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► **To cite this version:**

Mestapha Oumouni, Franck Schoefs, Bruno Castanier. Spatio-temporal modelling of degradation processes through stochastic Gamma and Gaussian processes. 27th European Safety and Reliability Conference (Esrel 2017), 2017, Portoroz, Slovenia. pp.789-796, 10.1201/9781315210469 . hal-02525524

HAL Id: hal-02525524

<https://univ-angers.hal.science/hal-02525524>

Submitted on 14 Aug 2020

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Spatio-temporal modelling of degradation processes through stochastic Gamma and Gaussian processes

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ABSTRACT: In this work, we propose a new meta-model to describe the evolution of the degradation of infrastructures for reliability purposes. The model takes into account both hazards, temporal and spatial. It is based on the classical homogenous or non-homogenous Gamma process, the scale parameter is modeled with log-normal random field in order to assess and incorporate the effect of the spatial variability and heterogeneity on the degradation process.

Simulation results are performed using synthetic data based on Monte-Carlo simulations to fit the model, the inference of their parameters is performed using the method of moments combined with curve fitting method. Computation and estimation of quantities of interests for reliability or maintenance studies, namely the failure time and remaining life-time are developed and illustrated by numerical examples.

1 INTRODUCTION

Mathematical models based on partial differential equations with stochastic parameters and data are extensively studied in Mechanical and Civil Engineering [1, 8, 2] to compute physical quantities varying in space and in time under uncertainty and spatial variability. However, their use in a reliability estimation context is faced with two major drawbacks.

Firstly, the approximation of such models to compute some physical quantities of interest can suffer from the curse of dimensionality, where it requires to solve a large number of deterministic problems. In particular when the variability and uncertainties are important [1, 6]. Secondly, health-monitoring data which are usually given by Non Destructive Techniques are not obviously linked to these models to update them with their associated parameters.

The Gamma process [12, 7, 5] is widely used model for modeling degradation process encountered in civil engineering. However, it model only temporal variability and assume a uniform degradation on the whole structure and does not incorporate heterogeneity and spatial variability through component. On the other hand, to construct a complete degradation model with accurate predictions, a large amount of data using destruc-

tive or nondestructive testing is required from a large amount of structures [12, 10, 9, 5] for predicting levels of the degradation with accuracy.

The major contribution detailed in this work is a new spatio-temporal random model based on Gamma process for predicting the degradation mechanism which takes into account both hazards, temporal and spatial. The temporal variability is modeled by the classical Gamma process and the spatial variability is modeled by a positive random field. This spatial field follows a log-normal distribution to describe the scaling parameter of the Gamma process. Under the assumption of stationarity satisfied by the random field, the spatial monitoring data of the component contributes in the parameters estimation to reduce uncertainty and increase the accuracy of the meta-model approach. Therefore, the method of moments based on the variogram curve fitting is used in the first stage to estimate the spatial parameters of the Gaussian field, in the second stage the method is reused to estimate temporal parameters.

The article is organized as follows, the spatio-temporal random field degradation model is presented in Section 2. The method of moments combined with curve fitting method are presented in Section 3 for identifying properties of the model in terms of statistical inference. Once the model is adjusted, Section 4 develops quantities of interests

which are useful in the reliability and maintenance analysis, Namely, the failure time defined in terms of degradation level passages and remaining useful life time defined as the time of inspection of the unit to failure. Finally, Section 5 presents a numerical example of one dimensional variability illustrating the proposed methodology for model validation. Once the model is adjusted, Quantities of interest are developed and illustrated by an analytic and sampling approach.

2 SPATIO-TEMPORAL DEGRADATION MODEL

We look for modeling the degradation process by a convenient spatio-temporal random field to take into account its aleatory evolution with time and space. A separable model is one simple spatio-temporal model obtained through the tensorial product between a merely stochastic process $(X_t)_{t \geq 0}$ and a spatial random field $Z(z)$, where z is the spatial variable. This class of separable random field is extensively used even in situations in which they are not always physically justifiable, since separability gives important computational and mathematic benefits.

