# Spatio-temporal geostatistics using gstat



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## 1 Introduction

Since gstat package version 1.0-0, a dependency of gstat on the R package spacetime was introduced, allowing the code in gstat to exploit spatio-temporal data structures from that package. This vignette describes the possibilities and limitations of the package for spatio-temporal geostatistics.

To understand some of the possibilities and limitations, some knowledge of the history of the software is needed. The original gstat software (Pebesma and Wesseling, 1998) was a standalone computer program written in around 25,000 lines of C code, and would do geostatistical modelling, prediction and simulation. The gstat R package (Pebesma, 2004) consisted mostly of an R interface to this C code, together with convenience functions to use R's modelling interface (formula's, see ?lm) and graphic capabilities (trellis graphics in package lattice to show cross variogram as matrix plots; interaction with variogram clouds using base plots).

Starting 2003, a group of programmers developed a set of classes and methods for dealing with spatial data in R (points, lines, polygons, grids), which was supported by the publications of the well-known ASDAR book (Bivand et al. 2008; see also <a href="http://www.asdar-book.org/">http://www.asdar-book.org/</a>) and helped convergence in the user community, with in 2011 over 2000 subscribers on the r-sig-geo mailing list. Package gstat was one of the first packages that adopted and benefited from these classes.

To realize a particular idea, writing code in C typically takes about 10-20 times as long as writing it in R. C code can be more efficient, gives more control over memory usage, but is also more error prone—mistakes in C code make an R session crash, something that is hard to do when writing R code.

The original C code of gstat (Pebesma and Wesseling, 1998) provides all kriging varieties (universal, ordinary, simple; univariable, or multivariable as in cokriging) for two- or three-dimensional data. When the spatial domain is constrained to two dimensions (and this might cover over 99% of the use cases!), the third dimension might be used to represent time. As such, the metric variogram model, which allows for geometric anisotropy definition in three dimensions, can be used for spatio-temporal kriging. When defining the three-dimensional variogram as the sum of 2 or more nested variogram (summed)

models, one can choose anisotropy coefficients for a single model such that this model is *effectively* zero in some directions, e.g. in space *or* in time; this allows one to approximate the so-called space-time sum model. It should be noted that at the C code there is no knowledge whether a third dimension represents space, or time. As such, particular characteristics of time cannot be taken care of.

Since the second half of 2010, the development of an R package spacetime started. It provides methods and classes for spatio-temporal data, and builds on the spatial data classes in sp and time series classes in xts. This document will explain how data in this form, and methods provided by this package, can be used for spatio-temporal geostatistics.

We will work with a data set with air quality (PM10) measurements over germany, taken from rural background stations available in the data sets provided by the European Environmental Agency.

```
> library(spacetime)
> rm(list = ls())
> data(air)
> ls()
[1] "DE_NUTS1" "rural"
```

# 2 Variography

## 2.1 Temporal autocorrelation and cross correlation

We will look into a subset of the data, ranging from 2005 to 2010, and remove stations that have only missing values in this period:

```
> rr = rural[,"2005::2010"]
> unsel = which(apply(as(rr, "xts"), 2, function(x) all(is.na(x))))
> r5to10 = rr[-unsel,]
> summary(r5to10)
Object of class STFDF
 with Dimensions (s, t, attr): (53, 1826, 1)
[[Spatial:]]
Object of class SpatialPoints
Coordinates:
coords.x1 6.28107 14.78617
coords.x2 47.80847 54.92497
Is projected: FALSE
proj4string:
[+init=epsg:4326 +proj=longlat +ellps=WGS84 +datum=WGS84 +no_defs
+towgs84=0,0,0]
Number of points: 53
[[Temporal:]]
     Index
                           ..1
        :2004-12-31
                             :2558
 Min.
                      Min.
 1st Qu.:2006-04-01
                      1st Qu.:3014
 Median :2007-07-01
                      Median:3470
```

```
:2007-07-01
Mean
                      Mean
                             :3470
3rd Qu.:2008-09-29
                      3rd Qu.:3927
Max.
       :2009-12-30
                      Max.
                             :4383
[[Data attributes:]]
                                                   NA's
  Min. 1st Qu. Median
                           Mean 3rd Qu.
                                           Max.
 0.560
                       16.260 20.330 269.100
                                                  21979
         9.275 13.850
```

Next, we will (rather arbitrarily) select four stations, which have the following labels:

```
> rn = row.names(r5to10@sp)[4:7]
> rn
```

#### [1] "DEBE056" "DEBE032" "DEHE046" "DENW081"

In the following, autocorrelation functions are computed and plotted. The resulting plot is shown in figure 1.

Auto- and cross correlations can be computed when a multivariate time series object is passed to acf:

```
> acf(na.omit(as(r5to10[rn,], "xts")))
```

The resulting plot is shown in figure 2. From these graphs one should be able to observe the following

- autocorrelations for lag 0 are always 1
- cross correlations for lag 0 are not always 1
- cross correlations can be asymmetric, meaning that when  $\rho_{AB}(h)$  is the correlation between  $Z(s_A, t)$  and  $Z(s_B, t + h)$ ,

$$\rho_{AB}(h) = \rho_{BA}(-h) \neq \rho_{AB}(-h)$$

with  $s_A$  and  $s_B$  the two stations between which a cross correlation is computed, and h the (directional!) lag between the series.

