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Spatial uncertainty assessment in modelling reference evapotranspiration at regional scale

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Abstract

Evapotranspiration is one of the major components of the water balance and has been identified as a key factor in hydrological modelling. For this reason, several methods have been developed to calculate the reference evapotranspiration (ET_0). In modelling

- reference evapotranspiration it is inevitable that both model and data input will present some uncertainty. Whatever model is used, the errors in the input will propagate to the output of the calculated ET₀. Neglecting information about estimation uncertainty, however, may lead to improper decision-making and water resources management. One geostatistical approach to spatial analysis is stochastic simulation, which draws alternative and equally probable, realizations of a regionalized variable. Differences between the realizations provide a measure of spatial uncertainty and allow to carry out
- an error propagation analysis. Among the evapotranspiration models, the Hargreaves-Samani model was used.

The aim of this paper was to assess spatial uncertainty of a monthly reference evapotranspiration model resulting from the uncertainties in the input attributes (mainly temperature) at regional scale. A case study was presented for Calabria region (southern Italy). Temperature data were jointly simulated by conditional turning bands simulation with elevation as external drift and 500 realizations were generated.

The ET_0 was then estimated for each set of the 500 realizations of the input variables, and the ensemble of the model outputs was used to infer the reference evapotranspiration probability distribution function. This approach allowed to delineate the areas characterized by greater uncertainty, to improve supplementary sampling strategies and ET_0 value predictions.





1 Introduction

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Reference evapotranspiration (ET_0) , defined as the potential evapotranspiration of a hypothetical surface of green grass of uniform height, actively growing and adequately watered, is one of the most important hydrological variables for scheduling irrigation systems, preparing input data for hydrological water-balance models, and calculating actual evapotranspiration for a region and/or a basin (Blaney and Criddle, 1950; Dyck, 1983; Hobbins et al., 2001a, b; Xu and Li, 2003; Xu and Singh, 2005; Gong et al., 2006).

The concept of reference evapotranspiration was introduced to study the evaporative demand of the atmosphere independently of crop type, crop development and management practices. As water is abundantly available at the reference evapotranspiring surface, soil factors do not affect ET_0 . The only factors affecting ET_0 are climatic attributes. Consequently, ET_0 is a climatic attribute and can be computed from weather data. ET_0 expresses the evaporating power of the atmosphere at a specific location and time of the year and does not consider the crop characteristics and soil factors (Allen et al., 1998).

A multitude of methods exists to estimate reference evapotranspiration, ET_0 , (e.g., Xu and Singh, 2002). The techniques for estimating ET_0 are based on one, or more, atmospheric variables, such as air temperature, solar or net total radiation and humidity,

or some measurement related to these variables, like pan evaporation (ET_{pan}). Some of these methods are accurate and reliable; others provide only a rough approximation. Most of the methods were developed in specific studies and work better when applied to the climate for which they were developed (Penman, 1948; Jensen, 1973).

The Penman-Monteith equation is widely recommended because of its detailed theoretical base and its accommodation of small time periods. However, the detailed climatological data required by the Penman-Monteith are not often available especially in developing nations. This lack of meteorological data was resolved by Hargreaves et al. (1985) who devised an easy approach for calculating ET₀. The Hargreaves equation (Hargreaves and Samani, 1985) requires only daily mean, maximum, and minimum air





temperature, usually available at most weather stations world-wide, and extraterrestrial radiation (Droogers and Allen, 2002). This method behaves best for weekly or longer predictions, although some accurate ET_0 daily estimations have been reported in literature (Hargreaves and Allen, 2003).

To improve the prediction capacity of ET₀ models for large areas, spatial data should be used as inputs because their continuous variation may reflect more appropriately the nature of the ET₀ in comparison to the measurements made only at a few weather station locations. When the input data are sparse or poorly correlated in space, their direct measurements could be supplemented by secondary information originating from other related attributes (Goovaerts, 1997). The estimation generally improves when additional and denser information is taken into consideration.

Evapotranspiration depends mainly on temperature, which in turn is strongly controlled by topographic attributes, longitude, latitude and distance from the coast. Temperature varies both in space and time, and it is generally well correlated with elevation.

Elevation can be regarded as completely known from accurate digital elevation models and can help to map temperature above the ground. A good example in which elevation was used as external drift to model temperature is reported in Hudson and Wackernagel (1994). Moreover, at regional scale, stationarity of temperature data cannot be assumed, so a non-stationary approach as kriging with external drift must be used.

