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A Bayesian Hierarchical Approach To Regional Frequency Analysis

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X - 2 RENARD: HIERARCHICAL REGIONAL ESTIMATION Abstract. Regional Frequency Analysis (RFA) has a long history in Hy-3 drology, and numerous distinct approaches have been proposed over the year 4 to perform the estimation of some hydrologic quantity at a regional level. 5 However, most of these approaches still rely on strong hypotheses that may limit their application and complicate the quantification of predictive un-7 certainty. The objective of this paper is to propose a general Bayesian hirarchical framework to implement RFA schemes that avoid these difficul-9 ties. The proposed framework is based on a two-level hierarchical model. The 10 first level of the hierarchy describes the joint distribution of observations. An 11 arbitrary marginal distribution, whose parameters may vary in space, is as-12 sumed for at-site series. The joint distribution is then derived by means of 13 an elliptical copula, therefore providing an explicit description of the spa-14 tial dependence between data. The second level of the hierarchy describes 15 the spatial variability of parameters using a regression model that links the 16 parameter values with covariates describing site characteristics. Regression 17 errors are modeled with a Gaussian spatial field which may exhibit spatial 18 dependence. This framework enables performing prediction at both gauged 19 and ungauged sites and, importantly, rigorously quantifying the associated 20 predictive uncertainty. A case study based on annual maxima of daily rain-21 fall demonstrates the applicability of this hierarchical approach. Although 22 numerous avenues for improvement can already be identified (amongst which 23 the inclusion of temporal covariates to model time variability), the proposed 24

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- $_{^{25}}\,$ model constitutes a general framework for implementing flexible RFA schemes
- ²⁶ and quantifying the associated predictive uncertainty.

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1. Introduction

1.1. Standard implementations of regional frequency analysis

The purpose of Regional Frequency Analysis (RFA) is to estimate the distribution of some hydrologic variable (e.g. annual maximum rainfall or runoff) using data from several sites. Compared with standard at-site frequency analysis, RFA attempts to improve the precision of estimates by sharing the information stemming from similar sites. Moreover, RFA enables estimation at ungauged or poorly gauged sites by transferring the information arising from neighboring gauging stations.

Numerous approaches have been proposed to implement RFA schemes. Amongst them, 33 the index flood method proposed by *Dalrymple* [1960] is still widely used in engineering 34 practice. It is based on a scale invariance hypothesis: within an homogeneous region, 35 distributions from all sites are assumed identical, up to a scale factor, termed the index 36 flood. The implementation of the index flood method can be summarized in three steps 37 [e.g., Hosking and Wallis, 1997]: (i) delineation of an homogeneous region; (ii) estimation 38 of the common regional distribution, based on scaled at-site data (i.e. divided by the index 39 flood); (iii) transfer of information to ungauged or poorly gauged site using a regression 40 model linking the index flood values with site characteristics. 41

⁴² The index flood method is widely used due to its ease of implementation and its robustness.
⁴³ However, its basic implementation is affected by several limitations:

• The delineation of homogeneous regions, where the scale invariance assumption holds, is far from obvious.

• The scale invariance assumption might simply be too restrictive is some cases.

In most cases, the index flood is defined as the mean or the median of at-site data,
but the physical reasons behind this choice are unclear.

• The regional distribution is estimated by pooling scaled at-site data, and treating them as is they were independent, which is rarely the case.

• Standard regression methods like ordinary least squares might be statistically inefficient because index flood values are statistics (as opposed to observations), and are therefore affected by estimation errors that may be dependent in space and whose properties may vary from site to site.

• The previous points make the quantification of the total predictive uncertainty challenging.

A wealth of research has been carried out over the years, either to improve the implementation of the index flood method, or to generalize it by abandoning some of its most restrictive assumptions (in particular, scale invariance). A non-exhaustive list of examples includes the work by *Ouarda et al.* [2001] on the concept of homogeneous region, the studies by *Stedinger* [1983] and *Hosking and Wallis* [1988] on estimating the regional distribution with spatially dependent data, or the extension to peak-over-threshold series of *Madsen and Rosbjerg* [1997] and *Ribatet et al.* [2007].

The transfer of information from gauged to ungauged site and the quantification of the associated predictive uncertainty has been a topic of particular attention [e.g., *Stedinger and Tasker*, 1985, 1986; *Reis et al.*, 2005; *Kjeldsen and Jones*, 2009a; *Micevski and Kuczera*, 2009]. Indeed, this transfer relies on a regression model linking the index flood values (estimated at gauged sites) and site characteristics. *Stedinger and Tasker* [1985, 1986] introduced the Generalized Least Squares (GLS) approach to account for both the het-

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⁷⁰ eroscedastic and spatially dependent nature of sampling errors (i.e., errors in estimating
⁷¹ the index flood at gauged sites) and the existence of regression errors. *Robson and Reed*⁷² [1999] and *Kjeldsen and Jones* [2009a] further generalized the GLS approach by consid⁷³ ering spatially dependent regression errors.

⁷⁴ Despite these advances, GLS-based transfer of information from gauged to ungauged site
⁷⁵ still relies on the following two-step procedure:

⁷⁶ 1. Local estimation: an index flood is first chosen (e.g. the at-site mean or median), ⁷⁷ and is estimated at each gauged site. This estimation is affected by sampling errors, ⁷⁸ which are spatially dependent and whose variance varies from site to site. Consequently, ⁷⁹ the covariance matrix $\hat{\Sigma}$ of sampling errors is also estimated;

⁸⁰ 2. Regional estimation: A regression model is estimated to link the index flood value ⁸¹ with site/catchment characteristics. Importantly, the estimation of the regression model ⁸² accounts for the existence of sampling errors in a GLS framework, and is hence performed ⁸³ conditionally on $\hat{\Sigma}$. Also note that estimating a regression model is not restricted to esti-⁸⁴ mating the regression coefficients, but also involves estimating the properties of regression ⁸⁵ errors, i.e. their covariance matrix $\hat{\Gamma}$. This is of primary importance since this matrix ⁸⁶ plays an important role in the predictive uncertainty at ungauged sites.

⁸⁷ Such a two-step procedure might be problematic in the context of quantifying the total ⁸⁸ predictive uncertainty. Indeed, estimates at step 2 are conditional on estimates at step 1. ⁸⁹ In particular, the covariance matrix $\hat{\Sigma}$ is an estimate, and may itself be in error (see e.g. ⁹⁰ the discussion in *Stedinger and Tasker* [1985] and *Kroll and Stedinger* [1998] on desirable ⁹¹ properties of $\hat{\Sigma}$). Such error may then propagates to step 2. Consequently, it would ⁹² be desirable to avoid separating the inference process in two separate steps: this can be

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⁹³ achieved using hierarchical models (see following section 1.2).

In addition to these issues related to the estimation procedure, the assumptions underlying 94 index flood approaches remain questionable. Indeed, the scale invariance assumption is 95 convenient because it merges all spatial variability into a single parameter (the index-flood parameter). However, this assumption forces the coefficient of variation of data to remain 97 constant throughout an homogeneous region: this might be too restrictive in some regions, or it might force to drastically reduce the spatial extent of the region to ensure the scale 99 invariance hypothesis is met. Several alternatives have therefore been proposed to move 100 beyond this restrictive framework, in particular region of influence approaches [Burn, 1990] 101 and recent developments [Kjeldsen and Jones, 2009b], normalized quantile regression [Fill 102 and Stedinger, 1998] or empirical Bayes procedures [Kuczera, 1982a, b, 1983] (see also the 103 discussion provided by *Griffis and Stedinger* [2007]). 104

1.2. Bayesian hierarchical models

An alternative approach, based on Bayesian hierarchical models, has been explored 105 more recently. In particular, Wikle et al. [1998] proposed a general hierarchical frame-106 work to describe the spatial variability of the distribution of some environmental variable. 107 The principle of a hierarchical model is to use several modeling layers. For instance, a 108 first layer may assume that the data follow some distribution with unknown parameters, 109 while a second layer may model the variability of those parameters in space, using some 110 regression model. This closely corresponds to the successive steps involved in the stan-111 dard implementation of RFA approaches. However, the main advantage of a hierarchical 112 model is that all unknown quantities can be inferred simultaneously, therefore accounting 113 for possible interactions between estimation errors made at different layers and yielding 114

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a more rigorous quantification of the predictive uncertainty. In other words, hierarchical
models allow describing the local variability of data *together* with their regional coherence,
without separating the inference process in several steps. Moreover, such models are more
general than the model underlying the index flood approach, since they do not require
assuming scale invariance.

Several applications of Bayesian hierarchical models in a hydrological context have been 120 proposed in the literature. For instance, *Cooley et al.* [2007] described the spatial variabil-121 ity of extreme rainfall, while Aryal et al. [2009] extended this description to both spatial 122 and temporal variability. Similarly, *Lima and Lall* used Bayesian hierarchical models to 123 describe daily rainfall occurrences [Lima and Lall, 2009] or runoff extremes [Lima and 124 Lall, 2010 in a regional context. In addition to these hydrological applications, similar 125 Bayesian hierarchical models have been used in other fields, e.g. for extreme wind speed 126 modeling [Coles and Casson, 1998; Casson and Coles, 1998, 2000]. 127

Despite improving the standard implementation of RFA approaches, all Bayesian hierar-128 chical models described above rely on an assumption of conditional independence: data 129 are assumed spatially independent given the values of their distribution's parameters. 130 This would be a valid assumption if most spatial covariation in the data was explained 131 by the spatial covariation in the parameters. However, it is questionable since spatial 132 dependence between data in the one hand and parameters in the other hand arise from 133 distinct processes, as noted by *Cooley et al.* [2007]: in a nutshell, data dependence can be 134 interpreted as *weather* spatial dependence, while dependence between parameters (also 135 termed process dependence) relates to *climate* spatial dependence. Weather should ex-136 hibit spatial dependence (at least within a short distance range), even if the climate were 137

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¹³⁸ perfectly known.

Examples of spatial hierarchical models explicitly accounting for intersite dependence in 139 the observations are very few. *Perreault* [2000] proposed such a model to detect a re-140 gional step-change in annual runoff. Alternatively, Micevski et al. [2006] and Micevski 141 [2007] proposed a Bayesian hierarchical regional flood model accounting for data depen-142 dence. However, in both cases, the explicit description of data dependence was rooted to 143 a particular distributional assumption: Gaussian data were assumed by *Perreault* [2000], 144 while *Micevski et al.* [2006] and *Micevski* [2007] used a mixture of log-normal distributions. 145 Unfortunately, Gaussian-related assumptions may be too restrictive for other hydrologic 146 variables or in other geographical contexts. 147

1.3. Objectives

¹⁴⁸ Building on previous work described in the preceding sections, this paper aims to derive ¹⁴⁹ a general Bayesian hierarchical framework for RFA. In particular, this framework should ¹⁵⁰ enable an explicit description of spatial dependence between data, without relying on ¹⁵¹ Gaussian-related distributional assumptions. This is achieved by means of the elliptical ¹⁵² copula family [*Genest and Favre*, 2007], which constitutes a convenient tool to model ¹⁵³ dependence in a highly dimensional and non-Gaussian context.

The paper is organized as follows. Section 2 describes the two-level hierarchical framework, with level 1 modeling the joint distribution of observations and level 2 modeling the variation of the distribution' parameters in space. Section 3 describes the inference of the hierarchical model and its use for prediction at both gauged and ungauged sites. Section 4 illustrates the application of the proposed framework for the regional estimation of

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¹⁵⁹ extreme rainfall. Avenues for further improvement are identified and discussed in section

¹⁶⁰ 5, before summarizing the main results in section 6.

