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# Compositions of pseudo-symmetric integrators with complex coefficients for the numerical integration of differential equations



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#### ABSTRACT

In this paper, we are concerned with the construction and analysis of a new class of methods obtained as double jump compositions with complex coefficients and projection on the real axis. It is shown in particular that the new integrators are symmetric and symplectic up to high orders if one uses a symmetric and symplectic basic method. In terms of efficiency, the aforementioned technique requires fewer stages than standard compositions of the same orders and is thus expected to lead to faster methods.

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# 1. Introduction

Given a differential equation

$$\dot{x} \equiv \frac{dx}{dt} = f(x), \qquad x(0) = x_0, \tag{1}$$

composition methods constitute a powerful technique to raise the order of a given integrator  $\psi_{\tau}$  applied to (1) with time-step  $\tau$ , as high as might be required, by considering expressions of the form

$$\phi_{\tau} = \psi_{\gamma_1 \tau} \circ \psi_{\gamma_2 \tau} \circ \cdots \circ \psi_{\gamma_5 \tau}, \tag{2}$$

where the coefficients  $\gamma_1, \gamma_2, \dots, \gamma_s$  are appropriately chosen so as to satisfy some *universal algebraic conditions* [1–3]. It is known in particular that if  $\psi_{\tau}$  is of order k, i.e. satisfies

$$\varphi_{\tau}(x_0) - \psi_{\tau}(x_0) = \mathcal{O}(\tau^{k+1}),$$

where  $\varphi_{\tau}$  denotes the exact flow of (1), then  $\phi_{\tau}$  will be at least of order k+1 (i.e., local error k+2) if the following two conditions are satisfied

(i) 
$$\sum_{i=1}^{s} \gamma_i = 1$$
 and (ii)  $\sum_{i=1}^{s} \gamma_i^{k+1} = 0$ . (3)

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Given that these two equations have no real solution for odd k and arbitrary s, a series of authors (e.g. [4,5]) suggested to start from a second-order method and to consider *symmetric compositions* only, i.e., schemes with coefficients satisfying the additional condition

$$\gamma_{s+1-i} = \gamma_i, \quad i = 1, \ldots, s.$$

This has led to so-called triple-jump compositions (s = 3,  $\gamma_3 = \gamma_1$ ) obtained by iterating the process described above to construct a sequence of symmetric methods with even orders (see, e.g., [1, pp. 44]).

In spite of its simplicity, the triple-jump rationale leads to inefficiencies for high orders as compared to methods obtained by solving directly the order conditions [1]. On top of this, it also suffers from the occurrence of negative time-steps, although this fact is not specific to triple-jump methods and concerns all composition or splitting methods of orders higher than two. This, of course, is a severe limiting factor for equations where the vector field (usually an operator) is not reversible, the prototypical example of which being the heat equation. To circumvent this difficulty, several authors have suggested to use complex time-steps (or complex coefficients) in the context of parabolic equations [6,7]. One indeed easily sees that, already for s = 2, solutions of equations (i)–(ii) exist in  $\mathbb{C}$ .

Generally speaking, suppose that  $\psi_{\tau}$  is an integrator of order k, denoted  $\mathcal{S}_{\tau}^{[k]}$  in the sequel for clarity, and consider the composition (2) with s=2,

$$\mathcal{S}_{\tau}^{[k+1]} = \mathcal{S}_{\gamma_1 \tau}^{[k]} \circ \mathcal{S}_{\gamma_2 \tau}^{[k]}. \tag{4}$$

Then, if the coefficients verify conditions (i)-(ii), that is to say if

$$\gamma_{1} = \bar{\gamma}_{2} \equiv \gamma = \frac{1}{2} + \frac{i}{2} \frac{\sin(\frac{2\ell+1}{k+1}\pi)}{1 + \cos(\frac{2\ell+1}{k+1}\pi)} \quad \text{for} \quad \begin{cases} -\frac{k}{2} \le \ell \le \frac{k}{2} - 1 & \text{if } k \text{ is even} \\ -\frac{k+1}{2} \le \ell \le \frac{k-1}{2} & \text{if } k \text{ is odd} \end{cases}$$
(5)

then (4) results in a method of order k+1, which can subsequently be used to generate recursively higher order composition schemes by applying the same procedure. The choice  $\ell=0$ ,

$$\gamma = \gamma^{[k]} := \frac{1}{2} + \frac{i}{2} \frac{\sin\left(\frac{\pi}{k+1}\right)}{1 + \cos\left(\frac{\pi}{k+1}\right)} = \frac{1}{2} + \frac{i}{2} \tan\left(\frac{\pi}{2(k+1)}\right) = \frac{1}{2\cos\left(\frac{\pi}{2(k+1)}\right)} e^{\frac{\pi}{2(k+1)}}$$
(6)

gives the solutions with the smallest phase. If we start with a symmetric method of order 2,  $S_{\tau}^{[2]}$ , and apply composition (4) with corresponding coefficients (6), we can construct the following sequence of methods:

$$\mathcal{S}^{[2]}_{\tau} \longrightarrow \mathcal{S}^{[3]}_{\tau} \longrightarrow \mathcal{S}^{[4]}_{\tau} \longrightarrow \mathcal{S}^{[5]}_{\tau} \longrightarrow \mathcal{S}^{[6]}_{\tau},$$

all of which have coefficients with positive real part [7]. The final method of order 6 involves 16 evaluations of the basic scheme  $\mathcal{S}^{[2]}_{\tau}$ . By contrast, there are composition methods of order 6 (both with real and complex coefficients) involving just 7 evaluations of  $\mathcal{S}^{[2]}_{\tau}$  [5,8]. It is thus apparent that this direct approach does not lead to cost-efficient high-order schemes

One should remark that the composition (4) does not provide a time-symmetric method, i.e.,  $S_{-\tau}^{[k+1]} \circ S_{\tau}^{[k+1]}$  is *not* the identity map, even if  $S_{\tau}^{[k]}$  happens to be symmetric. As we have mentioned before, symmetry allows to raise the order by two at each iteration by considering the triple-jump composition

$$S_{\tau}^{[2k+2]} = S_{\gamma_1 \tau}^{[2k]} \circ S_{\gamma_2 \tau}^{[2k]} \circ S_{\gamma_1 \tau}^{[2k]}$$
(7)

starting from a symmetric method. Apart from the real solution, the complex one with the smallest phase is

$$\gamma_1 = \frac{e^{i\pi/(k+1)}}{2^{1/(k+1)} - 2e^{i\pi/(k+1)}}, \qquad \gamma_2 = 1 - 2\gamma_1, \tag{8}$$

and symmetric methods up to order 8 with coefficients having positive real part are possible if one starts with a symmetric second-order scheme.<sup>1</sup> These order barriers have been rigorously proved in [8].

The simple third-order scheme (4) corresponding to k=2 has been in fact rediscovered several times in the literature [6,7,9–11]. In particular, it was shown in [10] that the method, when applied to the two-body Kepler problem, behaves indeed as a fourth-order integrator, the reason being attributed to the fact that the main error term in the asymptotic expansion is purely imaginary. In this note we elaborate further the analysis and provide a comprehensive study of the general composition (4), paying special attention to the qualitative properties the method shares with the continuous system (1). In addition, we show how it is possible to combine compositions and a trivial linear combination to raise the order, while still preserving the qualitative properties of the basic integrator up to an order higher than of the method itself.

<sup>&</sup>lt;sup>1</sup> It is actually possible to reach order 14 if, in the construction, one uses formula (7) alternatively with coefficients  $\gamma_1$ ,  $\gamma_2$  and coefficients  $\bar{\gamma}_1$ ,  $\bar{\gamma}_2$  [6].

# 2. Composition and pseudo-symmetry or pseudo-symplecticity

In what follows, we will assume for convenience that all values of x in (1) lie in a compact set K where the function f is *smooth*. Before starting the analysis, it is worth recalling the notions of adjoint method and symplectic flow.

