

# JACK KNIFING FOR SEMIVARIOGRAM VALIDATION <sup>(1)</sup>

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## ABSTRACT

The semivariogram function fitting is the most important aspect of geostatistics and because of this the model chosen must be validated. Jack knifing may be one the most efficient ways for this validation purpose. The objective of this study was to show the use of the jack knifing technique to validate geostatistical hypothesis and semivariogram models. For that purpose, topographical heights data obtained from six distinct field scales and sampling densities were analyzed. Because the topographical data showed very strong trend for all fields as it was verified by the absence of a sill in the experimental semivariograms, the trend was removed with a trend surface fitted by minimum square deviation. Semivariogram models were fitted with different techniques and the results of the jack knifing with them were compared. The jack knifing parameters analyzed were the intercept, slope and correlation coefficient between measured and estimated values, and the mean and variance of the errors calculated by the difference between measured and estimated values, divided by the square root of the estimation variances. The ideal numbers of neighbors used in each estimation was also studied using the jack knifing procedure. The jack knifing results were useful in the judgment of the adequate models fitted independent of the scale and sampling densities. It was concluded that the manual fitted semivariogram models produced better jack knifing parameters because the user has the freedom to choose a better fit in distinct regions of the semivariogram.

**Key words:** Semivariograms, stationarity, topography, scale of variation.

## RESUMO

### JACK KNIFING PARA VALIDAÇÃO DE SEMIVARIOGRAMAS

A função de ajuste do semivariograma é o aspecto mais importante da geoestatística, por esse motivo, o modelo escolhido deve ser validado. *Jack knifing* pode ser um dos métodos mais eficientes para esta finalidade. O objetivo deste estudo foi mostrar o uso da técnica *jack knifing* para validar modelos de hipótese geoestatística e de semivariograma. Para essa finalidade, foram analisados dados topográficos de seis campos de escalas distintas e de diferentes densidades de amostragem. Devido aos dados topográficos selecionados terem apresentado tendência em todos os campos, fato verificado pela ausência de patamar no semivariograma, a tendência foi removida através do ajuste de uma superfície de tendência pelo método dos quadrados mínimos. Os modelos de semivariograma foram ajustados por diferentes técnicas e validados pela comparação dos resultados do *jack knifing*. Os parâmetros de *Jack knifing* analisados foram a interseção, o coeficiente angular e o coeficiente de correlação para a regressão linear entre valores estimados e medidos, e a média e a variância dos erros calculados pela diferença entre valores medidos e estimados divididos pela raiz quadrada da variância da estimativa. Os resultados do *jack knifing* foram úteis no julgamento adequado dos modelos ajustados aos semivariogramas, independentemente da escala e da densidade de amostragem. Conclui-se que os modelos de semivariograma ajustados manualmente resultaram em melhores parâmetros de *jack knifing* devido à liberdade de escolha, por parte do usuário, do melhor ajuste em regiões distintas do semivariograma.

**Palavras-chave:** erro reduzido, estacionaridade, topografia, variação de escala.

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## 1. INTRODUCTION

Soil variability has always existed and if not taken into account when field work is involved, there is a risk to make wrong conclusions out of the data. If the soil variability is somewhat organized in the space and spatial dependence can be determined, the data must be analyzed using geostatistics. In this condition, it is possible to estimate values for the unsampled locations without bias and with minimum variance through the kriging interpolation technique. On the other hand, in order to make appropriate use of geostatistics, it is necessary assume that the measured data correspond to one realization of a continuous random function which exists in every point in the field (VIEIRA et al., 1983). For this reason, it is necessary that the data fit into some stationarity hypothesis (VIEIRA, 2000). Besides, the experimental semivariogram calculated will be a series of discrete data pairs of distances and semivariances to which a continuous mathematical function must be fitted. For this reason, it is commonly said that semivariogram function fitting is the most important aspect of geostatistics (McBRATNEY and WEBSTER, 1986) and the model chosen must be validated. Methods for fitting a model to the semivariogram are well documented in the literature (WOLLENHAUPT et al., 1997; GOTWAY, 1991; LEE, 1994; CARVALHO and VIEIRA, 2004). Jack knifing may be one the most efficient ways for this validation purpose (VIEIRA et al., 1983). Through this technique, a measured value is temporarily taken out of the data set, and then it is estimated using the semivariogram model fitted. Simultaneously with the estimated value, the estimation variance is also calculated in the kriging procedure. This procedure is successively repeated for every measured value and at the end it is possible to calculate errors using measured, estimated and estimation variance values whose values must remain within some specified statistical limits. This technique is also known as a cross validation or "leave-one-out" and has been used in some different applications (SCHECHTMAN and WANG, 2004; FORTES et al., 2004; ZHU et al., 2008, MELLO et al., 2008, PIRES and STRIEDER, 2006). The objective of this study was to show the use of the jack knifing technique to validate geostatistical hypothesis and semivariogram models.