The plot further more shows that for these four stations the asymmetry is not very strong, but that cross correlations are fairly strong and of a similar form of autocorrelations.

This kind of plot does not work very well in layouts of e.g.  $10 \times 10$  sub-plots; acf automatically chooses  $4 \times 4$  as the maximimum a single plot. To try this out, do a  $7 \times 7$  plot

```
> acf(na.omit(as(r5to10[4:10,], "xts")))
```

and note that here we see in the last figure (DESH & DESN04) a pair of plots with nearly no cross correlation. This might have to do with the spatial distance between these two stations:

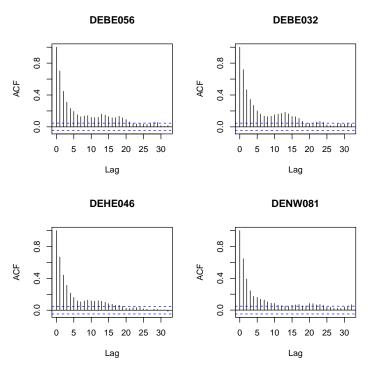


Figure 1: Autocorrelations for PM10; time lag unit in days.

> print(spDists(r5to10[4:10,]@sp), digits=3)

```
[,1]
             [,2] [,3] [,4] [,5] [,6] [,7]
[1,]
       0.0
            28.8
                   344
                        468
                              292
                                   235
[2,]
     28.8
             0.0
                   317
                        440
                              269
                                   231
                                         131
[3,] 343.6 317.3
                     0
                        150
                              309
                                   282
                                         285
[4,] 467.9 439.7
                   150
                          0
                              336
                                   432
                                         430
[5,] 291.7 268.5
                   309
                                   438
                        336
                                0
                                         362
[6,] 235.5 231.0
                   282
                        432
                              438
                                     0
                                         101
[7,] 134.6 130.9
                   285
                        430
                              362
                                   101
                                           0
```

(What is the spatial distance between stations DESH and DESN04?)

## 2.2 Spatial correlation, variograms

In the next steps, we will sample 100 time instances randomly,

```
> rs = sample(dim(r5to10)[2], 100)
```

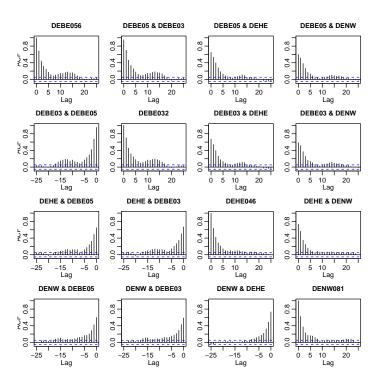


Figure 2: autocorrelations (diagional) and cross correlations (off-diagional) for the four stations selected; time lag unit in days.

we select these instances as a SpatialPointsDataFrame and add a time index to them. After this we bind them together in a single SpatialPointsDataFrame which has a time index ti:

```
> lst = lapply(rs, function(i) { x = r5to10[,i]; x$ti = i; x} )
> pts = do.call(rbind, lst)

Then, we can compute the pooled variogram
> library(gstat)
> v = variogram(PM10~ti, pts[!is.na(pts$PM10),], dX=0)
and plot it (figure 3):
> plot(v, fit.variogram(v, vgm(1, "Exp", 200, 1)))
> vmod = fit.variogram(v, vgm(1, "Exp", 200, 1))
> plot(v, vmod)

The fitted model is this:
> vmod
model psill range
```

Nug 3.471287 0.00000 Exp 73.655154 77.00846

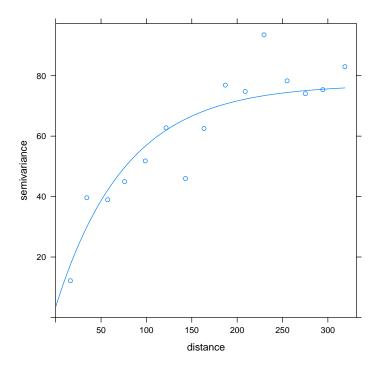


Figure 3: sample spatial variogram, averaged over 100 time randomly chosen time steps

One should note that the fit is rather poor, and not forget that we only have 53 stations selected. The time resolution is rich (1862 days) but the number of stations is small:

#### > dim(r5to10)

#### [1] 53 1826 1

We can fit a spatio-temporal variogram the usual way, by passing an object of class STFDF:

Alternatively, if this takes too long, a temporal subset can be taken, e.g. using the first 200 days:

taking random days from the full period will lead to the a wrong assumption that every time index increment reflect a constant lag increase. As an alternative, we will here load the precomputed S/T variogram:

#### > data(vv)

Plotting this object can be done in two ways, both shown in figure 4:

```
> plot(vv, ylab = "time lag (days)")
> plot(vv, map = FALSE, ylab = "time lag (days)")
```