The evolution in time models the intrinsic aleatory and the Gamma process is a natural candidate to catch this monotonous degradation. The spatial random field models the variability and uncertainty through the structure, and in many applications it is classically modeled by a second order stationary random field and given by a transformation of a Gaussian random field. Therefore, for simplicity and in order to construct a separable model, we assume that those source of randomness (time and space) are independent. Therefore, we consider a separable spatio-temporal random field to model the spatial variability in the degradation process.

Consider $\alpha(\cdot)$ to be a non-decreasing, right-continuous, real-valued function for $t \geq 0$ and vanishing at $t = 0$ and $\beta > 0$ a positive constant. A stochastic process $(X_t)_{t \geq 0}$ is said to be a Standard Gamma process with function shape $\alpha(\cdot)$ and identical scale parameter β if it satisfies the following properties

- $X_0 = 0$ with probability one,
- $X_{t+s} - X_t \sim Ga(\alpha(t+s) - \alpha(t), \beta)$,
- X_t has independent positive increments,

Where $Ga(\alpha(\cdot), \beta)$ is the Gamma distribution defined by the density function $f_{Ga}(x)$:

$$f_{Ga}(x) = \frac{\beta^{\alpha(\cdot)}}{\Gamma(\alpha(\cdot))} x^{\alpha(\cdot)-1} e^{-\beta x}, \text{ for each } x > 0,$$

and Γ is the classical Gamma function. The process X_t satisfies the following scaling property,

$$X_t \sim Ga(\alpha(\cdot), 1) \beta^{-1}, \text{ for each } \beta > 0. \quad (1)$$

The scaling property of the gamma process in (1) motivates us to consider the scale parameter to be a spatial random field to obtain a separable spatio-random filed model,

$$G_t(z) := X_t \beta(z)^{-1}, \quad (2)$$

where $\beta(\cdot)$ is spatial and positive random field which is assumed to be independent of X_t , and (without loss of generality we consider) $X_t \sim Ga(\alpha(\cdot), 1)$. The scaling property satisfied by the process X_t in (1) suggests to see (formally) the spatio-random field $G_t(z)$ as Gamma process with spatial random scale $\beta(\cdot)$.

In practice, it is difficult to verify and to find a positive distribution for a spatial or spatio-temporal random field. However, log-normal distribution occurs naturally as a limit distribution of physical processes, because the Central Limit Theorem applied to the product of positive independent random variables (number of measures > 30) ensures that the log normal distribution can occur. Therefore, we choose the Log-normal distribution for the random scale coefficient,

$$\beta(z) = e^{Y(z)},$$

where Y is the spatial Gaussian random field $Y(z)$ which defined in D a set in R^d with $d = 1, 2, 3$. The field Y is assumed to be homogenous (stationary field) and then completely defined [14] by its constant mean $\mu := \mathbb{E}[Y(\cdot)]$ and its stationary covariance function $cov(r) := \mathbb{E}[Y(z+r)Y(z)] - \mu^2$.

The Matérn model of the covariance functions are the commonly used covariance in the engineering applications for the Gaussian random field. There are defined by the following function:

$$cov(r) = \sigma^2 \frac{2^{1-\nu}}{\Gamma(\nu)} \left(\frac{\sqrt{2\nu}r}{l_c} \right)^\nu K_\nu \left(\frac{\sqrt{2\nu}r}{l_c} \right) \quad (3)$$

where the parameters σ^2 , ν , α and l_c are non-negative real numbers, σ^2 is the variance of Y , l_c is the correlation length, r is the Euclidean distance between two points, K_ν denotes the modified Bessel function of the second kind. The parameter $\nu > 0$ is a non-negative number that characterizes the degree of smoothness of cov which is related to the smoothness of the field Y . when $\nu = \frac{1}{2}$, cov coincides with the exponential covariance,

$$c(r) := e^{-r/l_c}.$$

this covariance is only Hölder continuous and so for the samples paths of Y . When $\nu \rightarrow \infty$, it approaches the gaussian covariance,

$$c(r) := e^{-r^2/(2l_c^2)},$$

which is an analytic function and so the samples paths of Y are also analytic.

