# The meuse data set: a tutorial for the **gstat** R package

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#### 1 Introduction

The meuse data set is a data set comprising of four heavy metals measured in the top soil in a flood plain along the river Meuse. The governing process seems that polluted sediment is carried by the river, and mostly deposited close to the river bank. This document shows a geostatistical analysis of this data set.

This tutorial introduced the functionality of the R package gstat, used in conjunction with package sp. Package gstat provides a wide range of univariable and multivariable geostatistical modelling, prediction and simulation functions, where package sp provides general purpose classes and methods for defining, importing/exporting and visualizing spatial data.

#### 2 R geostatistics packages

Package gstat (Pebesma, 2004) is an R package that provides basic functionality for univariable and multivariable geostatistical analysis, including

- variogram modelling, residual variogram modelling, and cross variogram modelling using fitting of parametric models to sample variograms
- $\bullet\,$  geometric anisotropy specfied for each partial variogram model
- restricted maximum likelihood fitting of partial sills
- variogram and cross variogram maps
- simple, ordinary, universal and external drift (co)kriging
- (sequential) Gaussian (co)simulation equivalents for each of the kriging varieties
- indicator (co)kriging and sequential indicator (co)simulation
- kriging in a local or global neighbourhood
- block (co)kriging or simulation for each of the varieties, for rectangular or irregular blocks

Other geostatistical packages for R usually lack part of these options (e.g. block kriging, local kriging, or cokriging) but provide others: e.g. package geoR and geoRglm (by Paulo Ribeiro and Ole Christensen) provide the model-based geostatistics framework described in Diggle et al. (1998), package fields (Doug Nychka and others) provides thin plate spline interpolation, covariance functions for spherical coordinates (unprojected data), and routines for network design optimization.

#### 3 Spatial data frames

Package gstat assumes that data are projected, i.e. they should not be provided as lattitude/longitude. As an example, we will look at the meuse data set, which is a regular data frame that comes with package gstat:

```
> library(gstat)
Loading required package: sp
> data(meuse)
> class(meuse)
[1] "data.frame"
> names(meuse)
               "y"
 [1] "x"
                          "cadmium" "copper"
                                               "lead"
                                                         "zinc"
                                                                    "elev"
 [8] "dist"
               "om"
                          "ffreq"
                                    "soil"
                                               "lime"
                                                         "landuse" "dist.m"
> coordinates(meuse) = ~x + y
> class(meuse)
[1] "SpatialPointsDataFrame"
attr(,"package")
[1] "sp"
> summary(meuse)
Object of class SpatialPointsDataFrame
Coordinates:
     min
            max
x 178605 181390
y 329714 333611
Is projected: NA
proj4string : [NA]
Number of points: 155
Data attributes:
    cadmium
                       copper
                                         lead
                                                          zinc
 Min.
        : 0.200
                  Min.
                         : 14.00
                                    Min.
                                           : 37.0
                                                     Min.
                                                            : 113.0
 1st Qu.: 0.800
                  1st Qu.: 23.00
                                    1st Qu.: 72.5
                                                     1st Qu.: 198.0
 Median : 2.100
                                    Median :123.0
                                                     Median : 326.0
                  Median : 31.00
 Mean
       : 3.246
                          : 40.32
                                    Mean :153.4
                  Mean
                                                     Mean
                                                            : 469.7
```

```
3rd Qu.: 3.850 3rd Qu.: 49.50 3rd Qu.:207.0 3rd Qu.: 674.5 Max. :18.100 Max. :128.00 Max. :654.0 Max. :1839.0
```

 elev
 dist
 om
 ffreq
 soil
 lime

 Min. : 5.180
 Min. : 0.00000
 Min. : 1.000
 1:84
 1:97
 0:111

 1st Qu.: 7.546
 1st Qu.: 0.07569
 1st Qu.: 5.300
 2:48
 2:46
 1: 44

 Median : 8.180
 Median : 0.21184
 Median : 6.900
 3:23
 3:12

 Mean : 8.165
 Mean : 0.24002
 Mean : 7.478

3rd Qu.: 8.955 3rd Qu.:0.36407 3rd Qu.: 9.000 Max. :10.520 Max. :0.88039 Max. :17.000