The adjoint method  $\psi_{\tau}^*$  of a given method is the inverse map of the original integrator with reversed time step  $-\tau$ , i.e.,  $\psi_{\tau}^* := \psi_{-\tau}^{-1}$ . A symmetric method satisfies  $\psi_{\tau}^* = \psi_{\tau}$  [1,12].

The vector field f in (1) is Hamiltonian if there exists a function H(x) such that  $f = J\nabla_x H(x)$ , where  $x = (q, p)^T$  and J is the basic canonical matrix. Then, the exact flow of (1) is a symplectic transformation,  $\varphi'_t(x)^T J \varphi'_t(x) = J$  for  $t \ge 0$  [13,14]. It then makes sense introducing the following definitions, taken from [15] and [16]:

#### **Definition 1.** Let $\psi_{\tau}$ be a smooth and consistent integrator:

1. it is pseudo-symmetric of pseudo-symmetry order q if for all sufficiently small  $\tau$ , the following relation holds true:

$$\psi_{\tau}^* = \psi_{\tau} + \mathcal{O}(\tau^{q+1}),\tag{9}$$

where the constant in the  $\mathcal{O}$ -term depends on bounds of derivatives of  $\psi$  on K.

2. it is *pseudo-symplectic* of *pseudo-symplecticity order* r if for all sufficiently small  $\tau$ , the following relation holds true when  $\psi_{\tau}$  is applied to a Hamiltonian system:

$$(\psi_{\tau}')^T J \psi_{\tau}' = J + \mathcal{O}(\tau^{r+1}), \tag{10}$$

where the constant in the  $\mathcal{O}$ -term depends on bounds of derivatives of  $\psi$  on K.

**Remark 1.** A symmetric method is pseudo-symmetric of any order  $q \in \mathbb{N}$ , whereas a method of order k is pseudo-symmetric of order  $q \ge k$ . A similar statement holds for symplectic methods.

As a first illustration of Definition 1, let us consider again a symmetric 2nd-order method  $S_r^{[2]}$  and form the composition

$$\psi_{\tau}^{[3]} = \mathcal{S}_{\nu\tau}^{[2]} \circ \mathcal{S}_{\bar{\nu}\tau}^{[2]}$$

with  $\gamma = \frac{1}{2} + i\frac{\sqrt{3}}{6}$ . Then, if the vector field f under consideration is real-valued, its real part

$$\Re(\psi_{\tau}^{[3]}) = \frac{1}{2} \left( \psi_{\tau}^{[3]} + \overline{\psi}_{\tau}^{[3]} \right) = \frac{1}{2} \left( \mathcal{S}_{\gamma\tau}^{[2]} \circ \mathcal{S}_{\bar{\gamma}\tau}^{[2]} + \mathcal{S}_{\bar{\gamma}\tau}^{[2]} \circ \mathcal{S}_{\gamma\tau}^{[2]} \right).$$

is a method of order 4 and pseudo-symmetric of pseudo-symmetry order 7. This result is a consequence of the fact that

$$(\psi_{\tau}^{[3]})^* = \overline{\psi}_{\tau}^{[3]}$$

and the following general statement, which lies at the core of the construction procedure described in this paper.

**Proposition 1.** Let  $\psi_{\tau}$  be any consistent smooth method for Eq. (1) and consider the new method

$$R_{\tau} = \frac{1}{2} \left( \psi_{\tau} + \psi_{\tau}^* \right).$$

Assume also that  $\psi_{\tau}$  is pseudo-symmetric of order q. Then  $R_{\tau}$  is of pseudo-symmetry order 2q+1. If  $\psi_{\tau}$  is furthermore of pseudo-symplecticity order r, then  $R_{\tau}$  is of pseudo-symplecticity order  $\min(2q+1,r)$ .

**Proof.** By assumption, there exists a smooth function  $(\tau, x) \mapsto \delta_{\tau}(x)$ , defined for all x in a compact set K and for all sufficiently small real  $\tau$ , such that

$$\psi_{\tau}^{*} = \psi_{\tau} + \tau^{q+1} \delta_{\tau} \quad \text{or} \quad \psi_{-\tau}^{-1} = \psi_{\tau} + \tau^{q+1} \delta_{\tau} \quad \text{or} \quad \psi_{\tau}^{-1} = \psi_{-\tau} + (-\tau)^{q+1} \delta_{-\tau}, \tag{11}$$

so that

$$R_{\tau} = \psi_{\tau} + \frac{1}{2} \tau^{q+1} \delta_{\tau}.$$

Composing the third relation of (11) from the left by  $\psi_{\tau}$ , we obtain

$$id = \psi_{\tau} \circ \psi_{-\tau} + (-\tau)^{q+1} \psi_{\tau}' \circ \psi_{-\tau} \cdot \delta_{-\tau} + \mathcal{O}(\tau^{2(q+1)}), \tag{12}$$

where the  $\mathcal{O}$ -term depends on bounds of the derivatives of  $\psi_{\tau}$  and  $\delta_{\tau}$  on K. Similarly, composing the second relation of (11) from the right by  $\psi_{-\tau}$ , we get

$$id = \psi_{\tau} \circ \psi_{-\tau} + \tau^{q+1} \delta_{\tau} \circ \psi_{-\tau}. \tag{13}$$

As a consequence, we have

$$\tau^{q+1}\delta_{\tau}\circ\psi_{-\tau}=(-\tau)^{q+1}\psi_{\tau}'\circ\psi_{-\tau}\cdot\delta_{-\tau}+\mathcal{O}(\tau^{2(q+1)}).$$

We are then in position to write

$$\begin{split} R_{\tau} \circ R_{-\tau} &= \left( \psi_{\tau} + \frac{1}{2} \tau^{q+1} \delta_{\tau} \right) \circ \left( \psi_{-\tau} + \frac{1}{2} (-\tau)^{q+1} \delta_{-\tau} \right) \\ &= \psi_{\tau} \circ \psi_{-\tau} + \frac{1}{2} (-\tau)^{q+1} \psi_{\tau}' \circ \psi_{-\tau} \cdot \delta_{-\tau} + \frac{1}{2} \tau^{q+1} \delta_{\tau} \circ \psi_{-\tau} + \mathcal{O}(\tau^{2(q+1)}) \\ &= \mathrm{id} + \mathcal{O}(\tau^{2(q+1)}), \end{split}$$

which proves the first statement. Now, if  $\psi_{\tau}$  is in addition of pseudo-symplecticity order r, then its adjoint  $\psi_{\tau}^*$  is also of pseudo-symplecticity order r, so that relation (11) leads to

$$J + \mathcal{O}(\tau^{r+1}) = (\partial_x \psi_{\tau}^*)^T J \partial_x \psi_{\tau}^* = (\psi_{\tau}' + \tau^{q+1} \delta_{\tau}')^T J (\psi_{\tau}' + \tau^{q+1} \delta_{\tau}')$$
  
=  $J + \mathcal{O}(\tau^{r+1}) + \tau^{q+1} \left( (\delta_{\tau}')^T J \psi_{\tau}' + (\psi_{\tau}')^T J \delta_{\tau}' \right) + \mathcal{O}(\tau^{2(q+1)}),$ 

which implies that

$$\tau^{q+1}\left((\delta_{\tau}')^T J \psi_{\tau}' + (\psi_{\tau}')^T J \delta_{\tau}'\right) = \mathcal{O}(\tau^{\min(2(q+1),r+1)}).$$

As an immediate consequence, we have that

$$(R'_{\tau})^{T}JR'_{\tau} = J + \mathcal{O}(\tau^{\min(2(q+1),r+1)})$$

which proves the second statement.  $\Box$ 

This result can be rendered more specific as follows:

**Proposition 2.** Let  $\mathcal{S}^{[2n]}_{\tau}$  be a smooth method of order  $2n \geq 2$  and pseudo-symmetry order  $q \geq 2n+1$ . Let us consider the composition method

$$\psi_{\tau}^{[2n+1]} = S_{\gamma_1 \tau}^{[2n]} \circ S_{\gamma_2 \tau}^{[2n]}, \tag{14}$$

where the coefficients  $\gamma_1$  and  $\gamma_2$  satisfy both relations  $\gamma_1 + \gamma_2 = 1$  and  $\gamma_1^{2n+1} + \gamma_2^{2n+1} = 0$ . Then the method

$$\hat{R}_{\tau} = \frac{1}{2} \left( \psi_{\tau}^{[2n+1]} + \overline{\psi}_{\tau}^{[2n+1]} \right) \tag{15}$$

is of order

$$\begin{cases}
2n+1 & \text{if } q = 2n+1, \\
2n+2 & \text{if } q \ge 2n+2
\end{cases}$$
(16)

when the vector field f in (1) is real, and of pseudo-symmetry order

$$\begin{cases} 2n+1 & \text{if } q=2n+1, \\ \min(q,4n+3) & \text{if } q \geq 2n+2. \end{cases}$$
 (17)

If in addition, f is a (real) Hamiltonian vector field and  $\mathcal{S}_{\tau}^{[2n]}$  is of pseudo-symplecticity order r, then  $\hat{R}_{\tau}$  is of pseudo-symplecticity order

$$\begin{cases}
\min(r, 2n+1) & \text{if } q = 2n+1, \\
\min(q, r, 4n+3) & \text{if } q \ge 2n+2.
\end{cases}$$
(18)

**Remark 2.** Note that in Proposition 2, one has necessarily  $q \ge 2n+1$ . This can be seen straightforwardly by a direct computation of  $\mathcal{S}_{-\tau}^{[2n]} \circ \mathcal{S}_{\tau}^{[2n]}(x)$  with  $\mathcal{S}_{\tau}^{[2n]}(x) = \varphi_{\tau}(x) + \tau^{2n+1}C(x) + \mathcal{O}(\tau^{2n+2})$ .

**Proof.** Noticing that  $\gamma_1$  and  $\gamma_2$  are complex conjugate and (1) is real, and taking into account that  $S_{\tau}^{[2n]}$  is of pseudo-symmetry order q, we have

$$\overline{\psi}_{\tau}^{[2n+1]} = S_{\gamma_{2}\tau}^{[2n]} \circ S_{\gamma_{1}\tau}^{[2n]} = \left( (S_{\gamma_{2}\tau}^{[2n]})^{*} + \mathcal{O}(\tau^{q+1}) \right) \circ \left( (S_{\gamma_{1}\tau}^{[2n]})^{*} + \mathcal{O}(\tau^{q+1}) \right) 
= (S_{\gamma_{2}\tau}^{[2n]})^{*} \circ (S_{\gamma_{1}\tau}^{[2n]})^{*} + \mathcal{O}(\tau^{q+1}) = (\psi_{\tau}^{[2n+1]})^{*} + \mathcal{O}(\tau^{q+1}).$$
(19)

Moreover, by construction,  $\psi_{\tau}^{[2n+1]}$  is at least of order 2n+1, so that

$$\psi_{\tau}^{[2n+1]} + \mathcal{O}(\tau^{2n+2}) = \overline{\psi}_{\tau}^{[2n+1]} = (\psi_{\tau}^{[2n+1]})^* + \mathcal{O}(\tau^{2n+2}), \tag{20}$$

and altogether

$$\hat{R}_{\tau} = R_{\tau} + \mathcal{O}(\tau^{\max(2n+2,q+1)}). \tag{21}$$

Now, since the pseudo-symmetry order of  $\psi_{z}^{[2n+1]}$  is at least 2n+1, the method

$$R_{\tau} = \frac{1}{2} \left( \psi_{\tau}^{[2n+1]} + (\psi_{\tau}^{[2n+1]})^* \right)$$

is, according to Proposition 1, of pseudo-symmetry order 4n + 3 and of pseudo-symplecticity order  $\min(4n + 3, q)$ . The first (16), second (17) and third (18) statements on orders then follow from (21).  $\Box$ 

In the Appendix we provide an alternative proof of Proposition 2 based on the Lie formalism, which allows us, in addition, to generalise the previous result on pseudo-symplecticity to other geometric properties the continuous system may possess (such as in volume preserving flows, isospectral flows, differential equations evolving on Lie groups, etc.).

Notice that, according with Proposition 2, if we start from n=1, that is to say from a basic symmetric  $(q=+\infty)$  and symplectic  $(r=+\infty)$  method of order 2, we get a method of order 4 that is pseudo-symmetric and pseudo-symplectic of order 7 just by considering the simple composition (14) and taking the real part of the output at each time step. If this technique is applied to a symmetric and symplectic method of order 4, i.e. with n=2, then  $\hat{R}_{\tau}$  is of order 6 and pseudo-symmetric and pseudo-symplectic of order 11.

Let us consider, in particular, the 4th-order symmetric scheme (7) with k=2 as basic scheme. Then, the resulting 6th-order integrator  $\hat{R}_{\tau}$  only requires the evaluation of 6 second-order methods  $\mathcal{S}_{\tau}^{[2]}$ , whereas the corresponding 6th-order scheme obtained by the triple-jump technique involves 9 evaluations. This number is reduced to 7 by considering general compositions of  $\mathcal{S}_{\tau}^{[2]}$  [8]. If we take this 6th-order composition of 7 schemes as the basic method  $\mathcal{S}_{\tau}^{[6]}$ , the resulting integrator of order 8,  $\hat{R}_{\tau}$ , involves the evaluation of 14  $\mathcal{S}_{\tau}^{[2]}$ , whereas 15 evaluations are required by pure composition methods. Notice that  $\hat{R}_{\tau}$  is pseudo-symmetric and pseudo-symplectic of order 15, so that for values of  $\tau$  sufficiently small, it preserves effectively the symmetry up to round-off error while the drift in energy for Hamiltonian systems is hardly noticeable.

# 3. Families of pseudo-symplectic methods

There is another possibility to increase the order, though, and it consists in applying the technique of Proposition 2 recursively. Thus, if denote by  $\hat{R}_{\tau}^{(1)} \equiv \hat{R}_{\tau}$  the method of Eq. (15), we propose to apply the following recurrence:

For 
$$i = 2, 3, ...$$

$$\Phi_{\tau}^{(i)} = \hat{R}_{\gamma^{[2i]}\tau}^{(i-1)} \circ \hat{R}_{\bar{\gamma}^{[2i]}\tau}^{(i-1)}$$

$$\hat{R}_{\tau}^{(i)} = \frac{1}{2} \left( \Phi_{\tau}^{(i)} + \overline{\Phi}_{\tau}^{(i)} \right)$$
(22)

where  $\gamma^{[2i]}$  is given by (6). Then, according with Proposition 2, it is possible to raise the order up to the pseudo-symmetry order of the underlying basic method  $\mathcal{S}^{[2n]}_{\tau}$ . Thus, in particular, the maximum order one can achieve by applying this technique to the basic symmetric method  $\mathcal{S}^{[2]}_{\tau}$  is 7, whereas if we start with a basic symmetric method of order 4,  $\mathcal{S}^{[4]}_{\tau}$ , the maximum order is 11. It is 15 from a symmetric method  $\mathcal{S}^{[6]}_{\tau}$  of order 6 and so on and so forth.

To give an assessment of the computational cost of the methods obtained by applying this type of composition, we notice that the computation of  $\Phi_{\tau}^{(i)}$  and  $\overline{\Phi}_{\tau}^{(i)}$  required to form  $\hat{R}_{\tau}^{(i)}$  by (22) at the intermediate stages can be done in parallel, whereas at the final stage it only requires taking the real part. Thus, the method of order 6 constructed recursively from  $S_{\tau}^{[2]}$  only requires the effective computation of 4 basic methods  $S_{\tau}^{[2]}$ .