## 2. MATERIAL AND METHODS

### Data

The following data sets used were: 1) A square field of 90m on each side named FIELD 1 sampled on a 2m grid with a total of 2500 points; 2) A triangular field named FIELD 2 of 110m by 220m, sampled on a 10m square grid, with a total of 164 points; 3) An approximately rectangular field named FIELD 3 measuring 90x250m, sampled on trapezoidal grid of 5m, with a total of 383 points; 4) An approximately rectangular field named FIELD 4, measuring 120x160m sampled on a 10m square grid with 302 data points; 5) An approximately rectangular field named FIELD 5, of approximately 35ha, sampled on a 50m square grid in a total of 146 data points; 6) A circular field named FIELD 6 of 77ha sampled on a square grid of 50m, at every one of the 322 points.

A summary about the six fields with grid information is shown in table 1. Topography was chosen as a data set to be analyzed because its form is easily verified in the field and its surface does not change with time allowing for the field validation of the results if necessary. These specific fields were chosen because they represent a very wide range of scales (from 0.81 to 77 hectares), of grid sampling spacing (from 2 to 50 m), of number of values (from 2500 to 146 values) and consequently of number of samples per hectare (from 3086.42 to 4.17). Because the topographical data showed very strong trend for all fields as it was verified by the absence of a sill in the experimental semivariograms, the trend was removed with a trend surface fitted by the minimum square deviation, according to VIEIRA (2000). The degree of the trend surface used in each case is listed in the last column of table 1.

### Detrending

The trend removal technique used is described in VIEIRA (2000). The presence of a trend is detected when the semivariogram does not have a stable sill. This condition violates the intrinsic hypothesis as it represents a field for which the mean value depends on the spatial position. The simplest trend removal consists in fitting a three dimensional surface to the data

**Table 1.** Area, grid, number, number of sample (N ha<sup>-1</sup>) and trend removed to sampling sites field 1 until 6

Site	Area	Grid	Number	Nha <sup>-1</sup>	Trend
	Ha	m			
Field 1	0.81	2	2500	3086.42	Parabolic
Field 2	1.21	10	164	135.54	Parabolic
Field 3	2.25	5	383	170.22	Cubic
Field 4	3.50	10	302	100.67	Parabolic
Field 5	35.00	50	146	4.17	Parabolic
Field 6	77.00	50	322	4.18	Linear

by the least squares and subtracting its values from the originals. For a parabolic trend surface, the equation is

$$Z^*(x, y) = A_0 + A_1X + A_2Y + A_3X^2 + A_4XY + A_5Y^2 \quad (1)$$

where  $Z(x, y)$  is the estimated trend surface,  $X$  and  $Y$  are the coordinate positions and  $A_0, A_1, A_2, A_3, A_4$  and  $A_5$  are the regression parameters determined by the least squares method. This surface is then subtracted from the originals generating a new variable which may be called Residuals.

$$Z_{res}(x, y) = Z(x, y) - Z^*(x, y) \quad (2)$$

The criteria for the choice of the degree for the detrending surface is the simplest surface that will produce a semivariogram with a stable sill. Thus, if a linear surface solves the stationarity problem producing a semivariogram with a sill, there no need to look for any other degree of a surface.

### Semivariogram

The semivariogram is, by definition:

$$\gamma(h) = 1/2 E\{Z(x_i) - Z(x_i + h)\}^2 \quad (3)$$

And can be estimated by:

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

where  $N(h)$  is the number of pairs of measured values  $Z(x_i), Z(x_i + h)$ , separated by a vector  $h$  (JOURNAL and HUIJBREGTS, 1978). The graph of  $\gamma^*(h)$  versus the corresponding values of  $h$ , called semivariogram, is a function of the vector  $h$ , and therefore it depends on both magnitude and direction of  $h$ . When the semivariogram is the same for all directions it is called isotropic. Many variables show anisotropic semivariograms depending on the dimensions of the field and of the nature of the variability. There are ways of transforming an anisotropic semivariogram (JOURNAL and HUIJBREGTS, 1978; BURGESS and WEBSTER, 1980) in order to reflect the variability in different directions. Jack knifing procedure can also be used to verify the distance range over which a semivariogram can be used before anisotropic effects may affect the results (VIEIRA, 2000).

