# The Nsolve Program 

Peter Buchholz<br>Informatik IV，University of Dortmund<br>D－44227 Dortmund，Germany<br>Email：peter．buchholz＠udo．edu

## 1 Short Introduction to Nsolve

Nsolve is a program to determine the stationary solution of hierarchical Markovian models exploiting a hierarchical Kronecker structure of the generator matrix．The model class contains stochastic automata networks（SANs）without functional transition rates and superposed generalized stochastic Petri nets （SGSPNs）which are both described by a hierarchical model with one state in the upper level．A hier－ archical model consists of a number of components（LLMs）and one high level model（HLM）．Detailed descriptions of the required matrix structure and some solution methods can be found in $[4,7,5,2]$ ．The different file formats which are used to describe hierarchical models are explained in a companion report ［8］．

Nsolve includes iterative solution techniques using the hierarchical Kronecker representation of the matrix and it includes as well iterative techniques using a sparse matrix representation．In the latter case，the flat matrix is first generated out of the hierarchical Kronecker representation．Observe that the implementation of the generation is not very efficient such that solution techniques for sparse matrices should only be used for models with a modest state space size $\left(\leq 10^{5}\right)$ ．

Nsolve requires the following files：
－$\langle$ name $\rangle$ ．conf configuration file（Description in 2）

- 〈 name＞0．mat HLM matrix（Description in［8］）
- 〈 name $\rangle$ ．spa HLM state space（Description in［8］）
- 〈 name $\rangle j$ ．mat LLM $j$ matrix（Description in［8］）

A file for each LLM has to exist．The number of LLMs can be found in the configuration file．
All files may include comments．Comments start with \＃as the first character in a line，then the rest of the line is interpreted as a comment！

The program is called with

```
Nsolve name
```


## 2 Configuration File

The configuration file includes the following values（plus possibly some comments）：

```
Number of LLMs (integer \geq0)
Solution method (integer \geq0)
Maximum number of iterations (integer \geq0)
\epsilon
\epsilon
\epsilon
Maximum CPU-time (real }\geq0.0
Relaxation-parameter (real \geq0.0)
Output overall vector (Boolean)
Output HLM vector (Boolean)
for all LLMs 1,..., Number of LLMs
Output LLM vector (Boolean)
```

```
Initial state HLM (integer \geq-2)
if initial state HLM \geq0 then
for all LLMs 1,...Number of LLMs
Initial state LLM (integer \geq0)
```


### 2.1 Interpretation of the parameters

Values in the configuration file are interpreted as follows.

## Number of LLMs

The number of components which form the model.

## Solution method

The currently available solution methods are summarized in table 1. The number in the first column of the table has to be used in the configuration file.

Maximal number of iterations
The solution stops, if the number of iterations has been reached
$\epsilon_{1}$
The solution stops, if the absolute value if the largest element in the residual vector becomes smaller than $\epsilon_{1}$. The residual vector is currently checked every 10 iterations in most methods. Observe that in some methods like the projection methods TFQMR and BICGSTAB the residual can only be estimated during computation. Thus, these methods may stop too early or too late. In any case, the true residual is computed after termination of the method.
$\epsilon_{2}$
The value is used for several purposes. In A/D-methods, an $A / D$ step for the HLM is performed if the residual norm projected onto the HLM state space exceeds $\epsilon_{2}$. For methods with ILUTH preconditioner, the value is interpreted as the threshold for keeping non-zero elements during the incomplete factorization.
$\epsilon_{3}$
The value is only used for A/D-methods where an A/D step for a LLM is performed, if the residual norm projected onto the LLM state exceeds $\epsilon_{3}$.

CPU time limit
The method stops, if the CPU time of the solution (excluding the time for matrix generation) exceeds the limit.

## Relaxation-parameter

Relaxation-parameter for the methods JOR, SJOR, SOR and BSOR. The value should be taken from $(0,2)$.

## Output parameters for solution vectors

Parameter values can be 0 ( $=$ no output) or 1 ( $=$ output). The first parameter is for the complete solution vector, the second for the solution vector of the HLM and then follow the values for all LLMs. Solution vectors are written in files. File name_overall.vec contains the complete vector, name0.vec HLM vector and name $j$.vec the vector for LLM $j$.

