# Spatial point processes: Theory and practice illustrated with R

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Lecture II, Feburary 17, 2011

# Contents of lecture II

- Properties for the Poisson point process.
- Maximum likelihood estimation for Poisson point processes.
- Investigating dependence between points using functional summary statistics.
- Relevant commands in spatstat (remember to load the package when a new session is started):
  - > library(spatstat)

#### Homogeneous Poisson processes

- The homogeneous Poisson process with intensity λ is defined by two properties:
  - N(A) is Poisson distributed with mean  $\lambda |A|$ , for all A.
  - Conditional on N(A) = n the n points are independent and uniformly distributed in A.
- The counts N(A) and N(B) are independent for disjoint A, B.
- We will use the homogeneous Poisson process as the basis for creating more complicated point process models.

#### Density for point processes

■ A point process **X** in the window *W* has density *f*(·) with respect to the unit rate Poisson process if

$$\mathbb{E}[h(\mathbf{X})] = \mathbb{E}[h(\mathbf{Y})f(\mathbf{Y})]$$
(1)

for all functionals h, where  ${\bf Y}$  is a unit rate Poisson process (i.e.  $\lambda=1).$ 

 $\blacksquare$  In particular the homogeneous Poisson process with intensity  $\lambda$  has density

$$f(\mathbf{x}) = e^{(1-\lambda)|W|} \lambda^{n(\mathbf{x})}.$$
 (2)

The maximum likelihood estimate  $\hat{\lambda}$  of the intensity  $\lambda$  is

$$\hat{\lambda} = \frac{n(\mathbf{x})}{|W|}$$

### Inhomogeneous Poisson processes

- The inhomogeneous Poisson process with intensity function  $\lambda(\cdot)$  is defined by two properties:
  - ▶ N(A) is Poisson distributed with mean  $\int_A \lambda(u) du$ , for all A.
  - ► Conditional on N(A) = n the n points are independent and identically distributed in A with density proportional to λ(·).
- The counts N(A) and N(B) are still independent for disjoint A, B.
- The density of an inhomogeneous Poisson process on W with intensity  $\lambda_{\theta}$  parametrised by  $\theta$  is

$$f(\mathbf{x}) = \exp\left(\int_{W} (1 - \lambda_{\theta}(u)) \,\mathrm{d}u\right) \prod_{i} \lambda_{\theta}(x_{i}). \tag{3}$$

#### Inhomogeneous Poisson processes

- For simulation a useful technique is independent thinning: Suppose that **X** is a Poisson process with intensity function  $\beta(\cdot)$ , and that each point of **X** is either deleted or retained, independently of other points. If the retention probability at location *u* is p(u), then the resulting process of retained points is Poisson, with intensity function  $\lambda(u) = \beta(u)p(u)$ , for all *u*.
- To simulate an inhomogeneous Poisson process with intensity function  $\lambda(\cdot)$  we start by simulating a homogeneous Poisson process with intensity  $\lambda_{\max} = \max_u \{\lambda(u)\}$ . Then a thinning is performed with retention probabilies  $p(u) = \lambda(u)/\lambda_{\max}$  to obtain the inhomogeneous process.

# Inhomogeneous Poisson processes

To simulate an inhomogeneous Poisson process we can supply the intensity as a function or a pixel image:

- > intenfun <- function(x, y) 200 \* x</pre>
- > plot(rpoispp(intenfun, lmax = 200), main = "")
- > intenim <- as.im(intenfun, dimyx = c(100, 100),

> plot(rpoispp(intenim), main = "")



The spatstat implementation requires the intensity function  $\lambda_{\theta}(u)$  to be loglinear in the parameter  $\theta$ :

$$\log \lambda_{\theta}(u) = \theta \cdot S(u) \tag{4}$$

where S(u) is a real-valued or vector-valued function of location u. The form of S is arbitrary so this is not much of a restriction.

In practice S(u) could be a function of the spatial coordinates of u, or an observed covariate, or a mixture of both.

- The fitting function is called ppm ('point process model') and is very closely analogous to the model fitting functions in R such as lm and glm.
- The statistic S(u) is specified by an R language formula, like the formulas used to specify the systematic relationship in a linear model or generalised linear model.
- The basic syntax is:
  - > ppm(X, ~trend)

where X is the point pattern dataset, and ~trend is an R formula with no left-hand side. This should be viewed as a model with log link, so the formula ~trend specifies the form of the logarithm of the intensity function.

