THE COMPUTATION OF STATIONARY DISTRIBUTIONS OF MARKOV CHAINS THROUGH PERTURBATIONS¹

JEFFREY J. HUNTER² Department of Mathematics and Statistics, Massey University Palmerston North, New Zealand

ABSTRACT

An algorithmic procedure for the determination of the stationary distribution of a finite, m-state, irreducible Markov chain, that does not require the use of methods for solving systems of linear equations, is presented. The technique is based upon a succession of m, rank one, perturbations of the trivial doubly stochastic matrix whose known steady state vector is updated at each stage to yield the required stationary probability vector.

Key words: Finite state Markov chains, perturbations, stationary probability vector.

AMS(MOS) subject classification: 60J10.

1. INTRODUCTION

In recent years, widespread attention has been given to the computation of stationary distributions of Markov chains. A variety of methods have been suggested and implemented. Before considering an alternative way for finding such distributions, it is of interest to give a brief survey of the techniques that have been employed.

Paige, Styan and Wachter (1975) presented a comprehensive survey of eight different algorithms involving a variety of procedures including the use of generalized inverses, rank reduction, least squares and power methods. Their recommendation was a direct method that involved transforming the singular set of stationary equations into a non-singular system using a rank one modification followed by Gaussian elimination with row pivoting. A further study by

Printed in the U.S.A. © 1991 The Society of Applied Mathematics, Modeling and Simulation

¹Received: January 1990, Revised: September 1990.

²This research was supported in part by the Air Force Office of Scientific Research under contract F49620 85 C 0144 through the Center for Stochastic Processes, Department of Statistics, University of North Carolina, Chapel Hill, North Carolina, 27514, USA.

Harrod and Plemmons (1984) provided another direct approach based upon the LU factorization using Gaussian elimination without pivoting.

Iterative techniques and approximation methods have been surveyed by Koury, McAllister and Stewart (1984). When the transition matrix is large and exhibits a nearly completely decomposable structure it is shown that a method of "aggregation" can be combined with point and block iterative techniques to produce methods which converge rapidly to the stationary probability vector.

Sheskin (1985) presented a partitioning algorithm that used a matrix reduction routine that partitions the transition matrix to create a sequence of smaller order transition matrices followed by a vector enlargement routine that enables the components of the steady state vector to be determined sequentially. A related procedure was developed by Grassmann, Taksar and Heyman (1985) using the theory of regenerative processes. They derived relationships between the steady state probabilities which are then used to develop a numerical algorithm to find these probabilities. Both of these latter two techniques appear to be, in effect, modifications of Gaussian elimination.

More recently, Meyer (1987), has utilized the concept of "stochastic complementation" whereby an irreducible chain is uncoupled into smaller irreducible chains whose stationary distributions can be coupled back together to produce the stationary distribution of the original chain.

In this paper, an entirely new approach involving the analysis of perturbed Markov chains is considered. In Hunter (1986), techniques for updating the stationary distribution of a finite irreducible Markov chain, following a rank one perturbation of its transition matrix, were presented. In this current paper, these techniques are utilized, to construct a general procedure for determining the stationary distribution of any finite irreducible Markov chain. A significant feature of the proposed algorithm is that at no stage does a system of linear equations have to be solved and consequently there is no reliance upon computer subroutines for matrix inversion or, the more generally accepted method of solution, Gaussian elimination with or without pivoting.

The basic idea is very simple. Suppose the steady state probability vector $\underline{\pi}'$ of an *m*-state irreducible Markov chain with given transition matrix P is required. Let P_0 be the transition matrix of another irreducible, *m*-state, Markov chain with known stationary probability vector $\underline{\pi}'_0$. By replacing, successively, the elements of each row of P_0 with the corresponding row elements as specified by P and recomputing the stationary probability vector of the resultant perturbed transition matrix, the vector $\underline{\pi}'_0$ can be transformed, in *m* stages, to $\underline{\pi}'$, by a series of *m* updates. As the irreducibility of a Markov chain is governed by the location of the positive entries in its transition matrix, to ensure the irreducibility of each perturbed Markov chain it is sufficient to commence with P_0 containing positive elements placed at least in the same position as those in P.

Consider starting with the trivial doubly stochastic matrix P_0 with each element having the value 1/m, so that $P_0 = \underline{e} \underline{e}'/m$, where $\underline{e}' = (1, 1, ..., 1)$ is a vector of ones. As can be easily shown, $\underline{\pi}'_0 = \underline{e}'/m$.

For i = 1, 2, ..., m, let e_i be the i^{th} elementary (column) vector with a one in the i^{th} position and zeros elsewhere. Let $p_i' = e_i' P$ be the i^{th} row of P and let

$$P_i = P_{i-1} + \underbrace{e}_i \underbrace{b}_i', \tag{1.1}$$

where

$$b_{i}' = p_{i}' - e_{i}'/m.$$

Let π'_i be the stationary probability vector associated with the Markov chain with transition matrix P_i , and, since $P_m = \sum_{i=1}^{m} e_i p'_i = P$, π'_m is in fact the required vector π' .