Starting from a symmetric second-order method  $S_{\tau}^{[2]}$ , say Strang splitting for instance, it is important to monitor the sign of the real part of all coefficients involved in the previous iteration. It is immediate to see that in the recursive construction

$$\mathcal{S}_{\tau}^{[2]} \rightarrow \hat{R}_{\tau}^{(1)} \rightarrow \hat{R}_{\tau}^{(2)} \rightarrow \hat{R}_{\tau}^{(3)}$$

envisaged in the recurrence (22), the basic method  $S_{\tau}^{[2]}$  is used with the following coefficients

$$\begin{split} &i=1:\gamma^{[2]},\quad \bar{\gamma}^{[2]}\\ &i=2:\gamma^{[4]}\gamma^{[2]},\quad \bar{\gamma}^{[4]}\gamma^{[2]},\quad \gamma^{[4]}\bar{\gamma}^{[2]},\quad \bar{\gamma}^{[4]}\bar{\gamma}^{[2]}\\ &i=3:\gamma^{[6]}\gamma^{[4]}\gamma^{[2]},\quad \gamma^{[6]}\bar{\gamma}^{[4]}\gamma^{[2]},\quad \gamma^{[6]}\gamma^{[4]}\bar{\gamma}^{[2]},\quad \bar{\gamma}^{[6]}\gamma^{[4]}\gamma^{[2]},\quad \bar{\gamma}^{[6]}\bar{\gamma}^{[4]}\gamma^{[2]},\quad \bar{\gamma}^{[6]}\bar{\gamma}^{[4]}\bar{\gamma}^{[2]},\quad \bar{\gamma}^{[6]}\bar{\gamma}^{[4]}\bar{\gamma}^$$

Given the expression of  $\gamma^{[k]}$  (see (6)), these coefficients have arguments of the form

$$\frac{\pi}{2} \sum_{i=1}^{1} \pm \frac{1}{2j+1} = \frac{\pi}{2} \left( \pm \frac{1}{3} \pm \frac{1}{5} \pm \dots \pm \frac{1}{2i+1} \right), \quad i = 1, 2, 3,$$

so that their maximum argument is

$$\frac{\pi}{2}\sum_{i=1}^{3}\frac{1}{2j+1}$$
.

For all the coefficients to have positive real parts, a necessary and sufficient condition is thus that

$$\sum_{i=1}^{3} \frac{1}{2j+1} \le 1.$$

It clearly holds for methods  $\hat{R}_{\tau}^{(1)}$ ,  $\hat{R}_{\tau}^{(2)}$  and  $\hat{R}_{\tau}^{(3)}$ , of respective orders 4, 6 and 7, since 1/3 + 1/5 + 1/7 = 71/105. Similarly, starting form a symmetric method of order 4 having real or complex coefficients with maximum argument  $\theta_4$ , the condition becomes

$$\frac{2\theta_4}{\pi} + \sum_{j=1}^4 \frac{1}{2j+3} = \frac{2\theta_4}{\pi} + \frac{1888}{3465} \le 1.$$

For instance, suppose that f in (1) can be split as  $f(x) = f_a(x) + f_b(x)$ , so that the exact  $\tau$ -flows  $\varphi_{\tau}^{[a]}$  and  $\varphi_{\tau}^{[b]}$  corresponding to  $f_a$  and  $f_b$ , respectively, can be computed exactly. Then, the following composition

$$S_{\tau}^{[4]} = \varphi_{b_{1}\tau}^{[b]} \circ \varphi_{a_{1}\tau}^{[a]} \circ \varphi_{b_{2}\tau}^{[b]} \circ \varphi_{a_{2}\tau}^{[a]} \circ \varphi_{b_{2}\tau}^{[b]} \circ \varphi_{b_{2}\tau}^{[a]} \circ \varphi_{b_{1}\tau}^{[a]} \circ \varphi_{b_{1}\tau}^{[b]}$$

$$(23)$$

with

$$b_1 = \frac{1}{10} - \frac{1}{30}i$$
,  $b_2 = \frac{4}{15} + \frac{2}{15}i$ ,  $b_3 = \frac{4}{15} - \frac{1}{5}i$  and  $a_1 = a_2 = a_3 = a_4 = \frac{1}{4}i$ 

provides a 4th-order symmetric scheme (see [6]). Taking (23) as basic method we get  $\max_{i=1,2,3} \operatorname{Arg}(b_i) = \arccos(4/5)$  so that

$$\frac{2\theta_4}{\pi} + \frac{1888}{3465} < 0.409666 + \frac{1888}{3465} < 0.96 < 1$$

and thus all methods

$$S_{\tau}^{[4]} \to \hat{R}_{\tau}^{(1)} \to \hat{R}_{\tau}^{(2)} \to \hat{R}_{\tau}^{(3)} \to \hat{R}_{\tau}^{(4)}$$

of respective orders 4, 6, 8, 10 and 11 obtained by the procedure (22) have all their coefficients with positive real parts. As far as the  $f_a$  part is concerned, the maximum argument is less than  $0.55\frac{\pi}{2}$ .

#### 4. Numerical experiments

In this section we illustrate the previous results on several numerical examples, comprising Hamiltonian systems and partial differential equations of evolution previously discretised in space.

#### 4.1. Harmonic oscillator

We first consider the simple harmonic oscillator, with Hamiltonian

$$H = T(p) + V(q) = \frac{1}{2}p^2 + \frac{1}{2}q^2.$$

If we denote by  $M_X(\tau)$  the exact matrix evolution associated with the Hamiltonians X = H, T and V, i.e.,  $(q(\tau), p(\tau))^T = M_X(\tau)(q(0), p(0))^T$ , then

$$M_H(\tau) = \begin{pmatrix} \cos(\tau) & \sin(\tau) \\ -\sin(\tau) & \cos(\tau) \end{pmatrix}, \qquad M_T(\tau) = \begin{pmatrix} 1 & \tau \\ 0 & 1 \end{pmatrix}, \qquad M_V(\tau) = \begin{pmatrix} 1 & 0 \\ -\tau & 1 \end{pmatrix},$$

respectively. We take as basic symmetric (and symplectic) scheme the leapfrog/Strang splitting:

$$S_{\tau}^{[2]} = M_T(\tau/2)M_V(\tau)M_T(\tau/2) \tag{24}$$

and compute the first three iterations in (22). In Table 1 we collect the main term in the truncation error for the resulting integrators  $\hat{R}_{\tau}^{(i)}$ , i=1,2,3. We also check their time-symmetry and the preservation of the symplectic character of the approximate solution matrix by computing its determinant (a 2 × 2 matrix A is symplectic iff det(A) = 1). One can observe that these results are in agreement with the previous estimates.

Next we take initial conditions (q, p) = (2.5, 0), integrate until the final time  $t_f = 10^4$  with  $\mathcal{S}_{\tau}^{[2]}$ ,  $\hat{R}_{\tau}^{(1)}$ , and  $\hat{R}_{\tau}^{(2)}$  and compute the relative error in energy along the evolution. The result is depicted in Fig. 1. We see that for  $\hat{R}_{\tau}^{(1)}$  and  $\hat{R}_{\tau}^{(2)}$  the error in energy is almost constant for a certain period of time, and then there is a secular growth proportional to  $\mathcal{O}(\tau^7)$ .

Table 1 Main term in the truncation error, degree of symmetry and symplecticity for schemes  $\hat{R}_r^{(i)}$  obtained from the basic leapfrog integrator for the simple harmonic oscillator.  $I_2$  stands for the 2  $\times$  2 identity matrix..

