

BOUNDS FOR THE COUPLING TIME IN QUEUEING NETWORKS PERFECT SIMULATION*

JANTIEN G. DOPPER, BRUNO GAUJAL AND JEAN-MARC VINCENT †

Abstract. In this paper, the duration of perfect simulations for Markovian finite capacity queuing networks is studied. This corresponds to hitting time (or coupling time) problems in a Markov chain over the Cartesian product of the state space of each queue. We establish an analytical formula for the expected simulation time in the one queue case which provides simple bounds for acyclic networks of queues with losses. These bounds correspond to sums on the coupling time for each queue and are either almost linear in the queue capacities under light or heavy traffic assumptions or quadratic, when service and arrival rates are similar.

Key words. Perfect simulation, Markov chain, Hitting time

AMS subject classifications. 60J10, 60J22, 65C40, 65C20, 68U20

1. Introduction. Markov chains are an important tool in modelling systems. Amongst others, Markov chains are being used in the theory of queueing systems, which itself is used in a variety of applications as performance evaluation of computer systems and communication networks. In modelling any queueing system, one of the main points of interest is sampling the behavior of the system in the long run. For an irreducible, ergodic (i.e. aperiodic and positive-recurrent) Markov chain with probability matrix P , this long run behavior follows the stationary distribution of the chain given by the unique vector π which satisfies the linear system $\pi = \pi P$. However, it may be hard to compute this stationary distribution, especially when the finite state space is huge which is frequent in queueing models. In that case, several approaches have been proposed to get samples of the long run behavior of the system.

The most classical methods are indirect. They consists in first computing an estimation of π and then sample according to this distribution (by classical methods such as p.d.f. inverse, rejection or aliasing).

Estimating π can be done through efficient *numerical iterative methods* solving the linear system $\pi = \pi P$. [10]. Even if they converge fast, they do not scale when the state space (and thus P) grows.

Another approach to estimate π is to use a *regenerative simulation* [4, 7] based on the fact that on a trajectory of a Markov chain returning to its original state, the frequency of the visits to each state is steady state distributed. This technique does not suffer from statistical bias but is very sensitive to the return time to the regenerative state. This means that the choice of the initial state is crucial but also that regenerative simulation complexity increases fast with the state space which is exponential in the number of queues. This can be partially overcome by using importance sampling [3] or semi-regenerative simulation [2]. However, the simulation times still typically exhibit a multiplicative behavior with the number of queues.

There also exist direct techniques to sample states of Markov chain according to its stationary distribution. The classical method has been *Monte Carlo simulation* for many years. This method is based on the fact that for an irreducible aperiodic finite

*This work was partially supported by the French ACI SurePath project and SMS ANR

†Jantien Dopper, Mathematical institute, Leiden University, NL (jgdopper@math.leidenuniv.nl) (this author was partially supported by a grant from INRIA), Bruno Gaujal and Jean-Marc Vincent, Laboratoire ID-IMAG, Mescal Project INRIA-UJF-CNRS-INPG, 51, avenue Jean Kuntzmann, F-38330 Montbonnot, France, e-mail : ({Bruno.Gaujal, Jean-Marc.Vincent}@imag.fr)

Markov chain with initial distribution $\pi^{(0)}$, the distribution $\pi^{(n)}$ of the chain at time n converges to π as n gets very large:

$$\lim_{n \rightarrow \infty} \pi^{(n)} = \lim_{n \rightarrow \infty} \pi^{(0)} P^n = \pi.$$

After running the Markov chain long enough, the state of the chain will not depend on the initial state anymore. However, the question is how long is long enough? That is, when is n sufficiently large so that $|\pi^{(n)} - \pi| \leq \epsilon$ for a certain $\epsilon > 0$? Moreover, the samples generated by this method will always be biased.

In 1996, Propp and Wilson[8] solved these problems for Markov chain simulation by proposing an algorithm which returns exact samples of the stationary distribution very fast. The striking difference between Monte Carlo simulation and this new algorithm is that Propp and Wilson do not simulate into the future, but go backwards in time. The main idea is, while going backwards in time, to run several simulations, starting with all $s \in S$ until the state at $t = 0$ is the same for all of them. If the output is the same for all runs, then the chain has coupled. Because of this coupling property and going backwards, this algorithm has been called Coupling From The Past (from now on: CFTP). A more detailed description of this algorithm will be presented in section 2.

When the coupling from the past technique is applicable, one gets in a finite time one state with steady-state distribution. Then one can use either a one long-run simulation from this state avoiding the estimation of the initial transient problem or replicate independently the CFTP algorithm to get a sample of independent steady-state distributed variables. The analysis of the choice could be done exactly as in [1]. The replication technique has been applied successfully in finite capacity queueing networks with blocking and rejection (very large state-space) [12]. The efficiency of the simulation allows also the estimation of rare events (blocking probability, rejection rate) is done in [11].

The aim of this paper is to study the simulation time needed to get a stationary sample for finite capacity networks. We show that for *monotone* systems CFTP scales very well with the state space explosion accompanying the increase in the number of queues.

More precisely, we study the coupling time τ of a CFTP algorithm (*i.e.* the number of steps needed to provide one sample). Our main interest is setting bounds on the expected coupling time. We first obtain exact analytical values for the expected simulation time for one M/M/1/C queue which serves as a building block for the following).

As for networks of queues, we show how upper bounds on the mean simulation time can be obtained as sums of the coupling times for all queues. One of the main result of this paper is to show that for acyclic networks with rejection in case of overflow,

$$\mathbb{E}\tau \leq \sum_i \frac{\Lambda}{\Lambda_i} \frac{C_i + C_i^2}{2},$$

where Λ is the global event rate in the network, Λ_i is the rate of events affecting Queue i and C_i is the capacity of Queue i . This result can be refined under light or heavy traffic assumptions to almost linear bounds in the capacities. All these bounds scale very well with the number of queues. This explains why perfect simulation of monotone queueing networks is so fast, especially when dealing with large scale

networks as in [11] where systems with up to 32 queues of capacity 30 (the state space is of size $31^{32} \approx 10^{47}$) are sampled over a classical desktop computer is less than 20 milli-seconds. This is good enough to estimate rare event probabilities.

The paper is organized as follows. We first introduce the coupling from the past algorithm in Section 2. Then we show general properties of the coupling time for open Markovian queueing networks in Section 3. We will investigate the M/M/1/c queue in Section 4 providing exact computation for the expected coupling time and the case of acyclic networks in Section 5 where bounds are derived, together with several experimental tests assessing their quality.

2. Coupling from the Past. Let $\{X_n\}_{n \in \mathbb{N}}$ be an irreducible and aperiodic discrete time Markov chain with a finite state space \mathcal{S} and a transition matrix $P = (p_{i,j})$. Let

$$\phi : \mathcal{S} \times \mathcal{E} \rightarrow \mathcal{S},$$

encode the chain, which means that it verifies the property $\mathbb{P}(\phi(i, e) = j) = p_{i,j}$ for every pair of states $(i, j) \in \mathcal{S}$ and for any e , a random variable distributed on \mathcal{E} . The function ϕ could be considered as a construction algorithm and e is the *innovation* for the chain. In the context of discrete event systems, e is an *event* and ϕ is the *transition function*. Now, the evolution of the Markov chain is described as a stochastic recursive sequence

$$X_{n+1} = \phi(X_n, e_{n+1}), \quad (2.1)$$

with X_n the state of the chain at time n and $\{e_n\}_{n \in \mathbb{N}}$ an independent and identically distributed sequence of random variables.

Let $\phi^{(n)} : \mathcal{S} \times \mathcal{E}^n \rightarrow \mathcal{S}$ denote the function whose output is the state of the chain after n iterations and starting in state $s \in \mathcal{S}$. That is,

$$\phi^{(n)}(s, e_{1 \rightarrow n}) = \phi(\dots \phi(\phi(s, e_1), e_2), \dots, e_n). \quad (2.2)$$

This notation can be extended to set of states. So for a set of states $A \subset \mathcal{S}$ we note

$$\phi^{(n)}(A, e_{1 \rightarrow n}) = \left\{ \phi^{(n)}(s, e_{1 \rightarrow n}), s \in A \right\}.$$

In the following, $|X|$ denotes the size of set X .

THEOREM 2.1 ([8]). *Let ϕ be a transition function on $\mathcal{S} \times \mathcal{E}$. There exists an integer l^* such that*

$$\lim_{n \rightarrow +\infty} \left| \phi^{(n)}(\mathcal{S}, e_{1 \rightarrow n}) \right| = l^* \text{ almost surely.}$$

The system *couple*s if $l^* = 1$. Then the *forward coupling time* τ^f defined by

$$\tau^f = \min\{n \in \mathbb{N}; \text{ such that } \left| \phi^{(n)}(\mathcal{S}, e_{1 \rightarrow n}) \right| = 1\},$$

is almost surely finite. The coupling property of a system ϕ depends only on the structure of ϕ . The probability measure on \mathcal{E} does not affect the coupling property, provided that all events in \mathcal{E} have a positive probability. Moreover, the existence of some pattern $e_{1 \rightarrow n_0}^*$ that ensures coupling, guarantees that τ^f is stochastically upper bounded by a geometric distribution

$$\mathbb{P}(\tau^f \geq k.n_0^*) \leq (1 - p(e_1^*).p(e_2^*) \dots p(e_{n_0}^*))^k; \quad (2.3)$$

where $p(e) > 0$ is the probability of event e .

