Eight lectures covering:

➤ *Introduction to modelling*, what is it, why is it useful?

➤ *Simulation techniques*, random number generation, analysis of results from simulation or measurements

➤ *Operational analysis*, performance bounds, balanced systems

➤ *Queueing theory*, Markov chains, single/multiple servers, bounded queues, queueing networks
Recommended books


— Simulation, random number generation, operational analysis, some basic queueing theory, substantial sections on designing and analysing experiments


— Operational analysis, basic queueing theory, small section on simulation


— A classic on queueing theory, much more emphasis on mathematical derivations
Why model?

- A manufacturer may have a range of compatible systems with different characteristics — which configuration would be best for a particular application?

- A system is performing poorly — what should be done to improve it? Which problems should be tackled first?

- Fundamental design decisions may affect performance of a system. A model can be used as part of the design process to avoid bad decisions and to help quantify a cost/benefit analysis.
A toy problem

<table>
<thead>
<tr>
<th>system</th>
<th>CPU time</th>
<th>disk time</th>
<th>total</th>
</tr>
</thead>
<tbody>
<tr>
<td>A</td>
<td>4.6</td>
<td>4.0</td>
<td>8.6</td>
</tr>
<tr>
<td>B1</td>
<td>5.1</td>
<td>1.9</td>
<td>7.0</td>
</tr>
<tr>
<td>B2</td>
<td>3.1</td>
<td>1.9</td>
<td>5.0</td>
</tr>
</tbody>
</table>

- A database running on Type A system is too slow
- Type B1 system available immediately, and a type B2 system in the future
- Which option is best:
  - Stick with A?
  - Change to B1 immediately?
  - Wait and change to B2?
- What factors affect the correct choice?
How can modelling help?

Typical performance questions we might ask include:

- How long will a database request wait before receiving CPU service?

- What is the utilization of the resource (CPU, disk, ...)? (Utilization is the proportion of time that the resource is busy.)

- What is the distribution of the number of requests queued at some time \( t \)? What is its mean, standard deviation, ...
Techniques

Many different approaches:

➢ *Measurement* — if the system already exists then maybe it can be changed and the effects analysed

➢ *Simulation* — construct a computer program that emulates the system under study and study that instead

➢ *Operational analysis* — analysis based on directly measured quantities and relationships between them: makes few assumptions about the system

➢ *Queueing theory* — stochastic processes, analytical models of queueing systems
Techniques (2)

Choosing between these techniques depends on . . .

➤ *Stage of development*: can only measure an existing system

➤ *Time available*: measurements or simulations can take a long time to complete. How easily can different trade-offs be evaluated?

➤ *Resources*: systems with which to experiment, people with the relevant skills, cost

➤ *Desired accuracy*: how do the assumptions made in analytic techniques effect the result? Are appropriate parameters and workloads used during experimental work?

➤ *Creditable*: will people believe (act on) the results?
Little’s result

Begin with a simple derivation of *Little’s Result* — a very general theorem relating the number of jobs in a system with the time they spend there.

\[
\text{e.g.: A disk server takes, on average, 10ms to satisfy an I/O request. If the request rate is 100 per second, then how many requests are queued at the server?}
\]
Little’s result (2)

\[ \alpha(t) = \text{number of arrivals in } (0, t) \]
\[ \delta(t) = \text{number of departures in } (0, t) \]
\[ N(t) = \alpha(t) - \delta(t) \]

The area \( \gamma(t) \) between the curves \( \alpha(t) \) and \( \delta(t) \) represents the total time all customers have spent in system in \((0, t)\).
**Little’s result (3)**

Let

\[ \lambda_t = \frac{\alpha(t)}{t} \]

\[ T_t = \frac{\gamma(t)}{\alpha(t)} \]

\[ \overline{N_t} = \frac{\gamma(t)}{t} \]

- \( \lambda_t \) — the average arrival rate during \((0, t)\)
- \( T_t \) — system time per customer averaged over all customers in \((0, t)\)
- \( \overline{N_t} \) — average number of customers in system during \((0, t)\)

Combining these:

\[ \overline{N_t} = \lambda_t T_t \]
Little’s result (4)

Assume the following limits exist

\[
\lambda = \lim_{t \to \infty} \lambda_t \\
T = \lim_{t \to \infty} T_t
\]

Then we have

\[
\bar{N} = \lambda T
\]

That is, the average number in the system equals the average arrival rate \( \times \) average time in system.

This is Little’s result. The proof makes no assumptions about the way that arrivals or departures are distributed, the queueing discipline or how many servers service the queue.

First proved by Little in 1961.
Applications of Little’s result

We can re-state this for any boundary of our queueing system.

We can split $T$ (average time in the system) into $W$ (average time spent waiting) and $\bar{x}$ (average time spent being served).

Similarly, we can split $N$ (average number in the system) into $N_q$ (average number in the queue) and $N_s$ (average number being served).

Applying Little’s result separately to the queue and to the server:

$$\overline{N_q} = \lambda W$$

$$\overline{N_s} = \lambda \bar{x}$$
Probability theory refresher

We write the probability that an event occurs as: \( P(\text{event}) \)

— e.g. \( P(X < 0.5) \): “the random variable \( X \) is less than 0.5”

\( P(a, b) \) is the *joint* probability that *both* events \( a \) and \( b \) occur

— e.g. \( P(X < 0.5, Y < 0.5) \): “both \( X \) and \( Y \) are less than 0.5”

\( P(a \mid b) \) is the *conditional* probability that event \( a \) occurs given that event \( b \) has occurred.

— e.g. \( P(X < 0.5 \mid Y < 0.5) \) “\( X \) is less than 0.5 given that \( Y \) is less than 0.5”

\[
P(a \mid b) := \frac{P(a, b)}{P(b)}
\]
CDF and PDF

We can describe the distribution from which a continuous random variable is drawn in two ways: using the cumulative distribution function (cdf) or the probability density function (pdf).

For example, for a random variable \( X \) uniformly distributed from 0 to 1, the cdf and pdf would be

\[
F_X(x) := \mathbb{P}(X \leq x) \quad f_X(x) := \frac{dF_X(x)}{dx}
\]

Note that

\[
F_X(x) = \int_{-\infty}^{x} f_X(y) \, dy , \quad \text{and} \quad \int_{-\infty}^{\infty} f_X(y) \, dy = 1.
\]
Expected value and moments

The *expected value* of $X$, $\mathbb{E}(X)$, is given by

$$\mathbb{E}(X) := \overline{X} := \int_{-\infty}^{\infty} x f_X(x) \, dx$$

Also called the average, mean or first moment of the distribution.

The \(n^{th}\) *moment* is defined as

$$\mathbb{E}(X^n) := \overline{X^n} := \int_{-\infty}^{\infty} x^n f_X(x) \, dx$$

Take care to distinguish between $\overline{X^n}$ and $\overline{X}^n$.

The \(n^{th}\) *central moment* is

$$\overline{(X - \overline{X})^n} = \int_{-\infty}^{\infty} (x - \overline{X})^n f_X(x) \, dx$$
Variance and std dev

The 2\textsuperscript{nd} central moment is known as the \textit{variance},

$$\sigma^2_X := (X - \bar{X})^2 = X^2 - (\bar{X})^2$$

From this, we define:

$$\sigma_X := \sqrt{\sigma^2_X} \quad \text{the standard deviation}$$

$$C_X := \frac{\sigma_X}{\bar{X}} \quad \text{the coefficient of variation}$$

Numerically higher values signify "more variable" data.

For example, a coefficient of variation of 5 might be considered large, and 0.2 considered small.
Example: uniform

The uniform distribution is used where a continuous random variable takes values in some bounded range and there is no reason to favour one value over another

\[ a = \text{lower limit} \]
\[ b = \text{upper limit, } b > a \]

\[ f(x) = \frac{1}{b - a} \]

\[ F(x) = \begin{cases} 
0 & \text{if } x < a \\
\frac{x - a}{b - a} & \text{if } a \leq x < b \\
1 & \text{if } b \leq x 
\end{cases} \]

The mean is \( \frac{1}{2}(a + b) \), the variance is \( \frac{1}{12}(b - a)^2 \).
Example: exponential

One scale parameter, $\lambda > 0$, with

$$f(x) = \lambda e^{-\lambda x}$$

$$F(x) = 1 - e^{-\lambda x}$$

The mean is $\frac{1}{\lambda}$, the variance is $\frac{1}{\lambda^2}$.
Consequently $\lambda$ may be viewed as the mean event rate.

The exponential distribution is the only continuous distribution with the Memoryless Property that

$$\mathbb{P}(X > t + s \mid X > t) = \mathbb{P}(X > s)$$

Intuitively, it may be used to model the distribution of inter-event times in which the time until the next event does not depend on the time already waited.
Discrete distributions

The previous examples have concerned \textit{continuous} random variables whose distributions have been defined by their \textit{cdf} or, equivalently, their \textit{pdf}.

