

# ECMM450 Stochastic Processes

## Simulation of Non-Homogeneous Poisson Processes

70054986

Department of Computer Science, University of Exeter  
(Dated: April 7, 2023)

This project is about Non-Homogeneous Poisson processes and how to simulate them. We will be reviewing the thinning algorithm of Lewis and Shedler (1979) for simulating NHPP.

### I. INTRODUCTION

This report will be loosely structured according to the following points:

1. Explanation of what is meant by a Non-Homogeneous Poisson Process (NHPP) accompanied by clear mathematical definition.
2. Review of the thinning algorithm of Lewis and Shedler (1979) for simulating NHPP. Short description of the algorithm explaining briefly why it works, and its main benefits compared to other approaches.
3. Simulation of occurrence of 1000 successive events  $t_1, t_2, \dots, t_{1000}$  for a homogeneous Poisson process having a rate of 8 events per year. Visual representation of the same.
4. Considering a NHPP that has a rate function that increases smoothly from 1 event per year at  $t = 0$  by 1% per year, i.e.  $\lambda(t) = (1.01)^t$ . Determining whether the thinning algorithm can be used to simulate this process from the previous homogeneous Poisson process data.
5. Code to perform the thinning algorithm and use it to find occurrence times for a NHPP having the rate function  $\lambda(t) = (1.01)^t$ .
6. Make a figure showing  $N(t)$  versus  $t$  for your NHPP simulation and compare it to what was shown in the figure for the homogeneous Poisson process. By integrating the rate function, add a line to your figure showing the expectation  $E[N(t)]$  versus  $t$ .

### II. DESCRIPTION AND MATHEMATICAL DEFINITION

The following section details what is meant by a Non-Homogeneous Poisson Process (NHPP) giving a clear mathematical definition.

A non-homogeneous Poisson process can be thought of as a generalization of the homogeneous Poisson process, in that, as opposed to its homogeneous counterpart where rate of occurrence of events is constant (denoted by  $\lambda$ ), here the rate is a function of time, denoted by  $\lambda(t)$ . Thus,

the number of occurrences in the interval  $(0, T]$  follows Poisson distribution  $Pois(\int_0^T \lambda(t) dt)$ . More formally, allowing the rate parameter to vary with time results in the following definition.

**Definition II.1** (Ross, 2009, p.339, Definition 5.4). The counting process  $N(t), t \geq 0$  is said to be a nonhomogeneous Poisson process with intensity function  $\lambda(t) \geq 0, t \geq 0$ , if

1.  $N(0) = 0$ .
2. The process has independent increments.
3.  $PN(t+h) - N(h) = 1 = \lambda(t)h + o(h)$ .
4.  $PN(t+h) - N(h) \geq 2 = o(h)$ .

where  $o(h)$  denotes higher order terms of  $h$

Time sampling an ordinary Poisson process (with constant rate  $\lambda$ ) results in a Non-Homogeneous Poisson process. Given  $\{N(t), t \geq 0\}$ , a Poisson process with rate  $\lambda$ , if the event occurring at time  $t$  is counted with probability  $p(t)$ , then  $\{N_c(t), t \geq 0\}$  is a Non-Homogeneous Poisson process. <sup>[Insert reference]</sup>

### III. REVIEW OF THE THINNING ALGORITHM

#### 1. Some other other algorithms

There exist other algorithms for simulating Poisson processes, some of which we will discuss now.

1. Time-scale transformation of a homogeneous Poisson process via inverse of the integrated rate function  $\Lambda(x)$
2. Generate intervals between the points individually
3. Order statistics from Poisson variates
4. Log-linear rate function

#### 2. The thinning algorithm

To construct a Non-Homogeneous Poisson process  $\{N(t), t \geq 0\}$ , with rate parameter  $\lambda(t)$ , over the interval  $(0, T]$ , the algorithm starts with a Non-Homogeneous

Poisson process  $\{N^*(t), t \geq 0\}$ , with rate parameter  $\bar{\lambda}(t)$  that dominates the set  $\lambda(t)$  for all  $t \in (0, T]$ , that is

$$\begin{aligned}\bar{\lambda}(t) &\geq \lambda(t) \forall t \in (0, T] \\ \bar{\lambda}(t) &= \sup_{t \in (0, T]} \lambda(t)\end{aligned}$$

Then, for all  $t$ , the point from the dominating NHPP is retained with probability  $\lambda(t)/\bar{\lambda}$ . The remaining points form a NHPP with rate parameter  $\lambda(t)$ . It is noted that since points are deleted independently, the number of points in  $\{N(x) : x \geq 0\}$  in any set of non-overlapping intervals are mutually independent.

---

**Algorithm 1:** (Lewis and Shedler, 1979, p.7, Algorithm 1) Simulation of an Inhomogeneous Poisson Process with Bounded Intensity Function  $\lambda(t)$ , on  $[0, T]$

---

**Input:**  $\lambda, T$   
Initialize  $n = m = 0, t_0 = s_0 = 0$ ,  
 $\bar{\lambda}(t) = \sup_{t \in (0, T]} \lambda(t)$ ;  
**while**  $s_m < T$  **do**  
    Generate  $u \sim \text{uniform}(0, 1)$ ;  
    Let  $w = -\ln(u)/\bar{\lambda}$ ;  
    Set  $s_{m+1} = s_m + w$ ;  
    Generate  $D \sim \text{uniform}(0, 1)$ ;  
    **if**  $D < \lambda(s_{m+1})/\bar{\lambda}$  **then**  
         $t_{n+1} = s_{m+1}$ ;  
         $n = n + 1$ ;  
    **end**  
     $m = m + 1$   
**end**  
**if**  $t_n \leq T$  **then**  
    **return**  $\{t_k\}_{k=1,2,\dots,n}$   
**else**  
    **return**  $\{t_k\}_{k=1,2,\dots,n-1}$   
**end**

---

### 3. Benefits over other algorithms

The paper discusses a few other methods with which to simulate NHPP. However, each method entails drawbacks with respect to computational efficiency. in its simplest implementation (insert reference), the thinning method obviates the need for numerical integration of the rate function, for ordering of points, and for generation of Poisson variates.

