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"Analyzing Field-Measured Soil Properties"

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ANALYZING FIELD-MEASURED SOIL-WATER PROPERTIES

D.R. Nielsen, Patricia M. Tillotson and S.R. Vieira

1. INTRODUCTION

It is a challenge to the world community of earth and geophysical scientists to develop a better technology for sampling the earth's crust. That thin mantle of soil has been managed for countless human generations with the primary objective being the production of food and fiber to meet the needs of the earth's inhabitants. In the past, that management has been judged on annual measurements of crop productivity, and not on measurements taken below the soil surface that could be used to signal the long term consequences of present-day management of soil and water resources. We are indebted to the pioneering works of R. A. Fisher and subsequent efforts of others that have and continue to afford conceptual frameworks to statistically judge and compare the merits of different management schemes or treatments, particularly those used for enhancing agricultural production. The selection of a preferred cultivar, an optimal fertilizer application, the best timing of an irrigation or the most effective soil fumigant has been accomplished in the agricultural sciences by analysis of variance procedures advocated by Fisher. Such "aggie" statistical procedures are invaluable, and it is not our intent to de-emphasize their importance now or in the future. Our intent here is to expand that conceptual framework to include a consideration of statistical analyses normally not included in the agricultural sciences. Such an expansion is fully justified when we wish to examine the changing quality of soils as well as that of water moving over and through them as a result of different management schemes. We believe more attention should be given to developing techniques to better monitor the soil environment, and at the same time we recognize that such development should, concomitantly, potentially enhance the efficiency of crop production.

Our objective of this presentation is to provide a qualitative review of statistical concepts not usually covered in "aggie" statistics, and to provide an opinion of the questions we believe future research shall answer in light of the kinds of efforts reported at the 1982 meeting of the European Geophysical Society. Our intentions preclude the identification of analytical prescriptions or algorithms to carry out various statistical procedures. We also shall not attempt to be rigorous in the identification of the fundamental underlying assumptions of each concept. The concepts have been known and used in other scientific disciplines for a relatively long time. If our presentation is successful, we would urge the reader to refer to the list of references to learn and fully appreciate the fundamentals

of each concept, and not to rely on our qualitative descriptions and examples.

2. SPATIAL AND TEMPORAL DEPENDENCE

Our reference to "aggie" statistics highlights the fact that most statistical analyses advocated in the agricultural sciences implicitly disregard the spatial coordinate at which an observation is made. For the most part, emphasis is given to the identification of an average value and its potential dispersion for a soil attribute within a given parcel of land, a regression of one attribute versus one or more other attributes relative to their magnitudes (not their coordinate positions on the landscape), and a difference between two mean values of a soil attribute that may exist for two parcels of land chosen more or less arbitrarily without regard to their spatial coordinate system. In fact, it is generally considered necessary or advantageous in "aggie" statistics to assume that observations are spatially independent of each other, and hence, a set of observations are reduced to their mean value and a measure of its uncertainty expressed in terms of an assumed probability density distribution estimated by a set of observations without regard to their spatial positions. We refer the reader to standard texts for such analyses.

Intuitively, we do not expect field observations of soil properties to be necessarily spatially independent. We would expect measurements made close together to yield nearly equal values, and measurements made some distance apart to yield values more correlated to each other. We would also expect a spatially repetitive behavior of soil observations as a result of cyclic tillage traffic and cropping patterns in cultivated fields, sequences of low and high topographical positions giving rise to cyclic locations of greater and lesser degrees of leaching, and sequences of soil mapping units not randomly located within a landscape owing to soil formation processes that are linked spatially to the coordinates of the soil surface. Because of the above expectations, it would appear advantageous to sample a field in a manner that would allow the detection of cyclic irregularities in relation to the size of the parcel of land being measured. Such expectations raise questions regarding the "proper" size of an observation, the "proper" distance between observations, the "proper" location of each observation and the "proper" number of observations. These questions are all relevant to the geostatistical concept which defines the size of the domain characterized by a single observation within a field soil. This differs markedly from the "aggie" statistical concept that defines for a given level of probability the accuracy and precision of an estimate of an average value within a field soil from a set of observations.

2.1 Spatial autocorrelation

A measure of the strength of the linear association between pairs of observations is useful in defining the separation distance between observations beyond which there is no correlation between pairs of

values. The autocorrelation coefficient, r , a function of the separation distance h , is a measure of that strength and is defined as

$$r_g(h) = \frac{\text{autocovar}[G(x), G(x+h)]}{\sqrt{\text{var}[G(x)] \text{var}[G(x+h)]}} \quad (1)$$

for a set of soil water content observations G taken along a transect in the x -direction*. For example, when observations are taken 1 unit apart, $r_g(1)$ is the value of the linear regression coefficient for $h = 1$ (lag 1) when values of $G(x+1)$ are plotted against values of $G(x)$. In other words, nearest neighbors are plotted against each other. Similarly, when $h = 2$, $r_g(2)$ is the value of the coefficient when values of G are plotted against other values observed a distance of 2 units away. Fig. 1 shows two examples. For example A, the autocorrelation coefficient decreases abruptly from 1 to zero within a lag of 1 (the smallest distance between observations). The value

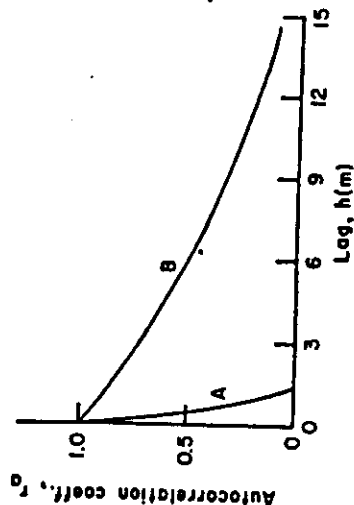


Fig. 1. Idealized autocorrelograms for average soil water content G observed at 1-m intervals along transects within fields A and B.

of r for example B approaches zero for lags in excess of 10. For example A, one concludes that the observations are spatially independent - that is, one cannot estimate the value of an observation from that of its nearest neighbor. Interpreting the results for example B where r decreases rather gradually as h increases, one concludes that

* Throughout this presentation, concepts expressed in one direction can be generalized to n directions.

the observations are spatially dependent. In other words, within a distance of about 10, it is possible to estimate from one neighbor the expected values of other neighbors. Soil scientists are beginning to use this concept to express the spatial dependence of field-measured soil properties and crop yields.

The functional relation $r_g(h)$ has been expressed with several empirical formulae. One of the most commonly used expressions is

$$r_g(h) = \exp(-h/\lambda) \quad (2)$$

where λ is selected in order that the sum of the measured deviations of r from the above expression is zero. Notice that the value of λ is equal to that distance h between measured values for which their correlation coefficient is $1/e$. λ is called the autocorrelation length or the scale of observation. A rather liberal interpretation of the significance of λ is that it represents the distance across the landscape characterized by a single observation within the field. In general, there are several ways of defining autocorrelation lengths. Important to our discussion here is that when sampling the field from which curve B (Fig. 1) was obtained, observations made at sampling intervals less than 10 units are somewhat unnecessary because they are related to each other. On the other hand, sampling the field from which curve A was obtained at sampling intervals greater than 1 unit does not allow meaningful interpolation between neighboring observations. It should be obvious that the functional relation between r and h depends upon the size of the sample, and that in general, the greater the sample size, the greater the value of the autocorrelation length λ . For any particular study, the investigator should consider the minimum distance between sampling locations ($h = 1$) in relation to the objectives of the experiment and the potential utility of the values of λ for each kind of observation.