### Models

Experimental semivariograms contain a set of discrete data points of distance and semivariance. A model must be fit to the experimental data with the objective of having semivariances available for every distance needed (GOTWAY, 1991). In order to be used to properly describe the spatial variability of any variable,

one of the requirements on the model is that the function used must be conditional positive definite (MCBRATNEY and WEBSTER, 1986). This condition will guarantee that the variances calculated will be positive. The main models which satisfy that condition and are adequate for use in geostatistical calculations are the spherical, the exponential and the Gaussian. On the equations bellow,  $C_0, C_1$ , and  $a$  represent the nugget effect, the structural variance and the range, respectively.

For the spherical model, usually symbolized as  $Sph(C_0, C_1, a)$ , the equation is:

$$\gamma(h) = C_0 + C_1 \left( 1.5 \frac{h}{a} - 0.5 \left( \frac{h}{a} \right)^3 \right), \quad h \leq a \quad (5)$$

$$\gamma(h) = C_0 + C_1, \quad h > a$$

The exponential model, symbolized as  $Exp(C_0, C_1, a)$ , the equation is:

$$\gamma(h) = C_0 + C_1 \left( 1 - \text{EXP} \left( \frac{-3h}{a} \right) \right) \quad (6)$$

The gaussian model, symbolized as  $Gau(C_0, C_1, a)$ , the equation is:

$$\gamma(h) = C_0 + C_1 \left( 1 - \text{EXP} \left( -3 \left( \frac{h}{a} \right)^2 \right) \right) \quad (7)$$

With the parameters fitted to the semivariogram the dependence ratio ( $DR$ ) can be calculated (ZIMBACK, 2001)

$$DR = \frac{C_1}{(C_0 + C_1)} \times 100 \quad (8)$$

The dependence ratio ( $DR$ ) represents the proportion of the semivariance which is structured. The smallest the  $DR$  value the weakest is the spatial dependence.

### Jack knifing

Estimating the semivariogram and associated parameters (nugget effect, range and sill) from a set of field measurements using the traditional estimator is a difficult task (SHAFER and VARLIEN, 1990). In order to verify if the semivariogram model adequately describes the spatial variability, there is a validation technique commonly known as jack knifing. This is a process of estimating known values by temporarily taking them out of the data set. The value taken out of the data set is estimated using the semivariogram model fitted, and a series of neighborhood sizes, generating an estimated value and an estimation variance. The process is then repeated for each measured value and at the end, there will be a set of measured,

estimated and estimation variances through which it is possible to calculate error parameters whose values must be within some statistically known limits. More details about the process can be found in JOURNAL and HUIJBREGTS (1978). There are some reports in the literature with some applications of the cross validation technique but using only the one-to-one graph of measured versus estimated values (FORTES et al., 2004; MELLO et al., 2005; MELLO et al., 2008). In this paper, six different jack knifing parameters are examined as criteria for judging the performance of semivariogram models, neighborhood of estimation and geostatistical hypothesis. All the estimations were made using ordinary kriging interpolation technique.

### The one-to-one Measured vs Estimated values

Using the  $N$  measured values,  $Z(x_i)$ , and the  $N$  values estimated through the jack knifing procedure,  $Z^*(x_i)$ , it is possible to make the graph known as the one-to-one and to calculate the linear regression between measured and estimated values. The regression will be:

$$Z^*(x_i) = a + b Z(x_i) \quad (9)$$

Where  $a$  is the intercept,  $b$  is the slope and  $r^2$  is the coefficient of determination between  $Z^*(x_i)$  and  $Z(x_i)$ .

Thus, if the estimation,  $Z^*(x_i)$ , is identical to the measured,  $Z(x_i)$ , for every one of the  $N$  points, then  $a$  is zero (0),  $b$  and  $r^2$  are equal to one (1.0), and the graph of  $Z(x_i)$  vs  $Z^*(x_i)$  would be a series of points on the one-to-one line. As the value of  $a$  depart from zero (0) to positive values, it is an indication that the estimator  $Z^*(x_i)$  is over estimating small values of  $Z(x_i)$  and under estimating large values. As the value of  $a$  gets negative the reverse happens. This way, the quality of the estimation may be assessed judging these parameters. The examination of the one-to-one scatter plot of measured versus estimated values is an important aid in judging the estimation performance but it only makes sense for the best selection of the other parameters (VIEIRA et al., 1983). Therefore, it is an useful technique but it needs the other parameters before a decision on neighborhood size and semivariogram parameters can be made.

### The reduced error

Remembering that through the kriging estimation of values,  $Z^*(x_i)$ , there is always a value of the estimation variance,  $\sigma_k^2(x_i)$ , corresponding to the uncertainty of the estimation process (VIEIRA et al., 1983), then it is possible to define the reduced error as:

$$RE(x_i) = \frac{Z^*(x_i) - Z(x_i)}{\sigma_k(x_i)} \quad (10)$$

The division by the square root of the estimation variance causes the reduced error,  $RE(x_i)$ ,

to be dimensionless which is a convenient situation for comparison between different variables.

