

# Prequel to Hawkes Processes: An Overview of Spatial, Temporal and Spatio-Temporal Point Processes and Some Simulations

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## **Abstract**

Events such as earthquake epicenters, crime patterns, forest wildfires, financial transactions, etc. often exhibit triggering and clustering behavior. The ability to capture events with such behavior gives Hawkes (or self-exciting) processes the potential to become a powerful predictive tool for a wide variety of applications. In this project, we give brief introductions, review definitions, discuss properties and applications of selected spatial and temporal point processes leading up to spatio-temporal self-exciting processes, and simulate some of the processes in 1D and 2D in hope of interested readers have the background knowledge to read and comprehend existing literature as well as explore the field further.

## **1 Introduction**

Real-world data are often spatial, temporal, or spatio-temporal in nature. Spatial data (e.g. soil properties, housing prices) often involve locations such as points and areas. Temporal data (e.g. sensor readings, stock prices) often involve times such as moments and intervals. Spatio-temporal data are data that relates to both locations and times. Examples of spatio-temporal data include forest inventories, remotely sensed images, earthquake epicenters, disease cases, map services and travel times, to name a few.

Various statistical models and methods have been developed for modeling spatial, temporal, or spatio-temporal data, and which models to use depend on questions of interest. For this project, we primarily focus on the point process models. Point process models are useful for describing phenomena that occurs at

random locations, times, or locations and times, and the questions of interests typically are: Does the rate for the occurrence of events vary with locations, times, or locations and times? Do events appear clustered? Do events trigger subsequent events?

Recall that spatial data can be broadly categorized into three types: geostatistical (point process) data, areal data, and point pattern data (Cressie, 2015), we are in the third category in which point pattern data are realizations of spatial point processes. Questions about point pattern data typically are: Is there clustering of events? Can we define a point process that captures the events? Examples of such models include Cox and cluster processes. On the other hand, when dealing with temporal data, *marked* point processes are sometimes used interchangeable with time series, and vice versa (Schoenberg, 2010). One major distinction is that in point processes, time intervals are treated as continuous, whereas in time series, they are treated as discrete. Examples of such models include (temporal) Poisson and Hawkes processes.

Hawkes processes are also known as self-exciting point process. More specifically, the original Hawkes processes are temporal, whereas the more recently developed self-exciting point processes have been extended from temporal Hawkes processes to account for both the spatial and temporal aspects of the data. The defining characteristic of Hawkes processes is that it ‘self-excites’. In other words, the occurrence of an event increases the occurrence of future events nearby in space and/or time, but the events don’t self-excite in perpetuity. In addition, given the history of such events, more recent events exert more influence on the rate at which events occur. For example, in seismology, an event can be an earthquake occurrence that causes aftershocks. In criminology, an event can be a gang rivalry that triggers retaliations following the gang crime. In both cases, the initial event can continue to spawn ‘offspring’ events and the ‘offspring’ events can spawn ‘offspring’ events of their own, but the spawns fade out eventually.

In addition to modeling earthquake epicenters (Ogata, 1988, 1998) and crime patterns (Mohler et al., 2011; Reinhart & Greenhouse, 2018), Hawkes processes have also been used in modeling events such as forest wildfires (Peng, et al., 2005), insurance claims (Stabile et al., 2010), financial transactions (Bauwens and Hautsch, 2009; Embrechts et al., 2011; Bacry et al., 2015), social network events (Zhao et al., 2015; Rizoïu et al., 2017), neuron activities (Johnson, 1996; Gerhard et al., 2017), and disease spread or transmission (Meyer et al., 2012; Meyer & Held, 2014), allowing them to find applications in a wide variety of fields such as emergency and disaster management, insurance, finance, social network, neuroscience, and epidemiology. Recent work has extended the use of self-exciting point processes to novel applications such as mass shootings (Boyd & Molyneux, 2021), COVID-19 transmission (Chiang et al., 2020), and gang violence (Park et al., 2021). However, there is still much work to be done which include computational advances to ease the burden of applying such models to bigger data sets (Holbrook et al., 2021), residual and model diagnostics,

multivariate Hawkes processes, methods that make the models more flexible and applicable, etc.

Given the flexibility and applicability of Hawkes processes, it is surprising to see that Hawkes processes have not gained enough attention from the machine learning communities, which would find their predictive capabilities beneficial. In addition, understanding Hawkes processes would benefit from knowing some of the relevant point processes (e.g. nonhomogeneous Poisson, Cox and cluster processes), which are often left out from graduate-level, introductory spatial statistics and stochastic processes courses. The objective of this project is then to give an overview of various types of point processes so that readers of interest have the necessary background knowledge to understand Hawkes and self-exciting processes.