2.2 Spatial cross-correlation

Instead of measuring only one kind of observation across a field, let us assume that two kinds are made: soil water content, G , as described above as well as the temperature, F , of a grain crop uniformly covering the field. The spatial cross-correlation coefficient r_c is defined by

$$r_c(h) = \frac{\text{covar}[F(x), G(x+h)]}{\sqrt{\text{var}[F(x)] \text{var}[G(x+h)]}} \quad (3)$$

where, in this case, F , the crop temperature and G , the soil water content, are each measured along a transect at positions x . Let us assume that as available soil water is depleted, evapotranspiration decreases and crop temperature, consequently, increases. For such a condition, crop temperature is inversely related to available soil water with the value of $r_c(0) < 0$. Equation (3) reduces to the linear

regression coefficient normally calculated using "aggie" statistics when $h = 0$ (a value of -0.8 for our example). Fig. 2 illustrates the use of equation (3) for other values of $h \geq 0$. In Fig. 2a, with the two distributions $F(x)$ and $G(x)$ overlapped a distance h , a linear regression analysis is calculated for the pairs $[F(x), G(x+h)]$. The results of such calculations of r_c are plotted in Fig. 2b for two hypothetical transects (one in field A and one in field B) having identical values of $r_c(0) = -0.8$. For field A, if the value of h is increased only slightly from 0, r_c rapidly approaches zero. On the

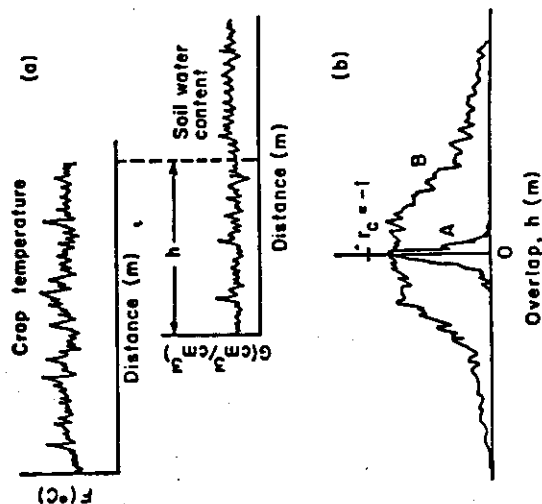


Fig. 2. Idealized cross-correlogram for crop temperature $F(x)$ and soil water content $G(x)$ along transects within fields A and B.

other hand, the results from field B show that crop temperatures measured at much greater distances from where the average soil water content observations were made remain significantly correlated. The general utility of such cross-correlation should be obvious. The area under each curve in Fig. 2b or the range of h over which the value of r_c remains near unity is an indication of the spatial distance

over which a linear relation exists between F and G . It allows the investigator to consider the spacing and size of one set of observation (F) with those of another (G), particularly when one set is difficult to obtain or relatively expensive. A related example would be the optimum choice of the size and spacing of observation pixels from overflight or satellite vehicles compared with size and spacing of those of ground observations.

2.3 Spectral analysis

In the paragraphs above, the interpretation of the correlation coefficients r_c and r_s of equations (1) and (3) were restricted to values of h for observations compared more or less in near vicinity to each other (i. e., for h much, much less than the width of the field being sampled). An opportunity to discern repetitious irregularities or cyclic patterns in soil or plant communities across a field exists with a spectral analysis that utilizes the function $r_s(h)$. We illustrate by assuming our measured distribution of average soil water content across a field $G(x)$ was taken where a crop had previously been grown along furrows 1 m apart. As a result of both plant extraction of soil water and infiltration occurring in 1-m cyclic patterns across the field, measured values of G will reflect local variations as well as a tendency toward a sinusoidal behavior having a 1-m period. A spectral analysis identifies this periodicity and can be calculated by

$$S(f) = 2 \int_0^p r_s(h) \cos(2\pi fh) dh \quad (4)$$

where f is the frequency equal to $1/p$ where p is the period. Fig. 3

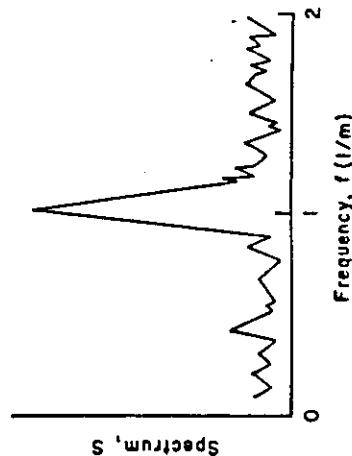


Fig. 3. Spectrogram $S(f)$ for a transect of soil water content observations taken normal to the direction of furrows.

illustrates the shape of the function $S(f)$ where it can be seen that most of the variance about the mean value of $G(x)$ is accounted for by observations of soil water content that reflect oscillations of wet and dry soil occurring, on the average, every 1 meter. Depending upon the kinds of repetitive features and processes that may be operating in a field, $S(f)$ may have a number of relative maxima that identifies their specific spatial occurrences. In other words, a spectral analysis is useful in partitioning the total variance of a set of observations among different frequencies and then assessing which of those frequencies has any significance for the field problem being studied.

If we extend the above example to conditions illustrated in Fig. 4, we see two additional periodicities - those greater water contents occurring every 2 meters owing to tractor tire compaction, and those occurring approximately every 10 meters associated with pre-plant border irrigation or some other kinds of previous traffic pattern. In this example, most of the total variance is accounted for by variations from the mean value occurring at periods of 1, 2 and 10 m. Had

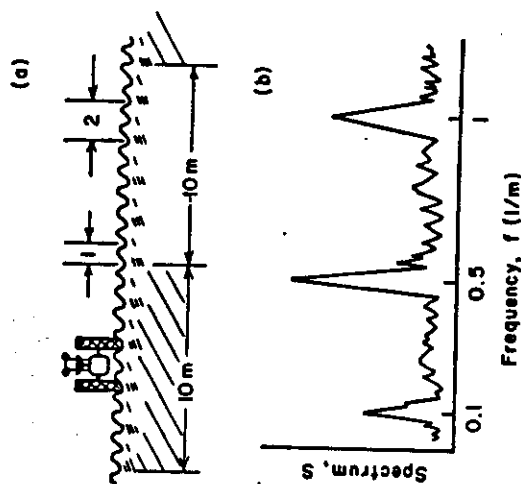


Fig. 4. Schematic diagram of furrows, tractor compaction and pre-plant irrigation causing cyclic variations of soil water content, and idealized spectrogram.

plant extraction, infiltration, compaction and pre-plant border irrigation had no influence on the spatial variations of average soil water content, $S(f)$ would have manifested no relative maxima.

2.4 Cospectral analysis

In a manner similar to the usage of $r(h)$ in spectral analysis, $r(h)$ is used to partition the total covariance for two sets of observations across a field. A cospectral analysis is made by

$$Co(f) = 2 \int_0^{\infty} \bar{r}_c(h) \cos(2\pi fh) dh \quad (5)$$

where $\bar{r}_c(h) = [r(h < 0) + r(h > 0)]/2$. Let us extend further our illustration given in Fig. 4 by assuming that a grain crop is growing in the field whose soil water content has cyclic distributions of soil water owing to furrows, compaction and pre-plant border irrigation. We have the distribution of crop temperatures $F(x)$ and the distribution of soil water contents $G(x)$ from which we calculate the cross-correlation coefficient $\bar{r}_c(h)$ from equation (3). Having calculated the average value $\bar{r}_c$ and integrating equation (5), we obtain $Co(f)$ depicted in Fig. 5. It is not surprising that three relative minima occur at periods of 1, 2 and 10 m. The area beneath the abscissa has a negative value and represents the total covariance between crop temperature and soil water content. A linear regression between $F(x)$ and $G(x)$ using "aggie" statistics would be negative indicating statistically that crop temperature is inversely related to

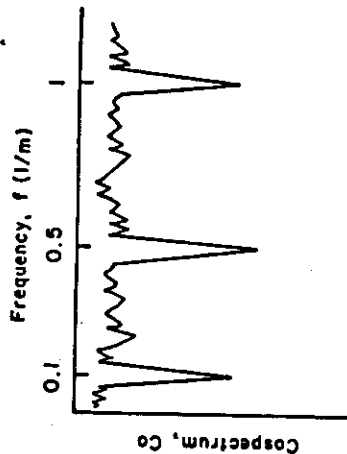


Fig. 5. Cospectrogram of crop temperatures $F(x)$ and soil water content $G(x)$ along a transect normal to the furrows shown schematically in Fig. 4a.

average soil water content. In this example, had pairs of observations of F and G been made in a spatially random manner across the field, the "aggie" linear regression correlation coefficient would have been approximately equal to $r(0)$ but such random samplings would not have identified the periodicity within both sets of observations.