3 METHOD OF MOMENTS FOR PARAMETERS INFERENCE

In order to obtain a complete and accurate degradation model for practical examples, statistical methods for parameters estimation of gamma processes and gaussian random field are required. A typical data set of $G_t(z)$ consists of inspection points in different increasing time t_j for $j = 0, \dots, N_t$, with the same period τ , where for each time t_j the inspection positions are given in uniform positions z_l for $l = 1, \dots, N_z$ with step size h . Here, we assume for simplicity that $\alpha(\cdot)$ follows a power law:

$$\alpha(t) = at^b, \quad (4)$$

for some unknown $a > 0$ and a known power $b = 1$.

3.1 Step 1: Spatial parameters

The Method of Moments (MOM) uses the benefit of the separability of the model to estimate the parameters of $G_t(z)$ in two steps. The first step of (MOM) consists in estimating the spatial parameters of the model. They are defined as the parameters of the second order stationary random field Y , i.e, the variance σ^2 , the correlation length l_c and the regularity parameter ν of the correlation function given in (3).

In order to obtain an estimate of these parameters, we use nonparametric estimate of the semi-variogram which is typically obtained through (MOM) of one or more realizations of Y and defined by [3]:

$$\hat{Y}_Y(h_t) = \frac{1}{2N_{h_t}} \sum_{z_i, z_j \in S_{h_t}} (Y(z_i) - Y(z_j))^2, \quad (5)$$

where S_{h_t} is the set of the points separated with distance h_t and N_{h_t} its cardinal.

For a fixed time instant t , the variogram of the field $\log(G_t(\cdot))$ equals the variogram of $Y(\cdot)$. Therefore, the empirical variogram \hat{Y}_Y is obtained by computing the empirical variogram of the random field $\log(G_t)$, the logarithm of G_t . Thus, in any fixed time t we get:

$$\hat{Y}_Y(h_t) = \frac{1}{2N_{h_t}} \sum_{z_i, z_j \in S_{h_t}} (\log(G_t(z_i)) - \log(G_t(z_j)))^2 \quad (6)$$

When M realizations of the degradation $G_t(z)$ are available, they can be used to improve the estimation of $\hat{Y}_Y(h_t)$ by empirical average.

Since the random field Y is stationary, the exact variogram is given by:

$$Y_Y(h_t) = \sigma^2 - \sigma^2 \text{cov}(h_t), \quad (7)$$

where the correlation function $\text{cov}(h)$ is given in (3). Therefore, we estimate the spatial parameters by minimizing the quadratic error (Least square method) between exact variogram and experimental one, thus σ^2 , l_c and ν are deduced by the following minimization problem.

$$\min_{\sigma, l_c, \nu > 0} \sum_{l=1}^{N_z} (\hat{Y}_Y(h_l) + \sigma^2 \text{cov}(h_l) - \sigma^2)^2 \quad (8)$$

Remark 3.1. *The classical least square method can be extended to a generalized least squares (GLS) method, where we minimize a weighted error, given by the correlation matrix R of the set $\{\hat{Y}_Y(h_l)\}_{l=1}^{N_z}$. A simplified approach of this generalization is the weighted method where the matrix R is diagonal with entries $R_l = \frac{2Y_Y(h_l)}{N_{h_l}}$ as suggested in [3] assuming a Gaussian law and non-correlation among $\{\hat{Y}_Y(h_l)\}_{l=1}^{N_z}$.*

We note when the space positions are not equidistant, estimation (5) is slightly modified for non uniform grid to compute the experimental variogram of Y from the logarithm of the field $G_t(\cdot)$, spatial mean is given on all pairs of points whose distance are between h and $h + \delta h$ for some threshold $\delta h > 0$ [3].

3.2 Step 2: Temporal parameters

The temporal parameters are defined as the parameters of the process X_t , i.e, the shape function $\alpha(t)$ defined in (4) and the deterministic contribution $\eta = e^\mu$ of the scale field β . Since we assume that the power b is known, we estimate then a and η . Note that η contains a contribution of the spatial mean of the random field Y . Let $\{G_{t_j}^i(z_l)\}$ be a sequence of independent and identically distributed (i.i.d.) simulations of the degradation model described in the previous section. Each i -th degradation process is observed at time t_j among N_t times and on location x_l among N_z locations.

Logarithm of increments of sample paths of $G_t(z)$ writes:

$$\xi_{j,l}^i := \log(\delta G_{t_j}^i(z_l)) = \log(X_{t_j}^i - X_{t_{j-1}}^i) + Y^i(z_l),$$

for $i \in [1, M]$, $j \in [1, N_t]$, $l \in [1, N_z]$.