2. A Bayesian hierarchical modeling framework

2.1. Data level

Let Y(s,t) denote the variable of interest at site s and time t. For instance, Y(s,t) may 161 represent the annual maximum daily rainfall at site s and year t. Let us further assume 162 that observations are available at sites $\tilde{s}_1, \dots, \tilde{s}_M$ and times t=1: T. The $(T \times M)$ observation 163 matrix is denoted by $\tilde{\boldsymbol{y}} = (\tilde{\boldsymbol{y}}(\tilde{s}_1), ..., \tilde{\boldsymbol{y}}(\tilde{s}_M))$, with $\tilde{\boldsymbol{y}}(\tilde{s}_i) = (\tilde{y}(\tilde{s}_i, t))_{t=1:T}$ denoting the time 164 series of observations at a given site \tilde{s}_k . The shorthand notation $\tilde{y}(t) = (\tilde{y}(\tilde{s}_i, t))_{i=1:M}$ is 165 also used to denote the (spatial) vector of observations at a given time t. Note that this 166 notation assumes that all sites share the same observation period. Although considering 167 non-concomitant observation periods would not affect the modeling hypotheses, it would 168 certainly create tedious notation and complicate the model implementation (see discussion 169 in section 5.1). 170

¹⁷¹ 2.1.1. At-site distribution

At an arbitrary site *s* within the area of study, and at any time *t*, Y(s, t) is assumed to be a realization from a given distribution whose parameters vary in space (see left-hand-side of Figure 1):

$$Y(s,t) \sim p(\boldsymbol{\theta}(s)) \tag{1}$$

RFA involves estimating the parameter vector $\boldsymbol{\theta}(s)$ at any (gauged or ungauged) site s. In order to avoid confusion with other parameters that will be introduced later on, the term "D-parameters" is systematically used to denote the parameters $\boldsymbol{\theta}(s) = (\theta_k(s))_{k=1:D}$ Author-produced version of the article published in Water Resources Research (2011) vol. 47, doi : 0.1029/2010WR010089. The original publication is available at http://www.agu.org/journals/wr/

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¹⁷⁸ of the parent distribution. Note that D-parameters are allowed to vary in space but not in ¹⁷⁹ time. An extension of the framework for allowing time-varying D-parameters is possible ¹⁸⁰ but is left for future work, since this paper mainly focuses on spatial aspects.

¹⁸¹ 2.1.2. Joint distribution

In a spatial context, the derivation of the likelihood of observations $\tilde{\boldsymbol{y}}$ requires knowing 182 the multivariate distribution of the spatial observation vector $\tilde{\boldsymbol{u}}(t)$ at any time t. An 183 assumption often made (explicitly or implicitly) by some regional estimation methods is 184 that the data are spatially independent. In this case, the multivariate pdf is simply equal 185 to the product of marginal pdfs in (1). Implications of this assumption are discussed 186 by e.g. Stedinger [1983], Hosking and Wallis [1988], Madsen and Rosbjerg [1997] and 187 Renard and Lang [2007]. In particular, these authors demonstrate that the variance (i.e. 188 the uncertainty) of estimated quantities is underestimated when data do not support the 189 independence assumption. 190

One of the main objectives of the framework presented in this paper is to overcome this limitation, by explicitly modeling spatial dependence. The approach taken to achieve this objective has to account for two important points: (i) it has to be applicable for an arbitrary choice of marginal distribution in (1); (ii) since regional analysis may involve hundreds of sites, it has to remain practical with high-dimensional data.

Point (i) above makes copulas a natural candidate to model dependence in the context of this framework. Indeed, the copula theory is based on the description of the dependence structure independently of marginal distributions [e.g., *Favre et al.*, 2004]. Point (ii) suggests focusing on the elliptical copula family, because of its ability to describe highdimensional datasets. A thorough description of the elliptical copula family can be found

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in the paper by *Genest and Favre* [2007], and a summary of the main characteristics of 201 two particular members (the Gaussian and the Student copulas) is given in Appendix A. 202 A *M*-dimensional elliptical copula is parameterized by a $(M \times M)$ symmetric dependence 203 matrix Σ describing pairwise dependences. Additional parameters η may be used for 204 some members of the family (e.g. the Student copula) to describe the strength of tail 205 dependence. The multivariate distribution is finally derived by combining this depen-20 dence structure with the marginal distributions given by equation (1) (see Appendix A 207 for details). An example of application of an elliptical copula (the Gaussian copula) to 208 regional frequency analysis is given by *Renard and Lang* [2007]. 209

Formally, it is assumed that the multivariate distribution of data from any set of M sites can be derived from an M-dimensional elliptical copula, with pairwise dependence matrix Σ , additional dependence parameters η and marginal distributions $\{p(\boldsymbol{\theta}(s_1)), ..., p(\boldsymbol{\theta}(s_M))\}$:

$$(Y(s_1,t),...,Y(s_M,t)) \sim EC_M(\Sigma, \eta, \{\theta(s_1),...,\theta(s_M)\})$$
(2)

The symbol EC_M stands for "M-dimensional Elliptical Copula", and analytical formulas for the corresponding joint pdf are given in Appendix A.

In order to simplify the model, it is further assumed that the dependence between data from two sites solely depends on the inter-site distance. An analogy can be drawn with the common treatment of stationary and isotropic spatial random fields in geostatistics [e.g., *Chiles and Delfiner*, 1999]. It follows that the elements of the pairwise dependence matrix Σ can be expressed as a function of the inter-site distance, parameterized by some vector ψ (see Figure 1):

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$$\Sigma(i,j) = \Psi(\|s_i - s_j\|; \psi)$$
(3)

Note that the inter-site distance is not necessarily Euclidean in the bi-dimensional xyspace (geographical coordinates). For instance, a distance in the tri-dimensional space xy+elevation may be useful if inter-site dependence is lower for sites having contrasted elevations. More generally, the distance may depend on covariates other than geographical coordinates [e.g., *Cooley et al.*, 2007; *Blanchet and Davison*, 2011]. However, for the sake of simplicity, this is not made explicit in the notation of equation (3). This topic is further discussed in section 5.8.

²²⁹ A valid dependence-distance function $\Psi(.; \psi)$ must ensure the positive-definiteness of the ²³⁰ dependence matrix Σ . Such a function can be chosen amongst the numerous covariogram ²³¹ models (e.g., exponential, spherical, Gaussian) existing in geostatistics. By analogy with ²³² the covariogram used in geostatistics, the dependence-distance model in (3) is termed ²³³ "dependogram". Note that a distinct naming convention is used because the elements ²³⁴ of the matrix Σ are not equal to the covariances between data pairs, but rather to the ²³⁵ dependence coefficients of the elliptical copula.

It is stressed that the use of an elliptical copula to model spatial dependence is motivated by practical considerations. Although the elliptical family is quite flexible (in particular, it encompasses tail-dependent and tail-independent models), there can be no guarantee that an elliptical copula will be able to model the data at hand. In particular, a multivariate distribution derived from an elliptical copula is not compatible with the family of multivariate extreme value distributions [*Mikosch*, 2005]. It follows that the use of the elliptical copula model in extrapolation to estimate low probabilities (e.g.,

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²⁴³ $Pr(Y(s_1,t) > u \cap ... \cap Y(s_M,t) > u)$ for some large u) is likely to yield inaccurate re-²⁴⁴ sults. However, the aim of the framework presented in this paper is not to estimate the ²⁴⁵ probability of extreme multivariate events, but rather to estimate the distribution of the ²⁴⁶ variable of interest at any site, while accounting for the existence of dependences at ob-²⁴⁷ served levels. The implications of the elliptical copula assumption are more thoroughly ²⁴⁸ discussed in section 5.2.

2.2. Process level

Section 2.1 described the construction of the multivariate distribution of data by specifying the following three components: (i) the at-site distribution (1); (ii) the elliptical copula (2) used to model spatial dependence; (iii) the dependogram (3). The second level of the hierarchical framework aims to describe the variation of the D-parameters in space, by means of a Gaussian spatial process whose mean depends on covariates describing the site (or catchment) characteristics (see right-hand-side of Figure 1).

²⁵⁵ 2.2.1. D-parameter regression model

Let $\theta(s) = (\theta_k(s))_{k=1:D}$ be the D-parameter vector. A regression model is used to describe the spatial variation of each D-parameter $\theta_k(s)$ as follows:

$$g_k(\theta_k(s)) = h_k(\boldsymbol{x}_k(s); \boldsymbol{\beta}_k) + \epsilon_k(s)$$
(4)

²⁵⁸ Equation (4) uses the following components:

1. One-to-one function g_k is termed the "link function" by analogy with generalized linear models [e.g., *Dobson*, 2001]. The identity function may be used in most cases. However, alternative functions might be useful for some D-parameters, e.g. a logarithm

function to ensure positivity or a logit function for a D-parameter comprised between zero and one.

264 2. Vector $\boldsymbol{x}_k(s)$ represents the set of covariates used to describe the characteristics 265 of site (or catchment) s (e.g., elevation, distance to sea, catchment size, etc.). As in 266 any regression framework, the choice of relevant covariates is of primary importance, and 267 preliminary analyzes are often useful e.g. to eliminate highly correlated covariates.

3. Function $h_k(.; \boldsymbol{\beta}_k)$ is the regression function, parameterized by a vector of regression parameters $\boldsymbol{\beta}_k$. The most common choice of regression function is the linear function $h_k(\boldsymbol{x}_k(s); \boldsymbol{\beta}_k) = \boldsymbol{x}_k(s)\boldsymbol{\beta}_k$, but alternative functions may be used.

4. $\epsilon_k(s)$ is the residual of the regression. It stems from the imperfect nature of the regression model, and is therefore termed "regression error". In the hierarchical modeling context described in this paper, regression errors are treated as latent variables, i.e. unobserved variables that need to be inferred.

275 2.2.2. Regression errors

Regression errors $\boldsymbol{\epsilon}_k = (\epsilon_k(s))_{s=1:M}$ are commonly assumed to be spatially independent 276 Stedinger and Tasker, 1985, 1986; Reis et al., 2005]. However, recent work by e.g., 277 Kieldsen and Jones [2009a] suggests that regression errors can be significantly dependent, 278 especially for nearby sites and/or regression models with poor predictive ability. Ignoring 279 spatial dependences between regression errors possibly affects regional frequency analysis 280 in two distinct ways: (i) it may impact the accuracy and/or precision of estimates for the 281 regression parameters; (ii) it may impact the predictive variance (i.e. the uncertainty in 282 quantities estimated at ungauged sites) by not taking advantage of the regression errors 283 estimated as nearby sites. Consequently, following Kjeldsen and Jones [2009a], regression 284

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errors are assumed to be a realization from a spatial Gaussian field with mean zero and covariance matrix Γ_k :

$$(\epsilon_k(s_1), \dots, \epsilon_k(s_M)) \sim N_M(0; \Gamma_k) \tag{5}$$

A covariogram model is used to relate pairwise covariances to inter-site distances:

$$\Gamma_k(i,j) = \Upsilon_k(\|s_i - s_j\|; \boldsymbol{v}_k) \tag{6}$$

In a Bayesian hierarchical context, the Gaussian distribution used to describe regression errors is termed the hyper-distribution. Parameters v_k defining the covariance matrix are termed the hyper-parameters (see Figure 1). Note that as in previous section 2.1.2, the inter-site distance is not necessarily Euclidean and can be interpreted in a wider sense.

²⁹² Moreover, note that regression errors $\epsilon_k(s)$ and $\epsilon_q(s)$ related to two distinct D-parameters ²⁹³ $\theta_k(s)$ and $\theta_q(s)$ will be described by independently using equations (5-6) twice. This ²⁹⁴ implies that an assumption of independence between regression errors $\epsilon_k(s)$ and $\epsilon_q(s)$ is ²⁹⁵ effectively made. This assumption will be further discussed subsequently (see sections 4 ²⁹⁶ and 5.3).