At time τ^f , all trajectories issued from all initial states at time 0 have collapsed in only one trajectory. Unfortunately, the distribution of X_{τ^f} is not stationary. In [6] an example is given that illustrates why it is not possible to consider that this process has the stationary regime.

In fact, this iteration scheme could be reversed in time as it is usually done in the analysis of stochastic point processes. For that, one needs to extend the sequence of events to negative indexes and build the reversed scheme on sets by

$$A_n = \phi^{(n)}(\mathcal{S}, e_{-n+1 \rightarrow 0}).$$

It is clear that the sequence of sets A_n is non-decreasing ($A_{n+1} \subset A_n$). Consequently, the system couples if the sequence A_n converges almost surely to a set with only one element. Almost surely, there exists a finite time τ^b , the *backward coupling time*, defined by

$$\tau^b = \min\{n \in \mathbb{N}; \text{ such that } |\phi^{(n)}(\mathcal{S}, e_{-n+1 \rightarrow 0})| = 1\}.$$

PROPOSITION 2.2 ([13]). *The backward coupling time τ^b and the forward coupling time τ^f have the same probability distribution.*

The main result of the backward scheme is the following theorem.

THEOREM 2.3 ([8]). *Provided that the system couples, the state when coupling occurs for the backward scheme, is steady state distributed.*

From this fact, a general algorithm (1) sampling the steady state can be constructed.

Algorithm 1 Backward-coupling simulation (general version)

```

for all  $s \in \mathcal{S}$  do
   $y(s) \leftarrow s$  {choice of the initial value of the vector  $y$ ,  $n = 0$ }
end for
repeat
   $e \leftarrow \text{Random\_event}$ ; {generation of  $e_{-n+1}$ }
  for all  $s \in \mathcal{S}$  do
     $y(s) \leftarrow y(\phi(s, e))$ ;
    { $y(s)$  state at time 0 of the trajectory issued from  $s$  at time  $-n + 1$ }
  end for
until All  $y(x)$  are equal
return  $y(x)$ 

```

The mean complexity c_ϕ of this algorithm is $c_\phi = O(\mathbb{E}\tau^b \cdot |\mathcal{S}|)$. The coupling time τ^b is of fundamental importance for the efficiency of the sampling algorithm. To improve its complexity, we could reduce the factor $|\mathcal{S}|$ and reduce the coupling time. When the state space is partially ordered by a partial order \prec and the transition function is monotone for each event e , it is sufficient to simulate trajectories starting from the maximal and minimal states [8]. Denote by MAX and MIN the set of maximal, respectively minimal elements of \mathcal{S} for the partial order \prec . The monotone version of algorithm (1) is given by algorithm (2). In this case, we need to store the sequence of events in order to preserve the coherence between the trajectories driven from $MAX \cup MIN$.

Algorithm 2 Backward-coupling simulation (monotone version)

```

n=1;
R[n]=Random_event;{array R stores the backward sequence of events }
repeat
  n=2.n;
  for all  $s \in MAX \cup MIIN$  do
     $y(s) \leftarrow s$  {Initialize all trajectories at time  $-n$ }
  end for
  for  $i=n$  downto  $n/2+1$  do
    R[i]=Random_event; {generates all events from time  $-n+1$  to  $\frac{n}{2}+1$ }
  end for
  for  $i=n$  downto 1 do
    for all  $s \in MAX \cup MIN$  do
       $y(s) \leftarrow \phi(y(s), R[i])$ 
      { $y(s)$  is the state at time  $-i$  of the trajectory starting in  $s$  at time  $-n$ }
    end for
  end for
until All  $y(s)$  are equal
return  $y(s)$ 
    
```

The doubling scheme (first step in the loop) leads to a mean complexity

$$c_\phi = O(\mathbb{E}\tau^b.(|MAX| + |MIN|)). \quad (2.4)$$

3. Open Markovian queueing networks. Consider an open network Q consisting of K queues Q_1, \dots, Q_K . Each queue Q_i has a finite capacity, denoted by C_i , $i = 1, \dots, K$. Thus the state space of a single queue Q_i is $\mathcal{S}_i = \{0, \dots, C_i\}$. Hence, the state space \mathcal{S} of the network is $\mathcal{S} = \mathcal{S}_1 \times \dots \times \mathcal{S}_K$. The state of the system is described by a vector $s = (s_1, \dots, s_K)$ with s_i the number of customers in queue Q_i . The state space is partially ordered by the component-wise ordering and there are a maximal state MAX when all queues are full and a minimal state when all queues are empty.

The network evolves in time due to exogenous customer arrivals from outside of the network and to service completions of customers. After finishing his service at a server, a customer is either directed to another queue by a certain routing policy or leaves the network. A routing policy determines to which queue a customer will go, taking into account the global state of the system. Moreover, the routing policy also decides what happens with a customer if he is directed to a queue with a buffer filled with C_i customers.

An event in this network is characterized by the movements of some clients between queues, modeling the routing strategy and the Poisson process defines the occurrence rate of the events. For example, consider the acyclic queueing network of figure 3.1, made of 4 queues and 6 events.

Since the network is open, clients are able to enter and leave the network. We assume that customers who enter from outside the network to a given queue arrive according to a Poisson process. Furthermore, suppose that the service times at server i are independent and exponentially distributed with parameter μ_i .

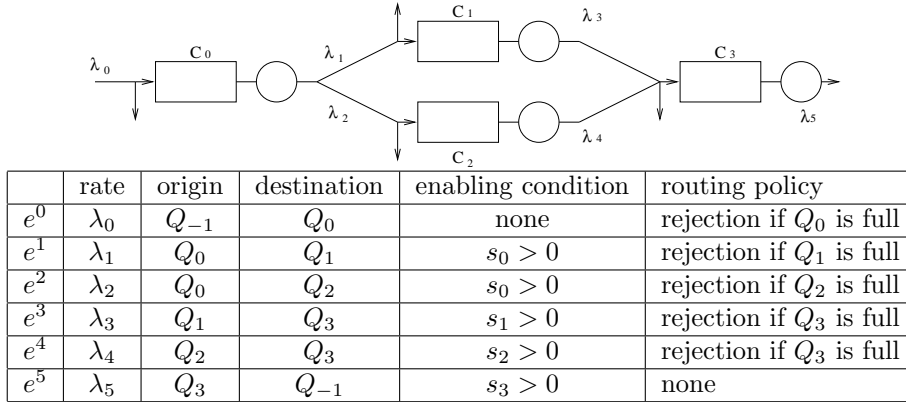


FIG. 3.1. Network with rejection

For example, for event e^1 (fig 3.1) we get

$$\phi(\cdot, e^1) : (s_0, s_1, s_2, s_3) \mapsto \begin{cases} (s_0 - 1, s_1 + 1, s_2, s_3) & \text{if } s_0 \geq 1 \text{ and } s_1 < C_1; \\ (s_0 - 1, s_1, s_2, s_3) & \text{if } s_0 \geq 1 \text{ and } s_1 = C_1(Q_1 \text{ full}); \\ (s_0, s_1, s_2, s_3) & \text{if } s_0 = 0(Q_0 \text{ empty}). \end{cases}$$

DEFINITION 3.1. An event e is monotone if $\phi(x, e) \leq \phi(y, e)$ for every x, y in \mathcal{S} with $x \leq y$.

It should be clear that event e^1 is monotone. Moreover usual events such as routing with overflow and rejection, routing with blocking and restart, routing with a index policy rule (eg Join the shortest queue) are monotone events [5, 12].

Denote by $\mathcal{E} = \{e_1, \dots, e_M\}$ the finite collection of events of the network. With each event e_i is associated a Poisson process with parameter λ_i . If an event occurs which does not satisfy the enabling condition the state of the system is unchanged.

To complete the construction of the discrete-time Markov chain, the system is uniformized by a Poisson process with rate $\Lambda = \sum_{i=1}^M \lambda_i$. Hence, one can see this Poisson process as a clock which determines when an event transition takes place. To choose which specific transition actually takes place, the collection \mathcal{E} of events of the network is randomly sampled with

$$p_i = \mathbb{P}(\text{event } e_i \text{ occurs}) = \frac{\lambda_i}{\Lambda}.$$

By construction, the following proposition should be clear.

PROPOSITION 3.2. The uniformized Markov chain has the same stationary distribution as the queueing network, and so does the embedded discrete time Markov chain.

Provided that events are monotone, the CFTP algorithm can be applied on queueing networks to build steady-state sampling of the network.

