

Lecture 7: Spatial Analysis II

GEOG413/613
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Spatial Analysis

- Point Pattern Analysis
 - Methods
 - Nearest Neighbor Analysis
 - Quadrat Analysis
 - Spatial Autocorrelation

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Descriptive Non-Spatial Statistics

- Mean, Standard Deviation, Skewness, Kurtosis

Descriptive Spatial Statistics

- Mean Centre, Weighted Mean Centre, Standard Distance, Relative Distance

Inferential Spatial Statistics and Analysis

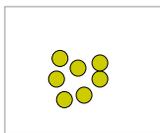
- Point Pattern Analysis (Nearest Neighbor Analysis, Quadrat Analysis)
- Cluster Analysis
- Spatial Autocorrelation
- Factor Analysis
- Spatial Interpolation

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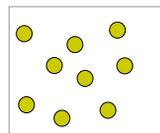
Examining Spatial Data

- ▶ How are the data points spatially distributed?

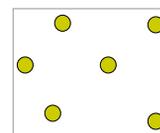
Clustered



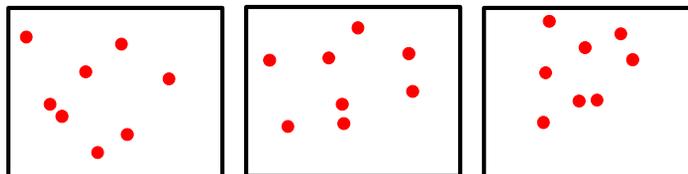
Random



Dispersed



- ▶ How do you know? Always test



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Point Pattern Analysis

- A set of quantitative tools for examining the spatial arrangement of point locations on the landscape as represented by a conventional map.
- Two methods are ***nearest neighbour analysis*** and ***quadrat analysis***.

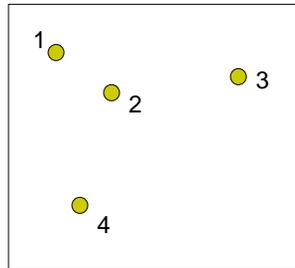
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Nearest Neighbour Analysis

- Distance of each point to its nearest neighbour is measured
- The average nearest distance for all points is then calculated
- Can compare results with expected average for a random distribution

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Nearest Neighbor Analysis



$$\begin{aligned}
 d_1 &= l_{12} \\
 d_2 &= l_{21} \\
 d_3 &= l_{32} \\
 d_4 &= l_{42} \\
 r_{obs} &= \frac{\sum_1^n d_i}{n}
 \end{aligned}$$

The Average Nearest Neighbor distance = r_{obs}

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Nearest Neighbor Analysis

- The average nearest neighbour distance is an absolute value
- It is a function of the units in which the distance is measured
- Problem
 - How can we compare data from different regions or studies?
 - Solution: **Standardized Nearest Neighbour Index**

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Nearest Neighbour Analysis

- The utility of the average nearest neighbour distance comes from comparing the index value for an observed pattern to the results produced from certain distinct point distributions
 - We can compare our results against values for random, clustered and dispersed distributions
 - Problem
 - How can we compare data from different regions or studies?
 - Solution: **Standardized Nearest Neighbour Index**

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Random Distribution

- For a random distribution, the average nearest neighbour distance is calculated as follows:

$$r_{rnd} = \frac{1}{2\sqrt{n/A}}$$

Where :

A is area of study region

n is number of points

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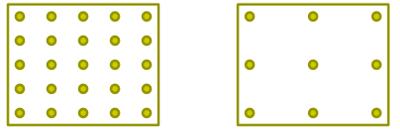
Maximum Dispersion Distribution

- If the distribution is perfectly uniform, the average nearest neighbour distance is calculated as follows:

$$r_{dsp} = \frac{1.07453}{\sqrt{n/A}}$$

Question:

Consider the two distributions below, assume that the area is the same. Do they have different r_{dsp} values? If so, which one will have a higher r_{dsp} value?