2.3. Remarks

In the hierarchical framework presented in previous sections, two distinct dependence structures are used: the dependogram (3) aims at describing dependence between observations, while the covariogram (6) aims at describing dependence between the parameters of their distribution (or more accurately, between the errors of the regression linking the D-parameters with catchment/site characteristics). As explained in the introduction, the former structure relates to weather spatial dependence, while the latter relates to climate

spatial dependence [see also the discussion in *Cooley et al.*, 2007]. These dependence 303 structures play distinct roles in the context of regional frequency analysis: climate spa-304 tial dependence allows transferring information from gauged to ungauged sites, while the 305 main effect of weather dependence is to diminish the information content of data collected 306 at nearby sites. Consequently, both dependence structures are likely to have an impact 307 on predictions and should therefore be accounted for. Note that the assumption of con-308 ditional independence frequently made in similar Bayesian hierarchical frameworks (see 309 introduction) corresponds to forcing the dependence matrix Σ in equation (3) to unity. 310 Moreover, the regression model of equation (4) yields several interesting particular cases 311 when the regression errors are forced to zero: 312

1. A purely local D-parameter (i.e. whose value is site-specific) can be obtained with the regression model $\theta_k(\tilde{s}) = \beta_k^{(\tilde{s})}$. This corresponds to introducing as many regression parameters $\beta_k^{(\tilde{s})}$ as there are sites, and use them to model a site effect. An obvious drawback of this approach is that the estimation at ungauged site *s* is not directly possible since the site effect $\beta_k^{(s)}$ is unknown.

2. A regional D-parameter (i.e. having an identical value for all sites within the region) corresponds to the regression model $\theta_k(\tilde{s}) = \beta_k$. This is a rather strong assumption, which may yield an underestimation of the predictive uncertainty. Indeed, in the absence of regression errors, the uncertainty in the estimation of θ_k is identical for all sites (gauged or not).

323 3. The strong assumption that parameter $\theta_k(\tilde{s})$ is identical for all sites can be relaxed 324 by using a regression model $g_k(\theta_k(\tilde{s})) = h_k(\boldsymbol{x}_k(\tilde{s}); \boldsymbol{\beta}_k)$. This corresponds to the "covariate 325 modeling" approach proposed by several authors [e.g., *Katz et al.*, 2002; *Maraun et al.*,

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³²⁶ 2009]. However, the possible underestimation of predictive uncertainty in the absence of ³²⁷ regression errors still holds.

Lastly, the Bayesian hierarchical framework shares several similarities with the GLS 328 approach proposed by Stedinger and Tasker [1985, 1986] and its latest developments by 329 *Kjeldsen and Jones* [2009a]. Indeed, GLS also uses two spatial dependence structures: (i) 330 the covariance matrix of sampling estimation errors (e.g., errors in estimating an index 331 flood at gauged sites), which results from the spatially dependent nature of data; (ii) the 332 covariance matrix of regression errors (termed model errors in the GLS approach). The 333 framework presented in this paper therefore borrows from GLS the objective of modeling 334 the spatial variability in the distinct sources of errors affecting RFA. However, it differs 335 in its implementation. GLS first estimates the spatial variability of sampling errors, and 336 then uses this estimation to fit the regression in a second step. By contrast, the proposed 337 framework performs the inference in a single step, which facilitates the quantification of 338 the total predictive uncertainty (see discussion in the introduction section). Moreover, the 339 regression is applied on each parameter, while in general GLS is rather applied to a single 340 hydrologic quantity (e.g. an index flood or a quantile, but see Tasker and Stedinger [1989] 341 and Griffis and Stedinger [2007] for exceptions). This is a more general approach, at least 342 in principle; however what level of model complexity can be identified given the limited 343 information content of data remains an open question (this will be further discussed in 344 section 5.5). 345

3. Estimation and prediction

3.1. Posterior distribution

- The application of the hierarchical framework described in section 2 requires estimating
- ³⁴⁷ the following unknown quantities:
- 348 1. The dependogram parameters ψ
- ³⁴⁹ 2. Additional parameters of the elliptical copula η
- 350 3. The regression parameters $\boldsymbol{\beta}_k$, for k = 1:D
- 4. The regression errors $\boldsymbol{\epsilon}_k$, for k = 1:D (latent variables)
- 5. The covariogram parameters \boldsymbol{v}_k , for k = 1:D (hyper-parameters)

The posterior distribution of these quantities, given observations \tilde{y} and covariates x, can be derived as follows:

$$p(\boldsymbol{\psi}, \boldsymbol{\eta}, \boldsymbol{\beta}_{k=1:D}, \boldsymbol{\epsilon}_{k=1:D}, \boldsymbol{\upsilon}_{k=1:D} | \tilde{\boldsymbol{y}}, \boldsymbol{x})$$

$$\propto p(\tilde{\boldsymbol{y}} | \boldsymbol{\psi}, \boldsymbol{\eta}, \boldsymbol{\beta}_{k=1:D}, \boldsymbol{\epsilon}_{k=1:D}, \boldsymbol{\upsilon}_{k=1:D}, \boldsymbol{x}) p(\boldsymbol{\psi}, \boldsymbol{\eta}, \boldsymbol{\beta}_{k=1:D}, \boldsymbol{\epsilon}_{k=1:D}, \boldsymbol{\upsilon}_{k=1:D} | \boldsymbol{x})$$
(7a)

$$= p(\tilde{\boldsymbol{y}}|\boldsymbol{\psi},\boldsymbol{\eta},\boldsymbol{\beta}_{k=1:D},\boldsymbol{\epsilon}_{k=1:D},\boldsymbol{x})p(\boldsymbol{\psi},\boldsymbol{\eta},\boldsymbol{\beta}_{k=1:D})p(\boldsymbol{\epsilon}_{k=1:D},\boldsymbol{v}_{k=1:D})$$
(7b)

$$= p(\tilde{\boldsymbol{y}}|\boldsymbol{\psi},\boldsymbol{\eta},\boldsymbol{\beta}_{k=1:D},\boldsymbol{\epsilon}_{k=1:D},\boldsymbol{x})p(\boldsymbol{\psi},\boldsymbol{\eta},\boldsymbol{\beta}_{k=1:D})p(\boldsymbol{\epsilon}_{k=1:D}|\boldsymbol{v}_{k=1:D})p(\boldsymbol{v}_{k=1:D})$$
(7c)

- ³⁵⁵ The following assumptions have been made to derive this posterior distribution:
- 1. Equation (7a) is a direct application of Bayes theorem.

2. Equation (7b) assumes: (i) the prior distribution does not depend on covariates \boldsymbol{x} ; (ii) prior independence between $\boldsymbol{\psi}, \boldsymbol{\eta}, \boldsymbol{\beta}_{k=1:D}$ in the one hand, and $\boldsymbol{\epsilon}_{k=1:D}, \boldsymbol{v}_{k=1:D}$ in the other hand. This assumption aims at isolating the hierarchical components of the model, i.e. $\boldsymbol{\epsilon}_{k=1:D}$ and $\boldsymbol{v}_{k=1:D}$. Moreover, it is noted that the likelihood of observations can be derived without using the hyper-parameters $\boldsymbol{v}_{k=1:D}$. Indeed, regression errors $\boldsymbol{\epsilon}_{k=1:D}$

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³⁶² suffice to apply the regression model (4) and hence to derive D-parameters (see detailed ³⁶³ derivation of the likelihood in following equation (9)).

- ³⁶⁴ 3. Equation (7c) is an application of conditional probability rules.
- ³⁶⁵ The posterior distribution (7) is made up of the following components:

1. The terms $p(\boldsymbol{\psi}, \boldsymbol{\eta}, \boldsymbol{\beta}_{k=1:D})$ and $p(\boldsymbol{v}_{k=1:D})$ are priors for the inferred parameters and hyper-parameters, respectively.

2. The term $p(\boldsymbol{\epsilon}_{k=1:D}|\boldsymbol{\upsilon}_{k=1:D})$ represents the hierarchical part of the model. Assuming regression errors $\boldsymbol{\epsilon}_k$ related to different D-parameters θ_k are mutually independent yields:

$$p(\boldsymbol{\epsilon}_{k=1:D}|\boldsymbol{\upsilon}_{k=1:D}) = \prod_{k=1}^{D} f_N(\boldsymbol{\epsilon}_k(\tilde{s}_1), \dots, \boldsymbol{\epsilon}_k(\tilde{s}_M)|\boldsymbol{0}, \boldsymbol{\Gamma}_k(\boldsymbol{\upsilon}_k))$$
(8)

where $\Gamma_k(\boldsymbol{v}_k)$ is the covariance matrix derived from the covariogram model (6) and 368 $f_N(\boldsymbol{z}|\boldsymbol{\mu},\boldsymbol{\Gamma})$ represents the pdf of a multivariate normal distribution with mean $\boldsymbol{\mu}$ and 369 covariance matrix Γ . Note that the assumption that there is no cross-correlation between 370 regression errors related to different D-parameters may be restrictive. This assumption 371 will be evaluated in the case study (section 4) and will be further discussed in section 5.3. 372 3. The term $p(\tilde{y}|\psi, \eta, \beta_{k=1:D}, \epsilon_{k=1:D}, x)$ is the likelihood of observations. Its derivation 373 requires further explanation. At a given time step t, the likelihood of (spatial) observations 374 $\tilde{\boldsymbol{y}}(t)$ can be computed as follows: 375

$$p(\tilde{\boldsymbol{y}}(t)|\boldsymbol{\psi},\boldsymbol{\eta},\boldsymbol{\beta}_{k=1:D},\boldsymbol{\epsilon}_{k=1:D},\boldsymbol{x}) = f_{EC}(\tilde{\boldsymbol{y}}(t)|\boldsymbol{\Sigma}(\boldsymbol{\psi}),\boldsymbol{\eta},\{[\theta_k(\boldsymbol{x}_k(\tilde{s}_1),\boldsymbol{\beta}_k,\epsilon_k(\tilde{s}_1))]_{k=1:D},...,[\theta_k(\boldsymbol{x}_k(\tilde{s}_M),\boldsymbol{\beta}_k,\epsilon_k(\tilde{s}_M))]_{k=1:D}\})$$
(9)

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In equation (9), $\Sigma(\psi)$ is the dependence matrix of the elliptical copula, derived from the dependogram model (3). For a given observation site \tilde{s}_j , the *k*th D-parameter $\theta_k(\boldsymbol{x}_k(\tilde{s}_j), \boldsymbol{\beta}_k, \epsilon_k(\tilde{s}_j))$ is derived by applying the regression equation (4), i.e.:

$$\theta_k(\boldsymbol{x}_k(\tilde{s}_j), \boldsymbol{\beta}_k, \epsilon_k(\tilde{s}_j)) = g_k^{-1} \left(h_k(\boldsymbol{x}_k(\tilde{s}_j), \boldsymbol{\beta}_k) + \epsilon_k(\tilde{s}_j) \right)$$
(10)

Lastly, $f_{EC}(\boldsymbol{z}|\boldsymbol{\Sigma}, \boldsymbol{\eta}, \{\boldsymbol{\theta}(s_1), ..., \boldsymbol{\theta}(s_M)\})$ represents the multivariate pdf of a *M*-dimensional vector \boldsymbol{z} derived from the elliptical copula with pairwise dependence matrix $\boldsymbol{\Sigma}$, additional dependence parameters $\boldsymbol{\eta}$ and marginal distributions $\{p(\boldsymbol{\theta}(s_1)), ..., p(\boldsymbol{\theta}(s_M))\}$, as detailed in Appendix A.

Assuming temporal independence between observations $\tilde{y}(t)$, the likelihood of the whole observation matrix \tilde{y} is simply obtained as follows:

$$p(\tilde{\boldsymbol{y}}|\boldsymbol{\psi},\boldsymbol{\eta},\boldsymbol{\beta}_{k=1:D},\boldsymbol{\epsilon}_{k=1:D},\boldsymbol{x}) = \prod_{t=1}^{T} p(\tilde{\boldsymbol{y}}(t)|\boldsymbol{\psi},\boldsymbol{\eta},\boldsymbol{\beta}_{k=1:D},\boldsymbol{\epsilon}_{k=1:D},\boldsymbol{x})$$
(11)

3.2. Inference

The posterior distribution (7) poses a computational challenge because its dimension 385 grows with the number of sites. This is due to the explicit modeling of regression errors 386 through latent variables. Consequently, the number of quantities to be inferred from the 387 posterior can amount to hundreds. This is typical of hierarchical models with latent vari-388 ables used to describe unobserved processes [e.g., Clark, 2005]. A Markov Chain Monte 389 Carlo (MCMC) sampler is used to address this difficulty. It is stressed that MCMC sam-390 pling from high-dimensional posteriors is challenging but by no means insurmountable. 391 Successful examples are provided by e.g. Crainiceanu et al. [2003]; Vrugt et al. [2008]; 392

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³⁹³ Thyer et al. [2009]; Reichert and Mieleitner [2009] in Hydrology, not to mention numer-³⁹⁴ ous applications in other fields [e.g., Clark, 2003, in ecology][Storz and Beaumont, 2002, ³⁹⁵ in genetics, etc.].