In our example of Figure 3.1 we ran the CFTP algorithm and produced samples of coupling time. The parameters used for the simulation are the following. Queues capacity : for all $i = 1, \dots, 4$, $C_i = 10$. Event rates: $\lambda_1 = 1.4$, $\lambda_2 = 0.6$, $\lambda_3 = 0.8$, $\lambda_4 = 0.5$ and $\lambda_5 = 0.4$. The global input rate λ_0 is varying. The number of samples used to estimate the mean coupling time is 10000. The result is displayed in Figure 3.2.

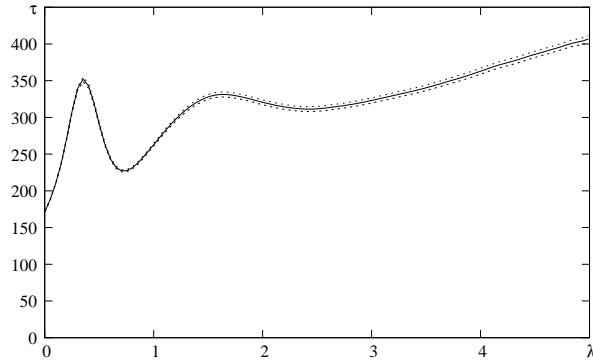


FIG. 3.2. The mean coupling time for the acyclic network of Figure 3.1 varies from 160 to 400 events when the input rate ranges from 0 to 4, with 95% confidence intervals.

This type of curve is of fundamental importance because the coupling time corresponds to the simulation duration and is involved in the simulation strategy (long run versus replication). These first results can be surprising because they exhibit a strong dependence on parameters values. The aim of this paper is now to understand more deeply what are the critical values for the network and to build bounds on the coupling time that are non-trivial.

Let N_i be the function from \mathcal{S} to \mathcal{S}_i with $N_i(s_1, \dots, s_K) = s_i$. So N_i is the number of customers in queue Q_i . As in section 2, τ^b refers to the *backward coupling time* of the chain, which is in case the coupling time from the past of the queueing network.

DEFINITION 3.3. Let τ_i^b denote the backward coupling time on coordinate i of the state space. Thus τ_i^b is the smallest n for which

$$\left| \left\{ N_i \left(\phi^{(n)}(s, e_{-n+1}, \dots, e_0) \right), s \in \mathcal{S} \right\} \right| = 1.$$

Because coordinate s_i refers to queue Q_i , the random variable τ_i^b represents the coupling time from the past of queue Q_i . Once all queues in the network have coupled, the CFTP algorithm returns one value and hence the chain has coupled. Thus

$$\tau^b = \max_{1 \leq i \leq K} \{ \tau_i^b \} \leq_{st} \sum_{i=1}^K \tau_i^b. \quad (3.1)$$

By taking expectation and interchanging sum and expectation we get:

$$\mathbb{E}[\tau^b] = \mathbb{E} \left[\max_{1 \leq i \leq K} \{ \tau_i^b \} \right] \leq \mathbb{E} \left[\sum_{i=1}^K \tau_i^b \right] = \sum_{i=1}^K \mathbb{E}[\tau_i^b] \quad (3.2)$$

It follows from Proposition 2.2 that τ^b and τ^f have the same distribution. The same holds for τ_i^f and τ_i^b . Hence $\mathbb{E}[\tau_i^b] = \mathbb{E}[\tau_i^f]$ and

$$\mathbb{E}\tau^b \leq \sum_{i=1}^K \mathbb{E}[\tau_i^f]. \quad (3.3)$$

The bound given in Equation 3.3 is interesting because $\mathbb{E}[\tau_i^f]$ is sometimes amenable to explicit computations, as shown in following sections. In order to de-

rive those bounds, one may provide yet other bounds, by making the coupling state explicit.

DEFINITION 3.4. *The hitting time $h_{j \rightarrow k}$ in a Markov chain X_n is defined as*

$$h_{j \rightarrow k} = \inf_{\mathbb{N}} \{n \text{ s.t. } X_n = k | X_0 = j\} \text{ with } j, k \in \mathcal{S}.$$

In the queueing framework, $h_{0 \rightarrow C_i}$ represents the number of steps it takes for queue Q_i to go from state 0 to state C_i . Now we consider queue Q_i out of the network and examine it independently. Suppose that $h_{0 \rightarrow C_i} = n$ for the sequence of events e_1, \dots, e_n . Because of monotonicity of ϕ we have

$$\phi^{(n)}(0, e_1, \dots, e_n) \leq \phi^{(n)}(s, e_1, \dots, e_n) \leq \phi^{(n)}(C_i, e_1, \dots, e_n) = 0,$$

with $s \in \mathcal{S}_i$. Hence, coupling has occurred. So $h_{0 \rightarrow C_i}$ is an upper bound on the forward coupling of queue Q_i . The same argumentation holds for $h_{C_i \rightarrow 0}$. Thus

$$\mathbb{E} [\tau_i^f] \leq \mathbb{E} [\min\{h_{0 \rightarrow C_i}, h_{C_i \rightarrow 0}\}]. \quad (3.4)$$

Hence,

$$\mathbb{E} \tau^b \leq \sum_{i=1}^K \mathbb{E} [\tau_i^f] \leq \sum_{i=1}^K \mathbb{E} [\min\{h_{0 \rightarrow C_i}, h_{C_i \rightarrow 0}\}] \leq \sum_{i=1}^K \min(\mathbb{E} h_{0 \rightarrow C_i}, \mathbb{E} h_{C_i \rightarrow 0}), \quad (3.5)$$

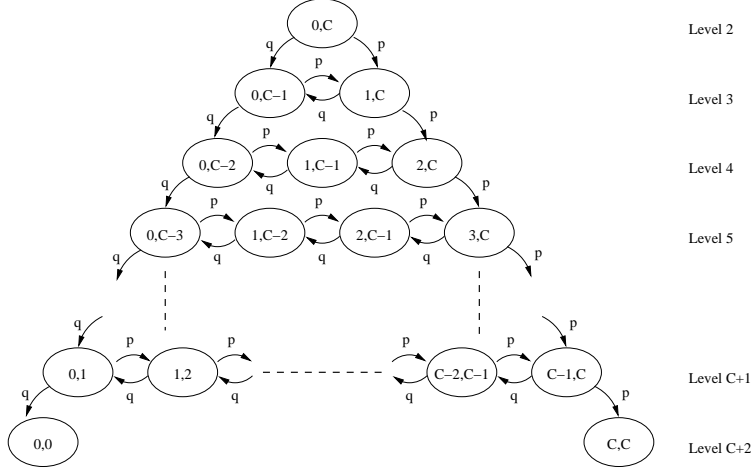
by Jensen's Inequality.

4. Coupling time in a M/M/1/C queue. The M/M/1/C queue is well known and there is no need to run simulations to get the distribution of its stationary distribution. However, the computation of hitting times provided here is new and will serve as a building block for the following section on networks.

In a M/M/1/C model, we have a single queue with one server. Customers arrive at the queue according to a Poisson process with rate λ and the service time is distributed according to an exponential distribution with parameter μ . In the queue there is only place for C customers. So the state space $\mathcal{S} = \{0, \dots, C\}$. If a customer arrives when there are already C customers in the queue, he immediately leaves without entering the queue. After uniformization, we get a discrete time Markov chain which is governed by the events e_a with probability $p = \frac{\lambda}{\lambda + \mu}$ and e_d with probability $q = 1 - p$. Event e_a represents an arrival and event e_d represents an end of service with departure of the customer.

In order to estimate the expectation of the coupling time from the past $\mathbb{E}[\tau^b]$ we use inequality 3.5. Since there is only one queue, the first two inequalities in 3.5 become equalities. Indeed, when applying forward simulation, the chain only can couple in state 0 or state C . This follows since for $r, s \in \mathcal{S}$ with $0 < r < s < C$ we have $\phi(r, e_a) = r + 1 < s + 1 = \phi(s, e_a)$ and $\phi(r, e_d) = r - 1 < s - 1 = \phi(s, e_d)$. So the chain cannot couple in a state s with $0 < s < C$. Furthermore we have $\phi(C, e_a) = C = \phi(C - 1, e_a)$ and $\phi(0, e_d) = 0 = \phi(1, e_d)$. Hence, forward coupling can only occur in 0 or C :

$$\mathbb{E} [\tau^b] = \mathbb{E} [\min\{h_{0 \rightarrow C}, h_{C \rightarrow 0}\}]. \quad (4.1)$$


 FIG. 4.1. Markov chain $X(q)$ corresponding to $H_{i,j}$

4.1. Explicit calculation of $\mathbb{E}[\tau^b]$. In order to compute $\min\{h_{0 \rightarrow C}, h_{C \rightarrow 0}\}$ we have to run two copies of the Markov chain for a M/M/1/C queue simultaneously (whose states are x and y respectively). One copy starts in state 0 and the other one starts in state C . We stop when either the chain starting in 0 reaches state C or when the copy starting in state C reaches state 0.