The outline of this project is as follows: In *Section 2*, we introduce, define and discuss properties and applications of counting processes, homogeneous and nonhomogeneous Poisson processes, Cox processes, cluster processes, Hawkes processes, and spatio-temporal self-exciting process. In *Section 3*, we wrap up the aforementioned point processes and discuss future work of Hawkes and self-exciting processes. In *Appendix*, we discuss in particular the thinning algorithm (acceptance-rejection method) and **spatstat** package of R that are used to simulate selected processes in 1D and 2D, respectively.

## 2 Introductions, Definitions and Properties

### 2.1 Counting Process

A counting process counts the occurrence (or number) of events over time, space, space-time or, in the most general sense, any metric space in which events occur and can be counted (Daley & Vere-Jones, 2003).

If we were to denote the time of arrival for customers at a super market, we would have a set of points in time in which we could count the number customers over some interval of time. On the other hand, if we consider the location of trees to occur at a point in space, then in some bounded region of the space, we could count the number of trees. An earthquake's epicenter would be a point in space but we can also capture when the point appears in time.

Let us restrict ourselves to the temporal domain so that interpretations are easier to follow. The counting process requires that 1) the number of events  $N(t)$  up to some time  $t$  to be greater than zero, 2) the number of events must be an integer, 3) the counts always increase, and 4) the number of events in specific time interval can be obtained by subtracting the number of events in previous interval from that in current interval. In addition, counting processes are independent, stationary, and homogeneous. In other words, the number of

events  $N(t)$  occurring in disjoint interval  $t$  are independent, the distribution of the number of events depends only on the length of the interval  $t$ , and the transition probability (i.e. the probability moving from one state to another state) between any two states at two times depends only on the difference between the states.

Before formally defining a counting process, first we need to define stochastic processes and point processes. A stochastic process is a collection of random variables indexed by time,  $t$ , space,  $s$  or space-time,  $(s, t)$ , but we restrict ourselves to the time domain again and follow with Obral (2016) to define a stochastic process and a point process

**Definition 2.1.1** (Stochastic Process) A stochastic process is a family of random variables indexed by time  $t$  and is defined as

$$\{X_t\}_{t \in T}.$$

**Definition 2.1.2** (Point Process) Let  $\{T_i\}_{i \in \mathbb{N}}$  be a sequence of non-negative random variables such that  $T_i < T_{i+1} \forall i \in \mathbb{N}$ , a point process on  $R^+$  is defined as

$$\{T_i\}_{i \in \mathbb{N}}.$$

A point process relies on the occurrence of an event occurring at a specific point in time. Stochastic processes, on the other hand, are more general. It can be related to a time interval, a waiting time, a state (e.g. blue or red) that changes over time, etc.

A counting process is then defined as follows

**Definition 2.1.3** (Counting Process) Let  $N(t)$  be the number of events up to some time  $t$  such that the values are nonnegative, integer valued, and nondecreasing, a stochastic process is said to be a counting process and is defined as

$$\{N(t), t \geq 0\}.$$

Let us look at a more explicit example, which we show in Figure 1. Suppose that  $N(t)$  counts the number of events up to some time  $t$  and events occur at times  $t = 0.1, 1, 1.5, 3, 5$ , etc. then  $N(2) = 3$  since events occurring at 0.01, 1, and 1.5 all occur in the time interval  $(0, 2]$ . Similiarly,  $N(4) = 4$  since 4 events occur in the time interval  $(0, 4]$ .

An alternative definition of a counting process is

**Definition 2.1.4** (Counting Process) Let  $\{T_i\}_{i \in \mathbb{N}}$  be a point process, a counting process associated with

