Applying Max-and-Smooth to the UKCP data

Brynjólfur Gauti Guðrúnar Jónsson

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Introduction

This document describes the implementation of the Max-and-Smooth algorithm for fast approximate Bayesian inference in spatial extreme value analysis of climate projections provided by the UKCP. The algorithm is specifically applied to Generalized Extreme Value (GEV) distributions and is implemented in C++ with R interfaces using Rcpp and RcppEigen as well as Stan.

Package Overview

The maxandsmooth R package provides tools for fast approximate Bayesian inference for spatial GEV models. The core of the package is implemented in C++ for efficiency, with R wrappers for ease of use.

Key features of the package include:

- Implementation of the Max-and-Smooth algorithm with Gaussian copula dependence
- Efficient C++ code using automatic differentiation and Eigen
- Spatial modeling of GEV parameters using Stan's efficient HMC sampler
- R interface for easy integration into existing extreme value analysis workflows

Algorithm Description

The Max-and-Smooth algorithm, as applied to spatial GEV models, consists of two main steps:

- 1. **Max Step**: Maximum likelihood estimation of GEV parameters at each spatial location using C++
- 2. **Smooth Step**: Spatial smoothing of the maximum likelihood estimates using a BYM2 model implemented in Stan

The algorithm treats the ML estimates as sufficient statistics for a latent Gaussian field, providing a fast approximation to full Bayesian inference for spatial extreme value models.

Code Structure

The package is organized into several key files:

- 1. src/gev_reverse_mode.cpp: Implements the Max step (maximum likelihood estimation for GEV) assuming a known Gaussian copula
- 2. Stan/stan smooth bym2.stan Implements the Smooth step using Stan

Max Step

The Max step involves computing location-wise maximum likelihood estimates (MLEs) for the GEV model parameters while accounting for spatial dependence through a Matérn-like Gaussian copula structure.

Data Structure and Model Specification

Let Y be an $n \times p$ matrix of observations where:

- Rows (i = 1, ..., n) represent temporal replicates
- Columns (j = 1, ..., p) represent spatial locations

The model combines GEV marginal distributions with a Gaussian copula:

1. **Marginal GEV distributions**: At each location j, observations follow a GEV distribution:

$$Y_{ij} \sim \text{GEV}(\mu_j, \sigma_j, \xi_j)$$

2. **Spatial dependence**: The dependence structure is captured by transforming the observations to standard normal using the probability integral transform:

$$Z_{ij} = \Phi^{-1}(F_{\mathrm{GEV}}(Y_{ij}|\mu_j,\sigma_j,\xi_j))$$

where $F_{\rm GEV}$ is the GEV CDF and Φ^{-1} is the standard normal quantile function.

3. **Matérn-like precision structure**: The transformed observations follow a multivariate normal distribution with precision matrix:

$$Q=(Q_{\rho_1}\otimes I_{n_2}+I_{n_1}\otimes Q_{\rho_2})^{\nu+1}$$

where:

- Q_{ρ} is the precision matrix of a standardized AR(1) process \otimes denotes the Kronecker product
- ν is a smoothness parameter
- The matrix is scaled to ensure unit marginal variances

Log-likelihood Function

The total log-likelihood combines the GEV marginal contributions and the Gaussian copula:

$$\ell(\theta|Y) = \sum_{i=1}^{p} \sum_{i=1}^{n} \ell_{\text{GEV}}(y_{ij}|\mu_{j}, \sigma_{j}, \xi_{j}) + \ell_{\text{copula}}(Z|Q)$$

where:

1. The GEV log-likelihood for a single observation is:

$$\ell_{\mathrm{GEV}}(y|\mu,\sigma,\xi) = -\log\sigma - (1+\frac{1}{\xi})\log(1+\xi\frac{y-\mu}{\sigma}) - (1+\xi\frac{y-\mu}{\sigma})^{-1/\xi}$$

2. The Gaussian copula log-likelihood is:

$$\ell_{\text{copula}}(Z|Q) = \frac{1}{2}\log|Q| - \frac{1}{2}Z^TQZ + \frac{1}{2}Z^TZ$$

where the last term accounts for the standard normal margins.