Fitting a Poisson model with an intensity that is log-linear in the cartesian coordinates, i.e.  $\lambda_{\theta}((x, y)) = \exp(\theta_0 + \theta_1 x + \theta_2 y)$ :

- > data(bei)
- > ppm(bei, ~x + y)

Nonstationary Poisson process

```
Trend formula: ~x + y
```

Fitted coefficients for trend formula: (Intercept) x y -4.7245290274 -0.0008031288 0.0006496090

In this case the fitted intensity function is

$$\lambda_{\theta}((x, y)) = \exp\left(-4.724529 + -0.000803 \, x + 0.00065 \, y\right).$$

To fit an inhomogeneous Poisson model with an intensity that is log-quadratic in the cartesian coordinates, i.e. such that  $\log \lambda_{\theta}((x, y))$  is a quadratic in x and y:

> ppm(bei, ~polynom(x, y, 2))

Nonstationary Poisson process

Trend formula: ~polynom(x, y, 2)

```
Fitted coefficients for trend formula:

(Intercept) polynom(x, y, 2)[x]

-4.275762e+00 -1.609187e-03

polynom(x, y, 2)[y] polynom(x, y, 2)[x<sup>2</sup>]

-4.895166e-03 1.625968e-06

polynom(x, y, 2)[x.y] polynom(x, y, 2)[y<sup>2</sup>]

-2.836387e-06 1.331331e-05
```

To fit a model with constant but unequal intensities on each side of the vertical line x = 500, the explanatory variable S(u) should be a factor with two levels, Left and Right say, taking the value Left when the location u is to the left of the line x = 500.

```
> ppm(bei, ~side(x))
```

Nonstationary Poisson process

```
Trend formula: ~side(x)
```

```
Fitted coefficients for trend formula:
(Intercept) side(x)right
-4.8026460 -0.2792705
```

For the rainforest data we may want to fit the inhomogeneous Poisson model with intensity which is a loglinear function of slope, i.e.

$$\lambda(u) = \exp(\beta_0 + \beta_1 Z(u))$$
(5)

where  $\beta_0, \beta_1$  are parameters and Z(u) is the slope at location u:

```
> slope <- bei.extra$grad
> ppm(bei, ~slope, covariates = list(slope = slope))
```

Nonstationary Poisson process

Trend formula: ~slope

Fitted coefficients for trend formula: (Intercept) slope -5.390553 5.022021

It might be more appropriate to fit the model

$$\lambda(u) = \beta Z(u) \tag{6}$$

where again Z(u) is the slope at u. Equivalently

$$\log \lambda(u) = \log \beta + \log Z(u). \tag{7}$$

There is no coefficient in front of the term  $\log Z(u)$  in (7), so this term is an 'offset'. To fit this model,

> ppm(bei, ~offset(log(slope)), covariates = list(slope = s Nonstationary Poisson process

Trend formula: ~offset(log(slope))

```
Fitted coefficients for trend formula:
(Intercept)
-2.427127
```

0.01 0.01

0.008

0.006

D.004

# Predicting the intensity function

The values of the intensity funcion can be calculated on a grid in the window for a fitted model via the function predict. Alternatively, we can simply plot the fitted object which calls the predict-function and plots the intensity function:

Fitted trend

# Variance of estimates

For Poisson models it is possible to calculate the variance-covariance matrix of  $\widehat{\theta}:$ 

> vcov(fit)

	(Intercept)	х	У
(Intercept)	1.854091e-03	-1.491267e-06	-3.528289e-06
x	-1.491267e-06	3.437842e-09	1.208410e-14
У	-3.528289e-06	1.208410e-14	1.338955e-08

This can be used to give standard deviations on the estimates:

i	$\theta_i$	$\operatorname{Var}(\widehat{\theta}_i)$	standard deviation
0	-4.724529	0.001854091	0.04305915
1	-0.0008031288	3.437842e-09	5.863311e-05
2	0.000649609	1.338955e-08	0.0001157132

#### Covariate effect plot

It is also possible to plot the 'effect' of a single covariate in the model. The command effectfun computes the intensity of the fitted model as a function of one of its covariates. This is chiefly useful if the model only has one covariate.

> fit <- ppm(bei, ~slope, covariates = list(slope = slope))
> plot(effectfun(fit, "slope"), main = "")



slope

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### Analysis of deviance

Analysis of deviance for **nested** Poisson point process models is done by using the function anova:

```
> fitnull <- update(fit, ~1)</pre>
> anova(fitnull, fit, test = "Chi")
Analysis of Deviance Table
Model 1: .mpl.Y ~ 1
Model 2: .mpl.Y ~ slope
  Resid. Df Resid. Dev Df Deviance P(>|Chi|)
1
      20507
                 18728
2
      20506
                 18346 1 382.25 < 2.2e-16 ***
Signif. codes: 0 âĂŸ***âĂŹ 0.001 âĂŸ**âĂŹ 0.01 âĂŸ*âĂŹ 0.0
Clearly, the model including slope is favored.
```

### Akaike Information Criterion

The homogeneous Poisson process is not nested in the model (6), in which intensity is proportional to slope. One possibility here is to use the Akaike Information Criterion **AIC** for model selection:

- > fit <- ppm(bei, ~offset(log(slope)), covariates = list(si > fitnull <- ppm(bei, ~1)</pre>
- > AIC(fit)
- [1] 42496.65

> AIC(fitnull)

[1] 42763.92

The smaller AIC favours the model (6) with intensity proportional to slope.

