

Componentwise accurate numerical methods for Markov-modulated Brownian motion

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Subtraction-free algorithms

Error analysis in an (imprecise) slogan

TL;DR: when you subtract two close numbers, you lose accuracy.

More precise: instead of a and b , a computer may store $a + \delta$ and $b + \varepsilon$; the number $(a + \delta) - (b + \varepsilon)$ may be at a large relative distance from $a - b$ (if a and b have the same sign).

So, **let's stop doing subtractions.**

Luckily, for many probabilities computations this is possible.

E.g., computing AB , for $A \geq 0, B \geq 0$.

Most subtractions come from M-matrices, but **we can avoid them!**

Regular M-matrices

A matrix A with sign pattern (possibly including zeros)

$$A = \begin{bmatrix} + & - & - & - \\ - & + & - & - \\ - & - & + & - \\ - & - & - & + \end{bmatrix}$$

is called **regular M-matrix** if there are $\mathbf{v} > \mathbf{0}$, $\mathbf{w} \geq \mathbf{0}$ such that $A\mathbf{v} = \mathbf{w}$.
E.g., $(-Q)\mathbf{1} = \mathbf{0}$ for the rate matrix of a CTMC

Attention! [Guo CH, 2013]

Not all M-matrices are regular! E.g., $\begin{bmatrix} 0 & 0 \\ -1 & 0 \end{bmatrix}$

GTH algorithm

For a regular M-matrix A , one can store its **off-diagonal** entries, \mathbf{v} and \mathbf{w} (**triplet representation**).

$$A = \begin{bmatrix} ? & - & - & - \\ - & ? & - & - \\ - & - & ? & - \\ - & - & - & ? \end{bmatrix} \quad A\mathbf{v} = \mathbf{w}$$

GTH-like algorithm [Grassmann et al '85, O'Conneide '93, Alfa et al '02...]

Given a triplet representation for $A \in \mathbb{R}^{n \times n}$, one can compute $B = A^{-1}$ subtraction-free, obtaining **perfect componentwise accuracy**:

$$|\tilde{b}_{ij} - b_{ij}| \leq O(n^3) \cdot |b_{ij}| \cdot \text{eps}.$$

Variants: LU factorization, left and right kernel, Perron vector.

Variant: $\mathbf{v}^\top A = \mathbf{w}^\top$

GTH algorithm

An example

$A = \begin{bmatrix} 1 & -1 \\ -1 & 1 + \varepsilon \end{bmatrix}$: can only compute inverse up to accuracy $\kappa(A) \approx \varepsilon^{-1}$.

$A = \begin{bmatrix} ? & -1 \\ -1 & ? \end{bmatrix}$ such that $A \begin{bmatrix} 1 \\ 1 \end{bmatrix} = \begin{bmatrix} 0 \\ \varepsilon \end{bmatrix}$: full accuracy possible.

Works especially well when dealing with different orders of magnitude.

Plan: triplet representation are a great idea — let's rewrite our matrix iterations to use them!

Obtaining triplets

Theorem [Nguyen P. '16 — or earlier?]

Given a regular M -matrix and its triplet representation partitioned as

$$\begin{bmatrix} A & B \\ C & D \end{bmatrix} \begin{bmatrix} v_1 \\ v_2 \end{bmatrix} = \begin{bmatrix} w_1 \\ w_2 \end{bmatrix},$$

one can obtain explicitly **without subtractions** triplet representations for its submatrices and Schur complements (ensorings):

$$\begin{aligned} Dv_2 &= w_2 - Cv_1, \\ (A - BD^{-1}C)v_1 &= w_1 - BD^{-1}w_2. \end{aligned}$$

Cyclic reduction

CR solves matrix equations of the form $R^2A - RB + C = 0$, with $A, C \geq 0$, and $B - A - C$ a regular M-matrix [Bini et al '01, BLM book]

Cyclic Reduction

$$\begin{aligned}A_0 &= A, B_0 = \hat{B}_0 = B, C_0 = C \\A_{k+1} &= A_k B_k^{-1} A_k, \\B_{k+1} &= B_k - A_k B_k^{-1} C_k - C_k B_k^{-1} A_k, \\C_{k+1} &= C_k B_k^{-1} C_k, \\\hat{B}_{k+1} &= \hat{B}_k - C_k B_k^{-1} A_k.\end{aligned}$$

At the end of the iteration, $R = C_0 \hat{B}_\infty^{-1}$, where $\hat{B}_\infty = \lim \hat{B}_k$.

Doubling algorithm

An unusual matrix iteration that can be seen as repeated censoring / Schur complementation:

$$\begin{bmatrix} E_{new} & G_{new} \\ H_{new} & F_{new} \end{bmatrix} = \begin{bmatrix} 0 & G \\ H & 0 \end{bmatrix} + \begin{bmatrix} E & 0 \\ 0 & F \end{bmatrix} \left(I - \begin{bmatrix} 0 & G \\ H & 0 \end{bmatrix} \right)^{-1} \begin{bmatrix} E & 0 \\ 0 & F \end{bmatrix}$$

Keep track of triplet representations at each step \implies
componentwise-accurate algorithms for fluid queues / Riccati equations.

