RESEARCH ARTICLE

Analysis of a large number of Markov chains competing for transitions

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We consider the behavior of a stochastic system composed of several identically distributed, but non independent, discrete-time absorbing Markov chains competing at each instant for a transition. The competition consists in determining at each instant, using a given probability distribution, the only Markov chain allowed to make a transition. We analyze the first time at which one of the Markov chains reaches its absorbing state. When the number of Markov chains goes to infinity, we analyze the asymptotic behavior of the system for an arbitrary probability mass function governing the competition. We give conditions for the existence of the asymptotic distribution and we show how these results apply to cluster-based distributed storage when the competition is handled using a geometric distribution.

Keywords: asymptotic analysis; competing Markov chains; cluster-based distributed systems; Markov chains, geometric distribution

1. Introduction

Competing Markov chains generally contend with each other for a set of resources, see for instance Fourneau (2008) and the references therein. The resulting process is then a multidimensional Markov chain based on the Cartesian product of the states spaces and on competition rules over resources. In this paper, the Markov chains do not compete for resources but for transitions. More precisely, we consider a stochastic system composed of n identically distributed, but non independent, discrete-time absorbing Markov chains competing at each instant for a transition. The competition consists in determining at each instant, using a given probability mass function with n values, the only Markov chain allowed to make a transition.

For this system, we analyze the absorption time Θ_n^{γ} at which one of the *n* Markov chains reaches its absorbing state, when the probability mass function is $\gamma(n)$. The distribution of this random variable has already been studied in Anceaume et al. (2010), in particular when the probability mass function $\gamma(n)$ handling the competition is uniform. In that case, we studied the asymptotic behavior of the system when the number *n* of Markov chains goes to infinity.

We propose here the study of the asymptotic behavior of the system when the number n of Markov chains goes to infinity, for an arbitrary probability mass function $\gamma(n)$ governing the competition. More precisely, we give conditions on the probability mass function $\gamma(n)$ governing the competition for the existence of a limiting distribution for Θ_n^{γ} . We apply these results to the

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case where the competition is governed by a geometric distribution and we study the effects of this distribution on a model of a large-scale cluster-based distributed storage.

The remainder part of the paper is organized as follows. In the next section, we describe the model, the notation and we give the transient state distribution of the global Markov chain composed of the *n* joined identically distributed local Markov chains. We also extend a result obtained in Anceaume et al. (2010). This result is a recurrence relation that allows us not only to compute the distributions of Θ_n^{γ} but also to compute the limiting distribution of Θ_n^{γ} when it exists. In Section 3, we study the asymptotic behavior of the system when *n* goes to infinity and we give conditions on the probability mass function $\gamma(n)$ governing the competition for the existence of the limiting distribution of Θ_n^{γ} . We also show how to compute this limiting distribution. We apply these results in Section 4 to the case where the probability mass function $\gamma(n)$ governing the competition is geometric. Section 5 is devoted to an application of these results to a model of a cluster-based distributed storage.

2. Transient Analysis

We consider a homogeneous discrete-time Markov chain $X = \{X_k, k \ge 0\}$ with a finite state space S composed of a set of transient states denoted by B and an absorbing state denoted by a. The associated transition probability matrix P can thus be partitioned as

$$P = \begin{pmatrix} Q \ v \\ 0 \ 1 \end{pmatrix},$$

where Q is the submatrix of dimension $|B| \times |B|$ containing the transitions between states of B. In the same way, v is the column vector with |B| entries representing the transitions from the transient states to the absorbing state. We suppose that the initial state lies in B, i.e. $\mathbb{P}\{X_0 \in B\} = 1$, and we denote by α the row vector of dimension |B| containing the initial probability distribution, i.e. for every $i \in B$, $\alpha_i = \mathbb{P}\{X_0 = i\}$. We denote by Θ_1 the total time the Markov chain X spends in B or equivalently the first instant at which the absorbing state a is reached. We have $\Theta_1 = \inf\{k \ge 0 \mid X_k = a\}$. The complementary cumulative distribution function of Θ_1 is easily derived as, see for instance Neuts (1981),

$$\mathbb{P}\{\Theta_1 > k\} = \mathbb{P}\{X_k \in B\} = \alpha Q^k \mathbb{1},$$

where $\mathbb{1}$ is the column vector of dimension |B| with all entries equal to 1. Since all the states of B are transient, the matrix I - Q is invertible and the expectation of Θ_1 is given by $E(\Theta_1) = \alpha(I - Q)^{-1}\mathbb{1}$, where I is the identity matrix.

Let us now consider, for $n \ge 1$, *n* Markov chains denoted by $X^{(1)}, \ldots, X^{(n)}$ stochastically equivalent to X, i.e. with the same state space S, the same transition probability matrix P and the same initial probability distribution α . These *n* Markov chains are in competition at each instant to make a transition using the probability mass function $\gamma(n) = (p_{1,n}, \ldots, p_{n,n})$.

From these *n* Markov chains, we construct a new Markov chain denoted by $Y = \{Y_k, k \ge 0\}$ as follows. The state space of *Y* is equal to S^n and $Y_k = (X_k^{(1)}, \ldots, X_k^{(n)})$. A transition in the Markov chain *Y* corresponds to a transition in only one of the Markov chains $X^{(1)}, \ldots, X^{(n)}$, all the rest staying in the same state. The Markov chain that makes the transition is chosen in accordance with the probability mass function $\gamma(n)$, which means that the Markov chain $X^{(\ell)}$ makes the transition with probability $p_{\ell,n}$. We suppose, without loss of generality that for any $\ell, 0 < p_{\ell,n} < 1$.

