1 Isotropic Covariance Functions

Let $\{Z(s)\}$ be a Gaussian process on \mathbb{R}^n , *i.e.*, a collection of jointly normal random variables Z(s) associated with *n*-dimensional locations $s \in \mathbb{R}^n$. The joint distribution of $\{Z(s)\}$ depends only on the means $\mu(s) = \mathsf{E}Z(s)$ and the covariances $C(s,t) = \mathsf{E}(Z(s) - \mu(s))(Z(t) - \mu(t))$.

The process is called *stationary* or *translation invariant* if the distribution wouldn't change under a rigid translation of the entire collection of locations, *i.e.*, if $\mu(s) = \mu(s+h)$ and C(s+h,t+h) = C(s,t) for all h; in this case $\mu(s) \equiv \mu$ is constant and C(s,t) = C(s-t,0) can only depend on the difference h = (s-t) between the two locations, so must be of the form $C(s,t) = C_0(s-t)$ for some function $C_0(h) = C(h,0)$ on \mathbb{R}^n . Not just any function $C_0(h)$ can be a covariance function; let's see what the choices are.

It's easy to see that the function C_0 must be even, i.e., must satisfy $C_0(h) = C_0(-h)$, since $C(s-t) = \mathsf{E}(Z(s) - \mu(s))(Z(t) - \mu(t)) = C(t-s)$. But more is true: if $\{s_j\}$ any collection of locations, then complex linear combinations $a^{\mathsf{T}}(Z-\mu) = \sum a_j(Z_j - \mu_j)$ of the centered random variables $Z_j = Z(s_j)$ (with means $\mu_j = \mu(s_j)$) must have nonnegative squared modulus $\mathsf{E} |\sum a_j(Z_j - \mu_j)|^2 = \sum a_jC(s_j - s_k)\bar{a}_k \geq 0$ for every set of complex numbers $\{a_j\} \subset \mathbb{C}$. A function $C_0(h)$ is called *positive semi-definite* if it always satisfies the inequality $\sum_{jk} a_jC(s_j - s_k)\bar{a}_k \geq 0$ for any locations s_j and complex numbers a_j ; this is equivalent to asking that C(h) = C(-h) for every $h \in \mathbb{R}^n$ and that $\sum a_jC(s_j - s_k)a_k \geq 0$ for all real numbers $a_j \in \mathbb{R}$. One way to get a symmetric positive semi-definite function $C_0(h)$ is by taking the Fourier transform

$$C_0(h) = \int_{\mathbb{R}^n} e^{ih \cdot \omega} G(\omega) \, d^n \omega$$

of any positive function $G(\omega)$ on \mathbb{R}^n or, more generally, of any finite positive measure $G(d\omega)$, because then

$$\sum_{jk} a_j C(s_j - s_k) \bar{a}_k = \int_{\mathbb{R}^n} \sum_{jk} (a_j e^{s_j \cdot \omega}) \overline{(a_k e^{s_k \cdot \omega})} G(d\omega)$$
$$= \int_{\mathbb{R}^n} \left| \sum_j a_j e^{s_j \cdot \omega} \right|^2 G(d\omega) \ge 0.$$

It turns out that this is the *only* way to get one— that every positive semidefinite function can be written in this form for some finite positive measure $G(d\omega)$, called the *spectral measure* (if $G(d\omega) = G(\omega) d\omega$ is absolutely continuous, $G(\omega)$ is called the *spectral density*). Known as "Bochner's Theorem," this result is really just the Fourier inversion formula in an unfamiliar setting:

$$G(\omega) = (2\pi)^{-n} \int_{\mathbb{R}^n} e^{-ih \cdot \omega} C_0(h) \, d^n h.$$

Since the process $\{Z(s)\}$ is real-valued, the spectral density $G(\omega) = G(-\omega)$ must be an even function and so we can write

$$C_0(h) = \int_{\mathbb{R}^n} \cos(h \cdot \omega) G(\omega) d^n \omega$$

$$G(\omega) = (2\pi)^{-n} \int_{\mathbb{R}^n} \cos(h \cdot \omega) C_0(h) d^n \omega$$

If the Gaussian process is also *isotropic*, or invariant under rotations, then $G(\omega) = g(|\omega|)$ must also be invariant under rotations and depend only on the length $r = |\omega|$ of the vector $\omega \in \mathbb{R}^n$. In this case we can simplify these integrals by transforming to polar coordinates.