The initial distribution
Different methods exist to define the initial distribution. The first value determines the type of the initial distribution. If the first value is larger or equal to zero, then it is assumed that a single initial state is defined. In this case the following values include the initial states of the LLMs. Observe that not every combination of states defines a correct state. If inconsistencies are detected, the program tries to repair the state and prints a warning. The distribution is initialized with 1.0 at the initial state and 0.0 elsewhere. Usually a uniform distribution is recommended as initial distribution because in a hierarchical model all states are reachable if the state space is generated appropriately (see [5]). A uniform distribution is generated if the first state value is set to -2 . In this case the following values are not interpreted. The last possibility to define the initial distribution is to read it from a file. In this case the value in the configuration file has to be set to -1 . The name of the file has to be name_overall.reach. The vector length has to be equal to the number of states in the state space of the hierarchical model. The vector is normalized to 1 before the iteration starts. With this feature result vectors of one method can be used as initial vectors of another method.

Apart from the parameters in the configuration file, some parameters exist in the source files of the different methods. However, modifications of these parameters requires a modification of the source code and recompilation of the program. We tried to set the parameters to reasonable values, but the choices might not be optimal for specific examples.

The methods NO _SOLUTION ( 0 ) and MAT_GEN (1) do not analyze the system. With NO_SOLUTION only the matrix structures are built to prove consistency of the model. MAT_GEN generates a flat sparse matrix which is written in the file name_overall.mat. The file format for the flat matrix can be found in [9].

## References

[1] P. Buchholz. Numerical solution methods based on structured descriptions of Markovian models. In G. Balbo and G. Serazzi, editors, Computer Performance Evaluation - Modelling Techniques and Tools, pages 251-267. Elsevier, 1992.
[2] P. Buchholz. A framework for the hierarchical analysis of discrete event dynamic systems. Habilitationsschrift, Fachbereich Informatik, Universität Dortmund, 1996. available upon request.
[3] P. Buchholz. An aggregation/disaggregation algorithm for stochastic automata networks. Probability in the Engineering and Informational Sciences, 11(2):229-253, 1997.
[4] P. Buchholz. An adaptive aggregation/disaggregation algorithm for hierarchical Markovian models. European Journal of Operational Research, 116(3):85-104, 1999.
[5] P. Buchholz. Hierarchical structuring of superposed GSPNs. IEEE Transactions on Software Engineering, 25(2):166-181, 1999.
[6] P. Buchholz. Projection methods for the analysis of stochastic automata networks. In B. Plateau, W. J. Stewart, and M. Silva, editors, Numerical Solution of Markov Chains (NSMC'99), pages 149-168. Prensas Universitarias de Zaragoza, 1999.
[7] P. Buchholz. Structured analysis approaches for large Markov chains. Applied Numerical Mathematics, 31(4):375-404, 1999.
[8] P. Buchholz. A matrix format for structured markovian models. working paper, 102010.
[9] P. Buchholz. The sparse matrix file format. working paper, 102010.
[10] P. Buchholz and T. Dayar. Block SOR for Kronecker structured representations. Linear Algebra and its Applications, 386:83-109, 2004.
[11] P. Buchholz and T. Dayar. Block SOR preconditioned projection methods for Kronecker structured Markovian representations. SIAM Journal on Scientific Computing, 26:1289-1313, 2005.
[12] P. Buchholz, J. Dunkel, B. Müller-Clostermann, M. Sczittnick, and S. Zäske. Quantitative Systemanalyse mit Markovschen Ketten. Teubner-Texte zur Informatik Band 8. Teubner, 1994.
[13] R. W. Freund and M. Hochbruck. On the use of two QMR for solving singular systems and applications in Markov chain modelling. Numerical Linear Algebra and Applications, 1:403-420, 1994.
[14] G. Horton and S. Leutenegger. A multi-level solution algorithm for steady state Markov-chains. ACM Performance Evaluation Review, 22:191-200, 1994.
[15] Y. Saad and K. Wu. DQGMRES: a direct quasi-minimal rsidual algorithm based on incomplete orthogonalization. Numerical Linear Algabra and Applications, 3:329-343, 1996.
[16] W. J. Stewart. Introduction to the numerical solution of Markov chains. Princeton University Press, 1994.
[17] E. Uysal and T. Dayar. Iterative methods based on splittings for stochastic automata networks. European Journal of Operational Research, 110(1):166-186, 1998.
[18] H. A. van der Vorst. Bi-CGSTAB: A fast and smoothly converging variant of Bi-CG for the solution of nonsymmetric linear systems. SIAM Journal on Scientific Computing, 13:631-644, 1992.