$\psi_{ au}$	$M_H( au)-\psi_ au$	$\psi_{\tau} \circ \psi_{-\tau} - I_2$	$\det (\psi_{ au})$
$\hat{R}_{ au}^{(1)}$	$\left(\begin{array}{cc} 0 & -\frac{1}{180} \\ -\frac{1}{120} & 0 \end{array}\right) \tau^5$	$\left(\begin{array}{cc} -\frac{1}{1728} & 0\\ 0 & -\frac{1}{1728} \end{array}\right) \tau^8$	$1 - \frac{1}{1728}\tau^8$
$\hat{R}_{\tau}^{(2)}$	$\left(\begin{array}{cc} 0 & 3.8 \times 10^{-5} \\ 5.1 \times 10^{-5} & 0 \end{array}\right) \tau^7$	$\left(\begin{array}{cc} 5.4 \times 10^{-6} & 0 \\ 0 & 5.4 \times 10^{-6} \end{array}\right) \tau^{8}$	$1 + 5.4 \times 10^{-6}  \tau^8$
$\hat{R}_{\tau}^{(3)}$	$\left(\begin{array}{ccc} 5.8 \times 10^{-9} & 0 \\ 0 & 5.8 \times 10^{-9} \end{array}\right) \tau^8$	$\left( \begin{array}{ccc} -1.1 \times 10^{-8} & 0 \\ 0 & -1.1 \times 10^{-8} \end{array} \right) \tau^8$	$1-1.1\times 10^{-8}\tau^8$

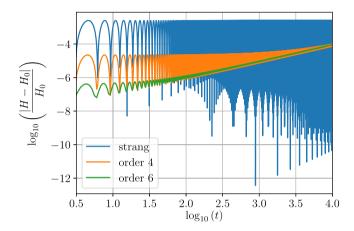


Fig. 1. Error in energy along the integration for the harmonic oscillator taking (24) as the basic integrator in the sequence (22).

#### 4.2. Kepler problem

Next, we consider the two-dimensional Kepler problem with Hamiltonian

$$H(q, p) = T(p) + V(q) = \frac{1}{2}p^{T}p - \mu \frac{1}{r}.$$
 (25)

Here  $q = (q_1, q_2), p = (p_1, p_2), \mu = GM, G$  is the gravitational constant and M is the sum of the masses of the two bodies. Taking  $\mu = 1$  and initial conditions

$$q_1(0) = 1 - e, q_2(0) = 0, p_1(0) = 0, p_2(0) = \sqrt{\frac{1 + e}{1 - e}},$$
 (26)

if  $0 \le e < 1$ , then the solution is periodic with period  $2\pi$ , and the trajectory is an ellipse of eccentricity e. Note that the gradient function must here be implemented carefully so as to be analytic for complex values of  $z = q_1^2 + q_2^2$ . Here, we define it using the following determination of the complex logarithm (analytic on the complex plane outside the negative real axis):

$$\forall (x,y) \in \mathbb{R}^2 \text{ s.t. } x + iy \notin \mathbb{R}_-, \qquad L(x+iy) = \log|x+iy| + 2i \arctan\left(\frac{y}{x+|x+iy|}\right).$$

As a consequence, the analytic continuation of the function  $1/r^3 = 1/(q_1^2 + q_2^2)^{3/2}$  writes

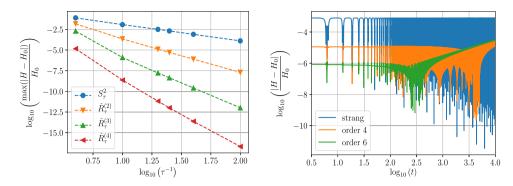
$$\exp\left(-\frac{3}{2}L(x+iy)\right),\,$$

where  $x=\Re(q_1^2+q_2^2)$  and  $y=\Im(q_1^2+q_2^2)$ . Here, as with the harmonic oscillator, we take as basic method the 2nd-order Strang splitting

$$S_{\tau}^{[2]} = \varphi_{\tau/2}^{[a]} \circ \varphi_{\tau}^{[b]} \circ \varphi_{\tau/2}^{[a]}, \tag{27}$$

where  $\varphi_{\tau}^{[a]}$  (respectively,  $\varphi_{\tau}^{[b]}$ ) corresponds to the exact solution obtained by integrating the kinetic energy T(p) (resp., potential energy V(q) in (25).

We take e = 0.6, integrate until the final time t = 20 with Strang and the schemes obtained by the recursion (22) with i = 1, 2, 3 for several time steps and compute the relative error in energy at the final time. Fig. 2 shows this error as



**Fig. 2.** Left figure: Relative error in energy vs. the inverse of the step size  $\tau$  after approximately 3.183 periods (t = 20) for the Kepler problem for the schemes obtained by the recurrence (22). Right figure: Evolution of this error along the integration.

a function of the inverse of the step size  $1/\tau$  to illustrate the order of convergence: order 2 for Strang, order 4 for  $\hat{R}_{\tau}^{(1)}$  and order 6 for  $\hat{R}_{\tau}^{(2)}$ . For  $\hat{R}_{\tau}^{(3)}$ , and contrary to what happens to the harmonic oscillator, the observed numerical order is higher than expected, varying between 7 and 8. We do not have at present a theoretical explanation for this phenomenon. Fig. 2 (right) depicts the time evolution of this error when the final time is  $t = 10^4$ .

# 4.3. The semi-linear reaction-diffusion equation of Fisher

Our third test-problem is the scalar equation in one-dimension

$$\frac{\partial u(x,t)}{\partial t} = \Delta u(x,t) + F(u(x,t)),\tag{28}$$

with periodic boundary conditions on the interval [0, 1]. Here F(u) is a *nonlinear* reaction term. For the purpose of testing our methods, we take Fisher's potential [17]

$$F(u) = u(1-u)$$

as considered for example in [8].

The splitting corresponds here to solving, on the one hand, the linear equation with the Laplacian (as  $f_a$ ), and on the other hand, the non-linear ordinary differential equation

$$\frac{\partial u(x,t)}{\partial t} = u(x,t)(1 - u(x,t)),$$

with initial condition  $u(x, 0) = u_0(x)$ , whose analytical solution is given by the well-defined (for small enough complex time t) formula

$$u(x,t) = u_0(x) + u_0(x)(1 - u_0(x)) \frac{(e^t - 1)}{1 + u_0(x)(e^t - 1)}.$$

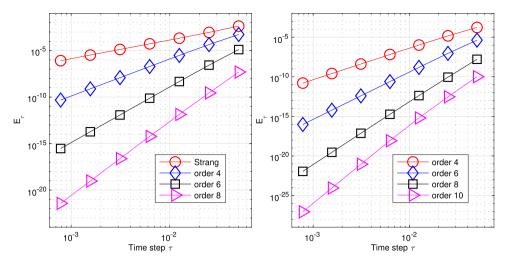
Here we aim to solve Eq. (28) with periodic boundary conditions on the interval [0, 1], and initial condition  $u_0(x) = \sin(2\pi x)$ . Numerically, the interval is discretised on a uniform grid, i.e.,  $x_j = j/N$ ,  $j = 0, \ldots, N-1$ ,  $N \in \mathbb{N}$ , and u(x, t) is approximated by Fourier pseudo-spectral methods. In this way we construct a vector  $\mathbf{u}$  with components  $(\mathbf{u})_j \approx u(x_{j-1}, t)$ ,  $j = 1, 2, \ldots, N$ . If we denote by  $\mathbf{u}_{\tau}$  the whole numerical solution computed by a certain integrator with step size  $\tau$  from t = 0 until the final time, and by  $\mathbf{u}_{\tau/2}$  the corresponding numerical solution computed by the same integrator with step size  $\tau/2$ , then the quantity  $E_{\tau} := \|\mathbf{u}_{\tau} - \mathbf{u}_{\tau/2}\|_{\infty}$  is a good indicator of the convergence order.

Numerical simulations were carried out in quadruple precision (with Intel Fortran) such that roundoff errors are suppressed. Fig. 3 shows the successive errors  $E_{\tau}$ , at final time T=10, of the methods obtained with the sequence (22) with the Strang splitting as the basic method  $\mathcal{S}_{\tau}^{[2]}$  (left) and the fourth order scheme  $\mathcal{S}_{\tau}^{[4]}$  given by (23) (right) with different time steps  $\tau_j=0.1/2^j,\ j=1,\ldots,7$ . One can clearly observe that the convergence order matches the previous analysis with a slightly better performance for the highest order, analogously to the Kepler problem. Fig. 4 shows the successive errors versus the number of basic integrators in each case.

