd, 2014).

An unbiased estimator of the IV component for X is derived from the multimodel mean of the IV estimates obtained for the G different chains respectively (Hingray and Blanchet, 2018, see also Appendix A, Eq.A.3 and Eq.A.6). When for each chain g, the trend model for Y can be expressed as a linear combination of L functions of time, it reads :

$$\widehat{\sigma}_{\eta,QE}^2 = \frac{2}{G(TM - L)} \sum_{g=1}^G \sum_{m=1}^M \sum_{k=1}^T \left\{ Y(g, m, t_k) - \widehat{\lambda}_{QE}(g, t_k) \right\}^2.$$
(8)

When the trend model is a simple linear function of time, L = 2.

An unbiased estimator of the sample model uncertainty variance $s_{\alpha}^{2}(t)$ under constraint $\sum_{g=1}^{G} \hat{\alpha}_{QE}(g,t) = 0$ is (Hingray and Blanchet, 2018, see also Appendix A, Eq.A.5):

$$\hat{s}_{\alpha,QE}^{2}(t) = \frac{1}{G-1} \sum_{g=1}^{G} \left\{ \hat{\alpha}_{QE}(g,t) \right\}^{2} - \frac{A(t,\mathcal{C})}{M} \hat{\sigma}_{\eta,QE}^{2}, \tag{9}$$

where $\hat{\alpha}_{QE}(g,t)$ is as previously an estimate of the deviation $\alpha(g,t)$ for model gexpressed as $\hat{\alpha}_{QE}(g,t) = \hat{\varphi}_{QE}(g,t) - \hat{\mu}_{QE}(t)$ with $\hat{\mu}_{QE}(t) = \frac{1}{G} \sum_{g=1}^{G} \hat{\varphi}_{QE}(g,t)$ the overall mean climate change response at t. A(t, C) depends on the projection lead time and on the different functions of time used for the trend model of Y (Hingray and Blanchet, 2018). It is positive and smaller than 1. The expression of A(t, C) is given in Appendix A when the trend model is a simple linear function of time (Eq. A.7).

An unbiased estimator of total variance for X(g, m, t) finally also reads: $\hat{\sigma}_{X,QE}^2(t) = \hat{s}_{\alpha,QE}^2(t) + \hat{\sigma}_{\eta,QE}^2$.

253 3 Simulations

 $_{254}$ In the following, we mainly focus on the precision of uncertainty estimators

obtained with STANOVA and QEANOVA methods for the change variable $X(g, m, t_e)$ obtained between some future prediction lead time t_e and the reference period t_C .

We construct via Monte Carlo simulations multiple synthetic MMEs where each MME is composed of $M \times G$ times series of raw projections Y(g, m, t), with M the number of members for each of the G chains (see Appendix C for details). For reasons of simplification, each MME is simulated based on the assumptions that the climate response function of each modeling chain g 8

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is a linear function of time and that the random deviations $\nu(g, m, t)$ due to internal variability come from independent and normally distributed random variables: $\nu(g, m, t) \sim \mathcal{N}(0, \sigma_{\nu}^2)$.

Each MME simulated for Y(g, m, t) allows deriving the corresponding time series for X(g, m, t). In the present analysis, for graphical simplification purposes, each MME for Y is constructed so that it leads, for X at time $t = t_e$, to a prescribed value of the Response-to-Uncertainty ratio [R2U(t_e)] and to a prescribed value of the fractional variance $F_{\eta}(t_e)$ due to internal variability (Appendix C).

In the following, simulations were produced over the 1960-2100 period for a 10-yr average climate variable (see illustrative MMEs in figure 1). Decadal time series were next aggregated to 20-yr time series for all analyses presented in section 4, in order to match with the usually preferred 20-yr temporal resolution in climate analyses. Y(g, m, t) time series are thus constituted from 7 values each (T = 140/20). The reference period is $t_C = 1990$.

For the sake of simplicity, a unique number of chains is considered in the following section (G = 5). The influence of the number of chains on the results is discussed in section 5.1. Most results are finally presented for prediction lead time $t_e = 2050$. t_e will be thus omitted in most following notations.

[FIGURE 1 HERE]

For each MME, both STANOVA and QEANOVA models are applied, allowing estimating the ratio $\mathbf{R}(\theta) = \hat{\theta}/\theta_{th}$ between the estimated value $\hat{\theta}$ and the theoretical (prescribed) value θ_{th} of different parameters θ : the grand ensemble mean response $\mu(t_E)$, the uncertainty variances $s_{\alpha}^2(t_E), \sigma_{\eta}^2(t_E), \sigma_X^2(t_E)$, the Response-to-Uncertainty ratio [R2U(t_E)] and the contribution to total variance of internal variability and model uncertainty, $F_{\eta}(t_E)$ and $F_{\alpha}(t_E)$.

We run the simulations for different combinations of prescribed R2U and F_{η} values and for different numbers of members M. For each combination, we simulate 1,000 MMEs. We subsequently estimate, for each parameter θ , the mean E[R(θ)] and the standard deviation SD[R(θ)] of the corresponding 1,000 ratios R(θ).

Application of a QEANOVA method requires fitting a trend model to the raw climate projections Y. A parametric model is usually used for convenience. For simplicity, we here also assume that the trend is a linear function of time for the whole transient period $[t_1, t_T]$ (the influence of the trend model on the results is discussed in section 5.2). The estimation is made using ordinary least squares.

Equations (5) and (8) give unbiased estimators of σ_{η}^2 . Respectively, equations (6) and (9) give unbiased estimators of s_{α}^2 . E[R(θ)] obtained for σ_{η}^2 and s_{α}^2 with both STANOVA and QEANOVA approaches are therefore expected to equal one whatever the ($M, F_{\eta}, R2U$) configuration. This actually also applies to all other STANOVA and QEANOVA estimators (not shown).

