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Rush-Larsen time-stepping methods of high order for stiff problems in cardiac electrophysiology

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Abstract

To address the issues of stability and accuracy for reaction-diffusion equations, the development of high order and stable time-stepping methods is necessary. This is particularly true in the context of cardiac electrophysiology, where reaction-diffusion equations are coupled with stiff ODE systems. Many research have been led in that way in the past 15 years concerning implicit-explicit methods and exponential integrators. In 2009, Perego and Veneziani proposed an innovative time-stepping method of order 2. In this paper we present the extension of this method to the orders 3 and 4 and introduce the Rush-Larsen schemes of order k (shortly denoted RL_k). The RL_k schemes are explicit multistep exponential integrators. They display a simple general formulation and an easy implementation.

The RL_k schemes are shown to be stable under perturbation and convergent of order k . Their Dahlquist stability analysis is performed. They have a very large stability domain provided that the stabilizer associated with the method captures well enough the stiff modes of the problem. The RL_k method is numerically studied as applied to the membrane equation in cardiac electrophysiology.

Keywords: stiff equations, explicit high-order multistep methods, exponential integrators, stability and convergence, Dahlquist stability

Subject classification: 65L04, 65L06, 65L20, 65L99

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Introduction

The general purpose of this article is the time integration of stiff reaction-diffusion equations, more precisely when a coupling with ODE systems occurs. In that framework, the matters of stability and accuracy together with the non linear nature of the equations are of first importance (as developed below). As a systemic example of these questions, we will consider the monodomain model in cardiac electrophysiology [3, 4, 5]. It has the general formulation: on the heart domain Ω and on the time interval $[0, T]$,

$$\frac{\partial v}{\partial t} = Av + f_1(v, w) + s(x, t), \quad \frac{\partial w}{\partial t} = f_2(v, w), \quad (1)$$

where A is a diffusion operator. The first unknown $v : \Omega \times [0, T] \rightarrow \mathbb{R}$ is the transmembrane-potential. The second (vector) unknown $w : \Omega \times [0, T] \rightarrow \mathbb{R}^N$ gathers variables describing the state of the cellular membrane : it incorporates ionic concentrations and gating variables. The source term $s(x, t)$ is an applied stimulation current. The reaction terms f_1 and f_2 model ionic currents and voltage across the cellular membrane, they are named ionic models. Ionic models originally have been developed by Hodgkin and Huxley [17] in 1952 for the squid axon. Several ionic models have been especially designed for cardiac cells, such as the Beeler and Reuter model [1], the Luo and Rudy models [22, 21] or the TNNP model [27]. A comprehensive list is available in [26].

Numerical simulations in cardiac electrophysiology have to face two difficulties. The first one is the nature of the reaction terms f_1 and f_2 in (1). It is non linear and the operation $(v, w) \rightarrow f_i(v, w)$ has a significant cost. For example, this operation for the TNNP model [27] involves the computation of 50 exponentials. These operations need to be performed at every node of the grid. They represent a large computational load. Their total amount needs to be maintained as low as possible. Fully implicit time-stepping methods (that require a non linear solver) therefore are avoided. The second difficulty is the stiffness. The management of the numerical instabilities thus is challenging since implicit methods are not affordable here. More precisely, the stiffness of the monodomain model (1) is caused by the presence of different space and time scales. The solutions of (1) display sharp wave-fronts. This is commonly coped with by resorting to very fine space and time grids, associated with high computational costs.

In this context, our strategy for the numerical resolution of (1) is to go towards high order methods, in order to have accurate simulations on much coarser space and time grids. A high order time-stepping method is required that fulfils these two conditions: it must have strong stability properties and it is explicit for the reaction terms. To this aim, we will focus in this paper on the time integration of stiff ODE systems,

$$\frac{dy}{dt} = f(t, y), \quad y(0) = y_0, \quad (2)$$

for which a reformulation of the following kind is available,

$$\frac{dy}{dt} = a(t, y)y + b(t, y), \quad y(0) = y_0. \quad (3)$$

The linear part $a(t, y)$ will be referred to as the stabilizer. The stabilizer will be inserted in the numerical scheme in order to stabilize the computations.

That approach is relevant in cardiac electrophysiology where stiffness is induced by the coexistence of fast and slow variables. The fast variables (gating variables) are described by equations in the ODE system in (1) of the form:

$$\frac{\partial w_i}{\partial t} = f_{2,i}(v, w) = \frac{w_\infty(v, w) - w_i}{\tau_i(v, w)},$$

that motivates the reformulation (3) with a diagonal stabilizer $a = \text{diag}(-1/\tau_i)$. More details are given in Section 4.1.

Exponential integrators are well suited in that framework. We refer to [23, 14, 11] for general reviews. They have been widely studied for the quasi-linear equations, $\partial_t y = Ay + b(t, y)$, see *e.g.* [12, 7, 13, 16, 28, 20]. The basic idea is to use the exact solution of the linearized equation in order to stabilize the numerical scheme. In general this implies to compute a matrix exponential. This is the supplementary cost associated to exponential integrators.

The targeted problem (3) displays a non constant linear part $a(t, y)$. Exponential integrators have been less studied in that case. Exponential integrators of Adams type for a non constant linear part have been first considered by Lee and Preiser [19] in 1978 and by Chu [2] in 1983. Recently, Ostermann *et al.* [15, 18] developed and analyzed the linearized exponential Adams method. The original problem (2) is reformulated after each time step as,

$$\frac{dy}{dt} = J_n y + c_n(t, y), \quad J_n = \partial_y f(t_n, y_n), \quad c_n(t, y) = f(t, y) - J_n y.$$

The Jacobian matrix J_n is used as the stabilizer with the following induced drawbacks. It requires the computation of matrix exponentials at every time steps. Moreover, when the fast variables of the system are known, stabilization can be performed only on these variables. Considering the full Jacobian as the stabilizer implies unnecessary computational efforts. To avoid these problems, an alternative is to set the stabilizer as a part or as an approximation of the Jacobian. This has been analyzed in [29], [25] and [6] for exponential Rosenbrock, exponential Runge-Kutta and exponential Adams type methods respectively. For exponential Adams type methods, equation (3) is reformulated after each time step as:

$$\frac{dy}{dt} = a_n y + c_n(t, y), \quad a_n = a(t_n, y_n), \quad c_n(t, y) = f(t, y) - a_n y.$$