Since the process X_t and the field Y are independent, the first two moments of $\xi_{j,l}^i$ are given by:

$$m_1 := \mathbb{E}[\xi_{j,l}^i] = \psi(a\tau_j) - \log(\eta) \quad (9)$$

$$m_2 := \text{var}[\xi_{j,l}^i] = \psi_1(a\tau_j) + \sigma^2, \quad (10)$$

where the function ψ is digamma function which is defined as the logarithmic derivative of gamma function Γ , and ψ_1 is the trigamma function defined as the derivative of ψ . The temporal grid being uniform, thus increments $(X_{t_j}^i - X_{t_{j-1}}^i)_{j=1}^{N_i}$ are independent and identically distributed. From MN_t realizations of those increments combined with spatial average we estimate m_1 and m_2 by the following sum:

$$m_1 \approx (MN_z N_t)^{-1} \sum_{i=1}^M \sum_{j=1}^{N_i} \sum_{l=1}^{N_z} \xi_{j,l}^i, \quad (11)$$

$$m_2 \approx (MN_z N_t)^{-1} \sum_{i=1}^M \sum_{j=1}^{N_i} \sum_{l=1}^{N_z} (\xi_{j,l}^i)^2 - m_1^2. \quad (12)$$

Therefore, an estimation of temporal parameters a and η is given by:

$$\begin{pmatrix} a \\ \eta \end{pmatrix} = f^{-1} \begin{pmatrix} m_1 \\ m_2 \end{pmatrix}, \quad (13)$$

where we define the function f by

$$f \begin{pmatrix} u \\ v \end{pmatrix} = \begin{pmatrix} \psi(u\tau) - \log(v) \\ \psi_1(u\tau) + \sigma^2 \end{pmatrix}, \quad (14)$$

with $\tau = t_j - t_{j-1}$ and m_1 et m_2 are estimated in (11, 12) and σ^2 is the estimate variance σ^2 obtained in equation (8).

Remark 3.2. When b is unknown, we can use the approach of pseudo maximum-likelihood method (PML) to estimate temporal parameters. It consists of maximizing the likelihood of the sequence of the increment $(\delta_i G, \&, \delta_{iN} G)$ on a given fixed spatial position z . These increments are conditionally independent and their likelihood is given by the product of the marginal density of the field $G_i(z)$. This marginal density is approximated in Section 4 (equation (15) and (16)).

Method of moments for temporal parameters can be extended when $b \neq 1$, the non-stationary Gamma process can be easily transformed to a stationary one by performing a monotonic transformation $u(t) = t^b$ on the time increments (see [12] for more details).

4 QUANTITIES OF INTEREST

4.1 Marginal density approximation

The marginal distribution noted by $f_t(\cdot)$ of the model $G_t(z)$ does not depend on the position z but only on the time t , because the spatial random field Y is homogenous. This distribution can be

constructed numerically by simulating paths of the field $G_t(z)$. It is useful in computation of some quantities of interest which are used in structural reliability and maintenance. We can also use this distribution as a pseudo-likelihood of $G_t(z)$ to conclude an estimation of the temporal parameters of the model.

Let $\xi(y)$ be the density of the standard gaussian random variable $N(0,1)$. By using the fact that the temporal paths of increments of the model $G_t(\cdot)$ are conditionally independent, we compute the distribution f_t for all $v > 0$ by the following form:

$$f_t(v) = \psi_t(v) \int_{\mathbb{R}} \exp(-v\eta \exp(\sigma y) + \alpha(t)\sigma y) \xi(y) dy, \quad (15)$$

where we set $\psi_t(v) := \frac{v^{\alpha(t)+\alpha(t)-1}}{\Gamma(\alpha(t))}$.

The integral in (15) has a transcendental form, thus we use Gauss-Hermite quadrature formula to approximate this marginal density f_t .