We therefore focus on the precision of estimators from STANOVA and QEANOVA approaches, i.e. on results obtained for $SD[R(\theta)]$, denoted SD below

307 for shortness. The smaller the standard deviation, the better the precision of

the estimator. We will additionally look at the ratio SD_{ST}/SD_{QE} , denoting the gain in precision obtained with the QEANOVA approach with respect to the STANOVA approach. The larger this ratio, the larger the gain.

The uncertainty analysis with the STANOVA method requires at least

two members for each chain. This is not the case for the QEANOVA method

which can be carried out with no more than one single member for each chain. In the following, we also present, for information, results obtained with the

³¹⁴ In the following, we also present, for information, result ³¹⁵ QEANOVA method in case of single member MMEs.

316 4 Results

In our case and whatever the $(M, F_{\eta}, R2U)$ configuration, the distribution of 317 the $R(\theta)$ ratios obtained respectively from the 1000 synthetic data sets was 318 always found to be roughly normal (not shown). In a first approximation, the 319 SD value of $R(\theta)$ thus determines the probability that the $R(\theta)$ ratio is outside 320 of any $[1 \pm e]$ interval, that is the probability that the estimated value of 321 the considered parameter is at least 100e% smaller or 100e% greater than the 322 theoretical value. For instance, the probability that the ratio $R(\theta)$ is outside the 323 [0.5, 1.5] interval is 1%, 32% or 62% when SD is 0.2, 0.5 or 1 respectively. When 324 the SD value is greater than 0.5, the probability for a large underestimation 325 (let say $R(\theta) < 0.5$) or for a large overestimation (let say $R(\theta) > 1.5$) is thus non

(let say $R(\theta) < 0.5$) or for a large overestimation (let say $R(\theta) > 1.5$) is thus non negligible (i.e. p > 0.32). When SD is greater than 1, a large underestimation

³²⁸ or overestimation of the considered parameter is very likely (i.e. p > 0.62).

329 4.1 Variance components, total and fractional variance

SD results obtained for variance parameters $(\sigma_{\eta}^2, s_{\alpha}^2, \sigma_X^2, F_{\eta}, F_{\alpha})$ are roughly independent on the R2U value prescribed for the simulation of the synthetic MMEs (not shown). They conversely depend on the value prescribed for F_{η} and on the number of members M. We here therefore only present results via the SD response functions in the (M, F_{η}) domain.

For the different variance parameters and both ANOVA methods, SD values are logically higher for small M numbers. Except for internal variability, they are also higher for large F_{η} values, highlighting the lower precision of the estimators in these configurations (figure 2). For all parameters, STANOVA SD values are also logically higher than QEANOVA SD values.

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[FIGURE 2 HERE]

For internal variability estimates, σ_{η}^2 , SD values do only depend on M. For configurations with $M \leq 3$, STANOVA SD values are greater than 0.4, highlighting the possibility for significant errors in σ_{η}^2 estimates. Whatever the (M, F_{η}) configuration, QEANOVA SD values are at least 3 times smaller than STANOVA SD values. QEANOVA estimates are thus expected to be much closer to the theoretical value.

For model uncertainty estimates s_{α}^2 , the precision gain obtained with QEANOVA 347 is at least 2.5. It is up to 4 for high F_{η} values (e.g. > 90%) or even more when 348 M is smaller than 5. Whatever the (M, F_{η}) configuration, SD values are much 349 larger than those obtained for internal variability. In the most unfavorable 350 configurations (very low M and very high F_{η} values), STANOVA SD values 351 are noticeably higher than 6! s_{α}^2 estimates are very likely to be largely over 352 or underestimated. In those cases, a misestimation can also be obtained with 353 QEANOVA but it would be somehow less critical. 354

³⁵⁵ Compared to s_{α}^2 and σ_{η}^2 , the potential error for total variance estimates, ³⁵⁶ σ_X^2 , is likely to be small for both methods. SD values are globally smaller than ³⁵⁷ 0.2 for QEANOVA. For STANOVA, they are smaller than 0.3 except in the low ³⁵⁸ M and high F_{η} configurations. This suggests that for any given dataset, errors ³⁵⁹ obtained for s_{α}^2 and σ_{η}^2 partly compensate. The gain in precision obtained with ³⁶⁰ QEANOVA is around 2.5, whatever the (M, F_{η}) configuration.

SD response functions obtained for fractional variances estimates, F_{η} and F_{α} , are very similar to those obtained for σ_{η}^2 and s_{α}^2 estimates respectively (not shown). Whatever the method, errors in F_{η} estimates are thus not expected to be very large. For F_{α} , large errors are conversely expected with STANOVA in the most unfavorable (M, F_{η}) configurations. The gain in precision obtained with QEANOVA is again at least 2.5.

³⁶⁷ 4.2 Mean response and Response-to-Uncertainty ratio

We focus here on the estimates obtained for the mean change response μ and for the Response-to-Uncertainty ratio R2U. For both ANOVA methods and both parameters, SD values are again logically higher for smaller M and/or higher F_{η} values. SD values additionally depend on the theoretical value prescribed for the Response-to-Uncertainty ratio R2U. SD values decrease when the theoretical value for R2U increases.