1.1 Polar Coordinates for Probabilists

Polar coordinates are a familiar tool in two-dimensional integrals, where the change of variables from $x \in \mathbb{R}^2$ to $r = \sqrt{x_1^2 + x_2^2}$ and $\theta = \arctan x_2/x_1$ (so $x_1 = r \cos \theta$, $x_2 = r \sin \theta$) and a change from d^2x to $r \, dr \, d\theta$ lead to simple expressions for the integrals of radial functions. Equivalently, we can let σ have a uniform probability distribution (denoted by $d\sigma$) over the unit circle $S^1 = \{x : x_1^2 + x_2^2 = 1\}$, and change variables from $x = (x_1, x_2)$ to (r, σ) , with $d^2x = dx_1 \, dx_2$ replaced by $2\pi r \, dr \, d\sigma$.

In three dimensions the first polar approach has its analogue in the Euler angles, while the second is simpler with uniform measure for σ on the unit sphere $S^2 \subset \mathbb{R}^3$, with $d^3x = dx_1 dx_2 dx_3$ replaced by $4\pi r^2 dr d\sigma$. Notice that $2\pi r$ and $4\pi r^2$ are the circumference of the circle and the area of the sphere of radius r, respectively. In any number n of dimensions the sphere S^{n-1} has area $2\pi^{n/2}r^{n-1}/\Gamma(n/2)$, and we can again again evaluate integrals in polar coordinates with the uniform probability distribution $d\sigma$ for $\sigma \in S^{n-1} \subset \mathbb{R}^n$, and $d^n x = \frac{2\pi^{n/2}}{\Gamma(n/2)}r^{n-1} dr d\sigma$. This makes it easy to compute integrals of radial functions; for functions that also depend on one or more of the components x_j , it is sometimes helpful to note that the squares $\{\sigma_j^2\}$ have a Dirichlet $\text{Di}(\frac{1}{2}, ..., \frac{1}{2})$ joint distribution, so each σ_j is distributed as the square root of a $\text{Be}(\frac{1}{2}, \frac{n-1}{2})$ random variable.

1.2 Evaluating $C_0(h)$

Switching to polar coordinates $r = |\omega| \ge 0$ and $\sigma = \omega/|\omega| \in S^{n-1}$ (where $d\sigma$ denotes the uniform probability measure on the unit sphere S^{n-1} in \mathbb{R}^n), and noting that the component $\sigma_h = \sigma \cdot h/|h|$ of $\sigma \in S^{n-1}$ in the direction h again has the same distribution as the square root of a $\mathsf{Be}(\frac{1}{2}, \frac{n-1}{2})$ random variable, writing ρ for |h|,

$$C_{0}(h) = \int_{\mathbb{R}^{n}} \cos(h \cdot \omega) g(|\omega|) d^{n} \omega$$

$$= \iint_{\mathbb{R}_{+} \times S^{n-1}} \cos(r\rho\sigma_{h}) g(r) \frac{2\pi^{n/2} r^{n-1}}{\Gamma(n/2)} dr d\sigma$$

$$= \int_{\mathbb{R}_{+}} \int_{0}^{1} \cos(r\rho\sqrt{u}) g(r) \frac{2\pi^{n/2} r^{n-1}}{\Gamma(n/2)} \frac{\Gamma(n/2)}{\Gamma(\frac{1}{2}) \Gamma(\frac{n-1}{2})} u^{1/2-1} (1-u)^{(n-1)/2-1} dr du$$