We consider m roots $\{y_j\}_{j=1}^m$ of the Hermite polynomial and their associated weights $\{w_j\}_{j=1}^m$. Thus, an approximation of f_t writes:

$$f_t^m(v) := \psi_t(v) \sum_{j=1}^m \exp(-v\eta \exp(\sigma y_j) + \alpha(t)\sigma y_j) w_j \quad (16)$$

The convergence of the sequence $f_t^m(v)$ for any fixed positive reel $v > 0$ is relatively fast, since the integrand is infinitely differentiable. However the norm of any m -derivative of this integrand depends on the value of the parameters $\alpha(t)$, σ and η . A large value of these parameters requires a large order m of the approximation in (16), in particular for large time t . Therefore, the order of the approximation m is established by the following stop criterion,

$$|f_t^m - f_t^{m-1}| \leq \varepsilon,$$

where $\varepsilon > 0$ is a convenient threshold value. Note that when $v \rightarrow \infty$ both $f_t(v)$ and $f_t^m(v)$ decreases to zero.

4.2 Distribution of failure time

The failure time T_F for a structural component is defined as the time at which the degradation path G_t first crosses a critical level g_F for any spatial location,

$$T_F = \inf\{t > 0; G_t(\cdot) \geq g_F\},$$

In what follows, the critical level g_F is assumed deterministic. For some simple path models, the distribution of T_F defined by $F_T(t) := P(T_F < t)$ can

be expressed in a closed form. However, this is not always possible and it can be numerically computed using Monte Carlo simulations by simulating paths of $G_t(z)$. The sample paths of G are monotonic since the spatial variability e^Y is positive, thus the failure time cumulative distribution function $F_T(T)$ satisfies:

$$F_T(t) = 1 - P(T_F > t) = 1 - \int_0^{g_F} f_t(z) dz, \quad (17)$$

where f_t is the marginal pdf of G_t given in (15). Therefore, using the approximation f_t^m with a convenient order m we approximate the distribution F_T ,

$$F_T^m(t) = 1 - \int_0^{g_F} f_t^m(v) dv, \quad (18)$$

The integral (18) can be computed accurately by any quadrature formula, for example the Legendre-Gauss quadrature.

The derivative of (17) and (18) with respect to the time t provides the probability density function of T_F and its approximation respectively.

The approximation of the cumulative density (18) requires a quite huge cost when t or σ^2 is large. In this case one can construct an estimation of F_T by generating a sufficiently large number of random sample paths of $G_t(z)$ with estimated parameters using Monte-Carlo (MC) simulations.

By fixing N_t desired times, N_z desired locations and M realisations of spatio-temporal paths $\{G_{t_i}^m(z_j)\}$ for $i=1 \& , N_t$, $j=1 \& , N_z$ and $m=1, \& , M$. The estimate \tilde{F}_T of F_T is given by,

$$\tilde{F}_T(t_i) := \frac{\sum_{j,m}^{N_z, M} \mathbb{I}_{\{G_{t_i}^m(x_j) \geq g_F\}}}{MN_z}, \quad (19)$$

where \mathbb{I}_A represents the characteristic function of the set A , i.e $\mathbb{I}_A(z) = 1$ if $z \in A$ and zero otherwise.

Note that since the random field G_t is homogeneous with respect to the spatial variables, the estimate \tilde{F}_T can be computed also using realizations of $G_t(z_p)$ fixed at any position z_p ,

$$\tilde{F}_T(t_i) \approx \frac{\sum_m^M \mathbb{I}_{\{G_{t_i}^m(z_p) \geq g_F\}}}{M}. \quad (20)$$

However the estimation (20) needs to use more MC simulations of $G_t(z)$ than (19) since by the ergodic property, the spatial average contributes to the convergence of \tilde{F}_T to F_T .

4.3 Remaining useful lifetime after inspection

In reliability analysis and survival studies, residual lifetime after inspection is a key indicator. In the maintenance decision analysis, the current

measured degradation is used to predict the remaining useful lifetime (RL) of the structure [11]. If t is the current time of inspection, the residual lifetime is defined by the random variable:

$$RL_t := \inf\{\tau > 0, G_{t+\tau} \geq g_F \mid G_t = g_t\},$$

where g_F is the critical level and g_t is the measured degradation at given time t , implicitly $g_t < g_F$. When we suppose that a component has survived to a given time t and we have not any information or measure about the current degradation path G_t , then a conditional reliability function gives an evaluation of the remaining lifetime:

$$R(\tau \mid T_F > t) := P(T_F \geq \tau + t \mid T_F > t) = \frac{\int_0^{g_F} f_{\tau+t}(y) dy}{\int_0^{g_F} f_t(y) dy}. \quad (21)$$