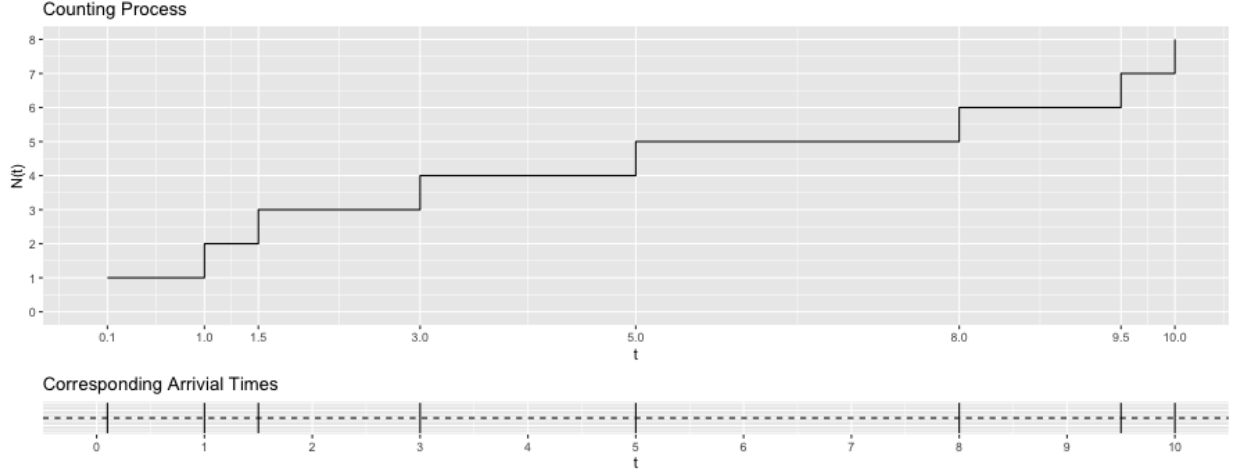


Figure 1: Counting Process

$\{T_i\}_{i \in \mathbb{N}}$  is defined as

$$N(t) = \sum_{i \in \mathbb{N}} I_{\{T_i \leq t\}}.$$

This equivalent definition may be more beneficial as we get into Poisson, cluster, etc. processes in the later sections because it is easier to see that for  $i \in \mathbb{N}$ , if  $T_i \leq t$ , then the indicator function  $I_{\{T_i \leq t\}}$  is equal to 1. Then, we sum up all the 1s for events which have occurred.

A useful corollary of counting process, which describes more formally some of the properties we stated above and will be helpful to understand as we move into Poisson processes and then beyond is

*Corollary 2.1.1* A counting process satisfies that

1.  $N(t) \geq 0$
2.  $N(t)$  is an integer
3. If  $t \leq t + h$ , then  $N(t) \leq N(t + h)$
4. If  $t < t + h$ , then  $N(t + h) - N(t)$  is the number of events occur in the interval  $(t, t + h]$ .

In other words, 1. An event has to occur for it to be counted. 2. We either count an event or we don't. There is no event that sort of occurs that results in decimal value. 3. Counts always increase because events don't disappear. Once we observe an event and count it, it remains in the counts. 4. The number of events in specific time interval can be obtained by subtracting the number of events in previous interval from that in the current interval.

## 2.2 Poisson Process

The homogeneous Poisson process (HPP) is one of the simplest yet most-widely used point processes (Baddeley et al., 2015). HPPs can be used to model the number of events such as bus arrivals at a bus stop, car accidents at a site, or the document requests on a web server over time. As we alluded to previously with counting processes, HPPs can also be considered over a space which is often taken to be a two-dimensional plane, such as the surface of Earth, or a three-dimensional volume, such as the interior of Earth.

Like counting processes, HPPs are also independent, stationary, and homogeneous. In addition, we assume that the numbers of events  $N(t)$  follow a Poisson distribution with a constant rate,  $\lambda$ , and the interarrival times between events,  $W$ , are exponentially distributed. We follow with Obral (2016) and Chen (2016) to formally define a Poisson process

**Definition 2.2.1** (Poisson Process) If the following conditions hold, a counting process  $\{N(t), t \geq 0\}$  is said to be a Poisson Process with constant rate (or intensity)  $\lambda > 0$

1.  $N(0) = 0$
2.  $N(t)$  has independent increments
3.  $P(N(t+h) - N(t) = 1) = \lambda h + o(h)$
4.  $P(N(t+h) - N(t) > 1) = o(h)$ ,

where the function of little  $o$   $o(h)$  is given as

$$\lim_{h \rightarrow 0^+} \frac{o(h)}{h} = 0.$$

In other words, 1. An event has to occur for it to be counted. 2. For any disjoint time intervals, the occurrence of an event does not affect the probability of the occurrence of one another event. 3.  $\lambda$  is the rate (i.e. events over time) at which points occur and is constant. 4. No more than 1 event can occur at the same location.

An alternative way to think of HPP is that it is a uniformly random process. If we were to take a realization of a HPP over some time interval  $(0, T]$  and ‘bin’ the number of events occurring in some set of equal intervals, then the histogram for the realization would resemble a realization of a uniform distribution over time 0 to  $T$ .