The resulting scheme is (see details in [15, 6]):

$$y_{n+1} = y_n + h [\varphi_1(a_n h) (a_n y_n + \gamma_1) + \varphi_2(a_n h) \gamma_2 + \dots + \varphi_k(a_n h) \gamma_k], \quad (4)$$

where γ_i are the coefficients of the Lagrange interpolation polynomial of $c_n(t, y)$ (in a classical k -step setting) and where the functions φ_j are given by

$$\varphi_0(z) = e^z, \quad \varphi_{j+1}(z) = \frac{\varphi_j(z) - 1/j!}{z}. \quad (5)$$

Independently, Perego and Veneziani [24] presented in 2009 an innovative exponential integrator of order 2, of a different nature:

$$y_{n+1} = y_n + h\varphi_1(\alpha_n h) (\alpha_n y_n + \beta_n), \quad (6)$$

also involving two constants α_n and β_n to be computed at each time step. It results in a schemes with a very simple definition. It is in particular simpler than the exponential Adams integrators (4). The essential difference with the previous approaches is that $\alpha_n \neq a(t_n, y_n)$. Precisely, the constants α_n and β_n are given by $\alpha_n = 3/2a_n - 1/2a_{n-1}$ and $\beta_n = 3/2b_n - 1/2b_{n-1}$ with $a_j = a(t_j, y_j)$ and $b_j = b(t_j, y_j)$.

In this paper we will study the general formulation (6). We will show that such schemes also exist at the orders 3 and 4, for an explicit definition of the two constants α_n and β_n . These schemes will be referred to as Rush-Larsen schemes of order k , shortly denoted RL_k . They will be shown to be stable under perturbation and convergent of order k . The Dahlquist stability analysis for the RL_k schemes is also performed. It is a practical tool that allows to dimension the time step h with respect to the variations of $f(t, y)$ in problem (2), see *e.g.* [10]. When considering varying stabilizers, the stability domain depends on how $f(t, y)y$ is decomposed in Equation (3), following [24]. The stability domains are numerically computed and shown to be much larger than in the absence of stabilization (*i.e.* when $a(t, y) = 0$) provided that $a(t, y)$ captures well enough the variations of $f(t, y)$. The performances of the RL_k method are evaluated for the membrane equation in cardiac electrophysiology. They are compared with the exponential Adams integrators (4). The two methods have a very similar robustness to stiffness. They both allow stable computations on coarse time grids.

At large time step, for the presented test case, the RL_3 and RL_4 schemes are more accurate, meanwhile with a simpler implementation and thus a lighter cost.

The paper is organized as follows. The RL_k schemes are derived in Section 1 and their numerical analysis is made in Sections 1 and 2. The Dahlquist stability analysis is in Section 3. The numerical results are presented in Section 4. The paper ends with the conclusion Section.

In the sequel h denotes the time step and $t_n = nh$ the associated time instants.

1 RL_k scheme definition and consistency

We start this section with the definition of the Rush-Larsen schemes.

Definition 1. The RL_k scheme is an explicit k -step method. It is defined with the formulation (6) for the following setting of α_n and of β_n : for the RL_2 scheme

$$\alpha_n = \frac{3}{2}a_n - \frac{1}{2}a_{n-1}, \quad \beta_n = \frac{3}{2}b_n - \frac{1}{2}b_{n-1},$$

for the RL_3 scheme

$$\begin{aligned}\alpha_n &= \frac{1}{12}(23a_n - 16a_{n-1} + 5a_{n-2}), \\ \beta_n &= \frac{1}{12}(23b_n - 16b_{n-1} + 5b_{n-2}) + \frac{h}{12}(a_nb_{n-1} - a_{n-1}b_n),\end{aligned}$$

and for the RL_4 scheme

$$\begin{aligned}\alpha_n &= \frac{1}{24}(55a_n - 59a_{n-1} + 37a_{n-2} - 9a_{n-3}), \\ \beta_n &= \frac{1}{24}(55b_n - 59b_{n-1} + 37b_{n-2} - 9b_{n-3}) \\ &\quad + \frac{h}{12}(a_n(3b_{n-1} - b_{n-2}) - (3a_{n-1} - a_{n-2})b_n),\end{aligned}$$

where $a_j = a(t_j, y_j)$ and $b_j = b(t_j, y_j)$.

A solution $y(t)$ of Equation (3) on a time interval $[0, T]$ is fixed. It is recalled that the scheme (6) is consistent of order k if:

- being given a time step h and a time instant $kh \leq t_n \leq T - h$,
- being given the numerical approximation y_{n+1} in (6) computed with $y_{n-j} = y(t_{n-j})$ for $j = 0 \dots k - 1$,

we have $|y_{n+1} - y(t_n + h)| \leq Ch^{k+1}$, for a constant C only depending on the problem (3) data a, b, y_0 and on T .

Proposition 1. Assume that the functions $a(t, y)$ and $b(t, y)$ in problem (3) are C^k regular. Moreover assume that $a(t, y)$ either is a diagonal matrix or a constant linear operator.

Then the RL_k scheme is consistent of order k .

Remark 1. In the case of a constant linear part $a(t, y) = A$, we always have $\alpha_n = A$. The definition of β_n also simplifies at the order 3 and 4,

$$\begin{aligned}RL_3 : \quad \beta_n &= \frac{1}{12}(23b_n - 16b_{n-1} + 5b_{n-2}) - \frac{h}{12}A(b_n - b_{n-1}). \\ RL_4 : \quad \beta_n &= \frac{1}{24}(55b_n - 59b_{n-1} + 37b_{n-2} - 9b_{n-3}) - \frac{h}{12}A(2b_n - 3b_{n-1} + b_{n-2}).\end{aligned}$$

Remark 2. The assumption “ $a(t, y)$ either is a diagonal matrix or a constant linear operator” in Proposition 1 has the following origin. To analyze the scheme consistency we will derive a Taylor expansion in h of the scheme formulation (6). That series is computed with the help of Taylor expansions in h for α_n and β_n .

Assume the simple form $\alpha_n = \alpha_0 + h\alpha_1$. We need to expand $\varphi_1(\alpha_n h)$ as a series in h . The function φ_1 is analytic on $\mathbb{C}$. However in the matrix case, the equality, $\varphi_1(M +$

$N) = \varphi_1(M) + \varphi'_1(M)N + \dots + \varphi_1^{(i)}(M)N^i/i! + \dots$ holds if M and N are commutative matrices. Therefore one cannot expand $\varphi_1(\alpha_n h)$ without the assumptions that α_0 and α_1 are commutative.