The unbiasedness condition of kriging estimation requires that:

$$\overline{RE} = E \{RE(x_i)\} = E \{(Z^*(x_i) - Z(x_i)) / \sigma_k(x_i)\} = 0 \quad (11)$$

The minimum variance condition requires that:

$$VAR(\overline{RE}) = E \{(Z^*(x_i) - Z(x_i)) / \sigma_k(x_i)\}^2 = 1 \quad (12)$$

These properties make this kind of error assessment a very valuable and easy to use tool for validation of geostatistical procedures. Because these errors have fixed reference values of 0 (zero) and 1 (one), respectively, and are dimensionless, their judgment and interpretation is much easier and allows comparison with other variables expressed in different units.

### The root mean square error (RMSE)

Another very powerful parameter of the jack knifing technique is the RMSE which can be calculated using

$$RMSE = \frac{1}{N} \sqrt{\sum_{i=1}^N [Z^*(x_i) - Z(x_i)]^2} \quad (13)$$

The disadvantage of this kind of error is that it does not have any standard to be compared with. Therefore the best results of the jack knifing technique will be obtained when the RMSE is minimum.

## 3. RESULTS AND DISCUSSION

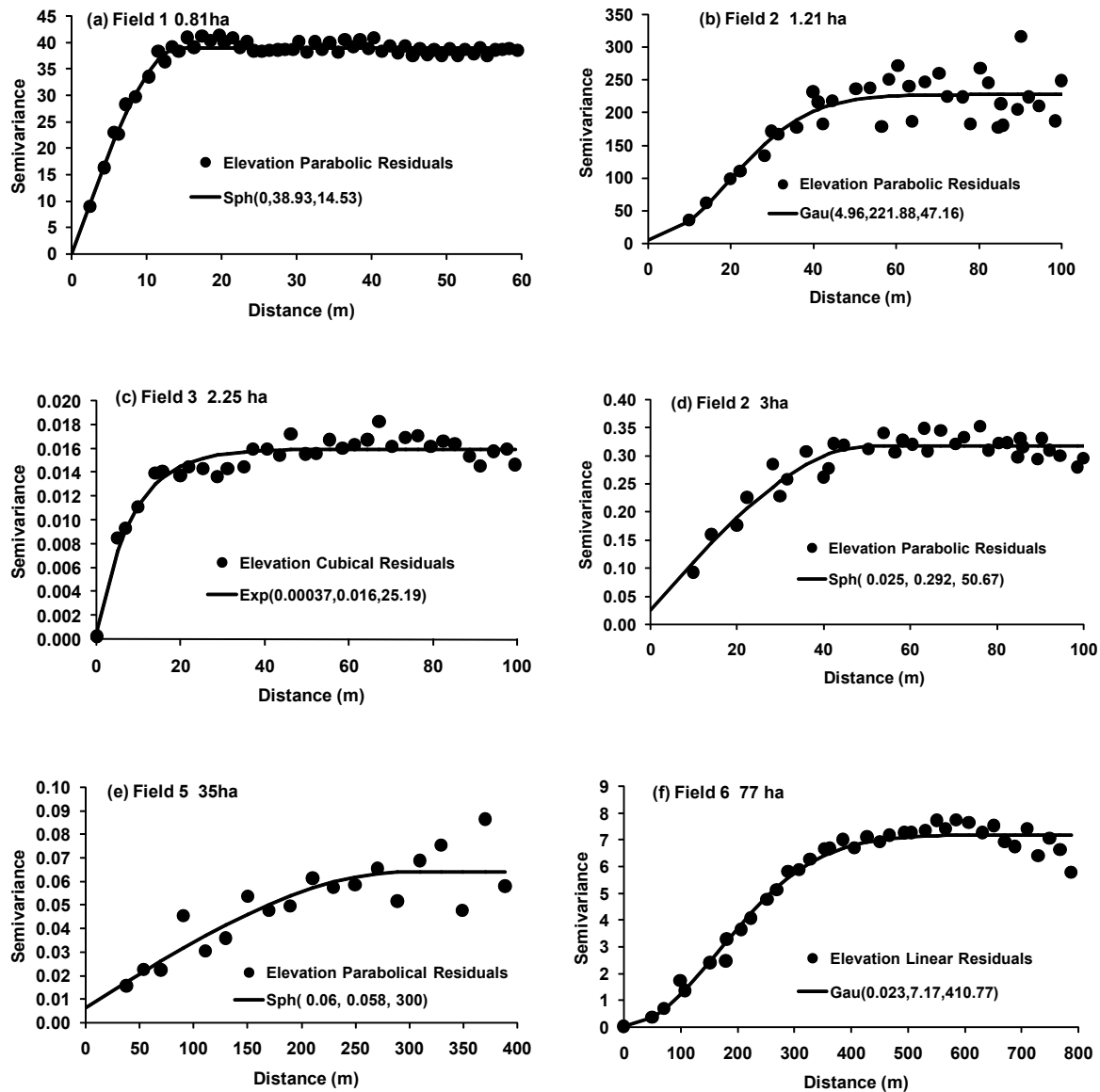
A summary about the data and the places from where they were sampled is shown in table 1, where it can be seen that the areas sampled range from 0.81 ha to 77 ha, and the sampling densities range from 3086 to 4.18 samples ha<sup>-1</sup>. Therefore, the areas sampled represent a very large range of field dimensions and topographic conditions, and for these reasons we hope that the results are adequate to evaluate the performance of the proposed method of validation. All the data sets presented very strong trends which had to be removed in order to satisfy the intrinsic hypothesis. The trend was removed by fitting with least square method a tri dimensional surface and subtracting it from the original data as described for VIEIRA et al. (2002). Last column in Table 1 identifies the kind of trend surface used to remove the trend of each of the data sets. The criteria used in the de trending process is to use the surface with the smallest degree that does the job of removing the trend. Thus, if the linear surface produces residuals whose semivariogram has a clear sill then there no need to try the parabolic surface because