The MCMC sampler used in this paper is made up of two stages. Stage one makes use of 396 an adaptive block Metropolis algorithm with univariate Gaussian jump distributions (i.e. 397 components of the parameter vector are updated one at a time). The adaption strategy 39 adjusts the jump variances to produce an adequate jump rate see *Renard et al.*, 2006, for 399 a detailed description]. This first sampler is used to perform a preliminary exploration 400 of the posterior distribution properties (notably in terms of posterior covariance). In 401 the second stage, a standard Metropolis sampler [Metropolis and Ulam, 1949; Metropolis 402 et al., 1953] is used, with a Gaussian jump distribution whose covariance matrix is speci-403 fied using the preliminary exploration performed at stage one. Convergence is assessed by 404 evolving four parallel chains and verifying that the Gelman-Rubin criteria [Gelman et al., 405 1995 are close to one for all inferred quantities. 406

⁴⁰⁷ Note that additional computing efficiency might be achieved in some cases by using con-⁴⁰⁸ jugate priors. However, this possibility is not investigated in this paper.

3.3. Prediction at gauged site

Once inference has been performed using the MCMC strategy outlined above, the next step is to use the model to predict some quantity of interest. In this section, focus is on prediction at a gauged site \tilde{s} . A typical quantity of interest is the *p*-quantile of the distribution of observations at site \tilde{s} , which can be directly derived from D-parameters $\theta(\tilde{s})$.

⁴¹⁴ In a Bayesian context, another interesting byproduct of the posterior distribution is the

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predictive distribution of a (future) observation. In general, if Θ denotes the vector of all parameters subject to inference, the pdf of this predictive distribution, evaluated at some value w, is mathematically defined as follows [Gelman et al., 1995]:

$$p(w|\tilde{\boldsymbol{y}}) = \int p(w|\boldsymbol{\Theta})p(\boldsymbol{\Theta}|\tilde{\boldsymbol{y}})d\boldsymbol{\Theta}$$
(12)

In practice, this integration is not performed analytically but is approximated using the MCMC replicates from the posterior (7). The approximation algorithm simply consists in generating a value from the at-site distribution (1) for each MCMC replicate. The resulting sample is a realization from the predictive distribution and can therefore be used to estimate its characteristics (mean, variance, probability interval, etc.).

Let us assume that MCMC sampling generated a set of n_{sim} replicates from the posterior distribution, $(\boldsymbol{\psi}^{(j)}, \boldsymbol{\eta}^{(j)}, \boldsymbol{\beta}_{k=1:D}^{(j)}, \boldsymbol{\epsilon}_{k=1:D}^{(j)}, \boldsymbol{v}_{k=1:D}^{(j)})_{j=1:n_{sim}}$. These replicates can be used to make a prediction at gauged site \tilde{s} using the following algorithm:

426 Do
$$j = 1 : n_{sim}$$

⁴²⁷ 1. compute D-parameters from the regression model (4), $\theta_k^{(j)}(\tilde{s}) = g_k^{-1} \left(h_k(\boldsymbol{x}_k(\tilde{s}), \boldsymbol{\beta}_k^{(j)}) + \epsilon_k^{(j)}(\tilde{s}) \right)$ ⁴²⁸ for k = 1 : D

⁴²⁹ 2. compute derived quantity $z^{(j)} = z(\boldsymbol{\theta}^{(j)}(\tilde{s}))$ (e.g., the T-year quantile) or generate a ⁴³⁰ value from the at-site distribution $w^{(j)} \sim p(\boldsymbol{\theta}^{(j)}(\tilde{s}))$

The samples $(z^{(j)})_{j=1:n_{sim}}$ and $(w^{(j)})_{j=1:n_{sim}}$ can be considered as realizations from the posterior distribution of the quantity of interest and from the predictive distribution, respectively.

3.4. Prediction at ungauged site

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The fundamental difference with the prediction approach presented in previous section 3.3 is that the regression error $\epsilon_k(s)$ has not been inferred for an ungauged site s. However, the properties of the hyper-distribution (5) have been inferred. Since the hyperdistribution is the joint distribution of a vector of regression errors, it can be used to indirectly infer the properties of the regression error $\epsilon_k(s)$ for the target ungauged site s. More precisely, the predictive distribution of the regression error $\epsilon_k(s)$, given observed data $\tilde{\boldsymbol{y}}$ and covariates \boldsymbol{x} , can be computed as follows:

$$p(\epsilon_{k}(s)|\tilde{\boldsymbol{y}},\boldsymbol{x}) = \int p(\epsilon_{k}(s),\epsilon_{k}(\tilde{s}_{1}),...,\epsilon_{k}(\tilde{s}_{M}),\boldsymbol{v}_{k}|\tilde{\boldsymbol{y}},\boldsymbol{x})d\epsilon_{k}(\tilde{s}_{1})...d\epsilon_{k}(\tilde{s}_{M})d\boldsymbol{v}_{k}$$
(13a)
$$= \int p(\epsilon_{k}(s)|\epsilon_{k}(\tilde{s}_{1}),...,\epsilon_{k}(\tilde{s}_{M}),\boldsymbol{v}_{k},\tilde{\boldsymbol{y}},\boldsymbol{x})p(\epsilon_{k}(\tilde{s}_{1}),...,\epsilon_{k}(\tilde{s}_{M}),\boldsymbol{v}_{k}|\tilde{\boldsymbol{y}},\boldsymbol{x})d\epsilon_{k}(\tilde{s}_{1})...d\epsilon_{k}(\tilde{s}_{M})d\boldsymbol{v}_{k}$$
(13a)

The second term in equation (13b) is the posterior distribution of $(\epsilon_k(\tilde{s}_1), ..., \epsilon_k(\tilde{s}_M), \boldsymbol{v}_k)$. 441 It can therefore be directly approximated using the MCMC replicates. The first term 442 represents the distribution of the regression error $\epsilon_k(s)$ for the target ungauged site s, 443 conditional on regression errors at gauged sites $\epsilon_k(\tilde{s}_1)...\epsilon_k(\tilde{s}_M)$ and hyper-parameters \boldsymbol{v}_k . 444 Following equation (5), the joint distribution of $\epsilon_k(s)$ and $\epsilon_k(\tilde{s}_1)...\epsilon_k(\tilde{s}_M)$ (conditional 445 on \boldsymbol{v}_k) is a multivariate Gaussian distribution with mean zero and covariance matrix 446 $\Gamma_k = \Gamma_k(\boldsymbol{v}_k)$ derived from the covariogram (6). Let us partition this covariance matrix as 447 follows: 448

$$\Gamma_k = \begin{pmatrix} \sigma_k^2 & \Lambda_k \\ \Lambda_k^t & \Omega_k \end{pmatrix}$$
(14)

where σ_k^2 is the marginal variance of regression errors, $\Omega_k = \Omega_k(\boldsymbol{v}_k)$ is the $M \times M$ covariance matrix of regression errors at gauged sites $\epsilon_k(\tilde{s}_1)...\epsilon_k(\tilde{s}_M)$, and $\Lambda_k = \Lambda_k(\boldsymbol{v}_k)$

(13b)

⁴⁵¹ is the 1 × *M* vector of covariances between $\epsilon_k(s)$ and $\epsilon_k(\tilde{s}_1)...\epsilon_k(\tilde{s}_M)$. Using a well-known ⁴⁵² formula for conditional Gaussian distributions, it follows that the conditional distribution ⁴⁵³ in equation (13b) is a univariate Gaussian distribution with mean $\mu_{k,cond}$ and variance ⁴⁵⁴ $\sigma_{k,cond}^2$:

$$p(\epsilon_k(s)|\epsilon_k(\tilde{s}_1), ..., \epsilon_k(\tilde{s}_M), \boldsymbol{v}_k, \tilde{\boldsymbol{y}}, \boldsymbol{x}) = N(\mu_{k,cond}; \sigma_{k,cond}^2)$$
$$\mu_{k,cond} = \boldsymbol{\Lambda}_k \boldsymbol{\Omega}_k^{-1}(\epsilon_k(\tilde{s}_1), ..., \epsilon_k(\tilde{s}_M))^t$$
$$\sigma_{k,cond}^2 = \sigma_k^2 - \boldsymbol{\Lambda}_k \boldsymbol{\Omega}_k^{-1} \boldsymbol{\Lambda}_k^t$$
(15)

⁴⁵⁵ The conditional distribution in equation (15) complements the regression to transfer ⁴⁵⁶ information from gauged to ungauged sites. Indeed, the vector Λ_k acts as a weight vector ⁴⁵⁷ favoring gauged sites nearby the target ungauged site *s*. If $\epsilon_k(s)$ is independent from all ⁴⁵⁸ $\epsilon_k(\tilde{s}_j)$'s, the conditional distribution is equal to the marginal distribution of regression ⁴⁵⁹ errors, i.e. a Gaussian distribution with mean $\mu_{k,cond} = 0$ and variance $\sigma_{k,cond}^2 = \sigma_k^2$. In ⁴⁶⁰ this case, the transfer of information to ungauged site *s* does not favor any particular ⁴⁶¹ gauged site \tilde{s} .

The explanations given above lead to the following algorithm for prediction at an ungauged site s:

464 Do
$$j = 1 : n_{sim}$$

465 1. compute
$$\boldsymbol{\Omega}_{k}^{(j)} = \boldsymbol{\Omega}_{k}(\boldsymbol{v}_{k}^{(j)}); \, \boldsymbol{\Lambda}_{k}^{(j)} = \boldsymbol{\Lambda}_{k}(\boldsymbol{v}_{k}^{(j)}); \, \boldsymbol{\sigma}_{k}^{2(j)} = \boldsymbol{\sigma}_{k}^{2}(\boldsymbol{v}_{k}^{(j)}).$$

466 2. compute
$$\mu_{k,cond}^{(j)}$$
 and $\sigma_{k,cond}^{(j)}$ according to equation (15).

3. generate regression error $\epsilon_k^{(j)}(s)$ from the conditional distribution (15) for k = 1 : D(Gaussian distribution with mean $\mu_{k,cond}^{(j)}$ and standard deviation $\sigma_{k,cond}^{(j)}$).

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469 4. compute D-parameters from the regression model (4), $\theta_k^{(j)}(s) = g_k^{-1} \left(h_k(\boldsymbol{x}_k(s), \boldsymbol{\beta}_k^{(j)}) + \epsilon_k^{(j)}(s) \right)$ 470 for k = 1 : D

5. compute derived quantity $z^{(j)} = z(\boldsymbol{\theta}^{(j)}(s))$ or generate a value from the at-site distribution $w^{(j)} \sim p(\boldsymbol{\theta}^{(j)}(s))$

4. Case study: Mediterranean extreme rainfall

The application of the Bayesian hierarchical approach is illustrated with a case study involving extreme rainfall data. The main objectives of this application are to demonstrate the feasibility of a Bayesian hierarchical approach and to assess the impact of some modeling hypotheses (in particular the choices of the data dependence model and the regression model). Note that a synthetic case study was also performed to verify the internal consistency of the modeling framework (not shown).

4.1. Data

Annual maxima from 87 series of daily rainfall are used in this study (Figure 2). Raingauges, whose elevations range from 1 to 1102 m., are located in the French Mediterranean area. This region is characterized by intense rainfall in autumn and is delimited by three mountainous areas: the Pyreneans (South-West), the Alps (East) and the Cevennes (Center). The thin lines in Figure 2 represent six homogeneous regions defined by *Pujol et al.* [2007].

