Non-Gaussian Geostatistical Modeling using 
(skew) t Processes

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Abstract

We propose a new model for regression and dependence analysis when addressing spatial data with possibly heavy tails and an asymmetric marginal distribution. We first propose a stationary process with t marginals obtained through scale mixing of a Gaussian process with an inverse square root process with Gamma marginals. We then generalize this construction by considering a skew-Gaussian process, thus obtaining a process with skew-t marginal distributions. For the proposed (skew) t process we study the second-order and geometrical properties and in the t case, we provide analytic expressions for the bivariate distribution. In an extensive simulation study, we investigate the...
use of the weighted pairwise likelihood as a method of estimation for the \( t \) process. Moreover we compare the performance of the optimal linear predictor of the \( t \) process versus the optimal Gaussian predictor. Finally, the effectiveness of our methodology is illustrated by analyzing a georeferenced dataset on maximum temperatures in Australia.

*Keywords:* Heavy-tailed processes; Hypergeometric functions; Multivariate skew-normal distribution; Scale mixing; Pairwise likelihood.

1 **Introduction**

The geostatistical approach models data coming from a limited number of monitoring stations as a partial realization from a spatial stochastic process (or random field) defined on the continuum space. Gaussian stochastic processes are among the most popular tools for analyzing spatial data because a mean structure and a valid covariance function completely characterized the associated finite dimensional distribution. Additionally, optimal prediction at an unobserved site depends on the knowledge of the covariance function of the process. Unfortunately, in many geostatistical applications, including climatology, oceanography, the environment and the study of natural resources, the Gaussian framework is unrealistic because the observed data have specific features such as negative or positive asymmetry and/or heavy tails.

The focus of this work is on non-Gaussian models for stochastic processes that vary continuously in the euclidean space, even if the proposed methodology can be easily extended to the space-time framework or to the spherical space. In particular, we aim to accommodate heavier tails than the ones induced by Gaussian processes and wish to allow possible asymmetry. In recent years, different approaches have been proposed in order to analyze these kind of data. Transformation of Gaussian (trans-Gaussian) processes is a general method to model non-Gaussian spatial data obtained by applying some nonlinear transformations to the original data (De Oliveira et al., 1997; Allcroft and Glasbey, 2003; De Oliveira, 2006). Then statistical
analyses can be carried out on the transformed data using any techniques available for Gaussian processes. However, it can be difficult to find an adequate nonlinear transformation and some appealing properties of the latent Gaussian process may not be inherited by the transformed process. A flexible trans-Gaussian process based on the Tukey $g - h$ distribution has been proposed in Xua and Genton (2017).

Wallin and Bolin (2015) proposed non-Gaussian processes derived from stochastic partial differential equations to model non-Gaussian spatial data. However this approach is restricted to the Matérn covariance model with integer smoothness parameter and its statistical properties are much less understood than those of the Gaussian process.

The copula framework (Joe, 2014) has been adapted in the spatial context in order to account for possible deviations from the Gaussian distribution. Even though which copula model to use for a given analysis is not generally known a priori, the copula based on the multivariate Gaussian distribution has gained a general consensus (Kazianka and Pilz, 2010; Masarotto and Varin, 2012; Gräler, 2014) since the definition of the multivariate dependence relies again on the specification of the correlation function. However, Gaussian copula could be too restrictive in some cases since it expresses a symmetrical and elliptical dependence.

Convolution of Gaussian and non-Gaussian processes is an appealing strategy for modeling spatial data with skewness. For instance, Zhang and El-Shaarawi (2010) proposed a Gaussian-Half Gaussian convolution in order to construct a process with marginal distributions of the skew-Gaussian type (Azzalini and Capitanio, 2014). Zareifard et al. (2018) developed bayesian inference for the estimation of a process with asymmetric marginal distributions obtained through convolution of Gaussian and Log-Gaussian processes. Mahmoudian (2017) proposed a skew-Gaussian process using the skew-model proposed in Sahu et al. (2003). The resulting process is not mean-square continuous and as a consequence it is not a suitable model for data exhibiting smooth behavior of the realization.

On the other hand, mixing of Gaussian and non-Gaussian processes is a useful strategy for modeling spatial data with heavy tails. For instance, Palacios and Steel (2006) and Zareifard and Khaledi (2013) proposed a (skew) Gaussian-Log-Gaussian
scale mixing approach in order to accommodate the presence of possible outliers for spatial data.

The $t$ distribution is a parametric model that is able to accommodate flexible tail behavior, thus providing robust estimates against extreme data and it has been studied extensively in recent years (Arellano-Valle and Bolfarine, 1995; Lange et al., 1989; Fonseca et al., 2008; Ferrari and Arellano-Valle, 1996; Arellano-Valle et al., 2012). Stochastic processes with marginal $t$ distributions have been introduced in Røislien and Omre (2006), Ma (2009), Ma (2010a) and DeBastiani et al. (2015), but as outlined in Genton and Zhang (2012), these models are not be identifiable when only a single realization is available (which is typically the case for spatial data).

In this paper, we propose a process with marginal $t$ distributions obtained though scale mixing of a standard Gaussian process with an inverse square root process with Gamma marginals. The latter is obtained through a rescaled sum of independent copies of a standard squared Gaussian process. Although this can be viewed as a natural way to define a $t$ process, the associated second-order, geometrical properties and bivariate distribution are somewhat unknown to the best of our knowledge. Some results can be found in Heyde and Leonenko (2005) and Finlay and Seneta (2006). We study the second-order and geometrical properties of the $t$ process and we provide analytic expressions for the correlation and the bivariate distribution. It turns out that both depend on special functions, particularly the Gauss hypergeometric and Appell function of the fourth type (Gradshteyn and Ryzhik, 2007). In addition, the bivariate distribution is not of elliptical type.

We then focus on processes with asymmetric marginal distributions and heavy tails. We first review the skew Gaussian process proposed in Zhang and El-Shaarawi (2010). For this process we provide an explicit expression of the finite dimensional distribution generalizing previous results in Alegría et al. (2017). We then propose a process with marginal distribution of the skew-$t$ type (Azzalini and Capitanio, 2014) obtained through scale mixing of a skew-Gaussian with an inverse square root process with Gamma marginals.

Our proposals for the $t$ and skew-$t$ processes have two main features. First, they allow removal of any problem of identifiability (Genton and Zhang, 2012),
and as a consequence, all the parameters can be estimated using one realization of the process. Second, the \( t \) and skew-\( t \) processes inherit some of the geometrical properties of the underlying Gaussian process. This implies that the mean-square continuity and differentiability of the \( t \) and skew-\( t \) processes can be modeled using suitable parametric correlation models as the Matérn model (Matérn, 1986) or the Generalized Wendland model (Gneiting, 2002; Bevilacqua et al., 2019).

For the \( t \) process estimation we propose the method of weighted pairwise likelihood (Lindsay, 1988; Varin et al., 2011; Bevilacqua and Gaetan, 2015) exploiting the bivariate distribution given in Theorem 2.3. In an extensive simulation study we investigate the performance of the weighted pairwise likelihood (\( wpl \)) method under different scenarios including when the degrees of freedom are supposed to be unknown. We also study the performance of the \( wpl \) estimation by assuming a Gaussian process in the estimation step with correlation function equal to the correlation function of the \( t \) process. It turns out that the Gaussian misspecified \( wpl \) (see Gouriéroux et al. (2017) with the references therein) leads to a less efficient estimator, as expected. However, the method has some computational benefits.

Additionally, we compare the performance of the optimal linear predictor of the \( t \) process with the optimal predictor of the Gaussian process. Finally we apply the proposed methodology by analyzing a real data set of maximum temperature in Australia where, in this case, we consider a \( t \) process defined on a portion of the sphere (used as an approximation of the planet Earth) and use a correlation model depending on the great-circle distance (Gneiting, 2013).

The methodology considered in this paper is implemented in the R package GeoModels (Bevilacqua and Morales-Oñate, 2018). The remainder of the paper is organized as follows. In Section 2 we introduce the \( t \) process, study the second-order and geometrical properties and provide an analytic expression for the bivariate distribution. In Section 3 we first study the finite dimensional distribution of the skew Gaussian process, and then we study the second-order properties of the skew-\( t \) process. In Section 4 we present a simulation study in order to investigate the performance of the (misspecified) \( wpl \) method when estimating the \( t \) process and the performance of the associated optimal linear predictor versus the optimal Gaussian
predictor. In Section 5, we analyze a real data set of maximum temperature in Australia. Finally, in Section 6, we give some conclusions. All the proofs has been deferred to the Appendix.

2 A stochastic process with t marginal distribution

For the rest of the paper, given a process $Q = \{Q(s), s \in A\}$ with $E(Q(s)) = \mu(s)$ and $Var(Q(s)) = \sigma^2$, we denote by $\rho_Q(h) = Corr(Q(s_i), Q(s_j))$ its correlation function, where $h = s_i - s_j$ is the lag separation vector. For any set of distinct points $(s_1, \ldots, s_n)^T$, $n \in \mathbb{N}$, we denote by $Q_{ij} = (Q(s_i), Q(s_j))^T$, $i \neq j$, the bivariate random vector and by $Q = (Q(s_1), \ldots, Q(s_n))^T$ the multivariate random vector. Moreover, we denote with $f_{Q(s)}$ and $F_{Q(s)}$ the marginal probability density function (pdf) and cumulative distribution function (cdf) of $Q(s)$ respectively, with $f_{Q_{ij}}$ the pdf of $Q_{ij}$ and with $f_Q$ the pdf of $Q$. Finally, we denote with $Q^*$ the standardized weakly stationary process, i.e., $Q^*(s) := (Q(s) - \mu(s))/\sigma$.

As outlined in Palacios and Steel (2006), given a positive process $M = \{M(s), s \in A\}$ and an independent standard Gaussian process $G^* = \{G^*(s), s \in A\}$, a general class of non-Gaussian processes with marginal heavy tails can be obtained as scale mixture of $G^*$, i.e. $\mu(s) + \sigma M(s)^{-\frac{1}{2}} G^*(s)$, where $\mu(s)$ is the location dependent mean and $\sigma > 0$ is a scale parameter. A typical parametric specification for the mean is given by $\mu(s) = X(s)^T \beta$ where $X(s) \in \mathbb{R}^k$ is a vector of covariates and $\beta \in \mathbb{R}^k$ but other types of parametric or nonparametric functions can be considered.

Henceforth, we call $G^*$ the ‘parent’ process and with some abuse of notation we set $\rho(h) := \rho_{G^*}(h)$ and $G := G^*$. Our proposal considers a mixing process $W_\nu = \{W_\nu(s), s \in A\}$ with marginal distribution $\Gamma(\nu/2, \nu/2)$ defined as $W_\nu(s) := \sum_{i=1}^\nu G_i(s)^2/\nu$ where $G_i, i = 1, \ldots, \nu$ are independent copies of $G$ with $E(W_\nu(s)) = 1$, $Var(W_\nu(s)) = 2/\nu$ and $\rho_{W_\nu}(h) = \rho^2(h)$ (Bevilacqua et al., 2018). If we consider a process $Y_\nu^* = \{Y_\nu^*(s), s \in A\}$ defined as

$$Y_\nu^*(s) := W_\nu(s)^{-\frac{1}{2}} G(s),$$

then, by construction, $Y_\nu^*$ has the marginal $t$ distribution with $\nu$ degrees of freedom.
with pdf given by:

\[
f_{Y^*_\nu}(y; \nu) = \frac{\Gamma\left(\frac{\nu+1}{2}\right)}{\sqrt{\pi \nu \Gamma\left(\frac{\nu}{2}\right)}} \left(1 + \frac{y^2}{\nu}\right)^{-\frac{\nu+1}{2}}.
\]  

(2.2)

Then, we define the location-scale transformation process \( Y_\nu = \{Y_\nu(s), s \in A\} \) as:

\[
Y_\nu(s) := \mu(s) + \sigma Y^*_\nu(s)
\]

(2.3)

with \( E(Y_\nu(s)) = \mu(s) \) and \( \text{Var}(Y_\nu(s)) = \sigma^2 \nu/(\nu - 2), \nu > 2 \).