the linear already did it. As shown in the last column of table 1, from the six fields studied, for one a cubic trend surface was used, for four the parabolic surface and for one the linear surface.

The parameters for the models fitted to the semivariograms are shown in Table 2 and the corresponding graphs of the semivariograms are shown in Figure 1 with the models fitted. It can be seen that

**Table 2.** Parameters of the models fitted to the semivariograms, nugget effect ( $C_0$ ), structural variance ( $C_1$ ), range ( $a$ ), coefficient of determination ( $r^2$ ), sum of squares deviation weighed (SQDP) and dependence ratio (DR%) to sampling sites field 1 until 6

Variable	Model	$C_0$	$C_1$	$a$	$r^2$	SQDP	DR (%)
Field 1	Spherical	0.00	38.93	14.53	0.9682	0.9553	100.00
Field 2	Gaussian	4.96	221.88	47.16	0.7400	727.68	97.81
Field 3	Exponential	0.00	0.02	25.19	0.9244	7.63E-07	97.66
Field 4	Spherical	0.02	0.29	50.67	0.8672	3.37E-04	92.13
Field 5	Spherical	0.01	0.06	300.00	0.7571	8.15E-05	90.38
Field 6	Gaussian	0.02	7.17	410.77	0.9723	0.0911	99.68



**Figure 1.** Semivariograms for topographic elevations in different sites (field 1 until 6). Sph stands for Spherical, Exp stands for exponential, Gau stands Gaussian. The numbers in parenthesis correspond to the parameters  $C_0$ ,  $C_1$  and  $a$  defined in equations (5), (6) and (7).

all semivariograms fit the intrinsic hypothesis (they all have a very well defined sill) and that the worst fitting found was for field 2, with  $r^2 = 0.7400$ . Otherwise, in general, the models fit quite well the experimental semivariograms. The dependence ratio shown in the last column of Table 2 indicates the very high degree of continuity that the residuals for the topographical data has. From semivariograms for the six fields, three of them were fitted to the spherical model, two to Gaussian and one to exponential model. Notice that the  $r^2$  values for the models fitted to the semivariograms (Table 2) are the lowest for field 2 and field 5 caused by the dispersion of values around the sill. Not much

importance should be given to this fact as the main portion of the semivariogram is the short distance (McBRATNEY AND WEBSTER, 1986). On the other hand, all semivariograms are very well fitted to their respective models for the distances smaller than the range.

The results of jack knifing for the five parameters ( $a$ ,  $b$ , and  $r^2$  for the regression 1:1, mean and variance of the reduced errors) used is shown in figure 2. The intercept values (Figure 2a) indicate that the semivariograms for all six surfaces produced good regression between measured and estimated values, as all of them, except for field 2 with four neighbors, approach a zero intercept. The slope

