



End of kriging? Or how tree-based Machine Learning Algorithms can be used to generate more accurate spatial predictions with combined geographical and feature space covariates

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Death to Kriging?



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Death to Kappa: birth of quantity disagreement and allocation disagreement for accuracy assessment

Robert Gilmore Pontius Jr & Marco Millones

Pages 4407-4429 | Received 27 Aug 2010, Accepted 20 Dec 2010, Published online: 17 Aug 2011

Download citation <http://dx.doi.org/10.1080/01431161.2011.552923>

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Abstract

The family of Kappa indices of agreement claim to compare a map's observed classification accuracy relative to the expected accuracy of baseline maps that can have two types of randomness: (1) random distribution of the quantity of each category and (2) random spatial allocation of the categories. Use of the Kappa indices has become part of the culture in remote sensing and other fields. This article examines five different Kappa indices, some of which were derived by the first author in 2000. We expose the indices' properties mathematically and illustrate their limitations graphically, with emphasis on Kappa's use of randomness as a baseline, and the often-ignored conversion from an observed sample matrix to the estimated population matrix. This article concludes that these Kappa indices are useless, misleading and/or flawed for the practical applications in remote sensing that we have seen. After more than a decade of working with these indices, we recommend that the profession abandon the use of Kappa indices for purposes of accuracy assessment and map comparison, and instead summarize the cross-tabulation matrix with two much simpler summary parameters: quantity disagreement and allocation

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Article

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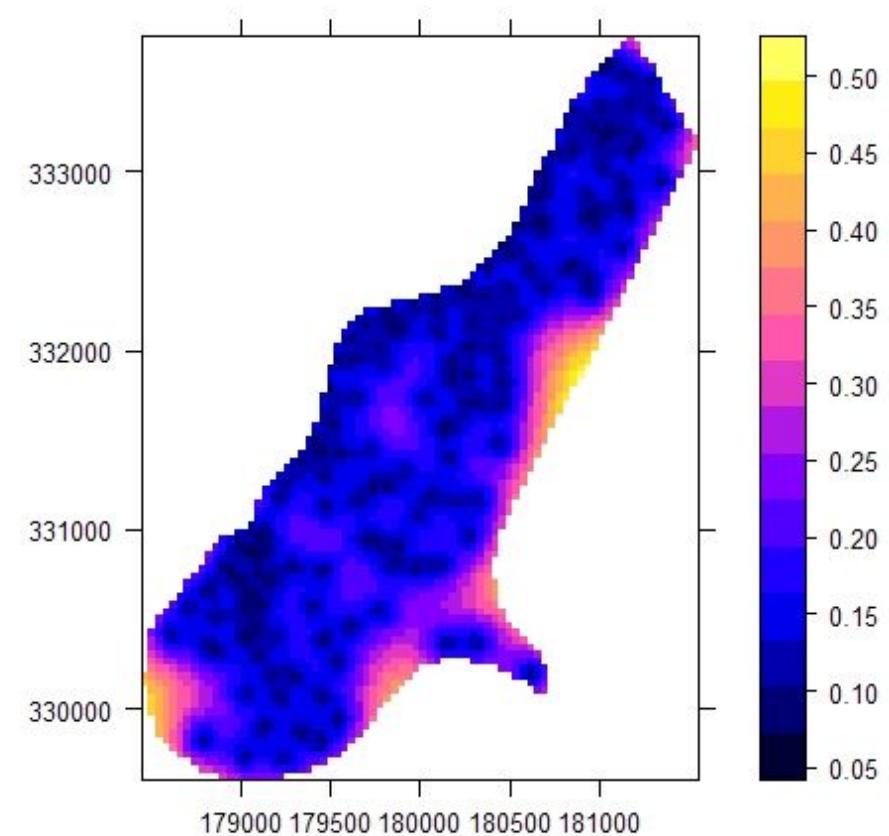
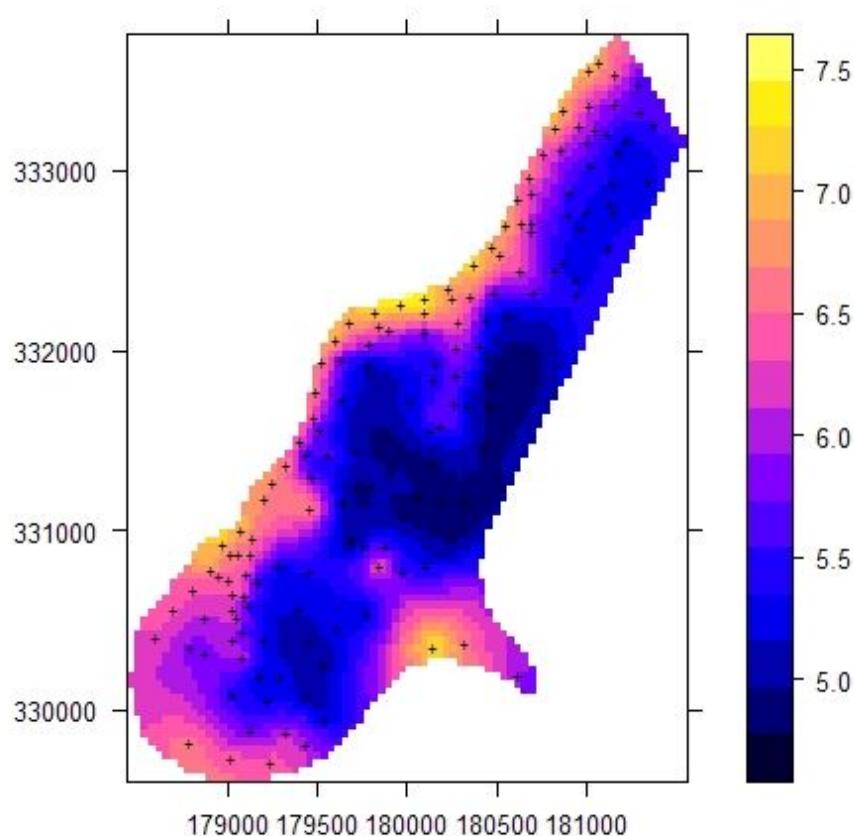
Robert Gilmore Pontius Jr et al.

International Journal of Remote Sensing

Published online: 4 Nov 2014



“Kriging” is basically a synonym for Geostatistics



Model-based geostatistics (state-of-the-art)



1. Determine distribution of the target variable and appropriate transformation (normal, log-normal, zero-inflated, Gamma, Poissonic ...)
2. Fit variogram (WLS, REML, ...), deal with multicollinearity (PCA?), non-stationary properties, support size, mixed effects...
3. Predict (mean values and uncertainty)
4. Validate predictions (mapping accuracy)



Regression-kriging

The Regression-kriging approach:

$$Y'(s) = m'(s) + e'(s) + \varepsilon$$

$$m'(s) = f[X(s)]; \quad e'(s) = f[h(s)|Y]$$

- First and second order stationarity
- Normal distribution (residuals)
- Anysotropy, multicollinearity...



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[PDF] researchgate.net

Case law

T Hengl, GBM Heuvelink, A Stein - Geoderma, 2004 - Elsevier
A methodological framework for spatial prediction based on **regression-kriging** is described and compared with ordinary kriging and plain regression. The data are first transformed using logit transformation for target variables and factor analysis for continuous predictors

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IOA Odeh, AB McBratney, DJ Chittleborough - Geoderma, 1995 - Elsevier

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About **regression-kriging**: from equations to case studies

[PDF] researchgate.net

T Hengl, GBM Heuvelink, DG Rossiter - Computers & geosciences, 2007 - Elsevier

This paper discusses the characteristics of **regression-kriging** (RK), its strengths and limitations, and illustrates these with a simple example and three case studies. RK is a spatial interpolation technique that combines a regression of the dependent variable on Cited by 435 Related articles All 10 versions Web of Science: 259 Cite Saved

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Two cultures: GLM vs RF

```
m = glm( zinc~log1p(dist)+ffreq, meuse,  
family=gaussian(link=log))
```

```
m = quantregForest(  
x=meuse@data[,c("dist","ffreq")],  
y=meuse$zinc)
```

Read more in: [Leo Braiman “Two cultures”](#)