The transition probability matrix of Y is detailed in Anceaume et al. (2010) where we give the transient distribution of the Markov chain Y. The first instant Θ_n^{γ} at which one of the n Markov

chains $X^{(1)}, \ldots, X^{(n)}$ gets absorbed when the probability mass function is $\gamma(n)$ is defined as

$$\Theta_n^{\gamma} = \inf\{k \ge 0 \mid \exists r \text{ such that } X_k^{(r)} = a\}.$$

When n = 1, we have $\gamma(1) = 1$ and, thus $\Theta_1^{\gamma} = \Theta_1$. We give below its distribution which has been obtained in Anceaume et al. (2010). We first introduce the set $S_{k,\ell}$ defined, for every $k \ge 0$ and $\ell \ge 1$, by $S_{k,\ell} = \{\underline{k} = (k_1, \ldots, k_\ell) \in \mathbb{N}^\ell \mid k_1 + \cdots + k_\ell = k\}$. For every $k \ge 0$ and $n \ge 1$, we have

$$\mathbb{P}\{\Theta_n^{\gamma} > k\} = \sum_{\underline{k} \in S_{k,n}} \frac{k!}{k_1! \cdots k_n!} \prod_{r=1}^n (p_{r,n})^{k_r} \alpha Q^{k_r} \mathbb{1}.$$
(1)

The computational complexity of $\mathbb{P}\{\Theta_n^{\gamma} > k\}$ using (1) is exponential. A way to reduce this complexity is provided by the following theorem. This theorem improves the one obtained in Anceaume et al. (2010) which has been obtained only for h = n.

Theorem 2.1: For every $k \ge 0$, $n \ge 2$ and h = 1, ..., n, we have

$$\mathbb{P}\{\Theta_{n}^{\gamma} > k\} = \sum_{\ell=0}^{k} \binom{k}{\ell} (p_{h,n})^{\ell} (1 - p_{h,n})^{k-\ell} \alpha Q^{\ell} \mathbb{1P}\{\Theta_{n-1}^{\gamma^{(1)}} > k - \ell\},$$
(2)

where the probability mass function $\gamma^{(1)}(n-1) = (p_{1,n-1}^{(1)}, \dots, p_{n-1,n-1}^{(1)})$ associated with $\Theta_{n-1}^{\gamma^{(1)}}$ and with the choice of h is defined, by

$$p_{r,n-1}^{(1)} = \frac{p_{r,n}}{1 - p_{h,n}}$$
 for $r = 1, \dots, h-1$ and $p_{r,n-1}^{(1)} = \frac{p_{r+1,n}}{1 - p_{h,n}}$ for $r = h, \dots, n-1$.

Proof: For every $k \ge 0$ and $n \ge 2$, we fix a value of h with $1 \le h \le n$. In Relation (1), we extract the index k_h in the multiple sum indexed by $\underline{k} \in S_{k,n}$. Then, we rename it ℓ and next, for h < n, we perform the changes of variables $k_{h+1} := k_h, \ldots, k_n := k_{n-1}$. We thus obtain

$$\mathbb{P}\{\Theta_{n}^{\gamma} > k\} = \sum_{\ell=0}^{k} \frac{(p_{h,n})^{\ell} \alpha Q^{\ell} \mathbb{1}}{\ell!} \sum_{\underline{k} \in S_{k-\ell,n-1}} \frac{k!}{k_{1}! \cdots k_{n-1}!} \prod_{r=1}^{h-1} (p_{r,n})^{k_{r}} \alpha Q^{k_{r}} \mathbb{1} \prod_{r=h}^{n-1} (p_{r+1,n})^{k_{r}} \alpha Q^{k_{r}} \mathbb{1}.$$

Multiplying and dividing the right-hand side respectively by $(k - \ell)!$ and $(1 - p_{h,n})^{k-\ell}$, we get

$$\mathbb{P}\{\Theta_{n}^{\gamma} > k\} = \sum_{\ell=0}^{k} \binom{k}{\ell} (p_{h,n})^{\ell} (1 - p_{h,n})^{k-\ell} \alpha Q^{\ell} \mathbb{1}$$
$$\times \sum_{\underline{k} \in S_{k-\ell,n-1}} \frac{(k-\ell)!}{k_{1}! \cdots k_{n-1}!} \prod_{r=1}^{h-1} \left(\frac{p_{r,n}}{1 - p_{h,n}}\right)^{k_{r}} \alpha Q^{k_{r}} \mathbb{1} \prod_{r=h}^{n-1} \left(\frac{p_{r+1,n}}{1 - p_{h,n}}\right)^{k_{r}} \alpha Q^{k_{r}} \mathbb{1}.$$

Using the definition of $\gamma^{(1)}(n-1)$, we obtain, from Relation (1) again,

$$\mathbb{P}\{\Theta_{n}^{\gamma} > k\} = \sum_{\ell=0}^{k} \binom{k}{\ell} (p_{h,n})^{\ell} (1-p_{h,n})^{k-\ell} \alpha Q^{\ell} \mathbb{1} \sum_{\underline{k} \in S_{k-\ell,n-1}} \frac{(k-\ell)!}{k_{1}! \cdots k_{n-1}!} \prod_{r=1}^{n-1} \left(p_{r,n-1}^{(1)} \right)^{k_{r}} \alpha Q^{k_{r}} \mathbb{1}$$
$$= \sum_{\ell=0}^{k} \binom{k}{\ell} (p_{h,n})^{\ell} (1-p_{h,n})^{k-\ell} \alpha Q^{\ell} \mathbb{1} \mathbb{P}\{\Theta_{n-1}^{\gamma^{(1)}} > k-l\},$$

which completes the proof.