2. GENERAL THEORY

The construction of the algorithm is based upon the following results.

Theorem 2.1. Let P_i be the transition matrix of a finite irreducible Markov chain with stationary probability vector π'_i .

- (a) $I P_i + t_i u_i'$ is non-singular if and only if $u_i' e_i \neq 0$ and $\pi_i' t_i \neq 0$.
- (b) Under the conditions of (a),

$$\pi_{i}^{\prime} = \alpha_{i}^{\prime} / \alpha_{i}^{\prime} e, \qquad (2.1)$$

where

$$\underline{\alpha}_{i}^{\prime} = \underline{u}_{i}^{\prime} \left[I - P_{i} + \underline{t}_{i} \underline{u}_{i}^{\prime} \right]^{-1}.$$

$$(2.2)$$

Proof. For (a) see Theorem 3.3 in Hunter (1982) and for (b) see Corollary 4.1.2 in Hunter (1982).

Theorem 2.2. If X is non-singular and $b' X^{-1}a \neq -1$, then

$$(X + \underline{a}, \underline{b}')^{-1} = X^{-1} - \frac{X^{-1} \underline{a}, \underline{b}' X^{-1}}{1 + \underline{b}' X^{-1} \underline{a}}.$$
(2.3)

Proof. This is the Sherman-Morrison formula. See Golub and Van Loan (1983), p. 3.

Suppose that, following the i^{th} perturbation, the stationary probability vector $\pi t_i'$ has been found for the Markov chain with transition matrix P_i , as given by (1.1), by using the procedure described by Theorem 2.1 (b) for suitable choices of t_i and u_i .

In Hunter (1986) it was shown that it is possible to find an expression for π'_{i+1} , associated with P_{i+1} , using the same procedure outlined in Theorem 2.1 (b), by choosing the t_{i+1} and u_{i+1} in such a way that $[I - P_{i+1} + t_{i+1} u'_{i+1}]^{-1}$ can be determined from the earlier deduced expression $[I - P_i + t_i u'_i]^{-1}$, without performing an additional matrix inversion.

For the particular situation under consideration, for i = 0, 1, ..., m-1, if $\underbrace{t}_{i+1} = \underbrace{e}_{i+1}$ and $\underbrace{u}_{i+1} = \underbrace{u}_i + \underbrace{b}_{i+1}$, where \underbrace{b}_i is given by (1.2), then, from (1.1),

$$I - P_{i+1} + \underbrace{t}_{i+1} \underbrace{u}_{i+1}'_{i+1} = I - P_i + \underbrace{e}_{i+1} \underbrace{u}_{i}'_{i},$$

$$= I - P_i + \underbrace{t}_{i} \underbrace{u}_{i}'_{i} + (\underbrace{t}_{i+1} - \underbrace{t}_{i}) \underbrace{u}_{i}'_{i}.$$
 (2.4)

Now if $[I - P_i + t_i u'_i]^{-1}$ exists, from the proof of Theorem 3.3 in Hunter (1982),

$$\underline{u}_{i}^{\prime}[I - P_{i} + \underline{t}_{i} \underline{u}_{i}^{\prime}]^{-1} = \frac{\underline{\pi}_{i}^{\prime}}{\underline{\pi}_{i}^{\prime} \underline{t}_{i}}.$$
(2.5)

Thus, using (2.3), (2.4) and (2.5), if $A_i \equiv [I - P_i + t_i u_i']^{-1}$ exists,

$$A_{i+1} = A_{i}[I + (\underbrace{t}_{i} - \underbrace{t}_{i+1}) \frac{\pi_{i}'}{\pi_{i}' \underbrace{t}_{i+1}}].$$
(2.6)

Equation (2.6) is ideally suited for recursive operations once an initial inverse $A_0 = [I - P_0 - t_0 u_0']^{-1}$ has been determined. However, because of the form of P_0 that has been selected, if $t_0 = e$ and $u_0' = e'/m$, no matrix inverse has to be computed since, in this instance,

$$I - P_0 + \underbrace{t}_{0} \underbrace{u}_{0}' = I - \underbrace{e}_{e} \underbrace{e}'/m + \underbrace{e}_{e} \underbrace{e}'/m = I.$$

Furthermore, using (2.2),

$$\underline{\alpha}_{0}' = \underline{u}_{0}' = \underline{e}'/m,$$

and, from (2.1),

 $\mathfrak{\underline{\pi}}_{0}^{\prime} = \mathfrak{\underline{e}}^{\prime}/m.$

The basic algorithmic procedure now follows:

Let
$$\underline{t}_{0} = \underline{e}_{i}, \underline{u}_{0}' = \underline{e}_{i}'/m, A_{0} = I, \underline{\pi}_{0}' = \underline{e}_{i}'/m.$$

For $i = 1, 2, ..., m$, let $\underline{t}_{i} = \underline{e}_{i}$, and $\underline{u}_{i}' = \underline{u}_{i-1}' + \underline{p}_{i}' - \underline{e}_{i}'/m.$

. .