Similar definitions apply to the case of \textit{discrete} random variables, $X$, taking values $x_i$ ($i \in I$). We have that

$$\sum_{i \in I} P[X = x_i] = 1$$

The \textit{expected value} of $X$ is

$$\mathbb{E}(X) := \sum_{i \in I} x_i P[X = x_i]$$

Similarly, for the other moments, where the integration becomes a summation over the set of possible values.
Example: Bernoulli

The *Bernoulli distribution* is used where a discrete random variable can take only two values, usually either 1 representing the *success* or 0 representing the *failure* of some operation.

\[
p = \text{probability of success, } 0 \leq p \leq 1
\]

\[
\mathbb{P}(X = x) = \begin{cases} 
1 - p & \text{if } x = 0 \\
p & \text{if } x = 1 \\
0 & \text{otherwise}
\end{cases}
\]

The mean is evidently \( p \) and the variance is \( p(1 - p) \).
Example: Binomial

The number of successes in a series of Bernoulli trials is a discrete random variable that has a distribution known as the \textit{Binomial distribution}.

\[ p = \text{probability of success, } 0 \leq p \leq 1 \]

\[ n = \text{number of trials, } n \in \mathbb{N}^+ \]

\[ \mathbb{P}(X = x) = \binom{n}{x} p^x (1 - p)^{n-x} \]

The mean is \( np \) and the variance is \( np(1 - p) \).

Note that \( \mathbb{P}(X = x) \) is the product of the number of ways that \( x \) successes can occur in \( n \) trials and the probability that exactly that pattern of successes and failures will occur.
A simple queueing system

We characterise queueing systems by:

- Arrival process
  \[ A(t) = P(\text{inter-arrival time} \leq t) \]

- Service process
  \[ B(x) = P(\text{service time} \leq x) \]

- Storage capacity available for waiting customers

- The number of servers/customers available

- The different kinds of arriving customers (big jobs, small jobs, ...)

- Queueing discipline used: FCFS, LCFS, priority, ...

- Defections, balking, bribing, ...
Queueing systems notation

The Kendall notation describes a single queueing system using the notation $A/B/m/k/l$ where:

- $A$ is the inter-arrival time distribution of customers
- $B$ is the service time distribution
- $m$ is the number of parallel servers
- $k$ is the limit on the customers in this system
- $l$ is the population size

If the population size or the limit on the queue length are not specified then they are assumed to be infinite.
Queueing notation (2)

Typical values for $A, B$ include:

- $M$ – exponential distribution
- $E_r$ – $r$ stage Erlangian distribution
- $D$ – Deterministic
- $G$ – General

Examples:

- $M/M/1$: exponential inter-arrival, exponential service, single server
- $M/E_r/1$: exponential inter-arrival, $r$ stage Erlang service, single server
- $M/G/1$: exponential inter-arrival, general service time, single server
- $M/M/K/K$: exponential inter-arrival, exponential service, $K$ servers and at most $K$ customers present
More generally, consider systems comprising multiple inter-connected service centres, forming a *queueing network*. Consider:

- the properties of each node
  - e.g. using Kendall notation

- the way in which jobs move between the nodes
  - e.g. the links that exist between nodes and the way in which a job leaving one node selects between these links

- the workload being analyzed
  - e.g. disk-server workload comprises a 20 : 1 mix of small/large requests
Queueing networks (2)

We can classify queueing networks as either

- **closed** networks in which a fixed set of jobs circulate between nodes, but no new jobs are introduced and no jobs leave the system
  — e.g. a computer system with a fixed number of terminals attached to it

- **open** networks in which jobs may enter and leave the system
  — e.g. the network on the previous slide: jobs arrive at 1 or 5 and leave from 1 or 3

Open networks may be further classified as **feed-forward** if a single job visits each node at most once.
Simulation

Simulation allows arbitrarily complex systems to be evaluated

➤ Able to capture the dynamic behaviour of systems

➤ Captures the dynamics of complex systems by imitation

➤ Tracks the evolution of the system over time

➤ Examples include communication network design, road traffic modelling, studying chemical reactions, fluid flow, etc.
Simulation (2)

Execution of simulation model consists of a series of state space changes.

These closely follow the chronological order of events in the system being modelled.

Consider simulation as a ‘set of equations’ describing evolution in time of system under study.

The ‘equations’ are ‘solved’ by following their evolution in time.

We *always* follow the *dynamic* evolution of the system, even if we only want a mean value — therefore, as well as techniques for implementing simulators, it is necessary to know how to analyse their results.

Simulation is of particular use when studying systems which are not in a steady state condition.
Types of simulation

Simulation studies may be classified as

➤ *Discrete state/event* simulation in which the state of the system is described by discrete variables
  
  — e.g. the number of jobs at different stages on a production line

➤ *Continuous state/event* simulation in which the state is described by continuous variables
  
  — e.g. the quantities of various chemical reagents in a vat

A similar distinction may be drawn between *discrete time* and *continuous time* simulations depending on whether the system state is defined at all times.

  — e.g. a simulation of the number of students attending these lectures would be discrete time
Types of simulation (2)

We will be concerned with *discrete event* simulation because it applies most naturally to computer systems in which state variables are generally discrete, e.g.

- the state of jobs in the system
- the number of jobs of different kinds
- the number or availability of devices
Pros and cons

The principal advantage of simulation is its extreme generality: the programmer may add arbitrary details. However,

➤ The design, coding and debugging of a simulation program is often time consuming and difficult to understand — it may even approach that of implementing the system and measuring it directly

➤ Generality can lead to complexity which can obscure understanding of the model — fine details may be irrelevant if the simulated workload is a poor estimate

➤ Execution of the simulation can be computationally expensive

➤ Statistical analysis of the output can be problematic — e.g. how long should the simulation run before taking an average of the results?
Events

Each event contains a time stamp identifying ‘when it occurs’ and denotes some change in the state of the system to be simulated e.g. ‘packet arrived’

Events are ordered in time in an event list

<table>
<thead>
<tr>
<th>Initialize the clock to 0</th>
</tr>
</thead>
<tbody>
<tr>
<td>Initialise the event list</td>
</tr>
<tr>
<td>WHILE termination criterion is not met</td>
</tr>
<tr>
<td>remove a smallest tuple (t,m) from the event list</td>
</tr>
<tr>
<td>update the clock to t</td>
</tr>
<tr>
<td>simulate the effect of transmitting m at time t</td>
</tr>
</tbody>
</table>

It is crucial that the simulator always selects the event with the least time stamp

Frequently most of the simulation time is spent maintaining the chronological order of the event list.
Physical systems

Concerned with the problem of representing so-called ‘physical systems’

A physical system consists of one or more physical processes

Each physical process operates independently, aside from interaction via messages (events)

All state changes are summarized in the event exchanges

Note the similarity with object-oriented styles of programming in which objects communicate solely by method invocations. Older object-oriented languages tended to refer to these as message-send operations.
An automatic car wash which is able to service one car at a time — is it viable to install a second car wash?

The model consists of an attendant (att), two car washes, (cw1, cw2), the entrance to the system (source) and the exit (sink)

- The inter-arrival time of cars that need to be washed is randomly distributed
- If both car washes are busy, an arriving car joins the queue
- When a car wash becomes idle then the car at the head of the queue is sent to it.
Example (2)

Events:

- **source→att**: car arrives in the system
- **att→cw**: car sent from the queue to the car wash
- **cw→att**: car wash becomes idle
- **cw→sink**: car departs the system

Note how the departure of a car is modelled by *two* events: one representing the car leaving the system and the other that signals to the attendant that the car wash is now idle.

Given that it takes *cw1* 8 minutes and *cw2* 10 minutes to wash a car, a *possible* sequence of events is ...
Sequence of events

<table>
<thead>
<tr>
<th>message</th>
<th>event</th>
<th>time</th>
<th>sender</th>
<th>receiver</th>
<th>content</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>-</td>
<td>0</td>
<td>cw1</td>
<td>att</td>
<td>idle</td>
</tr>
<tr>
<td>2</td>
<td>-</td>
<td>0</td>
<td>cw2</td>
<td>att</td>
<td>idle</td>
</tr>
<tr>
<td>3</td>
<td>1</td>
<td>6</td>
<td>source</td>
<td>att</td>
<td>car 1</td>
</tr>
<tr>
<td>4</td>
<td>2</td>
<td>6</td>
<td>att</td>
<td>cw1</td>
<td>car 1</td>
</tr>
<tr>
<td>5</td>
<td>3</td>
<td>11</td>
<td>source</td>
<td>att</td>
<td>car 2</td>
</tr>
<tr>
<td>6</td>
<td>4</td>
<td>11</td>
<td>att</td>
<td>cw2</td>
<td>car 2</td>
</tr>
<tr>
<td>7</td>
<td>5</td>
<td>12</td>
<td>source</td>
<td>att</td>
<td>car 3</td>
</tr>
<tr>
<td>8</td>
<td>6</td>
<td>14</td>
<td>cw1</td>
<td>sink</td>
<td>car 1</td>
</tr>
<tr>
<td>9</td>
<td>-</td>
<td>14</td>
<td>cw1</td>
<td>att</td>
<td>idle</td>
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<tr>
<td>10</td>
<td>7</td>
<td>14</td>
<td>att</td>
<td>cw1</td>
<td>car 3</td>
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<tr>
<td>11</td>
<td>8</td>
<td>17</td>
<td>source</td>
<td>att</td>
<td>car 4</td>
</tr>
<tr>
<td>12</td>
<td>9</td>
<td>19</td>
<td>source</td>
<td>att</td>
<td>car 5</td>
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<td>13</td>
<td>10</td>
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<td>cw2</td>
<td>sink</td>
<td>car 2</td>
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<td>-</td>
<td>21</td>
<td>cw2</td>
<td>att</td>
<td>idle</td>
</tr>
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<td>16</td>
<td>12</td>
<td>22</td>
<td>cw1</td>
<td>sink</td>
<td>car 3</td>
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<td>17</td>
<td>-</td>
<td>22</td>
<td>cw1</td>
<td>att</td>
<td>idle</td>
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<td>att</td>
<td>car 6</td>
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<td>20</td>
<td>15</td>
<td>30</td>
<td>cw1</td>
<td>sink</td>
<td>car 5</td>
</tr>
<tr>
<td>21</td>
<td>-</td>
<td>30</td>
<td>cw1</td>
<td>att</td>
<td>idle</td>
</tr>
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<td>22</td>
<td>16</td>
<td>30</td>
<td>att</td>
<td>cw1</td>
<td>car 6</td>
</tr>
</tbody>
</table>
Modelling stochastic systems

To represent the stochastic nature of the systems being modelled, simulation requires the use of random variables.