Therefore, let us rather consider a product Markov chain $X(q)$ with state space $\mathcal{S} \times \mathcal{S} = \{(x, y), x = 0, \dots, C, y = 0, \dots, C\}$. The Markov chain $X(q)$ is also governed by the two events e_a and e_d and the function ϕ is:

$$\begin{aligned} \psi((x, y), e_a) &= ((x + 1) \wedge C, (y + 1) \wedge C) \\ \psi((x, y), e_d) &= ((x - 1) \vee 0, (y - 1) \vee 0). \end{aligned}$$

Without any loss of generality, we may assume that $x \leq y$. This system corresponds with the *pyramid Markov chain* $X(q)$ displayed in Figure 4.1. The rest of this section is devoted to establishing a formula for the expected exit time of the pyramid. Although the technique used here (one step analysis) is rather classical, it is interesting to notice how this is related to random walks on the line (this also explains the shifted indexes associated to the levels of the pyramid).

Since we can only couple in 0 or C , this coupling occurs as soon as the chain $X(q)$ reaches states $(0, 0)$ or (C, C) . Define

$$H_{i,j} := \text{number of steps to reach state } (0, 0) \text{ or } (C, C) \text{ starting from state } (i, j)$$

with $(i, j) \in \mathcal{S} \times \mathcal{S}$. By definition, $\min\{h_{0 \rightarrow C}, h_{C \rightarrow 0}\} = H_{0,C}$. Now $H_{i,j}$ represents the hitting time of the coupling states $(0, 0)$ and (C, C) (also called absorption time) in a product Markov chain. Using a one step analysis, we get the following system of equations for $\mathbb{E}[H_{i,j}]$:

$$\begin{cases} \mathbb{E}[H_{i,j}] &= 1 + p\mathbb{E}[H_{(i+1) \wedge C, (j+1) \wedge C}] + q\mathbb{E}[H_{(i-1) \vee 0, (j-1) \vee 0}], & i \neq j, \\ \mathbb{E}[H_{i,j}] &= 0, & i = j \end{cases} \quad (4.2)$$

Two states (i, j) and (i', j') are said to be at the same level if $|j - i| = |j' - i'|$. In Figure 4.1 we can distinguish $C + 1$ levels. Because of monotonicity of ϕ , $|j - i|$

cannot increase. Hence, starting at a level with $|j - i|$, the chain will gradually pass all intermediate levels to reach finally the level with $|j - i| = 0$ in state $(0, 0)$ or (C, C) . Thus, starting in state $(0, C)$, the chain will run through all levels to end up at the level with $|j - i| = 0$. So, $H_{0,C} = \min\{h_{0 \rightarrow C}, h_{C \rightarrow 0}\}$. To determine $\mathbb{E}[H_{0,C}]$ we determine the mean time spent on each level and sum up over all levels.

A state (i, j) belongs to level m if $|j - i| = C + 2 - m$. Then state $(0, C)$ belongs to level 2 and the states $(0, 0)$ and (C, C) belong to level $C + 2$. To get from $(0, C)$ into either $(0, 0)$ or (C, C) , the chain $X(q)$ needs to cross all levels between the levels 2 and $C + 2$. Let T_m denote time it takes to reach level $m + 1$, starting in level m . Then

$$H_{0,C} = \sum_{m=2}^{C+1} T_m. \quad (4.3)$$

In order to determine T_m , we associate to each level m a random walk R_m on $0, \dots, m$ with absorbing barriers at state 0 and state C . In the random walk, the probability of going up is p and of going down is $q = 1 - p$. We have the following correspondence between the states of the random walk R_m and the states of $X(q)$ (see Figure 4.2).

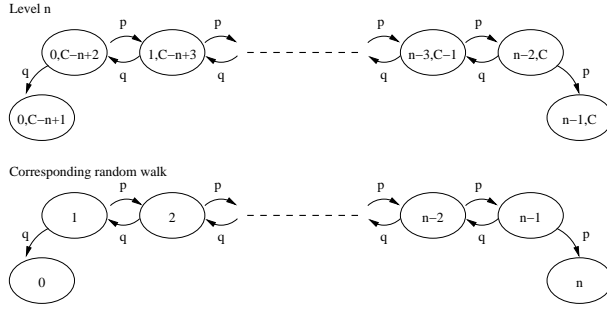


FIG. 4.2. Relationship between level m and random walk R_m .

- State 0 of R_m corresponds with state $(0, C - m + 1)$ of $X(q)$,
- State i of R_m corresponds with state $(i - 1, C - m + 1 + i)$ of the $X(q)$,
 $1 \leq i \leq m - 1$,
- State m of R_m corresponds with state $(m - 1, C)$ of $X(q)$.

Now the time spent on level m in $X(q)$ is the same as the time spent in a random walk R_m before absorption. Therefore, one can use the two following results on random walks in determining T_m , which are known as ruin problems (see for example [9]).

Let $\alpha_{i \rightarrow 0}^m$ denote the probability of absorption in state 0 of the random walk R_m starting in i . Then:

$$\alpha_{i \rightarrow 0}^m = \begin{cases} \frac{a^m - a^i}{a^m - 1}, & p \neq \frac{1}{2}, \\ \frac{m-i}{m}, & p = \frac{1}{2}, \end{cases} \quad (4.4)$$

where $a = q/p$.

Now, absorption occurs in R_m once the state 0 or C has been achieved.

LEMMA 4.1. Let \tilde{T}_i^m denote the mean absorption time of a random walk R_m

starting in i . Then:

$$\mathbb{E}[\tilde{T}_i^m] = \begin{cases} \frac{i-m(1-\alpha_{i \rightarrow 0}^m)}{q-p}, & p \neq \frac{1}{2}, \\ i(m-i), & p = \frac{1}{2}. \end{cases} \quad (4.5)$$

Now, let β_0^m (resp. β_m^m) denote the probability that absorption occurs in 0 (resp. m). Then

$$\beta_0^m = \sum_{i=0}^m \alpha_{i \rightarrow 0}^m \mathbb{P}(R_m \text{ starts in state } i), \quad (4.6)$$

and $\beta_m^m = 1 - \beta_0^m$. From the structure of the Markov chain $X(q)$ and the correspondence between $X(q)$ and the random walks, we have that (see Figure 4.2):

$$\mathbb{P}(\text{enter level } m+1 \text{ at } (0, C-m+1)) = \mathbb{P}(\text{absorption in } 0 \text{ in } R_m) = \beta_0^m.$$

Now one has:

$$\begin{aligned} \mathbb{E}[T_m] &= \mathbb{E}[\tilde{T}_1^m] \beta_0^{m-1} + \mathbb{E}[\tilde{T}_{m-1}^m] \beta_{m-1}^{m-1} \\ &= \mathbb{E}[\tilde{T}_{m-1}^m] + \left(\mathbb{E}[\tilde{T}_1^m] - \mathbb{E}[\tilde{T}_{m-1}^m] \right) \beta_0^{m-1}. \end{aligned} \quad (4.7)$$

4.1.1. Case $q = p = 1/2$. $\mathbb{E}[T_m]$ can be calculated explicitly for $p = \frac{1}{2}$. Since the random walk is symmetric, we have $\beta_0^m = \beta_n^m = \frac{1}{2}$. So:

$$\mathbb{E}[T_m] = \mathbb{E}[\tilde{T}_1^m] \beta_0^{m-1} + \mathbb{E}[\tilde{T}_{m-1}^m] \beta_{m-1}^{m-1} = m-1. \quad (4.8)$$

Hence,

$$\mathbb{E}[H_{0,C}] = \sum_{m=2}^{C+1} \mathbb{E}[T_m] = \sum_{m=2}^{C+1} m-1 = \frac{C^2+C}{2}.$$

LEMMA 4.2. For a $M/M/1/C$ with $\lambda = \mu$, $\mathbb{E}\tau^b = \frac{C^2+C}{2}$.

4.1.2. Case $p \neq \frac{1}{2}$. Since the random walks are not symmetric, one cannot apply the same reasoning as for the case $p = \frac{1}{2}$ to compute β_0^m . Entering the random walk R_m corresponds to entering level m in $X(q)$. Since we can only enter level m in the state $(0, C-m+2)$ or $(m-2, C)$ this means we can only start the random walk in state 1 or $m-1$. Therefore (4.6) becomes:

$$\begin{aligned} \beta_0^m &= \sum_{i=0}^m \alpha_{i \rightarrow 0}^m \mathbb{P}(R_m \text{ starts in state } i) \\ &= \alpha_{1 \rightarrow 0}^m \mathbb{P}(R_m \text{ starts in } 1) + \alpha_{m-1 \rightarrow 0}^m \mathbb{P}(R_m \text{ starts in } m-1) \\ &= \frac{a^m - a^{m-1}}{a^m - 1} + \frac{a^{m-1} - a}{a^m - 1} \beta_0^{m-1}. \end{aligned}$$

This gives the recurrence:

$$\begin{cases} \beta_0^m &= \frac{a^m - a^{m-1}}{a^m - 1} + \frac{a^{m-1} - a}{a^m - 1} \beta_0^{m-1} & m > 2; \\ \beta_0^2 &= 2. \end{cases} \quad (4.9)$$

Thus we obtain,

PROPOSITION 4.3. *For a M/M/1/C queue, using the foregoing notations,*

$$\mathbb{E}\tau^b = \mathbb{E}[H_{0,C}] = \sum_{m=2}^{C+1} \mathbb{E}[\tilde{T}_{m-1}^m] + \left(\mathbb{E}[\tilde{T}_1^m] - \mathbb{E}[\tilde{T}_{m-1}^m] \right) \beta_0^{m-1}, \quad (4.10)$$

with β_0^m defined by (4.9) and $\mathbb{E}[\tilde{T}_{m-1}^m]$ and $\mathbb{E}[\tilde{T}_1^m]$ defined by (4.5).