## IV. SIMULATION OF A HOMOGENEOUS POISSON PROCESS

Uniform random numbers are used to generate Poisson variates by using the following algorithm, that can be found in insert reference. Full implementation in Python can be found in the Appendix. Figure 1 shows the result of simulating a homogeneous Poisson process.

---

**Algorithm 2:** Simulation of a Homogeneous Poisson Process with Rate  $\lambda$ , on  $[0, T]$  insert reference to Yuanda Chen

---

**Input:**  $\lambda, N$   
Initialize  $n_0 = 0, t_0 = 0$ ;  
**while** *True* **do**  
    Generate  $u \sim \text{uniform}(0, 1)$ ;  
    Let  $w = -\ln(u)/\bar{\lambda}$ ;  
    Set  $t_{n+1} = t_n + w$ ;  
    **if**  $n + 1 > N$  **then**  
        **return**  $\{t_k\}_{k=1,2,\dots,n}$   
    **else**  
        Set  $n = n + 1$ ;  
    **end**  
**end**

---

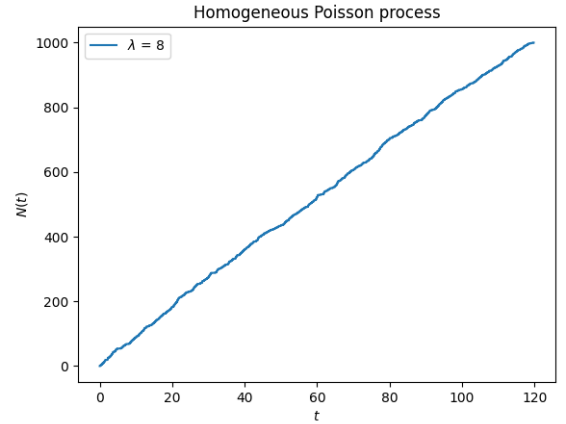


FIG. 1. Simulation of the time of occurrence of 1000 successive events for a homogeneous Poisson process having a rate of 8 events per year.  $N(t)$  denotes the number of events occurring in the time period  $[0, t]$ .

The observation is made that the graph appears relatively linear. This would be expected as the rate of occurrence of events remains constant over any interval of time  $(a, b)$  and hence the slope corresponding to the number of events over time (i.e. rate) also remains the same.

## V. NHPP WITH A SMOOTHLY INCREASING RATE FUNCTION

A NHPP is considered with a rate function that increases smoothly from 1 event per year at  $t = 0$  by 1% per year, i.e.  $\lambda(t) = (1.01)^t$ . Looking at the previous homogeneous Poisson process from IV, it can be seen that the maximum time taken to accumulate 1000 events is 140. Figure 2 shows this. Therefore, the maximum rate for the NHPP would be  $(1.01)^{140} = 4.027$ . Recall that the rate for the previous homogeneous Poisson process was 8. Hence, this is within the bounds of the original process. Thinning can be applied to simulate the NHPP from the

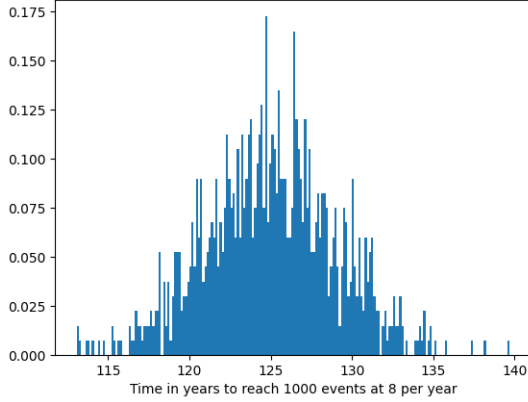


FIG. 2. Time in years to accumulate 1000 events at 8 per year

previous HPP data.

## VI. PERFORM THE THINNING ALGORITHM WITH DATA FROM HOMOGENEOUS POISSON PROCESS

Using the thinning algorithm, the data points from the previous homogeneous process are time-sampled to construct a Non-Homogeneous process.

## VII. GRAPHICAL COMPARISON OF HPP AND NHPP

Figure 3 showcases the cumulative events  $N(t)$  against time  $t$ . It is observed that, as opposed to homogeneous Poisson process, the line is not quite linear - the slope increases as time progresses. This can be expected as the rate parameter  $\lambda(t) = (1.01)^t$  also increases with time. Due to deletion of points from the original homogeneous process, it is also noted that the total number of points has decreased from 1000 to just over 300.

By integrating the rate function, we get

$$\begin{aligned} \int_0^t \lambda(t) dt &= \int_0^t 1.01^t dt \\ &= 1.01^t / \ln(1.01) + C \end{aligned}$$

Noting the boundary condition that at  $t = 0$ ,  $E[N(t)] = 0$ , we have  $C = \frac{1.01^0}{\ln(1.01)} = 100.499$ . Thus,  $E[N(t)] = \frac{1.01^t}{\ln(1.01)} - 100.499$ . Plotting this onto the earlier graph, Figure 4 is obtained. It can be seen that the orange line closely follows the blue stepped graph.