[Xue et al '12, Nguyen P. '15, Xue Li '16]

Markov-modulated Brownian motion [Asmussen '95, Karandikar

Kulkarni '95, Rogers '94]

- $\phi(t)$ continuous-time Markov chain with transition matrix $Q \in \mathbb{R}^{n \times n}$;
- $y(t)$ evolves as Brownian motion with drift $d_{\phi(t)}$ and variance $v_{\phi(t)}$.

The invariant density follows

$$\mathbf{p}''(x)V - \mathbf{p}'(x)D + \mathbf{p}(x)Q = \mathbf{0}.$$

$V, D \in \mathbb{R}^{n \times n}$ diagonal matrices containing v_i, d_i .

Invariant density and many properties can be computed using an **invariant pair**, i.e., $(X \in \mathbb{R}^{\ell \times \ell}, U \in \mathbb{R}^{\ell \times n})$ such that

$$X^2UV - XUD + UQ = 0.$$

[Rogers '94, Ivanovs, '10, Betcke Kressner '11, Gohberg et al '82]

Often, $U = \begin{bmatrix} I & \Psi \end{bmatrix}$.

Invariant pairs and Cyclic Reduction

How do we compute invariant pairs? **Cyclic reduction** gives a special one:

$$R^2/A - R/B + I/C = 0.$$

Plan: tinker with the problem to turn it into this form.

Discretizing transformation from eigenvalue properties: [Bini et al '10]

- we need a “continuous-time stable” pair: $\text{eig}(X) \subseteq$ left half-plane.
- CR produces a “discrete-time stable” one: $\text{eig}(R) \subseteq$ unit circle

So we make a change of variables $R = f(X)$.

$R = (I + X)(I - X)^{-1}$ **won't work**: cannot find $X = f^{-1}(R)$ subtraction-free.

Instead, we use $R = I + hX$, with h sufficiently small.

The algorithm

- 1 Choose h small enough so that

$$v_i + d_i h + q_{ii} h^2 > 0 \quad (*)$$

(and the subtraction is 'tame').

- 2 Set $A := \frac{1}{h^2} V \geq 0$, $B := 2\frac{1}{h^2} V + \frac{1}{h} D$, $C := \frac{1}{h^2} V + \frac{1}{h} D + Q \geq 0$.
- 3 Use subtraction-free CR on (A, B, C) to compute R .
- 4 Compute the off-diagonal of $X = h^{-1}(R - I)$.
- 5 Compute the left Perron vector μ of Q using the triplet $Q\mathbf{1} = \mathbf{0}$.
- 6 Compute the triplet $\mu(-X) = \frac{1}{h} \mu A_\infty \hat{B}_\infty^{-1}$, and obtain $\text{diag}(X)$.

Works whenever $v_i > 0$ for all i (positive variances).

Otherwise, we can't enforce (*)

Zero variances

- In $E_2 = \{i : v_i = 0, d_i \leq 0\}$, we can't obtain $v_i + d_i h + q_{ii} h^2 > 0$.
- We won't be able to choose $U = I$, because we only have enough stable eigenvalues to form an invariant pair of size $n - |E_2|$.

Solution to both problems: **shift** infinite eigenvalues [He et al '01].

From:

$$A = \begin{bmatrix} + & 0 \\ 0 & 0 \end{bmatrix}, \quad B = \begin{bmatrix} + & 0 \\ 0 & - \end{bmatrix}, \quad C = \begin{bmatrix} + & + \\ + & ?? \end{bmatrix}$$

move 2nd column “one matrix to the left” and change its sign:

$$\hat{A} = \begin{bmatrix} + & 0 \\ 0 & + \end{bmatrix}, \quad \hat{B} = \begin{bmatrix} + & - \\ 0 & M \end{bmatrix}, \quad \hat{C} = \begin{bmatrix} + & 0 \\ + & 0 \end{bmatrix}.$$

Shift \iff **differentiating** some of the equations.

See “index reduction” in ODE literature. [Kunkel Mehrmann '06]

Recovering the solution

Still, R is **not** the final solution.

The shift trick adds spurious zero eigenvalues: $R = \begin{bmatrix} + & 0 \\ + & 0 \end{bmatrix} \hat{B}_\infty^{-1}$.