# Simulation of fitted model

A fitted Poisson model can be simulated automatically using the simulate-function:

- > X <- simulate(fit)</pre>
- > plot(X, main = "")

#### Independence, regularity and clustering

Suppose that a point pattern appears to have constant intensity, and we wish to assess whether the pattern is Poisson. The alternative is that the point pattern is either regular or clustered.





The *Fry plot* is a scatterplot of the vector differences  $x_j - x_i$  between all pairs of distinct points in the pattern. It is computed by the command fryplot.



### Distances

Classical techniques for investigating interpoint interaction are *distance methods*, based on measuring the distances between points. Specifically we may consider

- **pairwise distances**  $s_{ij} = ||x_i x_j||$  between all distinct pairs of points  $x_i$  and  $x_j$  ( $i \neq j$ ) in the pattern;
- nearest neighbour distances t<sub>i</sub> = min<sub>j≠i</sub> s<sub>ij</sub>, the distance from each point x<sub>i</sub> to its nearest neighbour;
- empty space distances  $d(u) = \min_i ||u x_i||$ , the distance from a fixed reference location u in the window to the nearest data point.

These are availble in spatstat as the functions pairdist, nndist and distmap respectively.

# Empty space distances

- > data(swedishpines)
- > emp <- distmap(swedishpines)</pre>
- > plot(emp, main = "Empty space distances")
- > plot(swedishpines, add = TRUE)

Empty space distances



# Edge effects

Confining observations to a window W implies that the observed distance  $d(u, \mathbf{x}) = d(u, \mathbf{X} \cap W)$  to the nearest data point inside W, may be greater than the true distance  $d(u, \mathbf{X})$  to the nearest point of the complete point process  $\mathbf{X}$ .



# Empty space function F

Assuming X is stationary (statistically invariant under translations), we can define the cumulative distribution function of the empty space distance

$$F(r) = \mathbb{P}\left\{d(u, \mathbf{X}) \le r\right\}$$
(8)

where u is an arbitrary reference location. If the process is stationary then this definition does not depend on u.

The empirical cumulative distribution function (ecdf) of the observed empty space distances on a grid of locations u<sub>j</sub>, j = 1, ..., m,

$$F^*(r) = \frac{1}{m} \sum_j \mathbf{1} \left\{ d(u_j, \mathbf{x}) \le r \right\}$$
(9)

is a negatively biased estimator of F(r), as explained above.

# Edge-correction

 Typically edge-corrected estimates of F are weighted versions of the ecdf,

$$\widehat{F}(r) = \sum_{j} e(u_j, r) \mathbf{1} \{ d(u_j, \mathbf{x}) \le r \}$$
(10)

where e(u, r) is an edge correction weight designed so that  $\widehat{F}(r)$  is unbiased.

- The edge effect problem can also be regarded as a form of censoring (analogous to right-censoring in survival data), and a counterpart of the Kaplan-Meier estimator is available.
- Thus, assuming that the point process is homogeneous, we are able to compute an unbiased and reasonably accurate estimate of the empty space function F defined by (8).

#### Empty space function: Poisson case

Notice that d(u, X) > r if and only if there are no points of X in the disc b(u, r) of radius r centred on u. For a homogeneous Poisson process with intensity λ the probability that there are no points in this region is exp(-λπr<sup>2</sup>). Hence for a Poisson process

$$F_{\rho}(r) = 1 - \exp(-\lambda \pi r^2).$$
 (11)

■ Typically we compare *F*(*r*) with the value of *F<sub>p</sub>*(*r*) obtained by plugging in the estimated intensity *λ̂* = *n*(**x**)/area(*W*). Values *F̂*(*r*) > *F<sub>p</sub>*(*r*) suggest that empty space distances in the point pattern are shorter than for a Poisson process, suggesting a regularly space pattern; while values *F̂*(*r*) < *F<sub>p</sub>*(*r*) suggest a clustered pattern.

### The empty space function in spatstat

The function Fest computes estimates of F(r) using several edge corrections, and the benchmark value for the Poisson process.

```
> Fswed <- Fest(swedishpines)</pre>
```

> Fswed

```
Function value object (class \hat{a}\tilde{A}\tilde{Y}fv\tilde{a}\tilde{A}\tilde{Z})
for the function r -> F(r)
```

Entries:

id	label	description
r	r	distance argument r
theo	F[pois](r)	theoretical Poisson F(r)
cs	F[cs](r)	Chiu-Stoyan estimate of F(r)
rs	F[bord](r)	border corrected estimate of F(r)
km	F[km](r)	Kaplan-Meier estimate of F(r)
hazard	hazard(r)	Kaplan-Meier estimate of hazard funct

#### Plotting the empty space function

- > plot(Fswed)
- > plot(Fswed, cbind(km, rs, theo) ~ theo)



#### Plotting the empty space function

- > plot(Fswed, . ~ theo)
- > plot(Fswed, . theo ~ r)



# Thank your for your attention.

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