That difficulty vanishes if $a(t, y)$ is constant or a varying diagonal matrix.

The proof of Proposition 1 is based on the following result.

Lemma 1. With the same assumption as in Proposition 1, the scheme (6) is consistent of order k if:

$$\begin{aligned} k = 2 : \quad & \alpha_n = a_n + \frac{1}{2}a'_n h + O(h^2), \quad \beta_n = b_n + \frac{1}{2}b'_n h + O(h^2). \\ k = 3 : \quad & \alpha_n = a_n + \frac{1}{2}a'_n h + \frac{1}{6}a''_n h^2 + O(h^3), \\ & \beta_n = b_n + \frac{1}{2}b'_n h + \frac{1}{12}(a'_n b_n - a_n b'_n)h^2 + O(h^3). \\ k = 4 : \quad & \alpha = a_n + \frac{1}{2}a'_n h + \frac{1}{6}a''_n h^2 + \frac{1}{24}a'''_n h^3 + O(h^4), \\ & \beta = b_n + \frac{1}{2}b'_n h + \frac{1}{12}(a'_n b_n - a_n b'_n)h^2 + \left(\frac{1}{24}b'''_n + \frac{1}{24}(a''_n b_n - a_n b''_n) \right) h^3 + O(h^4). \end{aligned}$$

Where a'_n, a''_n, a'''_n and b'_n, b''_n, b'''_n denote the successive derivatives at time t_n of the functions $t \mapsto a(t, y(t))$ and $t \mapsto b(t, y(t))$.

Proof of lemma 1. By assumption the functions a and b in problem (3) are C^k regular. Therefore a solution y of problem (3) on a closed time interval $[0, T]$ is C^{k+1} regular. Its derivatives up to order $k+1$ can be bounded by constants only depending on the problem (3) data and on T . The Taylor expansion of y at time instant t_n is:

$$y(t_n + h) = y(t_n) + \sum_{j=1}^k \frac{s_j}{j!} h^j + O(h^{k+1}),$$

with $s_j = y^{(j)}(t_n)$. Using that $y' = ay + b$ we get

$$\begin{aligned} s_1 &= a_n y_n + b_n, \\ s_2 &= (a'_n + a_n^2) y_n + a_n b_n + b'_n, \\ s_3 &= (a''_n + 3a_n a'_n + a_n^3) y_n + b''_n + a_n b'_n + 2a'_n b_n + a_n^2 b_n, \\ s_4 &= (a'''_n + 4a''_n a_n + 3a_n'^2 + 6a'_n a_n^2 + a_n^4) y_n \\ &\quad + b'''_n + b''_n a_n + 3a''_n b_n + 5a'_n a_n b_n + 3a'_n b'_n + a_n^3 b_n + a_n^2 b'_n. \end{aligned}$$

A series expansion in h for α_n and for β_n is introduced as

$$\begin{aligned} \alpha_n &= \alpha_{n,0} + \alpha_{n,1}h + \dots + \alpha_{n,k-1}h^{k-1} + O(h^k), \\ \beta_n &= \beta_{n,0} + \beta_{n,1}h + \dots + \beta_{n,k-1}h^{k-1} + O(h^k). \end{aligned}$$

With the assumption that $a(t, y)$ is either constant or a diagonal matrix (see Remark 2), a Taylor expansion of the numerical solution y_{n+1} in (6) can be performed:

$$y_{n+1} = y(t_n) + \sum_{j=1}^k \frac{r_j}{j!} h^j + O(h^{k+1}).$$

A direct computation of the r_j gives,

$$r_1 = \alpha_{n,0} y_n + \beta_{n,0},$$

$$r_2 = (2\alpha_{n,1} + \alpha_{n,0}^2) y_n + 2\beta_{n,1} + \alpha_{n,0} \beta_{n,0},$$

$$r_3 = (6\alpha_{n,2} + \alpha_{n,0}^3 + 6\alpha_{n,0}\alpha_{n,1}) y_n + 3\alpha_{n,1}\beta_{n,0} + 6\beta_{n,2} + \alpha_{n,0}^2\beta_{n,0} + 3\alpha_{n,0}\beta_{n,1},$$

$$r_4 = (24\alpha_{n,0}\alpha_{n,2} + 24\alpha_{n,3} + 12\alpha_{n,1}\alpha_{n,0}^2 + 12\alpha_{n,1}^2 + \alpha_{n,0}^4) y_n \\ + 12\alpha_{n,2}\beta_{n,0} + 24\beta_{n,3} + 12\alpha_{n,0}\beta_{n,2} + 12\alpha_{n,1}\beta_{n,1} + 4\alpha_{n,0}^2\beta_{n,1} + 8\alpha_{n,0}\alpha_{n,1}\beta_{n,0} + \alpha_{n,0}^3\beta_{n,0}.$$

The condition to be consistent of order k is: $r_i = s_i$ for $1 \leq i \leq k$. Lemma 1 consistency conditions are obtained by solving recursively these relations. $\square$

Proof of Proposition 1. It is a direct and simple consequence of the backwards differentiation formula, that we first recall. Consider a real function f , its derivatives can be approximated as follows (with obvious notations). For the first derivative,

$$f'_n = \frac{f_n - f_{n-1}}{h} + O(h), \\ = \frac{1}{2h} (3f_n - 4f_{n-1} + f_{n-2}) + O(h^2), \\ = \frac{1}{6h} (11f_n - 18f_{n-1} + 9f_{n-2} - 2f_{n-3}) + O(h^3).$$

For the second derivative,

$$f''_n = \frac{1}{h^2} (f_n - 2f_{n-1} + f_{n-2}) + O(h), \\ = \frac{1}{h^2} (2f_n - 5f_{n-1} + 4f_{n-2} - f_{n-3}) + O(h^2).$$