4.1.3. Comparison between the cases $p = 1/2$ and $p \neq 1/2$.

PROPOSITION 4.4. *The coupling time in a M/M/1/C queue is maximal when the input rate λ and the service rate μ are equal.*

Proof. By definition, $\lambda = \mu$ corresponds to $p = q = 1/2$. The proof holds by induction on C . The result is obviously true when $C = 0$, because whatever q , $\mathbb{E}[H_{0,C}] = 0$.

For $C + 1$, let q be an arbitrary probability with $q > 1/2$ (the case $q < 1/2$ is symmetric). We will compare the expected time for absorption of three Markov chains. The first one is the Markov chain $X := X(1/2)$ displayed in Figure 4.1, with $q = p = 1/2$. The second one is the Markov chain $X' = X(q)$ displayed in Figure 4.1 and the last one X'' is a mixture between the two previous chains: The first C levels are the same as in X while the last level ($C + 1$) is the same as in X' .

The expected absorption time for the first C levels is the same for X and for X'' : $\sum_{m=2}^C \mathbb{E}T_m = \sum_{m=2}^C \mathbb{E}T'_m$. By induction, this is larger than for X' : we have $\sum_{m=2}^C \mathbb{E}T_m = \sum_{m=2}^C \mathbb{E}T''_m \geq \sum_{m=2}^C \mathbb{E}T'_m$. Therefore, we just need to compare the expected exit times out of the last level, namely $\mathbb{E}T_{C+1}$, $\mathbb{E}T'_{C+1}$ and $\mathbb{E}T''_{C+1}$, to finish the proof.

Let us first compare $\mathbb{E}T_{C+1}$ and $\mathbb{E}T''_{C+1}$. In both cases, the Markov chain enters level $C + 1$ in state $(0, 1)$ with probability $1/2$.

Equation (4.8) says that $\mathbb{E}T_{C+1} = C$ and Equation (4.5) gives after straightforward computations, $\mathbb{E}T''_{C+1} = 1/2 \frac{1-C(1-\alpha_{1 \rightarrow 0}^C)}{q-p} + 1/2 \frac{C-1-C(1-\alpha_{C-1 \rightarrow 0}^C)}{q-p} = \frac{C}{2q} \frac{a^C - a}{a^{C-1}} \leq C/(2q) < C = \mathbb{E}T_{C+1}$.

In order to compare $\mathbb{E}T'_{C+1}$ and $\mathbb{E}T''_{C+1}$, let us first show that β_0^m is larger than $1/2$, for all $m \geq 2$. This is done by an immediate induction on Equation (4.9). If $\beta_0^{m-1} \geq 1/2$, then $\beta_0^m \geq \frac{2a^m - a^{m-1} - a}{2(a^m - 1)}$. Now, $\frac{2a^m - a^{m-1} - a}{2(a^m - 1)} \geq 1/2$ if $2a^m - a^{m-1} - a \geq a^m - 1$, i.e. after recombining the terms, $(a-1)(a^{m-1} - 1) \geq 0$. This is true as soon as $q \geq 1/2$.

To end the proof, it is enough to notice that for the chain X' , time to absorption starting in 1, $\mathbb{E}\tilde{T}_1^{m'}$ is smaller than time to absorption starting in $m-1$, $\mathbb{E}\tilde{T}_{m-1}^{m'}$ for all m . The difference $\mathbb{E}\tilde{T}_{m-1}^{m'} - \mathbb{E}\tilde{T}_1^{m'}$ is

$$\frac{ma^m - ma^{m-1} + ma - m - 2a^m + 2}{p(a^m - 1)(a - 1)} = \frac{m(a-1) \left(\frac{a^{m-1} + 1}{2} - \frac{1+a+\dots+a^{m-1}}{m} \right)}{p(a^m - 1)(a - 1)} \geq 0,$$

by convexity of $x \mapsto a^x$. Finally,

$$\mathbb{E}T'_{C+1} = \beta_0^{C+1} \mathbb{E}\tilde{T}_1^{C+1'} + (1 - \beta_0^{C+1}) \mathbb{E}\tilde{T}_C^{C+1'} \leq \frac{1}{2} \mathbb{E}\tilde{T}_1^{C+1'} + \frac{1}{2} \mathbb{E}\tilde{T}_C^{C+1'} = \mathbb{E}T''_{C+1}.$$

□

4.2. Explicit Bounds. Equation (4.10) provides a quick way to compute $\mathbb{E}[H_{0,C}]$ using recurrence equation (4.9). However, it may also be interesting to get a simple closed form for an upper bound for $\mathbb{E}[H_{0,C}]$. This can be done using the last inequality in Equation (3.5) that gives an upper bound for $\mathbb{E}[H_{0,C}]$ amenable to direct computations.

$$\mathbb{E}[H_{0,C}] = \mathbb{E}[\min\{h_{0 \rightarrow C}, h_{C \rightarrow 0}\}] \leq \min\{\mathbb{E}[h_{0 \rightarrow C}], \mathbb{E}[h_{C \rightarrow 0}]\}. \quad (4.11)$$

The exact calculation of $\mathbb{E}[h_{0 \rightarrow C}]$ can be done using a one step analysis. Let F_i be the time to go from state i to 0. Then, $h_{0 \rightarrow C} = F_C$ and for all $i > 0$,

$$\mathbb{E}[F_i] = 1 + p\mathbb{E}[F_{(i+1) \wedge C}] + q\mathbb{E}[F_{i-1}]. \quad (4.12)$$

With an approach derived from [9] one can condition on the next event. Let T_i denote the time to go from state i to $i+1$. Then

$$\mathbb{E}[h_{0 \rightarrow C}] = \sum_{i=0}^{C-1} \mathbb{E}[T_i]. \quad (4.13)$$

To get an expression for T_i , with $0 < i \leq C$, we condition on the first event. Therefore let $\mathbb{E}[T_i|e]$ denote the conditional expectation of T_i knowing that the next event is e . Since $\mathbb{E}[T_i | e_a] = 1$ and $\mathbb{E}[T_i | e_d] = 1 + \mathbb{E}[T_{i-1}] + \mathbb{E}[T_i]$, conditioning delivers the following recurrent expression for the $\mathbb{E}[T_i]$:

$$\mathbb{E}[T_i] = \begin{cases} \frac{1}{p} + \frac{q}{p}\mathbb{E}[T_{i-1}], & 0 < i < C \\ \frac{1}{p}, & i = 0. \end{cases} \quad (4.14)$$

By induction one can show that $\mathbb{E}[T_i] = \frac{1}{p} \sum_{k=0}^i \left(\frac{q}{p}\right)^k$. Hence, $\mathbb{E}[T_i] = \frac{1 - \left(\frac{q}{p}\right)^{i+1}}{p - q}$ and from (4.13) it follows that

$$\mathbb{E}[h_{0 \rightarrow C}] = \sum_{i=0}^{C-1} \frac{1 - \left(\frac{q}{p}\right)^{i+1}}{p - q} = \frac{C}{p - q} - \frac{q \left(1 - \left(\frac{q}{p}\right)^C\right)}{(p - q)^2}. \quad (4.15)$$

By reasons of symmetry, we have

$$\mathbb{E}[h_{C \rightarrow 0}] = \frac{C}{q - p} - \frac{p \left(1 - \left(\frac{p}{q}\right)^C\right)}{(q - p)^2}. \quad (4.16)$$

The curves of $\mathbb{E}[h_{0 \rightarrow C}]$ and $\mathbb{E}[h_{C \rightarrow 0}]$ intersect in $C^2 + C$ when $p = q$. If $p > q$ then $\mathbb{E}[h_{0 \rightarrow C}] < \mathbb{E}[h_{C \rightarrow 0}]$ and because of symmetry, if $p < q$ then $\mathbb{E}[h_{0 \rightarrow C}] > \mathbb{E}[h_{C \rightarrow 0}]$. Since also $\frac{C^2 + C}{2}$ is an upper bound corresponding to the critical case $p = q$ on the mean coupling time $\mathbb{E}\tau^b$, as shown in Proposition 4.4, one can state:

PROPOSITION 4.5. *The mean coupling time $\mathbb{E}\tau^b$ of a $M/M/1/C$ queue with arrival rate λ and service rate μ is bounded using $p = \lambda/(\lambda + \mu)$ and $q = 1 - p$.*

$$\begin{aligned} \text{Critical bound:} & \quad \forall p \in [0, 1], \quad \mathbb{E}\tau^b \leq \frac{C^2 + C}{2}. \\ \text{Heavy traffic Bound:} & \quad \text{if } p > \frac{1}{2}, \quad \mathbb{E}\tau^b \leq \frac{C}{p - q} - \frac{q \left(1 - \left(\frac{q}{p}\right)^C\right)}{(p - q)^2}. \\ \text{Light traffic bound:} & \quad \text{if } p < \frac{1}{2}, \quad \mathbb{E}\tau^b \leq \frac{C}{q - p} - \frac{p \left(1 - \left(\frac{p}{q}\right)^C\right)}{(q - p)^2}. \end{aligned}$$