For the above example it is advantageous for our discussion to recognize an alternative situation where soil compaction occurs with relatively higher soil water contents. For this situation, it is possible that root growth is impeded, or because of ever-present root-rot microorganisms thriving in a wet, compacted soil environment, the crop roots are diseased. In either case, the crop temperature would be directly related to soil water content in zones of compacted soil, and inversely related in other locations in the field. For such a situation, equation (5) would give rise to a cospectral analysis illustrated in Fig. 6. Even though the total area is zero, the relative minima and a maximum show inverse correlations between F and G at spatial periods of 1 and 10 m, and a direct correlation at a period of 2 m. Had pairs of F and G been taken randomly across the field, a routine linear regression would have provided no enlightenment of the processes occurring inasmuch as its value would have been near zero.

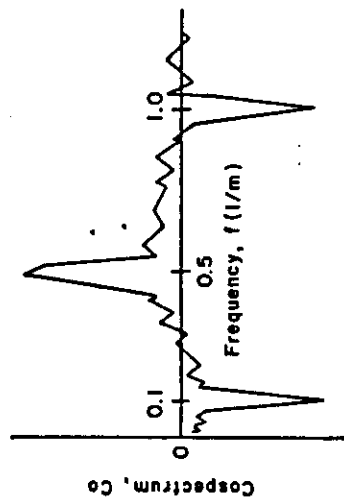


Fig. 6. Cospectrogram of crop temperatures F(x) and soil water content G(x) along a transect normal to the furrows shown in Fig. 4a when compaction decreases evapotranspiration.

2.5 Cospectral phase angles and coherence

The above examples of cospectral analysis compared two sets of observations whose periodicities were spatially equal for the same locations across the field. Soil attributes or processes that are

correlated for a particular spatial frequency may manifest this correlation at any phase angle of not necessarily zero. We illustrate this possibility in Fig. 7 where observations of crop temperature, T(x), and soil salinity, SS(x), have a periodicity of 1 m. We assume that crop temperatures are directly related to soil salinity and that soil salinity levels are smaller in the bottom of the furrows 1 m apart. For such conditions, each set of observations yields a relative maximum in its spectral analysis at a frequency of 1 as shown in the figure. However, because the sun's radiation is not received from a vertical direction, the maximum temperature of the crop will occur at a distance ϕ from where the highest soil salinity is observed. The phase angle ϕ illustrated in Fig. 7 is given by

$$\phi = \frac{1}{2\pi f} \tan^{-1} [Q(f)/Co(f)] \quad (6)$$

where Q(f) is the quadrature spectrum calculated in an equation similar to that of equation (5) where the cosine term has been replaced by

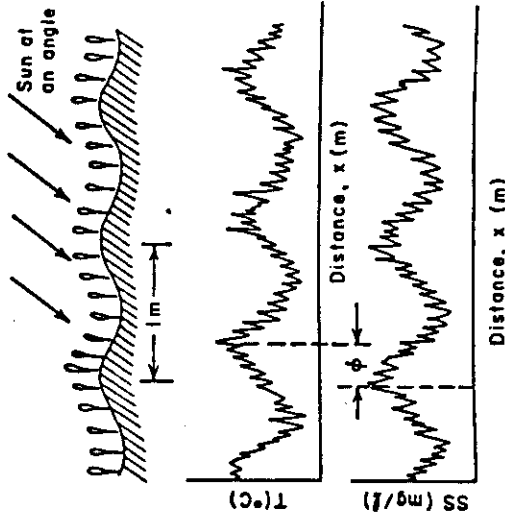


Fig. 7. Schematic diagram of observations of crop temperature T(x) and soil salinity SS(x) along a transect normal to furrows with the sun radiation being received from one side.

a sine term. Having calculated $r_c(h)$ for the two distributions $T(x)$ and $SS(x)$, $Q(f)$ can also be calculated from equation (5) if $r_c(h)$ is replaced by $r_c(h)$ where $r_c(h) = [r_c(h > 0) - r_c(h < 0)]/2$. The coherence of the cospectral analysis is given by

$$Ch(f) = \frac{Q^2(f) + Co^2(f)}{S_T(f)S_{SS}(f)} \quad (7)$$

where S_T and S_{SS} are calculated from equation (4) for T and SS , respectively. The coherence, whose values range between 0 and 1, is analogous to the r^2 -value in ordinary regression analysis of "aggie" statistics. It provides a measure of the certainty at which the phase angle is identified.

Spectral and cospectral analyses are potentially powerful tools for managing and increasing our knowledge of land resources. With them, we can spatially link observations of different physical, chemical, and biological phenomena. We can identify the existence and persistence of cyclic patterns across the landscape. In some cases, the cyclic behavior of soil attributes may be of more or equal importance than the average behavior. From a spectral analysis, some insights may be gained relative to the distances over which a meaningful selection of a particular size of a sensor should be based upon a knowledge of the potential periodicities to be manifested by such observations. And, with spectral analyses, it is possible to filter out trends across a field to examine more closely local variations, or vice versa. It should also be recognized that all of the above discussion from the beginning of this presentation could have had the time variable substituted for that of the distance variable.

2.6 Semivariograms and kriging

The spatial dependence of neighboring observations may also be expressed by the semivariogram $\gamma(h)$ estimated by

$$\gamma(h) = \frac{1}{2N(h)} \sum_{i=1}^{N(h)} [G(x_i) - G(x_i + h)]^2 \quad (8)$$

where $N(h)$ is the number of experimental pairs of observations $[G(x_i), G(x_i + h)]$ separated by a vector distance h . The shape of the semivariogram gives an indication of the spatial dependence of the soil physical properties. If for all values of h greater than zero γ remains essentially constant, it indicates that the observations are spatially independent. If for all values of h greater than zero, γ increases and approaches a constant value, it indicates that the observations are spatially dependent within a spatial area that can be characterized as a single domain. On the other hand, if as h increases, γ continues to increase, it indicates the area being sampled continually changes and is not comprised of a single domain. The

semivariogram describes the "variance structure" of a field, and is to geostatistics as the probability density distribution is to "aggie" statistics. Analytical interpretations of the shapes of variograms have proven especially helpful in the definition of soil mapping units as used in the context of soil morphology and classification.

Semivariograms, when the spatial dependence of observations exists, are useful to make interpretations between observed values and to identify improved future sampling schemes. Owing to the fact that the semivariogram gives the expected relation between pairs of observed neighbors, it is obvious for interpolation that different weights should be given to neighbor values depending upon their distance from the one to be interpolated. Hence,

$$G(x_0) = \sum_{i=1}^n \beta_i G(x_i) \quad (9)$$

where $G(x_i)$ is the estimated value at location x_i , β_i are the weights associated with each of the values G measured at locations x_i , and n is the number of locations. This interpolation method developed by G. Matheron, which he called kriging in honor of D. G. Krige is an optimum interpolator because it interpolates without bias and with minimum variance. And, because it allows the variance of the estimates to be estimated, it is extremely helpful for identifying improved sampling schemes. Kriging is becoming more common in geophysical studies, and in light of the presentation of Webster and Burgess, we shall not illustrate its usage. We point out, however, that through its usage for soil observations taken within and between furrows, or within and between rows of crop plants it is possible to construct contours of isovalues that are more meaningful than simplistic values of means calculated for the two positions - within the row and between the row, frequently repeated in agronomic journals.