An input to the simulation may have a stochastic distribution e.g. arrivals at a queue.

Simulation model needs to sample random variables from the given distribution to “re-create” the input process.

Random variables can be generated for a wide range of theoretical and empirical distributions.

This process requires only a sequence of independent random variables with a uniform distribution!
Random number generation

Many random number generators in use today are not particularly good.

It is important not to generate random numbers with an ad hoc method — complex algorithms do not necessarily generate random outputs.

Some operating systems include support for generating random numbers based on (e.g.) key strokes or network inter-arrival times — however, these mechanisms are not usually suited to generating large volumes of random data.

How can we generate a large random sequence in a deterministic manner?

The answer is that the sequence is not random, but appears random as far as can be determined from statistical tests. They are termed pseudo-random.
Random number generation
(2)

Important requirements include

➢ The algorithm should be fast
➢ The storage requirements should be low
➢ The random sequence should only repeat after a very long period
➢ The sequence generated should possess two important statistical properties: uniformity and independence

For ease of implementation and speed, most random number generators use integer representation over an interval $m$

Given a value in this range, a desired value in the range $(0, 1)$ can be obtained by dividing by $m$
Linear congruential method

Many successful random number generators are special cases of the following scheme, introduced by Lehmer in 1951

It is based on 4 magic numbers:

- The starting value $X_0 \geq 0$
- The multiplier $a \geq 0$
- The increment $c \geq 0$
- The modulus $m \geq X_0, m > a, m > c$

The desired sequence of random numbers, $X_n$ is then generated by setting

$$X_{n+1} = (aX_n + c) \mod m$$
Multiplicative congruential method

The special case where $c = 0$ yields a purely multiplicative method.

When $c = 0$ the period of the sequence is reduced but, to our advantage, generation is a little faster.

How do we choose the values?

The case $a = 1$ can be rejected immediately as this would mean that

\[ X_n = (X_0 + nc) \mod m \]

The case $a = 0$ is even worse.

So $a > 1$ is necessary.
Choice of values

Since the period can never be greater than $m$ this should be chosen to be as large as possible. The following values were chosen for efficient computation of random numbers on the IBM/370

\[
\begin{align*}
a &= 16807 \\
m &= 2147483647 \\
c &= 0 \\
X_0 &= 314159
\end{align*}
\]

Knuth has observed that the value $a$ should end in the digits $x21$ where $x$ is even e.g. 31415821 (Implementation problem: what about overflow?)
Tests for randomness

Various statistical tests exist.

These include tests for uniformity (such as the Kolmogorov-Smirnov test and Chi-Square test).

Runs tests, which examine the arrangement of numbers in a sequence (a run) to test the hypothesis of independence.

These tests frequently check for the number of “up runs”, the number of “down runs”, and the runs above and below the mean.

Autocorrelation tests check the correlation structure of the sequence of observations.
Random variable generation

We have a sequence of pseudo-random uniform variables. How do we generate variables from other distributions?

Random behaviour can be programmed so that the random variables appear to have been drawn from a particular probability distribution

If \( f(x) \) is the desired pdf, then consider the CDF

\[
F_X(x) = \int_{-\infty}^{x} f(x) \, dx
\]

This is non-decreasing and lies between 0 and 1.
Random variable generation
(2)

Given a sequence of random numbers $r_i$
distributed over the same range $(0, 1)$

Then the corresponding value $x_i$ such
that $r_i = F_X(x_i)$ is uniquely determined by

$$x_i = F_X^{-1}(r_i)$$

The sequence $x_i$ is randomly distributed with
the probability density function $f(x)$ and
cumulative distribution function $F_X(x)$. 
Uniform distribution \((a, b)\)

Consider the \textit{uniform} random variable on the interval \((a, b)\) where

- \(a, b\) are arbitrary real numbers
- \(a < b\)

The distribution function is

\[
F_X(x_i) = \frac{(x_i - a)}{(b - a)}
\]

for \(x_i\) in the interval \((a, b)\)

Given \(r_i\) on interval \((0, 1)\) the problem is that of obtaining \(x_i\)

Starting with \(r_i = \frac{(x_i - a)}{(b - a)}\) we find that \(x_i = (b - a)r_i + a\).
Exponential distribution

For the exponential distribution

\[ F_X(x_i) = 1 - e^{-\lambda x_i} \]

for positive \( x_i \)

Thus

\[ r_i = 1 - e^{-\lambda x_i} \]

\[ 1 - r_i = e^{-\lambda x_i} \]

\[ \ln(1 - r_i) = -\lambda x_i \]

\[ x_i = -\frac{1}{\lambda} \ln(1 - r_i) \]

Note that \( r_i \) has the same distribution as \( 1 - r_i \) so we might as well use

\[ x_i = -\frac{1}{\lambda} \ln(r_i) \]

Other random variables can be derived in a similar fashion.
Simulation performance measures

As the simulation itself is stochastic, so too are the observed outputs.

It is critical to realize that a simulation can only yield *estimates* for performance measures.

There will *always* be some statistical error in the estimates.

We can attempt to reduce the error by

- running the simulation for longer until sufficient samples have been taken.
- running the same simulation with a different pseudo-random number generator, and combining the results from multiple runs.
Utilization

The utilization is the proportion of time that a server is busy.

An estimate can therefore be obtained by taking the sum of the busy times of the server and dividing by $T$, the simulation length.

In the case of a $k$-server, the busy times can be estimated together and divided by $kT$.

There are two obvious ways to aggregate the busy times:

- Sample the system and observe whether servers are busy or idle
- When the server becomes busy the time is saved, and when it becomes idle again the difference in the two times is the busy period.
Throughput, queue length

A simple count is kept of the number of customers receiving service. The throughput is this value divided by the total time.

There are at least two ways to estimate queue length:

- estimate the queue length distribution, then use that to obtain the mean

- View the queue length as a function of time: the mean queue length is $1/T$ of integral of this function.
Queue length

- Each time the queue length changes the time $t_{i+1}$ is stored
- subtract previous recorded time $t_i$ from current time
- multiply by previous queue length $n_i$

Sum these areas to give mean queue length (for $M$ observations) by

$$\overline{N} = \frac{1}{T} \sum_{i=1}^{M} n_i(t_{i+1} - t_i)$$
Queueing time

Two obvious ways to obtain *queueing time*

- observe the queueing times and take their average
- simply use Little’s law
Statistical analysis of results

Each independent replication of a simulation experiment will yield a different outcome.

To make a statement about the accuracy we need to consider the distribution of our estimator.

A typical method to incorporate the variability of the estimates is to construct *confidence intervals*. 

Confidence intervals

Given some point estimate $p$ we produce a confidence interval $(p - \delta, p + \delta)$.

The “true” value is estimated to be contained within the interval with some chosen probability, e.g. 0.9, say.

The value $\delta$ depends on the confidence level — the greater the confidence required the larger the value of $\delta$.

Let $x_1, x_2, \ldots, x_n$ be the values of a random sample from a population determined by the random variable $X$.

Assume: either $X$ is normally distributed or $n$ is large then

*Law of large numbers*: $\overline{X} \approx$ normally distributed.
Confidence intervals (2)

Given the variance, \( \sigma \), of \( X \) the 100(1 - \( \alpha \))% confidence interval is

\[ \bar{x} \pm \delta \]

where

\[ \delta = \frac{z_{\alpha/2} \sigma}{\sqrt{n}} \]

and

\[ \bar{x} = \sum_{i=1}^{n} x_i \]

\( z_{\alpha} \) is defined to be the largest value of \( z \) such that \( \mathbb{P}(Z > z) = \alpha \) where \( Z \) is the standard normal random variable.

These values, \( z_{\alpha} \), can be taken from tables of the standard normal distribution.

For example, for a 95% confidence interval \( \alpha = 0.05 \) and \( z_{\alpha/2} = z_{0.025} = 1.96 \)
Using Student’s T

When we know neither $\mu$ nor $\sigma$ we use the observed sample mean $\bar{x}$ and sample standard deviation $s$ where

$$s^2 = \frac{1}{n-1} \sum_{i=1}^{n} (x_i - \bar{x})^2$$

If $n$ is large then we simply use $s$ for $\sigma$ as above.

If $n$ is small then we may use

$$\delta = \frac{t_{\alpha/2} s}{\sqrt{n}}$$

$t_{\alpha/2}$ is defined by $\mathbb{P}(T > t_{\alpha/2}) = \alpha/2$

$T$ has a Student-$t$ distribution with $n - 1$ degrees of freedom.

This is the more frequently used formula in simulation models.
Stopping rules

How do we know when a system has been run ‘long enough’ for performance measures to be realistic?

We also need to be aware of transient conditions at the start of the simulation.

We can also repeat the simulation several times to obtain many samples for the ensemble average.

These multiple runs are called *replications*.