In most quantitative ET_0 modelling with GIS, the calculation is very often assumed to produce an exact result, because most current GIS are intrinsically deterministic and cannot examine the impact of errors in input and output data. Knowing the quality of the model results is fundamental especially when they are used in spatial decision-making

and water management. The quality of model predictions essentially depends on three main factors: 1) quality of data; 2) quality of the model and 3) the way data and model interact (Burrough, 2001). Therefore it is very important to know how uncertainties in both model parameter and data propagate through the model. Model uncertainty is caused by: 1) the limitations in the mathematical models used to simulate the physical





system imposed by the simplifying assumptions and/or 2) parameter errors in regression models. Model uncertainty is generally difficult to quantify and a way of estimating the errors is to use retrospective validation with independent data. Data uncertainty is instead caused by measurement errors in the data, incomplete knowledge of spatial

- and temporal variations and heterogeneities at a spatial scale smaller than the sampling scale. To assess the model output error resulting from the uncertainties in the input attributes, a Monte Carlo analysis (Heuvelink, 1998) can be used, an analysis which consists in the generation of an adequate random input data set realizations and considers the joint distribution of all input variables. The model is then run for each
- single set of realizations of the input variables and the ensemble of model outputs is used to infer the output probability function. A single Monte Carlo simulation consists in model running at all locations of a fine grid covering the interest region. The simplest way to store error surfaces of interpolated input data in a GIS is to assume that all data are normally distributed and, then, that the error is correctly expressed by the standard
- ¹⁵ deviation. This method entails associating two numbers to each cell: the mean value and its standard deviation. However, a criticism to this approach is that spatial correlation is neglected and the spatially uncorrelated error is used for each realization. An alternative method consists in applying joint multivariate stochastic simulation (Gomez-Hernandez and Journel, 1992; Goovaerts, 1997), which is aimed at making predictions
- of cross-correlated variables. Such prediction is accomplished using a variogram matrix, which includes not only spatial autocorrelation, but also spatial cross-correlation between variables. The latter information is expected to improve the spatial prediction of ET₀ by reducing its uncertainty, when compared with traditional Monte Carlo simulations. Stochastic simulation actually allows to estimate cell-specific probability
- ²⁵ distribution functions which reflect the location of known data points and the spatial correlation structures of the variables. Although the stochastic simulation method is computer intensive, it has several advantages: unbiased predictions of model outputs; estimates of output uncertainties; assessment of error propagation in non-linear and complex models; estimation of probability of exceeding a critical threshold.





The objective of this paper was to assess spatial uncertainty of a monthly reference evapotranspiration model resulting from the uncertainties in the input attributes (mainly temperature data) at regional scale (Calabria region, southern Italy). In this study, it was focused on June mean reference evapotranspiration, and a subset of elevation as external drift variable at points on a 250-m square grid was used for simulation.

2 Materials and methods

The study case was the Calabria region located in the southern part of the Italian peninsula (Fig. 1) with an area of 15 080 km² and a coastline of 738 km on the Ionian and Tyrrhenian seas. In the North, it borders Basilicata region for 80 km. Calabria has an oblong shape with a length of 248 km, and a width ranging between 31 and 111 km. Although Calabria does not have many high summits, it is one of the most mountainous regions in Italy (Fig. 1): 42% of the land is mountainous, 49% hilly, and only 9% is flat. The maximum elevation is 2267 m a.s.l., while the average elevation is 597 m a.s.l.

Temperatures have been measured daily at 134 weather stations of the Italian Hydrographic Service (at present "Centro Funzionale Multirischi della Calabria" of the "Agenzia Regionale per la Protezione dell'Ambiente della Calabria" – Arpacal) during the period 1924–2009 (Fig. 1). Since it was focused on the month of June, only temperature data of this month were taken into account. Temperature time series having less than 30 years of observation were discarded and only temperature data from 42 weather stations were used. Some external stations were used to take into account the border effect. To map June temperatures, a subset of elevation at points on a 250-

m square grid from accurate digital elevation model was used in simulation as external drift variable.