**Remark 1:** A possible drawback for the Gamma process \( W_\nu \) is that it is a limited model due to the restrictions to the half-integers for the shape parameter. Actually, in some special cases, it can assume any positive value greater than zero. This feature is intimately related to the infinite divisibility of the squared Gaussian process \( G^2 = \{G^2(s), s \in A\} \) as shown in Krishnaiah and Rao (1961). Characterization of the infinite divisibility of \( G^2 \) has been studied in Vere-Jones (1997), Bapat (1989), Griffiths (1970) and Eisenbaum and Kaspi (2006). In particular Bapat (1989) provides a characterization based on \( \Omega \), the correlation matrix associated with \( \rho(h) \). Specifically, \( \nu > 0 \) if and only if there exists a matrix \( S_n \) such that \( S_n \Omega^{-1} S_n \) is an \( M \)-matrix (Plemmons, 1977), where \( S_n \) is a signature matrix, i.e., a diagonal matrix of size \( n \) with entries either 1 or -1. This condition is satisfied, for instance, by a stationary Gaussian random process \( G \) defined on \( A = \mathbb{R} \) with an exponential correlation function. The \( t \) process \( Y^*_\nu \) inherits this feature with the additional restriction \( \nu > 2 \). This implies that \( Y^*_\nu \) is well defined for \( \nu = 3, 4, \ldots \) and for \( \nu > 2 \) under non-infinite divisibility of \( G^2 \).

**Remark 2:** The finite dimensional distribution of \( Y^*_\nu \) is unknown to the best of our knowledge, but in principle, it can be derived by mixing the multivariate density associated with \( W^\nu_{\nu^2} \) with the multivariate standard Gaussian density. The multivariate Gamma density \( f_{W_\nu} \) was first discussed by Krishnamoorthy and Parthasarathy (1951) and its properties have been studied by different authors (Royen, 2004; Marcus, 2014). In the bivariate case, Vere-Jones (1967) showed that the bivariate Gamma distribution is infinite divisible, i.e., \( \nu > 0 \) in (A.2), irrespective of the correlation function. Note that this is consistent with the characterization given in Bapat (1989) since, given an arbitrary bivariate correlation matrix \( \Omega \), there
exists a matrix $S_2$ such that $S_2\Omega^{-1}S_2$ is a $M$-matrix. In Theorem 2.3 we provide the bivariate distribution of $Y_{\nu}^*$.

Note that, both $W_\nu$ and $G$ in (2.3) are obtained through independent copies of the ‘parent’ Gaussian process with correlation $\rho(h)$. For this reason, henceforth, in some cases, we will call $Y_{\nu}^*$ a standard $t$ process with underlying correlation $\rho(h)$.

In what follows, we make use of the Gauss hypergeometric function defined by (Gradshteyn and Ryzhik, 2007):

$$2F_1(a, b; c; x) = \sum_{k=0}^{\infty} \frac{(a)_k(b)_k x^k}{(c)_k k!}, \quad (2.4)$$

with $(s)_k = \Gamma(s + k)/\Gamma(s)$ for $k \in \mathbb{N} \cup \{0\}$ being the Pochhammer symbol and we consider the restrictions $a > 0$, $b > 0$, $c > 0$ and $x \geq 0$. If $c > a + b$ the radius of convergence of (2.4) is $0 \leq x \leq 1$ and, in particular (2.4) is convergent at $x = 1$ through the identity:

$$2F_1(a, b; c; 1) = \frac{\Gamma(c)\Gamma(c - a - b)}{\Gamma(c - a)\Gamma(c - b)}. \quad (2.5)$$

We also consider the Appell hypergeometric function of the fourth type (Gradshteyn and Ryzhik, 2007) defined as:

$$F_4(a, b; c, c'; w, z) = \sum_{k=0}^{\infty} \sum_{m=0}^{\infty} \frac{(a)_{k+m}(b)_{k+m}w^k z^m}{k!m!(c)_k(c')_m}, \quad |\sqrt{w}| + |\sqrt{z}| < 1.$$  

The following Theorem gives an analytic expression for $\rho_{Y_{\nu}^*}(h)$ in terms of the Gauss hypergeometric function.

**Theorem 2.1.** Let $Y_{\nu}^*$ be a standardized $t$ process with underlying correlation $\rho(h)$. Then:

$$\rho_{Y_{\nu}^*}(h) = \frac{\nu - 2}{2\Gamma^2\left(\frac{\nu}{2}\right)} \left[2F_1\left(\frac{1}{2}, \frac{1}{2}; \nu; \rho^2(h)\right) \rho(h)\right]. \quad (2.6)$$

The following Theorem depicts some features of the $t$ process. It turns out that nice properties such as stationarity, mean-square continuity and degrees of mean-square differentiability can be inherited from the ‘parent’ Gaussian process $G$. Further, the $t$ process has long-range dependence when the ‘parent’ Gaussian process has long-range dependence and this can be achieved when the correlation has some specific features. For instance, the generalized Cauchy (Gneiting and Schlather, 2004; Lim and Teo, 2009) and Dagum (Berg et al., 2008) correlation models can lead to
a Gaussian process with long range dependence. Finally, an appealing and intuitive feature is that the correlation of $Y_\nu^*$ approaches the correlation of $G$ when $\nu \to \infty$.

**Theorem 2.2.** Let $Y_\nu^*, \nu > 2$ be a standardized $t$ process with underlying correlation $\rho(h)$. Then:

a) $Y_\nu^*$ is also weakly stationary;

b) $Y_\nu^*$ is mean-square continuous if and only if $G$ is mean-square continuous;

c) Let $G$ $m$-times mean-square differentiable, for $m = 0, 1, \ldots$

• If $\nu > 2(2m + 1)$ then $Y_\nu^*$ is $m$-times mean-square differentiable;

• If $\nu \leq 2(2m + 1)$ then $Y_\nu^*$ is $(m - k)$-times mean-square differentiable if $2(2m - k + 1) < \nu \leq 2(2m - k + 3)$, for $k = 1, \ldots, m$.

d) $Y_\nu^*$ is a long-range dependent process if and only if $G$ is a long-range dependent process

e) $\rho_{Y_\nu^*}(h) \leq \rho(h)$ and $\lim_{\nu \to \infty} \rho_{Y_\nu^*}(h) = \rho(h)$.

One implication of Theorem (2.2) point c) is that the process $Y_\nu^*$ inherits the mean square differentiability of $G$ under the condition $\nu > 2(2m + 1)$. Otherwise, the mean square differentiability depends on $\nu$. For instance, if $G$ is one time mean square differentiable then $Y_\nu^*$ can be zero or one time differentiable depending if $\nu > 6$ or not.

**Remark 3:** A simplified version of the $t$ process in Equation (2.1), can be obtained assuming $W_\nu(s_i) \perp W_\nu(s_j), i \neq j$. Under this assumption, $Y_\nu^*$ is still a process with $t$ marginal distribution but, in this case, the geometrical properties are not inherited from the ‘parent’ Gaussian process $G$. In particular, it can be shown that the resulting correlation function exhibits a discontinuity at the origin and, as a consequence, the process is not mean-square continuous. A not mean-square continuous version of the $t$ process in Equation (2.1), can be obtained by introducing a nugget effect, i.e., a discontinuity of $\rho_{Y_\nu^*}(h)$ at the origin. This can be easily achieved by replacing $\rho(h)$ in (2.6) with $\rho^*(h) = 1$ if $h = 0$ and $\rho^*(h) = (1 - \tau^2)\rho(h)$ otherwise, where $0 \leq \tau^2 < 1$ represents the underlying nugget effect.
Since the $t$ process inherits some of the geometrical properties of the ‘parent’ Gaussian process, the choice of the covariance function is crucial. Two flexible isotropic models that allow parametrizing in a continuous fashion the mean square differentiability of a Gaussian process and its sample paths are as follows:

1. the Matérn correlation function \(^{[1986]}\)

\[
M_{\alpha,\psi}(h) = \frac{2^{1-\psi}}{\Gamma(\psi)} \left( \frac{|h|}{\alpha} \right)^\psi \mathcal{K}_\psi \left( \frac{|h|}{\alpha} \right),
\]

where $\mathcal{K}_\psi$ is a modified Bessel function of the second kind of order $\psi$. Here, $\alpha > 0$ and $\psi > 0$ guarantee the positive definiteness of the model in any dimension.

2. the Generalized Wendland correlation function \(^{[2002]}\), defined for $\psi > 0$ as:

\[
GW_{\alpha,\psi,\delta}(h) := \begin{cases} 
\int_0^1 u^{(\psi-1)(1-u)^\delta} du & \text{if } |h| < \alpha \\
0 & \text{otherwise}
\end{cases},
\]

and for $\psi = 0$ as:

\[
GW_{\alpha,0,\delta}(h) := \begin{cases} 
(1 - |h|/\alpha)^\delta & |h| < \alpha \\
0 & \text{otherwise}
\end{cases}.
\]

Here $\mathrm{B}(\cdot, \cdot)$ is the Beta function and $\alpha > 0$, $\delta \geq (d+1)/2 + \psi$ guarantee the positive definiteness of the model in $\mathbb{R}^d$.

In particular for a positive integer $k$, the sample paths of a Gaussian process are $k$ times differentiable if and only if $\psi > k$ in the Matérn case \(^{[1999]}\) and if and only if $\psi > k - 1/2$ in the Generalized Wendland case \(^{[2019]}\). Additionally, the Generalized Wendland correlation is compactly supported, an interesting feature from computational point of view \(^{[2013]}\), which is inherited by the $t$ process since $\rho(h) = 0$ implies $\rho_{Y^*_t}(h) = 0$.

In order to illustrate some geometric features of the $t$ process, we first compare the correlation functions of the Gaussian and $t$ processes using an underlying Matérn
model. In Figure 1 (left part) we compare $\rho_{Y^*}(\mathbf{h})$ when $\nu = 3, 7$ with the correlation of the ‘parent’ Gaussian process $\rho(\mathbf{h}) = \mathcal{M}_{1.5, \alpha^*}(\mathbf{h})$ where $\alpha^*$ is chosen such that the practical range is 0.2. It is apparent that when increasing the degrees of freedom $\rho_{Y^*}(\mathbf{h})$ approaches $\rho(\mathbf{h})$ and that the smoothness at the origin of $\rho_{Y^*}(\mathbf{h})$ is inherited by the smoothness of the Gaussian correlation $\rho(\mathbf{h})$ when $\nu = 7$ and if $\nu = 3$ then $\rho_{Y^*}(\mathbf{h})$ is not differentiable at the origin. On the right side of Figure 1 we compare a kernel nonparametric density estimation of a realization of $G$ and a realization of $Y^*_7$ (approximately 10000 location sites in the unit square) using $\rho(\mathbf{h}) = \mathcal{M}_{1.5, \alpha^*}(\mathbf{h})$.

In Figure 2 (a) and (b), we compare, from left to right, two realizations of $G$ with $\rho(\mathbf{h}) = \mathcal{M}_{0.5, \alpha^*}(\mathbf{h})$ and $\rho(\mathbf{h}) = \mathcal{M}_{1.5, \alpha^*}(\mathbf{h})$ where $\alpha^*$ is chosen such that the practical range is 0.2. In this case, the sample paths of $G$ are zero and one times differentiable. From the bottom part of Figure 2 (c) and (d) it can be appreciated that this feature is inherited by the associated realizations of $Y^*_7$.

![Figure 1](image1.png)

![Figure 2](image2.png)

Figure 1: Left part: comparison of $\rho_{Y^*}(\mathbf{h})$, $\nu = 3, 7$ with the correlation $\rho(\mathbf{h})$ of the ‘parent’ Gaussian process $G$ when $\rho(\mathbf{h}) = \mathcal{M}_{1.5, \alpha^*}(\mathbf{h})$ with $\alpha^*$ such that the practical range is 0.2. Right part: a comparison of a nonparametric kernel density estimation of realizations from $G$ and from the $t$ process $Y^*_7$.