Sixty raingauges are used for estimation, the remaining 27 series being used for model validation. Data are available over the period 1955-2004. Years with more than 15 days of missing values are treated as missing data. Since the treatment of missing data during estimation is not obvious, the 60 estimation series were chosen to ensure the completeness

of annual maxima series over the period 1955-2004 (this point is discussed in more details
in section 5.1).

4.2. Model specification

⁴⁹¹ **4.2.1.** Data level

Let $(y(\tilde{s}_i, t))_{i=1:60, t=1:50}$ denote the observed annual maxima. At any site *s*, the random variable (Y(s, t)) is assumed to follow a Generalized Extreme Value (GEV) distribution with parameters $(\mu(s), \lambda(s), \xi(s))$. The pdf of a GEV distribution with location parameter μ , scale parameter λ and shape parameter ξ is:

$$p(y|\mu,\lambda,\xi) = \frac{1}{\lambda} \left[1 - \xi \left(\frac{y-\mu}{\lambda} \right) \right]^{\frac{1}{\xi}-1} exp \left\{ - \left[1 - \xi \left(\frac{y-\mu}{\lambda} \right) \right]^{1/\xi} \right\}$$
$$\lambda > 0, \xi \neq 0, 1 - \xi \left(\frac{y-\mu}{\lambda} \right) > 0$$
(16)

The joint distribution of observations $\tilde{\boldsymbol{y}}(t)$ is derived using three distinct assumptions on the dependence structure. The first assumption corresponds to using a Gaussian copula [e.g., *Renard and Lang*, 2007] with dependence matrix $\boldsymbol{\Sigma}$:

$$(Y(\tilde{s}_1, t), \dots, Y(\tilde{s}_M, t)) \sim GCop_M(\Sigma, \{\mu(\tilde{s}_i), \lambda(\tilde{s}_i), \xi(\tilde{s}_i)\}_{i=1:M})$$
(17)

The second assumption corresponds to using a Student copula with dependence matrix 500 Σ and tail dependence coefficient ν :

$$(Y(\tilde{s}_1, t), \dots, Y(\tilde{s}_M, t)) \sim SCop_M(\boldsymbol{\Sigma}, \nu, \{\mu(\tilde{s}_i), \lambda(\tilde{s}_i), \xi(\tilde{s}_i)\}_{i=1:M})$$
(18)

Lastly, the third assumption corresponds to assuming spatial independence between data. It can be viewed as a special case of Equation (17) with the dependence matrix Σ

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⁵⁰³ forced to unity.

In both the Student and Gaussian copula cases, the dependogram is formulated as a weighted sum of two exponential functions (with $\psi_1 < \psi_2$) [e.g., *Kjeldsen and Jones*, 2009a]:

$$\Sigma(i,j) = \psi_0 exp(-\psi_1 \|\tilde{s}_i - \tilde{s}_j\|) + (1 - \psi_0) exp(-\psi_2 \|\tilde{s}_i - \tilde{s}_j\|)$$
(19)

507 4.2.2. Process level

The variation of D-parameters $(\mu(s), \lambda(s), \xi(s))$ in space is described using several modeling assumptions. First, an index-flood-like approach, assuming scale invariance, is used. This assumption induces strong constraints on the parameters of at-site distributions. More precisely, a GEV distribution complying with the scale invariance assumption can be reparameterized as follows [e.g., *Ribatet et al.*, 2007]:

$$Y(s,t) \sim GEV(\delta(s)\mu, \delta(s)\lambda, \xi)$$
(20)

Equation (20) states that (i) the shape parameter is constant throughout the region; (ii) the ratio between the location and the scale parameters is constant throughout the region, which implies that annual maxima from all sites have the same coefficient of variation. In other terms, the whole spatial variability of annual maxima is accounted for by the index flood parameter $\delta(s)$, with other parameters μ, λ, ξ being assumed regional.

Two distinct regression models are used in this index flood approach. The first model, \mathcal{M}_1 , does not use any covariate and simply describes the spatial variation of the index flood D-parameter $\delta(s)$ using a Gaussian spatial field:

$$log(\delta(\tilde{s}_i)) = \epsilon(\tilde{s}_i); (\epsilon(\tilde{s}_1), ..., \epsilon(\tilde{s}_M)) \sim N(0, \Gamma)$$
(21)

The second model, \mathcal{M}_2 , uses the raingauge elevation as covariate. Preliminary investigations indicate that the effect of elevation depends on the region. Consequently, the model uses region-specific relationships between the index flood D-parameter $\delta(\tilde{s}_i)$ and elevation:

$$log(\delta(\tilde{s}_i)) = \beta_0^{(j)} + \beta_1^{(j)} * elevation(\tilde{s}_i) + \epsilon(\tilde{s}_i)$$

(\epsilon(\tilde{s}_1), ..., \epsilon(\tilde{s}_M)) \simeq N(0, \mathbf{\Gamma}) (22)

where the superscript (j) is used to denote the region of site \tilde{s}_i . In order to ensure the identifiability of the model, the following additional constraint is applied:

$$\sum_{j=1}^{6} \beta_0^{(j)} = 0 \tag{23}$$

In a second step, the index-flood approach described above is made less restrictive by abandoning the scale invariance assumption. More precisely, the following at-site distribution is assumed:

$$Y(s,t) \sim GEV(\mu(s), \lambda(s), \xi) \tag{24}$$

⁵³⁰ Compared to the index-flood equation (20), equation (24) does not assume a constant ⁵³¹ ratio between the location and the scale parameters throughout the region. However, it ⁵³² still assumes a constant shape parameter. The spatial variation of the location $\mu(s)$ and ⁵³³ scale $\lambda(s)$ D-parameters is described using region-specific relationships with elevation: X - 30 RENARD: HIERARCHICAL REGIONAL ESTIMATION

$$log(\mu(\tilde{s}_i)) = \beta_0^{(j)} + \beta_1^{(j)} * elevation(\tilde{s}_i) + \epsilon_\mu(\tilde{s}_i)$$
$$(\epsilon_\mu(\tilde{s}_1), ..., \epsilon_\mu(\tilde{s}_M)) \sim N(0, \Gamma_\mu)$$

$$log(\lambda(\tilde{s}_i)) = \alpha_0^{(j)} + \alpha_1^{(j)} * elevation(\tilde{s}_i) + \epsilon_\lambda(\tilde{s}_i)$$
$$(\epsilon_\lambda(\tilde{s}_1), ..., \epsilon_\lambda(\tilde{s}_M)) \sim N(0, \Gamma_\lambda)$$
(25)

The model defined by equations (24)-(25) is noted \mathcal{M}_3 .

In all models \mathcal{M}_1 - \mathcal{M}_3 , the covariance matrixes Γ in equations (21), (22) and (25) are parameterized as follows:

$$\Gamma(i,j) = \sigma^2 \left[v_0 exp(-v_1 \| \tilde{s}_i - \tilde{s}_j \|) + (1 - v_0) exp(-v_2 \| \tilde{s}_i - \tilde{s}_j \|) \right]$$
(26)

Finally, vague priors are specified for all parameters by using uniform distributions with large support. The only exception is the shape parameter, for which a Gaussian prior with mean zero and standard deviation 0.3 is specified: this is similar to the "Geophysical prior" used by *Martins and Stedinger* [2000]. Prior specification is further discussed in section 540 5.6.

4.3. Estimation and prediction

In this section, inference is performed using the regression model \mathcal{M}_3 coupled with a Gaussian copula assumption for describing data dependence (equation (17)). This model is noted \mathcal{M}_3 -GCop.

⁵⁴⁵ 4.3.1. Parameter estimates

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Figure 3a shows the posterior pdf of the regional shape parameter ξ . The median value of about -0.12 corresponds to an heavier tail than the Gumbel distribution: this is consistent with previous studies in the same area [*Neppel et al.*, 2007]. Moreover, this estimation is quite precise: this is a consequence of the hypothesis that this parameter is constant throughout the region.

Figure 3b-c shows the posterior pdfs of a few regression errors, for both location (b) 551 and scale (c) regressions (*latent variables* in the vocabulary of hierarchical modeling). 552 Recall that those errors are assumed mutually independent in the inference framework (see 553 section 3). This assumption can be evaluated here by computing the correlation between 554 series of estimated regression errors $\hat{\epsilon}_{\mu}(\tilde{s}_1), ..., \hat{\epsilon}_{\mu}(\tilde{s}_M)$ and $\hat{\epsilon}_{\lambda}(\tilde{s}_1), ..., \hat{\epsilon}_{\lambda}(\tilde{s}_M)$ (maximum-555 posterior estimates are used). A correlation of about 0.52 is found, which suggests that 556 the assumption that both error processes are not cross-correlated might not be realistic. 557 This is further discussed in section 5.3. 558

The standard deviations of regression errors are represented in Figure 3d. Those terms (*hyper-parameters* in the vocabulary of hierarchical modeling) are of primary importance since they control the predictive uncertainty at ungauged sites. In this case study, the estimated hyper-standard deviations are similar for both location and scale regressions. They roughly correspond to a standard error of 15% in the regressions.

⁵⁶⁴ 4.3.2. Data and parameter dependences

Figure 4 shows the estimated dependence-distance relationships, for both the data dependence model (19) and the regression errors dependence model (26). The dependence between data (Figure 4a) is precisely estimated, and confirm that data are not spatially independent. Conversely, the dependence between regression errors (Figure 4b-c) appears Author-produced version of the article published in Water Resources Research (2011) vol. 47, doi : 0.1029/2010WR010089. The original publication is available at http://www.agu.org/journals/wr/

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⁵⁶⁹ more difficult to estimate, with large posterior intervals denoting a lower precision. How-⁵⁷⁰ ever, the strength of dependence is rather limited here, with the correlograms dropping ⁵⁷¹ to near-zero values at relatively short distances of about 10-20 km.

572 **4.3.3.** Prediction

The 100-year daily rainfall is predicted on a $40^{*}50$ grid (yielding $10^{*}10$ km cells) using 573 the procedure detailed in section 3.4. Since the highest raingauge has an elevation of only 574 1102 m, high-elevation areas from the Alps and the Pyreneans (whose maximal elevations 575 reach 4810 m and 3404 m, respectively) were excluded from the prediction (white areas 576 in Figure 5). The left panel of Figure 5 shows the 100-year daily rainfall estimated on 577 the grid using the posterior median. An area with higher $R_{0.99}$ values (reaching 400 mm 578 at some grid points) is located on the Cevennes mountain range: this area is well-known 579 to be affected by the highest rainfall intensities in France. Moreover, the rainfall quantile 580 abruptly drops to smaller values ($\approx 100-150 \text{ mm}$) downwind of the Cevennes. 581

The Bayesian framework used in this paper enables a direct assessment of the uncertainties affecting predictions. The right panel of Figure 5 shows the uncertainty in predicted 100year daily rainfall, measured by the posterior coefficient of variation. This uncertainty appears relatively uniform over the region, and mostly corresponds to 15-20% coefficients of variations. Note however the high-uncertainty area appearing on the foothill of the Alps (East): this is due to the low number of calibration sites in this region (only three sites), which does not enable a precise estimation of the elevation effect.

4.4. Impact of the data dependence model

In this section, the impact of the model used to describe spatial dependence between data is evaluated. To this aim, inference is performed with the three dependence models

⁵⁹¹ described in section 4.2.1, coupled with the regression model \mathcal{M}_3 . The three resulting ⁵⁹² models are noted \mathcal{M}_3 -Inde (independence assumption), \mathcal{M}_3 -GCop (Gaussian copula as-⁵⁹³ sumption) and \mathcal{M}_3 -SCop (Student copula assumption).