When the current degradation measure path of G_t is available, then by conditioning of the failure time would give a more accurate prediction than (21), i.e the probability that the unit survival after the time $t + \tau$ given its actual state $G_t = g_t$ at time t is:

$$P(RL_t > \tau) = P(G_{t+\tau} < g_F \mid G_t = g_t) = P(\delta_t G_t < g_F - g_t \mid G_t = g_t), \quad (22)$$

where we define the increment $\delta_t G_t := G_{t+\tau} - G_t$. If we note by $f_{\delta_t G_t \mid G_t}$ the conditional marginal density of the process $\delta_t G_t$ given the event $\{G_t = g_t\}$, then we compute the probability cumulative distribution $F_{RL} := P(RL_t \leq \tau)$ of the residual lifetime by:

$$P(RL_t \leq \tau) = 1 - P(\delta G_t < g_F - g_t \mid G_t = g_t) = 1 - \int_0^{g_F - g_t} f_{\delta_t G_t \mid G_t}(y) dy. \quad (23)$$

Using the independence of the increments of Gamma process, we compute the conditional density of $\delta_t G_t$ given $G_t = g_t$ by,

$$f_{\delta_t G_t \mid G_t}(u) = \frac{u^{\delta_t \alpha - 1} g_t^{\alpha(t) - 1} f_{\tau+t}(u + g_t)}{B(\delta_t \alpha, \alpha(t))(u + g_t)^{(\alpha(\tau+t) - 1)} f_t(g_t)} \quad (24)$$

where we set $\delta_t \alpha := \alpha(t + \tau) - \alpha(t)$, f_t is the marginal density of G_t given in (15) and $B(x, y) = \frac{\Gamma(x)\Gamma(y)}{\Gamma(x+y)}$ is the beta function. It follows that $F_{RL}(\tau)$ is given by,

$$F_{RL}(\tau) = 1 - \int_0^{g_F - g_t} f_{\delta_t G_t \mid G_t = g_t}(u) du. \quad (25)$$

By using an appropriate approximations of densities $f_{t+\tau}^m(u + g_t)$ and $f_t^m(g_t)$ we get $f_{\delta_t G_t \mid G_t = g_t}^m(u)$ an approximate of the conditional probability density in (24) and then an approximation of $F_{RL}(\tau)$.

Note that when $t = 0$ and $g_t = 0$, equation (24) is not definite and the cumulative probability function F_{RL} is the function F_T given in (17).

The derivative of (23) with respect to the variable τ provides the probability density function of the Residual lifetime RL .

5 NUMERICAL ILLUSTRATION

In this section we simulate the degradation model through Monte Carlo simulations to validate the inference process and the approximation of the quantities of interest.

We consider $M = 1, 10, 100$ sample paths of the field $(G^j)_{j=1}^M$, each trajectory G^j being simulated at N_t equidistant periods on the interval time $[0, 30]$ (in years) and N_z equidistant locations in one dimensional and two dimensions space.

The model is stationary with respect to time, i.e the shape parameter is linear $\alpha(t) = at$, the deterministic contribution of the scale random field is given by $\eta = e^\mu$ where $\mu = 2/3$ (mean of Y), $\sigma^2 = 0.6$ (variance of Y), $l_c = 1$ (correlation length) and $\nu = 2$ (smoothness parameter of Y).

The Gaussian random field is defined on the interval $[0, L]$ where $L = 100$. We use the method of circulant matrix [4] to simulate Y with exact discretization in N_z equidistant spatial positions.

In what follows, for simplicity, a fixed value of the smoothness parameter $\nu = 2$ is selected (i.e the paths of G are one time derivative in quadratic norm).

Figure 1 (left) plots one realization of the model $G_t(z)$ and compares (right) its experimental variogram with exact one of the Gaussian field Y . The experimental variogram is computed using the spatial trajectory of the logarithm $\log(G_t(z))$ at time $t = 30$.