This result shows that the computation of $\mathbb{P}\{\Theta_n^{\gamma} > k\}$ can be done using a simple recurrence with a polynomial complexity. The expectation of Θ_n^{γ} is then obtained by

$$E(\Theta_n^{\gamma}) = \sum_{k=0}^{\infty} \mathbb{P}\{\Theta_n^{\gamma} > k\}.$$
(3)

3. Asymptotic Analysis

This section is devoted to the analysis of the distribution on Θ_n^{γ} when n is large. This is generally the case in practice for large-scale distributed systems which are studied in the last section. For every $n \geq 1$ and $x \in \mathbb{R}$, we introduce the transform $F_n(x)$ defined by

$$F_n(x) = \sum_{k=0}^{\infty} \frac{x^k}{k!} \mathbb{P}\{\Theta_n^{\gamma} > k\}$$

An explicit expression of function F_n is given in the following theorem, which is proved in Anceaume et al. (2010).

Theorem 3.1: For every $n \ge 1$ and $x \in \mathbb{R}$, we have

$$F_n(x) = \prod_{\ell=1}^n \alpha e^{Qxp_{\ell,n}} \mathbb{1},\tag{4}$$

and, for every $k \ge 0$, $\mathbb{P}\{\Theta_n^{\gamma} > k\} = F_n^{(k)}(0)$, where $F_n^{(k)}$ is the k-th derivative of the function F_n with respect to x.

This result not only shows that $\mathbb{P}\{\Theta_n^{\gamma} > 0\} = 1$ as expected, but it also shows that, for every $n \ge 1$, we have $\mathbb{P}\{\Theta_n^{\gamma} > 1\} = \alpha Q \mathbb{1} = F'_n(0)$. It also gives access to an expression of $\mathbb{P}\{\Theta_n^{\gamma} > k\}$ for any k. Adopting this point of view, our strategy in order to compute $\lim_{n\to\infty} \mathbb{P}\{\Theta_n^{\gamma} > k\}$ is to compute $F(x) = \lim_{n\to\infty} F_n(x)$, an analytic function of x, so as to deduce the value $\lim_{n\to\infty} \mathbb{P}\{\Theta_n^{\gamma} > k\} = F^{(k)}(0)$. To pass to the limit in a clean fashion, we need the following. **Hypothesis (H)** – Limiting value of the powers sums of the $p_{\ell,n}$'s.

For any
$$k \ge 1$$
, the limit $V_k = \lim_{n \to \infty} \sum_{\ell=1}^n (p_{\ell,n})^k$ exists.

Important remark. The above assumption is harmless. Indeed, introducing the quantities $V_{n,k} = \sum_{\ell=1}^{n} (p_{\ell,n})^k$, it is clear that $0 \leq V_{n,k} \leq 1$ for any value of $n \geq 1$ and $k \geq 1$. Therefore, there exists a subsequence in n, say n_j with $n_j \to \infty$ as $j \to \infty$, such that $V_{n_j,k}$ has a limit as $j \to \infty$ for any $k \geq 1$. We are here merely assuming that the limit V_k is well defined

without referring to taking a subsequence in the original $V_{n,k}$'s. To give an example, suppose that the $p_{\ell,2n}$'s are uniformly distributed, i.e. $p_{\ell,2n} = 1/(2n)$, in which case $V_{2n,k} = 1/(2n)^{k-1} \to 0$ whenever $k \geq 2$, and $V_{2n,1} = 1$, while the $p_{\ell,2n+1}$'s are geometrically distributed with parameter b and truncation at step 2n+1, i.e. $p_{\ell,2n+1} = (1-b)^{\ell-1}b$, for $\ell \leq 2n$ and $p_{2n+1,2n+1} = (1-b)^{2n}$, in which case $V_{2n+1,k} \to b^k/(1-(1-b)^k)$. In that case it clearly does not make sense to study the whole sequence $V_{n,k}$ itself, and we need to separate the case when n is odd and the case when n is even. We assert here that this situation is generic, and that, up to extracting a subsequence, one may always assume that the original sequence $V_{n,k}$ itself possesses a limit V_k for any k.

With this assumption in mind, the following theorem gives the limit of the transform $F_n(x)$. This result was obtained in Anceaume et al. (2010) only in the case where the $p_{\ell,n}$'s are uniformly distributed.