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root RF vs OK

68cb598 34 minutes ago

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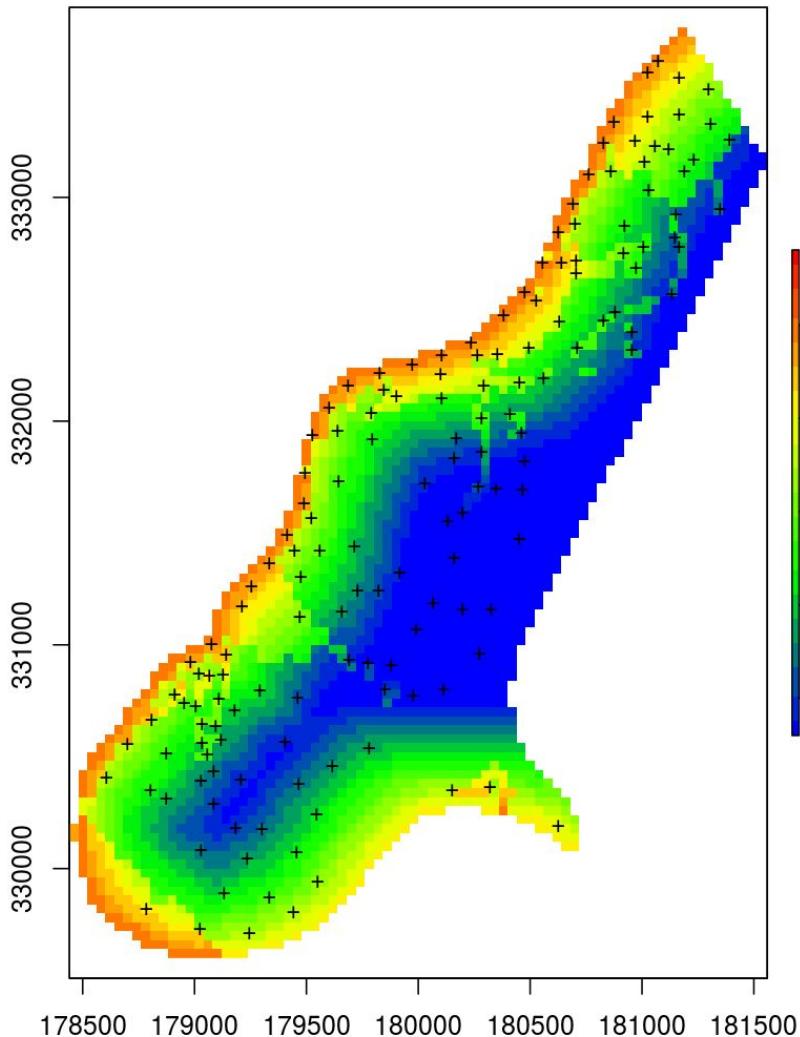
```
1 ## Comparison RF vs kriging
2 ## tom.hengl@isric.org
3
4 library(GSIF)
5 library(rgdal)
6 library(raster)
7 library(gstat)
8 library(randomForest)
9 library(quantregForest)
10 library(plotKML)
11 library(scales)
12 library(ranger)
13 library(RCurl)
14 #library(geoR)
15 leg = c("#0000ff", "#0028d7", "#0050af", "#007986", "#00a15e", "#00ca35", "#00f20d", "#1aff00", "#43ff00", "#6bff00", "#94ff00",
16 ## Load the Meuse data set:
17 demo(meuse, echo=FALSE)
18
19 ## compare GLM vs RF
20 m <- glm(zinc~log1p(dist)+ffreq, meuse, family=gaussian(link=log))
21 plot(m$fitted.values~m$y, asp=1)
22 abline(0,1)
23 rf <- quantregForest(x=meuse@data[,c("dist","ffreq")], y=meuse$zinc)
24 plot(rf$predicted~rf$y, asp=1)
25 abline(0,1)
26 meuse_grid_dalm_zinc <- predict(m, meuse_grid@data, type="response")
```

All examples shown are available via github

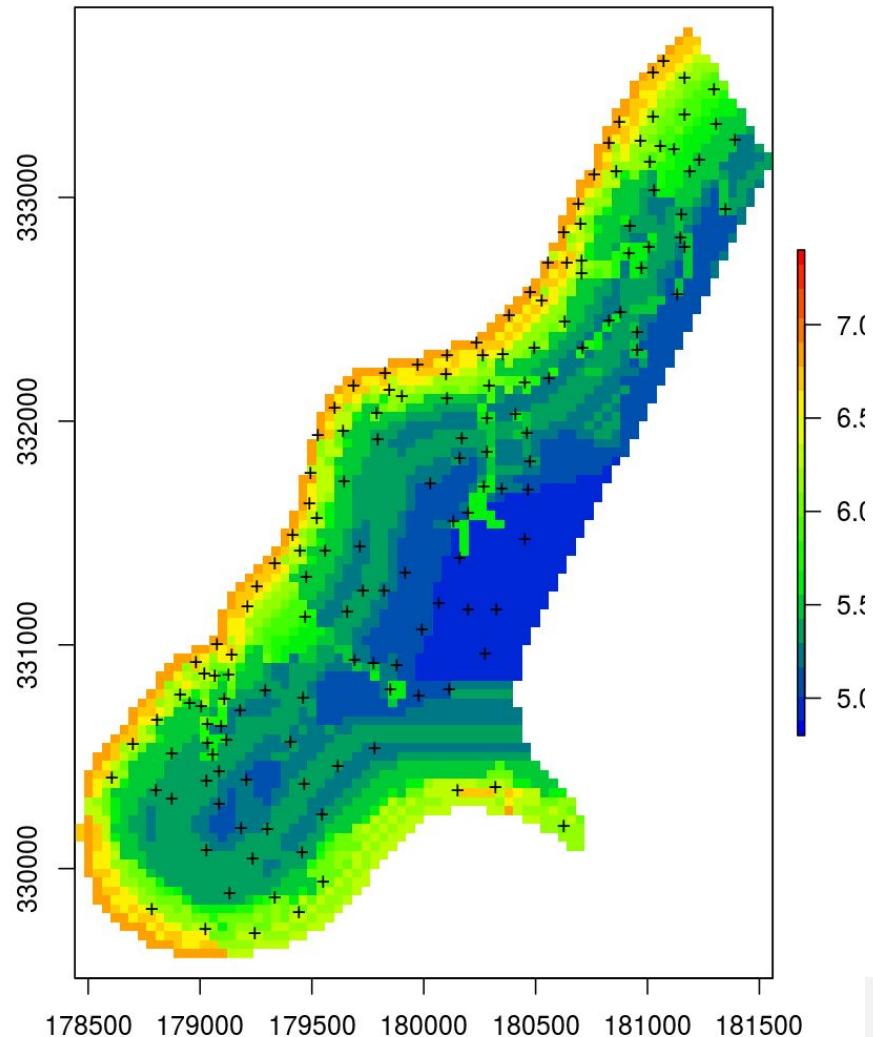
GLM vs RF (zinc meuse)



GLM



Random Forest





geoR

```
zinc.vgm <- likfit(zinc.geo, lambda=0,  
ini=c(var(log1p(zinc.geo$data)),500),  
cov.model="exponential")
```

```
zinc.ok <- krige.conv(zinc.geo, locations=locs,  
krige=krige.control(obj.m=zinc.vgm))
```

krige.conv: model with constant mean

krige.conv: performing the Box-Cox data transformation

krige.conv: back-transforming the predicted mean and variance

krige.conv: Kriging performed using global neighbourhood

Variogram (zinc meuse)

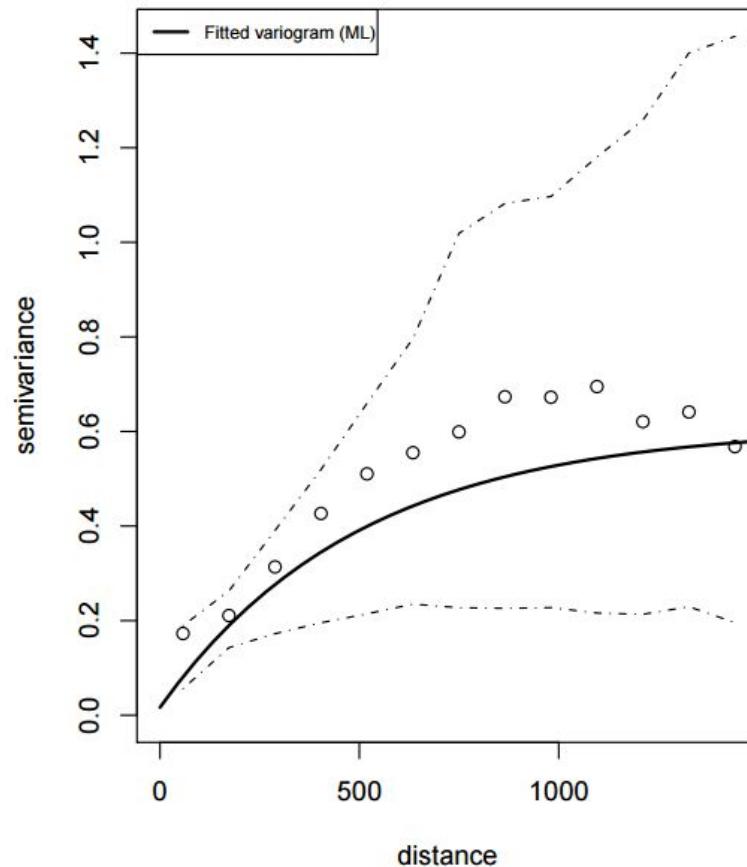
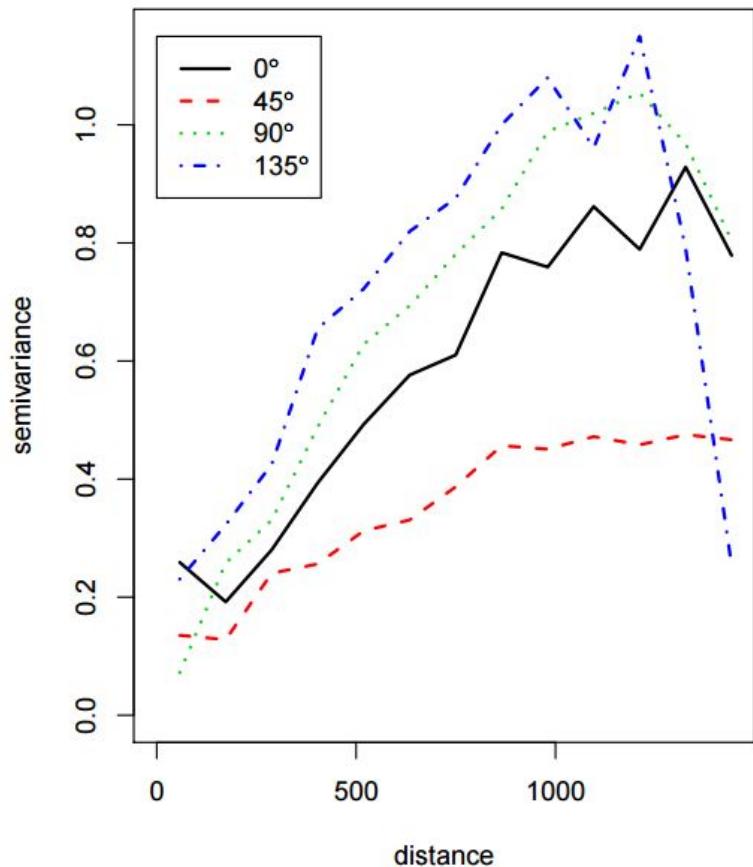