Method of moments (MOM) Step 1

The first step of MOM consists in estimating spatial parameters, variance σ^2 and correlation length

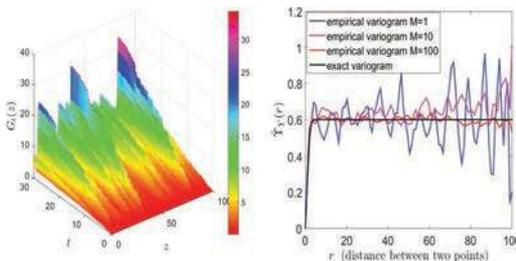


Figure 1. Left: example of one realization of $G_t(\cdot)$, right: variogram ($l_c = 1, \nu = 2, \sigma^2 = 0.6$).

Table 1. Parameters estimation of one dimensional spatial variability by MOM.

$N_z = 40$		
M	σ^2	l_c
1	0.714(0.114)	0.21(0.791)
10	0.5720(0.028)	0.81(0.19)
100	0.6048(0.0048)	0.890(0.1024)
$N_z = 100$		
M	σ^2	l_c
1	0.546(0.053)	1.186(0.186)
10	0.584(0.015)	0.934(0.065)
100	0.6009(0.009)	1.024(0.024)

l_c . The quality of the estimation is measured by the mean absolute errors which is given by the average of absolute difference between the exact parameter and 10 estimated values calculated across MOM.

Table 1 gives estimates of those parameters obtained by MOM where we minimize equation (8) with $\nu = 2$ and three values of M . (Mean absolute errors in brackets).

From Table 1, the estimation of the correlation length l_c depends strongly on the number of spatial positions N_z , obviously because l_c is very small with respect to the length $L = 100$. In contrast of l_c , the estimate of the variance σ^2 is largely acceptable even with small number of positions and with only one realization of $G_t(\cdot)$ ($M = 1, N_z = 40$). However, the estimation of quantities of interest depends strongly on the variance σ^2 , so an accurate estimate of σ^2 is needed to forecast reliable predictions.

Method of moments Step 2

Once spatial parameters are estimated, the second step of the MOM consists in estimating the temporal parameters, which are given by (13). An estimation of the variance σ^2 is inserted in (13) for each case of M and N_z . Table 2 summarizes estimations of temporal parameters a and η . Results show that their accuracy depends strongly on the total inspection times N_t and on the total number of positions N_z . An acceptable accuracy is reached for small number of realizations M ($M = 10$) when N_t and N_z are significantly large ($N_t = 60, N_z = 40$). In particular, estimate of η depends significantly on the number of locations N_z since it contains the stochastic contribution of the random field Y .

Figure 2 illustrate the convergence of the MOM for temporal parameters a and η , where we consider a database with size $M \times N_z \times N_t$ of the

Table 2. Parameters estimation of temporal variability using MOM.

$N_z = 40, N_t = 30$		
M	a	η
1	0.79(0.21)	1.487(0.46)
10	0.947(0.053)	2.157(0.21)
100	1.012(0.012)	1.986(0.038)
$N_z = 40, N_t = 60$		
M	A	η
1	0.902(0.098)	1.617(0.33)
10	1.027(0.027)	2.087(0.14)
100	1.01(0.009)	1.97(0.022)
$N_z = 100, N_t = 30$		
M	a	η
1	0.817(0.18)	2.377(0.43)
10	1.057(0.057)	1.837(0.11)
100	1.012(0.012)	2.015(0.068)
$N_z = 100, N_t = 60$		
M	a	η
1	1.01(0.1)	1.567(0.38)
10	1.029(0.029)	1.817(0.13)
100	1.01(0.009)	2.973(0.026)

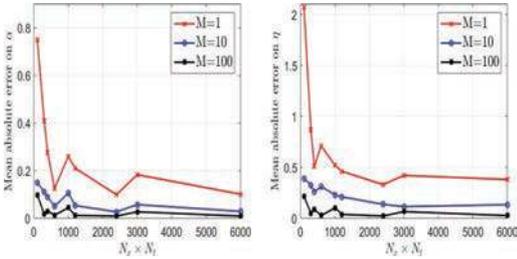


Figure 2. Mean absolute error on α (left) and η (right).

simulated model G . The mean absolute error on α and η is given as function of the product $N_z \times N_t$ where $N_z = 10, 40, 100$ and $N_t = 10, 30, 60$.