2.7 Cross-semivariograms and cokriging

In many field situations, one set of observations $G(x_i)$ may not be sampled sufficiently to yield interpolated values at other locations of acceptable accuracy. By considering the spatial correlation that may exist between that variable and another more frequently observed variable $H(x_i)$, cokriging may improve the precision of estimating the former. Cokriging relies not only upon the semivariogram but also on the cross-semivariogram estimated by

$$\gamma_c(h) = \frac{1}{2N(h)} \sum_{i=1}^{N(h)} [G(x_i) - G(x_i + h)][H(x_i) - H(x_i + h)] \quad (10)$$

where the pairs of values $[G(x_i), G(x_i + h)]$, $[H(x_i), H(x_i + h)]$ are separated by the vector h . An interpolated value of G at location x_0 is cokriged using

$$G(x_0) = \sum_{i=1}^n \beta_i^H(x_1) + \sum_{j=1}^m \beta_j^G(x_1) \quad (11)$$

where β_i^H and β_j^G are the weights associated with observations H and G , and n and m are the number of H and G used in the estimation of G at location x_0 . Values of H and G are measured while values of β_i^H and β_j^G are calculated from values of the semivariogram and cross-semivariogram given in equations (10) and (8), respectively. Cokriging is attractive when one of the two sets of observations correlated with each other is relatively inexpensive and abundant. We believe it has the potential of more precisely delineating boundaries between soil mapping units, especially when either or both G and H above are each functions of several soil properties. Present-day and future soil mapping units should be judged, in some degree, on the behavior of their semivariograms and cross-semivariograms.

3. DETERMINISTIC VERSUS STOCHASTIC EQUATIONS

Differential equations derived and solved for the description of soil processes have most often been based upon deterministic concepts regarding both their variables and parameters. Their derivations explicitly demanded that sufficient observations were available to identify the expected value of each term with more than sufficient precision deemed necessary for their accurate solution. Such equations, often used under strict laboratory conditions where observations of each term were reasonably precise and accurate, are now being questioned when their solutions are extended to natural field conditions. That questioning initially embraced both measurement error and sampling error. But with the continual development of instruments and methodology to assess soil properties *in situ* with greater accuracy, it is now recognized that the sampling error associated with spatial and temporal variations of those properties must be considered separately. That is, with any reasonably affordable present day sampling scheme, the spatial and temporal variance of field soil properties are sufficient to render estimates of their means highly unreliable. Hence, deterministic equations are giving way to mixed deterministic - stochastic and stochastic equations, with levels of probability of their solution defined.

3.1 Scaling

Scaling is part of a more general methodology known as fractional analysis which seeks to find partial solutions to physical problems which cannot be solved explicitly. Their complete solutions are unattainable owing to either some lack of understanding or the mathematical analysis is intractable. An example of such a problem is the simultaneous transport of water, solutes, heat and gases within an unsaturated field soil subjected to diurnal conditions. From a theoretical viewpoint, scaling is a process which reduces through dimensional or inspectional analysis the number of variables important in a given problem to the smallest number of variables which completely

describe the system. This reduction greatly simplifies the description of the system as well as provides descriptions of a great number of other systems having different values for common parameters. More than 25 years ago, E. E. Miller and R. D. Miller introduced scaling and the concept of similar media using inspectional analysis for describing the retention and transport of water through unsaturated soils. They and others have derived scale factors for soil water flow properties that have shown utility for structureless soils comprised of particles of different sizes having similar geometric shapes. Soil scientists have yet to address the scaling of solute, gas or heat movement through soils by dimensional or inspectional analysis, and further consideration of that for water movement appears justified.

3.2 Stochastic equations

More recently different kinds of scale factors, not necessarily related to those mentioned above, have been identified through regression analysis to aid in the quantitative description of field-measured soil-water functions or parameters required for the solution of deterministic equations. Such identification simplifies the description of the functions for a spatially variable field soil into one or more stochastic scale factors. Hence, the precision for which soil-water functions are known for a field can be ascertained from the probability density function and the spatial variance of these scale factors. Incorporating such scale factors with mathematical expressions of their statistical and geostatistical variances into the usual deterministic equations for soil water allows solutions and simulations of soil water and related processes to be calculated within prescribed levels of probability. The practicality of routinely using stochastic scale factors would be enhanced considerably if they could be adequately estimated through correlations of easily measured soil properties such as soil texture, porosity, bulk density, etc. potentially available for each soil mapping unit.

It is becoming clear that there is a need to analyze and simulate the behavior of field soils and agronomic regimes using equations in probabilistic viewpoints. It is also becoming obvious that a knowledge of the mean behavior of a field may be of less importance in some cases than that of its statistical or spatial variance. In this respect, solutions of deterministic equations calculated repeatedly using Monte Carlo procedures to identify realizations of their soil parameters are beneficial. Such repeated calculations theoretically correspond to a series of repeated field-measured values. Analytic solutions of deterministic equations containing random or regionalized variables are also gaining recognition. Stochastic differential equations beginning to be used in hydrology relating variations in the saturated hydraulic conductivity to the dispersion of solutes need to be explored for unsaturated soils. There also is the possibility of using transfer function models that treat the transformation of a soil profile input into an output without a knowledge or modelling of the mechanisms inside the profile. Such models have been used recently to investigate the transfer of solutes added to the soil surface to

greater depths in the soil profile. Physical, chemical and biological reactions of the solutes within the profile were explicitly ignored but implicitly included through estimated probabilities of the amount and rate at which the solutes arrived at a greater soil depth.

4. FUTURE RESEARCH

A dearth of properly designed field-measured observations of soil water properties precludes an adequate assessment of their spatial variance structure and the development of an efficient field technology to optimize the size and spacing of their measurement. Sampling schemes based upon "aggle" statistics are relatively inexpensive owing to fewer observations required compared with those based upon geostatistical concepts. Each statistical analysis has advantages and disadvantages. A particular sampling scheme should embrace the attributes of both "aggle" statistics and geostatistics. Unfortunately, present-day soil mapping units have been developed without sufficient regard to quantitative evaluations of the spatial variances of soil parameters. Future research needs to answer the question if present-day mapping units manifest commensurate spatial variance structures for each of their soil properties. Within each mapping unit, when should sampling schemes be 1-, 2-, or 3-dimensional? Should observations be equally spaced or should they be clustered in some manner? For 2-dimensional sampling schemes, are orthogonal configurations more, or less informative than triangular, pentagonal, and other configurations? Is there a future for the turning bands concept in soil water studies? Do spectral and cospectral analyses of soil water properties offer opportunities for improved soil water management? Are there particular frequencies associated with cultivation or with pedologic processes that should be more amenable to cospectral analysis? What is the future of kriging and cokriging in soil water studies? Does cokriging offer any substantive advantage over simply taking more observations of that parameter of primary interest? How much better an understanding of transport of solutes, heat and gases through field soils can be gained through dimensional and inspectional analysis? Are scaling factors correlated with soil properties and are they linked within or between soil mapping units?

Of paramount importance is to identify criteria for choosing deterministic rather than stochastic algorithms for ascertaining the behavior of water in field soils. The presentations that follow as well as those we envision during the next decade will make substantive contributions toward that identification.