We now consider the bivariate random vector associated with $Y^*_\nu$ defined by:

$$Y^*_{\nu,ij} = W_{\nu,ij}^{-\frac{1}{2}} \circ G_{ij}$$

where $\circ$ denotes the Schur product vector. The following Theorem gives the pdf of
Figure 2: Upper part: two realizations of the ‘parent’ Gaussian process $G$ on $[0, 1]^2$ with (a) $\rho(h) = M_{0.5, \alpha^*}(h)$ and (b) $\rho(h) = M_{1.5, \alpha^*}(h)$ (from left to right) with $\alpha^*$ such that the practical range is approximatively 0.2. Bottom part: (c) and (d) associated realizations of the $t$ process $Y^*_7$.

$Y^*_{\nu,ij}$ in terms of the Appell function $F_4$. It can be viewed as a generalization of the generalized bivariate $t$ distribution proposed in Miller (1968).

**Theorem 2.3.** Let $Y^*_{\nu}$, $\nu > 2$ be a standard $t$ process with underlying correlation
where 

\[
\rho(h) \text{ Then:}
\]

\[
f_{Y_{*ij}}(y_i, y_j) = \frac{\nu^\nu l_{ij}^{-(\nu+1)/2} \Gamma^2\left(\frac{\nu+1}{2}\right)}{\pi^{1/2}(1 - \rho^2(h))^{-(\nu+1)/2}} \frac{1}{l_{ij}} F_4\left(\frac{\nu + 1}{2}, \frac{\nu + 1}{2}, \frac{\nu}{2}, \frac{\nu}{2}; \frac{\nu^2 + 1}{2}; \frac{\rho^2(h)}{l_{ij}}, \frac{l_{ij}}{l_{ij}}\right) + \frac{\rho(h)y_i y_j \nu^{\nu+1} l_{ij}^{-\nu/2 - 1}}{2\pi(1 - \rho^2(h))^{-(\nu+1)/2}} F_4\left(\frac{\nu}{2} + 1, \frac{\nu}{2} + 1, \frac{\nu}{2}, \frac{\nu}{2}; \frac{\rho^2(h)}{l_{ij}}, \frac{l_{ij}}{l_{ij}}\right)
\]

(2.10)

where \(l_{ij} = [(y_i^2 + \nu)(y_j^2 + \nu)].\)

\[\]

Figure 3: Contour plots of the bivariate \(t\) distribution (2.10) when \(\rho(h) = 0.2, 0.9\) and \(\nu = 3, 9.\)

**Remark 4:** Note that \(f_{Y_{*ij}}(y_i, y_j)\) is defined for \(\nu > 2\) irrespectively of the correlation function since it is obtained from a bivariate Gamma distribution (see Remark 2). Moreover, when \(\rho(h) = 0\), according to (4.3) and using the identity \(\quad 2F_1(a, b; c'; 0) = 1\), we obtain \(F_4(a, b; c, c'; 0, 0) = 1\), and as a consequence,
\( f_{Y_{\nu ij}}(y_i, y_j) \) can be written as the product of two independent \( t \) random variables with \( \nu \) degrees of freedom. Thus, zero pairwise correlation implies pairwise independence, as in the Gaussian case. Figure (3) show the contour plots of (2.10) when \( \nu = 3, 9 \), and \( \rho(h) = 0.2, 0.9 \). It turns out that the bivariate \( t \) distribution is not elliptical and when increasing \( \nu \) the contour plots tend towards an elliptical form. Finally, the bivariate density of the process \( Y_\nu \) is easily obtained from (2.3):

\[
\frac{1}{\sigma^2} f_{Y_{\nu ij}} \left( \frac{y_i - \mu_i}{\sigma}, \frac{y_j - \mu_j}{\sigma} \right).
\] (2.11)

3 A stochastic process with skew-t marginal distribution

In this section we first review the skew-Gaussian process proposed in Zhang and El-Shaarawi (2010). For this process, we provide an explicit expression for the finite dimensional distribution generalizing previous results in Alegria et al. (2017). Then, using this skew-Gaussian process, we propose a generalization of the \( t \) process \( Y_\nu \) obtaining a new process with marginal distribution of the skew-\( t \) type (Azzalini and Capitanio, 2014).

Following Zhang and El-Shaarawi (2010) a general construction for a process with asymmetric marginal distribution is given by:

\[
U_\eta(s) := g(s) + \eta|X_1(s)| + \omega X_2(s), \quad s \in A \subset \mathbb{R}^d
\] (3.1)

where \( \eta \in \mathbb{R}, \omega > 0 \) and \( X_i i = 1, 2 \) are two independents copies of a process \( X = \{X(s), s \in A\} \) with symmetric marginals. The parameters \( \eta \) and \( \omega \) allow modeling the asymmetry and variance of the process simultaneously.

Zhang and El-Shaarawi (2010) studied the second-order properties of \( U_\eta \) when \( X \equiv G \). In this case, \( U_\eta \) has skew Gaussian marginal distributions (Azzalini and Capitanio, 2014) with pdf given by:

\[
f_{U_\eta(s)}(u) = \frac{2}{(\eta^2 + \omega^2)^{1/2}} \phi \left( \frac{(u - g(s))}{(\eta^2 + \omega^2)^{1/2}} \right) \Phi \left( \frac{\eta(u - g(s))}{\omega(\eta^2 + \omega^2)^{1/2}} \right)
\] (3.2)

with \( \text{E}(U_\eta(s)) = g(s) + \eta(2/\pi)^{1/2}, \text{Var}(U_\eta(s)) = \omega^2 + \eta^2(1 - 2/\pi) \) and with correlation

\[
\rho_\eta(h) = \frac{\text{Cov}(U_\eta(s), U_\eta(s+h))}{\sqrt{\text{Var}(U_\eta(s)) \text{Var}(U_\eta(s+h))}}
\]
function given:

\[ \rho_{\nu}(h) = \frac{2\eta^2}{\pi\omega^2 + \eta^2(\pi - 2)} \left( (1 - \rho^2(h))^{1/2} + \rho(h) \arcsin(\rho(h)) - 1 \right) + \frac{\omega^2\rho(h)}{\rho^2 + \eta^2(1 - 2/\pi)}. \]  

(3.3)

The following theorem generalizes the results in Alegría et al. (2017) and gives an explicit closed-form expression for the pdf of the random vector \( U_\eta \).

**Theorem 3.1.** Let \( U_\eta(s) = g(s) + \eta|X_1(s)| + \omega X_2(s) \) where \( X_i, i = 1, 2 \) are two independent copies of \( G \) the ‘parent’ Gaussian process. Then:

\[ f_{U_\eta}(u) = 2 \sum_{l=1}^{2^{n-1}} \phi_n(u - \alpha; A_l)\Phi_n(c_l; 0, B_l) \]  

(3.4)

where

\[ A_l = \omega^2\Omega + \eta^2\Omega_l \]

\[ c_l = \eta\Omega_l(\omega^2\Omega + \eta^2\Omega_l)^{-1}(u - \alpha) \]

\[ B_l = \Omega_l - \eta^2\Omega_l(\omega^2\Omega + \eta^2\Omega_l)^{-1}\omega^2 \Omega_l \]

\[ \alpha = [g(s)]_{i=1}^n \]

and the \( \Omega_l \)'s are correlation matrices that depend on the correlation matrix \( \Omega \).

Some comments are in order. First, note that \( f_U \) can be viewed as a generalization of the multivariate skew-Gaussian distribution proposed in Azzalini and Dalla-Valle (1996). Second, using Theorem (3.1), it can be easily shown that the consistency conditions given in Mahmoudian (2018) are satisfied. Third, it is apparent that likelihood-based methods for the skew-Gaussian process are impractical from computational point of view even for a relatively small dataset.

To obtain a process with skew-\( t \) marginal distributions (Azzalini and Capitanio, 2014), we replace the process \( G \) in (2.3) with the process \( U_\eta \). Specifically, we consider a process \( S_{\nu,\eta} = \{S_{\nu,\eta}(s), s \in A\} \) defined as

\[ S_{\nu,\eta}(s) := \mu(s) + \sigma W_{\nu}(s)^{-1/2} U_\eta(s) \]  

(3.5)

where \( W_{\nu} \) and \( U_\eta \) are supposed to be independent. In (3.1) we assume \( g(s) = 0 \) and \( \eta^2 + \omega^2 = 1 \). The pdf of the marginal distribution of \( S_{\nu,\eta}^* \) is given by:

\[ f_{S_{\nu,\eta}^*}(g) = 2f_{Y_{\nu}^*}(g; \nu)F_{Y_{\nu}^*}(s) \left( \eta g \sqrt{\frac{\nu + 1}{\nu + g^2}; \nu + 1} \right) \]  

(3.6)
with \( E(S^*_\nu,\eta(s)) = \frac{\sqrt{\Gamma\left(\frac{\nu-1}{2}\right)}\eta}{\sqrt{\pi}} \), and \( Var(S^*_\nu,\eta(s)) = \left[ \frac{\nu}{\nu-2} - \frac{\nu\Gamma\left(\frac{\nu-1}{2}\right)\eta^2}{\pi\Gamma^2\left(\frac{\nu}{2}\right)} \right] \).

If \( \eta = 0 \), (3.6) reduces to a marginal t density given in (2.2) and if \( \nu \to \infty \), (3.6) converges to a skew-normal distribution. Moreover, coupling (2.6) and (3.3) the correlation function of the skew-t process is given by:

\[
\rho_{S^*_\nu,\eta}(h) = a(\nu,\eta) \left[ 2F_1\left(\frac{1}{2} \; \frac{1}{2} \; \frac{\nu}{2} \; \frac{\nu^2}{2}\right) \left( (1 + \eta^2(1 - \frac{2}{\pi}))\rho_{\nu,\eta}(h) + \frac{2\eta^2}{\pi} \right) - \frac{2\eta^2}{\pi} \right],
\]

(3.7)

where \( a(\nu,\eta) = \frac{\pi(\nu-2)\Gamma^2\left(\frac{\nu}{2}\right)}{2\pi\Gamma^2\left(\frac{\nu}{2}\right)(1+\eta^2)-\eta^2(\nu-2)\Gamma^2\left(\frac{\nu}{2}\right)} \). Note that \( \rho_{S^*_\nu,\eta}(h) = \rho_{S^*_\nu,\eta}(h) \) that is, as in the skew-Gaussian process \( U_\eta \), the correlation is invariant with respect to positive or negative asymmetry and using similar arguments of Theorem 2.2 point e), it can be shown that \( \lim_{\nu \to \infty} \rho_{S^*_\nu,\eta}(h) = \rho_{U_\eta}(h) \).

Finally, following the steps of the proof of Theorem 2.2 it can be shown that properties a), b), c) and d) in Theorem 2.2 are true for the skew-t process \( S^*_\nu,\eta \).

Figure 4, left part, compares \( \rho_{S^*_7,0.9}(h) \) and \( \rho_{S^*_7,0}(h) = \rho_{Y^*_7}(h) \) with the underlying correlation \( \rho(h) = GW_{0.3,1,5}(h) \). The right part shows a realization of \( S^*_7,0.9 \).

**Figure 4:** From left to right: a) comparison between \( \rho_{S^*_7,0.9}(h) \); \( \rho_{S^*_7,0}(h) = \rho_{Y^*_7}(h) \) and the underlying correlation \( \rho(h) = GW_{0.3,1,5}(h) \); b) a realization from \( S^*_7,0.9(s) \).

### 4 Numerical examples

In this Section we analyze the performance of the *wpl* method when estimating the *t* process assuming \( \nu \) known or unknown. Following Remark 1 in Section 2,
we consider the cases when \( \nu > 2 \) and \( \nu = 3, 4, \ldots \). In the latter case, we give a practical solution for fixing the degrees of freedom parameter to a positive integer value through a two-step estimation.

We also compare the performance of the \( wpl \) using the bivariate \( t \) distribution (2.10) with a misspecified Gaussian standard and weighted pairwise likelihood. Finally, we compare the performance of the optimal linear predictor of the \( t \) process using (2.6) versus the optimal predictor of the Gaussian process.