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Clustered Distribution

- When all points lie in the same position (i.e. maximum clustering) the average nearest neighbour distance is 0

$$r_{cst} = 0$$

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Standardized Nearest Neighbor Index

- The Standardized Nearest Neighbor Index is computed as a ratio of r_{obs} to r_{rnd} , the expected average nearest neighbor distance for a random distribution

$$R = \frac{r_{obs}}{r_{rnd}}$$

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Standardized Nearest Neighbour Index

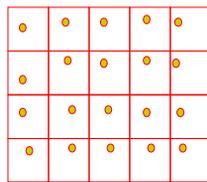
- To calculate the standardized nearest neighbour index for maximum dispersion:

$$R = \frac{r_{dsp}}{r_{rnd}}$$

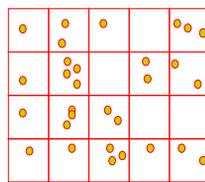
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Quadrat Analysis

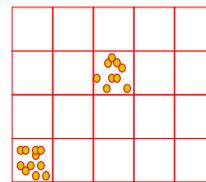
- Examines the frequency of points occurring in various parts of the study area.
- The point pattern arrangement in the study area is described with the aid of the frequency of points in a cell



Dispersed



Random



Clustered

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Quadrat Analysis

- In quadrat analysis, an index known as the variance-mean ratio (VMR) standardizes the degree of variability in cell frequencies relative to the mean of the cell frequency

$$VMR = \frac{Var}{Mean}$$

where n = number of points

m = number of cells

$Mean$ = mean cell frequency

Var = Variance of cell frequencies

$$Mean = \frac{n}{m}$$

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Quadrat Analysis

- Variance of Cell Frequencies

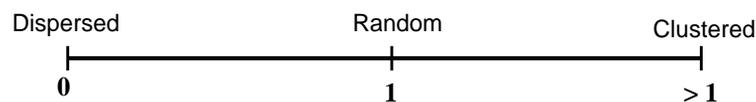
$$Var = \frac{\sum f_i x_i^2 - \frac{(\sum f_i x_i)^2}{m}}{m-1}$$

where f_i = frequency of cells with i cases
 x_i = number of cases per cell

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Quadrat Analysis

- Variance-Mean Ratio (VMR)
 - If each cell contains the same amount of points, then VMR = 0
 - If a point pattern is highly clustered with most cells containing no points, then VMR will be relatively large.
 - If the point pattern is perfectly random, then the mean cell frequency equals the variance of the cell frequency, and VMR = 1



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Quadrat Analysis

- Test of Significance
 - Applied to determine if distribution of points is random.
 - The test statistic used is chi-square:

$$X^2 = \text{VMR} (m - 1)$$

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Cluster Analysis

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Cluster Analysis

- To reduce data complexity by sorting the data into subsets (clusters) that share some common trait
- Achieve the reduction of observations by minimizing the within-group variation and maximizing the between group variation (i.e. the degree of association between two objects is maximal if they belong to the same group and minimal otherwise)

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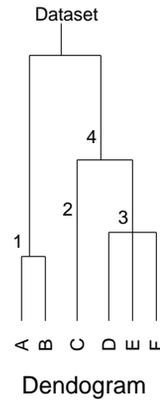
Clustering Analysis

- Searching for groups in the data in such a way that objects belonging to the same cluster resemble each other, whereas objects in different clusters are dissimilar.
- Two methods are partitioning and hierarchical clustering

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Cluster Analysis

- Hierarchical Methods
 - Can be **agglomerative** or **divisive**



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Cluster Analysis

- Hierarchical Methods
 - Identifies homogeneous groups of variables by using an algorithm that:
 - either starts with each observation in a separate cluster and combines clusters until only one is left (agglomerative),
 - or starts with the whole dataset and proceeds to divide it into successively smaller clusters (divisive).

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Cluster Analysis

- Partitioning Methods
 - Based on specifying an initial number of groups, and iteratively reallocating observations between groups until some equilibrium is attained
 - The most popular method of partitioning is the *k-means* method

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Cluster Analysis – Partitioning methods

- K-Means Clustering
 - Observations are classified as belonging to one of k groups
 - Group membership is determined by calculating the centroid for each group, then assigning each observation to the group with the nearest centroid

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Spatial Autocorrelation

- ▶ Nearest Neighbor Analysis, Quadrat Analysis and Cluster Analysis are limited as they only consider either the location or the attribute of the observations
- ▶ Spatial autocorrelation detects spatial patterns of a data distribution by considering both the **locations** and the **attributes** of the observations