For the third derivative,

$$f'''_n = \frac{1}{h^3} (f_n - 3f_{n-1} + 3f_{n-2} - f_{n-3}) + O(h),$$

With these formula, the consistency condition at order 3 on α_n in lemma 1 becomes,

$$\alpha_n = a_n + \frac{1}{2} a'_n h + \frac{1}{6} a''_n h^2 + O(h^3), \\ = a_n + \frac{1}{4} (3a_n - 4a_{n-1} + a_{n-2}) + \frac{1}{6} (a_n - 2a_{n-1} + a_{n-2}) + O(h^3) \\ = \frac{1}{12} (23a_n - 16a_{n-1} + 5a_{n-2}) + O(h^3).$$

We retrieve the definition of α_n for the RL₃ scheme. The same proof holds for β_n and extends at the other 4. $\square$

2 Stability under perturbation and convergence

We refer to [9, Ch. III-8] for the definitions of convergence and of stability under perturbation. For the analysis of time-stepping methods, it is commonly assumed that f in (2) is uniformly Lipschitz in its second variable y . That hypothesis will be replaced by assumptions based on the formulation (3). Precisely it will be assumed that,

$$a(t, y) \text{ is bounded, } a(t, y), b(t, y) \text{ are uniformly Lipschitz in } y. \quad (7)$$

The Lipschitz constants for a and b are denoted L_a and L_b respectively. The upper bound on $|a(t, y)|$ is denoted M_a .

Proposition 2. With the assumptions (7) the RL_k scheme is stable under perturbation.

Corollary 1. Assume that $a(t, y)$ and $b(t, y)$ are C^k regular and that $a(t, y)$ either is a diagonal or a constant matrix. In addition assume (7). Then the RL_k scheme is convergent of order k .

Stability under perturbation together with consistency implies convergence, see *e.g.* [9] or [6] where the current setting has been detailed. Therefore Corollary 1 is an immediate consequence of Propositions 1 and 2. Before to prove Proposition 2, definitions are needed.

Equation (2) is considered on $E = \mathbb{R}^N$ with the max norm $|\cdot|$. A final time $T > 0$ is considered. The space of $N \times N$ matrices is equipped with the operator norm $\|\cdot\|$ associated to $|\cdot|$. The space E^k is equipped with the max norm $|Y|_\infty = \max_{1 \leq i \leq k} |y_i|$ with $Y = (y_1, \dots, y_k)$. The RL_k scheme is defined with the mapping

$$s_{t,h} : Y = (y_1, \dots, y_k) \in E^k \longrightarrow s_{t,h}(Y) \in E,$$

with

$$s_{t,h} = y_k + h\varphi_1(\alpha_{t,h}(Y)h)(\alpha_{t,h}(Y)y_k + \beta_{t,h}(Y)),$$

in such a way that $y_{n+1} = s_{t,h}(y_{n-k+1}, \dots, y_n)$ in (6). The functions $\alpha_{t,h}$ and $\beta_{t,h}$ are given in Definition 1. For instance, $\alpha_{t,h}(Y)$ for the RL_3 scheme reads:

$$\alpha_{t,h}(Y) = \frac{1}{12}(23a(t, y_3) - 16a(t - h, y_2) + 5a(t - 2h, y_1)), \quad Y = (y_1, y_2, y_3).$$

A first way to prove the stability under perturbation is to show that $s_{t,h}$ is globally Lipschitz in Y . For this the derivative $\partial_Y s_{t,h}$ has to be analyzed. As developed in Remark 2, this will imply restrictions on $a(t, y)$: either diagonal or constant. A second way is to prove the two following stability conditions,

$$|s_{t,h}(Y) - s_{t,h}(Z)| \leq |Y - Z|_\infty (1 + Ch(|Y|_\infty + 1)), \quad (8)$$

$$|s_{t,h}(Y)| \leq |Y|_\infty (1 + Ch) + Ch \quad (9)$$

where C is a constant only depending on the data a, b, y_0 in Equation (3) and on the final time T . These are sufficient conditions for the stability under perturbation, as proved in

[6].

That second way will be used here, for it is more general and giving rise to less computations. The core of the proof is the following property of the RL_k scheme. For $Y = (y_1, \dots, y_k) \in E^k$:

$$s_{t,h}(Y) = z(t+h) \quad \text{for} \quad z' = \alpha_{t,h}(Y)z + \beta_{t,h}(Y), \quad z(t) = y_k. \quad (10)$$

It will be used together with the following Gronwall inequality (see [8, Lemma 196, p.150]). Suppose that $z(t)$ is a C^1 function and that there exist $M_1, M_2 > 0$ such that $|z'(t)| \leq M_1(t - t_0) + M_2$ for all $t \in [t_0, t_0 + h]$. Then,

$$\forall t \in [t_0, t_0 + h], \quad |z(t)| \leq e^{M_1(t-t_0)} (|z(t_0)| + M_2(t - t_0)). \quad (11)$$

Proof of Proposition 2. In this proof it is always assumed that $0 \leq h, t \leq T$. We will denote by C_i a constant only depending on the problem (3) data a and b and on T .

With the assumptions (7) and Definition 1, the function $\alpha_{t,h}$ is uniformly Lipschitz with a Lipschitz constant L_α . Moreover we have a uniform bound $\|\alpha_{t,h}\| \leq M_\alpha$. Since $b(t, y)$ is uniformly Lipschitz in y and since $0 \leq t \leq T$, there exists a constant K_b so that,

$$|b(t, y)| \leq K_b(1 + |y|). \quad (12)$$

Consider the RL_3 scheme,

$$\begin{aligned} |\beta_{t,h}(Y)|_\infty &\leq \frac{11}{3}K_b(1 + |Y|_\infty) + \frac{h}{12}M_a2K_b(1 + |Y|_\infty) \\ &\leq C_1(1 + |Y|_\infty). \end{aligned}$$

The same inequality holds for the RL_2 and RL_4 schemes. With (10) we have $s_{t,h}(Y) = z(t+h)$ and,

$$|z'| = |\alpha_{t,h}(Y)z + \beta_{t,h}(Y)| \leq M_\alpha|z| + C_1(1 + |Y|_\infty).$$

The initial state is $|z(t)| = |y_k| \leq |Y|_\infty$. With the Gronwall inequality (11) we obtain for $t \leq \tau \leq t + h$,

$$\begin{aligned} |z(\tau)| &\leq e^{M_\alpha h} (|Y|_\infty + hC_1(1 + |Y|_\infty)) \\ &\leq e^{M_\alpha h} (|Y|_\infty(1 + C_1h) + C_1h) \\ &\leq |Y|_\infty(1 + C_2h) + C_2h, \end{aligned} \quad (13)$$

by bounding the exponential with an affine function for $0 \leq h \leq T$. This gives the stability condition (9) for $\tau = t + h$.