### 4.1 Weighted pairwise likelihood estimation

Let \( (y_1, \ldots, y_n)^T \) be a realization of the \( t \) random process \( Y_\nu \) defined in equation (2.3) observed at distinct spatial locations \( s_1, \ldots, s_n, s_i \in A \) and let \( \theta = (\beta^T, \nu, \sigma^2, \alpha^T) \) be the vector of unknown parameters where \( \alpha \) is the vector parameter associated with the underlying correlation model. The method of \( wpl \), \( \text{(Lindsay, 1988; Varin et al., 2011)} \) combines the bivariate distributions of all possible distinct pairs of observations. The pairwise likelihood function is given by

\[
\text{pl}(\theta) := \sum_{i=1}^{n-1} \sum_{j=i+1}^{n} \log(f_{Y_{\nu,ij}}(y_i, y_j; \theta)) c_{ij}
\]

where \( f_{Y_{\nu,ij}}(y_i, y_j; \theta) \) is the bivariate density in (2.11) and \( c_{ij} \) is a nonnegative suitable weight. The choice of cut-off weights, namely

\[
c_{ij} = \begin{cases} 
1 & \|s_i - s_j\| \leq d_{ij} \\
0 & \text{otherwise}
\end{cases}
\]

for a positive value of \( d_{ij} \), can be motivated by its simplicity and by observing that the dependence between observations that are distant is weak. Therefore, the use of all pairs may skew the information confined in pairs of near observations \( \text{[Bevilacqua and Gaetan, 2015; Joe and Lee, 2009]} \). The maximum \( wpl \) estimator is given by

\[
\hat{\theta} := \arg\max_\theta \text{pl}(\theta)
\]

and, arguing as in \( \text{Bevilacqua et al., 2012; Bevilacqua and Gaetan, 2015} \), under some mixing conditions of the \( t \) process, it can be shown that, under increasing domain asymptotics, \( \hat{\theta} \) is consistent and asymptotically Gaussian with the
asymptotic covariance matrix given by $G_n^{-1}(\theta)$ the inverse of the Godambe information $G_n(\theta) := \mathcal{K}_n(\theta)J_n(\theta)^{-1}\mathcal{K}_n(\theta)$, where $\mathcal{K}_n(\theta) := \mathbb{E}[-\nabla^2 \text{pl}(\theta)]$ and $J_n(\theta) := \text{Var}[\nabla \text{pl}(\theta)]$. Standard error estimation can be obtained considering the square root diagonal elements of $G_n^{-1}(\hat{\theta})$. Moreover, model selection can be performed by considering two information criterion, defined as

$$\text{PLIC} := -2 \text{pl}(\hat{\theta}) + 2\text{tr}(\mathcal{K}_n(\hat{\theta})G_n^{-1}(\hat{\theta})),$$

$$\text{BLIC} := -2 \text{pl}(\hat{\theta}) + \log(n)\text{tr}(\mathcal{K}_n(\hat{\theta})G_n^{-1}(\hat{\theta}))$$

which are composite likelihood version of the Akaike information criterion (AIC) and Bayesian information criterion (BIC) respectively (Varin and Vidoni, 2005; Gao and Song, 2010). Note that, the computation of standard errors, PLIC and BLIC require evaluation of the matrices $\mathcal{K}_n(\hat{\theta})$ and $J_n(\hat{\theta})$. However, the evaluation of $J_n(\hat{\theta})$ is computationally unfeasible for large datasets and in this case subsampling techniques can be used in order to estimate $J_n(\theta)$ as in Bevilacqua et al. (2012) and Heagerty and Lele (1998). A straightforward and more robust alternative is parametric bootstrap estimation of $G_n^{-1}(\theta)$ (Bai et al., 2014). We adopt the second strategy in Section 5.

4.2 Performance of the weighted pairwise likelihood estimation

Following DiCiccio and Monti (2011) and Arellano-Valle and Azzalini (2013) we consider a reparametrization for the $t$ process by using the inverse of degrees of freedom, $\lambda = 1/\nu$. In the standard i.i.d case this kind of parametrization has proven effective for solving some problems associated with the singularity of the Fisher information matrix associated to the original parametrization. Here we consider two possible scenarios i.e. a $t$ process observed on a subset of $\mathbb{R}$ and $\mathbb{R}^2$.

1. We consider points $s_i \in A = [0, 1]$, $i = 1, \ldots, N$ and an exponential correlation function for the ‘parent’ Gaussian process. Then, according to Remark 1 in section 2 the $t$ process is well-defined for $0 < \lambda < 1/2$ and in this specific case all the parameters can be jointly estimated. We simulate, using Cholesky decomposition, 500 realizations of a $t$ process observed on a regular transect
We consider two mean regression parameters, that is, $\mu(s_i) = \beta_0 + \beta_1 u(s_i)$ with $\beta_0 = 0.5$, $\beta_1 = -0.25$ where $u(s_i)$ is a realization from a $U(0,1)$. Then we set $\lambda = 1/\nu$, $\nu = 3, 6, 9$ and $\sigma^2 = 1$.

As correlation model we consider $\rho(h) = M_{\alpha,0.5}(h) = e^{-|h|/\alpha}$ with $\alpha = 0.1/3$ and in the wpl estimation we consider a cut-off weight function with $d_{ij} = 0.002$. Table 1 shows the bias and mean square error (MSE) associated with $\lambda$, $\beta_0$, $\beta_1$, $\alpha$ and $\sigma^2$.

2. We consider points $s_i \in A = [0,1]^2$, $i = 1,\ldots,N$. Specifically, we simulate, using Cholesky decomposition, 500 realizations of a $t$ process observed at $N = 500$ spatial location sites uniformly distributed in the unit square. Regression, variance and (inverse of) degrees of freedom parameters have been set as in the first scenario. As an isotropic parametric correlation model, $\rho(h) = G W_{\alpha,0.4}(h)$ with $\alpha = 0.2$ is considered. In the wpl estimation we consider a cut-off weight function with $d_{ij} = 0.05$ and for each simulation we estimate with wpl, assuming the degrees of freedom are fixed and known.

We also consider the more realistic case when the (inverse of) degrees of freedom are supposed to be unknown. Recall that from Remark 1, $\nu$ must be fixed to a positive integer $\nu = 3, 4, \ldots$ in order to guarantee the existence of the $t$ process. A brute force approach considers different wpl estimates using a fixed $\lambda = 1/\nu$, $\nu = 3, 4, \ldots$ and then simply keeps the estimate with the best PLIC or BLIC. We propose a computationally easier approach by considering a two-step method. In the first step, we estimate all the parameters including $0 < \lambda < 1/2$ maximizing the wpl function. This is possible since the bivariate $t$ distribution is well defined for $0 < \lambda < 1/2$ (see Remark 4). In the second step $\nu$ is fixed equal to the rounded value of $1/\hat{\lambda}$ where $\hat{\lambda}$ is the estimation at first step. (If at the first step, the estimation of $1/\hat{\lambda}$ is lower than 2.5, then it is rounded to 3). Table 2 shows the bias and MSE associated with $\beta_0$, $\beta_1$, $\alpha$ and $\sigma^2$ when estimating with wpl, assuming (the inverse of) degrees of freedom 1) known and fixed, and 2) unknown and fixed using a two-step estimation and Figure 5 shows the boxplots of the wpl estimates for the case 1) and 2).
As a general comment, the distribution of the estimates are quite symmetric, numerically stable and with very few outliers for the three scenarios. In Scenario 1, the MSE of \( \lambda = 1/\nu \) slightly decreases when increasing \( \nu \). Moreover, in Table 2 it can be appreciated that only the estimation of \( \sigma^2 \) is affected when considering a two step estimation. Specifically, the MSE of \( \sigma^2 \) slightly increases with respect to the one-step estimation, \textit{i.e.}, when the degrees of freedom are supposed to be known.

| \( \lambda \) | 1/3 | 1/6 | 1/9 |
|--------------|-----|-----|-----|
| Bias         | MSE | Bias| MSE | Bias| MSE |
| \( \lambda \) | -0.01022 | 0.00321 | -0.01033 | 0.00215 | -0.00924 | 0.00172 |
| \( \beta_0 \) | -0.01466 | 0.00585 | -0.00559 | 0.00154 | 0.00028 | 0.00036 |
| \( \beta_1 \) | -0.000699 | 0.00062 | -0.00032 | 0.00049 | -0.00193 | 0.00049 |
| \( \alpha \) | -0.00213 | 0.0006 | -0.00278 | 0.0007 | -0.00189 | 0.0008 |
| \( \sigma^2 \) | -0.01064 | 0.07493 | -0.01064 | 0.07522 | -0.01064 | 0.07522 |

Table 1: Bias and MSE when estimating with \( wpl \) the \( t \) process with \( \lambda = 1/\nu \), \( \nu = 3, 6, 9 \) and exponential correlation function (Scenario 1).

| \( \lambda \) | 1/3 | 1/6 | 1/9 |
|--------------|-----|-----|-----|
| Bias         | MSE | Bias| MSE | Bias| MSE | Bias| MSE |
| \( \beta_0 \) | -0.00024 | 0.01261 | -0.00037 | 0.01271 | 0.00712 | 0.01135 | 0.00638 | 0.01153 | 0.00646 | 0.0115 |
| \( \beta_1 \) | -0.00702 | 0.00874 | -0.00721 | 0.00883 | 0.00419 | 0.00697 | 0.00419 | 0.00696 | -0.00089 | 0.00685 | -0.00098 | 0.00694 |
| \( \alpha \) | -0.00233 | 0.00058 | -0.00352 | 0.00058 | -0.00115 | 0.00059 | -0.00014 | 0.00061 | -0.00025 | 0.00056 | -0.00048 | 0.00058 |
| \( \sigma^2 \) | -0.00971 | 0.01341 | 0.01768 | 0.01745 | -0.00065 | 0.01142 | -0.00149 | 0.01148 | -0.00015 | 0.01657 |

Table 2: Bias and MSE when estimating with \( wpl \) the \( t \) process when the (inverse of) degrees of freedom (\( \lambda = 1/\nu \), \( \nu = 3, 6, 9 \)) are: 1) fixed and known, 2) unknown and fixed through a two-step estimation (Scenario 2).

### 4.3 Performance of the misspecified (pairwise) Gaussian likelihood estimation

Weighted pairwise likelihood estimation requires the evaluation of the bivariate distribution (2.10) \textit{i.e.} the computation of the Appell \( F_4 \) function. Standard statistical software libraries for the computation of the \( F_4 \) function are unavailable to the best of our knowledge. In our implementation we exploit the following relation with the
Figure 5: Boxplots of \( wpl \) estimates for \( \beta_0 = 0.5, \beta_1 = -0.25, \alpha = 0.2, \sigma^2 = 1 \) (from left to right) under Scenario 2 when estimating a \( t \) process with \( \lambda = 1/\nu, \nu = 6 \) when 1) \( \nu \) is assumed known, 2) \( \nu \) is assumed unknown and it is fixed to a positive integer through a two-step estimation.

Gaussian hypergeometric function \cite{BrychkovSaad2017}:

\[
F_4(a, b; c, c'; w, z) = \sum_{k=0}^{\infty} \frac{(a)_k(b)_k k^k}{k!(c')_k} \, _2F_1(a+k, b+k; c; w), \quad |\sqrt{w}| + |\sqrt{z}| < 1,
\]

truncating the series when the \( k \)-th generic element of the series is smaller than a fixed \( \epsilon \) and where standard libraries for the computation of the \( _2F_1 \) function can be used \cite{Pearsonetal2017}. Evaluation of the \( F_4 \) function can be time consuming depending of the speed of convergence of (4.3) and, as a consequence, if the number of location sites is large the computation of the \( wpl \) estimator can be computationally demanding.

An estimator that require smaller computational burden can be obtained by considering a misspecified \( wpl \). Specifically, if in the estimation procedure we assume a Gaussian process with mean equal to \( \mu(s) \), variance equal to \( \sigma^2 \nu/(\nu - 2) \) and correlation \( \rho_Y(h) \), then a Gaussian \( wpl \) only requires the computation of the Gaussian bivariate distribution and of the Gauss hypergeometric function in (2.6). Note that the misspecified Gaussian process matches mean, variance and correlation function of the \( t \) process. To avoid identifiability problems, we need a reparametrization of the variance \( i.e. \sigma_*^2 := \sigma^2 \nu/(\nu - 2) \). Then, maximization of the Gaussian \( wpl \) func-
tion leads to the estimation of $\mu(s)$, $\sigma^2$, $\nu$ and the parameters of the underlying correlation model $\rho(h)$.