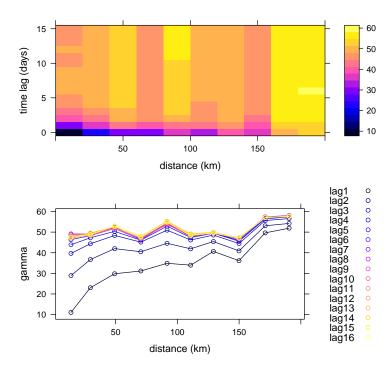


Figure 4: Spatio-temporal sample variogram map (top) and sample variograms for each time lag (bottom); both figures depict the information of object vv

### 2.3 Fitting a spatio-temporal variogram model

```
> ExpVgmMetric = function(x, s, t, nugget = 0) {
          h = sqrt(s^2 + (x[3] * as.numeric(t))^2)
+
          ifelse(h == 0, 0, nugget + x[1] * (1 - exp(-h / x[2])))
+
+
 }
 ExpFitFn = function(x, gfn, v, trace = FALSE, ...) {
          mod = gfn(x, v\$spacelag, v\$timelag, ...)
+
          resid = v\$gamma - mod
          if (trace)
                  print(c(x, MSE = mean(resid^2)))
          mean(resid^2)
+
 pars.M = optim(c(sill = 50, range = 100, anis = 50), ExpFitFn,
          gfn = ExpVgmMetric, v = vv, nugget = 10)
> pars.M$par
```

```
sill range anis
41.33730 83.17767 74.32199
```

The final model can be added, as fitted model to the sample variogram, and plotted by (figure 5):

```
> vv$model = ExpVgmMetric(pars.M$par, vv$spacelag, vv$timelag, nugget = 10)
> plot(vv)
```

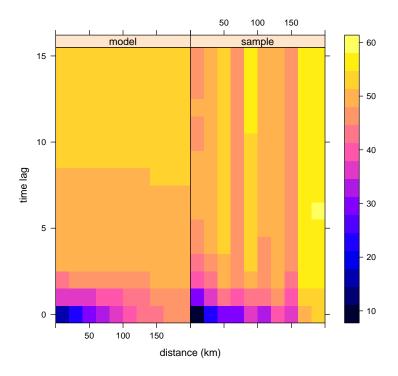


Figure 5: sample variogram map (right) and fitted metric model (left).

Now, let us try to fit and plot a separable model (figure 6):

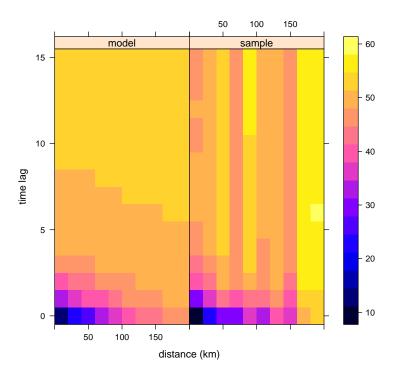


Figure 6: sample variogram map (right) and fitted seperable model (left).

A wireframe (3D) plot of the sample variogram can be obtained e.g. by

```
> library(lattice)
```

- > wireframe(gamma~spacelag\*timelag, vv, drape=TRUE,
- + col.regions=bpy.colors)

which is shown in figure 7.

# 3 Spatio-temporal prediction

The vignette in package spacetime gives an example of using the gstat function krigeST for spatio-temporal kriging of the Irish wind data. The variogram there needs to be specified as two separate variogram models, the product of which needs to be the target spatio-temporal variogram. The krigeST function uses global kriging, but only needs to invert the purely spatial and purely time covariance matrices.

For more generic spatio-temporal kriging where space is two-dimensional, one could use krige, defining the observations and prediction locations as three-dimensional data sets, see for an example

```
> demo(gstat3D)
```

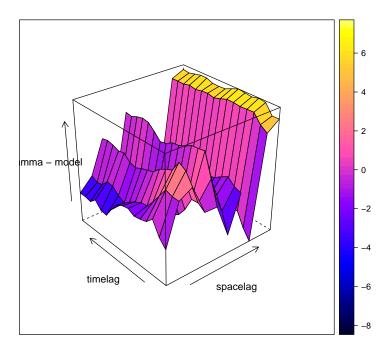


Figure 7: Wireplot of sample space-time variogram

It needs to be pointed out that in that case, the time (typically the third dimension) needs to be numeric, and three-dimensional anisotropy needs to be defined properly (see ?vgm).

In case the data set is too large for global kriging, one could try to use local kriging, and select data within some distance, or by specifying  $\mathtt{nmax}$  (the nearest n observations). In both cases, it is advisable to transform time such that one can use an isotropic variogram model in the three dimensions, as only in that case the nearest n observations correspond to the n most correlated observations.

An additional consideration is that in space-time, observations may not be regularly spaced. In some cases, the nearest n obervations may come from a single measurement location, which may lead to sharp jumps/boundaries in the interpolated values. This might be solved by using larger neighbourhoods, or by setting the omax in krige or gstat calls to the neighbourhood size to select per octant (this should be combined with specifying maxdist).

# References

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