For the RL_2 scheme $\beta_{t,h}$ is uniformly Lipschitz.

For the RL_3 scheme, consider $Y = (y_1, y_2, y_3)$ and $Z = (z_1, z_2, z_3)$ in E^3 . We have,

$$\begin{aligned} |\beta_{t,h}(Y) - \beta_{t,h}(Z)|_\infty &\leq \frac{11}{3}L_b|Y - Z|_\infty + \frac{h}{12}(|a(t, y_3)b(t-h, y_2) - a(t, z_3)b(t-h, z_2)| \\ &\quad + |a(t-h, y_2)b(t, y_3) - a(t-h, z_2)b(t, z_3)|) \end{aligned}$$

Let us bound the Lipschitz constant for a function of the type $F(Y) = a(\xi, y_2)b(\tau, y_3)$ for $0 \leq \tau, \xi \leq T$.

$$\begin{aligned} |F(Y) - F(Z)| &= |a(\xi, y_3)(b(\tau, y_2) - b(\tau, z_2)) + (a(\xi, y_3) - a(\xi, z_3))b(\tau, z_2)| \\ &\leq M_a L_b |Y - Z|_\infty + L_a |Y - Z|_\infty |b(\tau, z_2)|. \end{aligned}$$

With (12), this yields for $0 \leq \tau, \xi \leq T$ and for $Y, Z \in E^k$,

$$|F(Y) - F(Z)| \leq C_3 |Y - Z|_\infty (1 + |Z|_\infty),$$

As a result,

$$|\beta_{t,h}(Y) - \beta_{t,h}(Z)|_\infty \leq C_4 |Y - Z|_\infty (1 + |Z|_\infty)$$

The same inequality holds for the RL_4 scheme.

Finally consider $Y_1, Y_2 \in E^k$ and denote $\alpha_i = \alpha_{t,h}(Y_i)$, $\beta_i = \beta_{t,h}(Y_i)$. With the property (10), $s_{t,h}(Y_1) - s_{t,h}(Y_2) = (z_1 - z_2)(t + h)$ where z_i is the solution to,

$$z'_i = \alpha_i z_i + \beta_i, \quad z_i(t) = Y_{i,k}.$$

On the first hand, with the inequality (13), we have $|z_2(\tau)| \leq C_5(1 + |Y_2|_\infty)$ for $t \leq \tau \leq t + h$.

On the second hand, on $[t, t + h]$,

$$\begin{aligned} |(z_1 - z_2)'| &\leq |\alpha_1||z_1 - z_2| + |\alpha_1 - \alpha_2||z_2| + |\beta_1 - \beta_2| \\ &\leq M_\alpha |z_1 - z_2| + L_\alpha |Y_1 - Y_2|_\infty C_5(1 + |Y_2|_\infty) + C_4 |Y_1 - Y_2|_\infty (1 + |Y_2|_\infty) \\ &\leq M_\alpha |z_1 - z_2| + C_6 |Y_1 - Y_2|_\infty (1 + |Y_2|_\infty) \end{aligned}$$

The initial condition is $|(z_1 - z_2)(t)| = |Y_{1,k} - Y_{2,k}| \leq |Y_1 - Y_2|_\infty$. Then the Gronwall inequality (11) yields,

$$\begin{aligned} |(z_1 - z_2)(t + h)| &\leq e^{M_\alpha h} (|Y_1 - Y_2|_\infty + h C_6 |Y_1 - Y_2|_\infty (1 + |Y_2|_\infty)) \\ &\leq e^{M_\alpha h} |Y_1 - Y_2|_\infty (1 + C_6 h (1 + |Y_2|_\infty)). \end{aligned}$$

This last inequality implies the stability condition (8), again by bounding the exponential with an affine function for $0 \leq h \leq T$. $\square$

3 Dahlquist stability

For the general definition concerning the Dahlquist stability we refer to [10]. The background for the Dahlquist stability of exponential integrators with a general varying stabilizer $a(t, y)$ has been developed in [6], following the ideas of Perego and Veneziani [24]. Problem (2) is considered with the Dahlquist test function $f(t, y) = \lambda y$ that is decomposed in (3) as $f(t, y) = a(t, y)y + b(t, y)$ with,

$$a(t, y) = \theta \lambda, \quad b(t, y) = \lambda(1 - \theta)y.$$

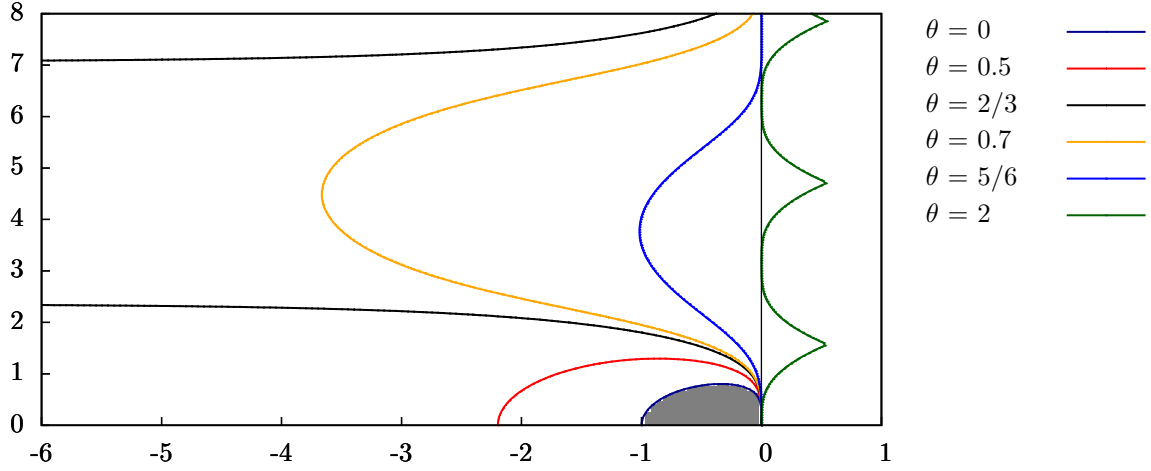


Figure 1: Stability domain D_θ for the RL_2 scheme for various values of θ . The stability domain for the particular case $\theta = 0$ (no stabilization) is in grey, corresponding to the Adams-Bashforth scheme of order 2.