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Statistical opportunities for analyzing spatial and temporal heterogeneity of field soils

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Key words: coherency, geostatistics, kriging, state-space, stochastic, time-series analysis

Abstract

Statistical techniques for analyzing data in the agricultural sciences have traditionally followed the pioneering efforts of R.A. Fisher who assumed that observations in the field were independent and identically distributed. Such techniques, proven useful in the past and still being used today for comparing the merits of different management practices or different treatments, are presently giving way to additional methods that are based upon observations being spatially or temporally correlated. It is physically more sensible to expect soil attributes to be correlated when they are measured at adjacent points in space or time. Spatially repetitive patterns of soil attributes for physical and biological processes occurring at distances of a few molecules to those of kilometers are also expected. Opportunities to use geostatistics, time series analyses, state-space models, spectral analyses of variance, lagged regression models and other alternative techniques for analyzing multidimensional random fields are available to enhance the understanding of agro-ecosystems. Approaches to modeling and fitting data using stochastic partial differential equations and scaling techniques also help reveal the underlying processes occurring in field soils. Inclusion of these alternatives in the development of an agro-ecological technology leads to improved sampling designs to better entire management units, rather than ascertaining the impact of particular, sometimes arbitrary treatments applied to a set of small plots using analysis of variance methods.

Introduction

Although agricultural scientists have conducted experiments throughout the world at an accelerated pace during the last half century, they have been less than successful explaining and managing biological processes that occur at different scales of space and time. The soil chemist, for example, understands the kinetics of chemical reactions within a homogeneous soil sample but cannot usually extend the results to a larger, more heterogeneous rhizosphere. Likewise, the microbiologist cannot easily account for shifts in microbial populations within the rhizosphere when the crop canopy experiences nonuniform, larger scale microclimate perturbations. And many agronomists or plant scientists never "trust their luck" with only one year's data, but rely on fertilizer trials made repetitiously during a period of several years. Too often

management unit. The investigators have the belief that an analysis of variance of the result stemming from their semi-empirical treatments will allow the development of superior management alternatives for agricultural production for sites of similar characteristics. Such alternatives are indeed developed with whole fields being treated or managed the same, but seldom are quantitative observations made without resorting to the use of additional treated plots to assess the appropriateness or impact of the management alternative.

The above experimental scenarios stem from the pioneering work of R.A. Fisher. Recognizing that the properties and attributes of our natural resources are heterogeneous and vary from one location to another, he developed, in the absence of computers, the concept of analysis of variance which quantified the impacts of treatments within prescribed limits of probability. In a sense, he was able to ignore where soils vary from location to location. The benefits of his pioneering statistical efforts have been enormous, but today's complex environmental problems need to be alternatively approached using different, more global methods. We must be able to monitor fields or regions at particular locations and times to ascertain the behavior of our natural resources integrated over different scales of space and time and to be able to ascertain the value of a plant or soil attribute at unsampled locations within prescribed limits of accuracy and precision.

The objective here is to briefly introduce statistical concepts that require and take advantage of experimental observations manifesting spatial and/or temporal dependence and to provide a few examples of the many available methods having potential relevance to microbial ecology of arable lands. We give here a perhaps overly simplistic presentation of the theory and its potential practice. A secondary objective is to suggest the advantages of stochastic equations and models over the more common deterministic formulations normally used in agricultural and environmental sciences.

Because of the spatially continuous nature of the landscape, soil formation processes and climatic zones, it is not unreasonable to expect that soil measurements taken in close proximity to each other would be correlated. It is also reasonable to expect spatial or temporal variations in one or more ecological attributes to be related to varia-

tions of other parameters or properties indigenous to the landscape. Using statistical techniques that rely on observations made at different scales being spatially or temporally dependent allows greater opportunity to explore and understand biological and physical processes occurring in arable lands.

An approach for examining spatial structure is time series analysis, a methodology used to study temporally and spatially correlated data. Both time domain and frequency domain analysis are treated in time series methodologies (Shumway, 1988). Time domain analysis expresses the observation variable in terms of a linear function of its past values and a random error, and is used for estimating equally spaced univariate and multivariate processes. This analysis, known as autoregressive modeling (Box and Jenkins, 1970), may not easily be used to estimate missing observations. Another analysis, known as state-space modeling introduced by Kalman (1960) and Kalman and Bucy (1961), yields two linear equations called observation and state equations which describe the observed series. Two different kinds of errors are recognized. The first is the observation error which might be due to measurement error or disturbances observed in the data not related to the variable under study. The second is the state error caused by the noise inherent in the variable under study. Frequency domain analysis assigns the variations of the observation variable to combinations of sine and cosine waves. Spectral analysis examines the series to find those frequencies that contribute the most to the total variance assuming an expected value or mean for the observation variable.

An approach utilizing the spatial structure of a data set is kriging which is an estimation technique for regionalized variables that exhibit significant spatial autocorrelation (see, for example Ripley, 1981). The theory of regionalized variables was developed by Matheron (1963) and has been applied in the field of mining in estimation of ore reserves and more recently has been applied to agronomic and soil water properties (Vieira *et al.*, 1983). Applications of this theory are now popularly known as geostatistics. The kriging estimate is a weighted sum of the observed data in which the weights are determined by using the spatial distribution of the observations. In addition to kriging Journal and Huijbregts (1978) developed the geostatistical cokriging technique of estimating a

variable that is undersampled by considering its spatial correlation with other variables that are more densely sampled (David, 1977). Cokriging was used by Vieira *et al.* (1983) to estimate agronomic properties. Many other applications of regionalized variable analyses or geostatistics are available and continue to be developed. We only introduce them here and begin with the definition of spatial dependence.

Quantification of spatial dependence

Consider a set of observations $Z(x_1), Z(x_2), \dots, Z(x_n)$ at locations $x_1, x_2, \dots, x_n$ where each location defines a point in 1-, 2-, or 3-dimensional space. The correlogram $r(h)$ of the observations is estimated by

$$r(h) = \left[\frac{\sigma^2}{n(h) - 1} \right] \sum_{i=1}^{n(h)} [Z(x_i) - \bar{Z}] \times [Z(x_i + h) - \bar{Z}] \quad [1]$$

where $n(h)$ is the number of pairs of observation points a distance h apart, $\bar{Z}$ the mean and σ^2 the variance. The correlogram has possible values from -1 to 1. Some idealized correlograms are given in Figure 1. In Figure 1A the value of $r(0)$ is unity while $r(h)$ gradually decreases to zero as h increases. Figure 2A shows autocorrelation of a nematode population calculated from 625 data points separated by a distance of 6.5 m (Alemi *et al.*, in manuscript). For large separation distances no correlation or spatial dependence exists. For a set of observations in 1 dimension the integral scale l defined by

$$l = \int_0^\infty r(h) dh \quad [2]$$

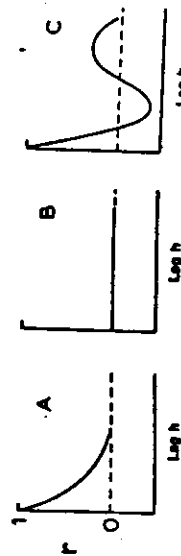


Fig. 1. Idealized correlograms. A is well-behaved and decreases monotonically to zero, B is for spatially independent or random observations, C is for observations that manifest repetitions, cyclic variations.

gives a measure of the distance h beyond which the pairs of observations are considered independent. The correlation length λ defined by

$$r = \exp(-h/\lambda) \quad [3]$$

gives still another measure beyond which the observations are considered independent. Numerically, the value of λ is that of h when r equals $1/e$.

Figure 1B is a correlogram manifesting a set of observations that are not correlated. In other words, for all separation distances, the observations are independent or may be said to be purely random. Figure 1C illustrates a set of observations whose values vary somewhat sinusoidally in space. The correlogram of an ideally sinusoidal set of observations would be a cosine curve having the same period as the observations and an amplitude of unity. In general, the correlogram shows fluctuations with the same kinds of periods as the original observations, except that all the fluctuations have been put in phase so that they reach a maximum value at $h = 0$. Figure 2B shows correlograms of water flow and its salt concentration measured monthly in a stream during a 6-year period. The autocorrelation functions indicate that both variables have strong cyclic behavior with a period of 12 lags (a yearly cycle).