Theorem 3.2: Under hypothesis (H), $F(x) = \lim_{n \to \infty} F_n(x)$ exists, whenever $|x| < \ln 2$, and the limit is uniform on compact subsets of $\{x \mid |x| < \ln 2\}$. Besides, we have the explicit value

$$F(x) = \exp\left(\sum_{m \ge 1} \sum_{k_1 \ge 1} \cdots \sum_{k_m \ge 1} \frac{(-1)^{m+1}}{m} \frac{\alpha Q^{k_1} \mathbb{1} \dots \alpha Q^{k_m} \mathbb{1}}{k_1! \dots k_m!} x^{k_1 + \dots + k_m} V_{k_1 + \dots + k_m}\right).$$
 (5)

Proof: Starting from Relation (4), we recover, expanding into power series in x, the value

$$\ln(F_n(x)) = \sum_{\ell=1}^n \ln\left(\alpha \, e^{Qxp_{\ell,n}} \mathbb{1}\right) = \sum_{\ell=1}^n \ln\left(1 + \sum_{k\geq 1} \frac{\alpha Q^k \mathbb{1}}{k!} x^k (p_{\ell,n})^k\right)$$

Hence, using the fact that $0 \leq \alpha Q^k \mathbb{1} \leq 1$ whenever $k \geq 0$, and deducing the bound

$$\left| \sum_{k \ge 1} \frac{\alpha Q^k \mathbb{1}}{k!} x^k (p_{\ell,n})^k \right| \le \sum_{k \ge 1} \frac{|x|^k}{k!} = e^{|x|} - 1 < 1,$$

whenever $|x| < \ln 2$, we may expand further and obtain

$$\ln (F_n(x)) = \sum_{\ell=1}^n \sum_{m \ge 1} \frac{(-1)^{m+1}}{m} \left(\sum_{k \ge 1} \frac{\alpha Q^k \mathbb{1}}{k!} x^k (p_{\ell,n})^k \right)^m$$

= $\sum_{\ell=1}^n \sum_{m \ge 1} \frac{(-1)^{m+1}}{m} \sum_{k_1 \ge 1} \cdots \sum_{k_m \ge 1} \frac{\alpha Q^{k_1} \mathbb{1} \dots \alpha Q^{k_m} \mathbb{1}}{k_1! \dots k_m!} x^{k_1 + \dots + k_m} (p_{\ell,n})^{k_1 + \dots + k_m}$
= $\sum_{m \ge 1} \sum_{k_1 \ge 1} \cdots \sum_{k_m \ge 1} \frac{(-1)^{m+1}}{m} \frac{\alpha Q^{k_1} \mathbb{1} \dots \alpha Q^{k_m} \mathbb{1}}{k_1! \dots k_m!} x^{k_1 + \dots + k_m} \left(\sum_{\ell=1}^n (p_{\ell,n})^{k_1 + \dots + k_m} \right).$

The above expansions clearly converge in any desirable sense whenever $|x| < \ln 2$ (say, for instance, uniformly in x on compact subsets of $\{x \mid |x| < \ln 2\}$). The existence of the limiting values V_k , together with the pointwise bound

$$\left|\frac{(-1)^{m+1}}{m}\frac{\alpha Q^{k_1}\mathbb{1}\dots\alpha Q^{k_p}\mathbb{1}}{k_1!\dots k_m!}\,x^{k_1+\dots+k_m}\left(\sum_{\ell=1}^n (p_{\ell,n})^{k_1+\dots+k_m}\right)\right| \leq \frac{|x|^{k_1+\dots+k_m}}{k_1!\dots k_m!},$$

a converging series whenever $|x| < \ln 2$, therefore provides the limit

$$\lim_{n \to \infty} \ln (F_n(x)) = \sum_{m \ge 1} \sum_{k_1 \ge 1} \cdots \sum_{k_m \ge 1} \frac{(-1)^{m+1}}{m} \frac{\alpha Q^{k_1} \mathbb{1} \dots \alpha Q^{k_p} \mathbb{1}}{k_1! \dots k_m!} x^{k_1 + \dots + k_m} V_{k_1 + \dots + k_m},$$

and the above convergence is uniform on compact subsets of $\{x \mid |x| < \ln 2\}$.

Armed with the above theorem, we deduce the limiting behavior of $\mathbb{P}\{\Theta_n^{\gamma} > k\}$.

Theorem 3.3: Under hypothesis (H), for every $k \ge 0$, we have

$$\lim_{n \to \infty} \mathbb{P}\{\Theta_n^{\gamma} > k\} = F^{(k)}(0).$$

Proof: The argument is standard. The function $F_n(x)$ being clearly analytic on the disk $\{z \in \mathbb{C} \mid |z| < \ln 2\}$, we may write, for any $0 < r < \ln 2$ and $k \ge 0$, the relation

$$\mathbb{P}\{\Theta_n > k\} = F_n^{(k)}(0) = \frac{k!}{2i\pi} \int_{|z|=r} \frac{F_n(z)}{z^{k+1}} dz.$$

Hence, using the above-mentioned uniform convergence of F_n towards F, we recover

$$\lim_{n \to \infty} \mathbb{P}\{\Theta_n > k\} = \lim_{n \to \infty} \frac{k!}{2i\pi} \int_{|z|=r} \frac{F_n(z)}{z^{k+1}} dz = \frac{k!}{2i\pi} \int_{|z|=r} \frac{F(z)}{z^{k+1}} dz.$$

On the other hand, in view of Relation (5), the function F is itself clearly analytic on the disk $\{z \in \mathbb{C} \mid |z| < \ln 2\}$, so we may write similarly

$$\frac{k!}{2i\pi} \int_{|z|=r} \frac{F(z)}{z^{k+1}} \, dz = F^{(k)}(0)$$

which completes the proof.

We denote by Θ^{γ} the random variable having as distribution the limiting distribution of Θ_n^{γ} . We then have from Theorem 3.3, under hypothesis (H), for every $k \ge 0$,

$$\lim_{n \to \infty} \mathbb{P}\{\Theta_n^{\gamma} > k\} = \mathbb{P}\{\Theta^{\gamma} > k\} = F^{(k)}(0).$$

The following two corollaries show how to compute recursively the distribution of Θ^{γ} . Again, note that this limiting distribution was obtained in Anceaume et al. (2010) only in the case where the $p_{\ell,n}$'s are uniformly distributed.