To investigate the performance of this kind of estimator, we consider 676 points on a regular spatial grid $A = [0,1]^2$ that is $(x_i, x_j)^T$ for $i, j = 1, \ldots, 21$ with $x_1 = 0, x_2 = 0.04, \ldots x_{26} = 1$ and we simulate, using Cholesky decomposition, 500 realizations of a $t$ process setting $\mu(s) = \mu = 0$, $\sigma^2 = 1$, $\nu = 3, 6, 9$ and underlying correlation function $\rho(h) = \mathcal{GW}_{0,\lambda}(h)$ with $\alpha = 0.2$. Then we estimate the parameters $\mu$, $\sigma^2$, $\alpha$ (assuming $\nu$ known and fixed) with wpl using the bivariate $t$ distribution (2.10) and with both misspecified Gaussian wpl and standard likelihood. In the wpl estimation we consider a cut-off weight function with $d_{ij} = 0.05$.

Table 3 shows the bias and MSE associated with $\mu$, $\alpha$, and $\sigma^2$ for the three methods of estimation. Note that, for comparison, the results of the variance parameter are reported in terms of the original parametrization. It is apparent that wpl with bivariate $t$ distribution show the best performance. In particular when $\lambda = 1/3$ the gains in terms of efficiency are considerable. However, when increasing the degrees of freedom the gains tends to decrease and when $\nu = 9$ the efficiencies of the three estimators are quite similar (see boxplots in Figure 6).

| $\lambda$ | Parameters | $WPL_T$ | $LG$ | $WPL_G$ |
|-----------|------------|---------|------|---------|
| $\lambda = 1/3$ | $\mu$ | 0.0045 | 0.0088 | 0.0036 | 0.0156 |
| | $\alpha$ | -0.0016 | 0.0003 | 0.0084 | 0.0009 |
| | $\sigma^2$ | -0.0019 | 0.0102 | -0.0577 | 0.0477 |
| $\lambda = 1/6$ | $\mu$ | -0.0057 | 0.0006 | -0.0053 | 0.0106 |
| | $\alpha$ | -0.0020 | 0.0003 | -0.0007 | 0.0003 |
| | $\sigma^2$ | -0.0105 | 0.0075 | -0.0150 | 0.0095 |
| $\lambda = 1/9$ | $\mu$ | -0.0029 | 0.0091 | -0.0038 | 0.0094 |
| | $\alpha$ | -0.0019 | 0.0003 | -0.0019 | 0.0003 |
| | $\sigma^2$ | -0.0075 | 0.0081 | -0.0089 | 0.0088 |

Table 3: Bias and MSE associated with $\mu$, $\alpha$ and $\sigma^2$ for wpl with bivariate $t$ distribution ($WPL_T$), standard misspecified Gaussian likelihood ($LG$) and wpl with bivariate misspecified Gaussian distribution ($WPL_G$) when $\lambda = 1/\nu$, $\nu = 3, 6, 9$. 

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Figure 6: Upper part: boxplots of $\mu$, $\alpha$, $\sigma^2$ for $wpl$ using $t$ bivariate distribution ($WPL_T$), standard misspecified Gaussian likelihood ($LG$) and $wpl$ using Gaussian misspecified bivariate distribution ($WPL_G$) when $\nu = 3$. Bottom part: The same boxplots with $\nu = 9$.

4.4 $t$ optimal linear prediction versus Gaussian optimal prediction

One of the primary goals of geostatistical modeling is to make predictions at spatial locations without observations. The optimal predictor for the $t$ process, with respect to the mean squared error criterion, is nonlinear and difficult to evaluate explicitly since it requires the knowledge of the finite dimensional distribution.

A practical and less efficient solution can be obtained using the optimal linear
prediction. Assuming known mean, correlation and the degrees of freedom of the t process, the predictor at an unknown location \( s_0 \) is given by:

\[
\hat{y}(s_0) = \mu(s_0) + c^T \nu R^{-1}_\nu (Y - \mu)
\]

where \( \mu = (\mu(s_1), \ldots, \mu(s_n))^T \), \( c_\nu = [\rho_{Y^\nu}(s_0 - s_j)]_{i=1}^n \) and \( R_\nu = [\rho_{Y^\nu}(s_i - s_j)]_{i,j=1}^n \), and the associated variance is given by:

\[
\text{Var}(\hat{y}(s_0)) = \sigma^2(1 - c^T \nu R^{-1}_\nu c_\nu).
\]

As an Associate Editor pointed out, this is equivalent to perform optimal Gaussian prediction with covariance function equal to \( \sigma^2 \rho_{Y^\nu}(h) \). Similarly, using (3.7), the optimal linear predictor of the skew-t process can be obtained.

We investigate the performance of (4.4) when compared with the Gaussian optimal predictor, assuming \( \rho(h) = GW_{0,2,\psi,4}(h) \), \( \psi = 0, 1, 2 \) as underlying correlation function, using cross-validation. With this goal in mind, we simulate 1000 realizations from a t process \( Y^\nu \) with \( \nu = 3, 7, 11 \) and a Gaussian process under the settings of Section 4.3 and for each realization, we consider 80% of the data for prediction and leave 20% as validation dataset.

For each model and for each realization, we compute the root mean square errors (RMSEs) that is:

\[
\text{RMSE}_l = \left( \frac{1}{n_l} \sum_{i=1}^{n_l} (y(s_{i,l}) - \hat{y}(s_{i,l}))^2 \right)^{\frac{1}{2}},
\]

where \( y(s_{i,l}), i = 1, \ldots, n_l \) are the observation in the \( l \)-th validation set and \( n_l \) is the associated cardinality (\( n_l = 135 \) in our example).

Finally we compute the empirical mean of the 500 RMSEs when the prediction is performed with the optimal linear predictor (4.4) i.e. using \( \rho_{Y^\nu}(h) \) with \( \nu = 3, 7, 11 \) and the optimal Gaussian predictor with \( \rho(h) \). Note that, from Theorem 2.2 (e), the prediction using \( \rho(h) \) can be viewed as the prediction using \( \rho_{Y^\nu}(h) \) when \( \nu \to \infty \).

In Table 4 we report the simulation results in terms of relative efficiency that is, for a given process and a given \( \psi = 0, 1, 2 \), the ratio between the mean RMSE of the best predictor and the mean RMSE associated with a competitive predictor. This implies that relative efficiency prediction is lower than 1 and it is equal to 1 in the best case. From Table 4, it can be appreciated that under the t process...
the prediction with $\rho_{Y^*_\nu}(h)$, for $\nu = 3, 7, 11$, performs overall better than the optimal Gaussian prediction using $\rho(h)$. As expected, the gain is more apparent when decreasing the degrees of freedom and increasing $\psi$. For instance, if $\nu = 3$ and $\psi = 2$ the loss of efficiency predicting with $\rho(h)$ is $19\%$ approximatively. It can also be noted that if $Y^*_3, Y^*_7$ or Gaussian are one or two mean squared differentiable ($\psi = 1, 2$), then the prediction using $\rho_{Y^*_3}(h)$ can be very inefficient. This is not surprising since from Theorem 2.2 (c), $Y^*_3$ is not mean square differentiable. Resuming, this numerical experiment study suggests that when predicting data exhibiting heavy tails the use of the correlation function $\rho_{Y^*_\nu}(h)$ should be preferred to the use of $\rho(h)$.

| $\mathcal{GW}_{0.2,\psi,4}(h)$ | Simulation from | $Y^*_3$ | $Y^*_7$ | $Y^*_11$ | Gaussian |
|-----------------------------|----------------|--------|--------|---------|---------|
| $\psi = 0$                  |                |        |        |         |         |
| $\rho_Y(h)$                 | 1              | 0.9896 | 0.9870 | 0.9838  |         |
| $\rho_{Y^*_7}(h)$           | 0.9904         | 1      | 0.9998 | 0.9987  |         |
| $\rho_{Y^*_11}(h)$          | 0.9877         | 0.9998 | 1      | 0.9996  |         |
| $\rho(h)$                   | 0.9836         | 0.9988 | 0.9996 | 1       |         |
| $\psi = 1$                  |                |        |        |         |         |
| $\rho_Y(h)$                 | 1              | 0.9399 | 0.9180 | 0.8789  |         |
| $\rho_{Y^*_7}(h)$           | 0.9503         | 1      | 0.9984 | 0.9895  |         |
| $\rho_{Y^*_11}(h)$          | 0.9327         | 0.9984 | 1      | 0.9966  |         |
| $\rho(h)$                   | 0.9095         | 0.9947 | 0.9969 | 1       |         |
| $\psi = 2$                  |                |        |        |         |         |
| $\rho_Y(h)$                 | 1              | 0.8827 | 0.8263 | 0.7109  |         |
| $\rho_{Y^*_7}(h)$           | 0.9223         | 1      | 0.9936 | 0.9511  |         |
| $\rho_{Y^*_11}(h)$          | 0.8834         | 0.9947 | 1      | 0.9817  |         |
| $\rho(h)$                   | 0.8169         | 0.9643 | 0.9848 | 1       |         |

Table 4: Relative mean RMSEs prediction efficiency over 500 runs for a $t$ process with $\nu = 3, 7, 11$ and a Gaussian process when predicting using $\rho_{Y^*_\nu}(h)$ for $\nu = 3, 7, 11$ and $\rho(h)$. The underlying correlation model is $\rho(h) = \mathcal{GW}_{0.2,\psi,4}(h)$ with $\psi = 0, 1, 2$.

5 Application to Maximum Temperature Data

In this section, we apply the proposed $t$ process to a data set of maximum temperature data observed in Australia. Specifically, we consider a subset of a global data set of merged maximum daily temperature measurements from the Global Sur-
face Summary of Day data (GSOD) with European Climate Assessment &Dataset (ECA&D) data in July 2011. The dataset is described in detail in Kilibarda et al. (2014) and it is available in the R package `meteo`. The subset we consider is depicted in Figure 7 (a) and consists of the maximum temperature observed on July 5 in 446 location sites, $y(s_i), i = 1, \ldots, 446$, in the region with longitude $[110, 154]$ and latitude $[-39, -12]$.

Figure 7: From left to right: a) spatial locations of maximum temperature in Australia in July 2011 and b) prediction of residuals of the estimated $t$ process.

Spatial coordinates are given in longitude and latitude expressed as decimal degrees and we consider the proposed $t$ process defined on the planet Earth sphere approximation $S^2 = \{ s \in \mathbb{R}^3 : ||s|| = 6371 \}$. The first process we use to model this dataset is a $t$ process:

$$Y_{\nu}(s) = \beta_0 + \beta_1 X(s) + \sigma Y_{\nu}^*(s), \quad s \in S^2$$

(5.1)

where $Y_{\nu}^*$ is a standard $t$ process. Here, $X(s)$ is a covariate called geometric temperature which represents the geometric position of a particular location on Earth and the day of the year (Kilibarda et al., 2014). As a comparison, we also consider a Gaussian process:

$$Y(s) = \beta_0 + \beta_1 X(s) + \sigma G(s), \quad s \in S^2$$

(5.2)
where $G$ is a standard Gaussian process. We assume that the underlying geodesically isotropic correlation function (Gneiting 2013; Porcu et al. 2016) is of the Matérn and generalized Wendland type. A preliminary estimation of the $t$ and Gaussian processes, including the smoothness parameters, highlights a multimodality of the (pairwise) likelihood surface, for both correlation models and a not mean-square differentiability of the process. For this reason, we fix the smoothness parameters and we consider the underlying correlation models $M_{\alpha,0.5}(d_{GC}) = e^{-d_{GC}/\alpha}$ and $GW_{\alpha,0.5}(d_{GC}) = (1 - d_{GC}/\alpha)^{5}$ where, given two spherical points $s_i = (\text{lon}_i, \text{lat}_i)$ and $s_j = (\text{lon}_j, \text{lat}_j)$, $d_{GC}(s_i, s_j) = 6371\theta_{ij}$, is the great circle distance. Here $\theta_{ij} = \arccos\{\sin a_i \sin a_j + \cos a_i \cos a_j \cos(b_i - b_j)\}$ is the great circle distance on the unit sphere with $a_i = (\text{lat}_i)\pi/180$, $a_j = (\text{lat}_j)\pi/180$, $b_i = (\text{lon}_i)\pi/180$, $b_j = (\text{lon}_j)\pi/180$.