| Nr. | Method | Explanation |
| :---: | :---: | :---: |
| Matrix structure test |  |  |
| 0 | NO SOLUTION | Matrix generation without solution |
| 1 | MAT_GEN | Flat matrix generation |
| Solution techniques for sparse matrices |  |  |
| 2 | LU | LU-Factorization ([16] pp. 66ff, [12] pp. 30ff) |
| 11 | POWER | Power-Method ([16] pp. 121ff) |
| 12 | JOR | JOR ([16] pp. 127f) |
| 13 | SOR | SOR ([16] pp. 128f) |
| 21 | POWER A/D | Power-Method with Aggregation/Disaggregation ([12], pp. 51ff) |
| 22 | JOR A/D | JOR with Aggregation/Disaggregation ([12], pp. 51ff) |
| 23 | SOR A/D | SOR with Aggregation/Disaggregation ([12], pp. 51ff) |
| 24 | KMS | KMS ([16], pp. 308) |
| 25 | ML | Multi-level method ([14]) |
| 31 | GMRES | GMRES ([16], pp. 198ff) |
| 32 | BICGSTAB | BiCGSTAB ([18])) |
| 33 | TFQMR | Transpose free variant of QMR ([13])) |
| 61 | POWER ILU0 | Power-Method with ILU0 Preconditioning ([16], pp. 145) |
| 62 | POWER ILUTH | Power-Method with ILUTH Preconditioning ([16], pp. 145) |
| 63 | GMRES ILU0 | GMRES with ILU0 Preconditioning ([16],pp. 205ff, 145) |
| 64 | GMRES ILUTH | GMRES with ILUTH Preconditioning ([16],pp. 205ff, 146) |
| 65 | BICGSTAB ILU0 | BICGSTAB with ILU0 Preconditioning |
| 66 | BICGSTAB ILUTH | BICGSTAB with ILUTH Preconditioning |
| 67 | TFQMR ILU0 | TFQMR with ILU0 Preconditioning |
| 68 | TFQMR ILUTH | TFQMR with ILUTH Preconditioning |
| Solution techniques for Kronecker representations |  |  |
| 111 | POWER | Power-Method ([12] pp. 203) |
| 112 | JOR | JOR ([1]) |
| 113 | SJOR | SJOR ([12] pp. 204f) |
| 114 | SOR | Real SOR ([17]) |
| 121 | POWER A/D | Power-Method with Aggregation/Disaggregation ([3, 4]) |
| 122 | JOR A/D | JOR with Aggregation/Disaggregation ([3, 4]) |
| 123 | SJOR A/D | SJOR with Aggregation/Disaggregation ([3, 4]) |
| 131 | GMRES | GMRES ([16], Spp 198ff) |
| 132 | DQGMRES | DQGMRES ([15]) |
| 133 | ARNOLDI | Arnoldi-Method ([16], pp. 191) |
| 134 | CGS | CGS ([16], pp. 222) |
| 135 | BICGSTAB | BiCGSTAB ([18])) |
| 136 | TFQMR | TFQMR ([13])) |
| 150 | PREPOWER | Power with Neumann preconditioning ([7]) |
| 151 | PREGMRES | GMRES with Neumann preconditioning ([7]) |
| 152 | PREARNOLDI | Arnoldi with Neumann preconditioning ([7]) |
| 154 | PRECGS | CGS with Neumann preconditioning ([6]) |
| 155 | PREBICGSTAB | BICGSTAB with Neumann preconditioning ([6]) |
| 156 | PRETFQMR | TFQMR with Neumann preconditioning([6]) |
| 157 | BSOR_BICGSTAB | BICGSTAB with Block SOR preconditioning ([11]) |
| 157 | BSOR_GMRES | GMRES with Block SOR preconditioning ([11]) |
| 157 | BSOR_TFQMR | TFQMR with Block SOR preconditioning ([11]) |
| 172 | STR_BSOR_T | Two level BSOR ([10]) |

Table 1: Overview of available solution methods