⁵⁹⁴ 4.4.1. Parameter estimates

The regional shape parameter is only moderately impacted by the data dependence model (Figure 6a). The posterior pdf of model \mathcal{M}_3 -SCop is slightly shifted and is more variable compared with models \mathcal{M}_3 -Inde and \mathcal{M}_3 -GCop. A similar observation can be made for the hyper-standard deviation of location regression errors (Figure 6b). Posterior pdfs for the hyper-standard deviation of scale regression errors (Figure 6c) have a similar variance but a slightly different mode.

⁶⁰¹ 4.4.2. Data and parameter dependences

Figure 7a shows the estimated dependence between data. The dependogram drops immediately to zero for \mathcal{M}_3 -Inde due to the spatial independence assumption. Dependograms for \mathcal{M}_3 -GCop and \mathcal{M}_3 -SCop are virtually indistinguishable. However, the Student copula also depends of an additional parameter controlling the strength of asymptotic dependence. The posterior pdf of this parameter is shown in Figure 7a, and corresponds to a rather limited tail dependence.

Stronger differences are observed for the dependence between regression errors (Figure 7b-c). In particular, the dependogram for \mathcal{M}_3 -GCop drops faster to zero for location regression errors (Figure 7b). Moreover, the choice of the data dependence model appears to impact the uncertainty in the estimation of the regression errors correlograms (for both location and scale regressions, Figure 7b-c): \mathcal{M}_3 -Inde yields wider posterior intervals than \mathcal{M}_3 -SCop and \mathcal{M}_3 -GCop.

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614 **4.4.3.** Prediction

Results described in previous sections 4.4.1 and 4.4.2 show that the impact of the model for data dependence is quite complex. A more integrated assessment can be made by comparing the prediction of 100-year quantiles obtained with the three models. Figure 8 maps the relative difference (in percent) between (i) \mathcal{M}_3 -Inde (top panels) and \mathcal{M}_3 -SCop (bottom panels) estimates in the one hand; and (ii) \mathcal{M}_3 -GCop estimates in the other hand. The Gaussian Copula is therefore considered as a benchmark model, and the two other dependence assumptions are compared to this benchmark.

Overall, assuming spatial independence between data yields only minor change in $R_{0.99}$ 622 estimates (top left panel). Exceptions are located in the Cevennes ($R_{0.99}$ values are \approx 623 10-15% higher), and on the foothill of the Alps ($R_{0.99}$ values are $\approx 10-35\%$ higher in the 624 North, and $\approx 10-35\%$ lower in the South). Simily, changes in the uncertainty of $R_{0.99}$ 625 estimates are minor (top right panel). However, uncertainty reductions dominate, with 626 decreases in the range 0-25%. This is consistent with the expected behavior when spatial 627 dependence is ignored (see discussion in section 2.1.2): the variance of estimates may be 628 underestimated, i.e. this decrease in uncertainty may be unduly optimistic. Note that 629 once again, the foothill of the Alps is an exception to this overall decrease. 630

⁶³¹ Overall, replacing the Gaussian copula model by a Student copula also yields minor change ⁶³² in $R_{0.99}$ estimates (bottom left panel), except in the Alps region where strong increases are ⁶³³ observed ($\approx 40\%$, culminating at $\approx 100\%$ for a couple of pixels). In terms of uncertainty ⁶³⁴ (bottom right panel), increases now dominate, but are still moderate (in the range 0-20%), ⁶³⁵ with the Alps showing an opposite behavior (uncertainty reductions in the range 20-40%).

4.5. Impact of the regression model

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In this section, the impact of the regression model is evaluated. Inference is performed with the three regression models described in section 4.2.2, coupled with the Gaussian copula model for data dependence. The three resulting models are noted \mathcal{M}_1 -GCop (scale invariance assumption, no elevation effect), \mathcal{M}_2 -GCop (scale invariance assumption with elevation effect) and \mathcal{M}_3 -GCop (elevation effect, scale invariance is not assumed).

⁶⁴¹ 4.5.1. Parameter estimates

The assumption of scale invariance notably impacts the shape parameter estimates 642 Figure 9a), with a marked shift between scale-invariant models \mathcal{M}_1 -GCop and \mathcal{M}_2 -GCop 643 in the one hand, and model \mathcal{M}_3 -GCop in the other hand. Conversely, the completeness 644 of the regression model seems to be the primary factor of influence for the hyper-standard 645 deviations (Figure 9b-c): model \mathcal{M}_1 -GCop, which ignores the elevation effect, yields a 646 markedly higher hyper-standard deviation than models \mathcal{M}_2 -GCop and \mathcal{M}_3 -GCop, which 647 use elevation as covariate. This is an expected result, since improving the regression model 648 results in decreasing the standard deviations of regression errors. Note that in Figure 9b-649 c, the posterior pdfs for \mathcal{M}_1 -GCop and \mathcal{M}_2 -GCop are related to the regression for the 650 index flood parameter $\delta(s)$, as defined in equation (21). Since this regression acts on both 651 the location and scale D-parameters (see Equation (20)), those pdfs are repeated in both 652 panels b and c. 653

⁶⁵⁴ 4.5.2. Data and parameter dependences

Figure 10a shows that data dependence is virtually identical with all regression models. This is consequence of using a copula formalism, with dependence being modeled independently of marginal distributions. Since the regression model only applies on marginal distributions, it does not impact the estimation of dependence between observed data.
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Conversely, the dependence between regression errors is strongly impacted by the choice 659 of a regression model (Figure 10b-c). The amount of dependence is markedly higher with 660 simple model \mathcal{M}_1 -GCop because the whole spatial variability in the distribution of an-661 nual maxima is accounted for by the regression errors. Models \mathcal{M}_2 -GCop and \mathcal{M}_3 -GCop 662 yield lower dependences, suggesting that the elevation effect explains a major part of 663 the spatial variability. This result is consistent with the findings of Kjeldsen and Jones 66 [2009a] who observed a decrease in the dependence between regression errors when the 665 regression model is improved. Lastly, model \mathcal{M}_3 -GCop yield higher uncertainties than 666 model \mathcal{M}_2 -GCop. This might be due to the estimation of two distinct regression models 667 for location and scale D-parameters in model \mathcal{M}_3 -GCop, while model \mathcal{M}_2 -GCop uses a 668 single regression acting on the index flood parameter $\delta(s)$. 669

670 **4.5.3.** Prediction

⁶⁷¹ Maps of 100-year quantiles are compared in a similar way to section 4.4.3. Figure 11 ⁶⁷² maps the relative difference (in percent) between (i) \mathcal{M}_1 -Gcop (top panels) and \mathcal{M}_2 -⁶⁷³ GCop (bottom panels) estimates in the one hand; and (ii) \mathcal{M}_3 -GCop estimates in the ⁶⁷⁴ other hand. The regression model \mathcal{M}_3 is therefore considered as a benchmark model, and ⁶⁷⁵ the two other regression models are compared to this benchmark.

The simple index-flood model \mathcal{M}_1 (which ignores the elevation effect) yields marked differences with the benchmark model \mathcal{M}_3 in terms of $R_{0.99}$ values (upper left panel): higher values are observed in the Cevennes (up to $\approx +100\%$) and in the Alps (up to $\approx +30\%$), while lower values are observed in the south-eastern corner of the domain (up to $\approx -$ 50%). Marked differences also appear for the uncertainty in $R_{0.99}$ estimates (upper right panel), with an overall increase exceeding $\approx 50\%$, except in the Alps where the uncer-

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tainty decreases (up to $\approx 50\%$). Note however that the difference between \mathcal{M}_1 and \mathcal{M}_3 682 uncertainties is not uniform and follows and interesting spatial pattern: uncertainties are 683 similar for both models nearby calibration sites, while \mathcal{M}_1 estimates become far more 684 uncertain than \mathcal{M}_3 estimates in poorly-gauged areas. This is a consequence of regression 685 errors showing a significant amount of spatial dependence with model \mathcal{M}_1 (see section 686 4.5.2 and Figure 10b-c). This spatial dependence improves the efficiency of the transfer of 68 information from gauged sites to nearby ungauged sites. However, when moving further 688 away from the calibration sites, this dependence vanishes and the uncertainty becomes 689 primarily controlled by the hyper-standard deviation of regression errors. This hyper-690 standard deviation is markedly higher for model \mathcal{M}_1 (see section 4.5.1 and Figure 9b-c), 691 which explains its higher uncertainty in poorly-gauged areas. 692

⁶⁹³ Qualitatively similar results are found for the index-flood model \mathcal{M}_2 (which includes an ⁶⁹⁴ elevation effect). However, differences with the model \mathcal{M}_3 in terms of $R_{0.99}$ values are ⁶⁹⁵ smaller (in the range $\approx \pm 20\%$, lower left panel). The uncertainty in $R_{0.99}$ estimates is ⁶⁹⁶ also larger with model \mathcal{M}_2 than with model \mathcal{M}_3 (lower right panel), with increases in the ⁶⁹⁷ range $\approx 30\text{-}40\%$ being quite evenly distributed in space.

⁶⁹⁸ 4.5.4. Assessment of the scale invariance assumption

The scale invariance assumption underlying models \mathcal{M}_1 and \mathcal{M}_2 can be appraised by computing the coefficients of variation (CV) of data from each site. As explained in section 4.2.2, scale invariance implies that the CV remains constant throughout the studied area. Figure 12 shows the empirical CV computed on each site (sites are reordered by increasing CV). The dashed horizontal lines show a 90% posterior interval of the CV resulting from model \mathcal{M}_2 -GCop. Numerous empirical CVs are well outside this interval, which casts

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 $_{705}$ doubts on the validity of the scale invariance assumption. Conversely, model \mathcal{M}_3 -GCop

⁷⁰⁶ (vertical bars) is able to track more closely the varying CVs at different sites.

4.6. Validation

The validation of predictions arising from the Bayesian hierarchical framework is per-707 formed by using tools and concepts borrowed from the field of probabilistic forecasting. 708 Such tools are of particular interest when the quantification of uncertainty is a primary 709 concern, which is the case in this paper. In particular, the reliability and the precision of 710 predictions are evaluated: reliability refers to the ability to derive predictive distributions 711 that are consistent with observations from validation sites, while precision refers to the 712 amount of uncertainty in predictions. Both concepts are complementary: predictions can 713 be reliable but not precise (uncertainties are large, but predictions remain consistent with 714 observations) or alternatively precise but not reliable (e.g. due to an underestimation of 715 uncertainties). An ideal predictive framework would yield predictions that are reliable 716 and as precise as possible. 717

⁷¹⁸ 4.6.1. Reliability of predictions

The reliability of predictions can be assessed by comparing annual maxima from the 719 27 validation sites with their predictive distribution (derived using the "ungauged site") 720 procedure described in section 3.4). More precisely, this comparison is based on the 721 predictive QQ-plot used by e.g. Dawid [1984], Gneiting et al. [2007], Laio and Tamea 722 [2007] or Thyer et al. [2009]. Let F_s be the cdf of the predictive distribution at site 723 s, and $\{Y(s,t)\}_{t=1:T_s}$ validation data at site s. Under the assumption that validation 724 data are realizations from the predictive distribution, the *p*-values $\{F_s(Y(s,t))\}_{t=1:T_s}$ are 725 realizations from a uniform distribution on [0;1]. This can be evaluated with a QQ-plot 726

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⁷²⁷ comparing the empirical cdf of p-values with the cdf of a uniform distribution. This ⁷²⁸ comparison can be performed individually for each validation site, or by pooling p-values ⁷²⁹ from all validation sites.