For the $t$ process the parameters were estimated using $wpl$ using the bivariate $t$ distribution with the two-step method described in Section 4 and using the weight function (4.2) with $d_{ij} = 150$ Km. It turns out that the estimation at the first step leads to fix $\nu = 4$ in the second step, irrespective of the correlation model. We also consider a Gaussian misspecified standard likelihood and $wpl$ estimation as described in Section 4.3 i.e. we estimate using the Gaussian process (5.2) with the $t$ correlation model fixing $\nu = 4$.

In addition, we compute the standard error estimation, PLIC and BLIC values through parametric bootstrap estimation of the inverse of the Godambe information matrix (Bai et al. 2014). For standard maximum likelihood we compute the standard errors as the square root of diagonal elements of the inverse of Fisher Information matrix (Mardia and Marshall 1984). The results are summarized in Table 5. Note that the regression parameters estimates are quite similar for the $t$ and Gaussian processes, irrespective of the correlation model. Furthermore, we note that the standard Gaussian process assigns lower spatial dependence and stronger variance compared to the other cases. Finally, for each correlation model, both the (pairwise) likelihood information criterion PLIC and BLIC select for the pairwise case the $t$ model and for the standard case the Gaussian model with $t$ correlation function.

Given the estimation of the mean regression and variance parameters of the $t$
Table 5: Estimates for the $t$ process using $wpl$ with bivariate $t$ distribution and misspecified (full and weighted pairwise) Gaussian likelihood and for the Gaussian process using standard likelihood with associated standard error (in parenthesis) and PLIC and BLIC values, when estimating the Australian maximum temperature dataset using two correlation models: $M_{\alpha,0.5}$ and $GW_{\alpha,0.5}$. Last three columns: associated empirical mean of RMSEs, MAEs and CRPSs.
process, the estimated residuals
\[ \hat{Y}_4^*(s_i) = \frac{y(s_i) - (\hat{\beta}_0 + \hat{\beta}_1 X(s_i))}{(\hat{\sigma}^2)^{\frac{1}{2}}} \quad i = 1, \ldots N \]
can be viewed as a realization of the process \( Y_4^* \). Similarly we can compute the Gaussian residuals. For the \( t \) process we use the \( \text{wpl} \) estimates obtained with the bivariate \( t \) distribution. Both residuals can be useful in order to check the model assumptions, in particular the marginal and dependence assumptions. In the top part of Figure 8 a qq-plot of the residuals of the Gaussian and \( t \) processes (from left to right) is depicted for the Matérn case. It can be appreciated that the \( t \) model overall fits better with respect the Gaussian model even if it seems to fail to model properly the right tail behavior. Moreover, the graphical comparison between the empirical and fitted semivariogram of the residuals (bottom part of Figure 8) highlights an apparent better fitting of the \( t \) model.

We want to further evaluate the predictive performances of Gaussian and \( t \) processes using RMSE and MAE as in Section 4.4. Specifically, we use the following resampling approach: we randomly choose 80% of the data to predict and we use the estimates previously obtained in order to compute RMSE and MAE values at the remaining 20% of the spatial locations. We repeat the approach for 2000 times and record all RMSEs and MAEs. Specifically, for each \( j \)–th left-out sample \((y_j^L(s_1), \ldots, y_j^L(s_K))\), we compute

\[ \text{RMSE}_j = \left[ \frac{1}{K} \sum_{i=1}^{K} \left( y_j^L(s_i) - \hat{Y}_j^L(s_i) \right)^2 \right]^{\frac{1}{2}} \]

and

\[ \overline{\text{MAE}}_j = \frac{1}{K} \sum_{i=1}^{K} |y_j^L(s_i) - \hat{Y}_j^L(s_i)|, \]

where \( \hat{Y}_j^L(s_i) \) is the optimal or the best linear optimal prediction for the Gaussian and \( t \) processes respectively. Finally, we compute the overall mean for both Gaussian and \( t \) processes and for both correlation models, that is RMSE = \( \sum_{j=1}^{2000} \text{RMSE}_j / 2000 \) and MAE = \( \sum_{j=1}^{2000} \overline{\text{MAE}}_j / 2000 \).

Additionally, to evaluate the marginal predictive distribution performance, we also consider, for each sample, the continuous ranked probability score (CRPS)
Figure 8: Upper part: Q-Q plot of the residuals versus the estimated quantiles in the Gaussian and t models ((a) and (b) respectively). Bottom part: Empirical semi-variogram (dotted points) of the residuals versus the estimated semivariogram (solid line) in the Gaussian and t models ((c) and (d) respectively). Distances are expressed in Km.

(Gneiting and Raftery [2007]). For a single predictive cumulative distribution function $F$ and a verifying observation $y$, it is defined as:

$$\text{CRPS}(F, y) = \int_{-\infty}^{\infty} (F(t) - 1_{[y, \infty)}(t))^2 dt.$$  

Specifically, for each $j-th$ left-out sample, we consider the averaged CRPS for the
Gaussian and $t$ distributions as:

$$\text{CRPS}_j = \frac{1}{K} \sum_{i=1}^{K} \text{CRPS}(F, y_j^i(s_i)) \quad F = F_G, F_{Y_4},$$

for $j = 1, \ldots, 2000$. In particular in the Gaussian case

$$\text{CRPS}(F_G, y_j^i(s_i)) = \sigma \left( \frac{y_j^i(s_i) - \mu(s)}{\sigma} \right) \left[ 2F_{G^*}(s) \left( \frac{y_j^i(s_i) - \mu(s)}{\sigma} \right) - 1 \right] + 2\sigma f_{G^*}(s) \left( \frac{y_j^i(s_i) - \mu(s)}{\sigma} \right) - \frac{\sigma}{\sqrt{\pi}},$$

and in the $t$ case with 4 degrees of freedom:

$$\text{CRPS}(F_{Y_4}, y_j^i(s_i)) = \sigma \left( \frac{y_j^i(s_i) - \mu(s)}{\sigma} \right) \left[ 2F_{Y_4^*}(s) \left( \frac{y_j^i(s_i) - \mu(s)}{\sigma} \right) - 1 \right] + 2 \left[ 4\sigma^2 + (y_j^i(s_i) - \mu(s))^2 \right] f_{Y_4^*}(s) \left( \frac{y_j^i(s_i) - \mu(s)}{\sigma} \right) - \frac{4\sigma B \left( \frac{1}{2}, \frac{7}{2} \right)}{3B^2 \left( \frac{1}{2}, 2 \right)},$$

where $\mu(s) = \beta_0 + \beta_1 X(s)$. We compute the CRPS in (5.4) and (5.5) plugging-in the estimates of the pairwise and standard (misspecified) likelihood estimation methods of $\beta_0$, $\beta_1$ and $\sigma$ using the R package \texttt{scoringRules} \cite{Jordan:2019}. Finally, we compute the overall mean for both Gaussian and $t$ processes and for both correlation models, that is $\text{CRPS} = \sum_{j=1}^{2000} \text{CRPS}_j / 2000$.

Table 5 reports the estimated RMSE, MAE and CRPS. As a general remark, the $t$ process outperforms the Gaussian process for the three measures of prediction performance irrespective of the method of estimation and for both correlation models. We point out that RMSE and MAE are computed using the optimal predictor in the Gaussian case and the linear optimal in the $t$ case. However, the RMSE and MAE results highlight a better performance for the $t$ process. In addition, a better RMSE and MAE for the Matérn correlation model with respect to the Wendland is apparent, irrespective of the type of process. The proposed $t$ process also leads to a clear better performance of the CRPS with respect to the Gaussian case. In this specific example, the use of the misspecified Gaussian \texttt{wpl} estimates leads to the best results in terms of RMSE and MAE. On the other hand, the best CRPS results are achieved by using the \texttt{wpl} estimates using the proposed $t$ bivariate distribution.

Finally, one important goal in spatial modeling of temperature data is to create a high resolution map in a spatial region using the observed data. In Figure 7 (b), we plot a high resolution map of the predicted residuals using the $t$ process with underlying Matérn correlation model estimated with \texttt{wpl}. 

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6 Concluding remarks

We have introduced a new stochastic process with \( t \) marginal distributions for regression and dependence analysis when addressing spatial with heavy tails. Our proposal allows overcoming any problem of identifiability associated with previously proposed spatial models with \( t \) marginals and, as a consequence, the model parameters can be estimated with just one realization of the process. Additionally, the proposed \( t \) process partially inherits some geometrical properties of the ‘parent’ Gaussian process, an appealing feature from a data analysis point of view. We have also proposed a possible generalization, obtaining a new process with the marginal distribution of the skew-\( t \) type using the skew-Gaussian process proposed in Zhang and El-Shaarawi (2010).

In our proposal, a possible limitation is the lack of amenable expressions of the associated multivariate distributions. This prevents an inference approach based on the full likelihood and the computation of the optimal predictor. In the first case, our simulation study shows that an inferential approach based on \( wpl \), using the bivariate \( t \) distribution given in Theorem 2.3, could be an effective solution for estimating the unknown parameters. An alternative less efficient solution that requires smaller computational burden can be obtained by considering a misspecified Gaussian \( wpl \) using the correlation function of the \( t \) process. In the prediction case, our numerical experiments show that the optimal linear predictor of the \( t \) process performs better than the optimal Gaussian predictor when working with spatial data with heavy tails.

Another possible drawback concerns the restriction of the degrees of freedom of the \( t \) process to \( \nu = 3, 4, \ldots \) under non-infinite divisibility of the associated Gamma process. This problem could be solved by considering a Gamma process obtained by mixing the proposed Gamma process with a process with beta marginals and using the results in Yeo and Milne (1991); however the mathematics involved with this approach are much more challenging.

The estimation of the skew-\( t \) process has not been addressed in this paper since the bivariate distribution in this case is quite complicated. In principle, after a
suitable parametrization, Gaussian misspecified \( wpl \) can be performed using (3.7) to estimate the parameters of the skew-t process. In this case an additional issue is the behavior of the information matrix when \( \eta = 0 \) (Arellano-Valle and Azzalini, 2013). Finally, a \( t \) process with asymmetric marginal distribution can also be obtained by considering some specific transformations of the proposed standard \( t \) process as in J. F. Rosco and Pewsey (2011) or under the two-piece distribution framework (Arellano-Valle et al., 2005) and this will be studied in future work.

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A Appendix

A.1 Proof of Theorem 2.1

Proof. Set \( R_\nu \equiv W_\nu^{-\frac{3}{2}} \). Then the correlation function of \( Y^*_\nu \) is given by

\[
\rho_{Y^*_\nu}(h) = \left( \frac{\nu - 2}{\nu} \right) \left( E(R_\nu(s_i)R_\nu(s_j))\rho(h) \right). \tag{A.1}
\]

To find a closed form for \( E(R_\nu(s_i)R_\nu(s_j)) \), we need the bivariate distribution of \( R_{\nu ij} \) that can be easily obtained from density of the bivariate random vector \( W_{\nu ij} \) given by \cite{bevilacqua2018}:

\[
f_{W_{\nu ij}}(w_i, w_j) = \frac{2^{-\nu+2} \nu^{\nu} (w_i w_j)^{\nu/2 - 1} \exp \left( -\frac{w_i w_j}{2(1 - \rho^2(h))} \right) \left( 2(1 - \rho^2(h)) \right)^{1 - \nu/2} \Gamma (\frac{\nu}{2}) (1 - \rho^2(h))^{\nu/2} \right) I_{\nu/2 - 1} \left( \frac{\nu \sqrt{\rho^2(h)w_i w_j}}{2(1 - \rho^2(h))} \right) \] \tag{A.2}

where \( I_\alpha(\cdot) \) denotes the modified Bessel function of the first kind of order \( \alpha \). \cite{jones1967} show the infinite divisibility of \( W_{\nu ij} \).