*a. Syntax* The argument of `\cite` may be a single *key*, or may consist of a comma-separated list of keys. The citation *key* may contain letters, numbers, the dash (-) character, or the period (.) character. New with natbib

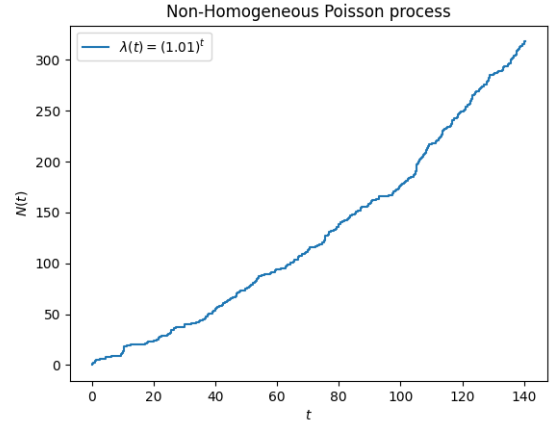


FIG. 3. Accumulated events following a NHPP with rate parameter  $\lambda(t) = 1.01^t$

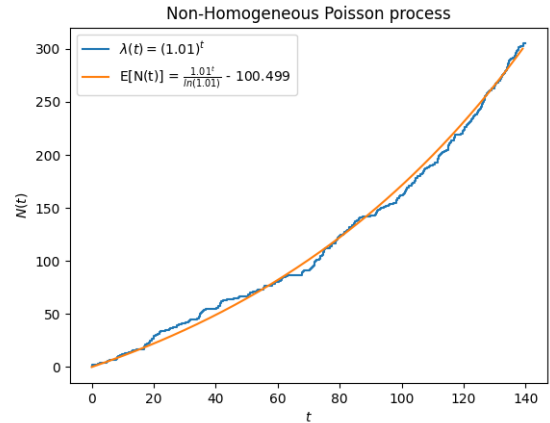


FIG. 4.  $E[N(t)]$ , denoted by  $\frac{1.01^t}{\ln(1.01)} - 100.499$  superimposed on the NHPP

8.3 is an extension to the syntax that allows for a star (\*) form and two optional arguments on the citation key itself. The syntax of the `\cite` command is thus (informally stated)

`\cite { key }, or`  
`\cite { optarg+key }, or`  
`\cite { optarg+key , optarg+key... },`  
 where *optarg+key* signifies

*key*, or  
 \**key*, or  
 [*pre*]*key*, or  
 [*pre*][*post*]*key*, or even  
 \* [*pre*][*post*]*key*.

where *pre* and *post* is whatever text you wish to place at the beginning and end, respectively, of the bibliographic reference (see Ref. [1] and the two under Ref. [2]). (Keep in mind that no automatic space or punctuation is applied.) It is highly recommended that you put the entire

*pre* or *post* portion within its own set of braces, for example: `\cite{[{text}]key}`. The extra set of braces will keep L<sup>A</sup>T<sub>E</sub>X out of trouble if your *text* contains the comma (,) character.

The star (\*) modifier to the *key* signifies that the reference is to be merged with the previous reference into a single bibliographic entry, a common idiom in APS and AIP articles (see below, Ref. [2]). When references are merged in this way, they are separated by a semicolon instead of the period (full stop) that would otherwise appear.

*b. Eliding repeated information* When a reference is merged, some of its fields may be elided: for example, when the author matches that of the previous reference, it is omitted. If both author and journal match, both are omitted. If the journal matches, but the author does not, the journal is replaced by *ibid.*, as exemplified by Ref. [2]. These rules embody common editorial practice in APS and AIP journals and will only be in effect if the markup features of the APS and AIP BibT<sub>E</sub>X styles is employed.

*c. The options of the cite command itself* Please note that optional arguments to the *key* change the reference in the bibliography, not the citation in the body of the document. For the latter, use the optional arguments of the `\cite` command itself: `\cite*[pre-cite][post-cite][key-list]`.

### 1. Example citations

By default, citations are numerical[3]. Author-year citations are used when the journal is RMP. To give a textual citation, use `\onlinecite{#1}`: Refs. 1 and 4. By default, the `natbib` package automatically sorts your citations into numerical order and “compresses” runs of three or more consecutive numerical citations. REVTeX provides the ability to automatically change the punctuation when switching between journal styles that provide citations in square brackets and those that use a superscript style instead. This is done through the `citeautoscript` option. For instance, the journal style `prb` automatically invokes this option because *Physical Review B* uses superscript-style citations. The effect is to move the punctuation, which normally comes after a citation in square brackets, to its proper position before the superscript. To illustrate, we cite several together [1, 2, 4–6], and once again in different order (Refs. [1, 2, 4–6]). Note that the citations were both compressed and sorted. Furthermore, running this sample file under the `prb` option will move the punctuation to the correct place.

When the `prb` class option is used, the `\cite{#1}` command displays the reference’s number as a superscript rather than in square brackets. Note that the location of the `\cite{#1}` command should be adjusted for the reference style: the superscript references in `prb` style must appear after punctuation; otherwise the reference must appear before any punctuation. This sample was written

for the regular (non-`prb`) citation style. The command `\onlinecite{#1}` in the `prb` style also displays the reference on the baseline.

## 2. References

A reference in the bibliography is specified by a `\bibitem{#1}` command with the same argument as the `\cite{#1}` command. `\bibitem{#1}` commands may be crafted by hand or, preferably, generated by BibT<sub>E</sub>X. REVTeX 4.2 includes BibT<sub>E</sub>X style files `apsrev4-2.bst`, `apsrmp4-2.bst` appropriate for *Physical Review* and *Reviews of Modern Physics*, respectively.

### 3. Example references

This sample file employs the `\bibliography` command, which formats the `apssamp.bbl` file and specifies which bibliographic databases are to be used by BibT<sub>E</sub>X (one of these should be by arXiv convention `apssamp.bib`). Running BibT<sub>E</sub>X (via `bibtex apssamp`) after the first pass of L<sup>A</sup>T<sub>E</sub>X produces the file `apssamp.bbl` which contains the automatically formatted `\bibitem` commands (including extra markup information via `\bibinfo` and `\bibfield` commands). If not using BibT<sub>E</sub>X, you will have to create the `thebibliography` environment and its `\bibitem` commands by hand.