The reference evapotranspiration was computed using the Hargreaves equation (Hargreaves and Samani, 1985), which can be expressed as:

$$\text{ET}_0 = 0.0023 R_a (T + 17.8) \sqrt{(T_{\text{max}} - T_{\text{min}})}.$$

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(1)

where ET_0 is the computed reference evapotranspiration (mm d⁻¹); R_a is the water equivalent of the extraterrestrial radiation (mm d⁻¹) computed according to Allen et al. (1998); T_{max} , T_{min} and T are the daily maximum, minimum and mean air temperature (°C), with T calculated as the average of T_{max} and T_{min} . 0.0023 is the original empirical coefficient proposed by Hargreaves and Samani (1985). The monthly mean reference evapotranspiration was obtained multiplying the result by 30 days.

Stochastic simulation of the input attributes

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In this paragraph only a very brief introduction to the algorithm used in the case study will be given; for a detailed presentation, interested readers should refer to ¹⁰ Goovaerts (1997), Chilès and Delfiner (1999), among others. Most geostatistics is based on the concept of a random function $Z(\mathbf{x})$, whereby the set of unknown values is regarded as a set of spatially dependent random variables $Z(\mathbf{x}_{\alpha})$. Each measurement of air temperature, $z(\mathbf{x}_{\alpha})$, at different location \mathbf{x}_{α} (\mathbf{x} is the location coordinates vector and α the sampling points = 1, ..., n) is interpreted as a particular realization of a random variable $Z(\mathbf{x}_{\alpha})$.

Geostatistical simulation, compared with an optimal procedure of estimation such as kriging, provides a more realistic means of evaluating the spatial variability of a variable (Castrignanò and Buttafuoco, 2004). Stochastic simulation results in a large number of equiprobable images, also called realizations, which honour the sample data and reproduce statistical characteristics and spatial features. Two types of simulations are available using geostatistics: unconditional and conditional. Unconditional simulations simply reproduce certain statistical measures (mean, variance, covariance function) of a variable without considering the observed data. Conditional simulations generate realizations that incorporate the correlation structure of the data, honour the data and reproduce some random component of variation, which is smoothed out by kriging.

There are several simulation techniques: the one used in this work is the turning bands method with external drift, which represents a reasonable trade-off between





quality and computing time. This method was chosen because the weather stations having temperature data with more than 30 years of observations were sparse and few (only 42). Elevation is additional and denser information easily available, which can improve the estimation. In the scope of simulation with external drift, the variable of interest Z(x) comprises deterministic and stochastic components, then it can represent the combination by the model:

$$Z(\mathbf{x}_{\alpha}) = m(\mathbf{x}) + \varepsilon(\mathbf{x})$$
 and $E[Z(\mathbf{x})] = m(\mathbf{x})$.

where $\varepsilon(x)$ is the stochastic component with zero mean and variogram $\gamma_{\varepsilon}(h)$ and m(x) is the drift which is usually modelled as a linear function of a smoothly varying secondary (external) variable y(x):

 $m(x) = a_0(x) + a_1(x) y(x).$

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The principle is to simulate a target variable using an auxiliary linear correlated variable known at the grid nodes of the result grid file. The auxiliary correlated variable in this case study was elevation because there is a good linear correlation between both

- ¹⁵ June mean minimum (-0.90) and June mean maximum (-0.87) temperature data. The value of the secondary variable (elevation) must be known at all primary data locations x_{α} ($\alpha = 1, ..., n$) and at all locations x_0 being estimated. Moreover, the secondary variable should vary smoothly in space to avoid instability of the kriging with external drift system (Goovaerts, 1997). The simulation algorithm generates a 2-D simulation from
- the 1-D simulations along the lines. The turning bands method is a powerful and useful mathematical operator (Christakos, 1987, 1992), however some authors (Deutsch and Journel, 1998) have criticised it because of the generation of artefacts in the simulated images. These artefacts are for the 3-D cases due to the limitation of a maximum of 15 lines which provide a regular partition of the 3-D space but there is no such limitation
- in 2-D. The turning band method consists in adding up a large number of independent 1-D simulations on the lines partitioning the plane. The value of the simulation at a point of the plane is the sum of the simulated values of the projected points on the different lines. In practice, the 2-D covariance is given and the 1-D covariance is obtained



(2)

(3)



by a deconvolution process. The simulations along each line are discretized so that the same simulated value at a point is assigned to the "band" perpendicular to the line and containing the point. Hence, the name turning "bands" given to the method. The only parameter of this method is the count of bands that has been fixed to 400 in this work. This was a good compromise to save computer time and obtain good results.

Moreover, the number of realizations was fixed to 500 because high accuracies are reached only when the number of runs is sufficiently large.