Compute
$$A_i = A_{i-1}[I + (\underbrace{t}_{i-1} - \underbrace{t}_{i})\underbrace{\pi}'_{i-1}/\underbrace{\pi}'_{i-1}\underbrace{t}_{i}].$$
 (2.7)

Compute

$$\alpha_i' = u_i' A_i. \tag{2.8}$$

Compute

$$\pi_{i}^{\prime} = \alpha_{i}^{\prime} / \alpha_{i}^{\prime} \epsilon .$$
(2.9)

Then $\pi' = \pi'_m$ is the stationary probability vector of the Markov chain with transition matrix $P = [p_{ij}].$

Since the elements of any stationary probability vector are always positive, $\pi i t_i$ and $\pi'_{i-1} t_i$ are both positive. Further, by induction, for $i = 0, 1, ..., m_i$

$$u_i' e_i = 1, \qquad (2.10)$$

so that the conditions of Theorem 2.1 and 2.2 are satisfied.

3. REFINEMENTS TO THE ALGORITHM

Although the procedure suggested in Section 2 will lead to the required stationary probability vector there are some modifications that, if employed, will lead to a more efficient procedure.

3.1. MODIFICATION TO THE π_i COMPUTATION

The ultimate aim of the algorithm is to determine $\pi = \pi m'$. Unless the stationary distributions of the intermediate perturbed Markov chains are required, some simplification can be effected by observing that (2.7) requires π'_{i-1} through its scaled version π'_{i-1}/π'_{i-1} . The scaling suggested by (2.9) is not required until the final step when i = m.

Thus, for i = 0, 1, ..., m - 1, let

$$\psi'_{i} \equiv \pi'_{i} / \pi'_{i} t_{i+1}. \tag{3.1}$$

Then, $\underbrace{\nu}_{0} = \underbrace{e}_{i}$ and for i = 1, 2, ..., m - 1,

$$\underline{v}_{i}^{\prime} = \underline{u}_{i}^{\prime} A_{i} / \underline{\omega}_{i}^{\prime} A_{i} \underbrace{e}_{i+1}.$$
(3.2)

If π'_i is required then it can be recovered simply. For i = 0, 1, ..., m-1,

$$\pi _{i}^{\prime} = \psi _{i}^{\prime} / \psi _{i}^{\prime} e . \qquad (3.3)$$

At the final step, compute $\pi m'$ using (2.8) and (2.9).

3.2. MODIFICATION TO THE A_i COMPUTATION

With the notation introduced in Section 3.1, it is easily seen that the early terms in the $\{A_i\}$ sequence are given, after simplification, as

$$A_0 = I,$$

$$A_1 = A_0 + (\underline{e} - \underline{e}_1) \underline{v}_0',$$
(3.4)

$$A_2 = A_1 + (\underline{e}_1 - \underline{e}_2) \underline{v}_1', \tag{3.5}$$

$$A_{3} = A_{2} + [(1 - v_{13}) \underbrace{e}_{1} + v_{13} \underbrace{e}_{2} - \underbrace{e}_{3}] \underbrace{v}_{2}', \qquad (3.6)$$

$$A_{4} = A_{3} + [\{1 - v_{14} - (1 - v_{13})v_{24}\} \underbrace{e}_{1} + (v_{14} - v_{13}v_{24}) \underbrace{e}_{2} + v_{24} \underbrace{e}_{3} - \underbrace{e}_{4}] \underbrace{v}_{3}', \qquad (3.7)$$

where $v_{ij} \equiv v_{ij}' e_{j}$, so that $v_{i,i+1} = 1$ for i = 1, 2, ..., m-1.

The above results provide motivation for the following theorem.

Theorem 3.1. For n = 0, 1, ..., m - 1,

$$A_{n+1} = A_n + B_n, \tag{3.8}$$

where

$$B_n = A_n(\underbrace{e}_n - \underbrace{e}_{n+1}) \underbrace{v}_n' \equiv \underbrace{b}_n \underbrace{v}_n'$$
(3.9)

with $\underset{c}{e}_{0} \equiv \underset{c}{e}$, so that $\underset{c}{b}_{0} = \underset{c}{e}_{-\underset{c}{e}_{1}}$ and for n = 1, 2, ..., m - 1,

$$b_{n} = b_{1n} e_{1} + \dots + b_{nn} e_{n} - e_{n+1}.$$
(3.10)

Proof. The theorem is obviously true by inspection, from (3.4) to (3.7), for n = 0, 1, 2, 3. Assume that (3.9) and (3.10) hold for n = 0, 1, ...k so that

$$A_{k+1} = A_k + B_k = A_{k-1} + B_{k-1} + B_k,$$

= ... = $I + \sum_{n=1}^{k} B_n.$ (3.11)

Hence

$$B_{k+1} = A_{k+1} (\underbrace{e}_{k+1} - \underbrace{e}_{k+2}) \underbrace{v}'_{k+1},$$