Implementation Details

The optimization is performed using automatic differentiation and the L-BFGS algorithm. Key implementation features include:

1. Parameter transformations:

$$(\psi, \tau, \phi) = (\log \mu, \log \sigma - \log \mu, \operatorname{logit}(\xi))$$

- 2. Efficient computation of the quadratic form Z^TQZ and log-determinant $\log |Q|$ by exploiting the Kronecker structure of the precision matrix
- 3. Probability integral transform using accurate approximations to the GEV CDF and normal quantile function
- 4. Automatic differentiation (using autodiff's reverse mode) for accurate gradient and Hessian computation

Smooth Step

The Smooth step performs Bayesian inference on the latent parameter fields using the maximum likelihood estimates from the Max step as noisy observations. We implement this using Stan's efficient Hamiltonian Monte Carlo sampler with a BYM2 (Besag-York-Mollié) spatial model.

Model Structure

Let $\hat{\eta}$ be the vector of maximum likelihood estimates from the Max step, arranged as:

$$\hat{\boldsymbol{\eta}} = (\hat{\psi}_1, \dots, \hat{\psi}_p, \hat{\tau}_1, \dots, \hat{\tau}_p, \hat{\phi}_1, \dots, \hat{\phi}_p)^T$$

where p is the number of spatial locations and $(\hat{\psi}, \hat{\tau}, \hat{\phi})$ represent the transformed GEV parameters.

Spatial Random Effects

For each parameter type $k \in \{\psi, \tau, \phi\}$, we decompose the spatial variation into structured and unstructured components following the BYM2 parameterization:

$$\eta_k = \mu_k \mathbf{1} + \sigma_k \left(\sqrt{\frac{\rho_k}{c}} \eta_k^{\text{spatial}} + \sqrt{1 - \rho_k} \eta_k^{\text{random}} \right)$$

where:

- μ_k is the overall mean
- $\sigma_k > 0$ is the marginal standard deviation
- $\rho_k \in [0,1]$ is the mixing parameter controlling the balance between spatial and unstructured variation
- c is a scaling factor that ensures the marginal variance of the spatial component is approximately 1
- η_k^{spatial} follows an intrinsic conditional autoregressive (ICAR) prior
- $\eta_k^{\text{random}} \sim \mathcal{N}(0, I)$ represents unstructured random effects

ICAR Prior Specification

The ICAR prior for the spatial component η_k^{spatial} is specified through its full conditional distributions:

$$\eta_{k,i}^{\text{spatial}} | \eta_{k,-i}^{\text{spatial}} \sim \mathcal{N}\left(\frac{1}{n_i} \sum_{j \sim i} \eta_{k,j}^{\text{spatial}}, \frac{1}{n_i}\right)$$

where $j \sim i$ indicates that locations i and j are neighbors, and n_i is the number of neighbors for location i. This is implemented efficiently in Stan through the sum of squared differences form:

$$p(\eta_k^{\text{spatial}}) \propto \exp\left(-\frac{1}{2} \sum_{i \sim j} (\eta_{k,i}^{\text{spatial}} - \eta_{k,j}^{\text{spatial}})^2\right)$$

with an additional soft sum-to-zero constraint implemented via $\sum_i \eta_{k,i}^{\text{spatial}} \sim \mathcal{N}(0, 0.001p)$.

Observation Model

The observation model links the MLEs to the latent field through a multivariate normal distribution:

$$\hat{\eta}|\eta \sim \mathcal{N}(\eta,Q_{\eta y}^{-1})$$

where $Q_{\eta y}$ is the precision matrix obtained from the negative Hessian in the Max step. To handle this efficiently in Stan, we:

- 1. Pre-compute the Cholesky factor L of $Q_{\eta y} = LL^T$
- 2. Store L in a sparse format using arrays of indices and values
- 3. Implement a custom log probability function that computes:

$$\log p(\hat{\eta}|\eta) = \frac{1}{2}\log|Q_{\eta y}| - \frac{1}{2}(\hat{\eta} - \eta)^TQ_{\eta y}(\hat{\eta} - \eta)$$

using the sparse Cholesky representation

Prior Distributions

We specify weakly informative priors:

$$\begin{split} & \sigma_k \sim \text{Exponential}(1) \\ & \rho_k \sim \text{Beta}(1,1) \\ & \mu_k \sim \text{flat} \end{split}$$

for each parameter type k. The exponential prior on σ_k provides weak regularization while ensuring positivity, while the uniform Beta prior on ρ_k allows the data to determine the balance between spatial and unstructured variation.

Posterior Inference

The model is fit using Stan's implementation of the No-U-Turn Sampler (NUTS), providing:

- 1. Posterior samples of the latent field η
- 2. Uncertainty quantification through the posterior distributions of μ_k , σ_k , and ρ_k 3. Decomposition of spatial variation through the posterior distributions of $\eta_k^{\rm spatial}$ and