The previous steps generate non-conditional realizations, which reproduce the given covariance function but do not honour the data. Conditioning is implemented in the software ISATIS[®] release 10.03 (www.geovariances.com) by kriging (Journel and Huijbregts, 1978; Chiles and Delfiner, 1999). The conditional simulation at location x_0 is given by:

$$Z_{\rm cs}(\boldsymbol{x}_0) = Z_{\rm s}(\boldsymbol{x}_0) + \sum_{L=1}^{N} \lambda_i^0 [Z(\boldsymbol{x}_i) - Z_{\rm s}(\boldsymbol{x}_i)].$$
(4)

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where $Z_{cs}(x_0)$ is the conditional simulation at x_0 ; $Z_s(x_0)$ the non-conditional simulation at x_0 ; $z(x_i)$ the experimental value at experimental location x_i ; $Z_s(x_c)$ the nonconditional simulation at experimental locations x_i ; λ_i^0 the kriging weight assigned at experimental location x_i when estimating at location x_0 ; and n the number of experimental locations for kriging. As the turning bands method is a Gaussian simulation technique, it requires a multi-Gaussian framework. Therefore, each variable has initially been transformed into a normal distribution with zero mean and unit variance, and the simulation results have subsequently been back-transformed to the raw distribution. Such procedure is known as Gaussian anamorphosis (Chilès and Delfiner, 1999; Wackernagel, 2003), and it is a mathematical function which transforms a variable with a Gaussian distribution into a new variable with any distribution. The Gaussian

²⁵ anamorphosis can be achieved using an expansion into Hermite polynomials $H_i(Y)$ (Wackernagel, 2003) restricted to a finite number of terms.





The joint turning bands simulation requires modelling and fitting a linear model of coregionalization (Goovaerts, 1997). The Linear Model of Coregionalization (LMC) is a quantitative measure of spatial correlation of the regionalized variable $z(x_i)$. It provides a method for modelling the direct and cross-variogram(s) of two or more variables so that the variables of any people linear combination of these variables is people as the time of the second second

⁵ that the variance of any possible linear combination of these variables is positive. Any experimental variogram is modelled as a combination of the same basic structures.

The aim was to build a model which described the major spatial features of the attributes under study. The models used can represent bounded or unbounded variation. In the former models, the variance (known as the sill variance) has a maximum at a finite lag distance (range) over which pairs of values are spatially correlated. In this case, to model the coregionalization of the two attributes (T_{max} and T_{min}), three (N(N + 1)/2) direct and cross variograms must be calculated and modelled jointly for the anamorphosed temperature data.

The best fitting function can be chosen by cross-validation, which checks the compatibility between the data and the model. It takes each data point in turn, removing it temporarily from the data set and using its neighbouring information to predict the value of the variable at its location. The estimate is compared with the measured value by calculating the experimental error, i.e. the difference between estimate and measurement, which can also be standardized by estimating the standard deviation. The goodness of fit was evaluated by the mean error (ME) and mean squared deviation (MSDR). The mean error, which proves the unbiasedness of estimate if its value is close to 0, is given by:

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$$\mathsf{ME} = \frac{1}{N} \sum_{i=1}^{N} [z^*(\boldsymbol{x}_i) - z(\boldsymbol{x}_i)].$$
(5)

where *N* is the number of observation points, $z^*(x_i)$ is the predicted value at location *i*, and $z(x_i)$ is the observed value at location *i*. The mean squared deviation ratio MSDR, which is the ratio between the squared errors and the kriging variance ($\sigma^2(x_i)$)





is expressed as:

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MSDR=
$$\frac{1}{N} \sum_{i=1}^{N} \frac{[z^{*}(x_{i})-z(x_{i})]^{2}}{\sigma^{2}(x_{i})}.$$

If the model for the variogram is accurate, the mean squared error should equal the kriging variance and the MSDR value should be 1.

Then each geostatistical multivariate simulation has been used as input to the ET₀ model. Probabilistic information has been extracted from the set of simulated images. By averaging the simulated values at each cell, two different maps have been produced: the map of the expected value at any given location (E-type or Expected-value estimate; Journel, 1983) and the one of its standard deviation. The uncertainty in model predictions has been quantitatively evaluated from the replicate stochastic images.

In sum, the proposed approach consisted in the following steps:

- 1. generating a set of input attributes (T_{min} and T_{max}) realizations a_i (i = 1, ..., 500) at nodes of a 250-m square grid using the joint stochastic simulation with elevation as external drift;
- for this set of inputs realizations a_i, computing the reference evapotranspiration using Eq. (1);
- 3. for each input and output attributes, computing average and standard deviation of the simulated values at each cell to produce the maps of the expected values at any given location and the ones of their standard deviation.