Another measure of spatial dependence is given by the variogram $\gamma(h)$ estimated by

$$\gamma(h) = \frac{1}{2n(h)} \sum_{i=1}^{n(h)} [Z(x_i + h) - Z(x_i)]^2 \quad [4]$$

Figure 3 provides an illustrative example of the several potential features of a variogram. Although $\gamma(0) = 0$, $\gamma(h)$ will appear to remain nonzero in many cases as h approaches zero as shown by the value C_0 commonly called the nugget. The value $h = a$, called the range, is the maximum separation distance for which pairs of observations remain

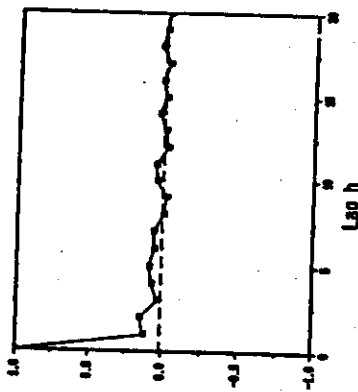


Fig. 2A. Autocorrelation function for observations of *Penstemon edorus* populations.

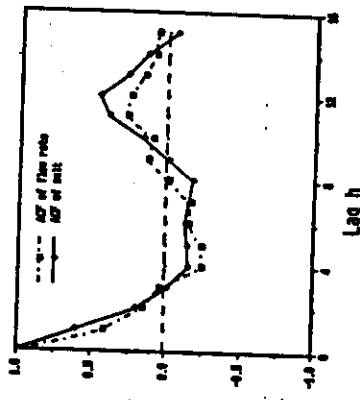


Fig. 2B. Autocorrelation functions of water flow and salt concentration.

correlated. For greater separation distances γ remains equal to C_0 and is called, the sill. Some variograms do not manifest one or more of the features C_0 , C or a . The use of the variogram provides an opportunity for interpolation and mapping schemes, for identifying new observation locations to improve estimates for a total region or subregion, and to ascertain the impact of the size of observation on the scale of the field problem being studied as well as to develop efficient sampling strategies.

We assume that the reader is relatively unfamiliar

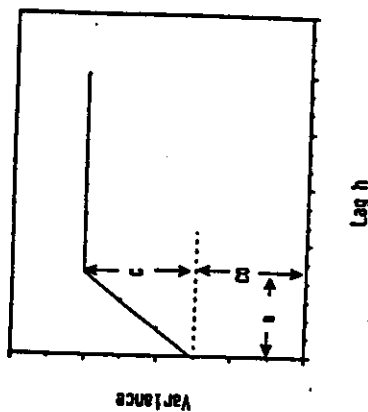


Fig. 3. Idealized variogram with range a , sill $(C + C_0)$ and nugget C_0 .

familiar with the concepts of time series and geostatistical analysis of data. Hence in the following section these methods are described briefly with real world examples.

Autoregressive integrated moving average (ARIMA) models

A stationary process Z_s is said to be autoregressive of order p , $AR(p)$ if

$$Z_s = \sum_{i=1}^p \phi_i Z_{s-i} + \omega_s \quad s = 1, 2, \dots \quad [5]$$

where ω_s is a white noise (independent and identically distributed normal variable) and ϕ_i are coefficients to be estimated. The moving average of order q is defined as

$$Z_s = \omega_s - \sum_{i=1}^q \theta_i \omega_{s-i} \quad s = 1, 2, \dots \quad [6]$$

where ω_s is again white noise and θ_i are coefficients to be estimated. The mixed autoregressive process of order p and moving average of order q , ARMA (p, q) is defined as

$$Z_s = \sum_{i=1}^p \phi_i Z_{s-i} - \sum_{i=1}^q \theta_i \omega_{s-i} + \omega_s \quad [7]$$

ARMA modeling can be applied to a nonstationary series, if the nonstationarity is removed for

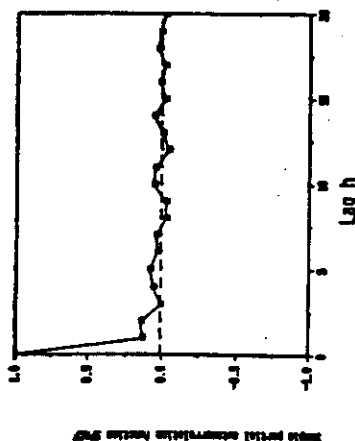


Fig. 4. Sample partial autocorrelation of *Paratrichodorus* series.

example by differencing the original series. If a difference of order d is applied, and an autoregressive model of order p and moving average of order q is used, the model is called autoregressive integrated moving average of orders p , d , and q or ARIMA (p, d, q).

The first step in ARIMA modeling is model identification. Autocorrelation and partial autocorrelation functions are the main tools in detection of the nonstationarity and model identification. For a complete description of the procedures for model identification and testing the adequacy of the model see Shumway (1988). The sample partial autocorrelation (SPAC) of a series of nematode

(*Paratrichodorus*) populations given in Figure 4 indicates that the SPAC is significant at lags one and two. Thus an AR(2) seems to be the appropriate model. The ACF of residuals (differences between the observed and estimated values in the series) is given in Figure 5 with no significant peak (ACF less than 1.96(S))²¹ where S is the length of the series). Hence, the AR(2) is an adequate model to describe the nematode series and can be written as

$$Z_s = 0.1068 Z_{s-1} + 0.1324 Z_{s-2} + \omega_s \quad [8]$$

We note that the nematode population at location s can be described in terms of its magnitude at the two previous locations ($s-1$) and ($s-2$).

State-space models

State-space models have been used to describe data in many different scientific endeavors. The multivariate version of the model is defined by the observation equation

$$Z_s = A_s \xi_s + v_s, \quad s = 1, 2, 3, \dots, S \quad [9]$$

where the p -dimensional vector $Z_s = (Z_{s1}, \dots, Z_{sp})$ is represented as a linear combination of unobserved q -dimensional vector $\xi_s = (\xi_{s1}, \dots, \xi_{sq})$, using a $p \times q$ design matrix A_s . The additive noise vector v_s is a $p \times 1$ vector of measurement error related to the observation vector Z_s . It is assumed that v_s are zero mean independent random vectors with covariance matrix R . The unobserved or so called state vector ξ_s satisfies the state equation

$$\xi_s = B \xi_{s-1} + \omega_s, \quad s = 1, 2, \dots, S \quad [10]$$

where B is a $q \times q$ transition matrix and ω_s is the state noise vector independent of v_s with mean zero and covariance matrix Q . The initial value of ξ_1 is assumed to be ξ_0 with mean vector μ and covariance matrix Σ .

In state-space modeling μ , Σ , B , Q , and R are estimated from the observed series Z_s , $s = 1, 2, \dots, S$ by an iterative procedure using Kalman filtering and expectation maximization (EM) algorithm given in Shumway and Stoffer (1982) and Shumway (1988). An example of application of a state-space model is given in Figure 6 where a series of cotton yield observations and that of a nematode population were used to estimate the cotton yield

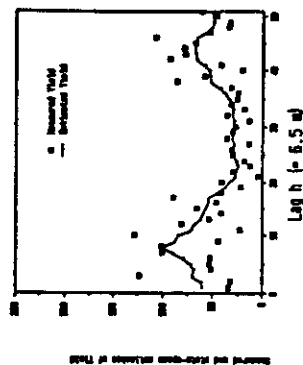


Fig. 6. Measured and state-space estimates of cotton yield.

series. The observation equation is

$$Z_s = A_s \xi_s + v_s \quad [11]$$

The state equation for the bivariate process using *Paratrichodorus* (P) and cotton yield (Y) series is

$$\begin{bmatrix} P \\ Y \end{bmatrix}_s = \begin{bmatrix} 0.92 & 0.23 \\ 0.01 & 0.94 \end{bmatrix} \begin{bmatrix} P \\ Y \end{bmatrix}_{s-1} + \begin{bmatrix} \omega_p \\ \omega_y \end{bmatrix} \quad [12]$$

with variance σ equal to 683 and 165, and those of ω equal to 4277 and 944, respectively. Using the standard deviation of variances, the magnitude of measurement and process errors can be calculated. The solid line given in the figure stems from equation [12] and tends to "smooth" the cotton yield data on the basis that there is a spatial correlation between cotton yield and nematode infestation.