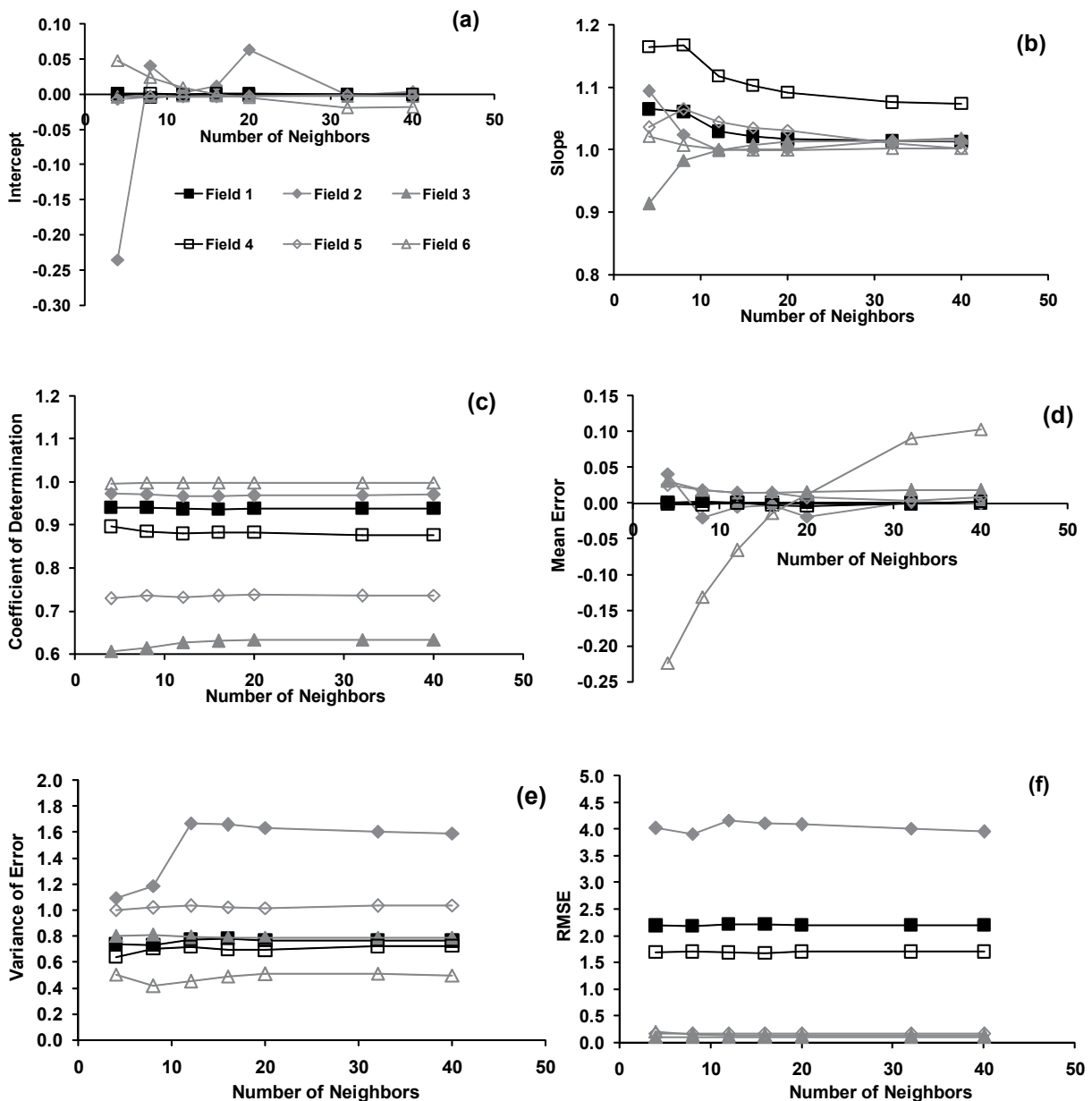


Figure 2. Results of jack knifing [intercept, slope, correlation coefficient, mean error, variance and root mean square error (RMSE)] for the six data fields.

of the regression line between measured and estimated values (Figure 2b) approaches the ideal situation ( $a=1$ ) for any neighborhood above 16 neighbors for all data sets. Field 4 was the only one for which this parameter was separated from the others and the cause for this has not been identified. The coefficients of determination (Figure 2c) showed a wide spread of values. In general, the values for this parameter approach the ideal value of 1.0, except for fields 3 and 5 which presented values around 0.7. The mean error (Figure 2d), except for field 6, are grouped around the ideal value of 0 (zero). The reason for the departure of the mean value for the reduced errors for field 6 is not known at this point. However, for 16 neighbors, all fields have a value of mean error very close to 0 (zero). The variance of the reduced errors (Figure 2e) should ideally approach the value of 1.0. In general all fields present values below 1.0 for the variance of the reduced errors, except for field 2 which presented values much above that level. Figure 2f shows a graph of the Root Mean Square Error (RMSE) between the measured and the estimated values. This parameter, although very robust, its values are somewhat arbitrary because it does not have any standards or ideal value to be compared with. An overall examination of the values of all parameters together shows that most of them approach the ideal values if a neighborhood of 16 values is used as it has been indicated by VIEIRA et al. (1983). The square grid sampling may be the explanation for this apparent coincidence around 16 neighbors for all of the data sets.

