

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

Input λ, t_{max}

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** ($t \leq t_{max}$)
 6. | Add $t_{vector} = c(t_{vector}, t)$
 7. **else**
 8. | **return** $\{t_k\}_{k=0,1,\dots}$
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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 $\mu = 0.5, \alpha = 0.7$, and $\beta = 0.5$. 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

Input $\mu, \alpha, \beta, \lambda, t_{max}$

1. Simulate a HPP using Algorithm 1
 2. Create a $\lambda(t)$ function where the function $= \mu + \sum_{i:T_i < t} \alpha e^{-\beta x}$
 3. Set $\lambda^* =$ 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** $\{t_k\}_{k=0,1,\dots}$
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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 3, **lambda** is set to be $= \text{rexp}(n = 1, \text{rate} = 1/100)$.

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. **r** specifies how far away cluster is from one another in radius, and **mu** gives the mean number of points per cluster. In the second figure of Figure 4, $\text{kappa} = 20, r = 0.05$, and $\text{mu} = 5$.