Spatial Autocorrelation I: Variograms

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DATA 589: Spatial Statistics



- 1. Overview
- 2. Impacts of autocorrelation
- 3. Detecting Spatial Autocorrelation
- 4. Applied Variogram Analysis

Overview



So far we have covered situations where locations were the variables of interest, and we studied these data using point processes

...but there are many situations where locations are arbitrary:



(Bell *et al.*, 2010) DATA 589: Spatial Statistics

Rainfall in Europe



(Ballabio et al., 2017)



(Feeney et al., 2022)

Overview





These data can't be treated as a point process because the locations of the points are (mostly) meaningless

...but they also can't be modelled using off the shelf techniques because they break an important assumption; that the data are IID.



Data are often collected by measuring quantities over space (e.g., abundances, concentrations, etc.).

When this is the case, spatial autocorrelation can arise when the variation between the values of the datapoints is affected by their spatial distance (i.e., data that are close together in space more similar than data collected further apart).

The underlying reason for this is that many of the drivers (e.g., environmental conditions, topography, geology, etc.) act at large spatial scales.

The result is that we can not assume that spatially collected data are IID. DATA 589: Spatial Statistics



What if we're studying the effect of rainfall on species diversity in the Amazon?



Source: Tadashi Fukami and Jes Coyle

Because rainfall is correlated in space, species diversity will also be correlated in space (if the relationship exists).

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Impacts of autocorrelation

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Anything that causes some data points to be more similar to each other than others can result in autocorrelation.

• **Time**: Data that are close together in time are more related.

• **Space**: Data that are close together in space are more related.



We will be focusing on spatial autocorrelation, but the ideas translate to other sources of autocorrelation.

Autocorrelation is important





C Chris Sorensen

Dr. Sam Wang, Neuroscientist —Princeton Election Consortium

"It is totally over. If Trump wins I will eat a bug." DATA 589: Spatial Statistics



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Nate Silver, Statistician —FiveThirtyEight.com

"Trump Is Just A Normal Polling Error Behind Clinton."

Autocorrelation is important





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Ignoring non-ind. \rightarrow overconfidence



Polls underestimated Trump in red states, Clinton in blue states

2016 election results vs. FiveThirtyEight's adjusted polling average by state



Are these polling errors independently distributed?

This same statistical issue that caused overly confident predictions of Clinton's 2016 victory can result in overconfidence in parameter estimates and predictions in regression models.

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SOURCE: DAVID WASSERMAN

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Sample size, n is the denominator when calculating SEs and CIs.

$$SE = \frac{\sigma}{\sqrt{n}} \qquad \qquad 95\%CI = \bar{x} \pm 1.96\frac{\sigma}{\sqrt{n}}$$

All else equal: $\uparrow n = \downarrow SE \& \downarrow CI$

But with autocorrelated data each new datapoint is related to a previously collected datapoint and does not bring a full independent datapoint worth of information (e.g., 90% autocorr. $\approx 10\%$ new info).

When data are autocorrelated $n_{\text{effective}} < n$, meaning SEs and CIs shrink faster than they should, resulting in a false sense of confidence.

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Spatially referenced measurements contain valuable information (e.g., soil organic carbon on a mountain in Mexico)



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(Fusaro et al., 2019)
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... but what if we want to know how much carbon is at a nearby, un-sampled location?

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Leveraging the information contained the spatial autocorrelation between samples allows us to make predictions to unsampled areas



(Fusaro et al., 2019)



Data that are collected across space will likely be autocorrelated, breaking the IID assumption.

Ignoring this autocorrelation results in biased, over-confident models, and (more importantly) does not leverage the full amount of information contained in the data.

During the rest of this course we will focus on ways to visualise, detect, and work with the autocorrelation contained in spatial data.

Detecting Spatial Autocorrelation



Scientists in many fields of research find themselves collecting data repeatedly over space, but the tools originally come from the field of geostatistics.

Can you think of data that are collected over space that might drive statistical innovation? Data where the ability to predict where things occur might be profitable?



The tools for working with spatial autocorrelation were developed for the goal of mapping mineral deposits.



We're going to work with a dataset on forest composition in Tatarstan, Russia from Zuur *et al.* (2007).

The variable of interest is a measure of boreality (\sim percent boreal species at a site).





Spatial autocorrelation can be difficult to see in a simple \times vs. y scatterplot (not designed for this purpose).





'Bubble plots' are an easy tool to quickly assess for autocorrelation.



Values are plotted in space, and sizes/colours are proportional to their values. The idea is to look for patterns.

Bubble plots are quick and easy to generate, but can be hard to read and are not particularly formal.