In order to investigate further the jackknifing potential for the validation of semivariogram models, four different models were fitted to the semivariogram for field 5 by different methods. The parameters fitted are shown in table 3 and on graph on figure 3. The four models were named Solver, Wrong, Sill 1 and Sill 2. The model Solver was fitted using the Solver technique in Excel to maximize the coefficient of determination. The

model called Wrong was purposely fitted with a wrong nugget effect value. The models Sill 1 and Sill 2 were fitted by trial and error by manually placing the sill value in different positions in order to provide information about the effect of the proper choice of the sill value on the jack knifing parameters. The results from the jack knifing for these models are shown in figure 4. The results indicate that if one single model had to be chosen it should be the model identified as Sill 2 with 16 neighbors as all the jack knifing parameters approach the ideal values with this choice. It can be clearly seen that the model identified as "wrong" had a very poor performance for all the jack knifing parameters. The above discussion illustrates the idea of using the jack knifing technique for fine tuning the semivariogram model fitting.

Because the data from field 1 was the one with the highest number (2500) and it was also the smallest field (highest density of samples, see Table 1) one set of jack knifing was calculated for this data set with 20 neighbors. The jack knifing parameters are shown in table 4. Except for the variance of the errors, all other parameters are very close to the ideal values. A graph with the measured versus estimated values for this calculation is plotted in figure 5 where it can be seen that there was a very good agreement between measured and estimated values.

#### 4. CONCLUSION

The technique shown in this work allowed to conclude that jack knifing may be a very helpful aid in the choice of the parameter models for the semivariogram. It is also possible to use jack knifing procedure for fine tuning the parameters fitted by running a sensitivity analysis with the jack knifing parameters. The jack knifing procedure was proven to discriminate very well between a representative model of the variability and a model which is not correct. The jack knifing procedure

**Table 3.** Parameter of models fitted (solver, wrong, sill1 and sill2) for jack knifing comparison, nugget effect ( $C_0$ ), structural variance ( $C_1$ ), range (a) and sum of squares deviation weighed (SQDP)

Model	$C_0$	$C_1$	a	SQDP
Solver	0.025	0.292	50.67	0.000337
Wrong	0.200	0.050	50.67	0.090506
Sill 1	0.040	0.300	50.67	0.000796
Sill 2	0.000	0.294	50.67	0.000934

**Table 4.** Jack knifing parameters for Field 1, using 20 neighbors

Parameter	Result
Number of neighbors	20
Intercept	0.0014
Slope	0.8661
r	0.9384
Mean error	0.0006
Variance	0.7661
Distance	9

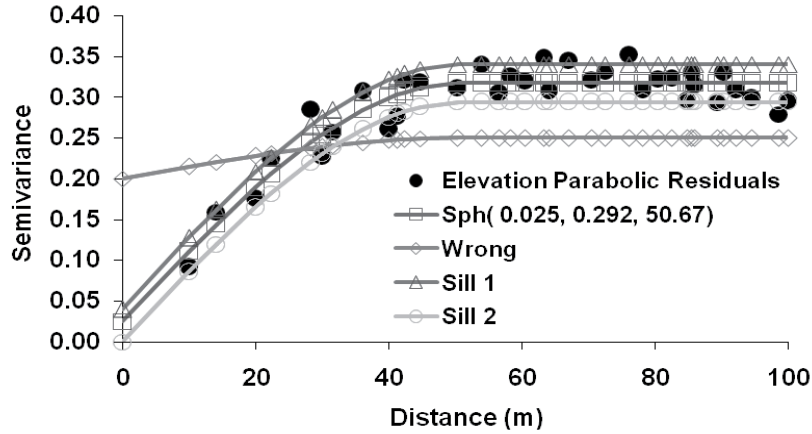


Figure 3. Models fitted to semivariogram of topographic data from FIELD 2.