Results are presented in Figure 3 for μ estimates and three theoretical 374 R2U values (0.2, 1, and 5). They illustrate typical configurations encountered 375 in climate analyses. A R2U value as high as 5 can be for instance typically 376 obtained in the 2050s for large scale temperature variables (e.g. global mean 377 temperature as in Hawkins and Sutton (2011)). A R2U value as low as 0.2 378 can be conversely obtained for noisy systems, e.g. regional climate variables 379 with a high natural variability (e.g. 20-yr average regional precipitation as in 380 Hingray and Saïd (2014)). 381

For μ estimates, STANOVA SD values obtained when R2U = 0.2 are often greater than 0.5 and even exceed 2.5 for the most unfavorable (M, F_{η}) configurations. When R2U = 1, SD values are much smaller. They are lower than 0.15 except for the low M and high F_{η} configurations. For larger prescribed R2U values, SD values tend to be negligible: they are always smaller than 0.03 when R2U = 5. QEANOVA SD values present a similar dependence on M, F_{η} and R2U. They are however again much smaller than STANOVA SD values

(roughly 2.5 times smaller whatever the M, F_{η} and R2U configuration).

390

[FIGURE 3 HERE]

For R2U estimates, the structure of SD response functions are very similar to those obtained for μ estimates (not shown). SD values are however systematically larger (from +20% for R2U=0.1 to more than +100% for R2U=5). For noisy systems and the most unfavorable (M,F_{η}) configurations, large errors are therefore expected with STANOVA. Errors are conversely expected to be low to moderate with QEANOVA estimates, which are at least 2.5 times more

³⁹⁷ precise.

398 5 Discussion

³⁹⁹ 5.1 Dependence to the setup and size of the projection ensemble

For the sake of simplicity, we considered that MMEs are composed of 5 mod-400 elling chains. This is much less than the number of chains currently available 401 in recent CMIP MMEs. We also considered that the same number of members 402 is available for all simulations chains. This is obviously never the case in real 403 MMEs which are mostly unbalanced. In real MMEs, a large number of model 404 combinations is also typically missing when the projections are produced with 405 chains composed of different models (e.g. not all GCM/RCM combinations are 406 available when projections are produced with different RCMs from different 407 GCM outputs). 408 Whatever the ANOVA method, estimates of the different uncertainty com-409

⁴¹⁰ ponents are expected to be more precise when more members are available, ⁴¹¹ when the number of chains is larger and/or when the number of missing model ⁴¹² combinations is smaller. In most of the cases however, a time series approach ⁴¹³ is still expected to produce more precise estimates of uncertainty components ⁴¹⁴ than a single time approach. This obviously results from the larger sample size ⁴¹⁵ used for the estimation which allows for a more precise estimate of the climate ⁴¹⁶ response and of the internal variability of each simulation chain.

As an illustration, the previous analysis was repeated with MMEs composed of 10 simulation chains instead of 5 (see Supplementary Material, Figure A1). As expected, the precision of all uncertainty components estimates is better with 10 chains (for both approaches) and QEANOVA estimates are more precise than STANOVA ones. In the present case, the gains in precision are actually roughly the same than those presented previously with 5 chains.

⁴²³ 5.2 The trend estimation issue in the QEANOVA approach

⁴²⁴ In the present case and for convenience, time series of the synthetic MMEs are
⁴²⁵ simulated based on the assumption that the climate response of each chain is
⁴²⁶ a linear function of time. The QEANOVA trend function used to estimate the

427 climate response of each chain is also assumed to be a linear function of time.

In real datasets of projections, the climate response likely evolves in a 428 more complex way, perhaps differently in different chains. In practice, the 429 precise form of the climate response is not known, and for small ensembles and 430 noisy systems it may be quite difficult to guess and/or fit. If the chosen trend 431 model were incorrect for some chain, the response function estimate could be 432 locally biased. Incorrect trend models could then lead to significant errors in 433 the estimation of the local effects of the chains, and additional uncertainty 434 would arise, particularly if internal variability is small. Because of this trend 435 estimation issue, the results of a time series approach can thus be sub-optimal. 436 The linear assumptions retained for practical reasons in Section 3, make the 437 QEANOVA results likely too optimistic since the assumed trend model is the 438 true one. Thus the precision gain with QEANOVA in Section 4 is likely to 439 overestimate the gain to be obtained with this method for real MMEs. 440

For real MMEs, non-linear models can be considered for the estimation of the trend (e.g. 4rd order polynomial used in Hawkins and Sutton (2009) and in Hingray and Said (2014) for temperature changes) and the risk of choosing non relevant trend models can be reduced with dedicated statistical tests and a visual check of the fit (if any expert knowledge can be used for instance to assess the climatic relevance of the estimated trend).

The trend estimation issue in a time series ANOVA approach, which may 447 be critical for some MME configurations, requires thus specific care. Notwith-448 standing, a time series approach is likely to present a significant advantage 449 in many cases, even when the trend is not linear, especially when internal 450 variability is large. For illustration, we considered the same set of MMEs as 451 previously but, instead of fitting a linear trend for estimating the climate re-452 sponse of each simulation chain, we fitted a 3rd order polynomial function. 453 Results are presented in Figure A2 of Supplementary Material. As expected, 454 whatever the configuration in the (M, F_{η}) domain, the precision of QEANOVA 455 estimates is lower with the polynomial trend model, especially for noisy MMEs 456 and for small ensembles. Indeed, with a polynomial trend, the degrees of free-457 dom are larger and in case of noisy systems, the local trend is often dominated 458 by internal variability which logically leads to a lower precision of the esti-459 mated climate response. Nevertheless, QEANOVA is still much more precise 460 than STANOVA. Note that a more relevant experiment would have been to 461 simulate synthetic MMEs with non linear trends for the true trends but such 462 a simulation is unfortunately not convenient and would have required to make 463 different arbitrary and non representative choices (e.g. shape and amplitude 464 of non linearity to be retained for each chain). The experiment discussed here 465 thus likely again underplays the trend estimation issue in real MMEs where 466 the true trends are not linear. Fitting higher order polynomials when the true 467 trend is linear is indeed expected to introduce little error especially when 468 the sample is large (the fit essentially yields the linear trend), but, as men-469 tioned previously, the opposite does not hold: fitting a linear trend when the 470 true trend is non linear (e.g. quadratic or cubic) will always yield local biases 471

⁴⁷² which may be large.