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Spatial Autocorrelation

- Spatial Autocorrelation – correlation of a variable with itself through geographic space.
 - The first law of geography: “everything is related to everything else, but near things are more related than distant things” – Waldo Tobler
 - If there is a systematic pattern in the spatial distribution of a variable, then there is spatial autocorrelation
 - *positive spatial autocorrelation* – neighbors are more alike
 - *negative autocorrelation* – neighbors are not alike/different
 - *no spatial autocorrelation* – Random distribution

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Spatial Autocorrelation

- Most statistical measures are based on the assumption that the values of observations in each sample are independent of one another (this is why we carry out random sampling)
- This independence may be violated by spatial autocorrelation. For example, positive spatial autocorrelation may violate the independence if the samples were taken from nearby areas
- Therefore it may lead to incorrect conclusions about relationships between variables
- Goals of spatial autocorrelation are
 - To measure the strength of spatial autocorrelation for a given distribution
 - To test the assumption of independence or randomness

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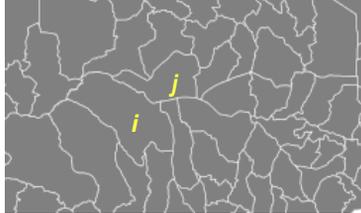
Moran's I Statistic

- Moran's I Statistic is a spatial autocorrelation technique that calculates the relationship between locations of observations (w_{ij}) the similarity between the attributes (c_{ij}) at those locations
- Given a set of features and an associated attribute, it evaluates whether the spatial data distribution is clustered, dispersed or random.
- It is applied to zones or points with continuous variables associated with them.

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Moran's I Statistic

If zone i is neighboring zone j



Similarity of Attribute c_{ij} $c_{ij} = (x_i - \bar{x})(x_j - \bar{x})$

Proximity of Location w_{ij} $w_{ij} = 1/d_{ij}$

Sample Variance s^2 $s^2 = \sum (x_i - \bar{x})^2$

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-
- Combines the measurement for attribute similarity and location proximity

$$I = \frac{\sum_{i=1}^n \sum_{j=1}^n w_{ij} c_{ij}}{s^2 \sum_{i=1}^n \sum_{i=1}^n w_{ij}}$$

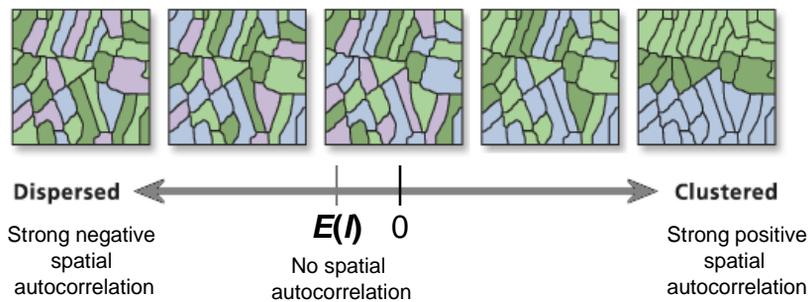
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- The numeric scale of Moran's I is anchored at the expected value of

$$E(I) = \frac{-1}{n-1}$$

- A Moran's I value equal to $E(I)$ demonstrates that the pattern has no spatial autocorrelation and hence can be considered random

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Testing for Significance

$$Z = \frac{I - E_I}{S^2}$$

Assuming a two-tailed test at the 0.05 significance level, the observed degree of spatial autocorrelation is significant if it is beyond the critical value of

$$Z = -1.96 \text{ or } +1.96$$

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Minimum Sample Size

- ▶ In statistics, it is often necessary to determine if our sample size is large enough to consider our results valid
- ▶ If the number of observations is too small the sample will not adequately represent the population
- ▶ Therefore, it is important to calculate the *minimum sample size* if there is spatial autocorrelation.
 - ▶ Generally the sample size will be larger than if spatial autocorrelation is not present
 - ▶ Different types of statistics have their methods for way calculating the *minimum sample size*

References

- David O'Sullivan, David Unwin, 2010
Geographic Information Analysis. Hoboken, NJ, John Wiley
- David W. S. Wong, Jay Lee, 2005. *Statistical Analysis of Geographic Information with ArcView GIS and ArcGIS*. Hoboken, NJ, John Wiley