Numerous examples of the use of the APS bibliographic entry types appear in the bibliography of this sample document. You can refer to the `apssamp.bib` file, and compare its information to the formatted bibliography itself.

## A. Footnotes

Footnotes, produced using the `\footnote{#1}` command, usually integrated into the bibliography alongside the other entries. Numerical citation styles do this[7]; author-year citation styles place the footnote at the bottom of the text column. Note: due to the method used to place footnotes in the bibliography, *you must re-run BibT<sub>E</sub>X every time you change any of your document’s footnotes.*

## VIII. MATH AND EQUATIONS

Inline math may be typeset using the `$` delimiters. Bold math symbols may be achieved using the `bm` package and the `\bm{#1}` command it supplies. For instance, a bold  $\alpha$  can be typeset as `$\bm{\alpha}$` giving  $\alpha$ . Fraktur and Blackboard (or open face or double struck) characters should be typeset using the `\mathfrak{#1}` and `\mathbb{#1}` commands respectively. Both are supplied

by the `amssymb` package. For example, `\mathbb{R}` gives  $\mathbb{R}$  and `\mathfrak{G}` gives  $\mathfrak{G}$ . In  $\text{\LaTeX}$  there are many different ways to display equations, and a few preferred ways are noted below. Displayed math will center by default. Use the class option `fleqn` to flush equations left.

Below we have numbered single-line equations; this is the most common type of equation in *Physical Review*:

$$\chi_+(p) \lesssim [2|\mathbf{p}|(|\mathbf{p}| + p_z)]^{-1/2} \begin{pmatrix} |\mathbf{p}| + p_z \\ px + ip_y \end{pmatrix}, \quad (1)$$

$$\left\{ 1234567890abc123\alpha\beta\gamma\delta1234556\alpha\beta\frac{1\sum_b^a}{A^2} \right\}. \quad (2)$$

Note the open one in Eq. (2).

Not all numbered equations will fit within a narrow column this way. The equation number will move down automatically if it cannot fit on the same line with a one-line equation:

$$\left\{ ab12345678abc123456abcde\alpha\beta\gamma\delta1234556\alpha\beta\frac{1\sum_b^a}{A^2} \right\}. \quad (3)$$

When the `\label{#1}` command is used [cf. input for Eq. (2)], the equation can be referred to in text without knowing the equation number that  $\text{\TeX}$  will assign to it. Just use `\ref{#1}`, where `#1` is the same name that used in the `\label{#1}` command.

Unnumbered single-line equations can be typeset using the `[, \]` format:

$$g^+g^+ \rightarrow g^+g^+g^+g^+ \dots, \quad q^+q^+ \rightarrow q^+g^+g^+ \dots$$

### A. Multiline equations

Multiline equations are obtained by using the `eqnarray` environment. Use the `\nonumber` command at the end of each line to avoid assigning a number:

$$\begin{aligned} \mathcal{M} = & \quad ig_Z^2(4E_1E_2)^{1/2}(l_i^2)^{-1}\delta_{\sigma_1,-\sigma_2}(g_{\sigma_2}^e)^2\chi_{-\sigma_2}(p_2) \\ & \times [\epsilon_j l_i \epsilon_i]_{\sigma_1} \chi_{\sigma_1}(p_1), \end{aligned} \quad (4)$$

$$\begin{aligned} \sum |M_g^{\text{viol}}|^2 = & \quad g_S^{2n-4}(Q^2) N^{n-2}(N^2 - 1) \\ & \times \left( \sum_{i < j} \right) \sum_{\text{perm}} \frac{1}{S_{12}} \frac{1}{S_{12}} \sum_{\tau} c_{\tau}^f. \end{aligned} \quad (5)$$

**Note:** Do not use `\label{#1}` on a line of a multiline equation if `\nonumber` is also used on that line. Incorrect cross-referencing will result. Notice the use `\text{#1}` for using a Roman font within a math environment.

To set a multiline equation without *any* equation numbers, use the `\begin{eqnarray*}`, `\end{eqnarray*}` format:

$$\begin{aligned} \sum |M_g^{\text{viol}}|^2 = & \quad g_S^{2n-4}(Q^2) N^{n-2}(N^2 - 1) \\ & \times \left( \sum_{i < j} \right) \left( \sum_{\text{perm}} \frac{1}{S_{12}S_{23}S_{n1}} \right) \frac{1}{S_{12}}. \end{aligned}$$

To obtain numbers not normally produced by the automatic numbering, use the `\tag{#1}` command, where `#1` is the desired equation number. For example, to get an equation number of (2.6'),

$$g^+g^+ \rightarrow g^+g^+g^+g^+ \dots, \quad q^+q^+ \rightarrow q^+g^+g^+ \dots \quad (2.6')$$

*a. A few notes on tags* `\tag{#1}` requires the `amsmath` package. Place the `\tag{#1}` command before the `\label{#1}`, if any. The numbering produced by `\tag{#1}` does not affect the automatic numbering in  $\text{\LaTeX}$ ; therefore, the number must be known ahead of time, and it must be manually adjusted if other equations are added. `\tag{#1}` works with both single-line and multiline equations. `\tag{#1}` should only be used in exceptional cases—do not use it to number many equations in your paper. Please note that this feature of the `amsmath` package is not compatible with the `hyperref` (6.77u) package.

