Other Covariance Functions and Anisotropy

Covariance Functions

Given the assumption that a Gaussian process is reasonable for the spatial process, a valid covariance function needs to be specified.

Up to this point, we have largely worked with isotropic covariance functions. In particular, the exponential covariance functions has primarily been used. However, a Gaussian process is flexible and can use any valid covariance function.

A valid covariance function $C(\mathbf{h})$, defined as $Cov(Y(\mathbf{s}), Y(\mathbf{s} + \mathbf{h}))$, for any finite set of sites $\mathbf{s_1}, \ldots, \mathbf{s_n}$ and a_1, \ldots, a_n should satisfy

$$Var\left[\sum_{i} a_{i} Y(\boldsymbol{s_{i}})\right] = \sum_{i,j} a_{i} a_{j} Cov(Y(\boldsymbol{s_{i}}), Y(\boldsymbol{s_{j}})) = \sum_{i,j} a_{i} a_{j} C(\boldsymbol{s_{i}} - \boldsymbol{s_{j}}) \ge 0$$

with strict inequality if all the a_i are not zero.

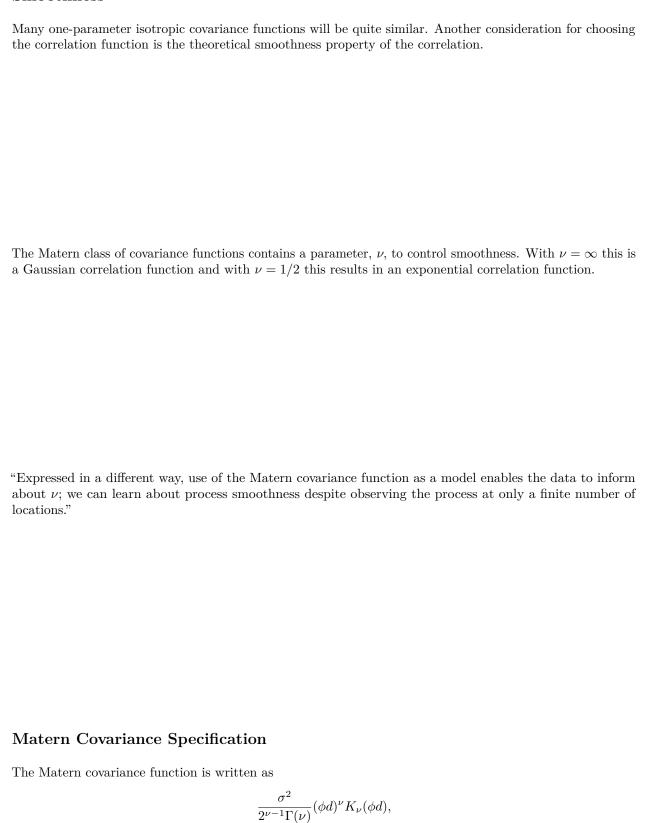
In other words, C(h) needs to be a positive definite function, which includes the following properties 1. $C(\mathbf{0}) \ge 0$ 2. $|C(h)| \le C(\mathbf{0})$

Constructing covariance functions

There are three approaches for building correlation functions. For all cases let C_1, \ldots, C_m be valid correlation functions:

- 1. Mixing: $C(\mathbf{h}) = \sum_{i} p_i C_i$ is also valid if $\sum_{i} p_i = 1$.
- 2. Products: $C(\mathbf{h}) = \prod_i C_i$
- 3. Convolution: $C_{12}(\mathbf{h}) = \int C_1(\mathbf{h} \mathbf{t})C_2(\mathbf{t})d\mathbf{t}$ this is based on a Fourier transform.

Smoothness



where $\Gamma()$ is a gamma function and K_{ν} is the modified Bessel function of order ν .

Anisotropy

Anisotropy means that the covariance function is not just a function of the distance ||h||, but also the direction.

Geometric anisotropy refers to the case where the coordinate space is anisotropic, but can be transformed to an isotropic space.

If the differences in spatial structure are directly related to two coordinate sets (lat and long), we can create a stationary, anistropic covariance function

Let

$$cor(Y(s+h), Y(s)) = \rho_1(h_u)\rho_2(h_x),$$

where $\rho_1()$ and $\rho_2()$ are proper correlation functions.

In general consider the correlation function,

$$\rho(\boldsymbol{h};\phi) = \phi_0(||L\boldsymbol{h}||;\phi)$$

where L is a $d \times d$ matrix that controls the transformation.