Fig. 5.15: Anisotropy (left) and variogram model fitted using the Maximum Likelihood (ML) method (right). The confidence bands (*envelopes*) show the variability of the sample variogram estimated using simulations from a given set of model parameters.

GSIF: correlate values with buffer distances



```
grid.dist0 <- buffer.dist(meuse["zinc"],  
meuse.grid[1], as.factor(1:nrow(meuse)))
```

```
dn0 <- paste(names(grid.dist0), collapse="+")  
fm0 <- as.formula(paste("zinc ~", dn0))
```

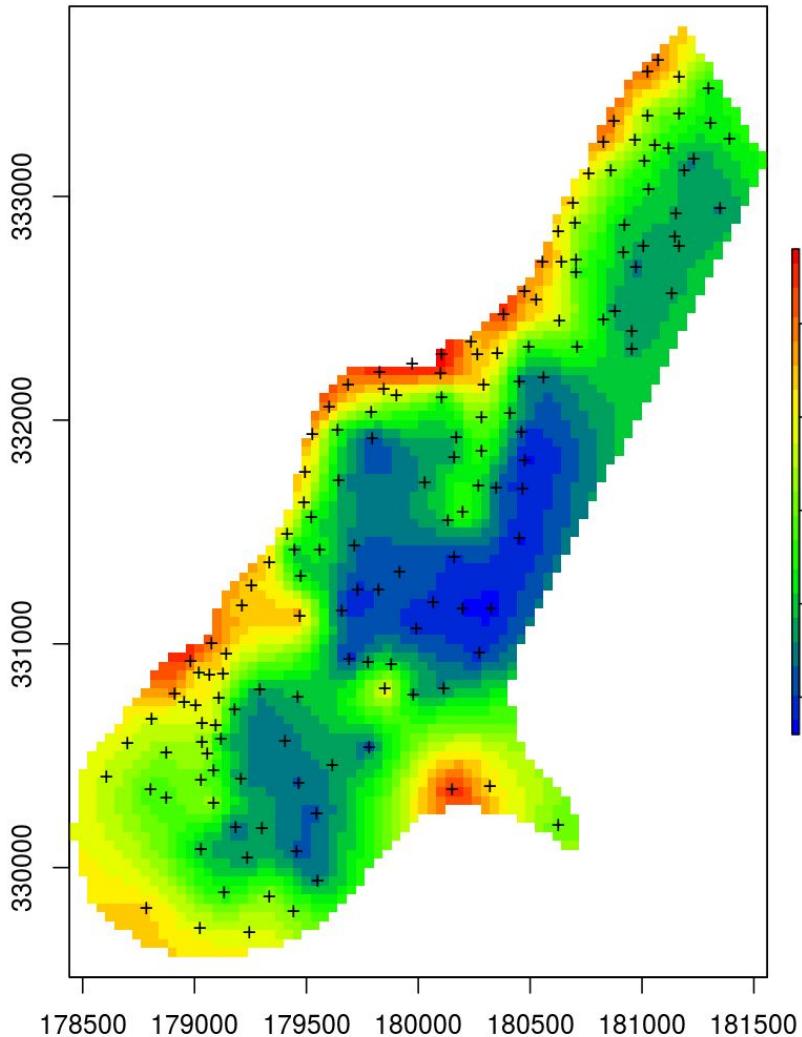
```
m0 <- fit.gstatModel(meuse, fm0, grid.dist0,  
method="ranger", rvgm=NULL)
```

```
rk.m0 <- predict(m0, grid.dist0)
```

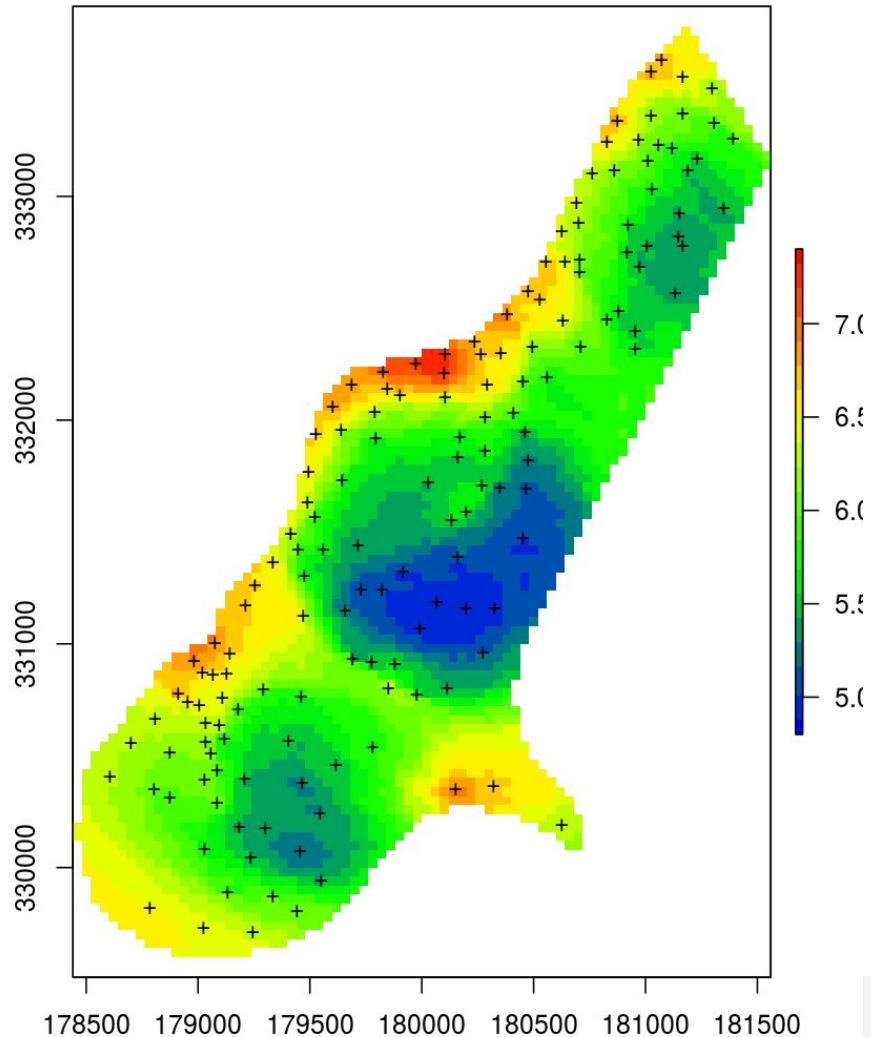
OK vs RF (zinc Meuse)



geoR (krige.conv)