Corollary 3.4: Under hypothesis (H) and if, for a fixed $h \ge 1$, we have $\lim_{n \to \infty} p_{h,n} = b \in (0,1)$, then, we have, for every $k \ge 0$,

$$\mathbb{P}\{\Theta^{\gamma} > k\} = \sum_{\ell=0}^{k} \binom{k}{\ell} b^{\ell} (1-b)^{k-\ell} \alpha Q^{\ell} \mathbb{1P}\{\Theta^{\gamma^{(1)}} > k-\ell\}.$$
(6)

Proof: In order to use Theorem 2.1, we have to check that hypothesis (H) is valid for distribution $\gamma^{(1)}(n-1)$. If hypothesis (H) is satisfied and if, for a fixed $h \ge 1$, $\lim_{n \to \infty} p_{h,n} = b \in (0,1)$,

then we have

$$V_{n-1,k}^{(1)} = \sum_{\ell=1}^{n-1} (p_{\ell,n-1}^{(1)})^k = \frac{1}{(1-p_{h,n})^k} \left(\sum_{\ell=1}^{h-1} p_{\ell,n} + \sum_{\ell=h+1}^n p_{\ell,n} \right) = \frac{V_{n,k} - (p_{h,n})^k}{(1-p_{h,n})^k},$$

so the limit $V_k^{(1)}$ exists and we have

$$V_k^{(1)} = \lim_{n \to \infty} V_{n-1,k}^{(1)} = \frac{V_k - b^k}{(1-b)^k}$$

Using Theorem 3.3 and taking the limit in Relation (2), we get Relation (6).

To illustrate this last result, consider for instance the distribution $\gamma(n) = (1/2, 1/2(n-1), \ldots, 1/2(n-1))$. Hypothesis (H) is clearly satisfied (we have $V_k = 1/2^k$, for every $k \ge 1$). Since $p_{1,n} = 1/2$, we take h = 1. We then have $\gamma^{(1)}(n-1) = (1/(n-1), \ldots, 1/(n-1))$, which is the uniform distribution. It has been shown in Anceaume et al. (2010) that $\mathbb{P}\{\Theta^{\gamma^{(1)}} > k\} = (\alpha Q \mathbb{1})^k$. We thus get from Relation (6)

$$\mathbb{P}\{\Theta^{\gamma} > k\} = \frac{1}{2^{k}} \sum_{\ell=0}^{k} \binom{k}{\ell} \alpha Q^{\ell} \mathbb{1}(\alpha Q \mathbb{1})^{k-\ell} = \alpha \left(\frac{1}{2}Q + \frac{\alpha Q \mathbb{1}}{2}I\right)^{k} \mathbb{1}$$

Corollary 3.5: Under hypothesis (H), if, for a fixed $h \ge 1$, we have $\lim_{n \to \infty} p_{h,n} = b \in (0,1)$ and if $\gamma^{(1)}(n-1) = \gamma(n-1)$, for every $n \ge 2$ then, we have $\mathbb{P}\{\Theta^{\gamma} > 0\} = 1$ and, for every $k \ge 1$,

$$\mathbb{P}\{\Theta^{\gamma} > k\} = \frac{1}{1 - (1 - b)^k} \sum_{\ell=0}^{k-1} \binom{k}{\ell} (1 - b)^\ell b^{k-\ell} \mathbb{P}\{\Theta^{\gamma} > \ell\} \alpha Q^{k-\ell} \mathbb{1}$$

Proof: Since $\mathbb{P}\{\Theta_n^{\gamma} > 0\} = 1$ for every $n \ge 1$, we have $\mathbb{P}\{\Theta^{\gamma} > 0\} = 1$. From Corollary 3.4 and since $\gamma^{(1)}(n) = \gamma(n)$ for every $n \ge 2$, we have

$$\mathbb{P}\{\Theta^{\gamma} > k\} = \sum_{\ell=0}^{k} \binom{k}{\ell} b^{\ell} (1-b)^{k-\ell} \alpha Q^{\ell} \mathbb{1}\mathbb{P}\{\Theta^{\gamma} > k-\ell\}.$$

Extracting the term containing $\mathbb{P}\{\Theta^{\gamma} > k\}$, which corresponds to index $\ell = 0$, from the right hand side, we get the desired relation.

Without any loss of generality, by renumbering the Markov chains, we take in the remainder part of the paper h = 1. This means, from Theorem 2.1, that the probability mass function $\gamma^{(1)}(n-1) = (p_{1,n-1}^{(1)}, \dots, p_{n-1,n-1}^{(1)})$ associated with $\Theta_{n-1}^{\gamma^{(1)}}$ is given, for $r = 1, \dots, n-1$, by

$$p_{r,n-1}^{(1)} = \frac{p_{r+1,n}}{1 - p_{1,n}}.$$
(7)