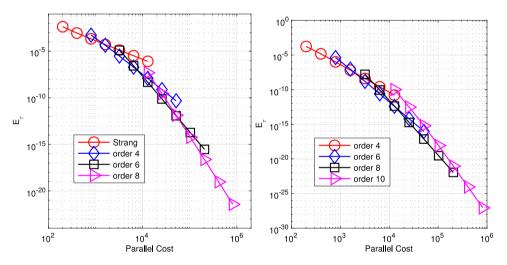
# 4.4. The semi-linear complex Ginzburg-Landau equation

Our final test problem is the complex Ginzburg–Landau equation on the domain  $(x, t) \in [-100, 100] \times [0, 100]$ ,

$$\frac{\partial u(x,t)}{\partial t} = \alpha \Delta u(x,t) + \varepsilon u(x,t) - \beta |u(x,t)|^2 u(x,t), \tag{29}$$



**Fig. 3.** Successive errors  $E_{\tau}$  versus time step  $\tau$  for Eq. (28) of the composition methods starting from the Strang scheme (left) and the fourth order scheme  $S_{\tau}^{[4]}$  (right).



**Fig. 4.** Successive errors  $E_{\tau}$  versus number of basic integrators for Eq. (28) of the composition methods starting from the Strang scheme (left) and the fourth order scheme  $S_{\tau}^{[4]}$  (right).

with  $\alpha = 1 + ic_1$ ,  $\beta = 1 - ic_3$  and initial condition  $u(x, 0) = u_0(x)$ . Here,  $\varepsilon$ ,  $c_1$  and  $c_3$  denote *real* coefficients. In physics, the Ginzburg–Landau appears in the mathematical theory used to model superconductivity. For a broad introduction to the rich dynamics of this equation, we refer to [18]. Here, we will use the values  $c_1 = 1$ ,  $c_3 = -2$  and  $\varepsilon = 1$ , for which plane wave solutions establish themselves quickly after a transient phase (see [19]). In addition, we set

$$u_0(x) = \frac{0.8}{\cosh(x - 10)^2} + \frac{0.8}{\cosh(x + 10)^2},$$

so that the solution can be represented in Fig. 5.

To apply the composition methods presented in previous sections, it seems natural to split equation (29) as

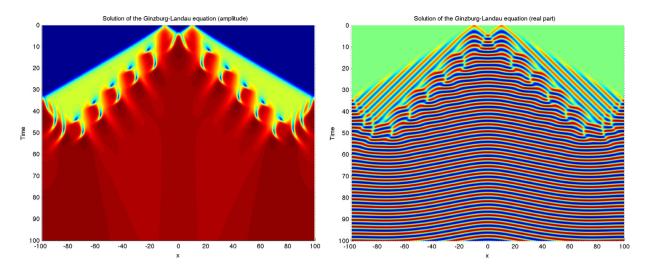
$$\frac{\partial u(x,t)}{\partial t} = (1+ic_1)\Delta u(x,t) + \varepsilon u(x,t),\tag{30}$$

whose solution is  $u(x, t) = e^{\varepsilon t} e^{t(1+ic_1)\Delta} u_0(x)$  for  $t \ge 0$ , and

$$\frac{\partial u(x,t)}{\partial t} = -(1-ic_3)|u(x,t)|^2 u(x,t) \tag{31}$$

with solution is for  $t \ge 0$ 

$$u(x,t) = e^{-(1-ic_3)\int_0^t M(x,s)ds} u_0(x) = e^{-\frac{\beta}{2}\log(1+2tM_0(x))} u_0(x).$$



**Fig. 5.** Colourmaps of the amplitude  $|u(x,t)|^2$  (left) and real part  $\Re(u(x,t))$  (right) of the solution of (29).

Here we have first solved the equation for  $M(x, t) := |u(x, t)|^2$ , given by

$$\frac{\partial M(x,t)}{\partial t} = -2M^2(x,t),$$

with solution

$$M(x, t) = \frac{M_0(x)}{1 + 2M_0(x)t}.$$

Considering t now as a complex variable with positive real part does not raise any difficulty for the first part, since  $e^{\varepsilon t}e^{t(1+ic_1)\Delta}$  is well-defined. More care has to be taken for the second part, since  $u\mapsto |u|^2u$  is *not* a holomorphic function, and this prevents us from solving (29) in its current form. As a consequence, we first rewrite (29) as a system for (v(x,t),w(x,t)) where  $v(x,t)=\Re(u(x,t))$  and  $w(x,t)=\Im(u(x,t))$ :

$$\begin{cases}
\frac{\partial v(x,t)}{\partial t} &= \Delta v(x,t) - c_1 \Delta w(x,t) + \varepsilon v(x,t) - (v^2(x,t) + w^2(x,t))(v(x,t) + c_3 w(x,t)) \\
\frac{\partial w(x,t)}{\partial t} &= c_1 \Delta v(x,t) + \Delta w(x,t) + \varepsilon w(x,t) - (v^2(x,t) + w^2(x,t))(-c_3 v(x,t) + w(x,t))
\end{cases} (32)$$

and now solve it for complex time  $t \in \mathbb{C}$  with  $\Re(t) \geq 0$ . Observing that

$$\begin{pmatrix} -1 & -c_3 \\ c_3 & -1 \end{pmatrix} = PD_3P^{-1} \quad \text{and} \quad \begin{pmatrix} 1 & -c_1 \\ c_1 & 1 \end{pmatrix} = PD_1P^{-1},$$

with

$$D_1 = \left( \begin{array}{cc} \alpha & 0 \\ 0 & \bar{\alpha} \end{array} \right), \qquad D_3 = \left( \begin{array}{cc} -\beta & 0 \\ 0 & -\bar{\beta} \end{array} \right), \qquad P = \left( \begin{array}{cc} i & 1 \\ 1 & i \end{array} \right) \qquad \text{and} \qquad P^{-1} = \left( \begin{array}{cc} -\frac{i}{2} & \frac{1}{2} \\ \frac{1}{2} & -\frac{i}{2} \end{array} \right),$$

system (32) can be rewritten as

$$\begin{cases}
\frac{\partial \tilde{v}(x,t)}{\partial t} &= \left(\alpha \Delta \tilde{v}(x,t) + \varepsilon \tilde{v}(x,t)\right) - \left(\beta \tilde{M}(x,t)\tilde{v}(x,t)\right) \\
\frac{\partial \tilde{w}(x,t)}{\partial t} &= \left(\bar{\alpha} \Delta \tilde{w}(x,t) + \varepsilon \tilde{w}(x,t)\right) - \left(\bar{\beta} \tilde{M}(x,t)\tilde{w}(x,t)\right)
\end{cases} (33)$$

where  $\tilde{M}(x, t) = 4i\tilde{v}(x, t)\tilde{w}(x, t)$  and where

$$\left(\begin{array}{c} \tilde{v} \\ \tilde{w} \end{array}\right) = \frac{1}{2} \left(\begin{array}{cc} -i & 1 \\ 1 & -i \end{array}\right) \left(\begin{array}{c} v \\ w \end{array}\right).$$

It is not difficult to see that the exact solution of the second part of (33) is given by

$$\begin{cases}
\tilde{v}(x,t) &= \tilde{v}_0(x)e^{-\frac{\beta}{2}\log(1+2t\tilde{M}_0(x))} \\
\tilde{w}(x,t) &= \tilde{w}_0(x)e^{-\frac{\bar{\beta}}{2}\log(1+2t\tilde{M}_0(x))}
\end{cases}$$
(34)

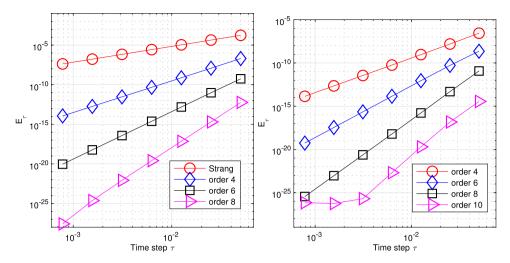


Fig. 6. Successive errors  $E_{\tau}$  versus time step  $\tau$  for Eq. (29) of the composition methods starting from the Strang scheme (left) and the fourth order scheme  $S_{\tau}^{[4]}$  (right).