Given a performance measure $L$ we can approach the ensemble process by

$$L = \lim_{n \to \infty} \sum_{i=1}^{n} \frac{L_i}{n}$$
Stopping rules (2)

Having repeated the experiment $n$ times, we can construct a confidence interval on our measure $L$.

Although large numbers of replications reduce the variance in the sample, each replication requires re-stabilizing the simulation.

Can we avoid this?

If we use only a single, long run, and break up our samples into $n$ blocks, each of these can form a sample $L_i$. 
Stopping rules (3)

What could go wrong with this?

Correlation between successive blocks could mean that we have biased samples.

If the block size is large then the correlation should be small.

Explicit techniques exist to estimate the correlation and obtain the block size.

The simulation can be stopped once the estimate of $L$ becomes stable, or once the confidence interval around $L$ becomes sufficiently narrow.
Operational Analysis

Based on the idea of *observation* rather than a probabilistic description of system behaviour.

It is also concerned with quantities ‘directly related’ to these observed quantities.

Operational analysis makes very weak assumptions about the system being modelled — unlike simulation which requires detailed system knowledge, or the techniques from queuing theory (in the next section) which depend extensively on the probability distributions involved.

We begin by examining some *fundamental quantities* and *operational laws*.
Operational analysis (2)

We examine a system for some time recording the customer arrivals and departures, and define the quantities of interest:

- $T$, the length of time we observe the system
- $A$, the number of arrivals observed
- $C$, the number of departures (or completions of service) observed
- $W$, the job-time product: the sum of the durations of all customers over the observation period

If the system is a single resource, then we can also measure:

- $B$, the time for which the resource was busy.
Operational analysis (3)

From these we can define the following quantities:

- **Arrival rate**, \( \lambda := \frac{A}{T} \)
  
  — the mean number of arrivals per unit time

- **Throughput**, \( X := \frac{C}{T} \).
  
  — the mean number of departures per unit time

- **Mean number of customers**, \( N := \frac{W}{T} \)
  
  — the job-time product in terms of \( N \) and \( T \)

- **Mean residence time**, \( R := \frac{W}{C} \)
  
  — the job-time product in terms of \( R \) and \( C \)
Operational analysis (4)

For a single resource, we can also define:

- **Utilization**, $U := \frac{B}{T}$
  - the proportion of the time that the resource is busy

- **Average service requirement**, $S := \frac{B}{C}$
  - the mean time that the resource spends for each departure
The utilization law

Our first “law” is just an algebraic identity

\[ U := \frac{B}{T} = \frac{C B}{T C} = X S \]

This is termed the *utilization law*.

For example, if the throughput \((X)\) is 5 departures/sec and the service demand \((S)\) is 0.1 sec/departure then the utilization \((U)\) is 50%.
Little’s law

Similarly, we can derive the familiar Little’s Law

\[ N := \frac{W}{T} = \frac{C}{T} \frac{W}{C} = XR \]

For example, if the throughput \((X)\) is 5 customers/sec and the mean residence time is 1 sec then the average number in the system is 5.

- Very weak assumptions about the system
- Applicable to a wide range of systems
- Can be applied recursively to subsystems and to individual resources
  — but take care that mutually-consistent values are used for \(X\) and \(R\); in particular, whether they apply to the queue, the server or the entire system
An example

Observe system for $T = 10$ sec

4 customers spend 10 s in the system
One customer spends 5 s in the system

Then we have $W = 4 \times 10 + 1 \times 5 = 45$ s.

If also $A = C = 5$ then

$X = C/T = 5/10 = 0.5$ customers per second

$\lambda = 5/10 = 0.5$ customers per second

$N = 45/10 = 4.5$ customers

$R = 45/5 = 9.0$ s per customer
A simple interactive system

Fixed number, $M$, of users logged on.
Customer is at the terminal whilst thinking.
The *think time*, $Z$, is the average time a user spends between receiving a prompt and responding.
A customer not thinking is somewhere inside the central subsystem.
simple interactive system (2)

We can use Little’s Law to relate some observable quantities in the central subsystem:

- $N$ is the number of customers in the central subsystem $0 \leq N \leq M$
- $X$ is the rate at which customers complete in the central subsystem
- $R$ is the average time a customer spends in the central subsystem (intuitively equivalent to “response time”)

If we observe that system throughput is 0.5 interactions per second and we find on average 7.5 users in the subsystem then

$$R = \frac{N}{X} = \frac{7.5}{0.5} = 15 \text{ s}$$

from Little’s Law.
simple interactive system (3)

We can also apply Little’s law to the entire system.

This is a closed system so the number of customers is fixed as $M$.

We can split the time spent during an interaction into the response time ($R$) and the think time ($Z$). The residence time is $R + Z$.

We consequently derive the interactive system version of Little’s Law:

$$M = X(R + Z)$$
simple interactive system (4)

With 10 users logged on, 5 s average think time and an average response time of 15 s,

\[ X = \frac{M}{R + Z} = \frac{10}{15 + 5} = 0.5 \text{ interactions/sec} \]

Under heavy load (M large or Z small)

\[ U \approx 1 \]

Using the Utilization Law the throughput
\[ X \approx \frac{1}{S} \] and hence

\[ R = \frac{M}{X} - Z \approx MS - Z \]

Thus, the response time grows approximately linearly with the number of users M.
Visit counts and forced flow

We now extend our notation to allow the modelling of multiple devices.

We will use subscripts $i = 1, 2, \ldots, K$ to identify each device, e.g. $X_i$ is the throughput at device $i$.

Assume that the service required by a customer is an inherent property of the customer not of the state of the system.

A visit count for a device is the number of completions at that device for every completion from the system

$$V_i := \frac{C_i}{C}$$

Where $C_i$ is the number of completions at device $i$.

Recall that in a feed-forward queueing network, $0 \leq V_i \leq 1$ because each customer visits a given device at most once.
The forced flow law

Since \( X = \frac{C}{T} \), we have that

\[
X_i = \frac{C_i}{T} = \frac{C_i}{C} \times \frac{C}{T} = V_i X
\]

the Forced Flow Law.

For example, if the throughput from the entire system is 20 customers per second and each customer visits a given device 3 times then the throughput of that device must be 60 completions per second.

If devices are load independent, then define the service demand a customer makes on a device \( i \) by

\[
D_i := V_i S_i
\]

— be careful to distinguish the service requirement \( (S_i) \) from the service demand \( (D_i) \)
Queue lengths at a server

Applying the utilization law at each device:

\[ U_i = X_i S_i = (XV_i)S_i = X(V_iS_i) = XD_i \]

Similarly, applying Little’s law at each device:

\[ N_i = X_i R_i \]

\( R_i \) is the residence time at device \( i \) and can be decomposed into the time spent queuing and the time spent in service, approximated by \( R_i^* \):

\[
R_i^* = N_i S_i + S_i \\
= R_i^* X_i S_i + S_i \\
= R_i^* U_i + S_i
\]
Queue lengths at a server (2)

Hence

\[ R_i^* = \frac{S_i}{1 - U_i} \]

So that

\[ N_i = X_i R_i^* \]

\[ = \frac{X_i S_i}{1 - U_i} \]

\[ = \frac{U_i}{1 - U_i} \]

Observe that

- \( N_i \) is zero when \( U_i \) is zero
- \( N_i \) grows rapidly without bound as \( U_i \) approaches one
Bottleneck analysis

A bottleneck in a system is a hindrance to movement or progress.

Given the forced flow assumption, at high loads system performance is determined by the device with the highest utilization: *the bottleneck*.

The ratio of the completion rates of any two devices is

\[
\frac{X_i}{X_j} = \frac{V_i X}{V_j X} = \frac{V_i}{V_j}
\]

Since \( U_i = X_i S_i \), we have a similar property for utilizations

\[
\frac{U_i}{U_j} = \frac{X_i S_i}{X_j S_j} = \frac{V_i S_i}{V_j S_j}
\]
Bottleneck analysis (2)

A system is *load independent* if

- \( V_i \) are intrinsic properties of customers,
- \( S_i \) are independent of the queue length at \( i \)

In such cases, the throughput and utilization ratios are the same for all loads.

This can be used to determine asymptotes for \( X \) and \( R \).

In general, \( U_i \leq 1 \) and \( X_i \leq \frac{1}{S_i} \).

A device \( i \) becomes *saturated* as \( U_i \to 1 \)

Thus, as \( U_i \to 1 \), we have that \( X_i \to \frac{1}{S_i} \): device \( i \) is working as fast as it can and consequently serves one customer every \( S_i \).
Bottleneck analysis (3)

We use the subscript $b$ to denote a device capable of saturating.

Since the utilization ratios are fixed, the device $i$ with the largest $V_i S_i$ product will be the first to achieve 100% utilization as $N$ increases:

$$V_b S_b = \max\{V_1 S_1, ..., V_K S_K\}$$

so the bottleneck is determined by both the device and workload (the $V_i$ and $S_i$) properties.
By the forced flow law

\[ X = \frac{X_b}{V_b} \]

So, as \( U_b \to 1 \) and \( X_b \to 1/S_b \)

\[ X_{\text{max}} = \frac{X_b}{V_b} \to \frac{1}{V_bS_b} \]
The total per-customer service required is

\[ R_{\text{min}} = \sum_{i=1}^{K} V_i S_i \]

\[ \Rightarrow X \leq \frac{N}{R_{\text{min}}} \]

\( R_{\text{min}} \) denotes the smallest possible value of mean response time, occurring when \( N = 1 \).
Maximum throughput (3)

If $k \leq K$ jobs always avoid each other then

$$X = \frac{k}{R_{\text{min}}} \leq \frac{1}{V_b S_b}$$

$$k \leq \frac{R_{\text{min}}}{V_b S_b} = \sum_{i=1}^{K} \frac{V_i S_i}{V_b S_b} = N^*, \quad \text{say}$$

So, beyond $N^*$ queueing is certain.
Maximum throughput (4)

- It stays below $1/(V_b S_b)$ because, at that point, a bottleneck will be operating at maximum utilization.