= $(I + B_0 + \sum_{n=1}^{k} B_n) (\underbrace{e}_{k+1} - \underbrace{e}_{k+2}) \underbrace{v}'_{k+1},$

implying that

$$b_{k+1} = [I + (e_{k-e_{1}})e_{k}' + \sum_{n=1}^{k} (\sum_{m=1}^{n} b_{mn}e_{m} - e_{n+1})v_{n}'](e_{k+1} - e_{k+2}).$$

Since $\mathfrak{E}'(\mathfrak{E}_{k+1}-\mathfrak{E}_{k+2})=0,$

$$\begin{split} b_{k+1} &= e_{k+1} - e_{k+2} + \sum_{n=1}^{k} \left(\sum_{m=1}^{n} b_{mn} e_{m} - e_{n+1} \right) (v_{n,k+1} - v_{n,k+2}), \\ &= e_{k+1} - e_{k+2} + \sum_{m=1}^{k} \left\{ \sum_{n=m}^{k} b_{mn} (v_{n,k+1} - v_{n,k+2}) \right\} e_{m} \\ &- \sum_{m=2}^{k+1} (v_{m-1,k+1} - v_{m-1,k+2}) e_{m}, \end{split}$$

showing that \underbrace{b}_{k+1} is linear combination of $\underbrace{e}_{1}, \ldots, \underbrace{e}_{k+2}$ with \underbrace{e}_{k+2} having coefficient -1. Thus (3.10) is true for n = k+1 and the theorem follows by induction.

Observe that

$$b_{k+1} = \sum_{m=1}^{k+1} b_{m,k+1} e_m - e_{k+2}, \qquad (3.12)$$

where

$$b_{1,k+1} = \sum_{n=1}^{k} b_{1,n} (v_{n,k+1} - v_{n,k+2}), \qquad (3.13)$$

$$b_{k+1,k+1} = v_{k,k+2}, (\text{since } v_{k,k+1} = 1), \tag{3.14}$$

and for $m = 2, \ldots, k$,

$$b_{m,k+1} = \sum_{n=m-1}^{k} b_{m,n} (v_{n,k+1} - v_{n,k+1}).$$
(3.15)

Note that

Thus, in the matrix B_n , all the entries in rows number n + 2, ..., m are zero. Obviously, this has considerable significance in the calculation of the matrices $A_i (i = 2, ..., m)$ as required in the algorithm.

The updating process, given by (2.7), can be replaced for i = 2, ..., m, by

$$A_i = A_{i-1} + B_{i-1}, \tag{3.17}$$

where

$$B_{i-1} - A_{i-1}(\underbrace{e}_{i-1} - \underbrace{e}_{i}) \underbrace{v}'_{i-1}, \qquad (3.18)$$

is such that only the first i rows of B_{i-1} are computed with the remaining entries set equal to zero.

Furthermore, from (3.16), some of the rows of B_n have a special form and do not require computation. In particular, the (n+1)th row is simply $-\underline{v}'_n$ while, from (3.14), the n^{th} row is $v_{n-1,n+1}$ times \underline{v}'_n .

Note also that the (n+1)st column of B_n is $\underset{n}{\flat}_n$ since $v_{n,n+1} = 1$.

There are also some other checks that can be applied.

Theorem 3.2. For i = 1, 2, ..., m,

$$A_{i\underline{e}}_{i} = \underline{e}, \qquad (3.19)$$

$$\frac{1}{m} \underbrace{e}' A_i = \underbrace{e}, \qquad (3.20)$$

$$\underline{e}'B_i = \underline{0}'. \tag{3.21}$$

Proof. Since $A_i = [I - P_i + t_i u_i]^{-1}$, equation (3.17) of Hunter (1982) implies that

$$A_{i, \overset{t}{\sim} i} = \frac{\underbrace{e}}{\underbrace{u'_{i, \overset{t}{\sim}}}} = \underbrace{e}_{\overset{t}{\sim}},$$

yielding (3.19) with $t_i = e_i$.

Equation (3.20) is obviously true when i = 1 since

$$\mathfrak{L}'A_1 = \mathfrak{L}'[I + (\mathfrak{L} - \mathfrak{L}_1)\mathfrak{L}'],$$
$$= \mathfrak{L}' + (m-1)\mathfrak{L}',$$
$$= m\mathfrak{L}'.$$

Thus, by induction, if (3.20) is true for $i = 1, 2, \dots, k$, from (2.7) and (3.1),

$$\frac{1}{m} \underbrace{\varepsilon}' A_{k+1} = \frac{1}{m} \underbrace{\varepsilon} A_k [I + (\underbrace{\varepsilon}_k - \underbrace{\varepsilon}_{k+1}) \underbrace{\upsilon}'_k],$$
$$= \underbrace{\varepsilon}' [I + (\underbrace{\varepsilon}_k - \underbrace{\varepsilon}_{k+1}) \underbrace{\upsilon}'_k],$$
$$= \underbrace{\varepsilon}' + (\underbrace{\varepsilon}' \underbrace{\varepsilon}_k - \underbrace{\varepsilon}' \underbrace{\varepsilon}_{k+1}) \underbrace{\upsilon}'_k = \underbrace{\varepsilon}'.$$

Thus (3.20) and hence also (3.21) follow.