Then, for each \( \nu > 2 \), the bivariate distribution of \( R_{\nu ij} \) is given by:

\[
f_{R_{\nu ij}}(r_{ij}) = \frac{2^{-\nu+2} \nu^{\nu} (r_i r_j)^{-\nu-1} \exp \left( -\frac{r_i r_j}{2(1 - \rho^2(h))} \right) \left( 2(1 - \rho^2(h)) \right)^{1 - \nu/2} \Gamma (\frac{\nu}{2}) (1 - \rho^2(h))^{\nu/2} \right) I_{\nu/2 - 1} \left( \frac{\nu \sqrt{\rho^2(h)r_i r_j}}{2(1 - \rho^2(h))} \right) \] \tag{A.3}

Using the identity \( {}_0F_1(b; x) = \Gamma(b) x^{(1-b)/2} I_{b-1}(2\sqrt{x}) \) and the series expansion of hypergeometric function \( {}_0F_1 \) in (A.3) we have

\[
E(R^a(s_i)R^b(s_j)) = \frac{2^{-\nu+2} \nu^{\nu} }{\Gamma^2 (\frac{\nu}{2}) (1 - \rho^2(h))^{\nu/2}} \int_{R^+} r_i^{-\nu+a-1} r_j^{-\nu+b-1} \exp \left( -\frac{r_i r_j}{2(1 - \rho^2(h))} \right) e^{-\frac{\nu (r_i^2 r_j)}{2(1 - \rho^2(h))}} r_i^a r_j^b dr_{ij}
\]

\[
= \frac{2^{-\nu+2} \nu^{\nu} }{\Gamma^2 (\frac{\nu}{2}) (1 - \rho^2(h))^{\nu/2}} \sum_{k=0}^{\infty} \int_{R^+} r_i^{-\nu+a-2k-1} r_j^{-\nu+b-2k-1} \exp \left( -\frac{r_i r_j}{2(1 - \rho^2(h))} \right) e^{-\frac{\nu (r_i^2 r_j)}{2(1 - \rho^2(h))}} r_i^a r_j^b dr_{ij}
\]

\[
= \frac{2^{-\nu+2} \nu^{\nu} }{\Gamma^2 (\frac{\nu}{2}) (1 - \rho^2(h))^{\nu/2}} \sum_{k=0}^{\infty} \frac{I(k) }{k! \left( \frac{\nu}{2} \right)^k} \left( \frac{\rho^2(h) \nu^2}{4(1 - \rho^2(h))} \right)^k \]

where, using Fubini’s Theorem

\[
I(k) = \int_{R^+} r_i^{-\nu+a-2k-1} \exp \left( -\frac{r_i^2}{2(1 - \rho^2(h))} \right) dr_i \int_{R^+} r_j^{-\nu+b-2k-1} \exp \left( -\frac{r_j^2}{2(1 - \rho^2(h))} \right) dr_j
\]
Using the univariate density $f_{R_n(s)}(r) = 2 \left( \frac{\nu}{2} \right)^{\nu/2} r^{-\nu-1} e^{-\frac{r^2}{2\nu}} / \Gamma \left( \frac{\nu}{2} \right)$, we obtain

$$I(k) = \Gamma \left( \frac{\nu-a}{2} + k \right) \Gamma \left( \frac{\nu-b}{2} + k \right) 2^{\nu-a+k-1} \nu^{\nu-k} + k \left( \frac{1 - \rho^2(h)}{\nu} \right) \nu^{\nu-k}$$

and combining equations (A.5) and (A.4), we obtain

$$E(R^a_i(s_i)R^b_i(s_j)) = \frac{2^{-\nu(a+b)/2} \nu(a+b)/2 (1 - \rho^2(h))^{(\nu-a-b)/2} \Gamma \left( \frac{\nu-a}{2} \right) \Gamma \left( \frac{\nu-b}{2} \right) \sum_{k=0}^{\infty} \frac{\left( \frac{\nu-a}{2} \right)_k \left( \frac{\nu-b}{2} \right)_k \rho^{2k}(h)}{\Gamma^2 \left( \frac{\nu}{2} \right)} F_1 \left( \frac{\nu-a}{2}, \frac{\nu-b}{2}; \rho^2(h) \right)$$

Then, using the Euler transformation, we obtain

$$E(R^a_i(s_i)R^b_i(s_j)) = \frac{2^{-\nu(a+b)/2} \nu(a+b)/2 (1 - \rho^2(h))^{(\nu-a-b)/2} \Gamma \left( \frac{\nu-a}{2} \right) \Gamma \left( \frac{\nu-b}{2} \right) \sum_{k=0}^{\infty} \frac{\left( \frac{\nu-a}{2} \right)_k \left( \frac{\nu-b}{2} \right)_k \rho^{2k}(h)}{\Gamma^2 \left( \frac{\nu}{2} \right)} F_1 \left( \frac{\nu-a}{2}, \frac{\nu-b}{2}; \rho^2(h) \right)$$

for $\nu > a$ and $\nu > b$. Finally, setting $a = b = 1$ in (A.6) and using it in (A.1) we obtain (2.6).

**A.2 Proof of Theorem 2.2**

Proof. If $G$ is a weakly stationary Gaussian process with correlation $\rho(h)$ then from (2.6) it is straightforward to see that $Y^*_\nu$ is also weakly stationary. Points b) and c) can be shown using the relations between the geometrical properties of a stationary process and the associated correlation. Specifically, the mean-square continuity and the $m$-times mean-square differentiability of $Y^*_\nu$ are equivalent to the continuity and $2m$-times differentiability of $\rho_{Y^*_\nu}(h)$ at $h = 0$, respectively (Stein, 1999). Recall that the correlation function of $Y^*_\nu$ is given by:

$$\rho_{Y^*_\nu}(h) = a(\nu) \left[ \gamma \left( \frac{1}{2}, \frac{\nu}{2}, \rho^2(h) \right) \right].$$

with $a(\nu) = (\nu-2)\Gamma^2 \left( \frac{\nu-1}{2} \right)$. Using (2.5) it can be easily seen that $\rho(0) = 1$ if and only if $\rho_{Y^*_\nu}(0) = 1$. Then $Y^*_\nu$ is mean-square continuous if and only if $G$ is mean-square continuous.

For the mean square differentiability, let $G$ $m$-times mean square differentiable. Using iteratively the $n$-th derivative of the $2F_1$ function with respect to $x$:

$$2F_1^{(n)}(a, b, c, x) = \frac{(a)_n(b)_n}{(c)_n} 2F_1(a + n, b + n, c + n, x), \quad n = 1, 2, \ldots$$

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and applying the convergence condition $c > b + a$ of identity \(2.5\), it can be shown that \(\rho_{Y^*}^{(2m)}(h)|_{h=0} < \infty\) if \(\nu > 2(2m + 1)\) and, as a consequence, \(Y^*\) is \(m\)--times mean square differentiable under this condition. On the other hand if \(\nu \leq 2(2m + 1)\) then \(Y^*\) is \(m-k\)--times mean-square differentiable if \(2(2(m-k)+1) < \nu \leq 2(2(m-k)+3)\), for \(k = 1, \ldots, m\).

For instance, let assume that \(G\) is 1--times mean square differentiable. This implies that \(\rho^{(i)}(h)|_{h=0} < \infty\), \(i = 1, 2\). Applying \((A.8)\) to \((A.7)\), the second derivative of \(\rho_{Y^*}^{(2)}(h)\) is given by:

\[
\rho_{Y^*}^{(2)}(h) = \frac{a(\nu)}{\nu(\nu+2)} \left[ (6 + 3\nu)_{2} F_{1} \left( 1.5, 1.5; 1 + \frac{\nu}{2}; \rho^{2}(h) \right) \rho(h) \{\rho^{(1)}(h)\}^{2} \right.
\]

\[
+ 9_{2} F_{1} \left( 2.5, 2.5; 2 + \frac{\nu}{2}; \rho^{2}(h) \right) \rho^{3}(h) \{\rho^{(1)}(h)\}^{2} + \nu(2 + \nu)_{2} F_{1} \left( 0.5, 0.5; \frac{\nu}{2}; \rho^{2}(h) \right) \rho^{(2)}(h) \left. \right]
\]

Then, applying the convergence condition of identity \(2.5\), \(\rho_{Y^*}^{(2)}(h)|_{h=0} < \infty\) if \(2 + \nu/2 > 5\) that is \(\nu > 6\). Therefore \(Y^*\) is 1--times mean square differentiable if \(\nu > 6\) and 0--times mean square differentiable if \(2 < \nu \leq 6\).

Point d) can be shown recalling that a process \(F\) is long-range dependent if the correlation of \(F\) is such that \(\int_{\mathbb{R}^+} |\rho_{F}(h)| d^n h = \infty\) (Lim and Teo, 2009). Direct inspection, using series expansion of the hypergeometric function, shows that \(\int_{\mathbb{R}^+} |\rho_{Y^*}(h)| d^n h = \infty\) if and only if \(\int_{\mathbb{R}^+} |\rho(h)| d^n h = \infty\) and, as a consequence, \(Y^*\) has long-range dependence if and only if \(G\) has long-range dependence.

Finally, note that if \(\nu > 2\) then \(a(\nu)_{2} F_{1} \left( \frac{1}{2}, 2; \frac{\nu}{2}; 0 \right) = a(\nu)\), \(a(\nu)_{2} F_{1} \left( \frac{1}{2}, \frac{1}{2}; \frac{\nu}{2}; 1 \right) = 1\) and \(a(\nu)_{2} F_{1} \left( \frac{1}{2}, \frac{1}{2}; 2; x^{2} \right)\) is not decreasing in \(0 \leq x \leq 1\). This implies \(a(\nu)_{2} F_{1} \left( \frac{1}{2}, \frac{1}{2}; \frac{\nu}{2}; x^{2} \right) \leq 1\) that is \(\rho_{Y^*}(h) \leq \rho(h)\). Moreover, \(\lim_{\nu \rightarrow \infty} a(\nu) = 1\) and using series expansion of the hypergeometric function:

\[
\lim_{\nu \rightarrow \infty} \rho_{Y^*}^{(2)}(h) = \rho(h).
\]
A.3 Proof of Theorem 2.3

Proof. Using the identity $\mathbf{0}_F(\nu; b; x) = \Gamma(b)x^{(1-b)/2}I_{b-1}(2\sqrt{x})$ and the series expansion of hypergeometric function $\mathbf{0}_F$, then under the transformation $g_i = y_i\sqrt{w_i}$ and $g_j = y_j\sqrt{w_j}$ with Jacobian $J((g_i, g_j) \to (y_i, y_j)) = (w_i w_j)^{1/2}$, we have:

$$f_{Y_i}(y_{ij}) = \int f_{G_{ij}|w_{ij}|g_{ij}|w_{ij}}f_{W_i}(w_{ij})dw_{ij}$$

$$= \frac{2^{-\nu}\nu^\nu}{2\pi^2 (\nu/2)} \int_{R^2_+} (w_i w_j)^{\nu/2-1} e^{-\gamma_2 R \gamma_2} w_{i+1}\gamma_i w_{j+1}\gamma_j [w_i y_i^2 + w_j y_j^2 - 2\rho(h)\sqrt{w_i y_i} y_j w_j] dw_{ij}$$

$$\times e^{-\gamma_2 R \gamma_2} \nu F_1\left(\frac{\nu^2 \rho^2(h) w_i w_j}{4(1-\rho^2(h))^2}\right) dw_{ij}$$

$$= \frac{2^{-\nu}\nu^\nu}{2\pi^2 (\nu/2)} \int_{R^2_+} (w_i w_j)^{\nu/2-1} e^{-\gamma_2 R \gamma_2} y_i^2 - 2\rho(h)\sqrt{w_i} y_i + \nu w_i e^{-\gamma_2 R \gamma_2} \gamma_2 w_j dw_{ij}$$

$$\times \sum_{k=0}^{\infty} \frac{1}{k!} \left(\frac{\nu^2 \rho^2(h) w_i w_j}{4(1-\rho^2(h))^2}\right)^k dw_{ij}$$