Finally, another possible limitation of a time series ANOVA approach is 473 when internal variability is assumed to be constant over time. This assumption 474 may be not really valid as suggested by the recent work of Olonscheck and 475 Notz (2017). As presented by Hingray and Saïd (2014) or Bracegirdle et al 476 (2014), this assumption is not necessarily required and can also be relaxed. 477 The internal variability can for instance be assumed to vary gradually from 478 one lead time to the other (Hingray and Saïd, 2014). It can also be estimated 479 from the variance of the residuals within a moving temporal window centered 480 on the targeted prediction lead time (Bracegirdle et al, 2014). 481

482 5.3 Toward a local-QEANOVA approach

The potential limitations discussed above could make the apparent higher attractiveness of a time series approach somehow reduced for some MMEs configurations (e.g. MMEs with complex trend shapes and low noise).

However, if difficult to exploit over the whole available simulation period, 486 the QEANOVA principles can be at least used locally, in the temporal neigh-487 borhood $\Omega(t)$ of each time point t. In such a local QEANOVA approach, the 488 climate response of any given chain g can be obtained for instance from lo-489 cal linear regression, in which the trend is assumed to be smooth so that the 490 climate response $\lambda(q,t)$ of the chain can be approximated by a straight line 491 in the neighborhood of t (Cleveland, 1979). The residuals estimated from all 492 data in $\Omega(t)$ can next be used to assess the local internal variability of the 493 projections. When the uncertainty analysis is produced for the climate change 494 variable, the local linear regression and the estimation of the local internal 495 variability is done for both the reference time t_c and the future lead time t_e . 496 In this configuration, note that the internal variability no longer has to be 497 assumed constant between both periods. 498

The theoretical developments of this local QEANOVA approach, given in 499 Hingray and Blanchet (2018) are summarized in Appendix B. For illustration, 500 the local QEANOVA approach was also applied to the different simulation 501 ensembles considered in the previous sections. The neighborhood $\Omega(t)$ was here 502 chosen to be composed from the 3 consecutive time steps centered around t. In 503 this context, the risk that the linear trend is dominated by natural variability 504 is likely important for noisy systems and small ensembles. This configuration 505 would have led to a risk of underestimating the natural variability but this is 506 not the case here thanks to the unbiased estimator derived for this variable 507 (see Appendix B). 508

Results of the local QEANOVA are presented for internal variability, model uncertainty and for the Response-to-Uncertainty ratio in figure 4. Results are not a surprise. For all uncertainty parameters, the precision obtained with the local QEANOVA is significantly higher than that obtained with STANOVA (confidence intervals are 2 to 3 times smaller for the local QEANOVA). The precision of the local QEANOVA is conversely smaller than that obtained with ⁵¹⁵ QEANOVA. These results again logically derive from the different sizes of the ⁵¹⁶ samples used by the three methods.

A local QEANOVA is thus a promising alternative to STANOVA. The risk 517 for a large misestimation of uncertainty components is indeed much lower. 518 519 Despite its smaller precision, a local QEANOVA can be also a good alternative to a time series approach. When the time series of the simulation chain presents 520 a low to moderate noise but has a complex and nonlinear climate response 521 for instance, a local QEANOVA actually reduces the risk of having a biased 522 estimate of the climate response, which may occur in a time series approach 523 because of the choice of an inappropriate trend function (see section 5.2). 524

525 [FIGURE 4 HERE]

526 5.4 Temporal variations of uncertainty component estimates

Whatever the estimation method, the uncertainty estimates are typically derived for the different prediction lead times of a given climate simulation period. The temporal behavior of these estimates is worth considering as it may be a relevant indicator of their precision. In most cases, the climate response of each simulation chain is indeed expected to evolve gradually with time. In most cases, the different uncertainty components of a given MME are thus also expected to evolve gradually with time.

Similarly, uncertainty components for model uncertainty and scenario uncertainty should start from zero when the climate variable considered for the analysis is a climate change variable. The climate change response of any simulation chain is indeed expected to start from zero when the prediction lead time corresponds to the control period.

Large temporal fluctuations of uncertainty estimates and/or large values 539 of model uncertainty for the control period should warn for the possible lack 540 of robustness of the analysis. The possibility to consider the temporal varia-541 tions of uncertainty estimates to assess their precision is illustrated for three synthetic data sets in Figure 5 for s_{α}^2 and σ_{η}^2 and in Figure 6 for F_{η} . The 542 543 three MMEs have been simulated following the simulation process described 544 in section 3. Confidence intervals (colored bounds) of uncertainty estimates 545 (white lines), obtained from Monte Carlo simulations similar to those carried 546 out in section 4, highlight the much lower precision of STANOVA, even when 547 the fractional variance due to internal variability is low. This is also clear from 548 the temporal variations of uncertainty estimates. When only 3 members are 549 available, the variations of STANOVA estimates are erratic and non-negligible 550 whatever the value of F_{η} . They are much smoother with a local QEANOVA 551 and they completely vanish with QEANOVA. As expected also, the model 552 uncertainty component estimated with QEANOVA starts from zero at the be-553 ginning of the century and then gradually increases. This is thus fully coherent 554 with the expected behavior of this uncertainty component (limit to zero for 555 the control period). This is not the case for the two other methods, especially 556

for STANOVA in the case of small ensembles and noisy systems (e.g. significant non zero values of model uncertainty can be achieved at time t = 2000in Figure 5, STANOVA, 2nd line).