For a fixed value of $n \geq 2$, the computation of the distribution of Θ_n^{γ} with a given probability mass function $\gamma(n)$ necessitates the computation of the distribution of $\Theta_{n-1}^{\gamma^{(1)}}$ with the probability mass function $\gamma^{(1)}(n-1)$ given by Relation (7). Let ε be a predetermined error tolerance. If we want to compute $\mathbb{P}\{\Theta_n^{\gamma} > k\}$ for every k such that $\mathbb{P}\{\Theta_n^{\gamma} > k\} > \varepsilon$, we need to determine an

integer K such that, for every i = 1, ..., n, $\mathbb{P}\{\Theta_i^{\gamma^{(n-i)}} > K\} \leq \varepsilon$, and then to compute, from Relation (2), for i = 1, ..., n, the values of $\mathbb{P}\{\Theta_i^{\gamma^{(n-i)}} > k\}$ for k = 0, ..., K - 1. Note that the probability mass function associated with $\Theta_i^{\gamma^{(n-i)}}$ is $\gamma^{(n-i)}(i) = (p_{1,i}^{(n-i)}, ..., p_{i,i}^{(n-i)})$ which is obtained from $\gamma^{(n-i-1)}(i+1)$ by normalizing its last *i* entries by $1 - p_{1,i+1}^{(n-i-1)}$, where we set $\gamma^{(0)}(n) = \gamma(n)$ and $p_{\ell,n}^{(0)} = p_{\ell,n}$. The following lemma will be used in the next theorem where we propose a value of K. An inequality between vectors is meant entrywise.

Lemma 3.6: For every $k \ge 1$, the vector function f(x) defined by $f(x) = (xQ + (1-x)I)^k \mathbb{1}$, for $x \in [0,1]$, is decreasing.

Proof: The function f is differentiable on the interval (0, 1) and its derivative f' is given by

$$f'(x) = k(xQ + (1-x)I)^{k-1}(Q\mathbb{1} - \mathbb{1}).$$

The matrix Q being substochastic, we have $Q\mathbb{1} - \mathbb{1} \leq 0$ with strict inequality for at least one entry. We thus have $f'(x) \leq 0$ which means that function f is decreasing on interval [0, 1]. \Box

For every $n \ge 1$, we introduce the numbers m_n defined by $m_n = \min_{i=1,...,n} p_{1,i}^{(n-i)}$. The two following theorems are improvements of the corresponding results obtained in Anceaume et al. (2010) since they do not use the fact that the $p_{\ell,n}$'s are decreasing with ℓ .

Theorem 3.7: For every $n \ge 1$, for every $\varepsilon \in (0, 1)$, we have

$$\max_{i=1,\dots,n} \mathbb{P}\{\Theta_i^{\gamma^{(n-i)}} > k\} \le \varepsilon \text{ for every } k \ge K,$$

where $K = \inf \left\{ k \ge 0 \mid \alpha \left(m_n Q + (1 - m_n) I \right)^k \mathbb{1} \le \varepsilon \right\}.$

Proof: For every $i = 1, \ldots, n$, we have

$$\mathbb{P}\{\Theta_{i}^{\gamma^{(n-i)}} > k\} = \sum_{\ell=0}^{k} \binom{k}{\ell} (p_{1,i}^{(n-i)})^{\ell} (1 - p_{1,i}^{(n-i)})^{k-\ell} \alpha Q^{\ell} \mathbb{1}\mathbb{P}\{\Theta_{i-1}^{\gamma^{(n-i+1)}} > k-\ell\} \\
\leq \sum_{\ell=0}^{k} \binom{k}{\ell} (p_{1,i}^{(n-i)})^{\ell} (1 - p_{1,i}^{(n-i)})^{k-\ell} \alpha Q^{\ell} \mathbb{1} \\
= \alpha \left(p_{1,i}^{(n-i)} Q + (1 - p_{1,i}^{(n-i)}) I \right)^{k} \mathbb{1} \\
\leq \alpha \left(m_{n} Q + (1 - m_{n}) I \right)^{k} \mathbb{1} \quad \text{(from Lemma 3.6).}$$

Note that matrix $m_n Q + (1 - m_n)I$ is substochastic, i.e. $(m_n Q + (1 - m_n)I)\mathbb{1} \leq \mathbb{1}$ with a strict inequality for at least one entry. This implies that $\alpha (m_n Q + (1 - m_n)I)^k \mathbb{1}$ is decreasing with k and that $\lim_{k \to \infty} \alpha (m_n Q + (1 - m_n)I)^k \mathbb{1} = 0$. So, for a fixed $\varepsilon \in (0, 1)$ and by definition of integer K we have, for every $i = 1, \ldots, n$,

$$\mathbb{P}\{\Theta_i^{\gamma^{(n-i)}} > k\} \le \varepsilon, \text{ for every } k \ge K,$$

which completes the proof.

In the same way, we obtain a similar result for the computation of the expected values $E(\Theta_i^{\gamma^{(n-i)}})$, for i = 1, ..., n, for which the truncation of the series (3) is needed.

Theorem 3.8: For every $n \ge 1$, for every $\varepsilon \in (0, 1)$,

$$0 \leq \max_{i=1,\dots,n} \left(E(\Theta_i^{\gamma^{(n-i)}}) - \sum_{k=0}^{L-1} \mathbb{P}\{\Theta_i^{\gamma^{(n-i)}} > k\} \right) \leq \varepsilon,$$
$$L = \inf\left\{ k \geq 0 \ \bigg| \ \frac{1}{m_n} \alpha (I-Q)^{-1} \left(m_n Q + (1-m_n) I\right)^k \mathbbm{1} \leq \varepsilon \right\}.$$

Proof: We introduce the notation

where

$$r_{i} = E(\Theta_{i}^{\gamma^{(n-i)}}) - \sum_{k=0}^{L-1} \mathbb{P}\{\Theta_{i}^{\gamma^{(n-i)}} > k\}.$$

