Appendix

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Simulation of a HPP

Recall that for a HPP, the interarrival times between events, W, are exponentially distributed. In this simulation of a HPP, after initializing the initial time, t, and the time vector, t_{vector} , we generate exponential random variables and use them to index the interarrival times between events.

Figure 2 as shown previously is a realization of a HPP with rates that are roughly constant at $\lambda = 10$.

Algorithm 1: Simulation of a HPP

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Input \lambda, t_{max}
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- 1. Initialize t, t_{vector}
- 2. while $(t \leq t_{max})$
- 3. Generate $u \sim U(0,1)$
- 4. Set t = t + w where $w = -log(u)/\lambda \sim exp(\lambda^* = \lambda)$
- 5. **if** $(t < t_{max})$
- 6. | Add $t_{vector} = c(t_{vector}, t)$
- 7. else
- 8. | **return** $\{t_k\}_{k=0,1,...}$

Simulation of a Hawkes Process

In this simulation of a Hawkes process, we use the thinning algorithm (or acceptance-rejection method) to simulate a temporal Hawkes process since it is one of the most popular choices for simulating both temporal and spatio-temporal NPP (Pasupathy, 2010). Broadly put, the thinning algorithm involves randomly deleting points from a point pattern. The process requires first simulating a HPP, creating a $\lambda(t)$ function and applying it to the HPP, and using $min(\lambda^*/\lambda, 1)$ as the accepting probability to randomly keep or 'thin' the points.

Figure 5 as shown previously is a realization of a Hawkes process with parameters such that $\mu = 0.5$, $\alpha = 0.5$, and $\beta = 0.7$. μ sets the background rate. α and β controls the shape and decay rate of the the exponentially decaying triggering function function.

Algorithm 2: Simulation of a Hawkes Process via Thinning Algorithm

Imput μ , α , β , λ , t_{max}

- 1. Simulate a HPP using Algorithm 1
- 2. Create a $\lambda(t)$ function where the function $= \mu + \sum_{i:T_i \leq t} \alpha e^{-\beta x}$
- 3. Set $\lambda^* = \text{apply the } \lambda(t)$ function to the HPP
- 4. Generate $u \sim U(0,1)$
- 5. if $(u < min(\frac{\lambda^*}{\lambda}, 1))$ where the accepting probability $= min(\lambda^*/\lambda, 1)$
- 6. | Keep the points
- 7. else
- 8. | "Thin" or reject the points and **return** $\{t_k\}_{k=0,1,...}$

Simulations of a HPP, NPP, Cox and Matern Cluster Process in 2D

All of the corresponding plots in 2D are created using the **spatstat** package of R (Baddeley & Turner, 2005).

HPP

The rpoispp function can be used to generate a random point pattern as a realization of a HPP or NPP. lambda controls the rate (or intensity) of a HPP and win sets the window in which the simulated pattern is observed. In Figure 3, we set lambda = 100 and win = square(1).

NPP

Similar to the above simulation, lambda can be set to control the intensity of a NPP. In the second figure of Figure 3, lambda = 400 * x * y.

Cox Process

Instead of setting lambda to be a deterministic function as in the NPP case, here we can set lambda to be a random function or variable. In Figure 4, lambda is set to be = rexp(n = 1, rate = 1/100).

Matern Cluster Process

Simulations of Matern cluster process are generated using the rMatClust function. Specifically, the process involves generating homogeneous Poisson parents and each parent gives rise to Poisson number of offspring uniformly distributed in a disc of radius r centered around the parent. kappa controls the intensity of the cluster centers and allows us to specify the number of clusters. \mathbf{r} specifies how far away cluster is from one another in radius, and $\mathbf{m}u$ gives the mean number of points per cluster. In the second figure of Figure 4, kappa = 20, r = 0.05, and mu = 5.