560 [FIGURE 5 HERE]

561 [FIGURE 6 HERE]

562 6 Conclusion

563 6.1 Precision of different methods for partitioning uncertainty

Two ANOVA approaches are currently used for partitioning model uncertainty and internal variability components in ensembles of climate projections: the popular Single Time approach (STANOVA), based on the data available for the considered projection lead time and time series approaches such as the QEANOVA approach considered here. For multiple synthetic ensembles, we compared the precision of uncertainty component estimates obtained with both approaches. The main findings are as follows:

In both cases, the precision is logically lower for noisy MMEs and/or small
 ensembles of members. The precision of the mean climate response and of
 the Response-to-Uncertainty ratio estimates are additionally lower when
 the theoretical (prescribed) Response-to-Uncertainty ratio is lower.

In many configurations, QEANOVA estimates are expected to be more precise than STANOVA ones. This does not depend on the setup and size of the MME considered (unbalanced MMEs, MMEs with missing chains, MMEs with more GCMs,...). The larger sample size used for the estimation is indeed expected to produce a better estimate of the climate response and of the internal variability of each simulation chain.

 Whatever the estimation method, the temporal variations of the uncertainty estimates should be considered as a relevant indicator of the precision of estimates. Large temporal fluctuations should call for caution and should warn for the likely low relevance of the uncertainty analysis.

⁵⁸⁵ Under the specific setup and assumptions of our analysis framework (e.g. ⁵⁸⁶ length of time series, number of simulations chains, linearity of trend, con-⁵⁸⁷ stancy of internal variability over time...), we find that:

• Simulated confidence intervals obtained with QEANOVA are at least 2.5 times smaller than those obtained with STANOVA.

• The precision of estimates is rather high for total uncertainty variance and for the multichain mean response (except for STANOVA for noisy MMEs and/or for MMEs with small ensembles). The two main features of any given ensemble of projections are thus likely to be rather correctly estimated in most cases. • This is obviously not the case for the individual uncertainty components. For model uncertainty variance for instance, the precision of estimates is low to very low for small ensembles and/or noisy MMEs. In the latter configurations, a large over- or underestimation of uncertainty components is thus very likely preventing any precise partition of uncertainty sources. This problem is particularly critical for STANOVA.

• STANOVA actually appears to be inappropriate in a large number of configurations as a result of the limited number of members, e.g. GCM runs, usually available for most modeling chains up to now. In the case of small ensembles and noisy MMEs, improving the relevance of uncertainty estimates requires using more data. A time series analysis approach as the QEANOVA method offers such an opportunity.

For a time series approach, the trend estimation can be an issue. For MMEs 607 where internal variability is small for instance, there is some risk to misrep-608 resent the forced trend. In such a case, a local QEANOVA approach offers 609 an interesting alternative as it is not subject to the same limitation as a full 610 time series approach. The precision of a local QEANOVA is additionally al-611 ways better than that of STANOVA. Note also that, similarly to a time series 612 approach, a local QEANOVA can be also applied when only a single member 613 is available for the chains. It can therefore make use of all climate experiments 614 produced in the recent years. This is another advantage over STANOVA which 615 requires multiple members of each chain. 616

617 6.2 Precision of uncertainty estimates for real ensembles of projections

Most ensembles of projections have potentially more uncertainty sources than 618 those considered in the present work. They may actually include scenario un-619 certainty, different components of model uncertainty (e.g. associated respec-620 tively to GCMs, Hydrological Models), of internal variability (with potentially 621 its large or small scale components) and, eventually, uncertainty associated to 622 interactions between models and/or scenarios (e.g. Hawkins and Sutton, 2011; 623 Vidal et al, 2016; Paeth et al, 2017). In all cases, and whatever the ANOVA 624 method used for the analysis, an idea of the precision of uncertainty compo-625 nents should be provided along with estimates of uncertainty components. 626

Results presented in the present work give already some benchmark for 627 simple ensembles of projections. They were obtained with moment estimators 628 of variance components. Likelihood-based estimation methods such as max-629 imum likelihood, restricted maximum likelihood or Bayesian methods could 630 (should) be used instead. They indeed allow for the estimation of confidence 631 or credibility intervals associated to each uncertainty component. They also 632 present next a number of interesting properties (e.g. Hartley and Rao, 1967; 633 Gelman, 2005). Conversely to method-of-moment based ANOVA approaches, 634 they can especially not produce negative estimates of variance components. 635 For the two ANOVA methods considered here, negative estimates are indeed 636 obtained for model uncertainty variance when the second terms in the right 637

hand sides of equations (6) or (9) are greater than the first terms. The probability of having a negative estimate for s_{α}^2 is actually directly related to the precision that can be achieved with the method for this component. For our work, this probability is roughly 2.5%, 16% or 31% when the SD value obtained for s_{α}^2 is 0.5, 1 or 2 respectively. It can be up to 50% in the most critical ($M_{\alpha}E_{\alpha}$ = 0.21) configurations

⁶⁴³ critical $(M, F_{\eta}, \text{R2U})$ configurations.

Likelihood-based estimation methods are also expected to use the data 644 more efficiently than moment-based estimation methods. They are for instance 645 easily applicable for unbalanced MMEs and/or MMEs with missing chains. 646 The work of Northrop and Chandler (2014) and of Evin et al (2018) give some 647 nice illustrations of their uses in a Bayesian context. The use of likelihood-648 based estimation methods is thus to be preferred to the use of moment-based 649 estimation methods. Even with such methods however, our results strongly 650 suggest that a full- or partial time series approach should be used instead of 651 a standard single time approach. 652

Author contributions. BH designed the analysis, developed the local QEANOVA and the synthetic simulations. JB derived the theoretical expressions for unbiased estimators of uncertainty components and wrote the ap-

pendixes. All authors contributed to write the manuscript and discuss results.

657 Acknowledgements We thank the three anonymous reviewers for their constructive sug-658 gestions which helped to significantly improve the content of our manuscript.