Enclosing display math within `\begin{subequations}` and `\end{subequations}` will produce a set of equations that are labeled with letters, as shown in Eqs. (6b) and (6a) below. You may include any number of single-line and multiline equations, although it is probably not a good idea to follow one display math directly after another.

$$\begin{aligned} \mathcal{M} = & \quad ig_Z^2(4E_1E_2)^{1/2}(l_i^2)^{-1}(g_{\sigma_2}^e)^2\chi_{-\sigma_2}(p_2) \\ & \times [\epsilon_i]_{\sigma_1} \chi_{\sigma_1}(p_1). \end{aligned} \quad (6a)$$

$$\left\{ abc123456abcde\alpha\beta\gamma\delta1234556\alpha\beta\frac{1\sum_b^a}{A^2} \right\}, \quad (6b)$$

Giving a `\label{#1}` command directly after the `\begin{subequations}`, allows you to reference all the equations in the `subequations` environment. For example, the equations in the preceding subequations environment were Eqs. (6).

#### 1. Wide equations

The equation that follows is set in a wide format, i.e., it spans the full page. The wide format is reserved for long equations that cannot easily be set in a single column:

$$\mathcal{R}^{(d)} = g_{\sigma_2}^e \left( \frac{[\Gamma^Z(3, 21)]_{\sigma_1}}{Q_{12}^2 - M_W^2} + \frac{[\Gamma^Z(13, 2)]_{\sigma_1}}{Q_{13}^2 - M_W^2} \right) + x_W Q_e \left( \frac{[\Gamma^\gamma(3, 21)]_{\sigma_1}}{Q_{12}^2 - M_W^2} + \frac{[\Gamma^\gamma(13, 2)]_{\sigma_1}}{Q_{13}^2 - M_W^2} \right). \quad (7)$$

This is typed to show how the output appears in wide format. (Incidentally, since there is no blank line between the `equation` environment above and the start of this paragraph, this paragraph is not indented.)

## IX. CROSS-REFERENCING

REVTeX will automatically number such things as sections, footnotes, equations, figure captions, and table captions. In order to reference them in text, use the `\label{#1}` and `\ref{#1}` commands. To reference a particular page, use the `\pageref{#1}` command.

The `\label{#1}` should appear within the section heading, within the footnote text, within the equation, or within the table or figure caption. The `\ref{#1}` command is used in text at the point where the reference is to be displayed. Some examples: Section I on page 1, Table I, and Fig. 5.

## X. FLOATS: FIGURES, TABLES, VIDEOS, ETC.

Figures and tables are usually allowed to “float”, which means that their placement is determined by L<sup>A</sup>T<sub>E</sub>X, while the document is being typeset.

Use the `figure` environment for a figure, the `table` environment for a table. In each case, use the `\caption` command within to give the text of the figure or table caption along with the `\label` command to provide a key for referring to this figure or table. The typical content of a figure is an image of some kind; that of a table is an alignment.

Insert an image using either the `graphics` or `graphix` packages, which define the `\includegraphics{#1}` command. (The two packages differ in respect of the optional arguments used to specify the orientation, scaling, and translation of the image.) To create an alignment, use the `tabular` environment.

The best place to locate the `figure` or `table` environment is immediately following its first reference in text;

TABLE I. A table that fits into a single column of a two-column layout. Note that REVTeX 4 adjusts the intercolumn spacing so that the table fills the entire width of the column. Table captions are numbered automatically. This table illustrates left-, center-, decimal- and right-aligned columns, along with the use of the `ruledtabular` environment which sets the Scotch (double) rules above and below the alignment, per APS style.

Left <sup>a</sup>	Centered <sup>b</sup>	Decimal	Right
1	2	3.001	4
10	20	30	40
100	200	300.0	400

<sup>a</sup> Note a.

<sup>b</sup> Note b.

this sample document illustrates this practice for Fig. 5, which shows a figure that is small enough to fit in a single column.

In exceptional cases, you will need to move the float earlier in the document, as was done with Table II: L<sup>A</sup>T<sub>E</sub>X’s float placement algorithms need to know about a full-page-width float earlier.

Fig. 6 has content that is too wide for a single column, so the `figure*` environment has been used.

The content of a table is typically a `tabular` environment, giving rows of type in aligned columns. Column entries separated by `&`’s, and each row ends with `\\`. The required argument for the `tabular` environment specifies how data are aligned in the columns. For instance, entries may be centered, left-justified, right-justified, aligned on a decimal point. Extra column-spacing may be specified as well, although REVTeX 4 sets this spacing so that the columns fill the width of the table. Horizontal rules are typeset using the `\hline` command. The doubled (or Scotch) rules that appear at the top and bottom of a table can be achieved enclosing the `tabular` environment within a `ruledtabular` environment. Rows whose columns span multiple columns can be typeset using the `\multicolumn{#1}{#2}{#3}` command (for example, see the first row of Table II).

Tables I, II, III, and IV show various effects. A table that fits in a single column employs the `table` environment. Table II is a wide table, set with the `table*` environment. Long tables may need to break across pages. The most straightforward way to accomplish this is to specify the `[H]` float placement on the `table` or `table*` environment. However, the L<sup>A</sup>T<sub>E</sub>X 2<sub>ε</sub> package `longtable` allows headers and footers to be specified for each page of the table. A simple example of the use of `longtable` can be found in the file `summary.tex` that is included with the REVTeX 4 distribution.

There are two methods for setting footnotes within a table (these footnotes will be displayed directly below the table rather than at the bottom of the page or in the bibliography). The easiest and preferred method is just to use the `\footnote{#1}` command. This will automatically enumerate the footnotes with lowercase roman letters. However, it is sometimes necessary to have multiple entries in the table share the same footnote. In this case,

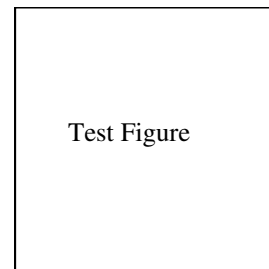


FIG. 5. A figure caption. The figure captions are automatically numbered.

Wide Test Figure

FIG. 6. Use the `figure*` environment to get a wide figure that spans the page in `twocolumn` formatting.

