

# Algorithms for nonnegative quadratic vector equations

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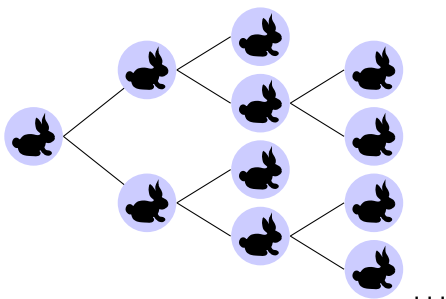
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18<sup>th</sup> Householder Symposium  
Tahoe City, CA 2011

# Markovian binary trees



MBTs model a colony of individuals that reproduce and die.

[Bean, Kontoleon, Taylor '04] [Hautphenne, Latouche, Rémiche '08]

## Simple example

$a = \mathbb{P}[\text{rabbit dies without spawning}]$

$b = \mathbb{P}[\text{rabbit spawns into two independent copies rabbit rabbit}]$

**Question:** starting from one individual, what is  $\mathbb{P}[\text{extinction}]$ ?

# A simple example

## Simple example

$a = \mathbb{P}$  [🐰 dies without spawning]

$b = \mathbb{P}$  [🐰 spawns into **two independent** copies 🐰 🐰 ]

**Question:** starting from one individual, what is  $\mathbb{P}$  [extinction]?

It is a nice elementary problem:  $x = \mathbb{P}$  [extinction]

$$x = a + b x^2$$

- Either 🐰 dies outright
- Or it spawns into two **independent** childs...  
... and the progenies of both die out

## The minimal solution

$x = a + bx^2$  has **two nonnegative solutions**. One is always 1, for  $a + b = 1$  (🐰 either reproduces or dies without)

Easy to prove that  $\mathbb{P}[\text{extinction}]$  is the **smaller** solution. Three cases:

**Subcritical**  $\mathbb{P}[\text{extinction}] = 1 > x_2$  (i.e. extinction = always)

**Supercritical**  $\mathbb{P}[\text{extinction}] = x_2 < 1$

**Critical** limit case: **1 double solution**

$\mathbb{P}[\text{extinction}] = 1$ , but needs an infinite time on average

# The vector case

Each 🐰 can be in  $N$  different states (e.g. age ranges)

$$a \in \mathbb{R}_+^N$$

$$a_i = \mathbb{P}[\text{🐰}_i \text{ dies}]$$

$$b \in \mathbb{R}_+^{N \times N \times N}$$

$$b_{ijk} = \mathbb{P}[\text{🐰}_i \text{ spawns into } \text{🐰}_j \text{ and } \text{🐰}_k]$$

$b$  contains  $N^3$  data!

Think to  $b$  as a **vector-valued bilinear form**

$$b : \mathbb{R}_+^N \times \mathbb{R}_+^N \rightarrow \mathbb{R}_+^N, \quad b(u, v) = \sum_{j,k} b_{ijk} u_j v_k$$

Our equation becomes

Markovian binary trees

$$x = a + b(x, x) \quad (\text{MBT})$$

# The classical algorithms

## Markovian binary trees

$$x = a + b(x, x) \quad (\text{MBT})$$

$e = \text{ones}(N, 1)$  is always a solution

$\mathbb{P}[\text{extinction}] = \text{minimal}$  nonnegative solution

Up to  $2^N$  nonnegative sol'ns, but there is always a **minimal one**:  
 $\hat{x}$  s.t.  $\hat{x} \leq x$  (component-by-component) for any other solution  $x$

**Subcritical or critical**:  $e$  is minimal, nothing to do

**Supercritical**: some other  $0 \leq \hat{x} \leq e$  is minimal: **how to compute it?**

# The classical algorithms

## Markovian binary trees

$$x = a + b(x, x) \quad (\text{MBT})$$

## Functional iterations [BKT '04]

$$x_{k+1} = a + b(x_k, x_k)$$

or something more elaborate, like

$$x_{k+1} = a + b(x_{k+1}, x_k)$$

i.e.