- It stays below the straight line $X = N/R_{min}$ because the throughput is limited by the number of customers in service.
Interactive response time

- **X** throughput
- **M** terminals
- Average think time **Z**
- Recall the interactive system version of Little’s law:

\[ R = \frac{M}{X} - Z \]

By considering a bottleneck device **b**:

\[ X \leq \frac{1}{V_b S_b} \]

\[ \Rightarrow R \geq M V_b S_b - Z \]

\[ \Rightarrow R \geq M V_i S_i - Z \quad \forall i \in \{1..K\} \]
The response time asymptote meets the horizontal axis at

\[ M_b = \frac{Z}{V_b S_b} \]

It intersects the minimum response time \( R_{\text{min}} \) at \( M_b^* \) (say) where

\[ M_b^* V_b S_b - Z = R_{\text{min}} \]
Interactive response time (3)

Thus

\[ M_b^* = \frac{R_{min} + Z}{V_1S_1} = N^* + M_b \]

When there are more than \( M_b^* \) terminals, queueing is *certain* to exist.
Summary

- The largest of the products $V_i S_i$ determines the bottleneck $b$.

- The sum of these products determines the smallest possible response time $R_{\text{min}}$.

- Queueing cannot be avoided when $N$ exceeds

$$N^* = \frac{R_{\text{min}}}{V_b S_b}$$

- Queueing cannot be avoided in an interactive system when the number of logged-on terminals exceeds

$$M^*_b = N^* + \frac{Z}{V_b S_b}$$
Example: interactive system

Suppose that $Z = 20$ s.

<table>
<thead>
<tr>
<th>No.</th>
<th>device</th>
<th>$S_i$ (s)</th>
<th>$V_i$</th>
<th>$D_i = V_iS_i$</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>CPU</td>
<td>0.05</td>
<td>20</td>
<td>1.00</td>
</tr>
<tr>
<td>2</td>
<td>disk</td>
<td>0.08</td>
<td>11</td>
<td>0.88</td>
</tr>
<tr>
<td>3</td>
<td>fast disk</td>
<td>0.04</td>
<td>8</td>
<td>0.32</td>
</tr>
</tbody>
</table>

$R_{\text{min}} = 2.20$

Question: Is a 8 second response time feasible with 30 users logged on? If not, what changes are required?
Example: interactive system (2)

\[ V_1 S_1 = 1s \quad \text{(bottleneck)} \]
\[ V_2 S_2 = 0.88s \]
\[ V_3 S_3 = 0.32s \]

\[ \Rightarrow R_{\text{min}} = \sum_{i=1}^{3} V_i S_i = 2.2s \]

For \( M = 30 \), the response time asymptote requires \( R \geq 30 \times 1 - 20 = 10 \) s.

So the answer is \textit{no}.
Example: interactive system (3)

To make a 8 second response time feasible, we need to speed up the CPU, so that the new service time obeys the condition

\[ MV_1 S'_1 - Z \leq 8 \]

or

\[ S'_1 \leq \frac{20 + 8}{30 \times 20} = .047 \text{ s} \]

which is a 7% speed up in the CPU.

Then \( V_1 S'_1 = 0.93 \) is still the largest product so the CPU is still the bottleneck.
**Example: interactive system (4)**

*Question:* Is a 10s response time feasible when 50 users are logged on? If not, how much CPU speedup is required?

If $S_1 \to 0$, the disk will become bottleneck

$$R \geq MV_2 S_2 - Z$$

For $M = 50$, this is

$$R \geq 50 \times 0.88 - 20 = 24 \text{ s}$$

so a 10s response time is not feasible with $M = 50$ and no amount of CPU speedup is capable of achieving it.
Balanced system bounds

Balanced system bounds provide *tighter* bounds at mid-range loads than bottleneck analysis.

A system is *balanced* if for any load the utilizations of all devices are equal.

Balanced systems exhibit the following important property

\[
U_i(N) = \frac{N}{N + K - 1}
\]

So the system throughput is given by

\[
X(N) = \frac{U_i}{D_i} = \frac{N}{N + K - 1} \times \frac{1}{D_i}
\]
Balanced system bounds (2)

For example,

\[ N = 1 \quad K = 2 \quad U_1 = U_2 = \frac{1}{2} \]
\[ N = 1 \quad K = 100 \quad U_1 = \ldots = U_{100} = \frac{1}{100} \]
\[ N = 100 \quad K = 2 \quad U_1 = U_2 = \frac{100}{101} \]
\[ N = 2 \quad K = 2 \quad U_1 = U_2 = \frac{2}{3} \]

We observe the system to determine

\( D_{\max} \) — maximum demand at any device

\( D_{\min} \) — minimum device demand

\( D_{\text{av}} \) — average demand at each device

\( D = D_{\text{tot}} \) — total demand across all devices

So, we have \( D_{\text{av}} = D/K \)
Balanced system bounds (3)

Consider imaginary balanced systems related to our system

**pess1:** balanced system with $K$ devices each with a demand of $D_{\text{max}}$

**opt1:** balanced system with $K$ devices each with a demand of $D_{\text{min}}$

The throughput of system **pess1** is

$$\frac{N}{N + K - 1} \times \frac{1}{D_{\text{max}}}$$

The throughput of system **opt1** is

$$\frac{N}{N + K - 1} \times \frac{1}{D_{\text{min}}}$$
Balanced system bounds (4)

So for the system being modelled we have

\[
\frac{N}{N + K - 1} \times \frac{1}{D_{\text{max}}} \leq X(N) \leq \frac{N}{N + K - 1} \times \frac{1}{D_{\text{min}}}
\]

And since \( N = XR \),

\[
(N + K - 1)D_{\text{max}} \geq R(N) \geq (N + K - 1)D_{\text{min}}
\]

This will give tighter bounds on the mid-range performance than the bottleneck bounds.
Balanced system bounds (5)

We can do even better by considering the best performance that the system can achieve which occurs when the load is spread out evenly among all the devices.

**opt2:** $D_{av}$ at each of the $K$ devices

$$X(N) = \frac{N}{N + K - 1} \times \frac{1}{D_{av}} = \frac{N}{D + (N - 1)D_{av}}$$
Balanced system bounds (6)

Now, what is the worst system subject to the constraints that $D$ and $D_{\text{max}}$ remain fixed?

**Answer:** place $D_{\text{max}}$ at as many devices as possible and 0 at the rest

**pess2:** $D_{\text{max}}$ at each of the $\frac{D}{D_{\text{max}}}$ devices

$$X(N) = \frac{N}{N + K - 1} \times \frac{1}{D_{\text{max}}} = \frac{N}{D + (N - 1)D_{\text{max}}}$$
Balanced system bounds (7)

\[
\frac{N}{D + (N - 1)D_{\text{max}}} \leq X(N) \leq \frac{N}{D + (N - 1)D_{\text{av}}}
\]

\[
D + (N - 1)D_{\text{max}} \geq R(N) \geq D + (N - 1)D_{\text{av}}
\]
Balanced system bounds (8)

As $N \to \infty$,

$$\frac{N}{D + (N - 1)D_{\text{max}}} \to \frac{1}{D_{\text{max}}}$$

Note that at high loads the bottleneck bounds are the limiting high bound.

The asymptotic bottleneck bound $\frac{1}{D_{\text{max}}}$ and the optimistic balanced bound intersect at $N^\dagger$ where

$$\frac{1}{D_{\text{max}}} = \frac{N^\dagger}{D + (N^\dagger - 1)D_{\text{av}}}$$

So that

$$N^\dagger = \frac{D - D_{\text{av}}}{D_{\text{max}} - D_{\text{av}}}$$
Markov processes

A Markov process is a family of random variables $X(t)$ indexed by a time parameter $t$ in which

- $X(t)$ denotes the state at time $t$
- The next state depends only on the current state

We are concerned with the particular case of Markov processes in which the state space is discrete, termed Markov chains

If we denote successive states as $x_1, x_2, \ldots$ then the Markov property says that

$$
P(X_{n+1} = x_{n+1} | X_1 = x_1, \ldots, X_n = x_n) = P(X_{n+1} = x_{n+1} | X_n = x_n)
$$
Other types of stochastic process

A Markov process is a particular kind of stochastic process

- **Birth death process** (BDP): a Markov process in which transitions occur only between neighbouring states
- **Random Walks**: refer to Probability 1A notes
- **Renewal Process**: $X(t)$ counts state transitions in $(0, t)$
Birth death processes

A *birth death process* is a special case of a continuous-time Markov chain in which we allow transitions only between neighbouring states.

The state space is the set of non-negative integers and we allow only a single birth or death per transition. Thus given $X_n = i$

$$X_{n+1} = \begin{cases} 
    i + 1 & \text{birth (or arrival)} \\
    i - 1 & \text{death (or departure)}
\end{cases}$$

We use $\lambda_i$ to represent the birth rate in state $i$, and $\mu_i$ to represent the death rate in state $n$. 

![Diagram of birth death processes](image-url)
Time dependent solution of BDP

We denote by $P_i(t)$ the probability of being in state $i$ at time $t$.

