GEOSTATISTICAL MODELLING OF ANTHROPIC ACTIVITY MARKERS IN THE PRESENCE OF BARRIERS.

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ABSTRACT KEYWORDS HIGHLIGHTS The analysis of chemical soil composition in archaeological domestic floors is becoming increasingly considered as an important topic for historical research. Mapping the distribution of certain elements allows us to understand the activities that were developed in these areas. This assumption comes from the idea that different social actions of production, consumption or distribution are the cause of variations observed in the material consequences detected through fieldwork. In this case, the variability on chemical soil composition is thought to be a valuable marker in order to detect, identify and analyse different activities in domestic contexts (Rondelli et al. 2014, Salisbury 2013, Middleton et al. 2010). Nonetheless, the reading of chemical differential concentrations in archaeological floors is not exempt of critical reflections about its limitations (Vyncke et al. 2011, Dore and López Varela 2010, Wells 2010, Terry et al. 2004).

Geostatistical methods are usually used to model the results of this kind of analyses and hence facilitate its interpretation. These techniques provide a set of statistical tools specifically designed for spatial problems, in which prediction is required over a region of interest where some observations have been taken. Predictions are based on an underlying statistical model that can take additional information into account as explanatory variables. In addition, the prediction error can be estimated based on propagation of uncertainty. One of the main limits of the use of geostatistical methods for that purpose is the assumption of a homogeneous, unrestricted space of analysis. This premise fail when we consider the spatial demarcations and topography that affect the distribution of our phenomena over the study region. The analysis of chemical residues in a domestic unit floors is a classic example of this kind of situation, in which the walls of the house affect to the diffusion of these elements.

In this work, we develop a methodology for overcoming this problem. We propose the use of cost-based distances to quantify the correlation between sampling locations. In this way, we present an interesting case study about the distribution of chemical elements in domestic floors within a controlled ethnographical context in Northern Gujarat (India) (Rondelli et al. 2014). This paper explores the relative spatial variability of these residues, taking into consideration spatial demarcations, to provide a method for the detection and interpretation of specific areas of activity. Our technique, therefore, might improve substantially the identification of both clustering patterns and different processes of floor maintenance and postdepositional dynamics considered as a background noise (Rondelli et al. 2014, Pecci et al. 2013, Barba 2007, Lloyd and Atkinson 2004).

2. ON THE USE OF NON-EUCLIDEAN DISTANCES IN GEOSTATISTICS

Geostatistics is a branch of statistics that encompasses the techniques that apply to geographical analysis. We owe its origins to the works of D.G. Krige (1951) and G. Matheron (1963) in the central decades of the 20th century. There are several applications of geostatistical methods in a very wide range of disciplines, with the common problem of modelling a Stochastic Process over a continuous spatial region from a partial group of observations. This process of interpolation is commonly assumed to be Gaussian, isotropic and intrinsically stationary (Cressie 1993). Geostatistical modelling is based on the principle of spatial dependence, this is near events are more related than distant ones. Nevertheless, what means near and how we calculate it?

Interpolation techniques assume that the correlation between the elements of a group of observations is a function of the Euclidean distance between them. In other words, stationarity is often accepted to mean that the spatial point process has constant intensity and uniform correlation depending only on the lag vector between pairs of points (Møller and Toftaker 2012). Considering the inherent irregularity of geographical terrain, either the presence of barriers or the difficulty to cross a region are presented as a major problem for this technical requirement. Imagine two locations at a given (Euclidean) distance such that they are significatively correlated, because of underlying relevant factors affecting both of them. Now put a barrier between them that blocks or absorbs the effect of the underlying factors. This obviously pulls the correlation down. Therefore, when some kind of barriers exist, the correlation depends on something other than this kind of metrics, which therefore cannot account for correlation by itself.

Methodologically, the first step in <u>classical</u> geostatistical processing is to fit the data and its empirical semivariogram function to a known parametric model. There is a variety of methods for estimating this correlation. Our approach here is to use maximum likelihood methods that fits the mean function value and the parameters of the semivariogram function. Once fitted, the main interest for analysing purposes is spatial prediction. Kriging is one of the most used approaches to this problem, in which a weighted average of the sample values is used to generate the prediction. This is, sample points near the prediction's location are given larger weights than those far away. Kriging determines these weights based on the semivariogram function, calculating them according to the value of the semivariogram, which is a function of the Euclidean distance (López-Quílez and Muñoz 2009). That seems to incur into the above mentioned error of assuming the validity of the spatial homogeneity premise. Thus, in certain cases, alternative measurements to Euclidean metrics, such as cost-based. Riemannian or pseudo-Euclidean ones, represent the distance argument r of the semivariogram function more naturally.