We remark a strong improvement for the estimation for $M > 1$, $N_z > 10$ and $N_t > 10$. Estimate accuracy of α is nearly independent of the spatial positions unlike η where accuracy depends on N_z and strongly on N_t . An acceptable accuracy is reached in the case of one component $M = 1$ when N_t and N_z are quite large ($N_t = 60, N_z = 40$). A large number of positions N_z does not improve the convergence of these temporal parameters unlike spatial parameters.

Once parameters estimation of the model is performed, we accurately simulate G at any position

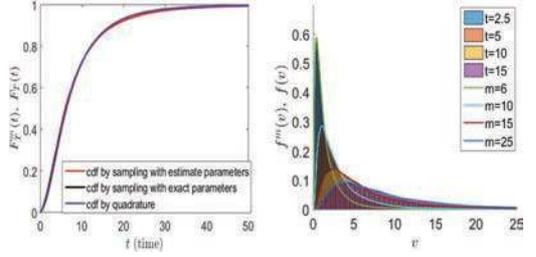


Figure 3. Left: estimate of the failure time distribution F_T , right: marginal density by sampling and quadrature approach.

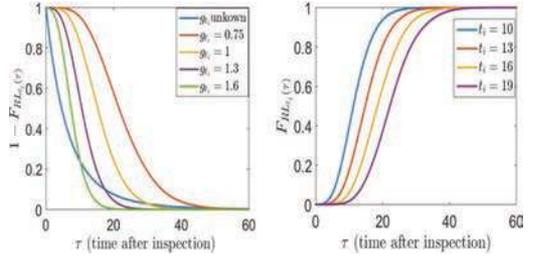


Figure 4. Left: Distribution of predictive time of inspection to failure, right: reliability and remaining lifetime function.

and time. Therefore, in order to estimate the marginal cumulative distribution F_T of failure time T_F by sampling method (19), we perform 10^3 MC simulations of $G_t(\cdot)$ discretized on $N_t = 100$ times and $N_z = 100$ positions. We plot in Figure 3 (left) the cdf F^T computed by sampling approach with exact and estimated parameters and by quadrature rule (18) with $m = 30$ Gaussian knots.

Figure 3 (right) compares marginal density of G_t by sampling and by quadrature given in (16) at several times t with a convergence criterion $\varepsilon \approx 10^{-3}$. We note that the required order of the quadrature rule increases with time t .

Figure 4 (left) illustrates the effect of conditioning the current state on the failure prediction at the current time of inspection $t_i = 10$ for several measured degradations levels from 0.75 to 1.6.

The survival function depends on the current state and gives more prediction than the reliability function (g_{t_i} is unknown, blue curve). The curves highlight how the reliability function underestimates or overestimates the time to failure under the value of g_{t_i} . In Figure 4 (right), we draw several failure curves of the remaining lifetime $F_{RL_{t_i}}(\tau)$ with several times of inspection t_i with the same condition state $g_{t_i} = 1.25$. This figures highlight an obvious result, the more the time for observing given condition state, the more the residual lifetime.

6 CONCLUSION

In this paper, we have developed a spatio-temporal degradation model that incorporates the spatial variability and heterogeneity across structural component. It is based on the basic gamma process with a scale parameter modeled with non-negative spatial random field. The temporal paths of the process are monotonic with conditionally independent increments, the positive random field scale follows a log-normal distribution as limit of independent positive variables.

The quantities of interest that are useful in reliability analysis and in maintenance, namely the distribution of failure time and the distribution of remaining useful lifetime are computed. A Method of moments is carried out to infer the spatial and temporal parameters of the model. Monte Carlo simulations illustrate the advantage of the proposed model. The advantages of the proposed model are that uncertainties are reduced and the accuracy of the inference is improved by exploiting with better manner the spatial data.

An interesting extension of the current model can consider the bivariate modeling based on the state dependent Gamma process.

ACKNOWLEDGEMENTS

The work reported in this article was supported by the Region Pays de la Loire de France through the regional project SI3M (Strategies of Identification of Meta-models for Structural Maintenance). The authors would like to thank the region for the support.

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