Probability of a birth in an interval $\Delta t$ when the system starts in state $i$ is assumed to be $\lambda_i \Delta t + o(\Delta t)$.

Probability of a death in $\Delta t$ when the system starts in state $i$ is $\mu_i \Delta t + o(\Delta t)$.

Probability of $>1$ event in $\Delta t$ is $o(\Delta t)$.

$o(\Delta t)$ denotes a quantity which becomes negligible when compared with $\Delta t$ as $\Delta t \rightarrow 0$.

To solve for $P_i(t)$ we write a set of difference equations called the Chapman Kolmogorov equations.
Chapman Kolmogorov equations

For $i \geq 1$:

$$
P_i(t + \Delta t) = P_i(t)(1 - \lambda_i \Delta t)(1 - \mu_i \Delta t)$$

$$+ P_{i+1}(t)(\mu_{i+1} \Delta t)(1 - \lambda_{i+1} \Delta t)$$

$$+ P_{i-1}(t)(\lambda_{i-1} \Delta t)(1 - \mu_{i-1} \Delta t)$$

$$+ o(\Delta t)$$

For $i = 0$:

$$
P_0(t + \Delta t) = P_0(t)(1 - \lambda_0 \Delta t)$$

$$+ P_1(t)(\mu_1 \Delta t)(1 - \lambda_1 \Delta t)$$

$$+ o(\Delta t)$$
Chapman Kolmogorov equations (2)

We derive differential-difference equations from these by dividing through by $\Delta t$ and taking the limit as $\Delta t \to 0$

$$\frac{dP_i(t)}{dt} = -(\lambda_i + \mu_i)P_i(t) + \mu_{i+1}P_{i+1}(t) + \lambda_{i-1}P_{i-1}(t)$$

for $i \geq 1$ and

$$\frac{dP_0(t)}{dt} = -\lambda_0P_0(t) + \mu_1P_1(t)$$

The time dependent solution is difficult for systems of interest, so we will study the stationary solution.
Stationary solution

We are interested in the long term probabilities after the system has reached an *equilibrium*. These probabilities are independent of the initial conditions.

System reaches equilibrium if, for all $i$,

$$\lim_{t \to \infty} P_i(t) = p_i \quad \text{exists}.$$ 

The quantities $p_i$ solve the Chapman Kolmogorov equations with $dP_i(t)/dt = 0$ so that

$$0 = -(\lambda_i + \mu_i)p_i + \mu_{i+1}p_{i+1} + \lambda_{i-1}p_{i-1}$$

$$0 = -\lambda_0 p_0 + \mu_1 p_1$$

Rewriting gives

$$p_{i+1} = \frac{\lambda_i + \mu_i}{\mu_{i+1}} p_i - \frac{\lambda_{i-1}}{\mu_{i+1}} p_{i-1} \quad (i \geq 1)$$

$$p_1 = \frac{\lambda_0}{\mu_1} p_0$$
Stochastic balance

Under steady-state conditions we require total flow into a state to equal total flow out of a state.

The total flow is the product of the steady state probabilities and the flow rates.

We enter state \( i \) at rate \( p_{i-1} \lambda_{i-1} + p_{i+1} \mu_{i+1} \).

We exit state \( i \) at rate \( p_i \lambda_i + p_i \mu_i \).

The equation

\[
p_{i-1} \lambda_{i-1} + p_{i+1} \mu_{i+1} = p_i \lambda_i + p_i \mu_i \quad (i \geq 1)
\]
equates the flow into and out of state \( i \).

This is called the *global balance* equation.
Stochastic balance (2)

The two equations

\[ p_i \lambda_i = p_{i+1} \mu_{i+1} \quad (i \geq 0) \]

describing flow from state \( i \) to state \( i + 1 \), and

\[ p_i \mu_i = p_{i-1} \lambda_{i-1} \quad (i \geq 1) \]

describing flow from state \( i \) to state \( i - 1 \) are the detailed balance equations.

Rewriting gives

\[ p_{i+1} = \frac{\lambda_i}{\mu_{i+1}} p_i \]

which gives us the product solution

\[ p_k = p_0 \prod_{i=0}^{k-1} \frac{\lambda_i}{\mu_{i+1}} \quad \text{for} \quad k \geq 1 \]

for \( p_k \ (k \geq 1) \) in terms of \( p_0 \).
Stochastic balance (3)

Since the sum of state probabilities must be unity,

\[ p_0 + \sum_{k=1}^{\infty} p_k = 1 \]
\[ p_0 + \sum_{k=1}^{\infty} p_0 \prod_{i=0}^{k-1} \frac{\lambda_i}{\mu_{i+1}} = 1 \]
\[ p_0 \left[ 1 + \sum_{k=1}^{\infty} \prod_{i=0}^{k-1} \frac{\lambda_i}{\mu_{i+1}} \right] = 1 \]

so that

\[ p_0 = \left[ 1 + \sum_{k=1}^{\infty} \prod_{i=0}^{k-1} \frac{\lambda_i}{\mu_{i+1}} \right]^{-1} \]
\[ p_k = p_0 \prod_{i=0}^{k-1} \frac{\lambda_i}{\mu_{i+1}} \]

which are known as the general flow balance equations.
The $M/M/1$ queue

The BDP maps well onto our domain of study — queueing systems.

Births represent arrivals to queue, deaths represent departures as customers finish service.

The $M/M/1$ queue is an infinite customer system, with infinite waiting room, and a state independent service rate.

This means that $\lambda_i = \lambda$ and $\mu_i = \mu$ for all $i$ and we can solve the balance equations as

$$p_k = p_0 \prod_{i=0}^{k-1} \frac{\lambda}{\mu}$$

$$= p_0 \left( \frac{\lambda}{\mu} \right)^k.$$
The $M/M/1$ queue (2)

Writing $\rho = \frac{\lambda}{\mu}$

$$p_0 = \frac{1}{1 + \sum_{k=1}^{\infty} \rho^k}$$

$$= \frac{1}{1 + \rho \sum_{k=0}^{\infty} \rho^k}$$

$$= \frac{1}{1 + \rho \left( \frac{1}{1-\rho} \right)}$$

$$= 1 - \rho$$

Consequently, the number in the system is geometrically distributed

$$p_k = (1 - \rho)\rho^k, \quad k = 0, 1, 2, \ldots$$

If $\rho > 1$, i.e. if $\lambda > \mu$ the system will not reach equilibrium.
The $M/M/1$ queue (3)

What is the average number of customers, $\bar{N}$, in the system?

\[
\bar{N} = \sum_{k=0}^{\infty} kp_k = \sum_{k=0}^{\infty} k(1 - \rho)\rho^k = (1 - \rho)\rho \frac{\partial}{\partial \rho} \left( \sum_{k=0}^{\infty} \rho^k \right) = (1 - \rho)\rho \frac{\partial}{\partial \rho} \left( \frac{1}{1 - \rho} \right) = \frac{\rho}{1 - \rho}
\]
The $M/M/1$ queue (4)

An arriving customer will find, on average $\overline{N}$ in the system, and will spend a time, say $\overline{T}$, in the system. During $\overline{T}$ there will be, on average $\lambda\overline{T}$ arrivals, leaving $\overline{N}$ customers in the queue. Thus

$$\overline{N} = \lambda\overline{T}$$

which is Little’s result restated. In our case

$$\overline{T} = \frac{\overline{N}}{\lambda} = \frac{\rho}{\lambda(1 - \rho)} = \frac{1}{\mu(1 - \rho)} = \frac{1}{\mu - \lambda}$$

which is the $M/M/1$ average response time.

Note that

$\gg \frac{1}{\mu}$ is the average service time

$\gg \rho$ is the utilization
Performance at high load

At high utilizations $\rho$ approaches one and the response time and queue lengths are unbounded.

Expected number in the system

\[ \rho = \frac{\lambda}{\mu} \]
Response time

Consider the effect on response time by increasing the utilization by a constant load of 0.1.

<table>
<thead>
<tr>
<th>utilization ($\rho$)</th>
<th>response time ($\bar{T}$)</th>
</tr>
</thead>
<tbody>
<tr>
<td>old</td>
<td>new</td>
</tr>
<tr>
<td>0.1</td>
<td>0.2</td>
</tr>
<tr>
<td>0.5</td>
<td>0.6</td>
</tr>
<tr>
<td>0.8</td>
<td>0.9</td>
</tr>
</tbody>
</table>

Predicting response times is very difficult at high loads.

Running systems at maximum utilization may please the providers, but it doesn’t please the users.
$M/M/1$ — an example

H.R Cutt barber shop, no appointment needed!

Customers served FCFS.

On Saturday mornings he is very busy and is considering hiring a helper.

Measures the (Poisson) arrival rate of customers to be 5 per hour.

Customers are prepared to wait, and he spends on average 10 min per cut.

What are the average number of customers in the shop, the average number of customers waiting? What percentage of time can a customer receive a cut without waiting?

He has 4 seats in his waiting room. What is the probability that an arriving customer will find no seat and have to stand?
**M/M/1 example solution**

The average number of customers in the shop:

\[ \lambda = 5 \text{ per hour}, \mu = 6 \text{ per hour} \]

So

\[ \rho = \frac{5}{6}, \text{ and hence } \bar{N} = 5. \]

Since the average number of customers in service is \( \rho \), the utilization, the average number of customers waiting is

\[ \bar{N}_q = \bar{N} - \rho = 4 \frac{1}{6} \]

How likely is the barber to be idle?

\[ p_0 = 1 - \rho = \frac{1}{6} \]

How often is no seat free?

\[ P(\text{no seat}) = P(N \geq 5) = \rho^5 \approx 0.402 \]
**$M/M/m — m$ servers**

This is just like the $M/M/1$ system, except that there are $m$ servers.