2.1 COST-BASED DISTANCES

Alternative measures to Euclidean distances have been largely tested in several disciplines. A multidimensional scaled reconfiguration of the spatial distribution has proved to be very useful in some cases, allowing to create a pseudo-Euclidean framework on which the analysis can be performed (Løland and Høst 2003; Negre 2015). A fast Fourier Transform has also been checked for integrating moving-average functions that may be used to generate a large class of valid, flexible variogram models. This transform allows us both to compute the cross-variogram on a set of discrete lags and to interpolate the cross-variogram for any continuous lag (Ver Hoef et al. 2004). In this same direction, recent works also propose the use Riemannian metrics associated to cost-based distances and Banach algebra using Kuratowski immersion (Muñoz 2012: 118). For its relative simple implementation, the use of cost-based distances directly into the covariance matrix of the Kriging, has proved to be a practical and competitive option for our research topic.

From a methodological perspective, the main goal of cost-based distances is to define the least cost path to reach a known point from each cell location in the original raster dataset. The calculation algorithms present the length of the irregular vectors formed by a spatial distribution using the shortest weighted distance is is the path with least accumulated cost. For allowing

these calculations, first we need to create a cost surface, which purpose is to assign an impedance value to each cell of a raster layer, that is, the ease with which it can be crossed. Formally, the resulting cost-weighted model can be defined as an f-function f, which describes for each cell of our model a real, positive value representing the difficulty to go through them, that is, its cost-weighted density. Therefore, the cost-weighted movement of a displacement that the point is a from this function, the cost of any path in A can be calculated as the integration of every cell in the model, which is gone through (Muñoz 2012: 55). Thus, the cost of a path α between points will be



These distances also maintain the same general properties than their metric counterparts (Waller and Gotway 2004: 321):

non-negativity



symmetry



triangle inequality



That implies that these measurements could be used into geostatistical functions. In spite of that, the results might be, in some cases, statistically irrelevant. The more homogeneous is the surface under study, the less significant are the changes with respect to the use of Euclidean measures. To obtain mathematical validity, the resultings covariance matrix of the observations must check thebe positive definitions of the covariance matrix of the observations. This condition requires that for any locations and complex set of coefficients , the function verify the next relationship:



where represents the cost-weighted distance between their arguments (Muñoz 2012: 201). Ultimately, and repovided verification of the above validation factors, a functional model be described providing the best linear unbiased prediction.

The adaptation of geostatistical analysis computation to these metrics is present in three major stages: empirical variogram computation, variogram model parameter fitting and the actual kriging prediction. Apart from observation data and prediction locations needed for standard kriging, we also need two Cost-Based distance matrices previously computed. One holdings the distances between observation points, a symmetric square matrix. The second matrix holdsing the distances between observation points and the prediction location(s), so it is an n (observations) x m (locations) sized matrix.

The empirical variogram is computed from the observation data only. It classifies pairs of observations into groups according to their distance, and then computes an estimator of the theoretical variogram value for that distance based on the differences between the observed values.

In order to make a cost-based empirical variogram it is enough to make the initial classification based on the cost-based distance values given in the corresponding matrix, rather than calculating Euclidean distances. Note that this modification produces a different grouping of observation pairs. Therefore, variogram estimates will be different. The variogram model parameter fitting is also made based on observation data only. It is typically accomplished through maximum likelihood methods, trying out many possible combinations iteratively and keeping the best. This implies computation of the covariance matrix for each combination being tested. All we need is to make sure that the covariance matrix is computed based on the cost-based distances provided by our previously computed matrix. Finally, there is the kriging prediction. At this point, the covariance model is assumed to be known. Here again, we need to make sure that the covariance matrix of the observations is computed with the cost-based distances. In addition, the covariance between observation points and prediction locations are to be computed in order to make predictions. So this is when the second of the cost-based distance matrices is to be used.