Figure 2: Homogeneous Poisson process (rate = 10)

Let us look at a realization of a HPP in time, which we show in Figure 2. First, we note that the cumulative number of points is growing at a constant linear rate. In addition, we can see that the histogram of rate appears roughly uniform; the rates are roughly constant at  $\lambda = 10$ . The algorithm used for simulating this HPP can be found in the Appendix section.

For an HPP, the numbers of events in any time interval  $N(t)$  are Poisson distributed. More formally, we can say that the number of events in any time interval  $(t, t + h]$ ,  $N((t, t + h])$ ,  $\sim \text{Pos}(\lambda \cdot h)$ . That is, for all  $t, h \geq 0$  and  $n = 0, 1, \dots$ ,

$$P(N(t + h) - N(t) = n) = \frac{(\lambda h)^n e^{-\lambda h}}{n!}.$$

Additionally, let us denote the interarrival times between events as  $W$ . For example, let  $T_0$  be the starting time of the process while  $T_1$  the time of the first occurrence of event and  $T_2$  the time of the second occurrence of event, then the elapsed time between the start of the process and first event is  $W_1$  and the elapsed time between the first event and second event is  $W_2$ . The interarrival times  $W$  are exponentially distributed. More formally, the interarrival times  $W$ ,  $\sim \exp(\frac{1}{\lambda})$ . That is, for rate  $\lambda > 0$ , the interarrival time  $W_i$   $i = 1, 2, \dots$ ,

$$P(W_1 > h) = P(N(h) = 0) = e^{-\lambda h}.$$

This is because the probability of the first point arriving after time  $h$  can be thought of as the probability that the first point does *not* arrive in the time interval  $(0, h]$ . Similarly,  $W_2$  is also  $\sim \exp(\frac{1}{\lambda})$  since

$$P(W_2 > h | W_1 = t) = P(N(t+h) - N(t) | N(t) - N(t^-) = 1) = P(N(t+h) - N(t) = 0) = P(N(h) = 0) = e^{-\lambda h}.$$

## 2.3 Nonhomogeneous Poisson Process

Assuming that the rate in which points occur is constant is often not realistic in practice. We may want a model that allows for more flexibility. The nonhomogeneous Poisson processes (NPPs) are a generalization of homogeneous Poisson processes that allow for the rate (or intensity)  $\lambda$  to vary as function of time  $t$  or space  $s$ .

We previously assumed that for the HPP the intensity  $\lambda$  is constant. If we have reasons to believe that the intensity is not constant, then we should model as a NPP instead. This would be the case if, as in the supermarket example, we have reasons to believe that the arrival rate of customers is higher during lunch time as compared to say, 2am, or as in the trees in a forest example, we speculate that environmental factors such as temperature, rainfall and light affect the spatial distribution of the trees.

Contrary then to the HPPs, NPPs are independent but not stationary nor homogeneous. Recall that HPP has stationary increments since the distribution of the numbers of events  $N(t)$  that occur in any interval of time  $t$  depends only on the length of the interval  $t$  but not the location of the interval  $t$ , NPP, in contrast, does not have stationary increments since the distribution of  $N(t)$  can change when shifted in  $t$ . Since stationary implies homogeneity, NPP is nonhomogeneous.

For HPP, we assume that the numbers of events in any time interval  $N(t)$  follow a Poisson distribution with a constant intensity  $\lambda$ . For NPP, we assume that  $N(t)$  follow a Poisson distribution too but with an intensity function  $\lambda(t)$  such that the intensity now varies with a function of time. This then leads to the following definition of a NPP (Obral, 2016; Chen, 2016)

**Definition 2.3.1** (Nonhomogeneous Poisson Process) If the following conditions hold, a counting process  $\{N(t), t \geq 0\}$  is said to be a nonhomogeneous Poisson Process with intensity function of time  $\lambda(t), t > 0$

1.  $N(0) = 0$
2.  $N(t)$  has independent increments
3.  $P(N(t+h) - N(t) = 1) = \lambda(t)h + o(h)$
4.  $P(N(t+h) - N(t) > 1) = o(h)$ .