Our example has involved only two observation variables — cotton yield and number of nematodes. State-space modeling easily handles several variables, does not require long series of observations, and is an excellent method to identify which local field properties or attributes tend to be related to the behavior of an observation variable of primary interest. Noting that the solid line in Figure 6 does not indicate a constant yield of cotton across the field implies that additional properties or attributes affecting crop yield are not accounted for in the bivariate state space model. The use of a multivariate model would potentially allow the identification of other properties or attributes that are responsible for the manner in which the cotton yield varies across the field.

Spectral analysis

Arable land is often cultivated with animals or machines in repeatable, cyclic patterns during one or more growing seasons. In addition, irrigations, fertilizers, agrochemicals, and crop residues are usually applied on or into the soil in cyclic patterns having various periodicities or frequencies depending upon the spacing between plant rows or the method of cultivation or irrigation. Under such conditions observations $Z(x_s)$ made in the field will yield a correlogram $r(h)$ that can be used to identify those portions of the spatial variance structure that are associated with sinusoidal patterns of specific periods or frequencies. The identification is made from a spectral analysis by calculating the power spectra from the correlogram in one or more directions. For one direction, the power spectra $f(v)$ is

$$f(v) = 2 \int_0^{\infty} r(h) \cos(2\pi v h) dh \quad [13]$$

where v is the frequency of the periodic observations. A graph of $f(v)$ reveals relative maxima at those values of v for which the observations manifest a sinusoidal behavior.

Soil surface temperature observations of a fallow, 6-ha field measured with an infrared thermometer on a 36 by 64 regular grid 2 days after a 2 cm rain are examined here to illustrate the use of spectral analysis (Bazza, 1985). Figures 7 and 8 show 3-dimensional plots of all 2304 observations viewed from the east and north, respectively. The most striking feature of the data is that variability is not the same across the field. An obvious feature is that variability in the E-W direction is completely stationary as compared with that in the other directions. The correlogram of the observations for the E-W direction indicated spatial independence while that for the N-S direction revealed a periodic behavior. The graph of $f(v)$ for the N-S direction (Fig. 9) shows a relative maximum of a frequency of about 0.06 corresponding to a period of about 18 lags. Figure 10 of average values of temperature of all 36 columns shows this periodic behavior about a mean of 6.6°C along the N-S direction. Also given in the figure is a graph of the salinity of the water used to irrigate the field for cotton production during the previous summer. Notice that there appears to be an inverse relation between the soil tem-

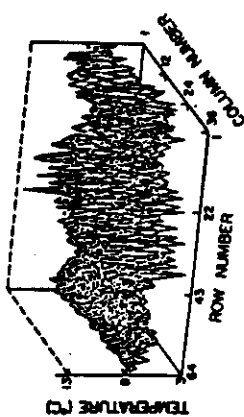


Fig. 7. Soil surface temperature observations viewed from the east.

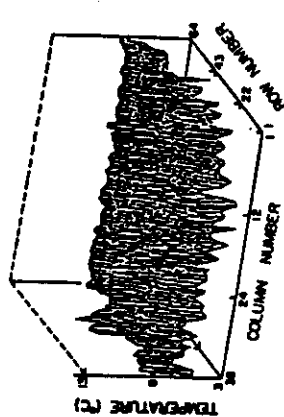


Fig. 8. Soil surface temperature observations viewed from the north.

perature and the salinity level of the irrigation water. The cross spectrum for two series Y_t and X_t can be written as

$$f_{xy}(v) = C_{xy}(v) - i q_{xy}(v) \quad [14]$$

where $f_{xy}(v)$ is cross spectrum between two series, $C_{xy}(v)$ and $q_{xy}(v)$ are called cospectrum and quad-

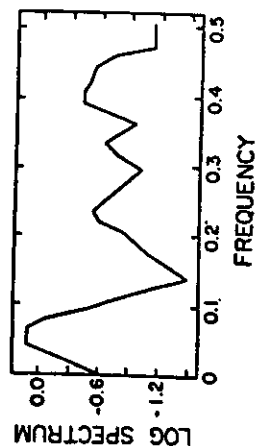


Fig. 9. Spectrum of surface soil temperature.

spectrum, respectively and i is the imaginary number. The cross-spectrum is the frequency dependent measure of the covariance between two series Y_t and X_t . The dependence between the two series Y_t and X_t with frequency can be explained by the coherence function between the two series. The coherence, defined by

$$\gamma_{xy}^2(v) = |f_{xy}(v)|^2 / f_x(v) f_y(v) \quad [15]$$

is the frequency dependent correlation between two series. If Y_t is an exact linear version of X_t then coherence becomes equal to unity at all frequencies.

After first calculating the cross-correlogram between the temperature versus salinity data already given as average values in Figure 10 in a manner analogous to equation (1) only for two variables. The cross-spectrum $f_{xy}(v)$ comes from equation [14]. The strong negative peak shown in Figure 11 indicates that the erratic small differences in temperature values across the field and salinity are spatially related at a frequency of about 0.06. At that frequency the magnitude of the coherence was nearly unity and indicates that the two series, temperature and salinity, were almost perfectly related. The reason behind the cyclic temperature behavior is probably due to the salinity causing lower soil water potential, lower evaporation and reduced hydraulic conductivity tending to prevent redistribution of soil water.

The concept of time stability of a spatially measured series as defined by Vachaud *et al.* (1985) has recently been expanded to include general linear transformations in time and to account for the occurrence of spatial scale dependency (Kachanoski and de Jong, 1988). Vachaud defined time dependency as the time invariant association between spatial location and classical statistical parametric

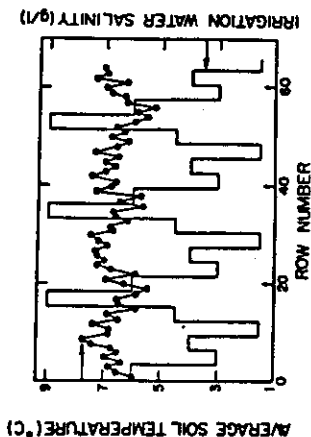


Fig. 10. Row averages of surface soil temperature and salt content of irrigation water.

values. For example, a given location in a field will tend to have a soil water content that remains in the same ranked order with those of other locations. In other words, a location that is relatively dry at one time compared to others will remain relatively dry at other times while a wet location will remain wet.

Let D represent the spatial domain of interest, j represent a horizontal spatial location vector in D , and $S_j(t)$ represent the cumulative soil water storage to depth L at location j and time t . If $s_j(t)$ is the realization of the variable $S_j(t)$ at location j and time t , the relative deviation from the expected or mean soil water storage $E[S_j(t)]$ at the same time across the spatial domain D is given by (Vachaud *et al.*, 1985):

$$\delta_j(t) = \frac{S_j(t) - E[S_j(t)]}{E[S_j(t)]} \quad [16]$$

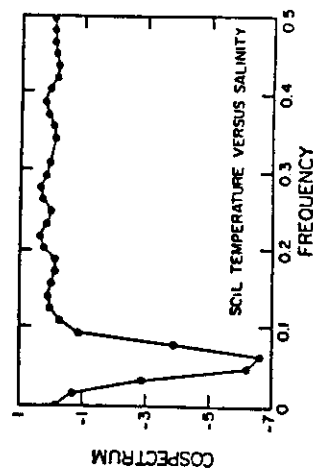


Fig. 11. Cross spectrum of surface soil temperature and salt content of irrigation water.

