

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 $\mu(s) = \mathbf{E}Z(s)$ and the covariances $C(s, t) = \mathbf{E}(Z(s) - \mu(s))(Z(t) - \mu(t))$.

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 $C(s - t) = \mathbf{E}(Z(s) - \mu(s))(Z(t) - \mu(t)) = C(t - s)$. But more is true: if $\{s_j\}$ any collection of locations, then complex linear combinations $a^\top(Z - \mu) = \sum a_j(Z_j - \mu_j)$ of the centered random variables $Z_j = Z(s_j)$ (with means $\mu_j = \mu(s_j)$) must have nonnegative squared modulus $\mathbf{E}|\sum a_j(Z_j - \mu_j)|^2 = \sum a_j C(s_j - s_k) \bar{a}_k \geq 0$ 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 $\sum_{jk} a_j C(s_j - s_k) \bar{a}_k \geq 0$ 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 $\sum a_j C(s_j - s_k) a_k \geq 0$ 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

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

It turns out that this is the *only* way to get one— that every positive semi-definite 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

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

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 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 $\text{Di}(\frac{1}{2}, \dots, \frac{1}{2})$ 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 $r = |\omega| \geq 0$ 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 $\text{Be}(\frac{1}{2}, \frac{n-1}{2})$ random variable, writing ρ for $|h|$,

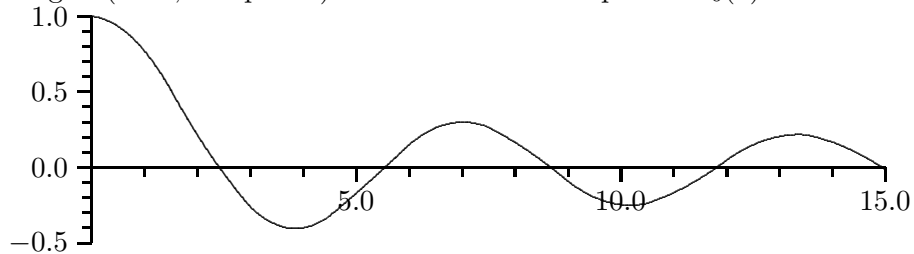
$$\begin{aligned}
 C_0(h) &= \int_{\mathbb{R}^n} \cos(h \cdot \omega) g(|\omega|) d^n \omega \\
 &= \iint_{\mathbb{R}_+ \times S^{n-1}} \cos(r\rho\sigma_h) g(r) \frac{2\pi^{n/2} r^{n-1}}{\Gamma(n/2)} dr d\sigma \\
 &= \int_{\mathbb{R}_+} \int_0^1 \cos(r\rho\sqrt{u}) g(r) \frac{2\pi^{n/2} r^{n-1}}{\Gamma(n/2)} \frac{\Gamma(n/2)}{\Gamma(\frac{1}{2}) \Gamma(\frac{n-1}{2})} u^{1/2-1} (1-u)^{(n-1)/2-1} dr du \\
 &= \int_0^\infty \rho (2\pi r/\rho)^{\nu+1} J_\nu(r\rho) g(r) dr, \quad \nu \equiv \frac{n}{2} - 1 \tag{1}
 \end{aligned}$$

$$\begin{aligned}
 &= \int_0^\infty (r\rho/2)^{-\nu} \Gamma(\nu+1) J_\nu(r\rho) \gamma(dr) \tag{2} \\
 &= \begin{cases} \int_0^\infty 2 \cos(r\rho) g(r) dr & \text{if } n = 1 \\ \int_0^\infty 2\pi r J_0(r\rho) g(r) dr & \text{if } n = 2 \\ \int_0^\infty \rho (2\pi r/\rho)^{3/2} J_{1/2}(r\rho) g(r) dr & \text{if } n = 3 \end{cases}
 \end{aligned}$$

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 $z \rightarrow \infty$.

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 :

$$\begin{aligned} C(\rho) &\approx \sum_j (2/r_j \rho)^\nu \Gamma(\nu + 1) J_\nu(r_j \rho) \gamma_j & (3) \\ &= \begin{cases} \sum_j \gamma_j \cos(r_j \rho) & \text{if } n = 1 \\ \sum_j \gamma_j J_0(r_j \rho) & \text{if } n = 2 \\ \sum_j \gamma_j \sqrt{\pi/2r_j \rho} J_{1/2}(r_j \rho) & \text{if } n = 3 \end{cases} \end{aligned}$$

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 $\sigma = h/|h| \in S^{n-1}$:

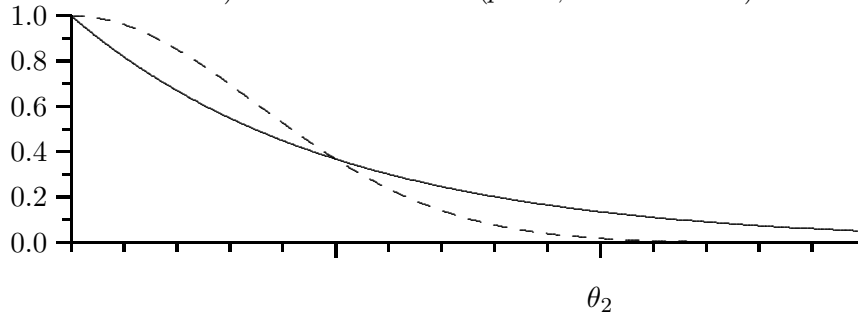
$$\begin{aligned} g(r) = G(\omega) &= \frac{1}{(2\pi)^n} \int \cos(-h \cdot \omega) C_0(h) d^n h \\ &= \frac{1}{(2\pi)^n} \iint_{\mathbb{R}_+ \times S^{n-1}} \cos(-r\rho\sigma_\omega) C(\rho) \frac{2\pi^{n/2} \rho^{n-1}}{\Gamma(n/2)} d\rho d\sigma \\ &= \int_0^\infty r(\rho/2\pi r)^{n/2} J_\nu(r\rho) C(\rho) d\rho, \quad \nu \equiv \frac{n}{2} - 1 & (4) \\ &= \begin{cases} \int_0^\infty \frac{2}{\pi} \cos(r\rho) C(\rho) d\rho & \text{if } n = 1 \\ \int_0^\infty (\rho/2\pi) J_0(r\rho) C(\rho) d\rho & \text{if } n = 2 \\ \int_0^\infty r(\rho/2\pi r)^{3/2} J_{1/2}(r\rho) C(\rho) d\rho & \text{if } n = 3 \end{cases} & (5) \end{aligned}$$

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

$$C(\rho|\theta, p) = \theta_1 \exp\{-|\rho/\theta_2|^p\}, \quad 0 < p \leq 2$$

Two notable covariograms in this family are the *exponential* ($p = 1$, solid below) and the *Gaussian* ($p = 2$, dashed below):

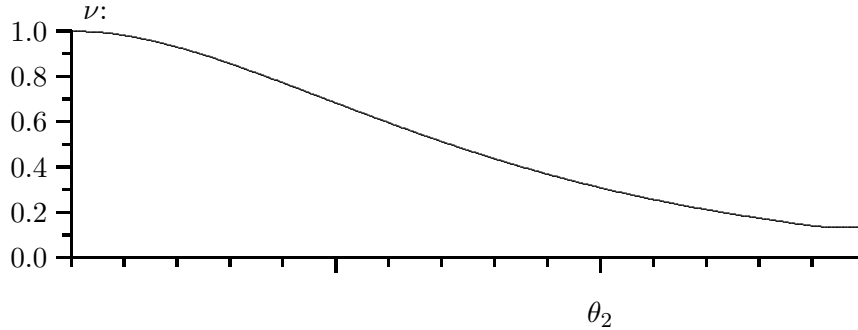


Notice that the exponential has a negative derivative at $z = 0$, so it falls off quickly at first, then slowly levels off, while the Gaussian has zero derivative near $z = 0$ then falls off very quickly. From (5) it follows that the exponential has spectral density function $g(r) = (\theta_1\theta_2^2/2\pi)/(1+r^2\theta_2^2)^{3/2}$,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

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

where $K_\nu(z)$ is the modified Bessel function of the third kind of order ν :



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 $\theta_3 \rightarrow \infty$), and many others. In n dimensions its spectral density function is

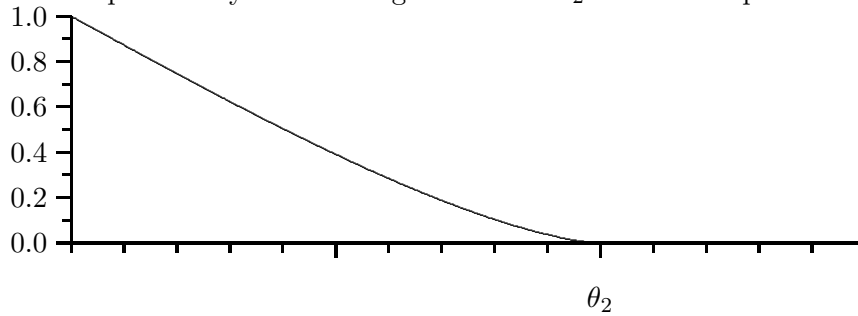
$$g(r) = \frac{\theta_1\theta_2^n}{\Gamma(\theta_3)\pi^{n/2}}(1+\theta_2^2r^2)^{-\theta_3-n/2},$$

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 $\theta_3 \rightarrow \infty$.

- Spherical

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

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 separated by a distance greater than θ_2 will be independent.



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 $E[\zeta(ds)] = 0$ and $E[\zeta(ds)^2] = ds$. The covariance is

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

with spectral density

$$\begin{aligned}
G(\omega) &= (2\pi)^{-n} \int e^{-i\omega \cdot h} C_0(h) dh \\
&= (2\pi)^{-n} \iint e^{-i\omega \cdot h} k(h-s) \overline{k(-s)} ds dh \\
&= (2\pi)^{-n} \left| \int e^{-i\omega \cdot x} k(x) dx \right|^2
\end{aligned}$$

so that an isotropic kernel may be computed from the spectral density as

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

or, in polar coordinates,

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

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 $\theta_3 \rightarrow \infty$) is available in any number of dimensions, with

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

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