A consequence of Theorem 3.2 is that the i^{th} column of A_i consists solely of unit elements while the sum of elements of each column of B_i is zero.

3.3. MODIFICATION TO THE α_i COMPUTATION

Although the $\{A_i\}(i=0,1,...,m)$ sequence plays an integral role in the procedure, the matrices A_i are required only to obtain the sequence of vectors $\underline{\alpha}'_i = \underline{u}'_i A_i$ and hence the vectors $\underline{\nu}'_i$. Thus is worth examining whether it is possible to dispense with explicit calculation of the A_i by deriving the $\{\underline{\alpha}'_i\}(i=1,2,...,m)$ sequence recursively.

Theorem 3.3. For i = 0, 1, 2, ..., m - 1,

$$\alpha'_{i+1} = \nu'_{i} - \epsilon' + p'_{i+1} A_{i+1}.$$
(3.22)

Proof. First observe that, from (2.8),

$$\underline{\alpha}_1' = \underline{u}_1' A_1 = \underline{p}_1' A_1,$$

so that (3.22) holds for i = 0 since $v_0' = e'$.

In general, for i = 1, 2, ..., m - 1, from (2.8),

$$\alpha'_{i+1} = \alpha'_{i+1}A_{i+1},$$

= $[\alpha'_{i} - (\alpha'/m) + \alpha'_{i+1}]A_{i+1}.$ (3.23)

Now

where $\alpha_{ij} = \alpha_{ij} \stackrel{\prime}{=} e_{j}$. But, from (3.19) and (2.10),

$$\alpha_{ii} = \underline{u}'_{i}A_{i}\underline{e}_{i} = \underline{u}'_{i}\underline{e}_{i} = 1, \qquad (3.25)$$

and, since from (3.2) and (2.8), $v_i' = \alpha_i'/\alpha_{i,i+1}$, (3.24) becomes

$$\underline{u}_{i}^{\prime}A_{i+1} = \underline{v}_{i}^{\prime}.$$
(3.26)

Equation (3.22) now follows from (3.23) upon substitution of (3.26) and (3.20).

Theorem 3.3 shows that in updating from α'_i to α'_{i+1} the term $p'_{i+1}A_{i+1}$ must be computed. The calculations of $\alpha'_1, \alpha'_2, \dots, \alpha'_i$ require, successively, p'_1, \dots, p'_i and for α'_{i+1} this is the first time p'_{i+1} , the (i+1)th row of P is involved.

Although, $p'_{i+1}A_{i+1}$ can be expressed in terms of A_i , as can be seen from the next theorem, very little advantage is gained since such terms are required for each i = 0, 1, ..., m-1.

Theorem 3.4. For i = 0, 1, ..., m - 1,

$$\underline{p}'_{i+1}A_{i+1} = \underline{p}'_{i+1}A_i + \underline{v}'_i - (\underline{p}'_{i+1}A_i \underline{e}_{i+1}) \underline{v}'_i.$$
(3.27)

Proof. For i = 0,

$$p_1'A_1 = p_1'[I + (\underline{e} - \underline{e}_1)\underline{e}'],$$
$$= p_1' + (p_1'\underline{e})\underline{e}' - (p_1'\underline{e}_1)\underline{e}',$$

and the result follows, since $\underset{l}{\mathcal{D}}'_{l} \underset{l}{\mathcal{C}} = 1, \underset{0}{\mathcal{U}}'_{0} = \underset{l}{\mathcal{E}}'$ and $A_{0} = I$.

In general, for i = 1, 2, ..., m - 1, from (2.7) and (3.1),

$$\underline{p}'_{i+1}A_{i+1} = \underline{p}'_{i+1}[A_i + A_i(\underline{e}_i - \underline{e}_{i+1})\underline{v}'_i].$$

Equation (3.27) follows since, from (3.19),

$$\underbrace{p'_{i+1}A_i \underbrace{e}_i = p'_{i+1} \underbrace{e}_i = 1.}$$

As a consequence of Theorem 3.3, it is suggested that (2.8) in the algorithm be replace by (3.22).

3.4. MODIFICATION TO THE $\pi m' = \pi'$ COMPUTATION

At the final step of the algorithm A_m can be computed and consequently $\pi m'$ derived as $\alpha m' / \alpha m' \in w$ where $\alpha m' = \mu m' A_m$. However, A_m need not be explicitly determined since, from Theorem 3.3 and 3.4,

$$a_{m}' = v_{m-1}' - e_{m}' + p_{m}'A_{m},$$

where $p_m'A_m = p_m'A_{m-1} + v_{m-1}' - (p_m'A_{m-1} e_m) v_{m-1}'$.

4. RECOMMENDED PROCEDURE

As a consequence of the refinements discussed in Section 3, it is suggested that the algorithm be constructed as follows:

2) Let $\alpha'_{1} = p'_{1}A_{1}$. 3) For i = 1, 2, ..., m - 2, compute (a) $\psi'_{i} = \alpha'_{i}/\alpha'_{i}e_{i+1}$, (b) $B_{i} = A_{i}(e_{i} - e_{i+1})\psi'_{i}$, (c) $A_{i+1} = A_{i} + B_{i}$, (d) $\alpha'_{i+1} = \psi'_{i} - e' + p'_{i+1}A_{i+1}$. 4) Let $\psi'_{m-1} = \alpha'_{m-1}/\alpha'_{m-1}e_{m}$.