We then have, for every $i = 1, \ldots, n$,

$$\begin{split} r_{i} &= \sum_{k=L}^{\infty} \mathbb{P}\{\Theta_{i}^{\gamma^{(n-i)}} > k\} \\ &= \sum_{k=L}^{\infty} \sum_{\ell=0}^{k} \binom{k}{\ell} (p_{1,i}^{(n-i)})^{\ell} (1 - p_{1,i}^{(n-i)})^{k-\ell} \alpha Q^{\ell} \mathbb{1P}\{\Theta_{i-1}^{\gamma^{(n-i+1)}} > k-\ell\} \\ &\leq \sum_{k=L}^{\infty} \sum_{\ell=0}^{k} \binom{k}{\ell} (p_{1,i}^{(n-i)})^{\ell} (1 - p_{1,i}^{(n-i)})^{k-\ell} \alpha Q^{\ell} \mathbb{1} \\ &= \sum_{k=L}^{\infty} \alpha \left(p_{1,i}^{(n-i)} Q + (1 - p_{1,i}^{(n-i)}) I \right)^{k} \mathbb{1} \\ &\leq \sum_{k=L}^{\infty} \alpha \left(m_{n} Q + (1 - m_{n}) I \right)^{k} \mathbb{1} \qquad (\text{ from Lemma 3.6}) \\ &= \alpha \left(I - (m_{n} Q + (1 - m_{n}) I) \right)^{-1} (m_{n} Q + (1 - m_{n}) I)^{L} \mathbb{1} \\ &= \frac{1}{m_{n}} \alpha (I - Q)^{-1} (m_{n} Q + (1 - m_{n}) I)^{L} \mathbb{1} \\ &\leq \varepsilon \text{ by definition of integer } L. \end{split}$$

which means that $\max_{i=1,\ldots,n} r_i \leq \varepsilon$.

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It is easily checked, from Relation (6), that the same result holds for the limiting expected value $E(\Theta^{\gamma})$. More precisely, if $\lim_{n \to \infty} p_{1,n} = b > 0$, then, for every $\varepsilon \in (0, 1)$, we have

$$0 \leq \left(E(\Theta^{\gamma}) - \sum_{k=0}^{J-1} \mathbb{P}\{\Theta^{\gamma} > k\} \right) \leq \varepsilon,$$

where $J = \inf \left\{ k \geq 0 \ \left| \ \frac{1}{b} \alpha (I-Q)^{-1} \left(bQ + (1-b) I \right)^k \mathbb{1} \leq \varepsilon \right\}.$

4. Geometric distribution

We suppose in this section that the probability mass function $\gamma(n)$ is the one of the geometric distribution with parameter $b \in (0,1)$, truncated at step n, i.e. $\gamma(n)$ is given, for $n \geq 2$ and $r = 1, \ldots, n-1$ by $p_{r,n} = (1-b)^{r-1}b$ and $p_{n,n} = (1-b)^{n-1}$. Clearly hypothesis (H) is satisfied and from Relation (7), we have $p_{1,n} = b$ and thus $\lim_{n \to \infty} p_{1,n} = b$. From Relation (7), we have $\gamma^{(1)}(n-1) = \gamma(n-1)$, for every $n \geq 2$. Thus, we have from Theorem 2.1, for every $n \geq 2$ and $k \geq 1$,

$$\mathbb{P}\{\Theta_n^{\gamma} > k\} = \sum_{\ell=0}^k \binom{k}{\ell} b^\ell \left(1-b\right)^{k-\ell} \alpha Q^\ell \mathbb{1P}\{\Theta_{n-1}^{\gamma} > k-\ell\}.$$
(8)

Moreover, Corollary 3.5 applies, i.e., for every $k \ge 1$,

$$\mathbb{P}\{\Theta^{\gamma} > k\} = \frac{1}{1 - (1 - b)^{k}} \sum_{\ell=0}^{k-1} \binom{k}{\ell} (1 - b)^{\ell} b^{k-\ell} \mathbb{P}\{\Theta^{\gamma} > \ell\} \alpha Q^{k-\ell} \mathbb{1}.$$
(9)