Figure 13 shows the site-specific predictive QQ-plots for each model (first five panels), then 730 compares the pooled-sites-predictive QQ-plots of the five models in a single plot (lower 731 right panel). Overall, the reliability of the predictive distributions appears acceptable 732 with all regression models, although some specific sites show signs of departure from the 733 1:1 line. Moreover, when all sites are pooled together, all models yield nearly-diagonal pp-734 plots (lower right panel), suggesting that all methods provide equally reliable predictions. 735 This result might appear surprising at first sight, given the differences between models 736 highlighted in previous sections 4.4 and 4.5. However, there is no contradiction in this 737 statement: in the framework of probabilistic prediction, several predictive distributions 738 can be equally reliable but distinct from each other. In particular, the fact that the overall 739 reliability is similar does not mean that, for a given site, the predictive distributions will 740 themselves be similar. This is illustrated in Figure 14, where the model \mathcal{M}_2 -GCop yields 741 a predictive distribution markedly different from the other four models. Moreover, the 742 probabilistic predictions may also differ in their precision (also termed "sharpness" in the 743 field of probabilistic prediction [see *Gneiting et al.*, 2007, for further discussion]). 744

⁷⁴⁵ 4.6.2. Precision of predictions

Given that all models were found equally reliable, one would favor the one yielding the most precise predictions. In order to evaluate the predictive precision, Figure 15 shows the posterior coefficient of variation of the estimated 100-year rainfall $R_{0.99}$, as a function of the distance between the validation site and the nearest calibration site. In general,

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⁷⁵⁰ index-flood models \mathcal{M}_1 -GCop and \mathcal{M}_2 -GCop appear less precise than models based on ⁷⁵¹ the regression \mathcal{M}_3 . Also note that model \mathcal{M}_2 -GCop yields very imprecise predictions ⁷⁵² (CV>0.3) for three validation sites, which are all located in the same region (South-⁷⁵³ Eastern Mediterranean coast).

Moreover, an apparent trend suggests that the \mathcal{M}_1 -predictions are more precise for val-754 idation sites located close to a calibration site than for isolated sites. For this model, 755 Figure 15 suggests that if a calibration site is available at about 10 km, the predictive 756 uncertainty is divided by almost two compared with isolated sites. This confirms an 757 observation previously made in section 4.5.3, and illustrates a positive outcome of mod-758 eling spatial dependence between regression errors. On the other hand, Figure 15 also 759 shows that when the regression model is improved (\mathcal{M}_3) , the predictive precision is as 760 good as model \mathcal{M}_1 nearby calibration sites, despite the fact that spatial dependence be-761 tween regression errors is negligible (see section 4.5.2). Moreover, \mathcal{M}_3 -predictions are 762 more precise for isolated sites. This important observation suggests that improving the 763 regression to make errors as spatially independent as possible is the most sensible strategy. 764 765

5. Discussion

The case study described in section 4 demonstrates the feasibility of a Bayesian hierarchical approach to regional frequency analysis. Despite this encouraging preliminary investigation, numerous avenues for improvement can already be identified. This section discusses the main issues to be addressed.

5.1. Treatment of missing data

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In the case study of section 4, calibration sites were selected in order to avoid missing 770 values in the calibration dataset. This was made for the sake of simplicity, but it cannot 771 be considered as an acceptable practical solution. Indeed, most regional analyses make 772 use of at-site data with different lengths. Restricting the calibration dataset to years 773 shared by all sites results in a loss of information. Moreover, data from poorly-gauged 774 sites still yield a substantial information on the at-site distribution. Combining this in-775 formation with the regional information transferred from nearby sites is likely to improve 776 the inference. 777

The treatment of missing data is challenging due to the explicit modeling of spatial de-778 pendence between data. Indeed, the likelihood equation (9) requires writing the joint dis-779 tribution of a (spatial) vector $\tilde{y}(t)$ at any time t. A first possibility would be to consider 780 that the size of vector $\tilde{\boldsymbol{y}}(t)$ varies with t, depending on the available data. Unfortunately, 781 this complicates the implementation of the model and increases its computational cost, 782 due to the manipulation and inversion of the dependence matrix Σ , whose dimension 783 would also vary with t in this case. An alternative and possibly more efficient approach 784 would be to consider models for missing data and data augmentation algorithms [see e.g. 785 Gelman et al., 1995]. 786

5.2. Sensitivity to the model for data dependence

As stressed in section 2.1.2, the use of an elliptical copula to model data dependence corresponds to a parametric assumption on the structure of dependence. As such, it might be an inappropriate model for some data, especially in the case of extreme data [*Mikosch*, 2005]. This is a crucial issue if the dependence model is to be used for estimating rare multivariate events (e.g., Pr(annual maxima from all raingauges larger than 100 mm)).

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The literature suggests that such estimations are highly sensitive to the choice of the dependence model [e.g. *Coles et al.*, 1999; *Renard and Lang*, 2007]. However, the impact of the dependence model on the estimation at ungauged site is not as clear. Indeed, prediction at ungauged site does not directly use the data dependence model (see section 3.4). Yet, the latter may indirectly influence predictions when inappropriate, by inducing bias in parameter estimates and/or inadequate quantification of uncertainties (posterior variance).

The preliminary investigation carried out in section 4.4 suggests that the choice of a 799 data dependence model is of second order importance compared with the choice of an 800 appropriate regression model to describe spatial variability. However, this result is based 801 on a single case study and can therefore not be considered as a generality. Consequently, 802 the sensitivity of predictions to the data dependence model needs to be further evaluated. 803 In particular, the impact of the spatial dependence model might be more pronounced for 804 datasets showing a higher level of asymptotic dependence. Including multivariate extreme 805 models recently proposed in the literature [e.g. Keef et al., 2009; Padoan et al., 2010] would 806 also constitute an improvement. 807

5.3. The role of spatial- and cross-dependence between regression errors

The framework developed in this paper follows the recent work by *Kjeldsen and Jones* [2009a] in assuming that regression errors are spatially dependent. The case study of section 4 illustrates the benefit of this assumption: when the regression model is poor and fails to capture key relationships with spatial covariates (model \mathcal{M}_1), spatial dependence between regression errors is significant and improves the predictive precision nearby calibration sites. However, this spatial dependence quickly vanishes when the regression

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model is improved (models \mathcal{M}_2 and \mathcal{M}_3). This illustrates the relationship between the 814 quality of the regression and the existence of spatial dependence in regression errors: in a 815 nutshell, spatial dependence acts as a surrogate for the spatial variability "missed" by the 816 regression model. This suggests that improving the regression to make regression errors 817 as spatially independent as possible is a better strategy than attempting to refine the 818 description of spatial dependence. However, in cases where no satisfying regression can 819 be uncovered, the explicit modeling of spatial dependence appears beneficial in terms of 820 predictive precision. 821

Moreover, cross-correlation between regression errors related to distinct D-parameters is neglected in this framework. Unfortunately, the case study suggests that such crosscorrelation exists (see section 4.3.1). Its impact on predictions is unclear at this stage and requires further evaluation.

An explicit modeling of cross-correlation was adopted by *Tasker and Stedinger* [1989], 826 in the case of spatially independent regression errors. This paper somehow considers 827 the opposite option: spatial dependence between regression errors is modeled, but their 828 cross-correlation is ignored. Which of these two options should be favored is likely case-829 specific, and depends on the relative strength of spatial- and cross-correlation. Moreover, 830 it is worth noting that both types of dependence can be constrained (to some extent) by 831 improving the model. In the one hand, improving the regression model decreases spatial 832 dependence as discussed above. In the other hand, cross-correlation can be decreased by 833 reparameterizing the marginal distributions: for instance, the GEV distribution could be 834 parameterized in terms of location, CV and shape parameters (instead of location, scale 835 and shape, see the discussion in *Stedinger and Griffis* [2011]), which may limit cross-836

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correlation between location and scale regression errors. Which of these two strategies 837 offers the most flexibility remains an open question, that will be addressed in future work. 838 Lastly, the third strategy would be to explicitly model both spatial- and cross-correlation 839 in regression errors. However, this would require a description of the dependence between 840 spatial random fields, which is a challenging task. Solutions might exist and be borrowed 841 from the field of geostatistics [e.g. using cokriging tools, see *Goovaerts*, 1998]. However, 842 improving the model to make either spatial- or cross-correlation negligible seems easier to 843 implement and might be an acceptable solution in many cases. 844

5.4. Validation procedures

As any predictive statistical model, a model constructed within the hierarchical frame-845 work presented in this paper needs to be thoroughly validated based on data that were 846 not used for estimation. In the case study of section 4, this was achieved by comparing 84 the predictive distribution (derived in an ungauged site context) with validation data. 848 Such a comparison assesses the overall reliability of the predictive distribution, and may 840 detect systematic predictive biases (e.g. the predictive mean is significantly smaller that 850 the mean of validation data). However, it might not be sufficient to assess the reliability 851 of predicted extreme quantiles, which are of primary interest in frequency analysis. Con-852 sequently, more specialized validation procedures are needed to open predicted extreme 853 quantiles to direct validation and scrutiny. The validation tools recently proposed by 854 Garavaglia et al. [2011] might be of particular interest in this respect. 855

5.5. Model complexity and identifiability

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The case study of section 4 compares several models differing in their flexibility (and, 856 as a consequence, in their complexity). In particular, models \mathcal{M}_1 and \mathcal{M}_2 are based on 857 the scale invariance assumption. This limits the complexity of the model by merging all 858 spatial variability into a single D-parameter (the index-flood parameter $\delta(s)$), the other 859 D-parameters being assumed constant over the region. However, this assumption seems 860 unrealistic in the studied region. The alternative model \mathcal{M}_3 describes the spatial varia-861 tions in two D-parameters, namely the location and the scale parameters, while the third 862 D-parameter (shape) remains constant. This is similar to the approach evaluated by Ste-863 dinger and Lu [1995], and it was found to yield reliable and more precise predictions in 864 this case study. Consequently, it would be tempting to go even further in terms of model 865 complexity, by using three distinct regression models for the location, scale and shape 866 parameters. 867

However, as in any statistical model, a trade-off between descriptive and predictive power 868 has to be found. Model complexity may come at the cost of reduced predictive ability. 869 Moreover, non-identifiability issues may arise if the information content of the data does 870 not support the inference of several spatial processes. Establishing an acceptable trade-off 871 between the flexibility of the model and the level of complexity that can be identified from 872 the data is a very challenging task. In particular, it is linked with the preceding discus-873 sion on validation procedures: introducing additional complexity is only beneficial if it 874 demonstrably improves predictive reliability and/or precision, yet the power of validation 875 procedures to detect model failures may be too limited, especially for extreme quantiles. 876 Consequently, in the case study presented in this paper, developing more efficient valida-877 tion tools seems to be a prerequisite before attempting to model spatial variations in the 878

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⁸⁷⁹ shape parameter.

Lastly, it is stressed that the high dimensionality of models resulting from the hierarchical framework presented herein does not necessarily lead to over-parameterized or nonidentifiable models. In the context of Bayesian hierarchical modeling, the issue of model complexity is far from obvious [see *Spiegelhalter et al.*, 2002, for a thorough discussion on this topic] and requires specialized criteria for model selection [e.g. *Cooley et al.*, 2007].

5.6. Prior specification

The case study of section 4 uses mostly non-informative priors. This is because the aim 885 of this case study is to understand some properties of the proposed inference approach, 886 rather than to seek an optimal estimation of rainfall extremes in this particular region. 887 In this context, using precise priors might exert a strong, case-specific leverage on the 888 conclusions. However, the specification of precise and accurate priors might indeed be 889 beneficial to the inference. In particular, the following investigations would be of interest: 890 • Prior specification depends on the particular choices of marginal distributions and 891 regression functions. It is therefore difficult to derive general specification guidelines. 892 In particular, the fact that D-parameters are not inferred directly, but are only defined 893 through the regression function (4) complicates prior specification. Methods for expressing 894 prior knowledge (either based on expertise or on external data not used for inference) in 895 the form of a proper prior distribution need to be developed. 896

• Using conjugate priors has the potential to ease computations by replacing (at least partly) MCMC sampling by explicit formulas. However, the choice of a conjugate family depends on the choice of marginal distributions and regression functions, making it difficult to provide general guidelines. The benefit of using conjugate priors could be

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⁹⁰¹ first evaluated for the hyper-parameters of the Gaussian hyper-distributions, for which ⁹⁰² conjugate families are known [e.g. *Gelman et al.*, 1995].

• The sensitivity of the inference to priors should be further evaluated, especially for ⁹⁰³ complex models that are more prone to identifiability issues (see discussion in section 5.5).