$$= \int_{0}^{\infty} \rho (2\pi r/\rho)^{\nu+1} J_{\nu}(r\rho) g(r) dr, \qquad \nu \equiv \frac{n}{2} - 1 \qquad (1)$$

$$= \int_{0}^{\infty} (r\rho/2)^{-\nu} \Gamma(\nu+1) J_{\nu}(r\rho) \gamma(dr)$$
(2)
=
$$\begin{cases} \int_{0}^{\infty} 2\cos(r\rho) g(r) dr & \text{if } n = 1 \\ \int_{0}^{\infty} 2\pi r J_{0}(r\rho) g(r) dr & \text{if } n = 2 \\ \int_{0}^{\infty} \rho(2\pi r/\rho)^{3/2} J_{1/2}(r\rho) g(r) dr & \text{if } n = 3 \end{cases}$$

where

$$J_{\nu}(z) = \frac{(z/2)^{\nu}}{\sqrt{\pi} \,\Gamma(\nu + 1/2)} \int_0^{\pi} \cos(z\cos\theta) \,\sin(\theta)^{2\nu} \,d\theta$$

is the Bessel function of the first kind of order ν (see Watson, 1944). Bessel functions aren't as familiar as sines and cosines, but they're common in engineering and physics and are in the standard C library, the GNU Scientific library (GSL), Maple and Mathematica, Matlab, *etc.*; see Abramowitz and Stegun (1964, Chapter 9) for details. Here's a plot of $J_0(z)$:



The plot of $J_0(z)$ looks a little like a sine or cosine, but falls off like $1/\sqrt{z}$ as $z \to \infty$.

The most general isotropic covariance is given in (2), with the absolutely continuous measure $g(r) \frac{2\pi^{n/2}}{\Gamma(n/2)} r^{n-1} dr$ replaced by an arbitrary positive finite measure $\gamma(dr)$ on $[0, \infty)$. Any isotropic covariance function may be approximated by one with a discrete spectral measure $\gamma(dr) = \sum \gamma_j \delta_{r_j}(dr)$ assigning mass γ_j to finitely many points r_j :

$$C(\rho) \approx \sum_{j} (2/r_{j}\rho)^{\nu} \Gamma(\nu+1) J_{\nu}(r_{j}\rho) \gamma_{j}$$

$$= \begin{cases} \sum_{j} \gamma_{j} \cos(r_{j}\rho) & \text{if } n = 1\\ \sum_{j} \gamma_{j} J_{0}(r_{j}\rho) & \text{if } n = 2\\ \sum_{j} \gamma_{j} \sqrt{\pi/2r_{j}\rho} J_{1/2}(r_{j}\rho) & \text{if } n = 3 \end{cases}$$

$$(3)$$

but a more common approach is to choose small parametric families of densities $g^{\theta}(r)$ or measures $g^{\theta}(dr)$.

We can recover the spectral density $g(r) = G(\omega)$ (for $r = |\omega|$) through the Fourier inversion formula, using polar coordinates with $\rho = |h| \in \mathbb{R}_+$ and $\sigma = h/|h| \in S^{n-1}$:

$$g(r) = G(\omega) = \frac{1}{(2\pi)^n} \int \cos(-h \cdot \omega) C_0(h) d^n h$$

$$= \frac{1}{(2\pi)^n} \iint_{\mathbb{R}_+ \times S^{n-1}} \cos(-r\rho \sigma_\omega) C(\rho) \frac{2\pi^{n/2} \rho^{n-1}}{\Gamma(n/2)} d\rho d\sigma$$

$$= \int_0^\infty r(\rho/2\pi r)^{n/2} J_\nu(r\rho) C(\rho) d\rho, \qquad \nu \equiv \frac{n}{2} - 1 \qquad (4)$$

$$= \begin{cases} \int_0^\infty \frac{2}{\pi} \cos(r\rho) C(\rho) d\rho & \text{if } n = 1 \\ \int_0^\infty (\rho/2\pi) J_0(r\rho) C(\rho) d\rho & \text{if } n = 2 \\ \int_0^\infty r(\rho/2\pi r)^{3/2} J_{1/2}(r\rho) C(\rho) d\rho & \text{if } n = 3 \end{cases}$$