NPPs have additional properties such as if the number of events in any time interval  $(t, t+h]$ , denoted as  $N((t, t+h])$ ,  $\sim \text{Pos}(\Lambda(t) = \int_t^{t+h} \lambda(v)dv)$ . That is, for all  $v, t, h \geq 0$  and  $n = 0, 1, \dots$ ,

$$P(N(t+h) - N(t) = n) = \frac{(\int_t^{t+h} \lambda(v)dv)^n e^{-\int_t^{t+h} \lambda(v)dv}}{n!},$$



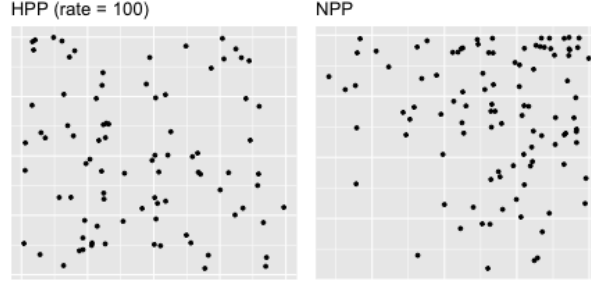


Figure 3: Left: HPP (rate = 100) Right: NPP (intensity =  $400xy$ )

where  $\lambda(v)$  again denotes a non-constant rate function.

Further, occurrence of the next point can be determined by utilizing the exponential distribution with

$$P(N(t, t+h] = 0) = e^{-\int_t^{t+h} \lambda(v) dv}.$$

## A Motivating Example

Before we delve further into other point processes, let us look at a motivating example, which we show in Figure 3. We demonstrate HPP and NPP in space so that visualizations are easier to look at and comprehend.

HPP in space is also called complete spatial randomness (CSR) (Baddeley et al., 2015). For HPP in space, the number of events in  $u$  with area  $|u|$ , denoted as  $N(u)$ ,  $\sim \text{Pos}(\lambda|u|)$ . The left figure of Figure 3 is a realization of a HPP with constant *rate* = 100. HPP points appear uniformly distributed in  $u$ .

For NPP in space, the number of events in  $u$ , denoted as  $N(u)$ ,  $\sim \text{Pos}(\int_u \lambda(v) dv)$ . The right figure of Figure 3 is a realization of a NPP with *intensity function* =  $400xy$ . NPP points are not uniformly distributed; they are distributed according to the intensity function of the process. In this example, the points appear to concentrate at the upper-right corner.

## 2.4 Cox and Cluster Process

Even more flexible models than NPP are Cox and cluster processes. Whereas, previously assumed independence between events which occur at a constant rate  $\lambda$  for the HPP or were independent but depend on an intensity function  $\lambda(t)$  for the NPP, we now discuss models that allow for the relaxation of this independence assumption. Additionally, Cox and cluster processes are mainly spatial processes so we are in the space domain.

Examples that are potentially better modelled as Cox processes than HPP include locations of emergent bramble cane (blackberry) plants and *Beilschmiedia* trees (Baddeley et al., 2015). In these examples, there appears to be some observable or unobservable spatial covariate (e.g. light, humidity, soil quality) or external factor that makes it more likely to observe the plants or trees preferentially in some areas as opposed to others.

Cox process (or doubly stochastic Poisson processes) can be defined as a Poisson process with a random intensity function; if the intensity surface  $\Lambda(u)$  is known, then it becomes a Poisson process with intensity function  $\Lambda(u)$  (Baddeley et al., 2015). This spatially varying intensity function of the Cox process is random since it relies on a set of random variables called random field to capture the covariates or external factors (Baddeley et al., 2015). When the intensity is relatively high in some area of space, points occur more frequently. When the intensity is low, fewer points occur. Since the intensity is random and changes based on a set of observed or potentially unobserved set of random variables, this accumulation of points sometimes appears to be clustered; this is one way to observe the dependence structure of the Cox process. Within the general class of Cox processes, there are also numerous specific models. Mixed Poisson process, the simplest example of Cox processes, involves generating a random variable  $\Lambda$  and, given the value of  $\Lambda$ , generating a Poisson process with intensity function  $\Lambda$ .

On the other hand, example that can be modelled as cluster processes includes seedlings and saplings of California redwood (Baddeley et al., 2015). In this example, unobserved ‘parent’ trees give rise to clusters of observed ‘offspring’ trees. It is also not hard to notice significant overlap between Cox processes and cluster processes since these processes tend to share a considerable amount of information in their constructions. Indeed, bramble canes can be modelled as cluster processes as well. While Cox processes can sometimes appear to exhibit clustering of points due to the underlying random field, cluster processes tend to make the clustering of points more explicit. For a cluster process, we first have some set of unobserved ‘parent’ points  $\mathbf{Y}$  generated by the process. Next, each ‘parent’ point  $y_i \in \mathbf{Y}$  gives rise to a random number of ‘offspring’ points  $z_{ij}$ . These ‘offspring’ points  $Z_{ij}$  then form a cluster process  $\mathbf{X}$  around the set of parents  $\mathbf{Y}$  and only the offspring points are observed. Within the general class of cluster processes, there are also numerous specific models which can be utilized depending on the choice of assumptions. Matern cluster processes, for one example, 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.