TABLE II. This is a wide table that spans the full page width in a two-column layout. It is formatted using the `table*` environment. It also demonstrates the use of `\multicolumn` in rows with entries that span more than one column.

Ion	$D_{4h}^1$		$D_{4h}^5$	
	1st alternative	2nd alternative	1st alternative	2nd alternative
K	$(2e) + (2f)$	$(4i)$	$(2c) + (2d)$	$(4f)$
Mn	$(2g)^a$	$(a) + (b) + (c) + (d)$	$(4e)$	$(2a) + (2b)$
Cl	$(a) + (b) + (c) + (d)$	$(2g)^a$	$(4e)^a$	
He	$(8r)^a$	$(4j)^a$	$(4g)^a$	
Ag		$(4k)^a$		$(4h)^a$

<sup>a</sup> The  $z$  parameter of these positions is  $z \sim \frac{1}{4}$ .

there is no choice but to manually create the footnotes using `\footnotemark[#1]` and `\footnotetext[#1]{#2}`. `#1` is a numeric value. Each time the same value for `#1` is used, the same mark is produced in the table. The `\footnotetext[#1]{#2}` commands are placed after the `tabular` environment. Examine the  $\LaTeX$  source and output for Tables I and IV for examples.

Video 1 illustrates several features new with REV $\TeX$ 4.2, starting with the `video` environment, which is in the same category with `figure` and `table`. The `\setfloatlink` command causes the title of the video to be a hyperlink to the indicated URL; it may be used with any environment that takes the `\caption` command. The `\href` command has the same significance as it does in the context of the `hyperref` package: the second argument is a piece of text to be typeset in your document; the first is its hyperlink, a URL.

*Physical Review* style requires that the initial citation of figures or tables be in numerical order in text, so don't cite Fig. 6 until Fig. 5 has been cited.

TABLE III. Numbers in columns Three–Five are aligned with the “d” column specifier (requires the `dcolumn` package). Non-numeric entries (those entries without a “.”) in a “d” column are aligned on the decimal point. Use the “D” specifier for more complex layouts.

One	Two	Three	Four	Five
one	two	three	four	five
He	2	2.77234	45672.	0.69
C <sup>a</sup>	C <sup>b</sup>	12537.64	37.66345	86.37

<sup>a</sup> Some tables require footnotes.

<sup>b</sup> Some tables need more than one footnote.

## ACKNOWLEDGMENTS

We wish to acknowledge the support of the author community in using REV $\TeX$ , offering suggestions and encouragement, testing new versions, ....

## Appendix A: Code

Below is the code for section IV

```
# Imports
```

TABLE IV. A table with numerous columns that still fits into a single column. Here, several entries share the same footnote. Inspect the  $\LaTeX$  input for this table to see exactly how it is done.

	$r_c$ (Å)	$r_0$ (Å)	$\kappa r_0$		$r_c$ (Å)	$r_0$ (Å)	$\kappa r_0$
Cu	0.800	14.10	2.550	Sn <sup>a</sup>	0.680	1.870	3.700
Ag	0.990	15.90	2.710	Pb <sup>b</sup>	0.450	1.930	3.760
Au	1.150	15.90	2.710	Ca <sup>c</sup>	0.750	2.170	3.560
Mg	0.490	17.60	3.200	Si <sup>d</sup>	0.900	2.370	3.720
Zn	0.300	15.20	2.970	Li <sup>b</sup>	0.380	1.730	2.830
Cd	0.530	17.10	3.160	Na <sup>e</sup>	0.760	2.110	3.120
Hg	0.550	17.80	3.220	K <sup>e</sup>	1.120	2.620	3.480
Al	0.230	15.80	3.240	Rb <sup>c</sup>	1.330	2.800	3.590
Ga	0.310	16.70	3.330	Cs <sup>d</sup>	1.420	3.030	3.740
In	0.460	18.40	3.500	Ba <sup>e</sup>	0.960	2.460	3.780
Tl	0.480	18.90	3.550				

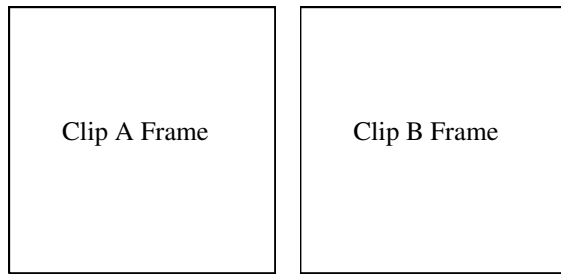
<sup>a</sup> Here's the first, from Ref. 2.

<sup>b</sup> Here's the second.

<sup>c</sup> Here's the third.

<sup>d</sup> Here's the fourth.

<sup>e</sup> And etc.



Video 1. Students explain their initial idea about Newton's third law to a teaching assistant. Clip (a): same force. Clip (b): move backwards.

```
import numpy as np
import matplotlib.pyplot as plt

# define hpp function
def gen_hpp(lmbda , N):
    '''
    param lmbda: rate parameter
    N: Number of events
    '''
    # inits
    t = [0]

    # begin loop
    while True:
        # generate uniform r.v. ~ Unif[0,1]
        u = np.random.uniform(0,1)
        # generate w ~ Exponential(lmbda)
        w = - np.log(u)/lmbda
        t.append(t[-1] + w)
        # exit condition
        if len(t) > N:
            # get time to event & count the
            # number of events
            return t , np.arange(len(t))

if __name__ == '__main__': # main namespace
    l , N = 8 , 1000
    # generate the time(s) to event(s)
    # AND count of events
    hpp_event_times , events = \
    gen_hpp(lmbda = 1 , N = N)
    print(events , hpp_event_times) # debug

    # Make plots
    fig , ax = plt.subplots()
    # step graph
    ax.step(hpp_event_times , events ,
            label = f"$\lambda$ = {l}" , lw = 0.5)
    ax.set_xlabel(r'$t$')
    ax.set_ylabel(r'$N(t)$')
    ax.set_title('Homogeneous Poisson process')
    ax.legend(loc='best')
    plt.show()
```