It is hard to imagine what $C_0(h)$ would look like for different choices of g(r); a simple approach is to take whatever symmetric functions G(u) whose Fourier transforms we can find, and see what we get. Here are some commonly used covariance families, in n = 2 dimensions; in each case $\theta_1 = C(0)$ is an overall level parameter and θ_2 is a distance scale parameter:

• Power family

$$C(\rho|\theta, p) = \theta_1 \exp\{-|\rho/\theta_2|^p\}, \ 0$$



Notice that the exponential has a negative derivative at z = 0, so it falls off quickly at first, then slowly levels off, while the Gaussian has zero derivative near z = 0 then falls off very quickly. From (5) it follows that the exponential has spectral density function g(r) = $(\theta_1 \theta_2^2 / 2\pi) / (1 + r^2 \theta_2^2)^{3/2}$, proportional to a bivariate Cauchy density function, while the Gaussian has spectral density $g(r) = (\theta_1 \theta_2^2 / 4\pi) \exp(-r^2 \theta_2^2 / 4)$, proportional to a normal density.

• Matérn

$$C(\rho|\theta) = \frac{2\,\theta_1}{\Gamma(\theta_3)} \left(\frac{\rho}{2\theta_2}\right)^{\theta_3} K_{\theta_3}(\rho/\theta_2)$$



The displayed plot has shape parameter $\theta_3 = 2$. The Matérn class is quite flexible and includes the exponential family (with $\theta_3 = \frac{1}{2}$), the Gaussian family (in the limit as $\theta_3 \to \infty$), and many others. In *n* dimensions its spectral density function is

$$g(r) = \frac{\theta_1 \theta_2^n}{\Gamma(\theta_3) \pi^{n/2}} (1 + \theta_2^2 r^2)^{-\theta_3 - n/2},$$

proportional to the familiar *n*-variate Student's *t* density function with $2\theta_3$ degrees of freedom and variance scale $\sigma^2 = 1/2\theta_2^2\theta_3$. This lends more insight into how the Matérn reduces to the exponential when $\theta_3 = 1/2$ and to the Gaussian when $\theta_3 \to \infty$.

• Spherical

$$C(\rho|\theta) = \begin{cases} \theta_1 \left[1 - \frac{2}{\pi} \left(\frac{\rho}{\theta_2} \sqrt{1 - (\frac{\rho}{\theta_2})^2} + \sin^{-1} \frac{\rho}{\theta_2} \right) \right] & \text{for } \rho < \theta_2 \\ 0 & \text{for } \rho \ge \theta_2 \end{cases}$$

The spherical covariance function is proportional to the area of intersection for two discs of diameter θ_2 with centers separated by distance ρ . In this model the Gaussian quantities Z_j and Z_k at loci s_j and s_k



This is not quite linear. Like the exponential, it has a negative slope at z = 0 and falls off rapidly at first; like the Gaussian, it falls off rapidly later and in fact reaches zero. The spectral density, while available in closed form, isn't illuminating; it's best to think of the spherical process as a convolution or moving average of Gaussian white noise, integrated at each locus over the surrounding ball of diameter θ_2 .