Alternatively, we can think of Cox process as a hierarchical model with two levels and cluster process as a hierarchical model with three levels (Geyer, 2020). For the Cox process, 1) there is some set of random variables that influences the intensity function and 2) based on this intensity function, we observe the set of

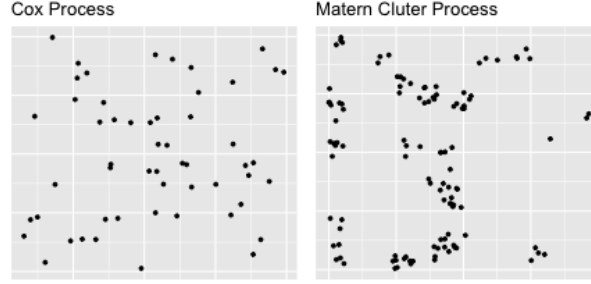


Figure 4: Left: Cox (intensity =  $\exp(n = 1, \text{rate} = 1/100)$ ) Right: Matern ( $\text{kappa} = 20, r = 0.05, \mu = 5$ )

points. For the cluster process, 1) there is some intensity function which can be random or not, 2) based on the intensity function, some set of ‘parent’ (or ‘center’) points are laid down which we often don’t observe, but 3) based on the location of the ‘parent’ points, some set of ‘offspring’ points are generated around the ‘parent’ points which we observe.

The left and right figure of Figure 4 is a realization of a mixed Poisson process with *intensity function* =  $\exp(1, 1/100)$  and a Matern cluster process with  $\text{kappa} = 20, r = 0.05, \mu = 5$ , respectively. Both Cox process and Matern cluster process points appear clustered, but the way the points cluster differ. Points in Cox process cluster accordingly to some specified distribution (i.e. a exponential distribution), whereas points in cluster process cluster in some defined area (i.e. a disc). The **spatstat** package of **R** used for simulating this HPP can be found in the Appendix section (Baddeley & Turner, 2005).

Having given some background on both the Cox and cluster processes, we follow with Coeurjolly (2015) and Daley & Vere-Jones (2003) now with their more formal definitions

**Definition 2.4.1** (Cox Process) Let  $\Lambda = (\Lambda(u))_{u \in S \subseteq \mathbb{R}^d}$  be a non-negative random field such that the values of  $\Lambda(u)$  is nonnegative and  $\Lambda(u)$  is a locally integrable function, where  $\Lambda$  is a random field means that  $\Lambda(u)$  is a random variable  $\forall u \in S$  and  $\Lambda(u)$  is a locally integrable function means that  $E(\Lambda(u))$  exists and is locally integrable with probability 1, if  $X \mid \Lambda \sim \text{Pos}(\Lambda)$ , then  $X$  is said to be a Cox process driven by  $\Lambda$  with intensity function  $\lambda(u) = E(\Lambda(u))$ . That is,

$$P(N(u) = n) = \frac{(\lambda(u))^n e^{-\lambda(u)}}{n!} = \frac{(E(\Lambda(u)))^n e^{-E(\Lambda(u))}}{n!},$$

where

$$\Lambda(u) \stackrel{\text{a.s.}}{=} \int_u^\infty \lambda(x) dx = \int_0^\infty \frac{x^n e^{-x} F_u(dx)}{n!}.$$

**Definition 2.4.2** (Cluster Process) Let  $x$  (‘parent’ points) be points in a point process  $N$  and replacing every

$x$  with a cluster of points  $N_x$  (‘offspring’ point) centered at  $x$  and assume also that each  $N_x$  is independent of one another, then the union of all the clusters forms a cluster process  $N_c$ . That is,

$$N_c = \bigcup_{x \in N} N_x.$$

Next, we follow with Baddeley et al. (2015) to state the following model assumptions for cluster processes

1. ‘Parent’ points follow a Poisson distribution.
2. Clusters are independent of one another.
3. Clusters are identically distributed, which means that clusters, when shifted, have the same distributions.
4. The locations of ‘offspring’ points of each parent point are independently and identically distributed.
5. The number of ‘offspring’ points of each parent point follows a Poisson distribution.
6. Clusters are isotropic, which means that the distribution of ‘offspring’ points for each parent point depends only on the distance between the ‘parent’ and the ‘offspring’.

Under assumption 1 - 4, we have a Neyman-Scott process. Under assumption 1 - 5, the cluster process is a Cox process. And finally, under assumption 1 - 6, we have a Matern or Thomas cluster process.