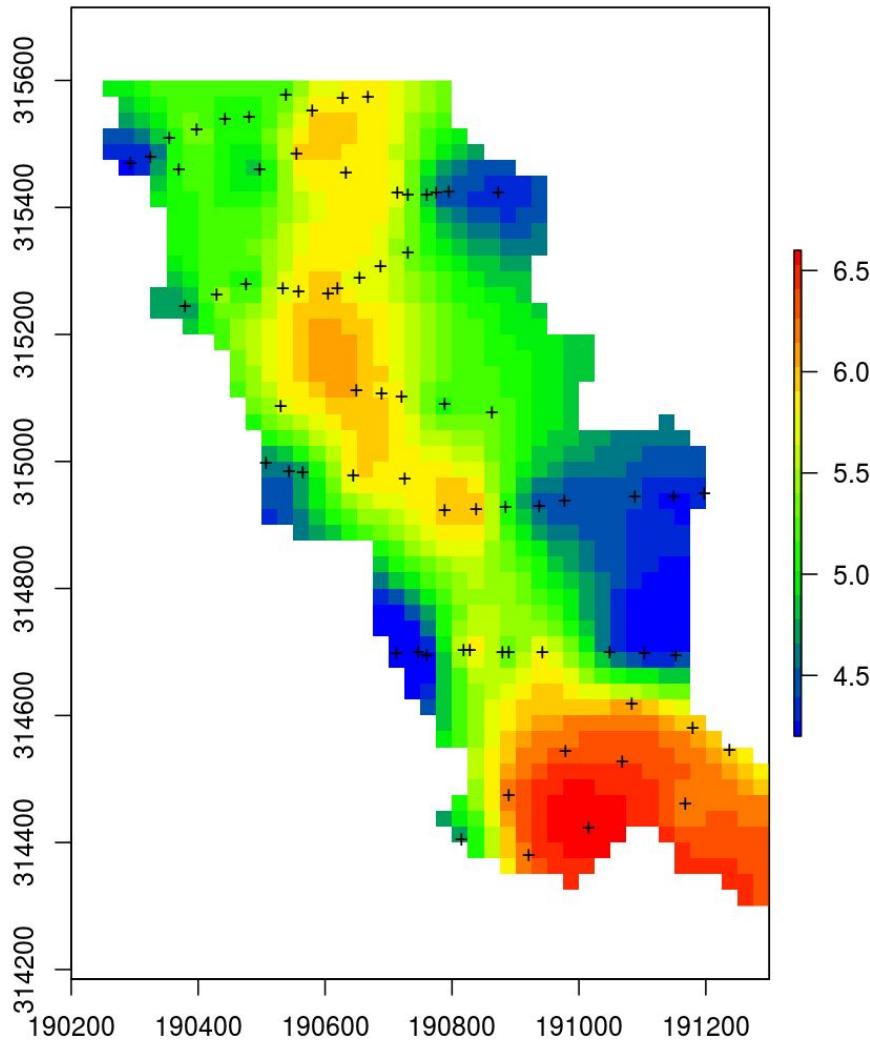
Random Forest



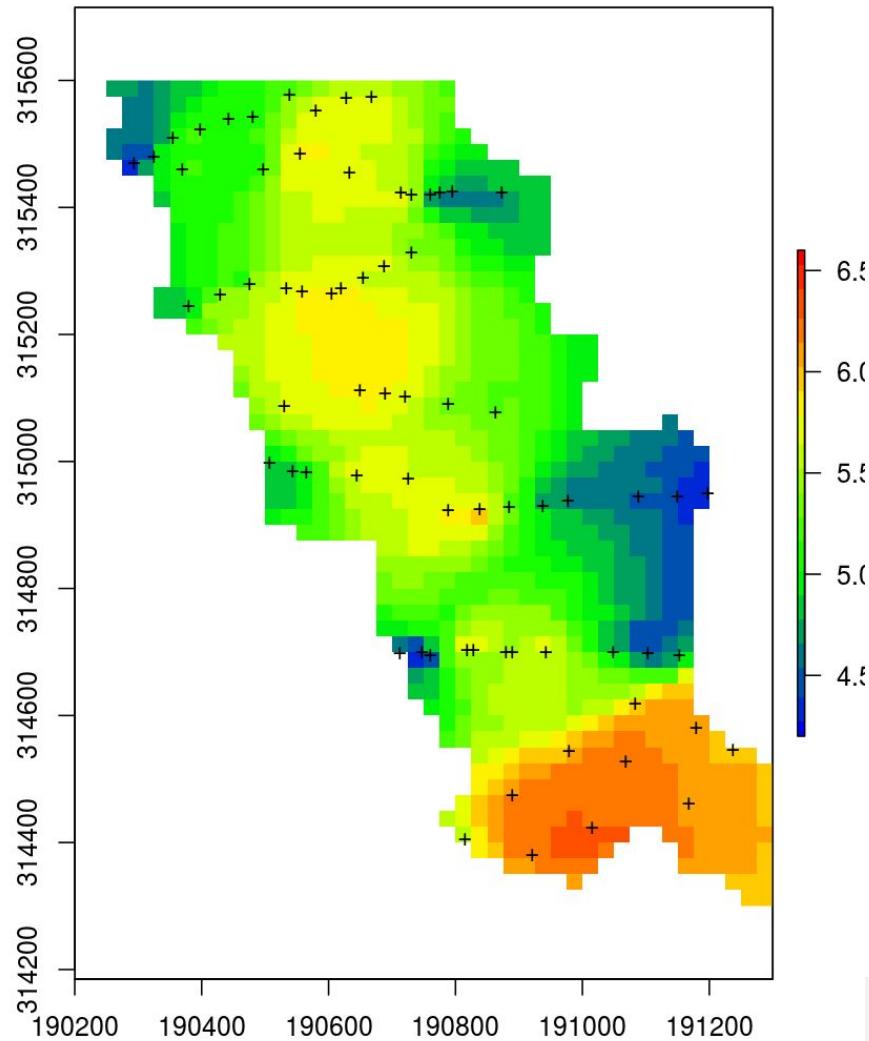
OK vs RF (Pb Gaul)



geoR (krige.conv)



Random Forest





sp predictions based on Random Forest

- ✓ No stationarity requirements.
- ✓ No Normal distribution requirements.
- ✓ No problems with choosing the right variogram (in fact, there is no need for a vgm at all).
- ✓ No (serious) problems with hot-spots.
- ✓ More complex distances can be added.



Random Forest complete list

Use both location-buffers + covariates:

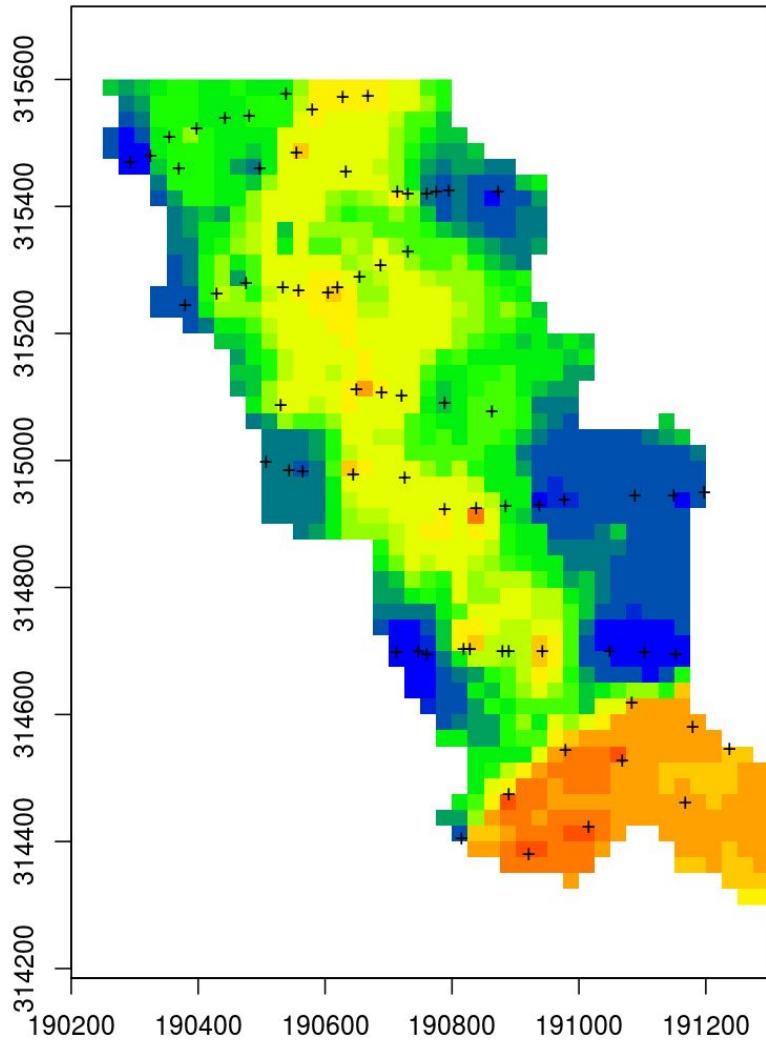
$$Y'(s) = f[h(s)|Y, X(s)]$$

- $h(s)|Y$ = buffer distances to points
- $X(s)$ = covariates / grids

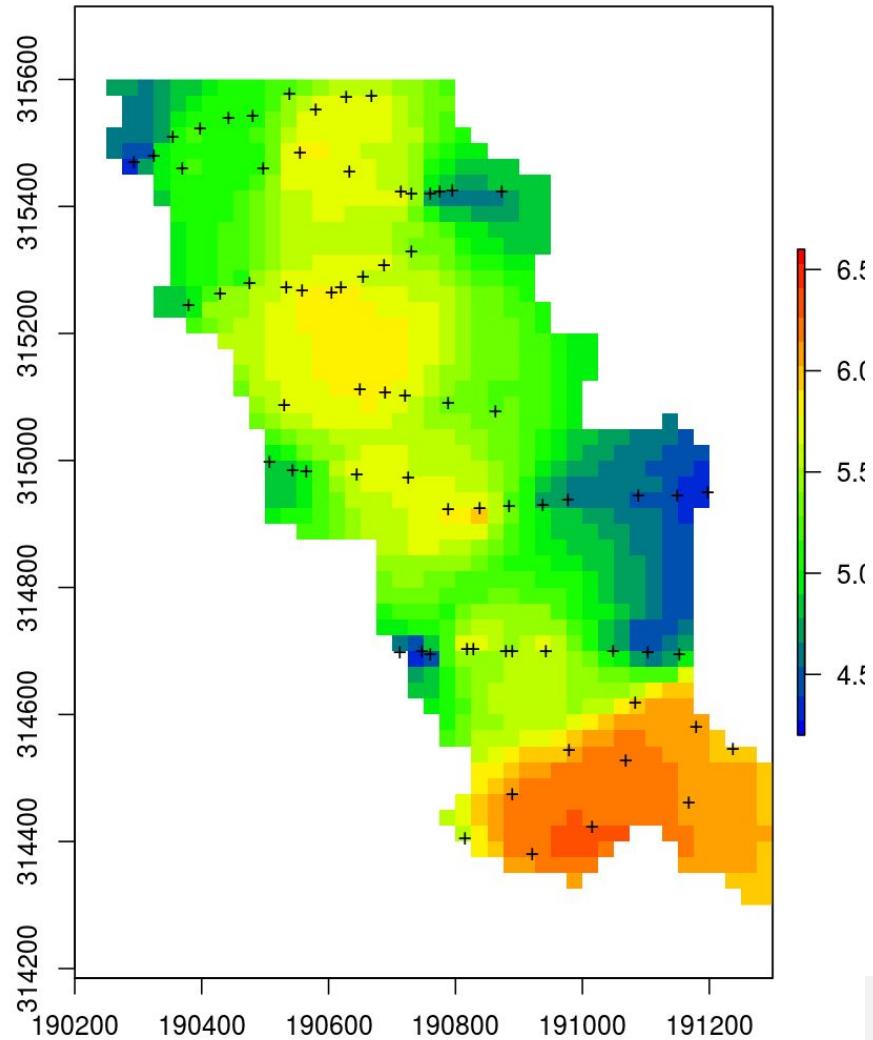
RF with covariates (Pb Gaul)

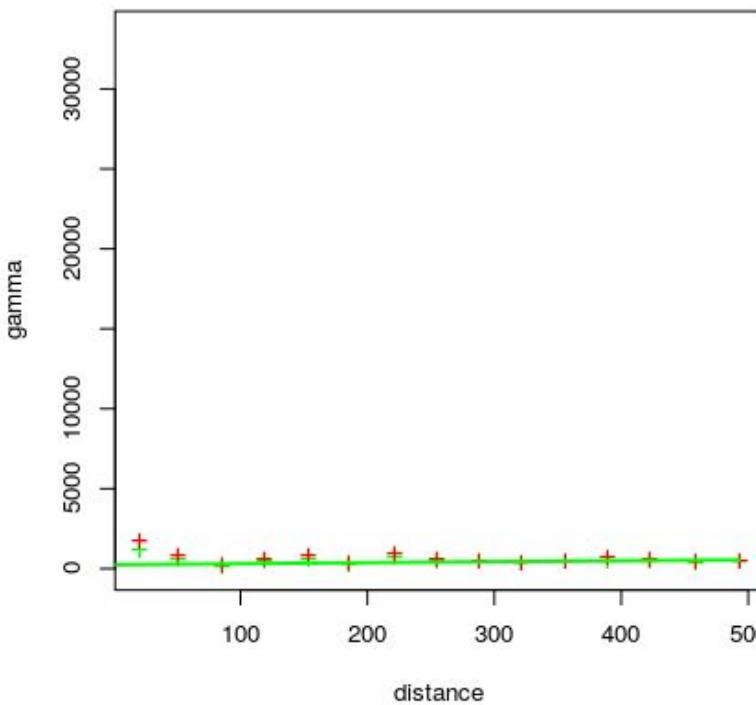
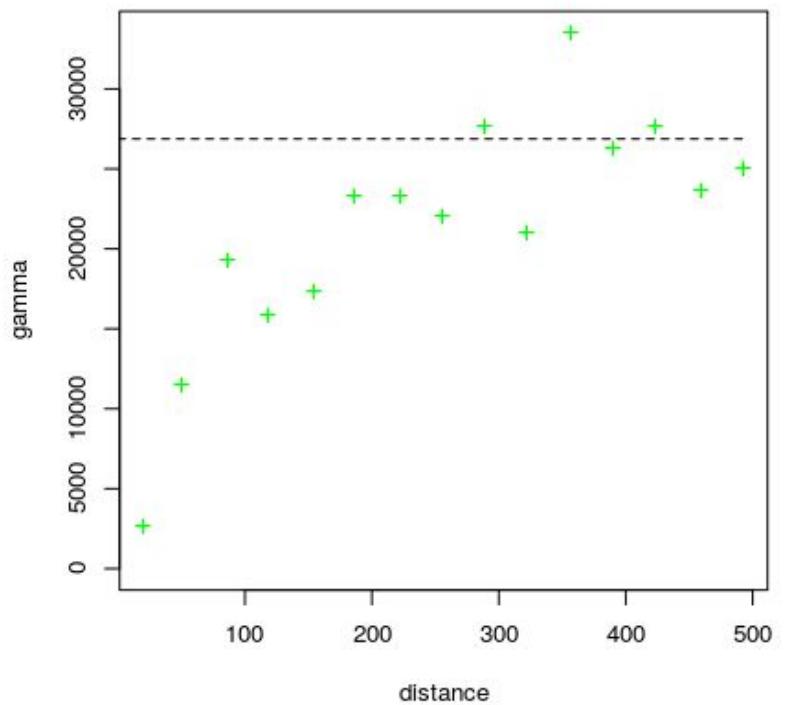
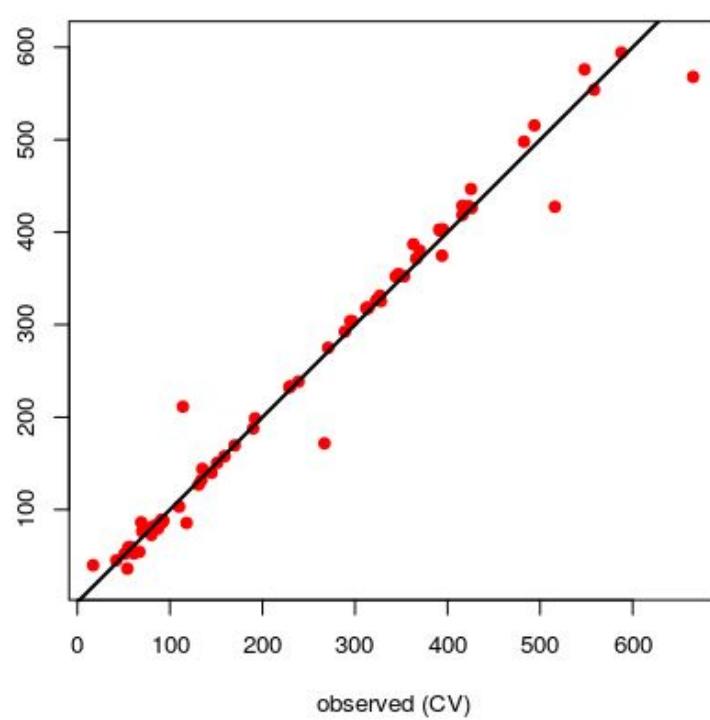
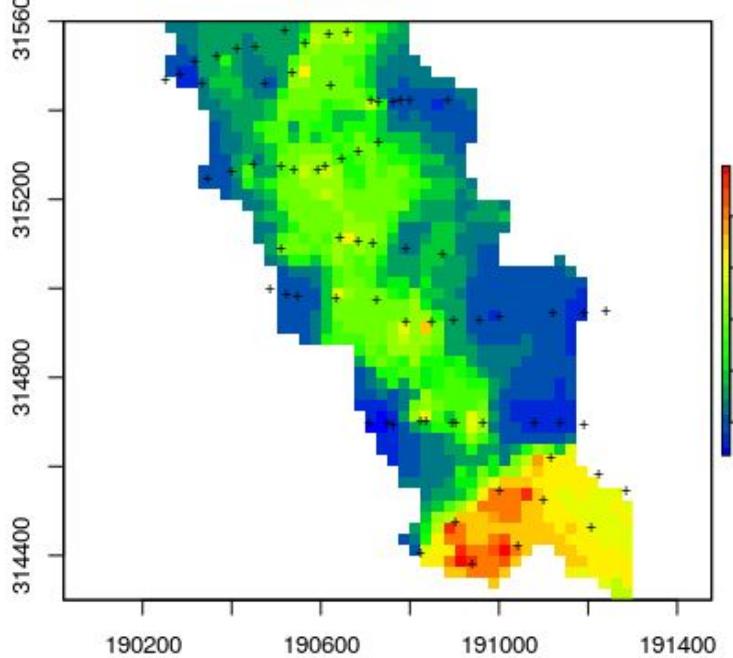