A variety of processes may be constructed similarly as kernel integrals of standard Gaussian white noise,

$$Z(h) = \int_{\mathbb{R}^n} k(h-s)\,\zeta(ds);$$

where "standard" means that $\mathsf{E}[\zeta(ds)] = 0$ and $\mathsf{E}[\zeta(ds)^2] = ds$. The covariance is

$$C_0(h) = \mathsf{E}[Z(0)\overline{Z(h)}] = \int_{\mathbb{R}^n} k(h-s) \,\overline{k(-s)} \, ds$$

with spectral density

$$G(\omega) = (2\pi)^{-n} \int e^{-i\omega \cdot h} C_0(h) \, dh$$

= $(2\pi)^{-n} \iint e^{-i\omega \cdot h} k(h-s) \, \overline{k(-s)} \, ds \, dh$
= $(2\pi)^{-n} \left| \int e^{-i\omega \cdot x} k(x) \, dx \right|^2$

so the kernel may be computed from the spectral density as

$$k(x) = (2\pi)^{-n/2} \int e^{i\omega \cdot x} G(\omega)^{1/2} d\omega$$

or, in polar coordinates,

$$k(\rho) = \int_0^\infty r^{\nu+1} \rho^{-\nu} J_{\nu}(r\rho) g(r)^{1/2} dr$$

=
$$\begin{cases} \int_0^\infty \sqrt{\frac{2}{\pi}} \cos(r\rho) \sqrt{g(r)} dr & \text{if } n = 1 \\ \int_0^\infty J_0(r\rho) r \sqrt{g(r)} dr & \text{if } n = 2 \\ \int_0^\infty J_{1/2}(r\rho) r^{3/2} \rho^{-1/2} \sqrt{g(r)} d\rho & \text{if } n = 3 \end{cases}$$

provided that the square root of the spectral density is the Fourier transform of a finite positive function, *i.e.*, is itself positive semidefinite. For the Matérn class, the root spectral density $\sqrt{g(r)} \propto (1 + \theta_2^2 r^2)^{-(\theta_3 + n/2)/2}$ will be another *n*-variate *t* density provided $\theta_3 > n/2$ and in this case, setting $\epsilon = (2\theta_3 - n)/4 > 0$, we find

$$k(\rho) = \frac{2\theta_1^{1/2} (2\rho\theta_2)^{-\epsilon - n/2}}{\Gamma(\epsilon + n/2)\sqrt{\Gamma(2\epsilon + n/2)}\pi^{n/4}} K_{\epsilon}(\rho/\theta_2)$$

leads to a moving-average kernel representation for the Matérn covariance class. In any number $n \geq 1$ of dimensions the restriction $\epsilon > 0$ entails $\theta_3 > n/2 \geq 1/2$, ruling out the exponential covariance, but the Gaussian covariance (the limiting case as $\theta_3 \to \infty$) is available in any number of dimensions, with

$$k(\rho) = \theta_1^{1/2} (\pi \theta_2^2/4)^{-n/2} e^{-2\rho^2/\theta_2^2}$$

References

- Abramowitz, M. and Stegun, I. A., eds. (1964), Handbook of Mathematical Functions With Formulas, Graphs, and Mathematical Tables, volume 55 of Applied Mathematics Series, Washington, D.C.: National Bureau of Standards.
- Cressie, N. A. C. (1993), *Statistics for Spatial Data*, New York, NY, USA: John Wiley & Sons.
- Le, N. D. and Zidek, J. V. (1992), "Interpolation with uncertain spatial covariances: A Bayesian alternative to kriging," *Journal of Multivariate Analysis*, 43, 351–374.
- Mardia, K. and Marshall, R. J. (1984), "Maximum likelihood estimation of models for residual covariance in spatial regression," *Biometrika*, 71, 135–146.
- Matérn, B. (1960), Spatial Variation, volume 49 of Meddelanden fran Statens Skogsforsningsinstitut, Stockholm: Statens Skogsforsningsinstitut, first edition, (second edition published by Springer-Verlag in 1986).
- Ripley, B. D. (1981), *Spatial Statistics*, New York, NY, USA: John Wiley & Sons.
- Yaglom, A. M. (1962), An Introduction to the Theory of Stationary Random Functions, New York: Dover, translated and Edited by Richard A. Silverman from 1952 Russian article in Uspekhi Matematicheskikh Nauk.