## 2.5 Hawkes Process

The final class of point process models we present are the Hawkes processes. Hawkes processes are also commonly known as self-exciting point processes. Like Cox and cluster process, the Hawkes process model allows for dependence between events; however, their dependence differs. In Hawkes processes, the occurrence rate of the events depends not only on time  $t$  but also past events  $\mathcal{H}_t^N$  up to some time  $t$ . This is the distinguishing feature of the Hawkes process as neither Cox nor cluster processes depends on the past history of events, which is also what makes Hawkes processes ‘self-excite’. It should also be noted that Hawkes processes can be powerful predictive models as they naturally capture triggering and clustering behavior (Reinhart, 2018).

As previously mentioned, examples of applications for Hawkes processes include locations of earthquake epicenters, locations of crimes, and locations of patients with a communicable disease. The major sharing

feature in each of these examples is that an occurrence of an event leads to an increase in the occurrence of subsequent events (i.e. ‘self-exciting’) in nearby time and/or space. In seismology, for example, a main quake is often followed by aftershocks. In gang crimes, one act of violence often features an act of retaliatory violence. In communicable diseases, one patient may infect others.

Since the intensity (or the expected rate in which points occur over some time and/or space) of the process is now as a function of past history, we refer to the expected rate as a conditional intensity function. Hawkes processes self-excite because of this conditional intensity function since conditioning on the past set of events, the expected rates at which points occur are expected to be higher when points have occurred in nearby time and/or space and then they gradually decline as the events get further away. This ‘self-exciting’ characteristic is captured by the summing part of the conditional intensity function, which we show in Definition 2.5.2.

Another characteristic is that more recent events exert more influence on the intensity but the intensity will fade out or decay until the next event. This ‘decaying’ characteristic is captured by the triggering part of the conditional intensity function, which again we show in Definition 2.5.2. The triggering function is also what makes the Hawkes process a self-exciting as well as cluster process as it allows for additional points to occur in nearby time and/or space before ultimately decaying back to the background rate. For example, an earthquake happens today may lead us believe that there is a higher chance of another earthquake that will occur tomorrow. However, an earthquake happened more than a year ago does not make us believe that there is a higher probability of another earthquake that will occur tomorrow.

A Hawkes process can be uniquely identified through its conditional intensity function (Daley & Vere-Jones, 2003), the formal definition of which we give below

**Definition 2.5.1** (Conditional Intensity Function) Let  $N(t)$  be the numbers of events  $N(t)$  that occur in any interval of time  $t$ , the conditional intensity function  $\lambda(t)$  with respect to the history of the process up to time  $t$ ,  $\mathcal{H}_t$ , is defined as

$$\lambda(t|\mathcal{H}_t) = \lim_{h \rightarrow 0^+} \frac{E(N(t, t+h)|\mathcal{H}_t)}{h},$$

where  $\mathcal{H}_t$  is the history prior to time  $t$ .

Having defined the conditional intensity function, we next more formally define the Hawkes process (Rizoiu, 2017; Reinhart, 2018)

**Definition 2.5.2** (Hawkes Process) A counting process  $\{N(t), t \geq 0\}$  associated with past events  $\{\mathcal{H}_t^N, t > 0\}$

is said to be a Hawkes process with conditional intensity function  $\lambda(t|\mathcal{H}_t^N), t > 0$  and takes the form

$$\lambda(t|\mathcal{H}_t^N) = \lambda_0(t) + \sum_{i:T_i < t} \phi(t - T_i),$$

where  $\lambda_0(t)$  is the base intensity function (or  $\mu$  the constant background rate),  $T_i < t$  are the events time occur before current time  $t$ ,  $\phi(\cdot)$  is the kernel function (or  $g(\cdot)$  the triggering function) through which intensity function depends on past events, and  $\mathcal{H}_t^N$  is the natural filtration (or simply  $\mathcal{H}_t$  the past history) which represents the internal history of  $N$  up to time  $t$ .

Typical choices of  $\phi(\cdot)$  include, for example, exponentially decaying function and power-law kernel (Rizoiu, 2017) and they take the forms of

$$\phi(x) = \alpha e^{-\beta x}$$

and

$$\phi(x) = \frac{\alpha}{(x + \beta)^{\eta+1}}.$$

Exponentially decaying function, which decays faster, has been applied to financial data (Embrechts et al., 2011). On the other hand, power-law kernel, which has a flatter tail, has been used in seismology (Ozaki, 1979) and social media (Rizoiu et al., 2017).