Below is the code for NHPP in section VI

```
# Imports
import numpy as np
import matplotlib.pyplot as plt

# define hpp function
def gen_nhpp(lmbda_bar = 8 , T = 140):
    '''
    param lmbda_bar: rate parameter that
    dominates the rate param of the NHPP
    T : maximum time
    '''
    # inits
    s = [0]
    t = [0]

    # begin loop
    while s[-1] < T:
        # generate uniform r.v. ~ Unif[0,1]
        u = np.random.uniform(0,1)
        # generate w ~ Exponential(lmbda)
        w = - np.log(u)/lmbda_bar
        s.append(s[-1] + w)
        # geerate D ~ Unif[0,1]
        D = np.random.uniform(0,1)
        # acceptance criterion
        if D < (1.01)*s[-1] / lmbda_bar:
            t.append(s[-1])

    if t[-1] > T:
        num_events = np.arange(len(t[:-1]))
        print(f'Breakpoint 1: the number of events \
is {num_events[-1]}, and the time taken to \
reach them is {t[:-1][-1]}')
        # get time to event & count the
        # number of events
        return t[:-1] , num_events
    else:
        num_events = np.arange(len(t))
        print(f'Breakpoint 2: the number of events is \
{num_events[-1]}, and the time taken to reach \
them is {t[-1]}')
        # get time to event & count the number
        # of events
        return t , num_events

# main namespace
if __name__ == '__main__':
    # generate the time(s) to event(s)
    # AND count of events
    nhpp_event_times , events = gen_nhpp()
    # Make plots
    fig , ax = plt.subplots()
    # step graph
    ax.step(nhpp_event_times , events ,
            label = f'$\lambda(t) = (1.01)^{t}$')
    # Integrating the rate function to get E[N(t)]
```



```
x = np.arange(140)
y = 1.01**x/np.log(1.01) - \
(1.01)**0/np.log(1.01)
ax.plot(x , y ,
label = r'E[N(t)] = $\frac{1.01^t}{\ln(1.01)}$ - 100.499')
# Auxiliaries
ax.set_xlabel(r'$t$')
ax.set_ylabel(r'$N(t)$')
ax.set_title('Non-Homogeneous \
Poisson process')
ax.legend(loc='best')
plt.show()
```

## Appendix B: Appendixes

To start the appendixes, use the `\appendix` command. This signals that all following section commands refer to appendixes instead of regular sections. Therefore, the `\appendix` command should be used only once—to setup the section commands to act as appendixes. Thereafter normal section commands are used. The heading for a section can be left empty. For example,

```
\appendix
\section{}
```

will produce an appendix heading that says “APPENDIX A” and

```
\appendix
```

```
\section{Background}
```

will produce an appendix heading that says “APPENDIX A: BACKGROUND” (note that the colon is set automatically).

If there is only one appendix, then the letter “A” should not appear. This is suppressed by using the star version of the appendix command (`\appendix*` in the place of `\appendix`).

## Appendix C: A little more on appendixes

Observe that this appendix was started by using

```
\section{A little more on appendixes}
```

Note the equation number in an appendix:

$$E = mc^2. \quad (C1)$$

### 1. A subsection in an appendix

You can use a subsection or subsubsection in an appendix. Note the numbering: we are now in Appendix C1.

Note the equation numbers in this appendix, produced with the subequations environment:

$$E = mc, \quad (C2a)$$

$$E = mc^2, \quad (C2b)$$

$$E \gtrsim mc^3. \quad (C2c)$$

They turn out to be Eqs. (C2a), (C2b), and (C2c).

- 
- [1] E. Witten, (2001), hep-th/0106109, and references therein
  - [2] See the explanation of time travel in R. P. Feynman, Phys. Rev. **94**, 262 (1954); The classical relativistic treatment of A. Einstein, Yu. Podolsky, and N. Rosen (EPR), *ibid.* **47**, 777 (1935) is a relative classic
  - [3] E. Beutler, in *Williams Hematology*, Vol. 2, edited by E. Beutler, M. A. Lichtman, B. W. Collier, and T. S. Kipps (McGraw-Hill, New York, 1994) Chap. 7, pp. 654–662, 5th ed.
  - [4] N. D. Birell and P. C. W. Davies, *Quantum Fields in Curved Space* (Cambridge University Press, 1982).
  - [5] J. G. P. Berman and J. F. M. Izrailev, Stability of nonlinear modes, Physica D **88**, 445 (1983).
  - [6] E. B. Davies and L. Parns, Trapped modes in acoustic waveguides, Q. J. Mech. Appl. Math. **51**, 477 (1988).
  - [7] Automatically placing footnotes into the bibliography requires using BibTeX to compile the bibliography.
  - [8] E. Beutler, in *Williams Hematology*, Vol. 2, edited by E. Beutler, M. A. Lichtman, B. W. Collier, and T. S. Kipps (McGraw-Hill, New York, 1994) 5th ed., Chap. 7, pp. 654–662.
  - [9] D. E. Knuth, in *Fundamental Algorithms*, The Art of Computer Programming, Vol. 1 (Addison-Wesley, Reading, Massachusetts, 1973) Section 1.2, pp. 10–119, 2nd ed., a full INBOOK entry.
  - [10] J. S. Smith and G. W. Johnson, Philos. Trans. R. Soc. London, Ser. B **777**, 1395 (2005).
  - [11] W. J. Smith, T. J. Johnson, and B. G. Miller, Surface chemistry and preferential crystal orientation on a silicon surface (2010), J. Appl. Phys. (unpublished).
  - [12] V. K. Smith, K. Johnson, and M. O. Klein, Surface chemistry and preferential crystal orientation on a silicon surface (2010), J. Appl. Phys. (submitted).
  - [13] U. Underwood, N. Net, and P. Pot, Lower bounds for wishful research results (1988), talk at Fanstord University (A full UNPUBLISHED entry).
  - [14] M. P. Johnson, K. L. Miller, and K. Smith, personal communication (2007).
  - [15] J. Smith, ed., *AIP Conf. Proc.*, Vol. 841 (2007).
  - [16] W. V. Oz and M. Yannakakis, eds., *Proc. Fifteenth Annual*, All ACM Conferences No. 17, ACM (Academic Press, Boston, 1983) a full PROCEEDINGS entry.
  - [17] Y. Burstyn, Proceedings of the 5th International Molec-