Let $A_1 = I + (e_1 - e_{-1}) e_1'$.

- 5) Let $\alpha'_m = 2 \upsilon'_{m-1} \varepsilon' + p mA_{m-1} (p mA_{m-1} \varepsilon_m) \upsilon'_{m-1}$.
- 6) Let $\pi' = \pi'_m = \alpha'_m / \alpha'_m \varepsilon$.

The order of the number of arithmetic operations (multiplication and division) required to determine π ' can be estimated as follows. The computation of the B_i and the π 'A_i have a dominant effect on the number of operations required. Since $A_i(\underline{e}_i - \underline{e}_{i+1})$ is effectively the difference of two columns of A_i , only mi operations are required to determine B_i , taking into consideration that B_i has only (i+1) non-zero rows, and, as a consequence of (3.21), that the elements of one row can be found from the other rows using the fact that each column sums to zero. On the other hand, for a general transition matrix, $\underline{p}'_i A_i$ will require m^2 operations, although this can be reduced to m(m-1) since the *i*th element of this row vector, $\underline{p}'_i A_i \underline{e}_i = \underline{p}'_i \underline{e} = 1$, (by using (3.19)). Since the other calculations required are relatively insignificant in comparison, the total number of operations is of the order of

$$\sum_{i=1}^{m-1} mi + \sum_{i=1}^{m} m(m-1) = 3m^2(m-1)/2, \text{ i.e. of order } 3m^3/2.$$

1)

To solve for the stationary distribution directly using Gaussian elimination requires of the order fo $4m^3/3$, while to solve directly using a matrix inversion routine requires of the order of $2m^3$ operations, (see Isaacson and Keller (1966)).

The procedure is, in effect, finding the stationary distribution of m different irreducible Markov chains and consequently the routine that has been developed offers much more information than other techniques currently available.

Although it has been suggested that the algorithm proceed row by row, there is no necessity to adhere to a strict sequential ordering of the rows. The procedure as outlined by (2.7), (2.8) and (2.9) can easily be adapted to such changes by altering the \underline{t}_i and \underline{u}'_i . A consequence of this is that the effect of changing selected transition probabilities upon the stationary distribution can easily be determined. (See also Hunter (1986)).

The procedure also offers the opportunity to utilize the structure of special transition matrices. For example, if the transition matrix of the chain is banded with $p_{ij} = 0$ for j < i - g and j > i + h, which occurs in some queueing models, the calculation of $p_i A_i$ will require at most (g+h)m operations and the algorithm will require on the order of only $m^3/2 + (g+h)m^2$ operations.

5. STRUCTURAL RESULTS

In Section 3.2 expressions for the first few terms of the $\{A_i\}$ sequence were derived. By using those terms and working through the first few steps of the algorithm it can be shown, that, following simplification, for i = 1, 2, 3,

$$\underline{\alpha}_{i}^{\prime} = (\mu_{i0} \underline{e}^{\prime} + \mu_{i1} \underline{p}_{1}^{\prime} + \dots + \mu_{ii} \underline{p}_{i}^{\prime}) / \mu_{ii}, \qquad (5.1)$$

$$\pi_{i}' = (\mu_{i0} \varepsilon' + \mu_{i1} p_{1}' + \dots + \mu_{ii} p_{i}') / (m\mu_{i0} + \mu_{i1} + \dots + \mu_{ii}),$$
(5.3)

where $\mu_{10} = 1 - p_{11}$,

$$\begin{split} \mu_{11} &= 1, \\ \mu_{20} &= (1-p_{11})(1-p_{22}) - p_{12}p_{21}, \\ \mu_{21} &= 1+p_{21}-p_{22}, \\ \mu_{22} &= 1-p_{11}+p_{12}, \end{split}$$

$$\begin{split} \mu_{30} &= (1-p_{11})(1-p_{22})(1-p_{33}) - p_{12}p_{23}p_{31} - p_{13}p_{21}p_{32} \\ &\quad -p_{13}(1-p_{22})p_{31} - (1-p_{11})p_{23}p_{32} + p_{12}p_{21}(1-p_{33}), \\ \mu_{31} &= p_{21}(1-p_{33}+p_{32}) + (1-p_{22})(1-p_{33}+p_{31}) + p_{23}(p_{31}-p_{32}), \\ \mu_{32} &= p_{31}(p_{12}-p_{13}) + p_{32}(1-p_{11}+p_{13}) + (1-p_{33})(1-p_{11}+p_{12}), \\ \mu_{33} &= (1-p_{11})(1-p_{22}+p_{23}) + p_{12}(p_{23}-p_{21}) + p_{13}(1-p_{22}+p_{21}). \end{split}$$