NA's : 2.000

landuse dist.m

W :50 Min. : 10.0

Ah :39 1st Qu.: 80.0

Am :22 Median : 270.0

Fw :10 Mean : 290.3

Ab : 8 3rd Qu.: 450.0

(Other):25 Max. :1000.0

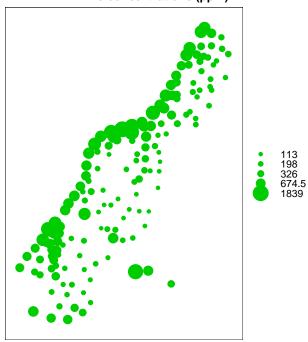
NA's : 1

> coordinates(meuse)[1:5, ]

x y

- [1,] 181072 333611
- [2,] 181025 333558
- [3,] 181165 333537
- [4,] 181298 333484
- [5,] 181307 333330
- > bubble(meuse, "zinc", col = c("#00ff0088", "#00ff0088"), main = "zinc concentrations (pp

#### zinc concentrations (ppm)



and note the following:

- the function coordinates, when assigned (i.e. on the left-hand side of an
   = or <- sign), promotes the data.frame meuse into a SpatialPoints DataFrame, which knows about its spatial coordinates; coordinates may
   be specified by a formula, a character vector, or a numeric matrix or data
   frame with the actual coordinates</li>
- 2. the function coordinates, when not assigned, *retrieves* the spatial coordinates from a SpatialPointsDataFrame.
- 3. the two plotting functions used, plot and bubble assume that the x- and y-axis are the spatial coordinates, and choose the aspect ratio of the axes such that one unit in x equals one unit in y (i.e., data are map data, projected).

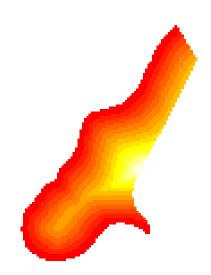
# 4 Spatial data on a regular grid

- > data(meuse.grid)
- > summary(meuse.grid)

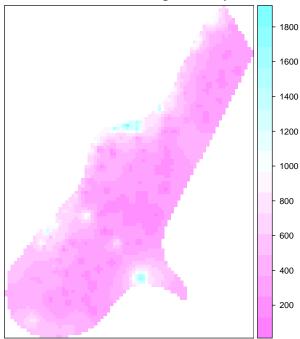
X	У	part.a	part.b	
Min. :178460	Min. :329620	Min. :0.0000	Min. :0.0000	
1st Qu.:179420	1st Qu.:330460	1st Qu.:0.0000	1st Qu.:0.0000	
Median :179980	Median :331220	Median :0.0000	Median :1.0000	
Mean :179985	Mean :331348	Mean :0.3986	Mean :0.6014	
3rd Qu.:180580	3rd Qu.:332140	3rd Qu.:1.0000	3rd Qu.:1.0000	

```
Max. :181540 Max. :333740 Max. :1.0000
                                                 Max. :1.0000
                soil
    dist
                               ffreq
Min. :0.0000 Min. :1.000
                               1: 779
1st Qu.:0.1193 1st Qu.:1.000
                                2:1335
Median :0.2715 Median :1.000
                               3: 989
Mean :0.2971 Mean :1.578
3rd Qu.:0.4402 3rd Qu.:2.000
Max.
      :0.9926 Max. :3.000
> class(meuse.grid)
[1] "data.frame"
> coordinates(meuse.grid) = ~x + y
> class(meuse.grid)
[1] "SpatialPointsDataFrame"
attr(,"package")
[1] "sp"
> gridded(meuse.grid) = TRUE
> class(meuse.grid)
[1] "SpatialPixelsDataFrame"
attr(,"package")
[1] "sp"
> image(meuse.grid["dist"])
> title("distance to river (red = 0)")
> zinc.idw = krige(zinc ~ 1, meuse, meuse.grid)
[inverse distance weighted interpolation]
> class(zinc.idw)
[1] "SpatialPixelsDataFrame"
attr(,"package")
[1] "sp"
> spplot(zinc.idw["var1.pred"], main = "zinc inverse distance weighted interpolations")
```

#### distance to river (red = 0)



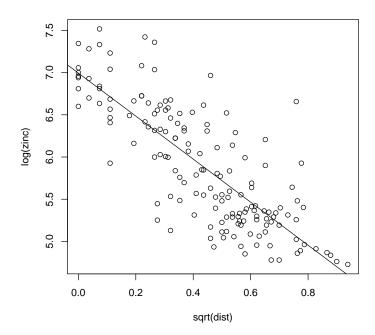
#### zinc inverse distance weighted interpolations



If you compare the bubble plot of zinc measurements with the map with distances to the river, it becomes evident that the larger concentrations are measured at locations close to the river. This relationship can be linearized by

log-transforming the zinc concentrations, and taking the square root of distance to the river:  $\,$ 

```
> plot(log(zinc) ~ sqrt(dist), meuse)
> abline(lm(log(zinc) ~ sqrt(dist), meuse))
```