In Figure 5, we show a realization of a Hawkes process with the exponentially decaying triggering function ( $\mu = 0.5, \alpha = 0.7, \beta = 0.5$ ). We can see that each time the event arrives, the intensity (conditional intensity function) increases and then it declines back to the background rate. When the next event arrives, the intensity jumps again and then declines. The intensity for this realization gradually increases over time too since the points of this simulation occur fairly regularly. The algorithm used for simulating this Hawkes process can be found in the Appendix section.

## 2.6 Spatio-Temporal Self-Exciting Process

Spatio-temporal self-exciting processes are an extension of temporal Hawkes processes. Recall that the conditional intensity function for temporal Hawkes processes take the form of

$$\lambda(t|\mathcal{H}_t) = \mu + \sum_{i:T_i < t} g(t - t_i),$$

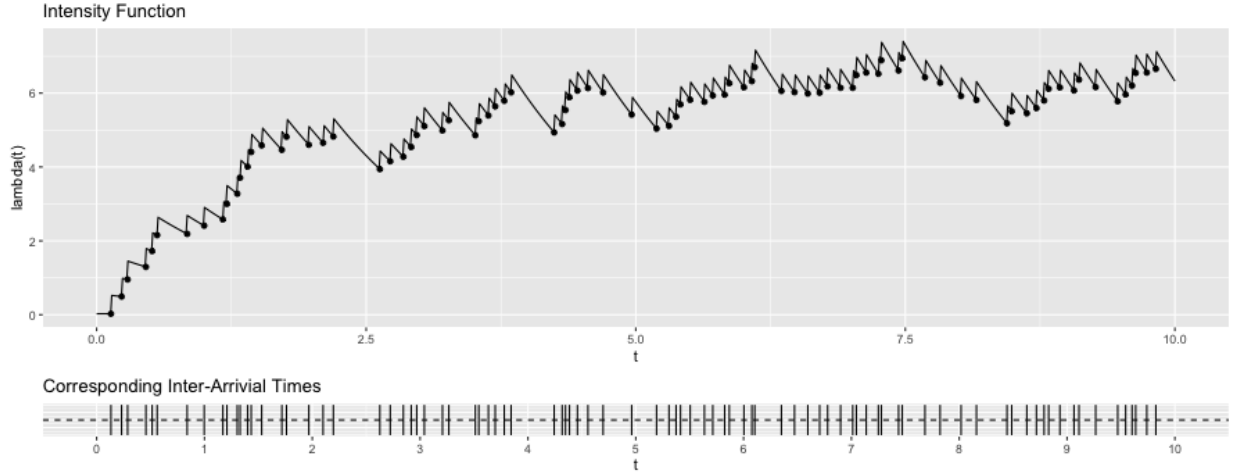


Figure 5: Hawkes Process

the conditional intensity function for spatio-temporal Hawkes processes take the form of

$$\lambda(t|\mathcal{H}_t) = \mu(s) + \sum_{i:T_i < t} g(s - s_i, t - t_i),$$

where  $s_i, i = 1, 2, \dots$  are the sequence of locations of events and  $t_i, i = 1, 2, \dots$  are the times of events. The triggering function is defined to be separable in space and time for simplicity. In addition, the background rate here can be taken to be a function of space, which is often a NPP or constant. Further, the triggering function here is now a function of space and time and, as a result, events that are closer to some given spatial and temporal location are more heavily influenced by the points that occur nearby in both time and space.

### 3 Conclusions and Discussion

In this project, we first introduce the counting process, which can be used to count the number of events over space, time, etc. Next, we introduce HPP, which is one of the simplest yet most widely used point processes for modeling the occurrence of events with constant arrival rate. Then, we relax the stationarity and homogeneity assumption of HPP to move on to NPP, which allows the intensity to vary with space or time. This allows the model to be more realistic in practice. We follow up with HPP and NPP in space for the purpose of visualization and comparison. Then, we further relax the independence assumption and introduce the Cox and cluster processes. Allowing the events to be dependent makes the models even more flexible, as a lot of these environmental and ecological applications rely on the dependence on covariates and/or external factors. Finally, we get to the Hawkes and self-exciting processes. With the intensity now depending on past

history, the models become a suitable choice for capturing events that exhibit both triggering and clustering behavior. Such models also have the potential to become a powerful predictive tool for a wide variety of applications. At this point, interested readers should have the background knowledge to comprehend Hawkes and self-exciting processes through additional, past and recent literature and ponder upon and explore areas for future work such as applications for larger dataset, residual and model diagnostics, etc. On my end, I really enjoy the month-long journey of learning the various point processes as well as witnessing their many practical and important uses for real-world data.

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