The general structure exhibited by (5.1), and hence also by (5.2) and (5.3), holds for all i = 1, 2, ..., m. [A proof by induction shows that if (5.1) and (5.2) hold for i = 1, ..., n then, since

$$p'_{n+1}A_{n+1} = p'_{n+1}[I + B_0 + \dots + B_n].$$

= $p'_{n+1}[I + b_0 v'_0 + \dots + b_n v'_n],$
= $p'_{n+1} + (p'_{n+1}b_0) v'_0 + \dots + (p'_{n+1}b_n) v'_n,$

using (3.22) with $i = n, \alpha'_{n+1}$ is a linear combination of $\nu'_0, \dots, \nu'_n, p'_{n+1}$, i.e. of $\varepsilon', p'_1, \dots, p'_{n+1}$. Furthermore, the coefficient of p'_{n+1} is unity whereby establishing the general structure of α'_{n+1} .]

Note also that, from (5.1) and (3.25), $\alpha'_i \stackrel{e}{\sim}_i = 1$, for i = 1, ..., m and thus

$$\mu_{i0} + \mu_{i1} p_{1i} + \ldots + \mu_{ii} p_{ii} = \mu_{ii}.$$
(5.4)

Further, for i = 1, 2 it can be shown, by direct verification, that

$$\mu_{i0} + \mu_{i1}p_{i,i+1} + \dots + \mu_{ii}p_{i,i+1} = \mu_{i+1,i+1},$$
(5.5)

which implies that

$$v_{ii} = v_{i}' e_{i} = \mu_{ii} / \mu_{i+1,i+1}, \qquad (5.6)$$

and

$$\alpha_{i,i+1} = \alpha_{i,i+1}' e_{i+1,i+1} / \mu_{ii}, \qquad (5.7)$$

results that it has not been possible to establish in general.

Let $(I - P)_i$ be the leading *ith* order principal submatrix of I - P formed by deleting all but the first *i* rows and columns then, for i = 1, 2, 3,

$$\mu_{i0} = det(I - P)_i. \tag{5.8}$$

For the special case when m = 3, with the notation used earlier in this section, it can be verified that

$$\underline{\alpha}_{i}^{\prime} = (\mu_{i1}, \mu_{i2}, \mu_{i3})/\mu_{ii}, \quad (i = 1, 2, 3), \tag{5.9}$$

$$\pi_{i}' = (\mu_{i1}, \mu_{i2}, \mu_{i3})/(\mu_{i1} + \mu_{i2} + \mu_{i3}), \quad (i = 1, 2, 3),$$
(5.11)

where

$$\mu_1 = 1 - p_{11} + p_{12} = \mu_{22},$$
$$\mu_{13} = 1 - p_{11} + p_{13},$$
$$\mu_{23} = \mu_{33} = 3\mu_{30}.$$

Observe that π_1' , π_2' and π_3' give, respectively, the stationary probability vectors of the Markov chains whose transition matrices are

p ₁₁	$p_{12}^{}$	p ₁₃			<i>p</i> ₁₂	p ₁₃			<i>p</i> ₁₁	<i>p</i> ₁₂	p ₁₃
1/3	1/3	1/3	,	p_{21}	p ₂₂	p_{23}	,	and	<i>p</i> ₂₁	p ₂₂	p ₂₃
1/3	1/3	1/3		1/3	1/3	1/3			p ₃₁	p ₃₂	p ₃₃

In examining (5.11), with i = 3, it can be shown that $\mu_{3j} = 3D_j$, (j = 1, 2, 3), where D_j is the determinant formed by striking out the *jth* row and *jth* column of I - P. This leads to an expression for the stationary probability vector of a general irreducible, three state, Markov chain as

$$\underline{\pi}' = (D_1, D_2, D_3) / \sum_{j=1}^3 D_j.$$
(5.12)

The natural extension of (5.12) for a general finite irreducible Markov chain is also true, such a result being attributed to Mihoc by Fréchet (1950) and rediscovered by Singer (1964).

Although the full details of a proof of the generalization of (5.12) using the techniques of this paper have not been worked out, it is conjectured that for an *m*-state chain $\mu_{mj} = mD_j$, (a result that holds for m = 2, 3), so that the procedures proposed in this paper appear to lead to an effective algorithmic construction of Mihoc's technique.

6. EXAMPLE

As an illustration of the perturbation procedure, we compute the stationary probabilities for the irreducible five-state Markov chain example used in Sheskin (1985) which was taken from Kemeny and Snell (1960) p. 199. The (1,1) entry has been changed to ensure that P is in fact a stochastic matrix, so that P is given by

$$P = \begin{bmatrix} 0.831 & 0.033 & 0.013 & 0.028 & 0.095 \\ 0.046 & 0.788 & 0.016 & 0.038 & 0.112 \\ 0.038 & 0.034 & 0.785 & 0.036 & 0.107 \\ 0.054 & 0.045 & 0.017 & 0.728 & 0.156 \\ 0.082 & 0.065 & 0.023 & 0.071 & 0.759 \end{bmatrix}$$

The following calculations were performed on a MacIntosh computer using the Microsoft Excel software spreadsheet. Although the program yields expressions to 14 significant figures, the results have been rounded to 3 decimal places.