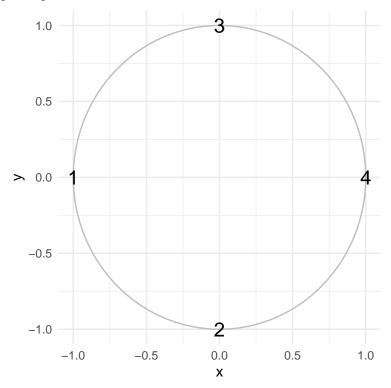
Let
$$Y(s) = \mu(s) + w(s) + \epsilon(s)$$
, and $Y(s) \sim N(\mu(s), \Sigma(\tau^2, \sigma^2, \phi, B))$, where $B = L^T L$.

The covariance matrix is defined as $\Sigma(\tau^2, \sigma^2, \phi, B) = \tau^2 I + \sigma^2 H((\boldsymbol{h}^T B \boldsymbol{h}^T)^{\frac{1}{2}})$, where $H((\boldsymbol{h}^T B \boldsymbol{h}^T)^{\frac{1}{2}})$ has entries of $\rho((\boldsymbol{h_{ij}}^T B \boldsymbol{h_{ij}}^T)^{\frac{1}{2}}))$ with $\rho()$ being a valid covariance function, typically including ϕ and $\boldsymbol{h_{ij}} = \boldsymbol{s_i} - \boldsymbol{s_j}$.

B is often referred to as a transformation matrix which rotates and scales the coordinates, such that the resulting transformation can be simplified to a distance.

Geometric Anisotropy Visual

• Consider four points positioned on a unit circle.



Now consider a set of correlation functions. For each, calculate the correlation matrix and discuss the impact of B on the correlation. Furthermore, how does B change the geometry of the correlation between points 1, 2, 3, and 4?

1.
$$\rho() = \exp(-\boldsymbol{h_{ij}}^T B \boldsymbol{h_{ij}}^T)^{\frac{1}{2}}))$$
, where $B = \begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix}$

2.
$$\rho() = \exp(-\boldsymbol{h_{ij}}^T B \boldsymbol{h_{ij}}^T)^{\frac{1}{2}}))$$
, where $B = \begin{pmatrix} 2 & 0 \\ 0 & 1 \end{pmatrix}$

3.
$$\rho() = \exp(-\boldsymbol{h_{ij}}^T B \boldsymbol{h_{ij}}^T)^{\frac{1}{2}}))$$
, where $B = \begin{pmatrix} 3 & 1 \\ 1 & 1 \end{pmatrix}$

1. $\rho() = \exp(-\boldsymbol{h_{ij}}^T I \boldsymbol{h_{ij}}^T)^{\frac{1}{2}})$

1.000	0.243	0.243	0.135
0.243	1.000	0.135	0.243
0.243	0.135	1.000	0.243
0.135	0.243	0.243	1.000

2. $\rho() = \exp(-\mathbf{h_{ij}}^T B \mathbf{h_{ij}}^T)^{\frac{1}{2}})$, where $B = \begin{pmatrix} 2 & 0 \\ 0 & 1 \end{pmatrix}$

1.000	0.177	0.177	0.059
0.177	1.000 0.135	0.135 1.000	0.177 0.177
0.059	0.177	0.177	1.0

3. $\rho() = \exp(-\mathbf{h_{ij}}^T B \mathbf{h_{ij}}^T)^{\frac{1}{2}})$, where $B = \begin{pmatrix} 3 & 1 \\ 1 & 1 \end{pmatrix}$

0.243	0.086	0.031
1.000	0.135	0.086
0.135	1.000	0.243
0.086	0.243	1.000
	1.000 0.135	1.000 0.135 0.135 1.000

The (effective) range for any angle η is determined by the equation

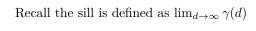
$$\rho(r_{\eta}(\tilde{\boldsymbol{h}}_{\eta}^{T}B\tilde{\boldsymbol{h}}_{\eta}^{T})^{\frac{1}{2}}) = .05,$$

where \tilde{h}_{η} is a unit vector in the direction η .

Okay, so if we suspect that geometric anisotrophy is present, how do we fit the model? That is, what is necessary in estimating this model?

- In addition to σ^2 and τ^2 we need to fit B.
- While B is a matrix, it is just another unknown parameter.
- To fit a Bayesian model we need a prior distribution for B. One option for the positive definite matrix is the Wishart distribution, which is a bit like a matrix-variate gamma distribution.

Sill, Nugget, and Range Anisotropy



Let \boldsymbol{h} be an arbitrary separation vector, that can be normalized as $\frac{\boldsymbol{h}}{||\boldsymbol{h}||}$

If $\lim_{a\to\infty} \gamma(a \times \frac{h}{||h||})$ depends on h, this is referred to as sill anisotropy.

Similarly the nugget and range can depend on h and give nugget anisotropy and range anisotropy