Using (3.462.1) of Gradshteyn and Ryzhik [2007], we obtain

$$I(k) = \int_{R^+} w_i^{\nu+1/2+k-1} e^{-\gamma_2 R \gamma_2} \left[-\gamma_2 R \gamma_2 w_i - \frac{\nu^2 \rho^2(h) w_i}{4(1-\rho^2(h))^2}\right] dw_i$$

$$= 2\left(\frac{y_i^2 + \nu}{1-\rho^2(h)}\right)^{-(\nu+1+k)} \Gamma(\nu + 1 + 2k) \int_{R^+} w_i^{\nu+1/2+k-1} e^{-\gamma_2 R \gamma_2} \left[-\gamma_2 R \gamma_2 w_i - \frac{\nu^2 \rho^2(h) w_i}{4(1-\rho^2(h))^2}\right] dw_i$$

$$\times D_{-(\nu+1+2k)}\left(-\frac{\rho(h) y_i y_j \sqrt{w_j}}{(1-\rho^2(h))(y_i^2 + \nu)}\right) dw_j$$

$$= 2\left(\frac{y_i^2 + \nu}{1-\rho^2(h)}\right)^{-(\nu+1+k)} \Gamma(\nu + 1 + 2k) A(k)$$

where $D_n(x)$ is the parabolic cylinder function. Now, considering (9.240) of Gradshteyn and Ryzhik [2007]:

$$D_{-(\nu+1+2k)}\left(-\frac{\rho(h) y_i y_j \sqrt{w_j}}{(1-\rho^2(h))(y_i^2 + \nu)}\right) = b_1 e^{-\gamma_2 R \gamma_2} \nu F_1\left(\frac{\nu+1}{2} + k; \frac{\mu^2 \rho^2(h) y_i^2 y_j}{4(1-\rho^2(h))(y_i^2 + \nu)}\right)$$

$$\times \left(1 + \frac{\rho(h) y_i y_j \sqrt{w_j}}{(1-\rho^2(h))(y_i^2 + \nu)}\right)$$

$$+ b_2 e^{-\gamma_2 R \gamma_2} \nu F_1\left(\frac{\nu+1}{2} + k; \frac{3\mu^2 \rho^2(h) y_i^2 y_j}{4(1-\rho^2(h))(y_i^2 + \nu)}\right).$$

\[A.10\]
where \( b_1 = \frac{2^{-(\nu+1)/2-k} \sqrt{k!}}{\Gamma\left(\frac{\nu}{2}+k+1\right)} \) and \( b_2 = \frac{2^{-\nu/2-k} \pi \rho(h) y_j y_j}{\Gamma\left(\frac{\nu}{2}+k+1\right) \sqrt{(1-\rho^2(h))(y_j^2+\nu)}} \). Replacing equations (A.11) in (A.10) and using (7.621.4) of Gradshteyn and Ryzhik (2007), we obtain

\[
A(k) = b_1 \int_{\mathbb{R}_+} w_j^{\nu/(2+k-1)} e^{-\frac{(y_j^2 + \nu)}{2(1-\rho^2(h))}} w_j \frac{\nu + 1}{2} + k \cdot \frac{\rho^2(h) y_j^2 y_j}{2(1-\rho^2(h))(y_j^2 + \nu)} \, dw_j
\]

\[
+ b_2 \int_{\mathbb{R}_+} w_j^{\nu/(2+k-1)} e^{-\frac{(y_j^2 + \nu)}{2(1-\rho^2(h))}} w_j \frac{\nu + 1}{2} + k \cdot \frac{\rho^2(h) y_j^2 y_j}{2(1-\rho^2(h))(y_j^2 + \nu)} \, dw_j
\]

\[
= b_1 \Gamma\left(\frac{\nu}{2} + k + 1\right) \left(\frac{y_j^2 + \nu}{2(1-\rho^2(h))}\right)^{-\frac{\nu}{2} - k - 1} 2F1\left(\frac{\nu + 1}{2} + k, \frac{\nu + 1}{2} + k; \frac{\rho^2(h) y_j^2 y_j^2}{(y_j^2 + \nu)(y_j^2 + \nu)}\right)
\]

\[
+ b_2 \Gamma\left(\frac{\nu}{2} + k + 1\right) \left(\frac{y_j^2 + \nu}{2(1-\rho^2(h))}\right)^{-\frac{\nu}{2} - k - 1} 2F1\left(\frac{\nu + 1}{2} + k, \frac{\nu + 1}{2} + k; \frac{\rho^2(h) y_j^2 y_j^2}{(y_j^2 + \nu)(y_j^2 + \nu)}\right)
\]

finally, combining equations (A.12), (A.10) and (A.9), we obtain

\[
f_{Y_j}(y_j) = \frac{\nu![(y_j^2 + \nu)(y_j^2 + \nu)]^{-\nu/2} 2\Gamma^2 \left(\frac{\nu}{2}\right) (1-\rho^2(h))^{-\nu/2}}{\pi^{\nu+2}(1-\rho^2(h))^{-\nu/2}} \sum_{k=0}^{\infty} \frac{\left(\frac{\nu}{2} + 1\right)^2}{k!} \left(\frac{\nu^2 \rho^2(h)}{(y_j^2 + \nu)(y_j^2 + \nu)}\right)^k
\]

\[
\times 2F1\left(\frac{\nu + 1}{2} + k; \frac{\nu + 1}{2} + k; \frac{\rho^2(h) y_j^2 y_j^2}{(y_j^2 + \nu)(y_j^2 + \nu)}\right)
\]

\[
+ \frac{\rho(h) y_j y_j \nu^{\nu/2} (y_j^2 + \nu)(y_j^2 + \nu)}{2\pi (1-\rho^2(h))^{-\nu/2}} \sum_{k=0}^{\infty} \frac{\left(\frac{\nu}{2} + 1\right)^2}{k!} \left(\frac{\nu^2 \rho^2(h)}{(y_j^2 + \nu)(y_j^2 + \nu)}\right)^k
\]

\[
\times 2F1\left(\frac{\nu}{2} + k + 1; \frac{\nu}{2} + k + 1; \frac{\rho^2(h) y_j^2 y_j^2}{(y_j^2 + \nu)(y_j^2 + \nu)}\right)
\]

and using (4.3) we obtain theorem 2.3

\[
\square
\]

A.4 Proof of Theorem 3.1

Proof. Consider \( \mathbf{U} = (U(s_1), \ldots, U(s_n))^T, \mathbf{V} = (|X_1(s_1)|, \ldots, |X_1(s_n)|)^T, \mathbf{Q} = (X_2(s_1), \ldots, X_2(s_n))^T \) where \( \mathbf{X}_k = (X_k(s_1), \ldots, X_k(s_n))^T \sim N_n(0, \Omega) \), for \( k = 1, 2 \), which are assumed to be independent. By definition of the skew-Gaussian process in (3.1) we have:

\[
\mathbf{U} = \mathbf{\alpha} + \eta \mathbf{V} + \omega \mathbf{Q}
\]

where, by assumption \( \mathbf{V} \) and \( \mathbf{Q} \) are independent. Thus, by conditioning on \( \mathbf{V} = \mathbf{v} \), we have \( \mathbf{U} | \mathbf{V} = \mathbf{v} \sim N_n(\mathbf{\alpha} + \eta \mathbf{v}, \omega^2 \Omega) \), from which we obtain

\[
f_{\mathbf{U}}(\mathbf{u}) = \int_{\mathbb{R}^n} \phi_n(\mathbf{u}; \mathbf{\alpha} + \eta \mathbf{v}, \omega^2 \Omega) f_{\mathbf{V}}(\mathbf{v}) d\mathbf{v}
\]

38
To solve this integral we need \( f_V(v) \), i.e., the joint density of \( V = (|X_1(s_1)|, \ldots, |X_1(s_n)|)^T \). Let \( X_k = (X_1, \ldots, X_n)^T = (X_1(s_1), \ldots, X_1(s_n))^T \) and \( V = (|X_1|, \ldots, |X_n|)^T \).

Additionally, consider the diagonal matrices \( D(l) = \text{diag}(l_1, \ldots, l_n) \), with \( l = (l_1, \ldots, l_n) \in \{-1, +1\}^n \), which are such that \( D(l)^2 \) is the identity matrix. Since \( l \circ v = D(l)v \) (the componentwise product) and \( X \sim N_n(0, \Omega) \), we then have

\[
F_V(v) = P_r(V \leq v) = P_r(|X| \leq v) = P_r(-v \leq X \leq v) = \sum_{l \in \{-1, +1\}^n} (-1)^{N_\omega} \Phi_n(D(l)v; 0, \Omega), \quad (N_\omega = \sum_{i=1}^n l_i \det\{D(l)\})
\]

\[
= \sum_{l \in \{-1, +1\}^n} \det\{D(l)\}\Phi_n(D(l)v; 0, \Omega)
\]

Hence, by using that

\[
\frac{\partial^n \Phi_n(D(l)v; 0, \Omega)}{\partial v_1 \cdots \partial v_n} = \det\{D(l)\}\Phi_n(D(l)v; 0, \Omega)
\]

we find that the joint density of \( V \) is

\[
f_V(v) = \sum_{l \in \{-1, +1\}^n} \left[ \det\{D(l)\}\right]^2 \phi_n(D(l)v; 0, \Omega)
\]

\[
= \sum_{l \in \{-1, +1\}^n} \phi_n(D(l)v; 0, \Omega), \quad (\left[\det\{D(l)\}\right]^2 = 1)
\]

\[
= \sum_{l \in \{-1, +1\}^n} |\det\{D(l)\}| \phi_n(v; 0, \Omega_l), \quad (\Omega_l = D(l)\Omega D(l) = (l_i l_j \rho_{ij}))
\]

\[
= \sum_{l \in \{-1, +1\}^n} \phi_n(v; 0, \Omega_l), \quad (|\det\{D(l)\}| = 1)
\]

\[
= 2 \sum_{l \in \{-1, +1\}^n \cap \Omega = \Omega} \phi_n(v; 0, \Omega_l)
\]

where the last identity is due to \( \Omega_{-\ell} = D(-\ell)\Omega D(-\ell) = D(\ell)\Omega D(\ell) = \Omega_l \) for all \( \ell \in \{-1, +1\}^n \); e.g. for \( n = 3 \) the sum must be performed on

\[
l \in \{ (+1, +1, +1), (+1, +1, -1), (+1, -1, +1), (-1, +1, +1) \}
\]

since

\[
-l \in \{ (-1, -1, -1), (-1, -1, +1), (-1, +1, -1), (+1, -1, -1) \}
\]

and both sets produce the same correlation matrices. The joint density of \( U \) is thus
given by

\[
f_U(u) = 2 \sum_{w \in \{-1, +1\}^n, w \neq -w} \int \phi_n(u; \alpha + \eta v, \omega^2 \Omega) \phi_n(v; 0, \Omega_l) dv
\]

\[
= 2 \sum_{w \in \{-1, +1\}^n, w \neq -w} \phi_n(u; \alpha, A_l) \int \phi_n(v; c_l, B_l) dv
\]

\[
= 2 \sum_{w \in \{-1, +1\}^n, w \neq -w} \phi_n(u; \alpha, A_l) \Phi_n(c_l; 0, B_l)
\]

where \( A_l = \omega^2 \Omega + \eta^2 \Omega_l, c_l = \eta \Omega_l A_l^{-1}(u - \alpha), B_l = \Omega_l - \eta^2 \Omega_l A_l^{-1} \Omega_l, \) and we have used the identity \( \phi_n(u; \alpha + \eta v, \omega^2 \Omega) \phi_n(v; 0, \Omega_l) = \phi_n(u; \alpha, A_l) \phi_n(v; c_l, B_l) \) which follows straightforwardly from the standard marginal-conditional factorizations of the underlying multivariate normal joint density.

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