659 Appendix A QEANOVA Estimates of uncertainty components

We summarize the theoretical developments of the QEANOVA approach to 660 achieve unbiased estimators of uncertainty components for a simplified con-661 figuration where the trend model is a simple linear function of time. The full 662 developments are given in Hingray and Blanchet (2018) for the general config-663 uration where the trend model is a linear combination of L functions of time. 664 Here, we first consider the case where the number of members differs from one 665 chain to the other; the simplified equations obtained when all chains have the 666 same number are then given. For the sake of conciseness, we omit the subscript 667 "QE" related to the QEANOVA approach. 668

Model. We first consider the raw projections Y(g, m, t) with M_g members for each of the G chains, assuming that, for all $t_s \leq t \leq t_f$:

$$Y(g,m,t) = \lambda(g,t) + \nu(g,m,t), \tag{A.1}$$

where $\lambda(g,t)$ is the trend model expressed as a linear function of time: $\lambda(g,t) = \Lambda_{g1} + \Lambda_{g2}(t-t_s)$ and where the $\nu(g,m,t)$ are independent and homoscedastic random variables (with variance $\sigma_{\nu_g}^2$). Second let consider the change variable at future prediction lead time $t \in [t_c, t_f]$:

$$X(g,m,t) = Y(g,m,t) - Y(g,m,t_c) = \alpha(g,t) + \eta(g,m,t),$$
(A.2)

where t_c is the reference period. We have thus $\alpha(g,t) = \Lambda_{g2}(t-t_c)$ and $\eta(g,m,t) = \nu(g,m,t) - \nu(g,m,t_c).$

Unbiased estimation of the model parameters for the raw variable Y. We discretize $[t_s, t_f]$ into T time steps (from $t_s = t_1$ to $t_f = t_T$) and write t_c as the Kth time step (i.e. $t_c = t_K$). We are interested in the future prediction lead time $t_k \in [t_K, t_T]$. Let consider the regression model (A.1) for a particular g. Unbiased estimators of the regression parameters $(\Lambda_{g1}, \Lambda_{g2})$ are given by the least square estimates

$$(\hat{A}_{g1}, \hat{A}_{g2})' = \mathbb{V} \mathbb{R}' \left(\frac{1}{M_g} \sum_{m=1}^{M_g} Y(g, m, t_1), \dots, \frac{1}{M_g} \sum_{m=1}^{M_g} Y(g, m, t_T)\right)'$$

where "" denotes the transpose, \mathbb{R} is the $T \times 2$ matrix of covariates whose *k*th row is $(1, t_k - t_1)$, for $1 \leq k \leq T$, and $\mathbb{V} = (\mathbb{R}'\mathbb{R})^{-1}$. Covariance matrix of the estimators $(\hat{A}_{g1}, \hat{A}_{g2})$ is given by $\widehat{\sigma_{\nu_g}^2} M_g^{-1} \mathbb{V}$ where $\widehat{\sigma_{\nu_g}^2}$ is an unbiased estimator of $\sigma_{\nu_g}^2$ given by

$$\widehat{\sigma_{\nu_g}^2} = \frac{1}{TM_g - L} \sum_{m=1}^{M_g} \sum_{k=1}^T \left\{ Y(g, m, t_k) - \hat{\lambda}_{QE}(g, t_k) \right\}^2$$
(A.3)

681 where L = 2 and $\hat{\lambda}_{QE}(g, t_k) = \hat{\Lambda}_{g1} - \hat{\Lambda}_{g2}(t_k - t_1)$.

In particular, an unbiased estimator of Λ_{g2}^2 is

$$\widehat{\Lambda_{g2}^2} = \widehat{\Lambda_{g2}^2} - \widehat{\sigma_{\nu_g}^2} M_g^{-1} V_{22}, \qquad (A.4)$$

where V_{22} is the element (2,2) of \mathbb{V} . Considering t_1, \ldots, t_T regularly spaced on $[t_s; t_f]$, we have $V_{22} = 12(T-1)/\{T(T+1)(t_T-t_1)^2\}$.

⁶⁶⁵ Unbiased estimation of the sample variance of the α 's in the change vari-⁶⁶⁶ able. Given (A.2) and (A.4), an unbiased estimator of the sample variance of ⁶⁶⁷ $\alpha(g, t_k)$, i.e. of $s_{\alpha}^2(t_k) = \frac{1}{G-1} \sum_{g=1}^G \{\alpha(g, t_k)\}^2$, is

$$\widehat{s_{\alpha}^{2}}(t_{k}) = s_{\hat{\alpha}}^{2}(t_{k}) - \frac{12}{T} \frac{T-1}{T+1} \left(\frac{t_{k}-t_{K}}{t_{T}-t_{1}}\right)^{2} \left(\frac{1}{G} \sum_{g=1}^{G} \frac{\widehat{\sigma_{\nu_{g}}^{2}}}{M_{g}}\right),$$

where

$$s_{\hat{\alpha}}^2(t_k) = \frac{(t_k - t_1)^2}{G - 1} \sum_{g=1}^G \hat{\Lambda}_{g2}^2.$$

When all GCMs have the same number of runs (M), this expression reduces to:

$$\widehat{s}_{\alpha}^{2}(t_{k}) = s_{\hat{\alpha}}^{2}(t_{k}) - \frac{A(t_{k}, \mathcal{C})}{M} \widehat{\sigma}_{\eta}^{2}, \qquad (A.5)$$

where $\hat{\sigma}_{\eta}^2$ is an unbiased estimator of internal variability variance for X

$$\widehat{\sigma}_{\eta}^{2} = \frac{2}{G} \sum_{g=1}^{G} \widehat{\sigma_{\nu_{g}}^{2}}$$
(A.6)

and where

$$A(t_k, C) = \frac{6(T-1)}{T(T+1)} \left(\frac{t_k - t_K}{t_T - t_1}\right)^2$$
(A.7)

692

⁶⁹³ Appendix B Estimates with a local QEANOVA approach

We here summarize the expressions of the different uncertainty estimators obtained with a local-QEANOVA approach. The full developments, similar to those presented in appendix A for the QEANOVA approach, are detailed in Hingray and Blanchet (2018).