Solution to remove them: switch to a different invariant pair:

$$(KRK^{-1})^2 KA + (KRK^{-1})KB + KC = 0.$$

$$\begin{bmatrix} R_{11} & 0 \\ R_{12} & 0 \end{bmatrix}^2 \begin{bmatrix} I & \Psi \\ 0 & I \end{bmatrix} A + \begin{bmatrix} R_{11} & 0 \\ R_{12} & 0 \end{bmatrix} \begin{bmatrix} I & \Psi \\ 0 & I \end{bmatrix} B + \begin{bmatrix} I & \Psi \\ 0 & I \end{bmatrix} C = 0.$$

New R is block-triangular \implies “reduced invariant pair” $(R_{11}, [I \ \Psi])$.

True but not obvious: all this can be done **subtraction-free**.

The (general) algorithm

- 1 Choose h small so that $h^{-2}v_i + h^{-1}d_i + q_{ii} > 0$ for all $i \notin E_2$.
- 2 Set $A := \frac{1}{h^2}V \geq 0$, $B := 2\frac{1}{h^2}V + \frac{1}{h}D$, $C := \frac{1}{h^2}V + \frac{1}{h}D + Q \geq 0$.
- 3 **Shift technique** to produce equivalent $\hat{A}, \hat{B}, \hat{C}$.
- 4 Use componentwise accurate CR on $(\hat{A}, \hat{B}, \hat{C})$ to compute \hat{B}_∞ .
- 5 **Use similarity** as in the previous slide to compute $(R_{11}, [I \ \Psi])$.
- 6 Compute the off-diagonal of $X = h^{-1}(R_{11} - I)$.
- 7 Compute the left Perron vector μ of Q using the triplet $Q\mathbf{1} = \mathbf{0}$.
- 8 Compute a triplet $\mathbf{v}^\top(-X) = \mathbf{w}^\top$ (**long formula omitted**).

Works **even with zero variances**.

Experiments: the competitors

- KK** [Karandikar Kulkarni '95]: compute eigenvalues explicitly.
- AS** [Agapie Sohraby '01]: iterative algorithm for span(stable eigenvalues), then orthogonal transformations.
- LN** [Latouche Nguyen '15]: (non-subtraction-free) Cyclic Reduction.
- QZ** QZ algorithm: orthogonal transformations — well-known linear algebra workhorse.
- NP** our new algorithm.

Problem	KK	AS	LN	QZ	NP
NP15	2.7e-12	2.5e-07	2.9e-13	1.8e-12	1.7e-16
NP15s	1.3e-12	2.2e-07	-	6.2e-13	1.8e-16
rand8	2.8e-15	1.5e-15	1.6e-15	2.4e-15	2.7e-16
rand8s	2.9e-15	1.8e-13	-	2.3e-15	3.1e-16
rand20	4.4e-15	9.6e-14	5.6e-15	4.8e-15	3.0e-16
rand20s	3.2e-15	3.0e-12	-	4.1e-14	1.1e-15
rand50	5.9e-15	4.0e-14	4.0e-14	5.6e-14	6.9e-16
rand50s	5.6e-14	1.2e-10	-	3.5e-14	5.2e-16
imb8	9.7e-12	1.9e-09	1.1e+00	7.1e-13	9.0e-13
imb8s	2.6e-14	1.3e-08	-	1.3e-12	1.1e-15
imb20	4.6e-11	2.1e-07	3.2e-04	1.1e-09	9.1e-12
imb20s	4.4e-12	6.9e-06	-	5.9e-12	4.0e-13
imb50	2.0e-10	9.8e-06	7.2e-01	1.0e-08	8.3e-10
imb50s	2.0e-10	3.3e-05	-	1.0e+00	2.6e-13

Table: Forward error $\frac{\|\tilde{X} - X\|}{\|X\|}$

Error on $U = [I \ \Psi]$

Problem	KK	AS	LN	QZ	NP
NP15s	2.3e-15	1.8e-11	-	2.8e-15	1.3e-16
rand8s	1.2e-14	3.7e-13	-	2.4e-15	2.5e-15
rand20s	7.1e-15	7.7e-11	-	6.7e-14	2.1e-15
rand50s	3.4e-14	3.5e-09	-	5.3e-14	4.7e-16
imb8s	8.3e-15	5.2e-09	-	1.1e-11	5.2e-15
imb20s	1.4e-10	1.9e-08	-	2.8e-11	4.0e-11
imb50s	6.9e-11	9.0e-09	-	1.0e-04	6.1e-08

Table: Forward error $\frac{\|\tilde{\Psi} - \Psi\|}{\|\Psi\|}$

Conclusions and open points

- Subtraction-free, **componentwise accurate** algorithm for MMBM. [Nguyen P. arXiv:1605.01482]
- There's also [Nguyen P. '15] for fluid queues.
- Similar to [Ramaswami '99] QBD construction but for MMBM.
- **Future plan:** remove the 2^k factor in the error for CR.
- Ideas from ODEs (index reduction, stability conditions) and linear algebra (shift technique, invariant pairs).

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Thanks for your attention!

Questions?