$$x_{k+1} = (I - b(\cdot, x_k))^{-1} a$$

$b(\cdot, x_k): \mathbb{R}_+^N \rightarrow \mathbb{R}_+^N$ : just a matrix

# The classical algorithms

## Markovian binary trees

$$x = a + b(x, x) \quad (\text{MBT})$$

Functional iterations [BKT '04]

Newton method [HLR '08]

$$x_{k+1} = (I - b(\cdot, x_k) - b(x_k, \cdot))^{-1} a$$

+ variants, e.g. [Hautphenne, Van Houdt '10]



# The classical algorithms

## Markovian binary trees

$$x = a + b(x, x) \quad (\text{MBT})$$

Functional iterations [BKT '04]

Newton method [HLR '08]

- When started from  $x_0 = 0$ , they converge monotonically:  
 $0 = x_0 \leq x_1 \leq x_2 \leq \dots \leq x^*$
- neat probabilistic interpretations:  
 $x_k = \mathbb{P}$  [extinction truncated to the  $k$ -th generation, or to a subtree]
- Become slower when **close to critical**:  
need more generations to capture the behaviour of the tree

# Deflation

Close to a double solution, and for Newton **double = trouble**

But one of these solutions  $x = e$  is known, we want to **deflate** it:

Set  $y := e - x$  **survival probability**; (MBT) becomes

The optimistic equation

$$y = \underbrace{(b(e - y, \cdot) + b(\cdot, e))}_{:= P_y} y = P_y y$$

Functional it's/Newton in this form: nothing changes, but...

# Perron vector-based algorithms

## The optimistic equation

$$y = \underbrace{(b(e - y, \cdot) + b(\cdot, e))}_{:= P_y} y = P_y y$$

New way to see the same equation:  $y$  is the Perron vector of a matrix depending (linearly) on  $y$  itself

$$y = PV(P_y) \tag{PE}$$

(+suitable **normalization** for the eig'vec:  $w^T \cdot \text{Residual} = 0$  for some  $w$ )

- Fixed point iteration based on (PE):  $y_{k+1} = PV(P_{y_k})$
- Newton's method

## Numerical experiments

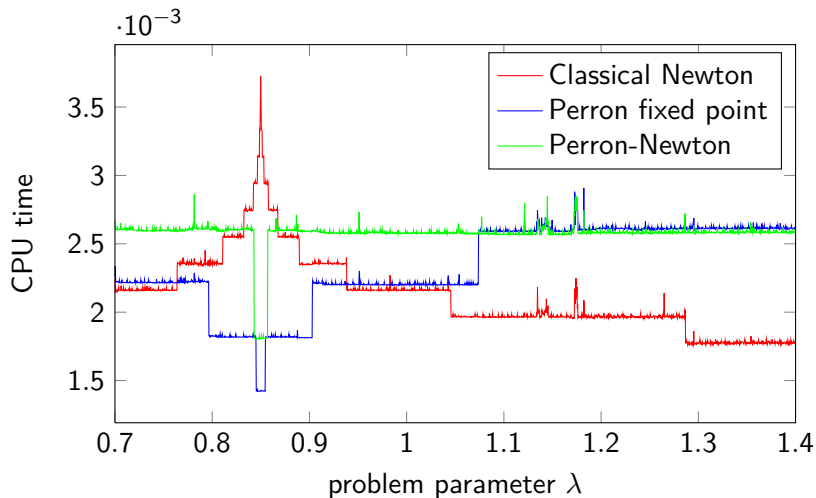


Figure: CPU time for a parameter-dependent problem [BKT '08, example 1]; lower=better