⁶⁹⁸ Model. A regression model is still considered to estimate the response function ⁶⁹⁹ $\lambda(g,t)$ for Y in Eq.A.1 but $\lambda(g,t)$ is assumed to be only locally linear in time, ⁷⁰⁰ in the neighborhoods of t_c and t_e respectively, i.e. on $[t_c - \omega, t_c + \omega]$ and ⁷⁰¹ $[t_e - \omega, t_e + \omega]$, where $t_e \in [t_c, t_f]$ is the future prediction lead time under ⁷⁰² consideration. $\lambda(g,t)$ can thus be expressed as

$$\lambda(g,t) = \begin{cases} \lambda_c(g,t) = \Lambda_{g1,c} + (t-t_c)\Lambda_{g2,c} \text{ for } t_c - \omega \le t \le t_c + \omega, \\ \lambda_e(g,t) = \Lambda_{g1,e} + (t-t_e)\Lambda_{g2,e} \text{ for } t_e - \omega \le t \le t_e + \omega. \end{cases}$$
(B.1)

The change variable X(g, m, t), for $t_e - \omega \le t \le t_e + \omega$, in Eq.A.2 is such that $\alpha(g, t) = (\Lambda_{g1,e} - \Lambda_{g1,c}) + (t - t_e)\Lambda_{g2,e}$. Each interval $[t_c - \omega, t_c + \omega]$ and $[t_e - \omega, t_e + \omega]$ is discretized into T^* regular

Each interval $[t_c - \omega, t_c + \omega]$ and $[t_e - \omega, t_e + \omega]$ is discretized into T^* regular periods of length $dt = 2\omega/(T^* - 1)$, with T^* odd, giving respectively the sequences t_1, \ldots, t_{T^*} and $t_{T^*+1}, \ldots, t_{2T^*}$. The values of Y for these different times are further considered to estimate the regression coefficients of the linear trend models in Eq.B.1. For the illustration given in section 5.3, $dt = \omega =$ 20yrs and $T^* = 3$.

711 Unbiased estimators of model uncertainty and internal variability variance.

- ⁷¹² Following Hingray and Blanchet (2018), an unbiased estimator of model un-
- ⁷¹³ certainty variance, i.e. of the sample variance of $\alpha(g, t_e)$, is

$$\widehat{s_{\alpha}^2}(t_e) = s_{\hat{\alpha}}^2(t_e) - \frac{4}{T} \left(\frac{1}{G} \sum_{g=1}^G \frac{\widehat{\sigma_{\nu_g}^2}}{M_g} \right)$$
(B.2)

where $T = 2T^*$ is the total number of time steps considered in the analysis and where $\widehat{\sigma_{\nu_g}^2}$ is an unbiased estimator of $\sigma_{\nu_g}^2$ given by

$$\widehat{\sigma_{\nu_g}^2} = \frac{1}{TM_g - 4} \sum_{m=1}^{M_g} \left[\sum_{j=1}^{T^*} \left\{ Y(g, m, t_j) - \hat{\lambda}_c(g, t_j) \right\}^2 \right]$$
(B.3)

$$+\sum_{j=T^*+1}^{2T^*} \left\{ Y(g,m,t_j) - \hat{\lambda}_e(g,t_j) \right\}^2 \right]$$
(B.4)

with $\hat{\lambda}_c(g,t_j) = \hat{A}_{g1,c} + \hat{A}_{g2,c}(t_j - t_c)$ and $\hat{\lambda}_e(g,t_j) = \hat{A}_{g1,e} + \hat{A}_{g2,e}(t_j - t_e)$ where $\hat{A}_{g1,c}, \hat{A}_{g2,c}, \hat{A}_{g1,e}$ and $\hat{A}_{g2,e}$ are the regression coefficients of the two linear models in Eq.B.1.

⁷¹⁹ When all GCMs have the same number of runs (M), the expression of ⁷²⁰ model uncertainty in Eq.B.2 reduces to

$$\widehat{s_{\alpha}^2}(t_e) = s_{\widehat{\alpha}}^2(t_e) - \frac{1}{MT^*} \widehat{\sigma}_{\eta}^2$$
(B.5)

where $\hat{\sigma}_{\eta}^2$ is an unbiased estimator of internal variability variance for X

$$\widehat{\sigma}_{\eta}^{2} = \frac{2}{G} \sum_{g=1}^{G} \widehat{\sigma_{\nu_{g}}^{2}}.$$
(B.6)

722 Appendix C Simulation of MMEs

Each MME is simulated for Y(g, m, t) assuming that, for all $t_s \leq t \leq t_f$:

$$Y(g,m,t) = \lambda(g,t) + \nu(g,m,t), \tag{C.1}$$

where the climate response function $\lambda(g,t)$ is a linear function of time and where the $\nu(g,m,t)$ are independent and homoscedastic random variables (with variance $\sigma_{\nu_g}^2$). For convenience we further assume that $\lambda(g,t)$ can be decomposed as $\lambda(g,t) = w(t) + d(g,t)$, for $t = t_1, \ldots, T$ where the mean climate response w(t) of the *G* chains and the deviations d(g,t) of chain *g* are linear functions of time, expressed as: $w(t) = B + P.(t - t_1)/(t_e - t_C)$ and $d(g,t) = D(g).(t - t_1)/(t_e - t_C)$ with the constraint $\sum_{g=1}^{G} D(g) = 0$.