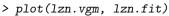
### 5 Variograms

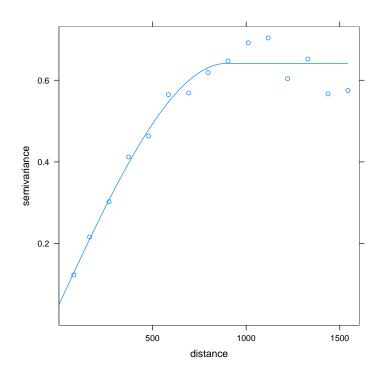
Variograms are calculated using the function variogram, which takes a formula as its first argument: log(zinc)~1 means that we assume a constant trend for the variable log(zinc).

```
> lzn.vgm = variogram(log(zinc) ~ 1, meuse)
> lzn.vgm
```

	np	dist	gamma	dir.hor	dir.ver	id
1	57	79.29244	0.1234479	0	0	var1
2	299	163.97367	0.2162185	0	0	var1
3	419	267.36483	0.3027859	0	0	var1
4	457	372.73542	0.4121448	0	0	var1
5	547	478.47670	0.4634128	0	0	var1
6	533	585.34058	0.5646933	0	0	var1
7	574	693.14526	0.5689683	0	0	var1
8	564	796.18365	0.6186769	0	0	var1
9	589	903.14650	0.6471479	0	0	var1
10	543	1011.29177	0.6915705	0	0	var1

```
11 500 1117.86235 0.7033984
                                  0
                                           0 var1
12 477 1221.32810 0.6038770
                                           0 var1
                                  0
13 452 1329.16407 0.6517158
                                  0
                                           0 var1
14 457 1437.25620 0.5665318
                                  0
                                           0 var1
15 415 1543.20248 0.5748227
                                           0 var1
> lzn.fit = fit.variogram(lzn.vgm, model = vgm(1, "Sph", 900, 1))
> lzn.fit
 model
             psill
                      range
   Nug 0.05066243
                     0.0000
   Sph 0.59060780 897.0209
```

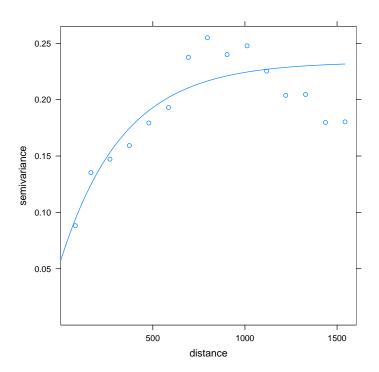




Instead of the constant mean, denoted by ~1, we can specify a mean function, e.g. using ~sqrt(dist) as a predictor variable:

```
> lznr.vgm = variogram(log(zinc) ~ sqrt(dist), meuse)
> lznr.fit = fit.variogram(lznr.vgm, model = vgm(1, "Exp", 300,
+ 1))
> lznr.fit

model    psill    range
1    Nug  0.05712231    0.0000
2    Exp  0.17641559    340.3201
> plot(lznr.vgm, lznr.fit)
```

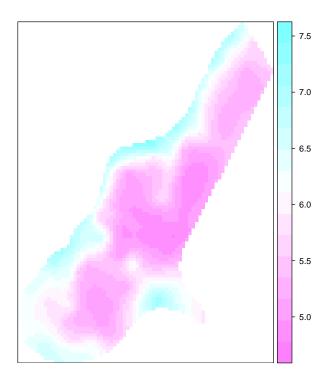


In this case, the variogram of residuals with respect to a fitted mean function are shown. Residuals were calculated using ordinary least squares.

## 6 Kriging

```
> lzn.kriged = krige(log(zinc) ~ 1, meuse, meuse.grid, model = lzn.fit)
[using ordinary kriging]
```

> spplot(lzn.kriged["var1.pred"])

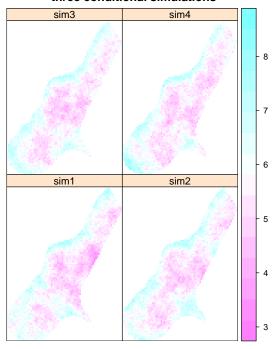


### 7 Conditional simulation

```
> lzn.condsim = krige(log(zinc) ~ 1, meuse, meuse.grid, model = lzn.fit,
+ nmax = 30, nsim = 4)

drawing 4 GLS realisations of beta...
[using conditional Gaussian simulation]
> spplot(lzn.condsim, main = "three conditional simulations")
```