For all $k$, $\lambda_k = \lambda$, but now the service rate is a function of the number of customers in the system

$$
\mu_k = \begin{cases} 
k\mu & \text{if } 0 < k \leq m \\
2m & \text{if } k > m
\end{cases}
$$

For an equilibrium distribution to exist we require that $\frac{\lambda}{m\mu} < 1$. 

![Diagram of $M/M/m$ system]
**$M/M/1/K$ — finite capacity**

The system can hold up to $K$ customers.

Now for $k \geq K$, $\lambda_k = 0$ and for $k > K$, $p_k = 0$.

Using the equations from the $M/M/1$ queueing system, but limiting the summation, and again writing $\rho = \frac{\lambda}{\mu}$,

\[
p_k = p_0 \rho^k \quad \text{for} \quad k \leq K
\]

\[
p_0 = \frac{1}{1 + \sum_{k=1}^{K} \rho^k}
\]

\[
= \frac{1}{1 + \frac{\rho - \rho^{K+1}}{1 - \rho}}
\]

\[
= \frac{1 - \rho}{1 - \rho^{K+1}}
\]

Note that $p_0$ is greater than in the $M/M/1$ case.

For this system with a finite state space an equilibrium distribution always exists whatever the arrival and departure rates.
$M/M/1//N$ — finite population

Single server, unbounded queue capacity and a population of $N$ customers.

We solve this system by modifying the $\lambda_k$ to model the arrival rate.

Instead of having an arrival rate for the population as a whole, we assign an arrival rate to each customer, say $\lambda$.

If there are no customers in the system, then all of them are eligible to be born, so that

$$\lambda_0 = N\lambda.$$

As we have more customers in the system, we have less eligible to be born. So that,

$$\lambda_k = (N - k)\lambda, \quad \text{for} \quad 0 \leq k \leq N.$$

With a single server the service rate is constant

$$\mu_k = \mu, \quad \text{for} \quad k \geq 1.$$
An application of this system is to model a link in a telephone network.

Such a link contains $m$ circuits each of which carries a single call.

Suppose that calls arrive at the link according to a Poisson process of rate $\lambda$.

If there is a free circuit the call is connected and holds the circuit for an exponentially distributed length of time, with mean $\frac{1}{\mu}$.

At the end of this holding period the call terminates and the circuit is freed.

If there are no free circuits then the call is lost.
\( M/M/m/m \)

\[ \lambda_k = \begin{cases} \lambda & k < m \\ 0 & k \geq m \end{cases} \]

\[ \mu_k = k\mu \quad \text{for} \quad 1 \leq k \leq m \]

The flow balance equations give for \( k \leq m \)

\[ p_k = p_0 \prod_{i=0}^{k-1} \frac{\lambda_i}{\mu_{i+1}} \]

\[ = p_0 \prod_{i=0}^{k-1} \frac{\lambda}{(i + 1)\mu} \]

\[ = p_0 \left( \frac{\lambda}{\mu} \right)^k \frac{1}{k!} \]
\( M/M/m/m \) (2)

Solving for \( p_0 \) gives

\[
\sum_{k=0}^{m} p_k = p_0 \sum_{k=0}^{m} \left( \frac{\lambda}{\mu} \right)^k \frac{1}{k!} = 1
\]

\[
\Rightarrow \ p_0 = \left[ \sum_{k=0}^{m} \left( \frac{\lambda}{\mu} \right)^k \frac{1}{k!} \right]^{-1}
\]

The probability that an arriving call finds all circuits occupied, \( p_m \), is called the *loss probability* for the telephone link. Thus,

\[
p_m = p_0 \left( \frac{\lambda}{\mu} \right)^m \frac{1}{m!}
\]

\[
= \left( \frac{\lambda}{\mu} \right)^m \frac{1}{m!} \left[ \sum_{k=0}^{m} \left( \frac{\lambda}{\mu} \right)^k \frac{1}{k!} \right]^{-1}
\]

This expression for the loss probability is known as *Erlang’s formula*. 
BDM extensions

First we relax our constraints on the arrival process distribution.

We want to model systems in which the coefficient of variation of the interarrival time is less than one.

Consider a system in which a birth occurs in two stages.

Each stage has an exponentially distributed residence time.
BDM extensions (2)

If the desired birth rate is $\lambda$, then let each stage have a rate $2\lambda$.

The average time to get through the combined birth process will be

$$\bar{\tau} = \frac{1}{2\lambda} + \frac{1}{2\lambda} = \frac{1}{\lambda}.$$ 

Since each stage has exponentially distributed residence times, the variance of each stage is

$$\sigma^2_{\text{single}} = \frac{1}{(2\lambda)^2}.$$ 

The two stages are independent, so the variance of $\tau$, the time to get through both stages is

$$\sigma^2_\tau = \frac{1}{(2\lambda)^2} + \frac{1}{(2\lambda)^2} = \frac{1}{2\lambda^2}.$$
BDM extensions (3)

So the coefficient of variation is

$$C_r = \frac{\sqrt{\frac{1}{2\lambda^2}}}{\frac{1}{\lambda}} = \frac{1}{\sqrt{2}}.$$  

In general, if we use $r$ stages each with rate $r\lambda$ we get an average time through all stages of $\frac{1}{\lambda}$ and a coefficient of variation of $\frac{1}{\sqrt{r}}$.

The distribution that describes this $r$-stage process is the Erlangian distribution, denoted $E_r$. 
Example, $M/E_2/1$

Allow the state of the process to represent the number of stages remaining to be served.

An arrival increases the number of stages remaining to be served by 2 and occurs at rate $\lambda$.

A departure from a stage reduces the number of stages to be served by 1 and occurs at rate $2\mu$. 

\[
M = E^2 = 1
\]
Parallel Servers

Combining stages in series reduces the coefficient of variation.

If, instead, we combine them in parallel with a probability $\alpha_i$ of choosing the $i^{th}$ parallel stage we get a service distribution with coefficient of variation larger than 1.

![Diagram of parallel servers with probabilities $\alpha_1$ and $\alpha_2$ and service rates $\mu_1$ and $\mu_2$.]

The coefficient of variation is given (see Kleinrock) by

$$C_r = \frac{2 \sum_{i=1}^{r} \frac{\alpha_i}{\mu_i^2}}{\left( \sum_{i=1}^{r} \frac{\alpha_i}{\mu_i} \right)^2} - 1 \geq 1.$$
Queueing Networks

We have seen the solution of several queueing systems of interest.

In general we want to solve a system of such queues representing a real world performance problem e.g. a distributed computing system.

We represent the system under study as a network of connected queues.

Customers move (instantaneously) between service centres where they receive service.
Model definition

➢ *customers*: typically these represent programs or data packets etc

➢ *service centres*: the resources in the system e.g. disks, CPU, transmission links

➢ *service time distributions*: may vary according to customer type and visit

➢ *load dependence*: multi-processor systems have load dependent service rates

➢ *waiting lines and scheduling*: may have limited capacity and various scheduling algorithms

➢ *customer types*: multiple customer classes may exist
Open Queueing Networks

Allow customers to arrive from an external source, with rate $\gamma$.

Customers may leave the network on completion.

Assume we have $N$ nodes, each a single server queue with infinite waiting room.

Each server $i$ has exponential service time with mean $1/\mu_i$.

Customers arrive as a Poisson stream at node $i$ at rate $\gamma_i$.

A customer completing at node $i$ moves to node $j$ with probability $q_{ij}$ for $(i, j = 1, 2, \ldots, N)$.

Note that

$$\sum_{j=1}^{N} q_{ij} \leq 1$$
Open Queueing Networks (2)

A job leaves the network from node $i$ with probability

$$q_{i0} = 1 - \sum_{j=1}^{N} q_{ij}.$$  

The probabilities $q_{ij}$ are called the *routing probabilities*.

Written as an $N \times N$ matrix this is called the *routing matrix* $Q = (q_{ij})$.

An open network with parameters $\gamma_i, \mu_i$ and $Q$ is called a *Jackson network*.

The system state is $(k_1, k_2, \ldots, k_N)$, where $k_i$ is the number of jobs present at node $i$. 
Steady state solution

Let $\lambda_j$ be the average number of arrivals to node $j$ per unit time.

On average the departure count per unit time is therefore $\lambda_j$.

A fraction $q_{ji}$ go to node $i$.

The rate of traffic from node $j$ to node $i$ is thus $\lambda_j q_{ji}$. 

\[ \begin{align*} 
\gamma_i & \rightarrow \lambda_i \\
\lambda_1 q_{1i} & \rightarrow \lambda_i \\
\lambda_2 q_{2i} & \rightarrow \lambda_i \\
\lambda_N q_{Ni} & \rightarrow \lambda_i \\
\lambda_i q_{i0} & \rightarrow \lambda_i \\
\lambda_i q_{i1} & \rightarrow \lambda_i \\
\lambda_i q_{i2} & \rightarrow \lambda_i \\
\vdots & \rightarrow \lambda_i \\
\lambda_i q_{iN} & \rightarrow \lambda_i 
\end{align*} \]
Traffic equations

Adding together all contributions,

$$\lambda_i = \gamma_i + \sum_{j=1}^{N} \lambda_j q_{ji} \quad i = 1, 2, \ldots, N.$$ 

These are known as the *traffic equations*.