where  $\tilde{M}_0(x)$  is now defined as  $\tilde{M}_0(x) := 4i\tilde{v}_0(x)\tilde{w}_0(x)$ . Note that here, by convention, the logarithm refers to the principal value of  $\log(z)$  for complex numbers: if  $z = (a+ib) = r\mathrm{e}^{\mathrm{i}\theta}$  with  $-\pi < \theta \leq \pi$ , then

$$\log z := \ln r + i\theta = \ln |z| + i \arg z = \ln(|a+ib|) + 2i \arctan\left(\frac{b}{a+\sqrt{a^2+b^2}}\right).$$

Since  $\log(z)$  is not defined for  $z \in \mathbb{R}^-$ , this means that the solution  $(\tilde{v}(x,t), \tilde{w}(x,t))$  is defined only as long as  $1+2\tilde{M}_0(x)t \notin$  $\mathbb{R}^-$ . Finally, the solution (v(x,t),w(x,t)) is of the form

$$\begin{cases} v(x,t) &= v_0(x) \frac{(\mathrm{e}^{-\beta L(x,t)} + \mathrm{e}^{-\bar{\beta} L(x,t)})}{2} - w_0(x) \frac{(\mathrm{e}^{-\beta L(x,t)} - \mathrm{e}^{-\bar{\beta} L(x,t)})}{2i} \\ w(x,t) &= v_0(x) \frac{(\mathrm{e}^{-\beta L(x,t)} - \mathrm{e}^{-\bar{\beta} L(x,t)})}{2i} + w_0(x) \frac{(\mathrm{e}^{-\beta L(x,t)} + \mathrm{e}^{-\bar{\beta} L(x,t)})}{2} \end{cases}$$

where  $L(x, t) := \log(1 + 2t\tilde{M}_0(x)) = \log(1 + 2tM_0(x))$  with  $M_0(x) = v_0^2(x) + w_0^2(x)$ . Denoting  $V = (v_1, \dots, v_N) \in \mathbb{R}^N$  and  $W = (w_1, \dots, w_N) \in \mathbb{R}^N$ , we eventually have to numerically solve the following system:

$$\begin{cases} \dot{V} = AV - c_1AW + \varepsilon V - G(V + c_3W) \\ \dot{W} = c_1AV + AW + \varepsilon W - G(-c_3V + W) \end{cases}$$

where G is the diagonal matrix with  $G_{i,i} = v_i^2 + w_i^2$ .

Eq. (29) is solved with periodic boundary conditions on the interval [-100, 100]. Now, in the previous example, the interval is discretised on a uniform grid, i.e.,  $x_i = j/N$ ,  $j = 0, \dots, N-1$ ,  $N \in \mathbb{N}$  with N = 512, and u(x, t) is approximated by Fourier pseudo-spectral methods. The successive errors  $E_{\tau} := \|\mathbf{u}_{\tau} - \mathbf{u}_{\tau/2}\|_{\infty}$  are shown also here to confirm the convergence order. Fig. 6 shows the successive errors, at final time T = 10, of the schemes obtained by applying the sequence (22) from the basic Strang splitting and the fourth-order scheme (23) with  $\tau_i = 0.1/2^j$ ,  $j = 1, \dots, 7$ . The observed order of convergence matches the previous analysis with a slightly better performance for the highest order. Fig. 7 shows the successive errors versus the number of basic integrators.

# Acknowledgements

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#### **Appendix**

In this Appendix we provide an alternative proof of Proposition 2 via Lie formalism. This allows us not only to gain some additional insight into the structure of the methods, but also to generalise the result on pseudo-symplecticity to other properties of geometric character, very often related to Lie groups, the differential equation may possess.

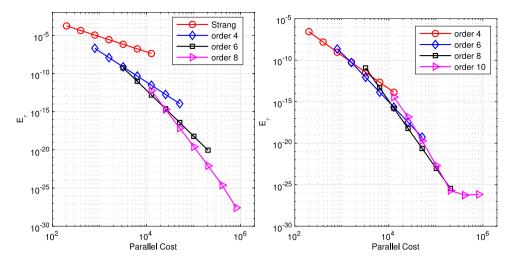


Fig. 7. Successive errors  $E_{\tau}$  versus number of basic integrators for Eq. (29) of the composition methods starting from the Strang scheme (left) and the fourth order scheme  $S_{z}^{[4]}$  (right).

To begin with, if  $\varphi_{\tau}$  is the exact flow of Eq. (1), then for each infinitely differentiable map g, the function  $g(\varphi_{\tau}(x))$ admits an expansion of the form [14,20]

$$g(\varphi_{\tau}(x)) = \exp(\tau F)[g](x) = g(x) + \sum_{k>1} \frac{\tau^k}{k!} F^k[g](x),$$

where F is the Lie derivative associated with f,

$$F = \sum_{i>1} f_i(x) \frac{\partial}{\partial x_i}.$$
 (35)

Analogously, for a given integrator  $\psi_{\tau}$  one can associate a series of linear operators so that

$$g(\psi_{\tau}(x)) = \exp(Y(\tau))[g](x),$$
 with  $Y(\tau) = \sum_{j \ge 1} \tau^j Y_j$ 

for all functions g [21]. The integrator  $\psi_{\tau}$  is of order k if

$$Y_1 = F$$
 and  $Y_i = 0$  for  $2 \le j \le k$ .

For the adjoint integrator  $\psi_{\tau}^* = \psi_{\tau}^{-1}$ , one clearly has

$$g(\psi_{\tau}^*(x)) = \exp(-Y(-\tau))[g](x).$$

This shows that  $\psi_{\tau}$  is symmetric if and only if  $Y(\tau) = \tau Y_1 + \tau^3 Y_3 + \cdots$ , and in particular, that symmetric methods are of even order.

An integrator  $S_{\tau}^{[2n]}$  of order  $2n \geq 2$  can be associated with the series

$$\Phi(\tau) = \exp\left(\tau F + \tau^{2n+1} N_{2n+1} + \tau^{2n+2} N_{2n+2} + \cdots\right) \tag{36}$$

for certain operators  $N_k$ . Then, the adjoint method  $(\mathcal{S}_{\tau}^{[2n]})^*$  has the associated series

$$\Phi^*(\tau) = \exp\left(\tau F + \tau^{2n+1} N_{2n+1} - \tau^{2n+2} N_{2n+2} + \cdots\right).$$

In consequence,  $\mathcal{S}_{\tau}^{[2n]}$  is pseudo-symmetric of order  $q \geq 2n+1$ . (i) Let us analyse first the case q > 2n+1. Then,  $N_{2n+2} = \cdots = N_q = 0$  in (36) and the series of operators associated with the composition  $\psi_{\tau}^{[2n+1]} = \mathcal{S}_{\gamma_1 \tau}^{[2n]} \circ \mathcal{S}_{\gamma_2 \tau}^{[2n]}$  is

$$\Psi(\tau) = \Phi(\gamma_2 \tau) \Phi(\gamma_1 \tau) \equiv \exp(V(\tau)).$$

where  $V(\tau)$  can be formally determined by applying the Baker-Campbell-Hausdorff formula [13] as

$$\begin{split} V(\tau) &= (\gamma_1 + \gamma_2)\tau F + (\gamma_1^{2n+1} + \gamma_2^{2n+1})\tau^{2n+1} N_{2n+1} \\ &+ \frac{1}{2}(\gamma_2\gamma_1^{2n+1} - \gamma_1\gamma_2^{2n+1})\tau^{2n+2} [F,N_{2n+1}] + (\gamma_1^{2n+3} + \gamma_2^{2n+3})\tau^{2n+3} N_{2n+3} + \mathcal{O}(\tau^{2n+4}). \end{split}$$