2.2 METHODOLOGICAL OVERVIEW

2.2.1 Computing the cost-surface

The first operation is to encode the spatial heterogeneity of the working area into one *cost-surface*. This imply some modelling decisions and assumptions, which are not technical but scientific in nature. In our case, we assume that the soil is homogeneous except for the areas with solid sunken structures. These structures will completely interrupt the continuity of the area, blocking the dispersion of substances.

This conceptual model yields a cost-surface with a constant value of 1 everywhere, except over the structures where it takes an infinite value. In practice, any value larger than the diameter of the region will suffice. Alternatively, can be more practical to work in the inverted scale of a conductivity surface. In this case, the values would be simply 1 for regular conductivity and 0 for no conductivity, or infinite cost. Any of these alternative surfaces can be easily produced from a digital representation of the region with any GIS software, or with other spatially capable software like (R Core Team, 2015).

In our case, we imported the ESRI shapefiles describing the geometry of the structures into spatial classes defined in the package (Pebesma and Bivand, 2005). Then we used functions from the rasterthe function rasterize() from the raster package (Hijmans, 2015) to produce a discretized surface with constant value 1 over the region of interest, and over the solid structures. the function to determine which raster cells fall within the boundaries of the solid structures. These raster cells are then assigned a value of 0. We used a resolution of about 520 pixels/m.

2.2.2 Computing the cost-based distances

The cost/conductivity surface is the object representing our model of the region, and from where the distances between locations can be computed. Specifically, two matrices of cost-based distances are required: one matrix with distances among the observations, and one with the distances between each observation to each one of the prediction locations.

We used the centroids of the conductivity raster cells as prediction locations to simplify mapping, although any set of prediction locations can be used. The computation of the distance matrices can be also be performed using a GIS software or right within . López-Quílez and Muñoz (2009) use the first approach with the help of a specific script (Muñoz 2015b) for GRASS GIS (GRASS Development Team, 2010). For this study, we used the package geoRcb (Muñoz 2015a) (van Etten, 2015) instead.

This package provides the function `distmatGen()`, which automatically computes the two cost-based matrices given the coordinates of the observations and the conductivity surface. Internally, it leverages the package gdistance (van Etten, 2015) for efficient computation of least-cost paths, while taking care of all the technical details. Ultimately, the costbased distance is computed using the This package provides the function , which computes the transition value from every raster cell to every other adjacent cell. This value is computed after a user-specified function of the corresponding conductivity values. Here we used the function , so that the transition value between two cells any of which with conductivity of 0 remains 0. We considered adjacent all the 16 cells that can be reached by one move of the king or the knight pieces in the game of Chess. The transition values are corrected for diagonal connections. These settings provide the maximum accuracy for the subsequent calculations of distances based on a discretized surface. Finally, the function -uses the transition values to compute least-cost distances between two sets of points. It is based on the well-known Dijkstra's algorithm for finding shortest paths between nodes in a network.

2.2.3 Using cost-based geostatistical algorithms

3 The geoRcb 💆 package -(Muñoz, 2015a) is an extension extends 2015) that provide alternative version of some functions, in order to make them capable of working with cost-based distances. Specifically, the 3 feature an additional argument functions and which takes a symmetric matrix of distances between observation locations. These functions are respectively used to compute empirical variograms and fit variogram models. Finally, the alternative function 3 performs the cost-based spatial prediction through conventional kriging by taking the required distance matrices as the additional arguments dd.dists.mat and 3

2.2.4 Presentation of results

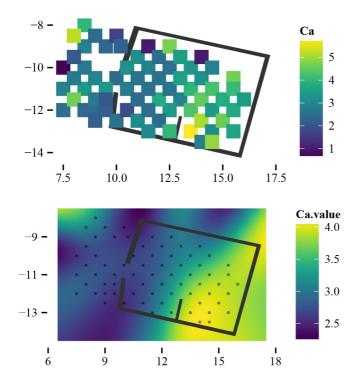
We use the viridis color palette (Garnier, 2015) for all the maps in the present paper. This palette is perceptually-uniform and is also designed to be perceived by readers with the most common form of color blindness.