(6)

3 Results and discussion

The summary statistics of mean maximum (T_{max}) and mean minimum temperature (T_{min}) data (Celsius degrees) for June and elevation data (m a.s.l.). are reported in Table 1. The assumption of normal distribution was not accepted for both mean maximum (T_{max}) and mean minimum temperature (T_{min}) data at a probability level p > 0.10, and the data distributions showed long negative tails. Therefore, before conducting joint Gaussian simulation, we applied a Gaussian transformation to T_{max} and T_{min} data. No anisotropy was evident in the maps of the 2-D variograms (not shown) to a maximum lag distance of 100 km. A nested isotropic LMC (Table 2) was fitted to all the experimental direct and cross-variograms of the Gaussian transformed variables. The LMC (Table 2) includes three different structures: a nugget effect, an exponential structure with a practical range of 30 000 m, and a spherical structure with a range of 60 000 m. The goodness of fit was evaluated by a cross-validation test and the results in terms of mean error (ME) and mean squared deviation (MSDR). ME ranged between -0.02

- and -0.07, while MSDR between 0.92 and 0.93. The optimal values for the two statistics are 0 for ME and 1 for MSDR, then the multivariate model of spatial correlation was unbiased and reproduced the experimental variance adequately. The sum of the eigenvalues at each spatial scale provides an estimate of the variance at that scale (Table 2). The nugget was about 50% of total variance (2.18), while the contribution
- of the shorter range component (30 000 m) of variation to the total variance was about 40% and the contribution of the longer range component (60 000 m) was 10%. The nugget effect component represents the unstructured spatial variation. It is mainly due to measurement errors and to the spatial variation at a scale lesser than the minimum distance of sampling. Moreover, in this study, the nugget ratio is due to the sparse and
- ²⁵ limited number of sampling locations. The variation at shorter scale (40%) is probably related to the local orographic characteristics of the region, while the longer scale of variation (10%) could be related to large scale factors.





The above LMC was used to generate the 500 simulations of T_{min} and T_{max} at nodes of a 250-m square grid. Afterwards, the expected values of T_{min} and T_{max} and their standard deviation were computed and mapped: Fig. 2 and Fig. 3 present a way to treat the jointly simulated images of the two variables, by calculating the mean and the standard deviation respectively of the 500 simulations at each grid node, and then mapping the results for each variable. The mean maps (Figs. 2 and 3) show the complexity in spatial distribution of temperatures. The maps of the standard deviation (Figs. 2 and 3), obtained by post-processing the simulations, have allowed to assess the uncertainties of non-Gaussian variables and to overcome the drawback of kriging variance of its independence from actual sample values. From a visual inspection, it shows clearly

- ¹⁰ Independence from actual sample values. From a visual inspection, it shows clearly how the uncertainty distributions of temperature are mostly related to the density of the sample data (Fig. 1). Figures 2 and 3 also show, as expected, that lower values of mean temperatures are estimated in correspondence to the mountainous areas (Sila Massif, Serre Chain and Aspromonte Massif) (Fig. 1), which have also the higher uncertainties (birther values of standard deviation). Figures 2 also show a large standard deviation.
- tainties (higher values of standard deviation). Figure 2 also shows a large area of the region (the northern part) characterized by high values of standard deviation. These high values occurred in an area with high variability in elevation and consequently in temperature data at short distance.

The 500 realizations of temperature data were used as input for ET_0 model and then the maps of the mean and standard deviation of ET_0 (Fig. 4) were obtained in a similar way to the maps of Fig. 2 and Fig. 3. A visual inspection shows clearly where the uncertainty in ET_0 is high. The simulation has shown how the previous uncertainties in input variables can affect the predictions of ET_0 model. In particular, one can note that there are extended areas characterised by high uncertainties localised on the Sila

²⁵ Massif, the Aspromonte Massif, the Serre Chain and the north-western portion of the region. The areas characterised by medium-high values of ET₀ present not very high values of standard deviation, and therefore less uncertainty.

This approach has demonstrated that it is possible to produce maps of uncertainty, which are more useful than the simple extrapolations of estimation points. Of course, it





is important to evaluate how well the model approximates reality, i.e. model uncertainty. If the model has high uncertainty, a difference in model output may not indicate a real change and could thus be meaningless.