When $\theta = 1$, the method is exact and thus A -stable.

When $\theta \simeq 1$, the exact linear part of $f(t, y)$ in Equation (2) is well approximated by $a(t, y)$. The stability domain depends on θ . It is denoted D_θ . At a fixed value of θ , it is given by the modulus of a stability function, with the same definition as for multistep methods, see *e.g.* [10]. That stability function has been numerically computed pointwise on a grid inside the complex plane $\mathbb{C}$.

Order 2 Rush-Larsen

The stability domain for the RL_2 scheme has been analyzed in [24]. The situation for this scheme is interesting and we reproduced the results on Figure 1.

- If $0 \leq \theta < 2/3$ the stability domain D_θ is bounded. Its size increases with θ , starting from the Adams-Bashforth scheme of order 2 stability domain when no stabilization occurs ($\theta = 0$).
- If $\theta = 2/3$, D_θ contains the negative real axis: the method is $A(0)$ stable. The domain boundary is asymptotically parallel to the real axis so that the method is not $A(\alpha)$ stable.
- If $\theta > 2/3$, the stability domain is located around the y -axis: the method is $A(\alpha)$ stable. The angle α increases with θ , it goes to $\pi/2$ as $\theta \rightarrow 1^-$.

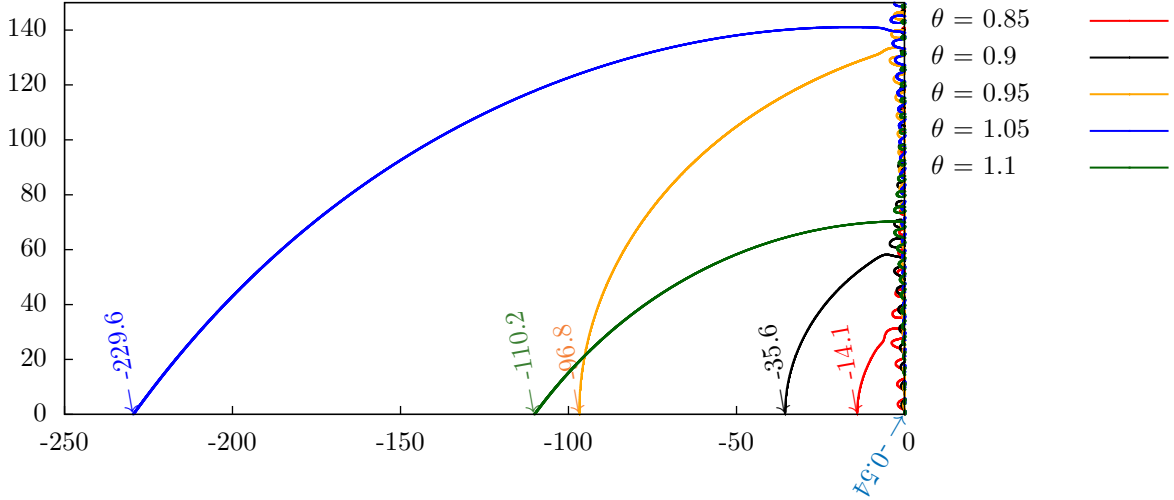


Figure 2: Stability domain D_θ for the RL_3 scheme. In the particular case $\theta = 0$ (no stabilization, corresponding to the Adams-Bashforth scheme of order 3), the stability domain crosses the x -axis at $x \simeq -0.54$ (dark blue arrow).

Rush-Larsen of order 3 and 4

The situation is different for the Rush-Larsen methods of order 3 and 4. The stability domains D_θ for various values of θ have been numerically computed and depicted on Figures 2 and 3.

Excepted for the case $\theta = 1$, the stability domains are always bounded: the schemes are not A -stable. However, this stability domain for values of $\theta \simeq 1$ are much larger than the $D_{\theta|_{\theta=0}}$ stability domain when no stabilization occurs (corresponding to the Adams-Bashforth scheme of order 3 or 4). For the order 3 case, the stability domain for $\theta = 0.85$ is 25 times wider on the left than $D_{\theta|_{\theta=0}}$, and for $\theta = 1.05$ it is 400 times wider. For the order 4 case, $D_{\theta|_{\theta=1.05}}$ is almost 300 times wider on the left than $D_{\theta|_{\theta=0}}$.

4 Numerical results

In this section are presented numerical experiments in order to investigate the performances of the RL_k method. It will be compared to the exponential integrator of Adams type of order k , shortly denoted EAB_k , defined by Equation (4). The EAB_k schemes have been numerically studied in [6] as compared to several classical methods. It had been shown to be a good candidate for the resolution of the stiff membrane equation in cardiac electrophysiology.

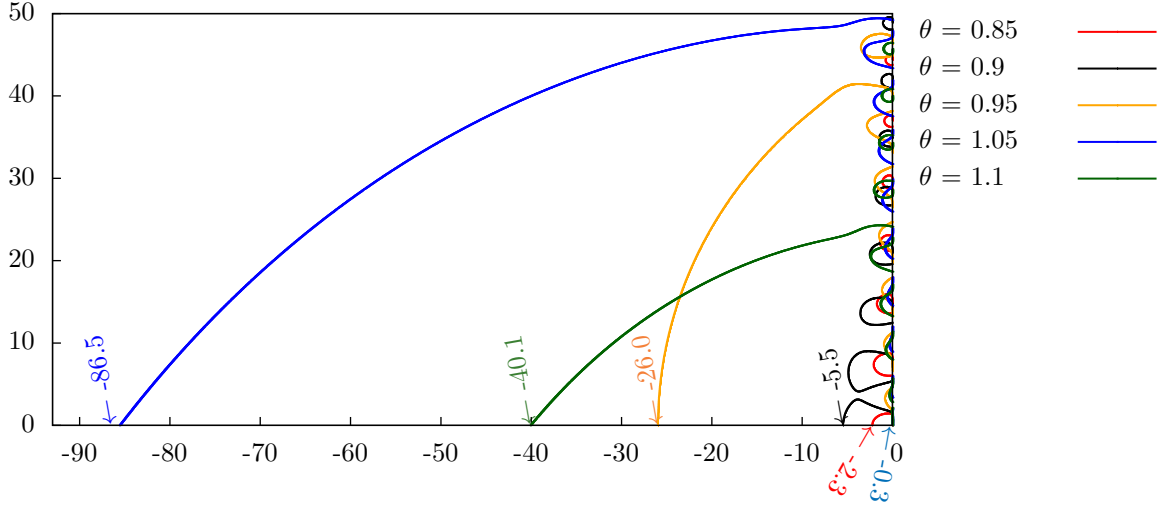


Figure 3: Stability domain D_θ for the RL_4 scheme. In the particular case $\theta = 0$ (no stabilization, corresponding to the Adams-Bashforth scheme of order 3), the stability domain crosses the x -axis at $x \simeq -0.3$ (dark blue arrow).