According to Vachaud *et al.*, a small temporal variation in $\delta_j(t)$ (i.e. time independence) is an indication of time stability, thus

$$\delta_j(t) \approx \delta_j(t) \quad [17]$$

where t_1 and t_2 represent two different measurement times. Substitution of equation [16] into equation [17] and taking the variables $S_j(t)$ and $S_j(t)$ in lieu of the observations $s_j(t)$ and $s_j(t)$ respectively, gives:

$$S_j(t) \approx \frac{E[S_j(t)]}{E[S_j(t)]} S_j(t) \quad [18]$$

Thus time stability as defined in equation [18] implies a linear relationship between soil water storage at time t_1 and t_2 across all spatial locations.

Since spatial variation is often scale dependent, it follows that time stability may also be scale dependent. For example, localized surface runoff may significantly alter the spatial variation of soil water storage on a small scale, but the changes may be insignificant compared to large scale variations. Other factors, such as climatic variations may affect time stability at large scales.

The persistence of a spatial pattern at different spatial scales was examined by Kachanoski and de Jong (1988) by expanding on the concept that time stability requires a linear relationship between successive time samplings. The presence of a significant linear relationship at different spatial scales is tested using spatial coherence analysis introduced by equation [15]. In this case, the coherence function estimates the proportion of the spatial variance of $S_j(t)$ which can be explained by the

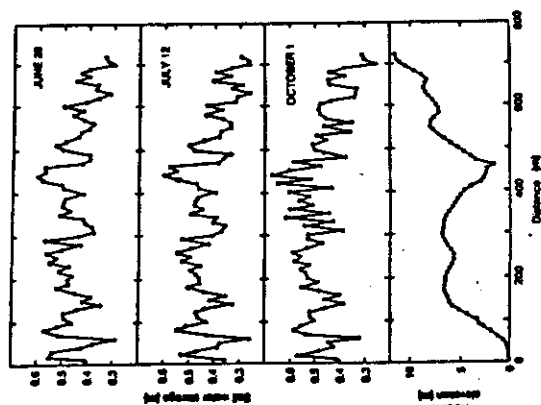


Fig. 12. Measured soil water storage and relative elevation for the study site.

spatial variance of $S_v(\cdot)$, as a function of spatial scale. Thus, the coherency function is a direct measurement of time stability as a function of spatial scale. A more detailed discussion has been given by Kachanoski *et al.* (1983a) which in turn is a summary of the direct Fourier method of spectrum estimation (Brillinger, 1981).

The soil of the study area of Kachanoski and de Jong was a moderately fine textured, moderately calcareous Typic Haploboroll resting on a gently rolling to undulating topography. Neutron probe access tubes were installed in a fallow field every 10 m in a 720 m long transect in a north-south direction. Soil water measurements were taken at all locations with depth on June 28, July 12 and October 1. Elevation measurements were taken at each location and 10 m east and west of each location using a rod and surveying level. The grid of elevation readings around each soil water sampling location was used to estimate local surface curvature by a least squares surface analysis. No significant precipitation was recorded during the first two sampling dates, (June 28-July 12), thus this period is characterized by drying. The

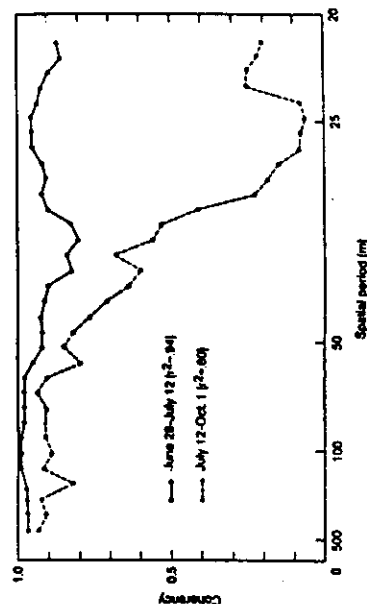


Fig. 13. Coherency spectra for the drying (June 28-July 12) and recharge (July 12-Oct. 1) periods.

able carbon and nutrient sources, using coherency analyses should be able to infer spatial scales of hydrologic microbial and crop community processes.

Geostatistics

An example of the variogram given in Figure 3 for the nematode populations measured by Alemi *et al.* (1988) on a 6.5 m grid across an agricultural field is given in Figure 14. The values of the range, sill and nugget are 32.5 m (5 lags), 8500 and 7000, respectively. We may use this variogram to kriging or

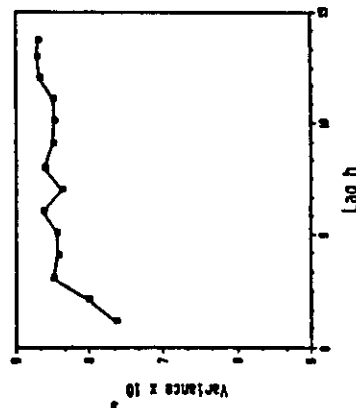


Fig. 14. Average semivariogram of *Pseudocerosus*.

estimate the nematode population at all locations across the agricultural field. The variogram can be used for estimating nematode values Z^* at locations x_0 where measurements have not been made. The procedures lead not only to the estimated values Z^* , but also an estimate of $\text{var}(Z - Z^*)$ which indicates the reliability of the estimate. "Error maps" can be prepared to reveal areas within the field where the variance is highest and presumably where additional sampling should be made.

In the kriging procedure, an estimate Z^* is assumed to be a linear function of the known values Z already measured at N other locations

$$Z^*(x_0) = \sum_{i=1}^N \lambda_i Z(x_i) \quad [19]$$

by choosing the weight factors λ_i such that the expected value and variance of $[Z^*(x_0) - Z(x_0)]$ are zero and a minimum, respectively. The variogram showing the spatial variance structure of Z is used to estimate the values of λ_i . Cokriging is a similar procedure that takes advantage of the spatial structure of the covariance between two variables such as cotton yield and nematode. Still another procedure, block kriging, allows interpolation between averages of values centered at locations instead of individual observations for each location. Universal kriging provides for combination of a deterministic trend across the field to coexist with random variations while disjunctive kriging can be used to identify areas within a field having values expected to be above or below a

threshold magnitude. Hence, there are many attractive kriging and cokriging analyses that can be used to develop and evaluate alternative sampling grids or locations relative to a specific management system.

The excellent comprehensive reference of Journel and Huijbregts (1978) on geostatistics and that of Warnick *et al.* (1986) applied to soil science provide essential details and conditions omitted from the above introductory presentation. Many additional statistical opportunities exist when one recognizes that spatial dependence also depends upon direction and that more than one variable can be treated simultaneously.

Stochastic equations

It is not possible to model a system relative to what is "a typical response to management practices" or what is "a typical pattern in ecology" without understanding the spatial and temporal variations for both small and large scales. "Models are being constructed, refined, elaborated, tinkered with and displayed with little or no effort to link them to the real world," Pielou (1981). The standard approach to field scale or plot scale problems has been to construct analytical or numerical solutions of the classical differential equations, assuming that the parameters take on known values which can be easily evaluated by a few observations. Although there have been major advances in numerical modeling techniques, standard modeling approaches are giving way to stochastic methods which incorporate the effects of natural or anthropogenically-induced variability. Stochastic models are more realistic representations of field situations than deterministic models and they provide better predictions owing to the probabilistic nature of the data incorporated. Because stochastic models can be examined in terms of the variability of the model parameters, data collection can be improved to emphasize those parameters having the greatest effects on the results.

During the past decade, stochastic analyses of groundwater (Baker *et al.*, 1978) and unsaturated flow (e.g., Yeh *et al.*, 1985) have been accelerating in number and kind of applications. Nielsen *et al.* (1987) recently reviewed conceptual alternatives for modeling water and solute transport in the vadose

zone. Promising areas for further research in microbial ecology of arable lands await similar development.

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