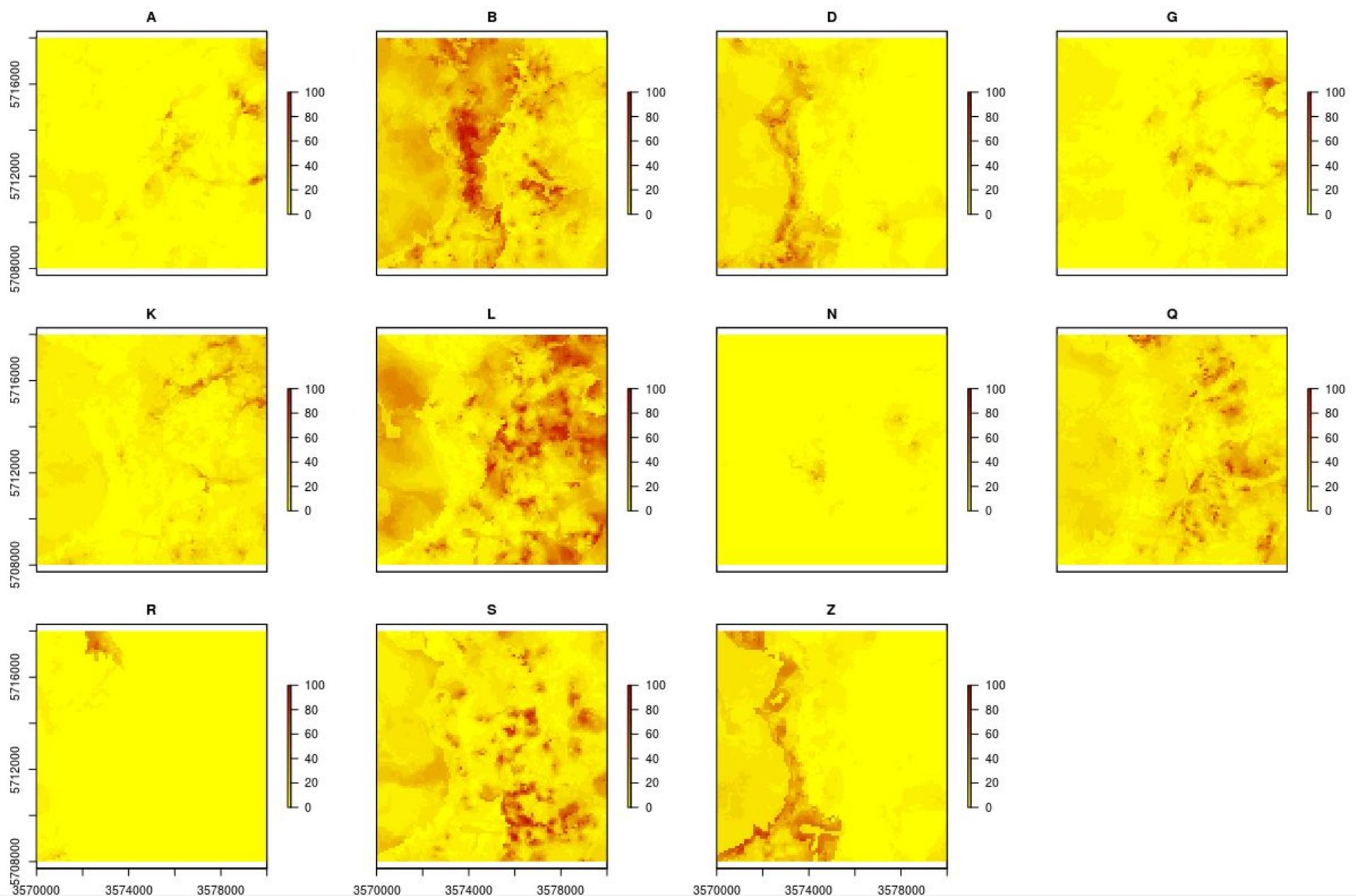
Random Forest + cova



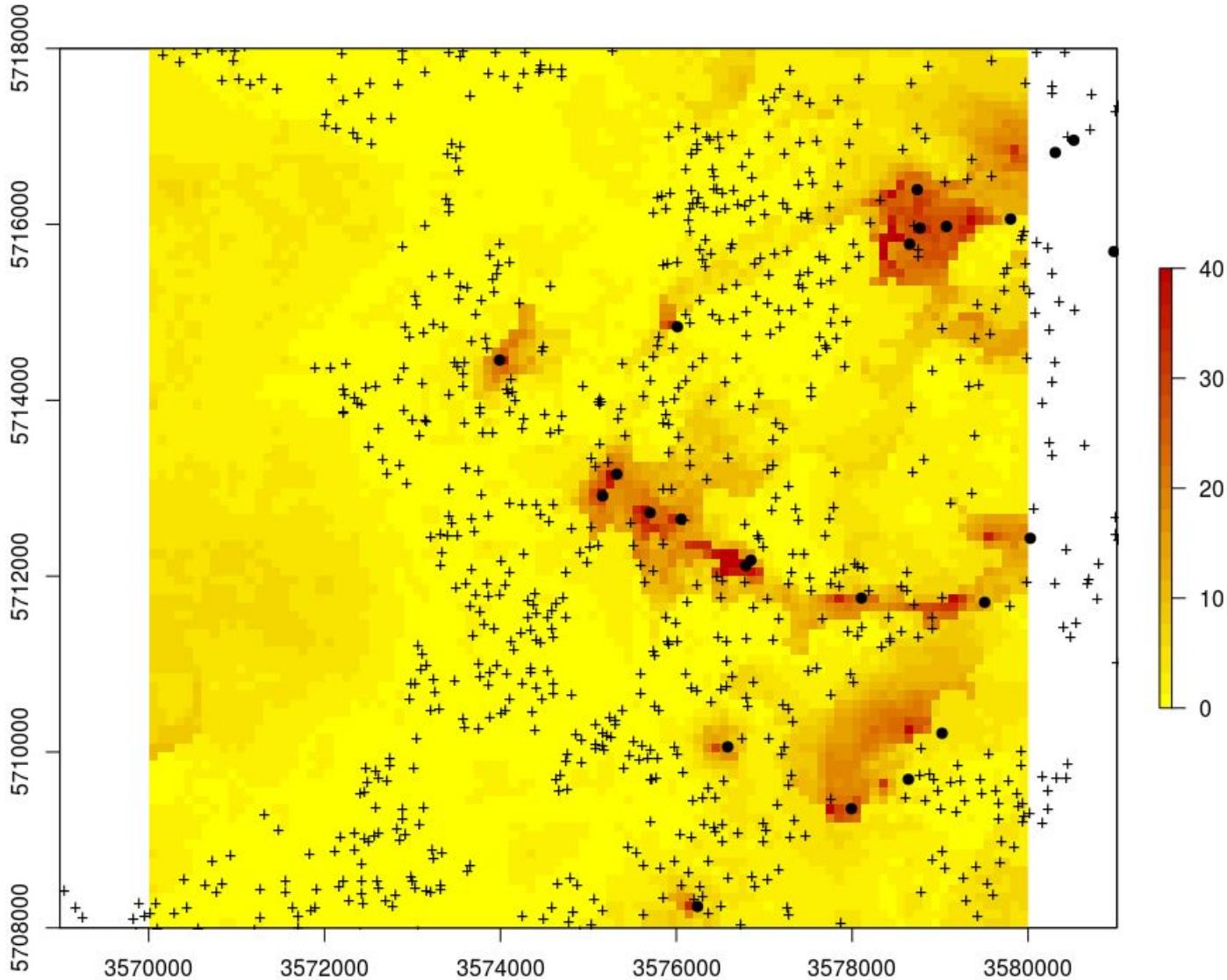
Random Forest