•

α' ₁	=	(1.000,	0.202,	0.182,	0.197,	0.264),
Ľ 1′	=	(4.950,	1.000,	0.901,	0.975,	1.307),
≈′2		(1.277,	1.000,	0.202,	0.244,	0.403),
v 2′	=	(6.309,	4.939,	1.000,	1.203,	1.991),
α'_3	=	(1.593,	1.232,	1.000,	0.302,	0.571),
v′₃	=	(5.275,	4.079,	3.311,	1.000,	1.890),
¢ 4′	=	(1.717,	1.295,	0.935,	1.000,	0.705),
2'4	=	(2.434,	1.836,	1.326,	1.418,	1.000),
¢ ′₅	=	(0.841,	0.573,	0.237,	0.459,	1.000).

From the above calculations, the stationary probabilities for each successive perturbed

π_1'	=	(0.542,	0.109,	0.099,	0.107,	0.143),
π'_2	=	(0.409,	0.320,	0.065,	0.078,	0.129),
π'3	=	(0.339,	0.262,	0.213,	0.064,	0.122),
π'4	=	(0.304,	0.229,	0.165,	0.177,	0.125),
π'_{5}	=	(0.270,	0.184,	0.076,	0.148,	0.322).

transition matrix can easily be calculated, yielding the following results:

7. FINAL COMMENTS

The initial choice of P_0 as e e'/m ensures that it is possible to start with an irreducible Markov chain whose stationary distribution is easily found without having to compute a matrix inverse or to solve a general set of linear equations. The fact that every element of P_0 is non-zero leads to a sequence of matrices A_1, A_2, \ldots that are "dense". Is it possible to start with a different Markov chain, say one that is relatively sparse, whose stationary distribution is well known and such that, for the early recursions, the equivalent sequence A_1, A_2, \ldots retains such a sparsity property?

The periodic Markov chain with entries $p_{ii+1}^{(0)} = 1, (i = 1, 2, ..., m-1)$, and $p_{m1}^{(0)} = 1$ is a potential candidate for P_0 , whose stationary probability vector is also $\pi_0' = \underline{e}'/m$. Even if \underline{t}_0 and \underline{u}_0 can be specified so that $A_0 = [I - P_0 + \underline{t}_0 \underline{u}_0']^{-1}$ has a simple structure much care would be required in carrying out any sequential row modification with this P_0 . For example if, for the specified P transition matrix, $p_{12} = 0$ then state 2 is never reached in the Markov chain with transition matrix P_1 violating the required irreducibility property of P_1 .

The major advantage in choosing $P_0 = e e'/m$ is that the irreducibility of each P_i transition matrix is guaranteed at each step of the procedure.

REFERENCES

- FRECHLET, M., Recherches Théoriques Modernes sur le Calcul des Probabilités, Vol. 2, 2nd ed., Gauthier-Villars, Paris, 1950.
- [2] GOLUB, G.H., and VAN LOAN, C.F., Matrix Computations, The Johns Hopkins University Press, Bailtimore, 1983.

[3]	GRASSMAN, W.K., TAKSAR, M.I., and HEYMAN, D.P., Regenerative analysis and
	steady state distributions for Markov chains, Oper. Res. 33 (1985), pp. 1107-1116.

- [4] HARROD, W.J. and PLEMMONS, R.J., Comparison of some direct methods for computing stationary distributions of Markov chains, SIAM J. Sci. Statist. Comput., 5 (1984), pp. 453-469.
- [5] HUNTER, J.J., Generalized inverses and their application to applied probability problems, Linear Algebra Appl. 45 (1982), pp. 157-198.
- [6] HUNTER, J.J., Stationary distributions of perturbed Markov chains, Linear Algebra Appl. 82 (1986), pp. 201-214.
- [7] ISAACSON, E. and KELLER, H.B., Analysis of Numerical Methods, John Wiley and Sons, New York, 1966.
- [8] KEMENY, J.G. and SNELL, J.L., Finite Markov Chains, Van Nostrand, Princeton, 1960.
- [9] KOURY, J.R., MCALLISTER, D.F. and STEWART, W.J., Iterative methods for computing stationary distributions of nearly completely decomposable Markov chains, SIAM J. Alg. Disc. Meth., 5 (1984), pp. 164-186.
- [10] MEYER, Jr., C.D., Uncoupling Markov chains and the Simon-Ando theory of nearly reducible systems, Centre for Research in Scientific Computation, North Carolina State Univ., Tech. Rep. 10018701, 1987.
- [11] PAIGE, C.C., STYAN, G.P.H. and WACHTER, P.G., Computation of the stationary distribution of a Markov chain, J. of Statist. Comput. Simulation, 4 (1975), pp. 173-186.
- [12] SHESKIN, T.J., A Markov chain partitioning algorithm for computing steady state probabilities, Oper. Res. 33 (1985), pp. 228-235.
- [13] SINGER, A., The steady state probabilities of a Markov chain as a function of the transition probabilities, Oper. Res. 12 (1964), pp. 498-499.