4 Conclusions

⁵ The results of a spatial uncertainty analysis have shown that the prediction quality depends on the uncertainties of the data used in the analysis; therefore map makers should convey the accuracy of the maps they produce (Heuvelink, 1998). A complete characterisation of the accuracy of spatial data should also include the spatial correlation of the attributes used for estimation and stored in a GIS. In the past, a single root mean squared error was sufficient to assess spatial accuracy, but now it is no longer sufficient, and much more information should be provided to characterise the quality of a map.

The objective of this study, however, was not to validate the model or assess the errors associated with the model type and coefficients, but rather to evaluate how the variability of inputs affects uncertainty of model prediction. By definition, a model is an approximation of reality and some models describe reality better than others. Therefore, the choice of model plays an important role in error prediction. In this paper, however, it was assumed that an appropriate model was selected and that the model errors were associated only with the spatial variation of the input attributes.

- In order to obtain realistic values of the model output uncertainty, when the model outputs are supposed to be spatially correlated, it is critically important to model and assess spatial correlations of input variables (Heuvelink and Pebesma 1999). Ignoring spatial correlation between input variables, as in the traditional Monte Carlo approach, implies modelling input variables as white noise. In this case, all uncertainty in ET₀ predictions might vanish after mapping with a dense point grid. The required density
- ²⁵ predictions might vanish after mapping with a dense point grid. The required density depends on the estimate precision level, and it is paramount to model spatial correlation correctly to separate input error from model uncertainty. The two types of variation





are quite different. Spatial variation refers to the deterministic variation of ET_0 or a single realization of the input attributes, whereas uncertainty refers to ET_0 distribution for a single point obtained from the ensemble of Monte Carlo simulations. The approach would put more emphasis on the quality of the input data and on how the input uncertainties may have a considerable impact on prediction uncertainty. Looking at the standard deviation map of ET_0 (Fig. 4), only a weak spatial pattern can be distinguished; the errors do not appear correlated with the estimates of ET_0 , but with the density of weather stations.

Finally, it is worth pointing out the consequences of estimation uncertainty in the context of decision-making and water resources management. There is currently some reluctance to perform error recognition, possibly because of the greater analysis required. However, studying uncertainty leads to increased understanding of the roles played by the different input attributes, which also allows to evaluate the relative costs and benefits of using different scenarios.

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e and elevation data (m a.s.l.).				
$T_{\rm max}$	$T_{\rm min}$	Elevation		
26.9	15.8	366.4		
2.5	3.0	386.5		
30.9	19.6	1315.0		
28.7	18.0	516.0		
27.9	17.1	242.0		
26.0	15.2	15.0		
20.1	7.7	2.0		
-1.0	-1.4	1.1		
3.3	4.2	3.4		
	ata (m a T _{max} 26.9 2.5 30.9 28.7 27.9 26.0 20.1 -1.0 3.3	ata (m a.s.l.).		

Table 1. Basic statistics of mean maximum (T_{max}) and mean minimum temperature (T_{min}) data (Celsius degrees) for June and elevation data (m a.s.l.).



Table 2. Fitted linear model of coregionalization of anamorphosed T_{max} (G T_{max}) and T_{min} (G T_{min}). The coregionalization matrices, the eigenvalues, and the corresponding percentage of variance explained by each eigenvector for the three basic structures are reported.

Variable	G T _{max}	G T _{min}		
(1) Nugget effect				
G T _{max}	0.5613			
G T _{min}	0.2752	0.5784		
Eigenvalue:	0.8452	0.2945		
	(74.16%)	(25.84%)		
<i></i>				
(2) Exponential model (Pr. Range = 30000m)				
G T _{max}	0.3108			
G T _{min}	0.3589	0.5168		
Eigenvalue:	0.7872	0.0403		
	(95.13%)	(4.87%)		
(3) Spherical model (Range = 60 000 m)				
G T _{max}	0.2204			
G T _{min}	0.0037	0.0001		
Eigenvalue:	0.2205	0.0000		
	(100.00%)	(0%)		







Fig. 1. Digital elevation model (right) of the study area and location (left) of the weather station (pushpins).







Fig. 2. Mean and standard deviation maps of simulations for the mean minimum temperature $(T_{\min}).$



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Fig. 3. Mean and standard deviation maps of simulations for the mean maximum temperature $(T_{\rm max}).$



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Fig. 4. Mean and standard deviation maps of simulations for ET₀.

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