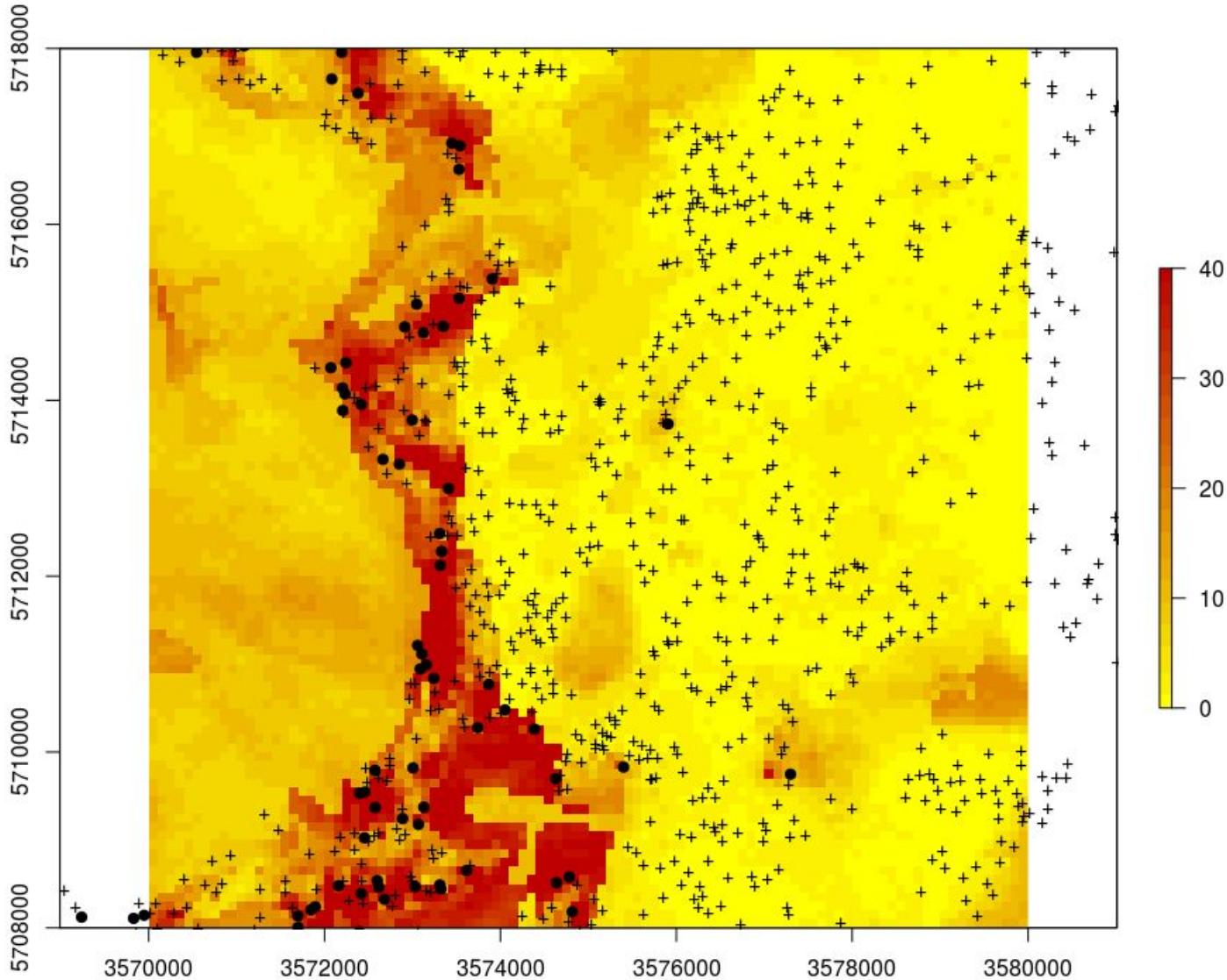




Soil type “G”



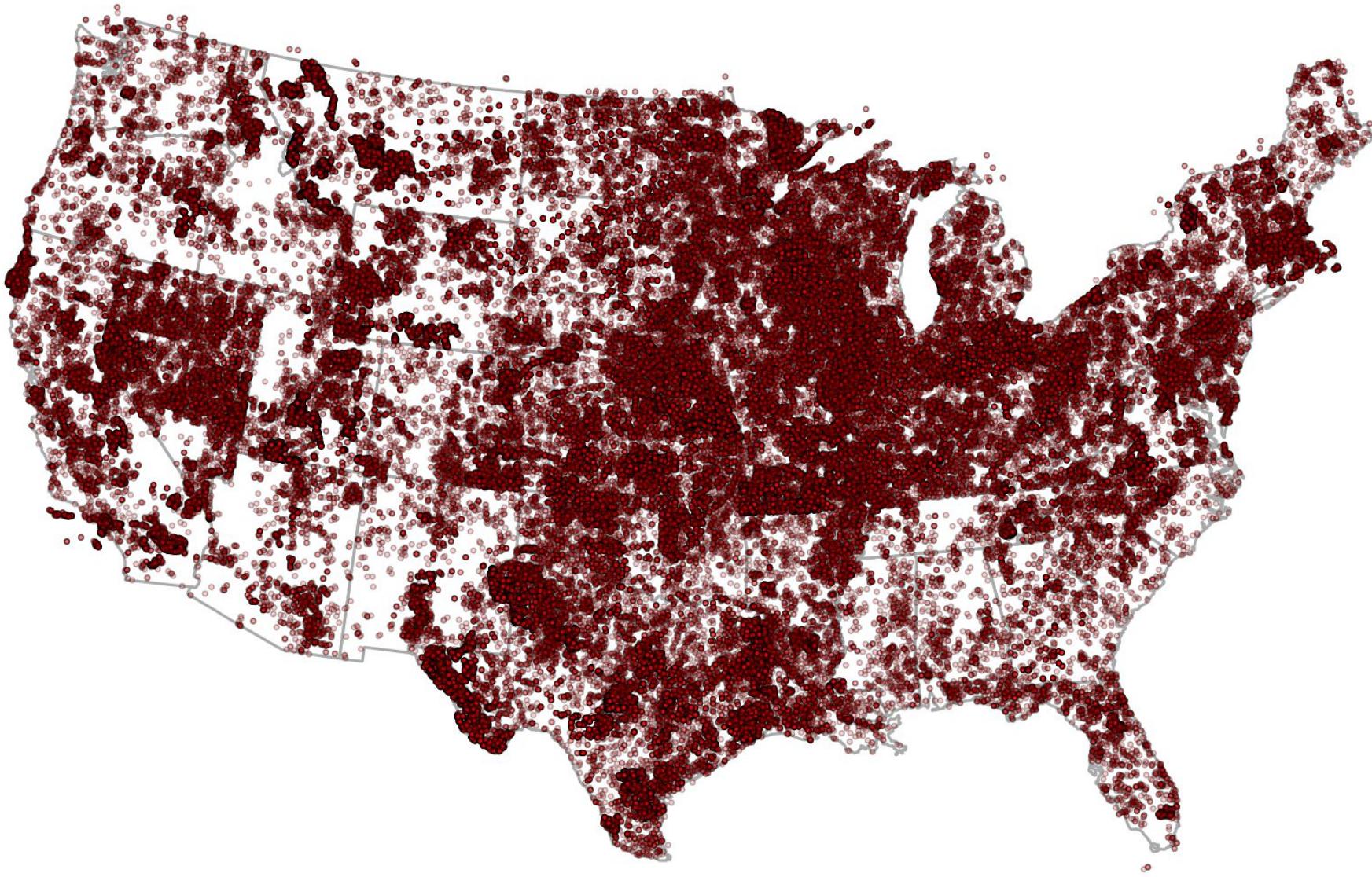
Soil type “D”