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Moran's I



Moran's I is a correlation coefficient that measures the overall spatial autocorrelation of a data set (think of it as \sim weighted covariance):

$$I = \frac{N}{W} \frac{\sum_{i} \sum_{j} w_{ij}(x_i - \bar{x})(x_j - \bar{x})}{\sum_{i} (x_i - \bar{x})^2}$$

N is the number of spatial units indexed by i and j; x is the variable of interest and \bar{x} is the mean of x; w_{ij} is a matrix of spatial weights and W is the sum of all w_{ij} .

Values of I usually range from -1 to +1.



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Many R packages for calculating Moran's I

```
library(ape)
library(fields)
#Vector of spatial coordinates
coords = cbind(data$x, data$y)
#Matrix of distances for the weights
w = fields::rdist(coords)
#Calculate Moran's I
ape::Moran.I(data$Bor, w = w)
$observed
[1] -0.03019649
$expected
[1] -0.001879699
$sd
[1] 0.001368412
$p.value
[1] 3.991055e-95
```



The *p*-value tells us we have significant spatial autocorrelation.



Moran's I can be a useful tool for identifying the presence of autocorrelation and is quite popular.

The challenge is how to act on this information (i.e., lets you know if you have autocorrelation, but doesn't help in modelling it)?



Moran's I is also very sensitive to how you define the weights:

"The idea is to construct a matrix that accurately reflects your assumptions about the particular spatial phenomenon in question. A common approach is to give a weight of 1 if two zones are neighbors, and 0 otherwise, though the definition of 'neighbors' can vary. Another common approach might be to give a weight of 1 to k nearest neighbors, 0 otherwise. An alternative is to use a distance decay function for assigning weights. Sometimes the length of a shared edge is used for assigning different weights to neighbors. The selection of spatial weights matrix should be guided by theory about the phenomenon in question." — Wikipedia



Semi-variograms are functions describing the degree of spatial dependence in a spatial stochastic process, $Z(\mathbf{s})$.

Semi-variance, $\gamma(h)$, is a measure of the degree of similarity between pairs of points separated by distance h, given by

$$\gamma(h) = \frac{1}{2V} \iint_V \left[f(M+h) - f(M) \right]^2 dV,$$

where M is a point in the spatial field V, f(M) is the value at point M (in arbitrary units); h is the separation distance (in e.g., meters or km); and the double integral is over 2 dimensions.

Semi-variograms properties



- $\gamma(h) \ge 0$, since it is the expectation of a square.
- Since $Z(\mathbf{s}_1) Z(\mathbf{s}_1) = 0$, $\gamma(0)$ is always 0.
- If the process is stationary, in the limit where $h \to \infty$ $\gamma_s(h) = var(Z(\mathbf{s})).$
- If a stationary process has no spatial dependence, the semi-variogram is constant everywhere except at the origin, where it is zero.
- The semi-variogram might be discontinuous at the origin (the height of the jump at the origin is referred to as nugget).

For a discussion of these properties see Bachmaier & Backes (2011).



To obtain the semi-variogram for a given $\gamma(h)$, all pairs of points at that exact distance, h, would need to be sampled. In practice, this is impossible, so the empirical semi-variogram is used instead.

For values of $Z(\mathbf{s})$ for all pairs separated by distance h, $\gamma(h)$ is estimated as:

$$\hat{\gamma}(h\pm\delta) := \frac{1}{2|N(h\pm\delta)|} \sum_{(i,j)\in N(h\pm\delta)} |z_i - z_j|^2$$

where δ is some bin width.



$$\hat{\gamma}(h\pm\delta) := rac{1}{2|N(h\pm\delta)|} \sum_{(i,j)\in N(h\pm\delta)} |z_i - z_j|^2$$



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```
# Import the data
data <- read.csv("Datasets/Boreality.csv")</pre>
# Spatial data frame of boreality
DATA <- data.frame(Z_s = data$Bor,
                    x = data$x,
                    v = data$v)
# Define coordinates
sp::coordinates(DATA) <- c("x","y")</pre>
# object of class = "SpatialPointsDataFrame"
# Calculate empirical variogram
vg <- gstat::variogram(Z_s ~ 1, data = DATA)
# object of class = "gstatVariogram"
plot(vg)
```



Autocorr. in the boreality data in R



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Applied Variogram Analysis



Annoni *et al.* (2019) used a semi-variograms and Moaran's I to examine regional economic resilience in the EU, and to what extent GDP is influenced by neighbours.





Weighting GDP by Moran's I showed how regions influenced each other.



Core a periphery GDP zones



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Autocorrelation contains a lot of information about a spatial process.

There are many methods for visualising and working with spatially correlated data (e.g., bubble plots, Moran's I).

Semi-variograms are useful tools for visualising spatial autocorrelation, are objective, and have a long, robust history.

Usefully, the shape of a dataset's empirical variogram can also provide clues on how to best model the autocorrelation in the data, which we will cover next lecture.

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