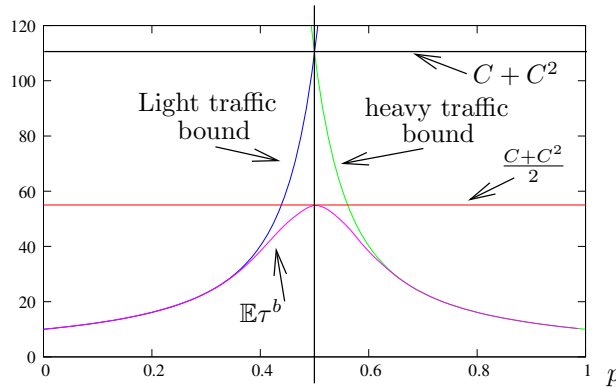


FIG. 4.3. Expected coupling time in an $M/M/1/10$ queue when q varies from 0 to 1 and the three explicit bounds given in Proposition 4.5

Figure 4.3 displays both the exact expected coupling time for a queue with capacity 10 as given by Equation (4.10) as well as the three explicit bounds given in Proposition 4.5. Note that the bounds are very accurate under light or heavy traffic ($q \leq 0.4$ and $q > 0.6$). In any case, the ratio is never larger than 1.2.

5. Coupling in acyclic queueing networks. This section is dedicated to the effective computation of a bound of the coupling time in acyclic networks. In the acyclic network given in Figure 3.1, the coupling time has a peak at $\lambda_0 = 0.4$, as one can see in Figure 3.2. This corresponds to the case when the input rate and service rate in Queue 3 are equal. This should not be surprising regarding the result for a single queue, which says that the coupling time is maximal when the rates are equal. Then a second peak occurs around $\lambda_0 = 1.4$ when coupling in Queue 0 is maximal. The rest of the curve shows a linear increase of the coupling time which may suggest an asymptotic linear dependence in λ_0 . In this part, an explicit bound on the coupling time which exhibits these two features will be derived.

The first result concerns an extension of inequality (3.5) to distributions. The second part shows how the results for a single $M/M/1/C$ queue can be used to get an effective computation of bounds for acyclic networks on queues.

In the following, the queues Q_0, \dots, Q_K are numbered according to the topological order of the network. Thus, no event occurring in queue Q_i has any influence on the state of queue Q_j as long as $i > j$.

5.1. Computation of an upper bound on the coupling time. Here, an acyclic network of $M/M/1/C$ queues with an arbitrary topology and Bernoulli routings is considered. The events here are of only two types: exogenous arrivals (Poisson with rate γ_i in queue i) and routing of one customer from queue i to queue j after service completion in queue i (with rate μ_{ij}). Queue $K+1$ is a dummy queue representing exits: routing a customer to queue $K+1$ means that the customer exits the network forever. In case of overflow, the new customer trying to enter the full queue is lost. The service rate at queue i is also denoted $\mu_i = \sum_{j=0}^{K+1} \mu_{ij}$.

Let us introduce new random variables. $\tau^b(s_j = x)$ is the backward coupling time of the network, over the set of all initial states with the j -th coordinate equal to x .

Namely,

$$\tau^b(s_j = x) = \min \{n \text{ s.t. } |\phi^{(n)}(\mathcal{S} \cap \{s_j = x\}, e_{-n+1}, \dots, e_0)| = 1\}.$$

$\tau_i^b(s_j = x)$ is the backward coupling time on coordinate i given $s_j = x$:

$$\tau_i^b(s_j = x) = \min \{n \text{ s.t. } |N_i(\phi^{(n)}(\mathcal{S} \cap \{s_j = x\}, e_{-n+1}, \dots, e_0))| = 1\}.$$

It should be obvious that $\tau^b(s_j = x) \leq_{st} \tau^b$ and for all i , $\tau_i^b(s_j = x) \leq_{st} \tau_i^b$. We also have the same notions for forward coupling times:

$$\tau^f(s_j = x) = \min \{n \in \mathbb{N}; \text{ s.t. } |\phi^{(n)}(\mathcal{S} \cap \{s_j = x\}, e_{1 \rightarrow n})| = 1\},$$

$\tau_i^f(s_j = x)$ being defined in the same manner, and for hitting times:

$$h_{C_i \rightarrow 0}(s_j = x) = \min \{n \in \mathbb{N}; \text{ s. t. } \phi^{(n)}(\mathcal{S} \cap \{s_i = C_i, s_j = x\}, e_{1 \rightarrow n}) \in \mathcal{S} \cap \{s_i = 0\}\}.$$

Now, sweeping the list of queues in the topological order, one can construct a sequence of backward simulations in the following way.

First simulate the queueing system from the past up to coupling of queue 0. The number of steps is by definition τ_0^b . Queue Q_0 has coupled in a random state X_0 . Then, run a second backward simulation up to coupling of Queue Q_1 given $s_0 = X_0^0$. This simulation takes $\tau_1^b(s_0 = X_0^0)$ steps and the state at time $t = 0$ is X_1^1 for Q_1 and X_0^1 for Q_0 .

This construction goes on up to the backward simulation up to coupling of Queue Q_K given $s_0 = X_0^{K-1}, s_1 = X_1^{K-1}, \dots, s_{K-1} = X_{K-1}^{K-1}$. The last simulation takes $\tau_i^b(s_0 = X_0^{K-1}, s_1 = X_1^{K-1}, \dots, s_{K-1} = X_{K-1}^{K-1})$ steps and the coupling state of Q_K is X_K^K .

LEMMA 5.1. *One has $\tau^b \leq_{st} \sum_{i=0}^K \tau_i^b(s_0 = X_0^{i-1}, \dots, s_{i-1} = X_{i-1}^{i-1})$, and for all i , (X_0^i, \dots, X_i^i) is steady state distributed for Q_0, \dots, Q_i . Furthermore, for all i ,*

$$\tau^b \leq_{st} \sum_{i=0}^K h_{C_i \rightarrow 0}(s_0 = X_0^{i-1}, \dots, s_{i-1} = X_{i-1}^{i-1}). \quad (5.1)$$

Proof. From the previous sequence of backward simulations one can construct a single one by appending them in the reverse order: the backward simulation for Queue Q_K preceded by the simulation of Q_{K-1} , and so forth up to the simulation of Q_0 . This is a backward simulation of the system (the last state is (X_0^K, \dots, X_i^K)). This construction is illustrated in the case of two queues in tandem in Figure 5.1.

A straightforward consequence, using acyclicity, is that (X_0^i, \dots, X_i^i) is steady state distributed for Q_0, \dots, Q_i for all i .

Furthermore, one gets in distribution

$$\begin{aligned} \tau^b &\leq_{st} \sum_{i=0}^K \tau_i^b(s_0 = X_0^{i-1}, \dots, s_{i-1} = X_{i-1}^{i-1}) = \sum_{i=0}^K \tau_i^f(s_0 = X_0^{i-1}, \dots, s_{i-1} = X_{i-1}^{i-1}) \\ &\leq_{st} \sum_{i=0}^K h_{C_i \rightarrow 0}(s_0 = X_0^{i-1}, \dots, s_{i-1} = X_{i-1}^{i-1}), \end{aligned}$$

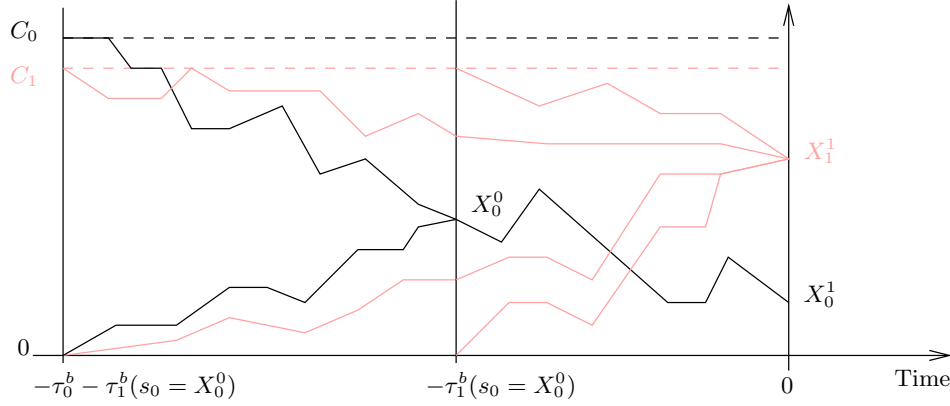


FIG. 5.1. The trajectories of the state in Q_0 are in black while the trajectories for Q_1 are in the lighter color. Starting at time $-\tau_0^b - \tau_1^b (s_0 = X_0^0)$, the state of Q_0 has coupled in X_0^0 at time $-\tau_1^b (s_0 = X_0^0)$. From then on, Q_0 stays coupled and Q_1 couples at some time before 0.

by independence of the variables, given the initial states X^{i-1} .