# Numerical experiments

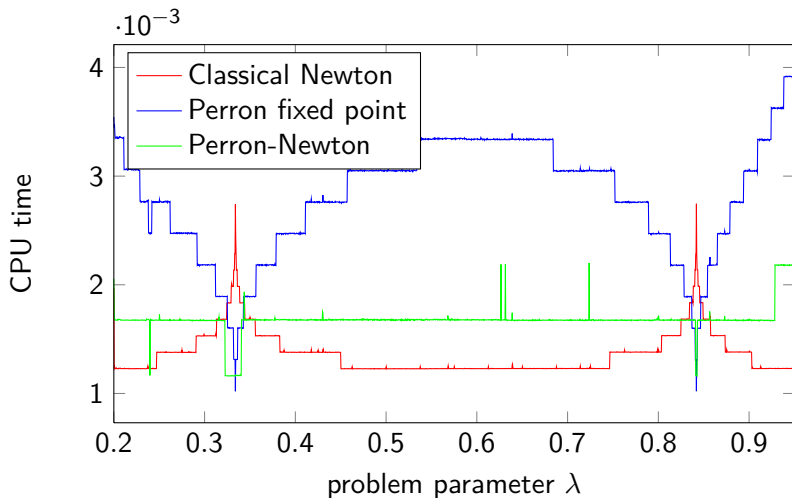


Figure: CPU time for a parameter-dependent problem [BKT '08, example 2]; lower=better

# Convergence results

- Convergence is not monotonic
- Convergence is not guaranteed for very far-from-critical problems

Theorem [Meini, P., SIMAX 2011]

- Explicit formula for the Jacobian of the Perron iteration
- For a special normalization choice,  
if **problem**  $\rightarrow$  **critical** then  $\rho(\text{Jac}) \rightarrow 0$

Thus, locally convergent for close-to-critical  
with speed that **tends to superlinear**

Theorem [Bini, Meini, P., NLAA (to appear)]

When the algorithm converges, it converges to the right solution  $\hat{x}$

# Applicability

We may ensure applicability even when strict positivity/irreducibility assumptions do not hold:

- 1 **deflate** away entries  $i$  s.t.  $\hat{x}_i = 0$ : they can be determined in  $O(N^3)$  from the nonzero pattern of  $a$  and  $b$
- 2 all  $P_y$  have the same nonzero pattern; if they are **reducible**, we may split the problem into two subproblems

(as with linear equations; **idea**: if  $P_y = \begin{bmatrix} P_{11} & P_{12} \\ 0 & P_{22} \end{bmatrix}$ , we can solve for the second block alone and back-substitute)

# A unifying framework

Why is this problem interesting?

$$Mx = a + b(x, x)$$

$$XCX - AX - XD + B = 0 \quad (\text{Nonsym. Riccati})$$

$$PX^2 + QX + R = 0 \quad (\text{QBD equation})$$

$$\begin{cases} Ix = (Py) \cdot x + e \\ Iy = (Qx) \cdot y + e \end{cases} \quad (\text{Transport theory})$$

With a bit of  $\text{vec}(\cdot)$ , several matrix equations can be reduced to (MBT)

Although no known  $(x = e)$  solution  $\rightarrow$  no PV-based algorithms

Open problem

Can we recover something similar from partial information (e.g., one known eigenpair of  $X$ )? Would carry over to many matrix equations



## Common aspects

- **minimal solution**  $x_* \geq 0$ , i.e.,  $x_* \leq x$  for any other solution  $x$
- functional iterations and Newton's method exhibit **monotonic convergence**:  $0 = x_0 \leq x_1 \leq x_2 \leq \dots \rightarrow x_*$
- **close-to-critical problems**: when close to a double solution, convergence is slower and more unstable

Common framework to work with several equations from different applications [P., to appear (LAA)]

Advantages:

- **unified proofs**: clear hypotheses, role of strict positivity of  $x_*$   
no matrix structure or spectral properties needed
- **unified algorithms**: take an algorithm for one equation, apply it to the others

**Example** a Newton variant [Hautphenne, Van Houdt '10] useful for the transport theory eqn

# Conclusions

## Open questions

- Understand doubling methods (SDA/Cyclic Reduction) in this framework:

If we try to construct doubling for (MBT) we get **Newton** instead; are the two related?

- Shift technique + what happens to spectral properties?
- Perron-based algorithms without a “full” known solution  $x = e$

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- Understand doubling methods (SDA/Cyclic Reduction) in this framework:

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- Shift technique + what happens to spectral properties?
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Thanks for your attention!