A necessary and sufficient condition for the existence of an equilibrium distribution is that

$$\rho_i := \frac{\lambda_i}{\mu_i} < 1$$

where $\lambda_i$ is the solution of the traffic equations.
**Jackson’s Theorem**

Let \( p(k_1, k_2, \ldots, k_N) \) be the steady state equilibrium distribution. Then *Jackson’s Theorem* states that

\[
p(k_1, k_2, \ldots, k_N) = p_1(k_1)p_2(k_2) \cdots p_N(k_N)
\]

where \( p_i(k_i) \) is the equilibrium distribution that there are \( k_i \) jobs in an \( M/M/1 \) queue with traffic intensity \( \rho_i \).

Jackson’s theorem has some important implications.

- The numbers of jobs at the various nodes are independent.

- Node \( i \) behaves as if subjected to a Poisson stream with rate \( \lambda_i \).

Jackson’s theorem may be generalized to the case where node \( i \) has \( n_i \) servers and so the nodes behave as independent \( M/M/n_i \) queues.
Closed Queueing Networks

Frequently used to model systems at high load or where a limited, constant number of jobs is admitted to service.

No external arrivals or departures.

Now the routing probabilities satisfy

\[ \sum_{j=1}^{N} q_{ij} = 1, \quad i = 1, 2, \ldots, N \]

The number of jobs in the system is always a constant, denoted by \( K \).

The states of the system, described by the vector \((k_1, k_2, \ldots, k_N)\), thus satisfy the constraint,

\[ \sum_{i=1}^{N} k_i = K. \]
Closed Queueing Networks (2)

The state space, $S$, is then finite. The number of states is

$$\binom{K + N - 1}{N - 1}$$

The traffic equations become

$$\lambda_i = \sum_{j=1}^{N} \lambda_j q_{ji}, \quad i = 1, 2, \ldots, N.$$  

With a finite state space there always exists an equilibrium distribution.

Analogous to Jackson’s theorem for the open network case it may be shown that

$$p(k_1, k_2, \ldots, k_N) = \frac{1}{G} r_1(k_1)r_2(k_2)\cdots r_N(k_N)$$

where $r_i(k_i)$ is the probability that there are $k_i$ jobs in an $M/M/1$ queue with traffic intensity given by a solution to the traffic equations.
Open vs closed networks

The normalization constant $G$ has to be determined by

$$G = \sum_{s\in S} r_1(k_1)r_2(k_2)\cdots r_N(k_N)$$

obtained by summing over all states $s = (k_1, k_2, \ldots, k_n)$ in the state space $S$.

With closed networks need to compute the normalization constant $G$ — a computationally hard problem.

The constraint $\sum_{i=1}^N k_i = K$ means that the numbers of jobs in the various queues are no longer independent.

For instance, consider the extreme case where all $K$ jobs are at one node.
Derivation of the size of $S$

We require to show that

$$p(K, N) := \binom{K + N - 1}{N - 1}$$

is the number of (ordered) partitions of a positive integer $K$ into $N$ integer summands

$$K = \sum_{i=1}^{N} k_i.$$

**Proof**

Consider $K + N - 1$ boxes aligned in a row and select $N - 1$ of these boxes (without replacement) which can be done in $p(K, N)$ ways. Place a “/” symbol in each of the boxes and a “1” in each of the other boxes. The boxes now represent an (ordered) partition of $K$ into $N$ groups of “1” which when added together give the $k_i$ summands.
Aside: application to balanced systems

Recall that in a balanced system with $N$ jobs and $K$ devices the common utilization at each device is given by

$$U_i(N) = \frac{N}{N + K - 1} \quad \text{for } i = 1, \ldots, K.$$

But using the function $p(K, N)$ we can see that, by symmetry, the utilization is given by

$$U_i(N) = 1 - \frac{p(N, K - 1)}{p(N, K)}$$

$$= 1 - \left( \frac{N+K-2}{K-1} \right) \left( \frac{N+K-1}{k-1} \right)$$

$$= 1 - \frac{(N + K - 2)!}{(K - 2)!N!} \cdot \frac{N!(K - 1)!}{(N + K - 1)!}$$

$$= 1 - \frac{K - 1}{N + K - 1} \cdot \frac{N}{N + K - 1}.$$
The $M/G/1$ queue

It is usually easier to justify the memoryless property for arrivals than for service times.

For arrivals, we can appeal to asymptotic results on the merging of large numbers of independent streams to help justify the memoryless property for arrivals.

For service times, it is easy to think of examples where the service times have a quite different distribution to the exponential. For example, the service times might be constant corresponding to certain packet lengths in a communication network.

This leads to an interest in the $M/G/1$ queue with general service times given by

CDF $B(x) = \mathbb{P}(\text{service time} \leq x)$. 
(Lack of) Markov property

With general service times we no longer find that \( X(t) \), the number of customers in the system at time \( t \), has the Markov property.

This follows since the future evolution of \( X(t) \) now depends not just on the number present but on the remaining service time of the customer (if any) currently in service.

Recall, that in the \( \cdot/M/\cdot \) case the remaining service time always has the same memoryless distribution whenever we observe the queue.
Embedded Markov Chain

It would be possible to formulate a model for the $M/G/1$ queue using a state variable with two components $(n, x)$ where $n$ is the number present and $x$ is the remaining service time, if any, of the customer in service. This augmented model does have the Markov property and can be analyzed directly.

Instead, it is possible to pick out a discrete set of times where the Markov property holds and build a model on this discrete time Markov Chain. Such a set of times is given by $t_i$ ($i = 1, 2, \ldots$) where $t_i$ is the time of the $i^{th}$ departure from the queue. There is no remaining service time to worry us at these time instants.

Thus, $X(t_i)$, $i = 1, 2, \ldots$ is a Markov Chain embedded in the stochastic process $X(t)$. 

_Compurer Systems Modelling, Nov 2001_
Performance measures

The determination of a full description of the $M/G/1$ model is possible but difficult. Instead, we shall look at some steady state performance measures.

Let $1/\mu$ be the mean service time of a customer in the $M/G/1$ queue, obtained from the CDF of the service time distribution $B(\cdot)$, say.

Then the mean queueing time, $\overline{T_q}$, of a customer before it receives service is given by

$$\overline{T_q} = \overline{N_q} \frac{1}{\mu} + \rho \overline{x}$$

where $\overline{N_q}$ is the average number of customers waiting in the queue at the time of arrival, $\overline{x}$ is the average remaining service time of the customer, if any, in service and $\rho = \lambda/\mu$ is the traffic intensity which is also the utilization of the server.
Remaining service time

A result from renewal theory is that $\bar{x} = \mu s^2 / 2$.

Notice that this involves the 2nd central moment of the service time distribution, $s^2$.

For the exponential case, $s^2 = 1/\mu^2$ so that $\bar{x} = 1/(2\mu)$ as might be intuitively expected.
Performance measures (2)

From Little’s law,

\[ N_q = \lambda T_q \]

and so

\[ T_q = \frac{1}{\mu} + \rho x \]

\[ = \frac{\rho x}{1 - \rho} \]

\[ = \frac{\rho \mu s^2}{2(1 - \rho)} \]

\[ = \frac{\lambda s^2}{2(1 - \rho)}. \]
Performance measures (3)

Let $C_s$ be the coefficient of variation of the service time distribution then

\[ C_s^2 = \frac{s^2}{(\mathbb{E}(s))^2} - 1 \]

where $\mathbb{E}(s) = 1/\mu$ so

\[ \frac{s^2}{\mathbb{E}(s)} = \frac{(1 + C_s^2)}{\mu^2} \]

Hence,

\[ T_q = \frac{\lambda(1 + C_s^2)}{\mu^2 2(1 - \rho)} \]

\[ = \frac{\rho(1 + C_s^2)}{2\mu(1 - \rho)} . \]
Pollaczek-Khintchine formula

Consider now the total time, \( \bar{T} \), for a customer to pass through the system given by their waiting time in the queue and their own service time.

Thus,

\[
\bar{T} = \bar{T}_q + \frac{1}{\mu}
\]

\[
= \frac{1}{\mu} \left( 1 + \frac{\rho(1 + C^2_s)}{2(1 - \rho)} \right) .
\]

Using Little’s law for the entire system we can now find, \( \bar{N} \), the mean number of customers in an \( M/G/1 \) queueing system by

\[
\bar{N} = \lambda \bar{T}
\]

\[
= \rho + \frac{\rho^2(1 + C^2_s)}{2(1 - \rho)}
\]

This is known as the \textit{Pollaczek-Khintchine} (PK) formula.
PK formula

The Pollaczek-Khintchine formula tells us that the mean number of customers is determined not only by the mean interarrival and mean service times but also by the coefficient of variation of the service time distribution, $C_s$.

There are several cases.

- $C_s = 0$: this is the case of constant service times. For example, in ATM networks where the cells (that is, the packets) are of fixed length (53 bytes).

- $C_s < 1$: this is the case where the variability is less than in the case of exponential service times, thus the $M/M/1$ model will be conservative in its performance estimates.
PK formula (2)

$C_s \approx 1$: this is where the $M/M/1$ model works best and many systems correspond to this model. For example, batch jobs on a mainframe.

$C_s > 1$: this is the case where the $M/G/1$ model is required. An example, is the observed packet lengths in Internet traffic. The distribution of packet sizes (and hence service times) is often found to be bimodal with many small packets and many longer packets of length determined by the MTU.