5.7. Modeling variability in time

Since this paper mainly focuses on spatial variability, at-site distributions were assumed 905 to be constant in time (see equation (1)). A natural extension of the framework would 906 be to allow temporal variations in at-site distributions, by including covariates varying 907 with time. The simplest application would be to include a linear trend in one (or possibly 908 several) D-parameter(s). The covariate could be in this case the time itself, or a large-scale 900 climatic index (e.g. NAO or PDO indexes). This would be an important development 910 since the strong natural variability of at-site data (in particular, extreme data) limits the 911 detectability of climatic effects. Studying the impact of climate variability/change at a 912 regional scale is likely to improve this assessment [Renard et al., 2008; Aryal et al., 2009]. 913

5.8. Application to runoff or other hydrological variables

Although the hierarchical approach presented in this paper can in principle be applied to runoff data, it is likely to yield sub-optimal predictions if standard dependence models are used. This is because runoff data are structured by the hydrologic network. Consequently, euclidean distances (e.g. between gauging stations or between catchment centroids) are in general a sub-optimal predictor of intersite dependences. Moreover, predictions should be constrained to ensure the consistency between estimates at upstream and downstream nested catchments. Several authors proposed specialized approaches to model network

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dependences [e.g. *Gottschalk*, 1993a, b; *Sauquet et al.*, 2000; *Sauquet*, 2006; *Skoien et al.*, 2006]. The applicability of such approaches within the Bayesian hierarchical framework presented in this paper will be evaluated in future work.

More generally, the choice of a relevant distance to explain spatial dependences is a chal-924 lenging task, but is also a promising avenue to improve predictions. This choice heavily 925 depends on the spatial properties of the hydrological variable. For rainfall, a xy-Euclidean 926 distance (as used in the case study of section 4) may seem reasonable at first sight, al-927 though accounting for elevation may improve predictions. Alternatively, consider apply-928 ing the hierarchical approach presented in this paper to snow depth values: a meaningful 929 dependence-distance relationship is very unlikely to be derived if elevation is neglected in 930 the definition of the inter-site distance. Recent work by Blanchet et al. [Blanchet et al., 931 2009; Blanchet and Lehning, 2010; Blanchet and Davison, 2011] addresses this issue, and 932 proposes practical solutions to derive meaningful distances. Those solutions might be 933 extended to hydrological variables other than snow depth. 934

6. Conclusion

Regional frequency analysis has a long history in Hydrology, and is widely applied 935 in practice to estimate the distribution of a hydrologic variable at ungauged or poorly 936 gauged sites. However, despite numerous methodological developments over the years 937 [e.g. Stedinger, 1983; Stedinger and Tasker, 1985, 1986; Hosking and Wallis, 1988; Mad-938 sen and Rosbjerg, 1997; Reis et al., 2005; Ribatet et al., 2007; Kjeldsen and Jones, 2009a], 939 most implementations still rely on several hypotheses that complicate the quantification 940 of predictive uncertainty; moreover, unrealistic hypotheses may yield unreliable predic-941 tions. The objective of this paper was therefore to propose a general Bayesian hierarchical 942

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⁹⁴³ framework to overcome some of these limitations, and to evaluate its applicability based ⁹⁴⁴ on a case study. This framework builds on previous work by several authors [in particular, ⁹⁴⁵ Wikle et al., 1998; Perreault, 2000; Micevski et al., 2006; Micevski, 2007; Cooley et al., ⁹⁴⁶ 2007; Lima and Lall, 2009; Aryal et al., 2009; Lima and Lall, 2010], who explored the ⁹⁴⁷ usefulness of spatial and temporal Bayesian hierarchical models. The main features of the ⁹⁴⁸ proposed framework are the following:

⁹⁴⁹ 1. At-site data can be modeled with any distribution, with the sole requirement that ⁹⁵⁰ the same distribution (but with different parameters) is used for all sites.

⁹⁵¹ 2. Intersite dependence is explicitly modeled by means of an elliptical copula.

3. The variation of parameters in space is described with a regression model linking
 parameter values and covariates.

4. Regression errors are modeled by means of a Gaussian spatial field, which allows
 transferring estimates from gauged to ungauged sites, while quantifying the associated
 predictive uncertainty.

A case study based on extreme rainfall data in Mediterranean France demonstrated the applicability of the framework and its reliable estimation of predictive uncertainty. Although numerous improvements remain to be implemented (see discussion section), these encouraging results warrant further research to develop this framework. In particular, the inclusion of temporal covariates would allow describing the evolution of hydrologic variables in both space and time. Importantly, this improved understanding of hydrologic variability would be achieved with a rigorous quantification of the associated uncertainties.

Appendix A: The Gaussian and Student Elliptical copulas

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- ⁹⁶⁴ The following notation is used:
- 1. $F_1(y), ..., F_H(y)$ are the marginal cdfs of a *H*-dimensional random vector $(Y_1, ..., Y_H)$.
- ⁹⁶⁶ 2. $f_1(y), ..., f_H(y)$ are the corresponding marginal pdfs.
- $_{967}$ 3. $\phi(y)$ is the cdf of the standard normal distribution.
- $_{968}$ 4. $\psi(y)$ is the corresponding pdf.
- 5. $\tau_{\nu}(y)$ is the cdf of the Student's t distribution with ν degrees of freedom
- 970 6. $t_{\nu}(y)$ is the corresponding pdf
- $\Phi_{\Sigma}(y_1, ..., y_H)$ is the joint cdf of a *H*-dimensional Gaussian distribution, with mean
- $_{\rm 972}$ $\,$ 0 and covariance matrix Σ
- 973 8. $\Psi_{\Sigma}(y_1, ..., y_H)$ is the corresponding joint pdf
- 974 9. $\Theta_{\Sigma,\nu}(y_1,...,y_H)$ is the joint cdf of a *H*-dimensional Student distribution, with mean
- $_{975}$ 0, covariance matrix Σ and ν degrees of freedom
- ⁹⁷⁶ 10. $T_{\Sigma,\nu}(y_1,...,y_H)$ is the corresponding joint pdf
- ⁹⁷⁷ The Gaussian and the Student copulas build the joint cdf of the random vector ⁹⁷⁸ $(Y_1, ..., Y_H)$ as follows:

$$F_{Gaussian}(y_1, \dots, y_H) = \Phi_{\Sigma}(u_1, \dots, u_H)$$
(A1)

$$F_{Student}(y_1, ..., y_H) = \Theta_{\Sigma, \nu}(v_1, ..., v_H)$$
(A2)

⁹⁷⁹ where $u_i = \phi^{-1}(F_i(y_i))$ and $v_i = \tau_{\nu}^{-1}(F_i(y_i))$.

⁹⁸⁰ The corresponding joint pdfs can be obtained by differentiating the above cdfs:

$$f_{Gaussian}(y_1, ..., y_H) = \frac{(\prod_{i=1}^H f_i(y_i))}{(\prod_{i=1}^H \psi(u_i))} \times \Psi_{\Sigma}(u_1, ..., u_H)$$
(A3)

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$$f_{Student}(y_1, ..., y_H) = \frac{(\prod_{i=1}^H f_i(y_i))}{(\prod_{i=1}^H t_\nu(v_i))} \times T_{\Sigma,\nu}(v_1, ..., v_H)$$
(A4)

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X - 60 RENARD: HIERARCHICAL REGIONAL ESTIMATION

Figure 1. Schematic of the hierarchical modeling framework

Figure 2. Raingauges location. The thin lines represent six homogeneous regions defined by *Pujol et al.* [2007].

Figure 3. Posterior pdfs of some inferred quantities: (a) regional shape parameter ξ ; (b) errors ϵ_{μ} in the regression for the location D-parameter (only 7 distributions are shown for readability); (c) errors ϵ_{λ} in the regression for the scale D-parameter (d) hyper-standard deviations σ_{μ} (solid black line) and σ_{λ} (dashed red line), corresponding to the standard deviations of location/scale regression errors.

Figure 4. Estimated dependence structures. (a) Data dependence. The dots represent estimated pairwise dependences for all available pairs of sites; (b) dependence in regression errors ϵ_{μ} for the location D-parameter; (c) dependence in regression errors ϵ_{λ} for the scale D-parameter.

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Figure 5. Estimation of the 100-year daily rainfall $R_{0.99}$. Left: $R_{0.99}$ point-estimates (posterior median); Right: uncertainty in estimating $R_{0.99}$ (measured by the posterior coefficient of variation).

Figure 6. Posterior pdfs of some inferred quantities with three distinct models for data dependence. (a) Regional shape parameter ξ ; (b) hyper-standard deviations σ_{μ} (standard deviations of location regression errors); (c) hyper-standard deviations σ_{λ} (standard deviations of scale regression errors).

Figure 7. Estimated dependence structures with three distinct models for data dependence. Solid black line = \mathcal{M}_3 -Inde; dashed red line = \mathcal{M}_3 -GCop; dotted blue line = \mathcal{M}_3 -SCop. Shaded areas represent 90% posterior intervals. (a) Data dependence; (b) dependence in regression errors ϵ_{μ} for the location D-parameter; (c) dependence in regression errors ϵ_{λ} for the scale D-parameter.

Figure 8. Impact of the data dependence model on predictions of $R_{0.99}$. Maps show relative differences with the benchmark model \mathcal{M}_3 -GCop. Top panels = \mathcal{M}_3 -Inde; bottom panels = \mathcal{M}_3 -SCop. Left = $R_{0.99}$ point-estimates (posterior median); Right = uncertainty in estimating $R_{0.99}$ (measured by the posterior coefficient of variation).

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Figure 9. Posterior pdfs of some inferred quantities with three distinct regression models. (a) Regional shape parameter ξ ; (b) hyper-standard deviations σ_{μ} (standard deviations of location regression errors); (c) hyper-standard deviations σ_{λ} (standard deviations of scale regression errors).

Figure 10. Estimated dependence structures with three distinct regression models. Dotted blue line = \mathcal{M}_1 -GCop; solid black line = \mathcal{M}_2 -GCop; dashed red line = \mathcal{M}_3 -GCop. Shaded areas represent 90% posterior intervals. (a) Data dependence; (b) dependence in regression errors ϵ_{μ} for the location D-parameter; (c) dependence in regression errors ϵ_{λ} for the scale D-parameter.

Figure 11. Impact of the regression model on predictions of $R_{0.99}$. Maps show relative differences with the benchmark model \mathcal{M}_3 -GCop. Top panels = \mathcal{M}_1 -GCop; bottom panels = \mathcal{M}_2 -GCop. Left = $R_{0.99}$ point-estimates (posterior median); Right = uncertainty in estimating $R_{0.99}$ (measured by the posterior coefficient of variation).

Figure 12. Evaluation of the scale invariance hypothesis. Points = empirical coefficients of variation of data; dashed horizontal lines = coefficient of variation estimated from \mathcal{M}_2 -GCop (90% posterior interval); vertical bars = coefficient of variation estimated from \mathcal{M}_3 -GCop (90% posterior intervals). Figure 13. Predictive QQ-plots at validation sites for each of the five studied models.

The lower right panel corresponds to merging data from all validation sites together.

Figure 14. Comparison of predictive distributions obtained at one particular validation site.

Figure 15. Posterior coefficients of variation of $R_{0.99}$ estimated at validation sites, as a function of the distance between the validation site and the nearest calibration site.









Ν





(a)

(b)

(c)







Posterior CV of R_{0.99}









Relative change in R_{0.99} (%)





Relative change in

R_{0.99} uncertainty (%)


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1.0 1.0 1.0 0.8 0.8 0.8 Pseudo-correlation 0.6 0.6 Correlation 0.6 Correlation 0.4 0.4 0.4 0.2 0.2 0.2 0.0 0.0 0.0 60 50 50 30 60 20 30 40 50 60 0 10 20 30 40 0 10 20 40 0 10

> Distance (km) (b)

Distance (km) (a) Distance (km) (c)



Relative change in R_{0.99} (%)





Relative change in

R_{0.99} uncertainty (%)



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