#### three conditional simulations



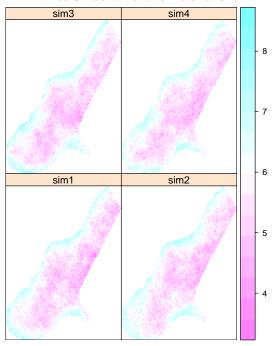
For UK/residuals:

```
> lzn.condsim2 = krige(log(zinc) ~ sqrt(dist), meuse, meuse.grid,
+ model = lznr.fit, nmax = 30, nsim = 4)
```

drawing 4 GLS realisations of beta...
[using conditional Gaussian simulation]

> spplot(lzn.condsim2, main = "three UK conditional simulations")

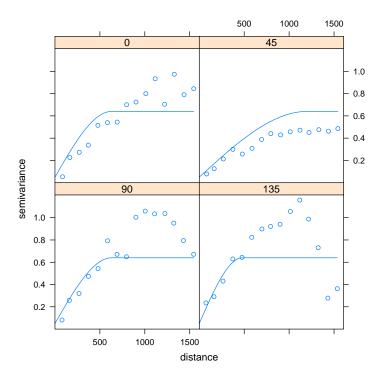




## 8 Directional variograms

The following command calculates a directional sample variogram, where directions are binned by direction angle alone. For two point pairs, Z(s) and Z(s+h), the separation vector is h, and it has a direction. Here, we will classify directions into four direction intervals:

```
> lzn.dir = variogram(log(zinc) ~ 1, meuse, alpha = c(0, 45, 90,
+ 135))
> lzndir.fit = vgm(0.59, "Sph", 1200, 0.05, anis = c(45, 0.4))
> plot(lzn.dir, lzndir.fit, as.table = TRUE)
```



Looking at directions between 180 and 360 degrees will repeat the image shown above, because the variogram is a symmetric measure:  $(Z(s) - Z(s + h))^2 = (Z(s + h) - Z(s))^2$ .

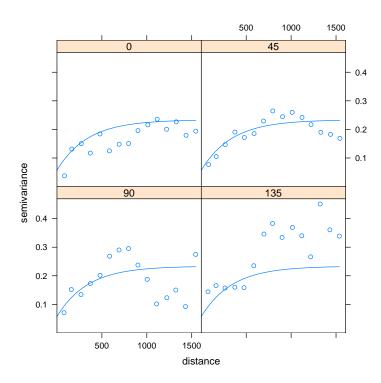
The first plot gives the variogram in the zero direction, which is North; 90 degrees is East. By default, point pairs are assigned to the directional variorgram panel with their nearest direction, so North contains everything between -22.5 and 22.5 degrees (North-West to North-East). After classifying by direction, point pairs are binned by separation distance class, as is done in the usual omnidirectional case.

In the figure, the partial sill, nugget and model type of the model are equal to those of the omnidirectional model fitted above; the range is that in the direction with the largest range  $(45^o)$ , and the anisotropy ratio, the range in the 135 direction and the range in the 45 direction, estimated "by eye" by comparing the 45 and 135 degrees sample variograms. Gstat does not fit anisotropy parameters automatically.

We do not claim that the model fitted here is "best" in some way; in order to get to a better model we may want to look at more directions, other directions (e.g. try alpha = c(22, 67, 112, 157)), and to variogram maps (see below). More elaborate approaches may use directions in three dimentions, and want to furnter control the direction tolerance (which may be set such that direction intervals overlap).

For the residual variogram from the linear regression model using sqrt(dist) as covariate, the directional dependence is much less obvious; the fitted model here is the fitted isotropic model (equal in all directions).

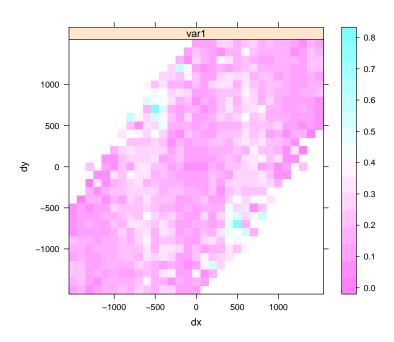
```
> lznr.dir = variogram(log(zinc) \sim sqrt(dist), meuse, alpha = c(0, + 45, 90, 135))
```



## 9 Variogram maps

Another means of looking at directional dependence in semivariograms is obtained by looking at variogram maps. Instead of classifying point pairs Z(s) and Z(s+h) by direction and distance class separately, we can classify them jointly. If  $h=\{x,y\}$  be the two-dimentional coordinates of the separation vector, in the variogram map the semivariance contribution of each point pair  $(Z(s)-Z(s+h))^2$  is attributed to the grid cell in which h lies. The map is centered around (0,0), as h is geographical distance rather than geographical location. Cutoff and width correspond to some extent to map extent and cell size; the semivariance map is point symmetric around (0,0), as  $\gamma(h)=\gamma(-h)$ .

```
> vgm.map = variogram(log(zinc) ~ sqrt(dist), meuse, cutoff = 1500,
+ width = 100, map = TRUE)
> plot(vgm.map, threshold = 5)
```