Here  $[\cdot,\cdot]$  denotes the usual Lie bracket. Clearly, the order of  $\psi_{ au}^{[2n+1]}$  is 2n+1 if

$$\gamma_1 + \gamma_2 = 1, \qquad \gamma_1^{2n+1} + \gamma_2^{2n+1} = 0,$$
 (37)

so that  $\gamma_2 = \bar{\gamma}_1 \equiv \gamma$  is given by Eq. (5) (with k = 2n). In that case we can write

$$V(\tau) = \tau F + \tau^{2n+2} G(\tau),$$
 with  $G(\tau) = \sum_{i=0}^{\infty} \tau^i G_i,$ 

whereas for the adjoint method one has

$$\Psi^*(\tau) = \exp(-V(-\tau)) = \exp\left(\tau F + \tau^{2n+2}\widetilde{G}(\tau)\right), \quad \text{with} \quad \widetilde{G}(\tau) = \sum_{i=0}^{\infty} (-1)^{i+1} \tau^i G_i.$$

In particular,  $G_0 = \frac{1}{2}(\gamma_2\gamma_1^{2n+1} - \gamma_1\gamma_2^{2n+1})[F, N_{2n+1}]$ , and  $G_1 = (\gamma_1^{2n+3} + \gamma_2^{2n+3})N_{2n+3}$ . The series  $\Psi(\tau)$  can also be written as

$$\Psi(\tau) = \exp\left(\frac{\tau}{2}F\right) \exp W(\tau) \exp\left(\frac{\tau}{2}F\right),$$

where  $W(\tau)$  is determined by applying the symmetric BCH formula [13] as

$$W(\tau) = \tau^{2n+2}G(\tau) + \frac{1}{24}\tau^{2n+4}[F, [F, G(\tau)]] + \mathcal{O}(\tau^{4n+4})$$
  
=  $\tau^{2n+2}G_0 + \tau^{2n+3}G_1 + \tau^{2n+4}(G_2 + \frac{1}{24}[F, [F, G_0]]) + \mathcal{O}(\tau^{2n+5}).$ 

By the same token,

$$\Psi^*(\tau) = \exp\left(\frac{\tau}{2}F\right) \exp\left(-W(-\tau)\right) \exp\left(\frac{\tau}{2}F\right).$$

Consider now the method

$$R_{\tau} = \frac{1}{2} \left( \psi_{\tau}^{[2n+1]} + (\psi_{\tau}^{[2n+1]})^* \right). \tag{38}$$

Clearly, its associated series of operators,

$$\mathcal{R}(\tau) = \frac{1}{2}\Psi(\tau) + \frac{1}{2}\Psi^*(\tau),$$

can be expressed as

$$\mathcal{R}(\tau) = \exp\left(\frac{\tau}{2}F\right)\mathcal{Y}\,\exp\left(\frac{\tau}{2}F\right),$$

where

$$y = \frac{1}{2} \exp(W(\tau)) + \frac{1}{2} \exp(-W(-\tau)).$$

By expanding, we have

$$y = I + \frac{1}{2}(W(\tau) - W(-\tau)) + \frac{1}{4}(W^{2}(\tau) + W^{2}(-\tau)) + \cdots,$$

but

$$W(\tau) - W(-\tau) = 2\tau^{2n+3} \sum_{i=0}^{\infty} \tau^{2i} z_{2i} \equiv 2\tau^{2n+3} Z(\tau),$$

with  $z_0 = G_1$ ,  $z_2 = G_3 + \frac{1}{24}[F, [F, G_1]]$ , etc. In general,  $z_{2i}$  is a linear combination of the operators  $\{F, N_{2n+1}, N_{2n+2}, \ldots\}$  and their nested Lie brackets. In addition,  $W^2(\tau) + W^2(-\tau) = \mathcal{O}(\tau^{4n+4})$ , so that we can write

$$\mathcal{Y} = I + \tau^{2n+3}Z + \mathcal{O}(\tau^{4n+4}) = \exp(\tau^{2n+3}Z) + \mathcal{O}(\tau^{4n+4})$$

and

$$\mathcal{R}(\tau) = \exp\left(\frac{\tau}{2}F\right) \exp\left(\tau^{2n+3}Z\right) \exp\left(\frac{\tau}{2}F\right) + \mathcal{O}(\tau^{4n+4}),$$

whence the following statements follow at once:

• Method (38) is of order 2n + 2, since  $\mathcal{R}(\tau) = \exp(\tau F) + \mathcal{O}(\tau^{2n+3})$ .

- Since  $Z(\tau)$  only contains even powers of  $\tau$  (up to  $\tau^{q+1}N_{q+1}$ ), then  $e^{\frac{\tau}{2}F}e^{\tau^{2n+3}Z}e^{\frac{\tau}{2}F}$  is a symmetric composition and  $R_{\tau}$  is pseudo-symmetric of order  $\min(q, 4n + 3)$ .
- Let us suppose that scheme (38) is applied to a Hamiltonian system and that  $\mathcal{S}_{\tau}^{[2n]}$  is of pseudo-symplecticity order r. Since Z is an operator in the free Lie algebra generated by  $\{F, N_{2n+1}, N_{2n+2}, \ldots\}$ , clearly the composition  $e^{\frac{\tau}{2}F}e^{\tau^{2n+3}Z}e^{\frac{\tau}{2}F}$  is symplectic (at least up to terms  $\mathcal{O}(\tau^r)$ ). As a matter of fact, this can be extended to any geometric property the differential equation (1) has: volume-preserving, unitary, etc., as long as the basic scheme  $\mathcal{S}_{\tau}^{[2n]}$  preserves this property up to order r.

Finally, in view of (19)–(21) and recalling that  $q \ge 2n+2$ , the same considerations apply if we take the complex conjugate instead of the adjoint, i.e., to the scheme

$$\hat{R}_{\tau} = \Re(\psi_{\tau}^{[2n+1]}) = \frac{1}{2} \left( \psi_{\tau}^{[2n+1]} + \overline{\psi}_{\tau}^{[2n+1]} \right). \tag{39}$$

(ii) We analyse next the case q=2n+1. Then  $N_{2n+2}\neq 0$  in (36) and, if  $\gamma_1$  and  $\gamma_2$  verify equations (37), then  $V(\tau)$  read

$$V(\tau) = \tau F + \tau^{2n+2} V_0 + \mathcal{O}(\tau^{2n+3})$$

with

$$V_0 = (\gamma_1^{2n+2} + \gamma_2^{2n+2})N_{2n+2} + \frac{1}{2}(\gamma_2\gamma_1^{2n+1} - \gamma_1\gamma_2^{2n+1})[F, N_{2n+1}].$$

Notice that, whereas  $\gamma_1^{2n+2} + \gamma_2^{2n+2}$  is a real number,  $\gamma_2 \gamma_1^{2n+1} - \gamma_1 \gamma_2^{2n+1}$  has non-vanishing real and imaginary parts. In any event, the same procedure as in the previous case can be carried out, leading to the conclusion that method (38) is still of order 2n + 2.

The situation is different, however, for method (39), since relations (19)–(21) do not provide further information. We have to analyse instead

$$\Re(\Psi(\tau)) = \exp\left(\frac{\tau}{2}F\right)\Re\left(\exp W(\tau)\right)\exp\left(\frac{\tau}{2}F\right),$$

with  $W(\tau) = \tau^{2n+2}V_0 + \mathcal{O}(\tau^{2n+3})$ . Noting that

$$\Re(\exp W(\tau)) = I + \Re(W(\tau)) + \mathcal{O}(\tau^{4n+4}) = I + \tau^{2n+2}\Re(V_0) + \mathcal{O}(\tau^{2n+3})$$

then we can write

$$\Re(\Psi(\tau)) = \exp\left(\frac{\tau}{2}F\right) \, \exp\!\left(\tau^{2n+2}\Re(V_0) + \mathcal{O}(\tau^{2n+3})\right) \, \exp\left(\frac{\tau}{2}F\right) + \mathcal{O}(\tau^{4n+4}).$$

In consequence,  $\hat{R}_{\tau}$  is a method of order 2n+1, pseudo-symmetric of order 2n+1 and pseudo-symplectic of order  $\min(r, 2n+1)$ .

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