- ular Beam Epitaxy Conference, Santa Fe, NM (2004), (unpublished).
- [18] B. Quinn, ed., *Proceedings of the 2003 Particle Accelerator Conference, Portland, OR, 12-16 May 2005* (Wiley, New York, 2001) albeit the conference was held in 2005, it was the 2003 conference, and the proceedings were published in 2001; go figure.
  - [19] A. G. Agarwal, Proceedings of the Fifth Low Temperature Conference, Madison, WI, 1999, *Semiconductors* **66**, 1238 (2001).
  - [20] R. Smith, Hummingbirds are our friends, *J. Appl. Phys.* (these proceedings) (2001), abstract No. DA-01.
  - [21] J. Smith, *Proc. SPIE* **124**, 367 (2007), required title is missing.
  - [22] T. Terrific, *An  $O(n \log n / \log \log n)$  Sorting Algorithm*, Wishful Research Result 7 (Fanstord University, Computer Science Department, Fanstord, California, 1988) a full TECHREPORT entry.
  - [23] J. Nelson, TWI Report 666/1999 (Jan. 1999) required institution missing.
  - [24] W. K. Fields, ECE Report No. AL944 (2005) required institution missing.
  - [25] Y. M. Zalkins, e-print arXiv:cond-mat/040426 (2008).
  - [26] J. Nelson, U.S. Patent No. 5,693,000 (12 Dec. 2005).
  - [27] J. K. Nelson, M.S. thesis, New York University (1999).
  - [28] É. Masterly, *Mastering Thesis Writing*, Master's project, Stanford University, English Department (1988), a full MASTERSTHESIS entry.
  - [29] S. M. Smith, Ph.D. thesis, Massachusetts Institute of Technology (2003).
  - [30] S. R. Kawa and S.-J. Lin, *J. Geophys. Res.* **108**, 4201 (2003), DOI:10.1029/2002JD002268.
  - [31] F. P. Phony-Baloney, *Fighting Fire with Fire: Festooning French Phrases*, PhD dissertation, Fanstord University, Department of French (1988), a full PHDTHESIS entry.
  - [32] D. E. Knuth, *Seminumerical Algorithms*, 2nd ed., *The Art of Computer Programming*, Vol. 2 (Addison-Wesley, Reading, Massachusetts, 1981) a full BOOK entry.
  - [33] J. C. Knvth, *The programming of computer art*, Vernier Art Center, Stanford, California (1988), a full BOOK-LET entry.
  - [34] R. Ballagh and C. Savage, Bose-einstein condensation: from atomic physics to quantum fluids, proceedings of the 13th physics summer school (World Scientific, Singapore, 2000) cond-mat/0008070.
  - [35] R. Ballagh and C. Savage, Bose-einstein condensation: from atomic physics to quantum fluids, in *Proceedings of the 13th Physics Summer School*, edited by C. Savage and M. Das (World Scientific, Singapore, 2000) cond-mat/0008070.
  - [36] W. Opechowski and R. Guccione, Introduction to the theory of normal metals, in *Magnetism*, Vol. IIa, edited by G. T. Rado and H. Suhl (Academic Press, New York, 1965) p. 105.
  - [37] W. Opechowski and R. Guccione, Introduction to the theory of normal metals, in *Magnetism*, Vol. IIa, edited by G. T. Rado and H. Suhl (Academic Press, New York, 1965) p. 105.
  - [38] W. Opechowski and R. Guccione, Introduction to the theory of normal metals, in *Magnetism*, Vol. IIa, edited by G. T. Rado and H. Suhl (Academic Press, New York, 1965) p. 105.
  - [39] J. M. Smith, *Molecular dynamics* (Academic, New York, 1980).
  - [40] V. E. Zakharov and A. B. Shabat, Exact theory of two-dimensional self-focusing and one-dimensional self-modulation of waves in nonlinear media, *Zh. Eksp. Teor. Fiz.* **61**, 118 (1971), [*Sov. Phys. JETP* **34**, 62 (1972)].
  - [41] J. M. Smith, in *Molecular Dynamics*, edited by C. Brown (Academic, New York, 1980).
  - [42] D. D. Lincoll, Semigroups of recurrences, in *High Speed Computer and Algorithm Organization*, Fast Computers No. 23, edited by D. J. Lipcoll, D. H. Lawrie, and A. H. Sameh (Academic Press, New York, 1977) 3rd ed., Part 3, pp. 179–183, a full INCOLLECTION entry.
  - [43] A. V. Oaho, J. D. Ullman, and M. Yannakakis, On notions of information transfer in VLSI circuits, in *Proc. Fifteenth Annual ACM*, Boston, 1982, All ACM Conferences No. 17, edited by W. V. Oz and M. Yannakakis, ACM (Academic Press, New York, 1983) pp. 133–139, a full INPROCEEDINGS entry.
  - [44] L. Manmaker, *The Definitive Computer Manual*, Chips-R-Us, Silicon Valley, silver ed. (1986), a full MANUAL entry.