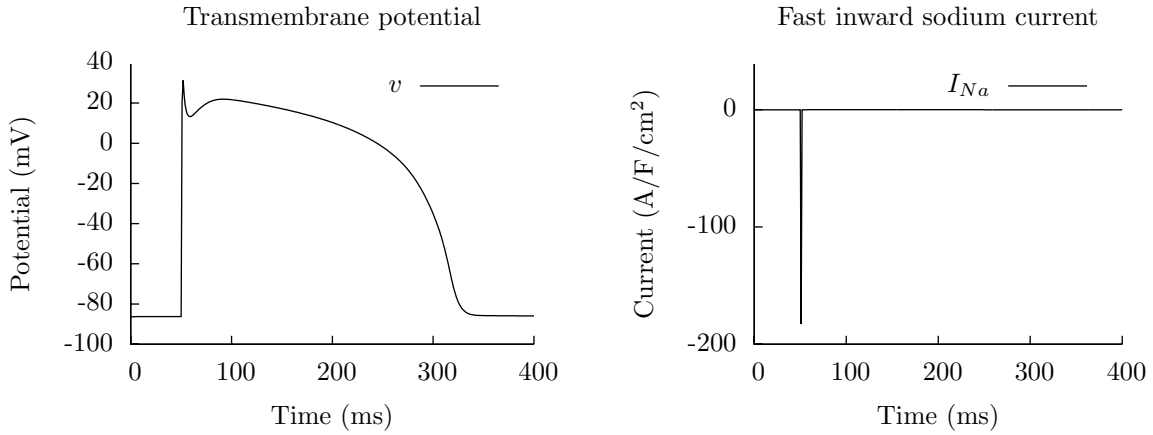


Figure 4: TNNP model illustration. Left, cellular action-potential: starting at a (negative) rest value, the membrane potential $v(t)$ has a stiff depolarization followed by a plateau and a repolarization to the rest value. Right, depolarization is induced by an ionic sodium current I_{Na} , with obvious large stiffness.

4.1 The membrane equation

We consider a class of models in cardiac electrophysiology. As illustrated on Figure 4, these models display a stiff behaviour characterized by the presence of heterogeneous time scales. The models used to simulate the electrical activity of cardiac cells are ODE

systems of the form, see [17, 1, 21, 27]:

$$\begin{aligned}\frac{dw_i}{dt} &= \frac{w_{\infty,i}(v) - w_i}{\tau_i(v)}, & \frac{dc}{dt} &= g(w, c, v), \\ \frac{dv}{dt} &= -I_{ion}(w, c, v) + I_{st}(t),\end{aligned}\tag{14}$$

where $w = (w_1, \dots, w_p) \in \mathbb{R}^p$ is the vector of the gating variables, $c \in \mathbb{R}^q$ is a vector of ionic concentrations or other state variables, and $v \in \mathbb{R}$ is the transmembrane-potential. The four functions $w_{\infty,i}(v)$, $\tau_i(v)$, $g(w, c, v)$ and $I_{ion}(w, c, v)$ are given reaction terms. The function $I_{st}(t)$ is a source term. It represents a stimulation current. Problem (14) reformulates into problem (3) form with:

$$a(t, y) = \begin{pmatrix} -1/\tau(v) & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix}, \quad b(t, y) = \begin{pmatrix} w_{\infty}(v)/\tau(v) \\ g(y) \\ -I_{ion}(y) + I_{st}(t) \end{pmatrix}, \tag{15}$$

for $y = (w, c, v) \in \mathbb{R}^N$ ($N = p + q + 1$) and where $-1/\tau(v)$ the $p \times p$ diagonal matrix with diagonal entries $(-1/\tau_i(v))_{i=1\dots p}$. The resulting matrix $a(t, y)$ is diagonal.

We will consider two such models: the BR model (Beeler and Reuter [1]) and the TNNP model [27] for human cardiac cells.

4.2 Convergence

No theoretical solution are available for the chosen application. A reference solution y_{ref} for a reference time step h_{ref} is computed with the Runge-Kutta scheme of order 4 to analyze the convergence properties of the RL_k scheme. Numerical solutions y are computed to y_{ref} for coarsest time steps $h = 2^p h_{ref}$ for increasing p .

Any numerical solution y consists in successive values y_n at the time instants $t_n = nh$. On every interval (t_{3n}, t_{3n+3}) the polynomial $\bar{y}$ of degree at most 3 so that $\bar{y}(t_{3n+i}) = y_{3n+i}$, $i = 0 \dots 4$ is constructed. On $(0, T)$, $\bar{y}$ is continuous and piecewise polynomial of degree 3. Its values at the reference time instants nh_{ref} are computed. This provides a projection $P(y)$ of the numerical solution y on the reference grid. Then $P(y)$ can be compared with the reference solution y_{ref} . The numerical error is defined by,

$$e(h) = \frac{\max |v_{ref} - P(v)|}{\max |v_{ref}|}, \tag{16}$$

where the potential v is the last and stiffest component of y in Equation (14). The numerical convergence graphs for the BR model are plotted on Figure 5. All the schemes display the expected asymptotic behaviour $e(h) = O(h^k)$ as $h \rightarrow 0$ in Corollary 1.