3. CHEMICAL RESIDUES ANALYSIS ON ARCHAEOLOGICAL FLOORS IN THE PRESENCE OF BARRIERS

3.1 CONTEXT AND DATA DESCRIPTION

The data used in this methodological presentation was obtained from a fieldwork campaign carried out in a domestic compound in Jandhala, a

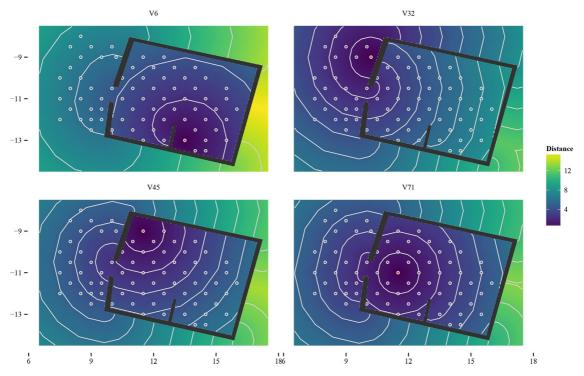
village in the Patan district of North Gujarat, India. The compound is a nonmechanised farmer's residence built using local materials and traditional techniques. Several researchers of the North Gujarat Archaeological Project (NoGAP) participated on taking chemical soil samples in one of the domestic units of the compound between 2009 and 2010. A complete ethnoarchaeological study was developed in order to define the types and cycles of activities carried out in the compound, as well as their spatial distribution within the courtyard, veranda inner and Micromorphological floor samples were recovered during these works in order to obtain, among others, the chemical signatures of the average indicators of all the residues that fell on the floor since the foundation of the house, 10 years before sampling. Theses analyses were carried out with Inductively Coupled Plasma - Atomic Emission Spectrometry (ICP-AES), providing quantitative data on chemical elements. Finally, a complete study of the use of space based on floor chemical analysis was published (Rondelli et al. 2014). This is not the main goal of our work to analyse the specific use of space in this case study, but to confront the previous geostatistical results with the ones provided using our technique.



For testing purposes, we want to model a chemical spatial distribution map of Calcium residues in one of the domestic units of the Jandhala compound. Seventy measurements have been made in various points distributed homogopusly over the house and its courtyard and veranda. The Calcium values from a minimum of 0.66 to a maximum of 5.70, being its their mean value 3.11 and its their median 2.95 pure 1 shows the spatial arrangement and values of the measurements. The data presents a Normal or Gaussian distribution of values, which stands for a non-random process (Barceló 2007: 45-49)

The walls of the house are considered non-transparent barriers for the diffusion of chemical residues on the archaeological floors. For this reason, cost-based distances betweeneach sampling points to everywhere else have

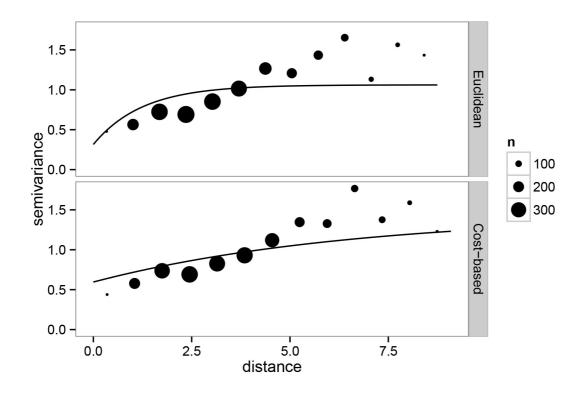
been calculated in order to create a symmetric matrix of distances between observation locations. Here, several examples of this calculations we show the cost-based distance maps to four selected points, which show how these distances honour the geometry of the region:



3.2 VARIOGRAM AND FITTED MODEL

The We fitted an exponential variogram model using both Euclidean and cost-based distances is exponential. We choose to estimate the nugget effect, which may account for measurement error, for example. The main difference being that the cost-based approach yields a larger range. This is usually the case, as incorporating the geometry into the analysis helps making sense of what was before interpreted as unstructured noise. Also the nugget is higher in the estimated cost-based variogram, which will result in increased smoothing.