For graphical simplification purposes, each MME for Y is constructed so that the parameters $(\sigma_{\nu}^2, P, D(g), g = 1, ..., G)$ lead, for the change variable X at time $t = t_e$, to a prescribed value of the Response-to-Uncertainty ratio [R2U(t_e)] and to a prescribed value of the fractional variance $F_{\eta}(t_e)$ due to internal variability.

For X, we have by definition $\varphi(g,t) = \lambda(g,t) - \lambda(g,t_C)$ and $\eta(g,m,t) = \nu(g,m,t) - \nu(g,m,t_C)$. We have thus for the change variable $\mu(t) = w(t) - w(t_C)$, $\alpha(g,t) = d(g,t) - d(g,t_C)$ and in turn $\mu(t) = P(t - t_C)/(t_e - t_C)$ and $\alpha(g,t) = D(g)(t - t_C)/(t_e - t_C)$.

The theoretical values for $\mu(t_e)$ and $s_{\alpha}^2(t_e)$ are thus as follows : $\mu(t_e) = P$ and $s_{\alpha}^2(t_e) = \mathbb{V}ar(D(g))$. Fixing P to 1, we thus simply require in turn

$$\sigma_X^2(t_e) = \frac{1}{[R2U(t_e)]^2};$$
 (C.2)

$$\sigma_{\nu}^{2}(t_{e}) = \frac{1}{2}\sigma_{\eta}^{2}(t_{e}) = \frac{1}{2}F_{\eta}(t_{e}).\sigma_{X}^{2}(t_{e}); \qquad (C.3)$$

$$\mathbb{V}ar(D(g)) = s_{\alpha}^{2}(t_{e}) = (1 - F_{\eta}(t_{e})) \cdot \sigma_{X}^{2}(t_{e}).$$
(C.4)

For each MME simulation, the deviations of the different chains, $D(g), g = 1, \ldots, G$, are obtained from a sample of G realizations in a normal distribution. These realizations are scaled so that their mean is zero and their variance corresponds to the prescribed value $s_{\alpha}^{2}(t_{e})$.

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Fig. 1 Three synthetic multimodel multimember ensembles of time series projections (MMEs) obtained via simulation with different F_{η} ratios and M numbers (left: $F_{\eta} = 0.1$, M = 3; middle $F_{\eta} = 0.5$, M = 3; right: $F_{\eta} = 0.5$, M = 9). Time series over the 1960-2100 period for the raw projections Y(g, m, t) of a 10-yr average climate variable. Cases on the left and in the middle are illustrative of typical configurations encountered in climate analyses: "small" number of members available for each GCM (M=3) with relatively low (resp. high) contribution of IV to total variance (case 1: $F_{\eta} = 0.1$; case 2: $F_{\eta} = 0.5$). The case on the right is similar as the middle case but in a configuration with a (unrealistic) "large" number of members (M = 9).



Fig. 2 Precision of QEANOVA (left) and STANOVA (middle) estimates for uncertainty variance components, and gain in precision between STANOVA SD values and QEANOVA SD values (right)(gain = ratio between STANOVA SD values and QEANOVA SD values). Top: Internal Variability, middle: Model Uncertainty, bottom: Total Uncertainty. SD values are given as a function of the fraction of total variance explained by internal variability (F_{η}) for a few representative values of number of members M: (1), 2, 3, 5, 10, 20. Results are presented for a theoretical ratio R2U = 1. For the sake of clarity, the upper limit of the figures for s_{α}^2 is truncated to 2. The highest values are greater than 10. Figures of the ratios between SD values obtained with both approaches are truncated to 5.



Fig. 3 Precision of QEANOVA (left) and STANOVA (middle) estimates for mean change response estimates (μ). Gain in precision between STANOVA and QEANOVA in the right panels. Results are presented for different values of the theoretical response-to-uncertainty ratio R2U (top: R2U = 0.2, middle: R2U = 1, bottom: R2U = 5). For details see caption of Figure 2.



Fig. 4 Precision of uncertainty estimates obtained with the local QEANOVA approach. Results for internal variability variance $(\sigma_{\eta}^2, \text{top})$, model uncertainty variance $(s_{\alpha}^2, \text{middle})$, response-to-uncertainty ratio (R2U, bottom). Left: SD values, middle: gain in precision between local QEANOVA and QEANOVA approaches (ratio between local QEANOVA SD values), right: gain in precision between STANOVA and local QEANOVA SD values), right: gain in precision between STANOVA and local QEANOVA SD values). Results for a theoretical ratio R2U = 1. For details see caption of Figure 2.



Fig. 5 Temporal variations and 90% confidence intervals of model uncertainty and internal variability estimates for the three reference datasets presented in Fig. 1 (top: $F_{\eta} = 0.1$, M = 3; middle $F_{\eta} = 0.5$, M = 3; bottom: $F_{\eta} = 0.5$, M = 9). Left: with STANOVA; middle: with local QEANOVA where $\Omega(t)$ is composed of the 3 consecutive time steps centered around t; right: with QEANOVA. s_{α}^2 in red, σ_{η}^2 in blue. The blue and red lines correspond to the theoretical variance components used for the generation of the reference dataset. The variance component estimates (white curves) obtained with each ANOVA approach are presented with their 90% confidence intervals.



Fig. 6 Temporal variations and 90% confidence intervals of uncertainty estimates for the fractional variance due to internal variability (F_{η} ratio) for the three reference datasets presented in Fig. 1. Left: with STANOVA; middle: with local QEANOVA; right: with QEANOVA. The blue lines correspond to the theoretical fractional variance used for the generation of the reference dataset. See Fig. 5 for caption details.