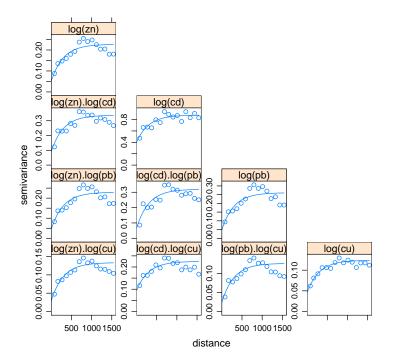
The threshold assures that only semivariogram map values based on at least 5 point pairs are shown, removing too noisy estimation.

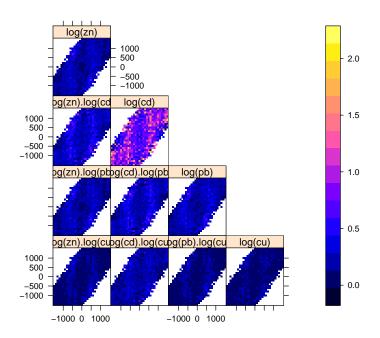
#### 10 Cross variography

Fitting a linear model of coregionalization.

```
> g = gstat(NULL, "log(zn)", log(zinc) ~ sqrt(dist), meuse)
> g = gstat(g, "log(cd)", log(cadmium) ~ sqrt(dist), meuse)
> g = gstat(g, "log(pb)", log(lead) ~ sqrt(dist), meuse)
> g = gstat(g, "log(cu)", log(copper) ~ sqrt(dist), meuse)
> v = variogram(g)
> g = gstat(g, model = vgm(1, "Exp", 300, 1), fill.all = TRUE)
> g.fit = fit.lmc(v, g)
> g.fit
log(zn) : formula = log(zinc) ~ sqrt(dist) ; data dim = 155 x 12
log(cd) : formula = log(cadmium) ~~ sqrt(dist) ; data dim = 155 x 12
log(pb) : formula = log(lead) `~`sqrt(dist) ; data dim = 155 x 12
log(cu) : formula = log(copper) `~ `sqrt(dist) ; data dim = 155 x 12
variograms:
                   model
                              psill range
log(zn)[1]
                     Nug 0.05141798
                     Exp 0.17556219
                                       300
log(zn)[2]
                     Nug 0.39996573
log(cd)[1]
log(cd)[2]
                     Exp 0.47893816
                                       300
```

```
Nug 0.04770893
log(pb)[1]
                                         0
log(pb)[2]
                     Exp 0.21323027
                                       300
log(cu)[1]
                     Nug 0.04577523
log(cu)[2]
                     Exp 0.07827374
                                       300
log(zn).log(cd)[1]
                     Nug 0.09190848
                                         0
log(zn).log(cd)[2]
                                       300
                     Exp 0.24542024
log(zn).log(pb)[1]
                     Nug 0.04528367
                                         0
log(zn).log(pb)[2]
                     Exp 0.18407011
                                       300
log(cd).log(pb)[1]
                     Nug 0.06425412
                                         0
                                       300
log(cd).log(pb)[2]
                     Exp 0.25525359
log(zn).log(cu)[1]
                     Nug 0.02912806
                                         0
log(zn).log(cu)[2]
                     Exp 0.10438748
                                       300
log(cd).log(cu)[1]
                     Nug 0.09441635
                                         0
log(cd).log(cu)[2]
                      Exp 0.13073936
                                       300
log(pb).log(cu)[1]
                     Nug 0.02369778
                                         0
                                       300
log(pb).log(cu)[2]
                      Exp 0.10267516
> plot(v, g.fit)
> vgm.map = variogram(g, cutoff = 1500, width = 100, map = TRUE)
 plot(vgm.map, threshold = 5, col.regions = bpy.colors(), xlab = "",
      ylab = "")
```





### References

- Diggle, P.J., J.A. Tawn, R.A. Moyeed, 1998. Model-based geostatistics. Applied Statistics 47(3), pp 299-350.
- Pebesma, E.J., 2004. Multivariable geostatistics in S: the gstat package. Computers & Geosciences 30: 683-691.
- Wackernagel, H., 1998. Multivariate Geostatistics; an introduction with applications, 2<sup>nd</sup> edn., Springer, Berlin, 291 pp.