□

Let us now consider a new circuit with one difference from the original one: all queues are replaced by infinite queues, except for queue Q_i which stays the same. In the following, all the notations related to this new network will be expressed by appending the ∞ symbol to all variables corresponding to this new circuit.

The new circuit up to Queue i is product form and using Burke's Theorem, the input stream in Queue i is Poisson. The rate of the input stream in queue i is given by ℓ_i , the solution of the flow equations:

$$\ell_i = \sum_{j < i} \ell_j \frac{u_{ji}}{\mu_j} + \gamma_i.$$

The network is said to be *stable* for Queue i as soon as $\ell_i < \mu_i$. We assume stability for all i in the following.

One can construct a sequence of backward simulations for the new network in the same way as for the original network. This provides the quantities ${}^\infty X_j^{i-1}$, ${}^\infty \tau_i^b (s_0 = {}^\infty X_0^{i-1}, \dots, s_{i-1} = {}^\infty X_{i-1}^{i-1})$, ${}^\infty \tau_i^f (s_0 = {}^\infty X_0^{i-1}, \dots, s_{i-1} = {}^\infty X_{i-1}^{i-1})$, and ${}^\infty h_{C_i \rightarrow 0} (s_0 = {}^\infty X_0^{i-1}, \dots, s_{i-1} = {}^\infty X_{i-1}^{i-1})$.

The monotony property given above implies that $X_i^j \leq_{st} {}^\infty X_i^j$ and

$$h_{C_i \rightarrow 0} (s_0 = X_0^{i-1}, \dots, s_{i-1} = X_{i-1}^{i-1}) \leq_{st} {}^\infty h_{C_i \rightarrow 0} (s_0 = {}^\infty X_0^{i-1}, \dots, s_{i-1} = {}^\infty X_{i-1}^{i-1}).$$

The next step is to build yet another model. This third model is made of a single $M/M/1/C_i$ queue with three types of events, arrivals of customers with rate ℓ_i (provided that the number of customers is smaller than C_i), departures with rate μ_i (provided that the number of customers is positive) and null events (with no effect on the queue) with rate $\Lambda - \ell_i - \mu_i$.

For this isolated model, let us introduce the uniformizing probabilities $p = \ell_i / \Lambda$, $q = 1 - p$ and $d = (\Lambda - \ell_i - \mu_i) / \Lambda$. Let F_k be the time to go from state k to state 0

in the isolated system. A one step analysis gives

$$\begin{aligned}\mathbb{E}[F_k] &= 1 + d\mathbb{E}[F_k] + \frac{\ell_i}{\Lambda}\mathbb{E}[F_{(k+1)\wedge C_i}] + \frac{\mu_i}{\Lambda}\mathbb{E}[F_{(k-1)}] \\ &= \frac{1}{1-d} + p\mathbb{E}[F_{(k+1)\wedge C_i}] + q\mathbb{E}[F_{(k-1)\vee 0}].\end{aligned}$$

We get the same equation as (4.14) except for the additional constant which is now $\frac{1}{1-d}$ instead of 1, so that the solution is the same as before up to a multiplicative factor of $\frac{1}{1-d} = \frac{\Lambda}{\ell_i + \mu_i}$. Using Equation (4.16), one gets

$$\mathbb{E}[F_{C_i}] = \frac{\Lambda}{\ell_i + \mu_i} \left(\frac{C_i}{q-p} - \frac{p(1 - (\frac{p}{q})^{C_i})}{(q-p)^2} \right). \quad (5.2)$$

LEMMA 5.2. *Under the foregoing notations and assumptions,*

$${}^\infty h_{C_i \rightarrow 0}(s_0 = {}^\infty X_0^{i-1}, \dots, s_{i-1} = {}^\infty X_{i-1}^{i-1}) = F_{C_i},$$

in distribution.

Proof. First, using Lemma 5.1 for the new network with infinite queues (except for Q_i), the state $({}^\infty X_0^{i-1}, \dots, {}^\infty X_{i-1}^{i-1})$ is steady state distributed. Using Burke's Theorem, this implies that the input stream in queue Q_i is Poisson with rate ℓ_i , when one runs a simulation starting in any state in $\mathcal{S} \cap \{s_i = C_i, s_j = {}^\infty X_j^{i-1}, j < i\}$.

Now, during this simulation, one can couple the addition, subtraction et null events for queue Q_i in isolation and for Q_i in the complete network of infinite queues, all of them having the same laws. This implies that the state of queue Q_i in both systems is the same under that coupling. Hence, they reach 0 at the same time: ${}^\infty h_{C_i \rightarrow 0}(s_0 = {}^\infty X_0^{i-1}, \dots, s_{i-1} = {}^\infty X_{i-1}^{i-1}) = F_{C_i}$ in distribution. \square

We are ready to put everything together in expectation.

$$\mathbb{E}\tau^b \leq_{st} \sum_i \mathbb{E}[h_{C_i \rightarrow 0}(s_0 = X_0^{i-1}, \dots, s_{i-1} = X_{i-1}^{i-1})] \quad (5.3)$$

$$\leq \sum_i \mathbb{E}[{}^\infty h_{C_i \rightarrow 0}(s_0 = {}^\infty X_0^{i-1}, \dots, s_{i-1} = {}^\infty X_{i-1}^{i-1})] \quad (5.4)$$

$$\leq \sum_i \mathbb{E}[F_{C_i}]. \quad (5.5)$$

The sequence of inequalities may not hold in distribution because the variables X^i and thus $h_{C_i \rightarrow 0}(s_0 = X_0^{i-1}, \dots, s_{i-1} = X_{i-1}^{i-1})$ are not independent.

Using (5.2),

$$\mathbb{E}\tau^b \leq \sum_i \frac{\Lambda}{\ell_i + \mu_i} \left(\frac{C_i}{q-p} - \frac{p(1 - (\frac{p}{q})^{C_i})}{(q-p)^2} \right).$$

The result of this part is summarized in the following theorem.

THEOREM 5.3. *In an acyclic stable network of $K + 1$ $/M/1/C_i$ queues with Bernoulli routing and losses in case of overflow, the coupling time from the past sat-*

isfies in expectation,

$$\mathbb{E}[\tau^b] \leq \sum_{i=0}^K \frac{\Lambda}{\ell_i + \mu_i} \left(\frac{C_i}{q-p} - \frac{p(1 - (\frac{p}{q})^{C_i})}{(q-p)^2} \right) \leq \sum_{i=0}^K \frac{\Lambda}{\ell_i + \mu_i} (C_i + C_i^2). \quad (5.6)$$

Note that this bound on the expectation is ultimately linear in the rate of any event in the system. This behavior is also noticeable for $\mathbb{E}[\tau^b]$ itself.

5.2. Some numerical experiments. In the construction of the bound given in Theorem 5.3, several factors may be responsible for the inaccuracy of the bound.

1. The first factor is the replacement of the max by the sum. We believe that it may be a hard task to get rid of this first approximation because of the intricate dependencies between the queues. Furthermore, experiments reported below show that this may not even be possible in many cases (see Figure 5.2.b).

2. Another factor which may increase the inaccuracy of our bounds is the fact that most events change the states of several queues at the same time, while the bound given here disregards this. In the network studied here, this may add a factor 2 between the true coupling time and the bound given in Theorem 5.3.

3. The most important factor which jeopardizes the quality of the bound is the load issue. If one queues has an heavy load (larger than 1), the bound provided by Equation (5.6), also called the light traffic bound in Proposition 4.5 is very bad (as seen in Figure 4.3). So far we have not been able to come up with a better bound for queues with large loads. However, when all queues have a small load (smaller than one, and even more so when the load is smaller than 2/3), the bound tends to be more accurate. This is further verified in the experiments reported below.

Computations for the network displayed in Figure 3.1 are reported in Figure 5.2. We have used the following parameters. The input rate is $\lambda_0 = 0.4$. the rates of the other events are $\lambda_1 = 1.4$, $\lambda_2 = 0.6$, $\lambda_3 = 0.8$, $\lambda_4 = 0.5$. The number of simulation runs is 10000. The capacity C is the same in all queues, and we let it vary from 1 to 20. The service rate in the last queue λ_5 takes three values, respectively 0.2, 0.6 and 0.4.

In the first case (Fig. 5.2.a), $\lambda_5 = 0.2$ so that queue Q_3 has a load larger than one. Figure 5.2 displays the bound given by formula (5.6) as well as the mean coupling time computed over 10000 simulation runs. As hinted before, the bound is indeed very bad for this system. A ratio larger than 10 w.r.t the true coupling time is reached when $C = 5$. It should also be noticed that our bound is convex in C while the coupling time does not seem to be so.