4.3 Stability

In [26] has been evaluated the stiffness of the BR and TNNP models along one cellular electrical cycle (as depicted on Figure 4). The largest negative real part of the eigenvalues

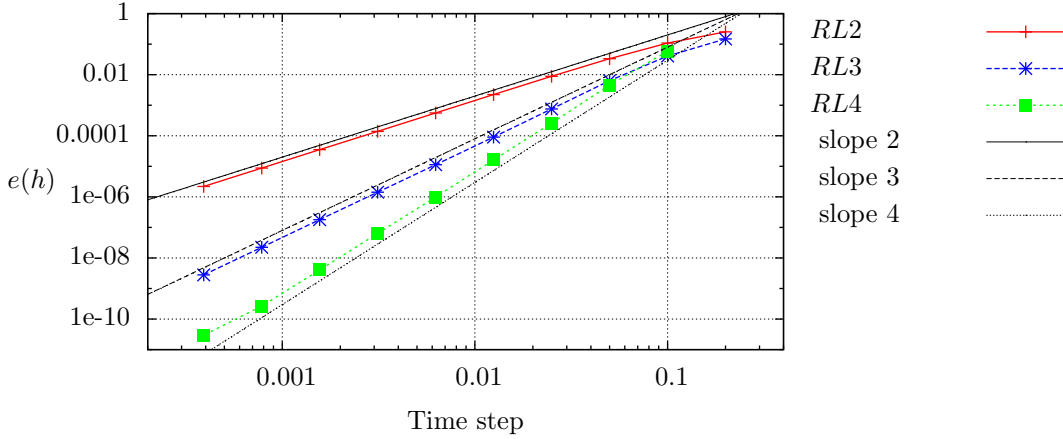


Figure 5: Relative error $e(h)$ (definition (16)) as a function of the time step h for the RL_k schemes, for $k = 2$ to 4 and in Log/Log scale.

of the Jacobian matrix during this cycle is of -1170 and -82 for the TNNP and BR models respectively. The TNNP model thus is 15 times stiffer than the BR model ($15 \simeq 1170/82$). Robustness to stiffness for the RL_k scheme is evaluated by comparing the critical time step for these two models. The critical time step Δt_0 is defined as the largest time step such that the numerical simulation runs without overflow for $h < \Delta t_0$. The results are presented in Table 1.

Table 1: Critical time step Δt_0 for the RL_k and EAB_k schemes

method	RL_2	RL_3	RL_4	EAB_2	EAB_3	EAB_4
BR	0.323	0.200	0.149	0.424	0.203	0.123
TNNP	0.120	0.148	0.111	0.233	0.108	$7.56 \cdot 10^{-2}$

An excellent robustness to stiffness can be observed. The critical time step is divided by 2.7, 2.0 and 1.3 for $k = 2, 3$ and 4 respectively. A comparison with the EAB_k schemes shows that the two scheme display a robustness to stiffness of same order.

For a method that is not $A(\alpha)$ stable, it is expected for the critical time step to be divided by 15 in case of an increase of stiffness of magnitude 15. This is not observed here, though the RL_k scheme is not $A(\alpha)$ stable. The reason for this is that the ODE system (14) is only partially stabilized by (15). Loss of stability is induced by the non-stabilized part, whose eigenvalues are less modified between the BR and the TNNP models.

4.4 Accuracy

The RL_k scheme is here compared to the EAB_k scheme in terms of accuracy. This comparison is done by computing the relative error $e(h)$ in Equation (16). The two BR and TNNP models are considered. We recall than the TNNP model is stiffer by a factor 15. The results are collected in Tables 2 and 3.

Table 2: Relative error $e(h)$ (eq. (16)) for the BR model.

h	RL ₂	RL ₃	RL ₄	EAB ₂	EAB ₃	EAB ₄
0.2	0.251	0.147	-	0.284	0.516	-
0.1	0.107	$4.07 \cdot 10^{-2}$	$5.86 \cdot 10^{-2}$	$9.26 \cdot 10^{-2}$	$9.17 \cdot 10^{-2}$	0.119
0.05	$3.35 \cdot 10^{-2}$	$6.34 \cdot 10^{-3}$	$4.58 \cdot 10^{-3}$	$2.31 \cdot 10^{-2}$	$1.09 \cdot 10^{-2}$	$8.96 \cdot 10^{-3}$
0.025	$8.88 \cdot 10^{-3}$	$7.57 \cdot 10^{-4}$	$2.61 \cdot 10^{-4}$	$5.39 \cdot 10^{-3}$	$1.17 \cdot 10^{-3}$	$4.33 \cdot 10^{-4}$

Table 3: Relative error $e(h)$ (eq. (16)) for the TNNP model.

h	RL ₂	RL ₃	RL ₄	EAB ₂	EAB ₃	EAB ₄
0.1	0.177	0.305	0.421	0.351	0.530	-
0.05	$7.39 \cdot 10^{-2}$	$4.54 \cdot 10^{-2}$	$4.61 \cdot 10^{-2}$	$9.01 \cdot 10^{-2}$	$5.59 \cdot 10^{-2}$	$8.93 \cdot 10^{-2}$
0.025	$2.21 \cdot 10^{-2}$	$6.53 \cdot 10^{-3}$	$5.96 \cdot 10^{-3}$	$2.14 \cdot 10^{-2}$	$7.34 \cdot 10^{-3}$	$8.34 \cdot 10^{-3}$
0.0125	$5.75 \cdot 10^{-3}$	$8.05 \cdot 10^{-4}$	$3.21 \cdot 10^{-4}$	$5.11 \cdot 10^{-3}$	$7.62 \cdot 10^{-4}$	$3.70 \cdot 10^{-4}$

For the RL₂ and the EAB₂ schemes, the accuracies are very close, the EAB₂ scheme being slightly more accurate for the BR model. For the orders 3 and 4, a non negligible difference is observed between the RL and EAB schemes. The RL scheme is more accurate at large time steps. For smaller time steps, accuracies are almost the same. This means that RL and EAB schemes are equivalent in terms of accuracy considering the asymptotic convergence region, but outside this region, RL scheme is more precise.

Conclusion

We introduced in this paper two new ODE solvers that we named Rush-Larsen of order 3 and 4. They are explicit multistep exponential integrators. Their general Definition (6) is very simple inducing an easy implementation. We provided a convergence and stability under perturbation analysis of these two schemes. We also performed their Dahlquist stability analysis: they are not $A(0)$ stable but display a very large stability domain for sufficiently precise stabilization. The numerical properties of the schemes are analyzed for a complex and realistic stiff application. The RL _{k} schemes are as stable as exponential integrators of Adams type, allowing simulations at large time step. On the presented example, the RL _{k} schemes moreover are more accurate for $k = 3$ and 4 when considering large time steps. They are shown to be robust to stiffness both in terms of stability and of accuracy.

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