5. Application to cluster-based distributed storage

A cluster-based distributed storage peer-to-peer system guarantees durable access to large scale applications such as file sharing, streaming, or video-on-demand. It is achievable by harnessing the very large storage space globally provided by the many unused or idle nodes connected to the network. A common approach to handle these nodes is by having nodes that are close to each other according to a given proximity metric to self-organize into clusters. Specifically, each object (e.q. data stream, file) is divided into k equal size fragments, and recoded into a potentially unlimited number of independent check blocks through a rateless-erasure coding (also called Fountain) scheme (e.g. Luby (2002)). Fundamental property of erasure coding is that one may recover an initial object by collecting k' distinct check blocks generated by different sources, with k' slightly greater than k. During the coding phase, each check block c_i is generated by (i) choosing a degree d_i from a particular degree distribution, (ii) randomly choosing d_i distinct input symbols (called neighbors of c_i) among the k input symbols, and (iii) combining the d_i neighbors into a check block c_i by performing a bitwise XOR operation. The key idea of the decoding process is to build the Tanner graph based on the set of received check blocks. Upon receipt of check blocks, the decoder (i) finds any check block c_i with degree equal to one, (ii) removes the edge between c_i and k_i in the Tanner graph, and *(iii)* executes a bitwise XOR operation between k_i and any remaining check block c_r that has k_i as neighbor, and remove the edge between c_r and k_i . These steps are repeated until all k input symbols are successfully recovered. To guarantee the success of the decoding, the degree distribution is designed so that as few as possible check blocks are needed to ensure minimum redundancy among them, and the average degree is as low as possible to reduce the average number of symbol operations to recover the original data. This amounts to generating check blocks so that in average no more than 1/4 of them are of degree one to start the decoding and to prevent a too high amount of redundancy among these check blocks, 1/2 of them are of degree 2 so that combined with degree 1 check blocks they allow to cover a large proportion of input blocks, and 1/8 of them are of degree 3 so that the decoding process is unlikely to get stuck. The repartition of the other check blocks classically shows a steep decline, i.e. $1/2^i$ of them are of degree *i*. These check blocks are disseminated to the nodes of the system so that all the nodes that receive degree 1 check blocks self-organize into a cluster, those that receive degree two check blocks self organize into another cluster, and so on. Nodes can freely join and leave a cluster. For scalability and reliability reasons the number of nodes in a cluster is constrained. When the cluster size undershoots m nodes, then new check blocks are generated so that new nodes will join the cluster. Similarly when it exceeds M then generation of check blocks is suspended. In this model, $\gamma(n) = (p_{1,n}, \ldots, p_{n,n})$, where $p_{\ell,n} = 1/2^{\ell}$, for $\ell = 1, \ldots, n-1$ and $p_{n,n} = 1/2^{n-1}$, represents the distribution that allows to choose, at each instant (that corresponds to a join or leave event), a given cluster. Θ_n^{γ} then represents the first instant at which a cluster reaches its minimal (m) or maximal (M) size.

We model the effects of join and leave events using a homogeneous discrete-time Markov chain denoted by $X = \{X_n, n \ge 0\}$, which represents the evolution of the number of nodes in a cluster. This Markov chain is depicted in Figure 5 in which q = 1 - p and $p \in (0, 1)$. The transition



Figure 1. Markov chain model of one cluster

probability p means that a new peer joins the cluster while the transition probability q means that a peer leaves the cluster. The transition from state m + 1 to the absorbing state expresses the fact that the cluster has reached its minimal size m and that the coding has to be activated. In the same way the transition from state M - 1 to the absorbing state means that the cluster has reached its maximal size M and that the coding has to be suspended. The initial distribution α that we consider is the unit row vector e_j where $j = \lfloor (M + m)/2 \rfloor$. This means that $X_0 = j$ with probability 1. The matrix Q which gives the transitions between the transient states of Xis thus a tri-diagonal matrix where non-zero entries are $Q_{i,i+1} = p$ and $Q_{i,i-1} = q = 1 - p$. As explained above, the probability mass function $\gamma(n)$ is a truncated geometric distribution with parameter b = 1/2, i.e. given by $p_{r,n} = 1/2^r$, for every $r = 1, \ldots, n-1$ and $p_{r,n} = 1/2^{n-1}$. The distribution of Θ_n is then given by (8) and its limiting behavior is given by (9).

For the numerical evaluations, we have chosen p = 1/2. With this value, we easily get $E(\Theta_1 | X_0 = i) = (i - m)(M - i)$. We have also chosen m = 4 and M = 16 which implies that the number of transient states is equal to 11, that $X_0 = 10$ and $E(\Theta_1) = 36$. By the way, it is easily checked that $\alpha(I - Q)^{-1} = (1, 2, 3, 4, 5, 6, 5, 4, 3, 2, 1)$. For an error tolerance fixed to $\varepsilon = 10^{-4}$, the obtained values for the truncation steps K, L and J used to compute respectively the distribution of Θ_n^{γ} , its expectation and the expectation of Θ^{γ} are K = 550 and L = J = 787. Note that, since $m_n = 1/2$ is independent of n, so are K and L and since $m_n = b = 1/2$, we have L = J.

Figure 2 shows the distributions of Θ_n^{γ} and Θ^{γ} for different values of n. It is worth noting that the limiting distribution is reached very quickly for small values of n.

This result shows that the application designer does not need to determine the system size n (for $n \ge 5$) in order to calibrate the periods at which he must generate or suspend the coding process. For instance, for every $n \ge 5$, we have $\mathbb{P}\{\Theta_n^{\gamma} > 100\} \in [0.0893, 0.0928)$. This proves the scalability of this approach. This argument is confirmed in Table 1 which shows the expected values of Θ_n^{γ} and Θ^{γ} for different values of n.

$E(\Theta_1^{\gamma})$	$E(\Theta_2^{\gamma})$	$E(\Theta_3^{\gamma})$	$E(\Theta_4^{\gamma})$	$E(\Theta_5^{\gamma})$	$E(\Theta_6^{\gamma})$	$E(\Theta_7^{\gamma})$	$E(\Theta_8^{\gamma})$	$E(\Theta^{\gamma})$
36	42.2046	47.4027	50.6644	52.0177	52.3487	52.3928	52.3960	52.3961



Figure 2. From bottom to the top: $\mathbb{P}\{\Theta_1^{\gamma} > k\}$, $\mathbb{P}\{\Theta_2^{\gamma} > k\}$, $\mathbb{P}\{\Theta_3^{\gamma} > k\}$, $\mathbb{P}\{\Theta_4^{\gamma} > k\}$, $\mathbb{P}\{\Theta_5^{\gamma} > k\}$, $\mathbb{P}\{\Theta^{\gamma} > k\}$ as functions of k.



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