In the scatter plot, the horizontal patterns correspond to probberved values. Otherwise, the differences are negligible. The area the observations, the cost-based approach has a larger prediction error due to its increased estimation of the nugget (i.e. short-range variance). In the main area, the prediction errors are practically the same with both appreparates. Behind the walls, the Euclidean prediction error is unrealistically low.



	Euclidean	Cost_based
Nugget	0.32	0.60
Partial sill	0.75	0.85
phi	1.25	6.53
Pract. range	3.75	19.56

Once again, cost-based distance explains the correlation between locations better than Euclidean distance. A natural question to ask is how different the values from the two types of distances are. The following plot regarding pointwise comparison of predictions allows observe the high variability of the calculation depending on the method.

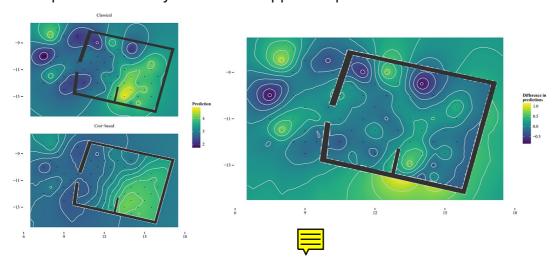
3.3 COMPARISON OF KRIGING ESTIMATES AND PREDICTED ERRORS

In the following image, it can be observed the maps with the prediction for each location using Euclidean and Cost-based calculations under the same scale, and their difference. It is interesting to compare both results. There are three remarkable aspects to be noted. First, the highest differences in prediction happen in the outer unsampled area, specially just across the wall from the most extreme observations. The Euclidean prediction is strongly influenced by these values, in contrast to the average value predicted by the cost-based approach. Second, the cost-based prediction is clearly smoother, as expected given the parameter estimates of the variogram model. This produces important differences in the neighbourhoods of the most extreme measurements. Finally, the walls of the house modify the prediction even in the sampled area. For example, in the corner near the central inner wall, across from the highest measurement, the Euclidean method predicts higher, undoubtedly under the influence of the observation.

On the contrary, in the veranda, the Euclidean predictions are lower, influeced by the observations in the interior.

In summary, taking into account the heterogeneities of the working region through the use of the cost-based distance leads to differences in predictions that can be split into these three main aspects. We believe that the cost-based approach produces more reliable results, as it avoids assuming certain unrealistic assumptions about the homogeneity of the region.

The walls of the house modify the prediction as well as the method per se, as the map of predictions differences shows. Taking into account only the house and the veranda areas, cost-based approach is highly more accurate The area of Calcium concentration is better defined through our approach, as well as the false positive areas, such as the external perimeter of the house or in the veranda. The chemical representation of this element over the floor is perfectly demarcated close to the interior wall of the house, just in the south-east corner of it. That allows a better interpretation of the spatial distribution of this element and, especially, a better integration with other geochemical residues. Archaeological reading of chemical signatures in domestic floors is highly related to the combination of diverse elements, such as different combinations of Potassium, Calcium, Phosphorus. Magnesium and Strontium for deposited food remains, living room sediments, enclosed spaces and burning areas (Rondelli et al. 2014: 488). Calcium has long been regarded as good indicator of human activity, usually correlated with Strontium, suggesting similar chemical pathways. Usually, these combined elements are interpreted as good indicators of covered, enclosed spaces. They should, then, serve as indicators of the integrity of anthropogenic deposits (Middleton et al. 1996: 679). In the presence of several geochemical indicators related at the same time, a perfect delimitation of their distribution areas result crucial for a valid inference. That plus of accuracy is what our approach provides.



4. DISCUSSION

The most important aspect of this work lies in the general methodology for overcoming the geostatistical restriction on the homogeneity of the prediction region. In the study of chemical residues on archaeological floors, that limitation is a major problem in order to obtain significant interpretation of the sampling data. Thus, more accurate spatial distribution map of soil composition elements benefits from this methodology, since walls in

domestic contexts are relevant restrictions in its diffusion. Furthermore, the possibility of applying geostatistical techniques enables us to obtain results based on statistical models, providing reliable predictions together with estimations of uncertainty, which commonly used deterministic methods cannot provide. Altogether, all this information allows the correct interpretation of the archaeological data, both the distribution of the chemicals and the postdepositional effects over them, linked to the intensity of the uncertainty measure, among other factors.

To conclude,

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