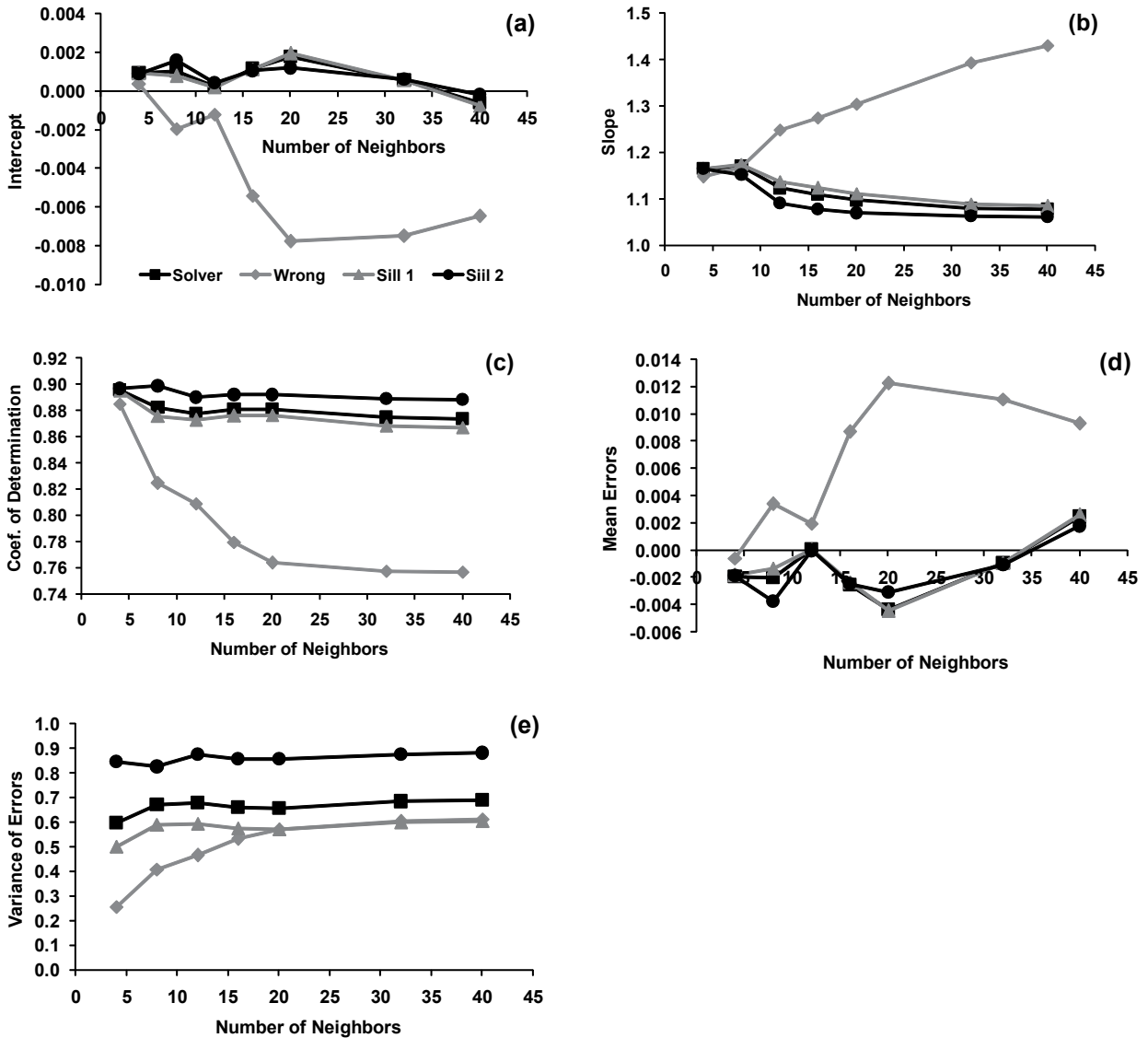
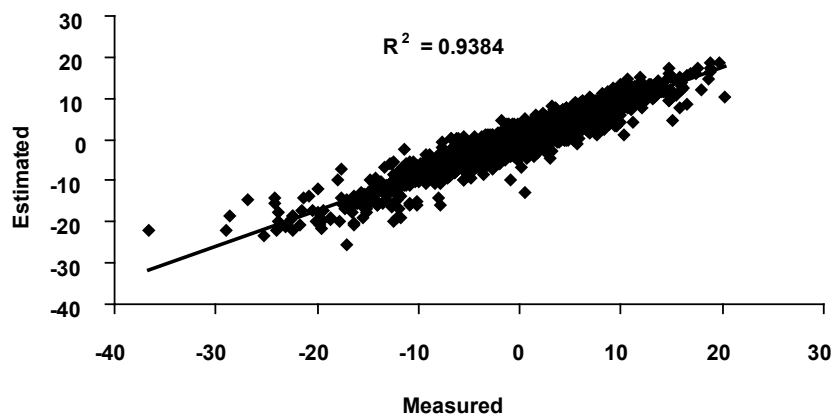


Figure 4. Results of jack knifing [intercept, slope, correlation coefficient, mean error, variance and root mean square error (RMSE)] corresponding to residuals of parabolic trend of topographical heights for field 2.





**Figure 5.** One-to-one graph for measured versus estimated values of elevation parabolic residuals for field 1 with jack knifing using 20 neighbors.

proposed in this paper also helps in identifying the best neighborhood size for the kriging interpolation. It does not seem possible to pick one single jack knifing parameter for this analysis as a judgment of all parameters seems to be a more secure decision tool.

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