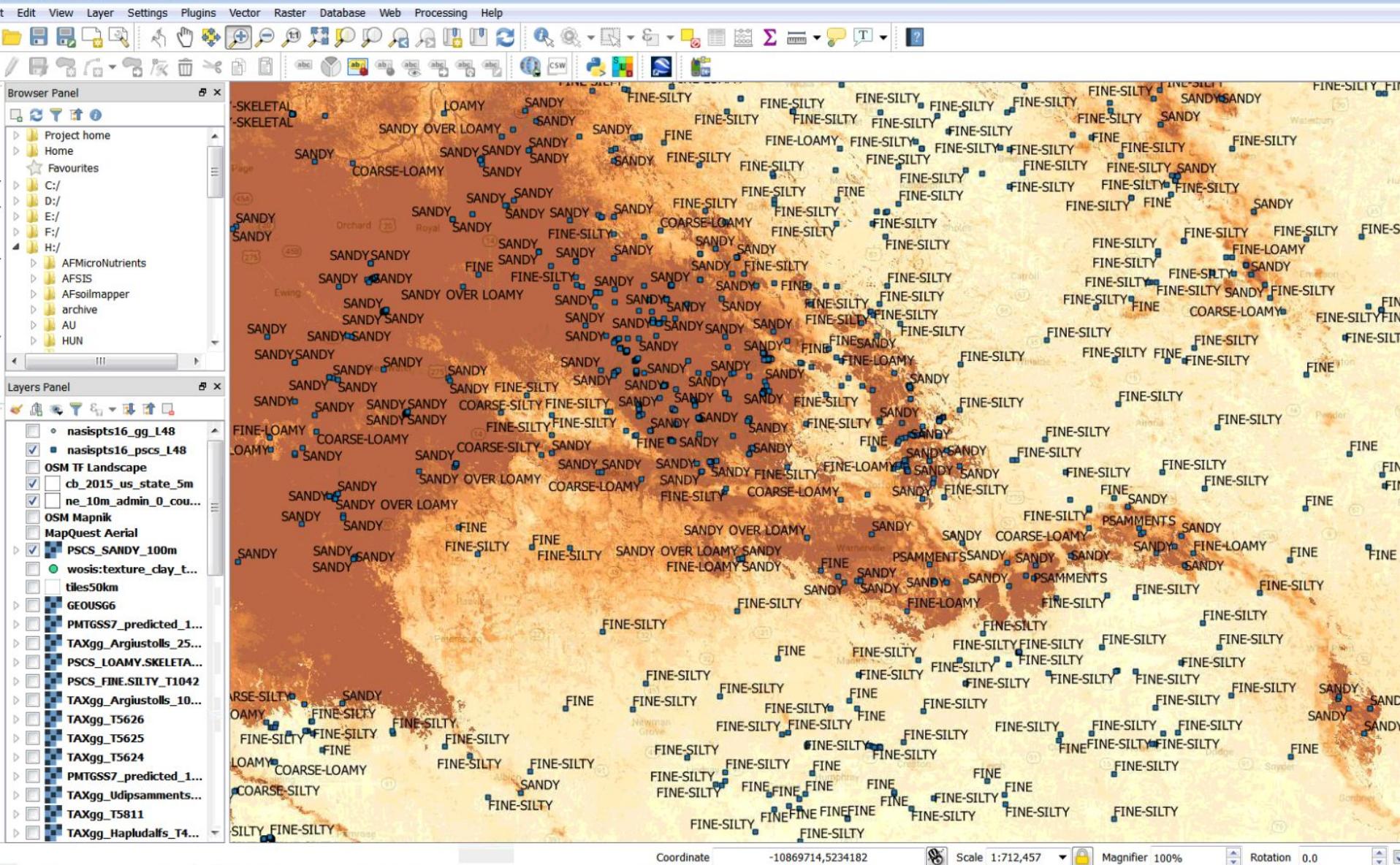




USA data

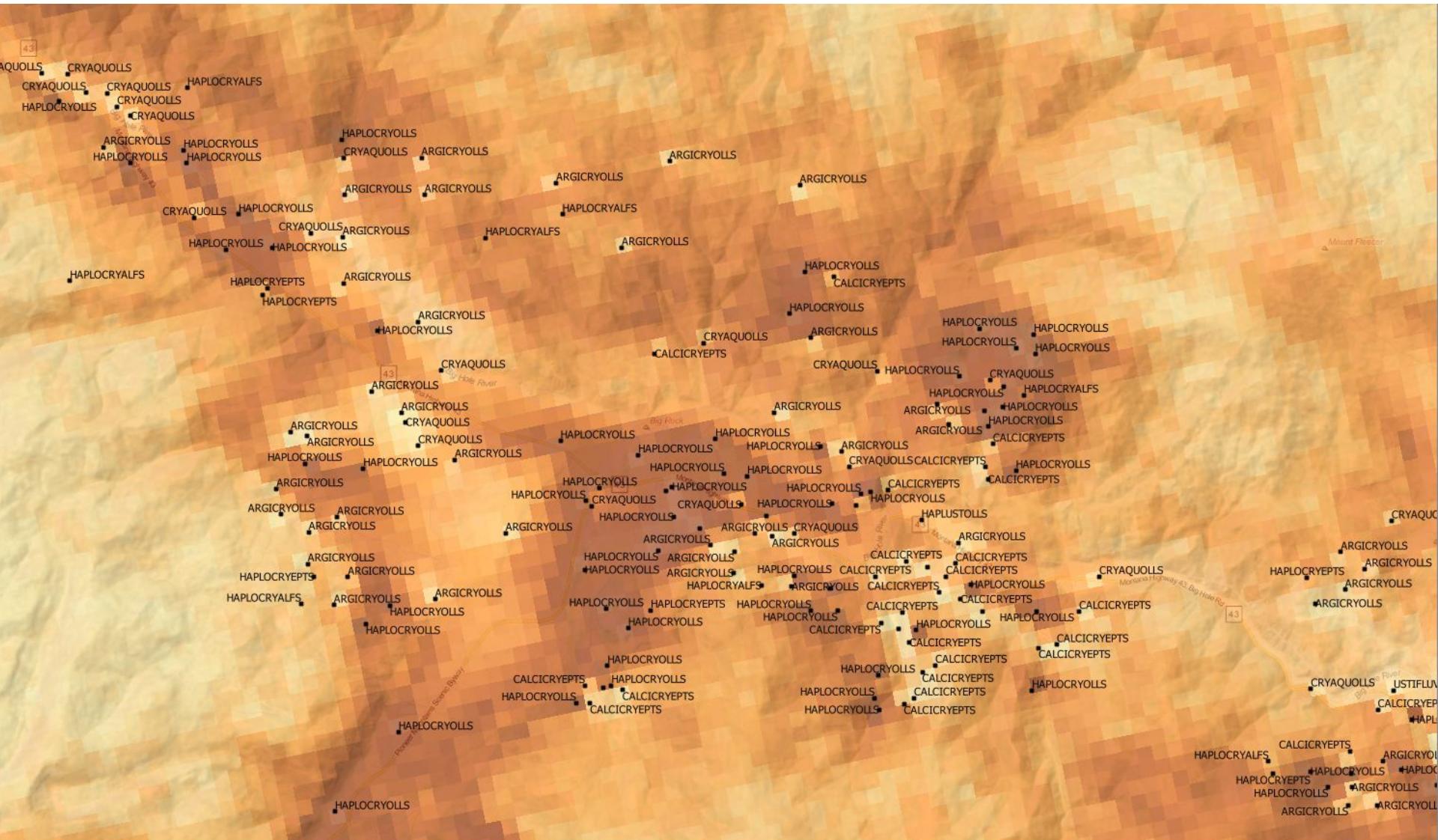
NASIS points (N=327,041)







Kriging patterns



Coordinate

-12571418,5758361



Scale

1:107,419

Rotation 0.0



Render

EPSG:3857 (OTF)





If machine learning is so
efficient in generating
spatial predictions, do we
still need kriging?

Random forest
Topic

Kriging
Topic

+ Add comparison

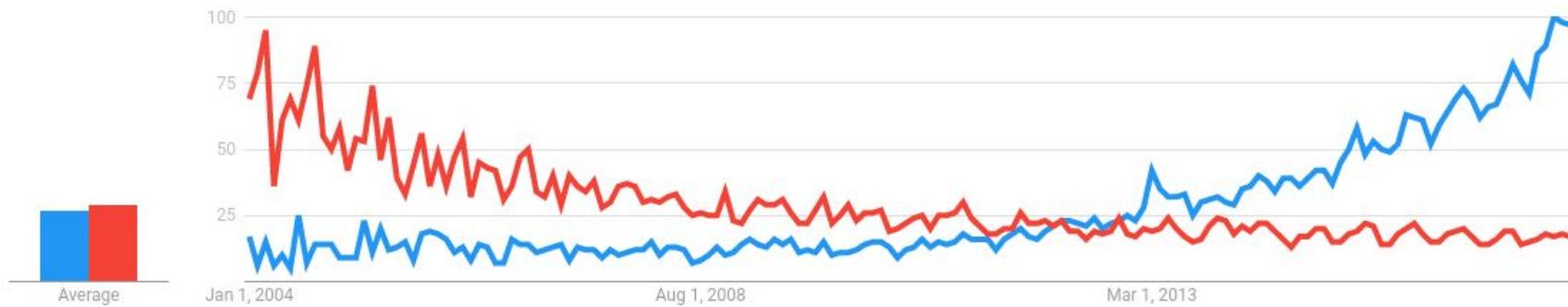
Worldwide ▾

2004 - present ▾

All categories ▾

Web Search ▾

Interest over time ?



Interest by region ?





A generic framework for sp

In essence, three types of covariates:

$$Y'(s) = f[h(s) | Y, X_R(s), X_P(s)]$$

- $h(s) | Y$ = buffer distances to points
- $X_R(s)$ = reflectances
- $X_P(s)$ = physical and chemical processes



Data-driven modeling (MLA)

1. Any target variable distribution is fine.
2. Multicollinearity – no problem.
3. A lot of covariates – no problem.
4. Complex non-linear relationships – great!

MLA = Hyper-parametric non-linear
(nested) models, usually requiring
intensive computing



However

- Computing intensity of MLA can be **MASSIVE**
- MLA usually **very very sensitive to artifacts in the input data** (even few typos can shift all predictions); most importantly **it requires high quality sampling**
- There are still many things unsolved (spatial simulations? spatial clustering? spacetime distances?)



Conclusions #1

Kriging?

Probably not needed any more.



Conclusions #2

Spatial statistics?

**Probably needed more
than ever!**



Conclusions #3

MLA is magic?

**Yes, but also very
sensitive on data quality
and still many many
issues to be solved.**

Computing

The Extraordinary Link Between Deep Neural Networks and the Nature of the Universe

Nobody understands why deep neural networks are so good at solving complex problems. Now physicists say the secret is buried in the laws of physics.

by Emerging Technology from the arXiv September 9, 2016



In the last couple of years, deep learning techniques have transformed the world of artificial intelligence. One by one, the abilities and techniques that humans once imagined were uniquely our own have begun to fall to the onslaught of ever more powerful machines. Deep neural networks are now better than humans at tasks such as face recognition and object recognition. They've mastered the ancient game of Go and thrashed the best human players.

But there is a problem. There is no mathematical reason why networks arranged in layers should be so good at these challenges.

Mathematicians are flummoxed. Despite the huge success of deep



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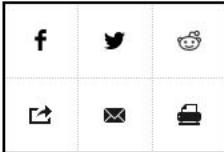
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How the Computer Beat the Go Master

As a leading go player falls to a machine, artificial intelligence takes a decisive step on the road to overtaking the natural variety

By Christof Koch on March 19, 2016

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South Korean professional Go player Lee Sedol is seen on a TV screen during the

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