In the second case (Fig. 5.2.b), $\lambda_5 = 0.6$, and all queues have a load smaller that 2/3. Figure 5.2.b shows the bound provided by (5.6) and the true coupling time computed by simulation runs. Both curves appear to be almost linear in C (this is true for the bound: when q/p is small, $\mathbb{E}H_{C_i,0}$ is almost linear in C_i) and the ratio is smaller than 1.3. In that case, the curve $\max_{i \in \{0, \dots, K\}} \mathbb{E}H_{C_i,0}$ is also displayed and is below the actual coupling time. This is to be related with the first item in the comments above.

The last case (Fig. 5.2.c) is for $\lambda_0 = 0.4$, so that Q_3 has load exactly one. This would correspond to the maximal coupling time for Q_3 if it were alone. Figure 5.2.c displays the backward coupling time and the bound provided by Equation (5.6). For queue Q_3 , we use a bound in $C_3 + C_3^2$ which is a bad approximation because of the loss of the factor 2 when compared with the bound for isolated queues. Note that the total gap has a ratio which is almost 2. In that case both the coupling time and the bound exhibit a convex behavior w.r.t. C .

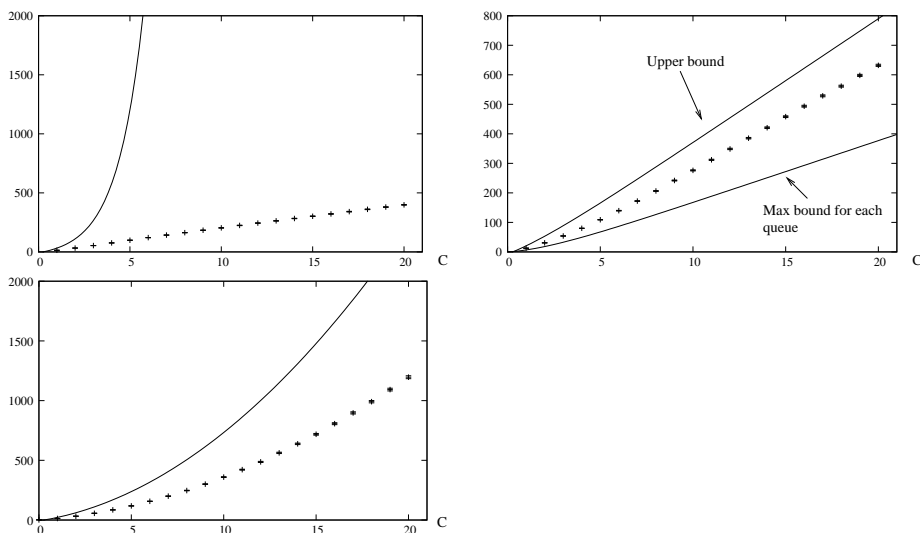


FIG. 5.2. The capacity C varies from 1 to 20 in all queues. The upper left figure displays the coupling time (dots) with 95% confidence intervals, and the bound given by Equation (5.6) when Queue Q_3 is unstable ($\lambda_5 = 2/10$). In the upper right figure, are given the bound in Equation (5.6), the mean coupling time (dots) with 95% confidence intervals and the maximum over Equations (5.6) for all queues, when Queue Q_3 is stable ($\lambda_5 = 6/10$). The lower figure displays the coupling time (dots) with 95% confidence interval and the bound given by Equation (5.6) when Queue Q_3 is barely unstable ($\lambda_5 = 4/10$)

A ratio smaller than 2 is indeed interesting because efficient perfect simulation algorithm use a doubling window technique to reduce the complexity and their running time (see Equation (2.4)) so that our bound gives a good estimation of the mean running time of the algorithms.

One should also note that, on a practical point of view, most actual networks which require stationary performance evaluations have small loads.

5.2.1. Extension to more general networks. Actually, extensive simulation runs over many examples show that the bound given in Theorem 5.3 is robust and also holds for more general networks with blocking and with circuits. While we have only been able to show that the light traffic bound holds for each queue, we conjecture that the heavy traffic bound and the critical bound should also hold. This would yield an overall quadratic bound: $\mathbb{E}[\tau^b] \leq \sum_{i=0}^K \frac{\Lambda}{\ell_i + \mu_i} O(C_i^2)$, for any monotone Markovian network of queues with a finite state space. Furthermore under light or heavy traffic in all queues, the bound should rather be linear: $\mathbb{E}[\tau^b] \leq \sum_{i=0}^K \frac{\Lambda}{\ell_i + \mu_i} O(C_i)$.

To illustrate this conjecture, we have run simulations for the network displayed in Figure 3.1 with the following parameters. The rates are $\lambda_0 = 0.4$, $\lambda_1 = 1.4$, $\lambda_2 = 0.6$, $\lambda_3 = 0.8$, $\lambda_4 = 0.5$. The capacity is fixed to 10 in all queues and we let λ_5 (the service rate in Q_3) vary from 0 to 4. As long as $\lambda_5 < 0.4$, Q_3 is unstable and our proven bound (B_1) is poor. As soon as λ_5 is large enough our bound becomes acceptable. In Figure 5.3, note that both the bound and the coupling time τ have a linear asymptotic growth in λ_5 . The Figure also displays the heavy traffic bound B_2 and the critical bound B_3 . Should these two bounds hold, the minimum of B_1, B_2, B_3 (in bold in the figure) would provide a remarkable bound on the coupling time, up to an additional constant. This issue is the subject of our current investigations.

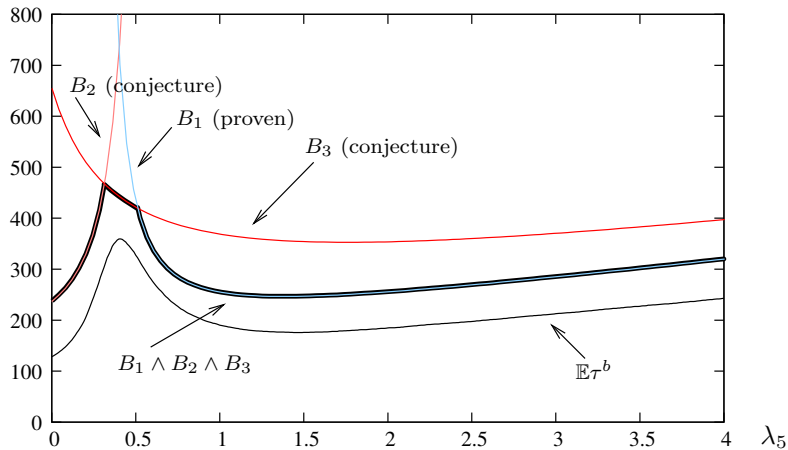


FIG. 5.3. This figure displays the actual coupling time $\mathbb{E}\tau^b$ for the network of Figure 3.1, when the service rate of the last queue ranges from 0 to 5, together with the proven light traffic bound B_1 , the conjectured heavy traffic bound B_2 , the conjectured critical bound B_3 and the minimum of the three bounds.

Acknowledgments. The authors would like to thank Jérôme Vienne who partially designed the `psi2` perfect simulation software which was used to run all the simulations presented here.

REFERENCES

- [1] C. ALEXOPOULOS AND D. GOLDSMAN, *To batch or not to batch?*, ACM Trans. Model. Comput. Simul., 14 (2004), pp. 76–114.
- [2] J.M. CALVIN, P.W. GLYNN, AND M.K. NAKAYAMA, *The semi-regenerative method of simulation output analysis*. submitted.
- [3] ———, *Importance sampling using the semi-regenerative method*, in Proceedings of the Winter Simulation Conference, vol. 1, 2001, pp. 441–450.
- [4] M. CRANE AND D.L. IGLEHART, *Simulating stable stochastic systems, iii: Regenerative processes and discrete-event simulation*, Operation Research, 23 (1975), pp. 33–45.
- [5] P. GLASSERMAN AND D.D. YAO, *Monotone Structure in Discrete-Event Systems*, Wiley InterScience, Series in Probability and Mathematical Statistics, 1994.
- [6] O. HÄGGSTRÖM, *Finite Markov Chains and Algorithmic Applications*, Cambridge University Press, 2002.
- [7] S.G. HENDERSON AND P.W. GLYNN, *Regenerative steady-state simulation of discrete-event systems*, ACM Trans. Model. Comput. Simul., 11 (2001), pp. 313–345.
- [8] J. PROPP AND D. WILSON, *Exact sampling with coupled Markov chains and applications to statistical mechanics*, Random Structures and Algorithms, 9 (1996), pp. 223–252.
- [9] S. M. ROSS, *Probability models*, Academic Press, 2003.
- [10] W.J. STEWART, *Introduction to the Numerical Solution of Markov Chains*, Princeton, 1994.
- [11] J.-M. VINCENT, *Perfect simulation of monotone systems for rare event probability estimation*, in Winter Simulation Conference, Orlando, dec 2005.
- [12] ———, *Perfect simulation of queueing networks with blocking and rejection*, in Saint IEEE conference, Trento, 2005, pp. 268–271.
- [13] J.-M. VINCENT AND C. MARCHAND, *On the exact simulation of functionals of stationary markov chains*, Linear Algebra and its Applications, 386 (2004), pp. 285–310.