

Three approaches for the design of adaptive time-splitting methods

Mechthild Thalhammer
Leopold–Franzens Universität Innsbruck, Austria

Workshop on Advances in Mathematical Modelling and
Numerical Simulation of Superfluids
Rouen, France, August 2017

Theme

Splitting methods. Time integration of **nonlinear evolution equations** by **exponential operator splitting methods**

$$\begin{cases} u'(t) = F(u(t)) = A(u(t)) + B(u(t)), & t \in (0, T), \\ u(0) \text{ given.} \end{cases}$$

Areas of application.

- Schrödinger equations (Quantum mechanics)
- Damped wave equations (Nonlinear acoustics)
- Parabolic equations (Pattern formation)
- Kinetic equations (Plasma physics)

Main theme

Local error control. Use of **local error control** to adjust time stepsize

$$\tau_{\text{optimal}} = \tau_{\text{current}} \cdot \min\left(\alpha_{\text{max}}, \max\left(\alpha_{\text{min}}, \sqrt[p+1]{\alpha \cdot \frac{\text{tol}}{\text{err}_{\text{local}}}}\right)\right)$$

in general enhances **reliability** and **efficiency** of time integration.

Question. How to construct **estimators for local error** in the context of splitting methods?

Approches. Different approaches rely on

- embedded splitting methods (with OTHMAR KOCH),
- defect-based a posteriori local error estimators (with HARALD HOFSTÄTTER, OTHMAR KOCH, WINFRIED AUZINGER),
- associated approximations with negligible additional cost (with SERGIO BLANES, FERNANDO CASAS).

Exponential operator splitting methods for nonlinear evolution equations

Exponential operator splitting methods

Splitting methods. For **nonlinear evolution equations** of form

$$\begin{cases} u'(t) = F(u(t)) = A(u(t)) + B(u(t)), & t \in (0, T), \\ u(0) \text{ given,} \end{cases}$$

determine **approximations** at time grid points $0 = t_0 < \dots < t_N \leq T$ with associated stepsizes $\tau_{n-1} = t_n - t_{n-1}$ for $n \in \{1, \dots, N\}$ by recurrence

$$u_n = \mathcal{S}_F(\tau_{n-1}, u_{n-1}) \approx u(t_n) = \mathcal{E}_F(\tau_{n-1}, u(t_{n-1})).$$

Splitting methods rely on presumption that corresponding subproblems are solvable in accurate and efficient manner

$$\begin{aligned} v'(t) &= A(v(t)), & w'(t) &= B(w(t)), \\ v(t) &= \mathcal{E}_A(t, v(0)), & w(t) &= \mathcal{E}_B(t, w(0)). \end{aligned}$$

High-order splitting methods are cast into following format with suitably chosen real (or complex) coefficients

$$\mathcal{S}_F(\tau, \cdot) = \mathcal{E}_B(b_s \tau, \cdot) \circ \mathcal{E}_A(a_s \tau, \cdot) \circ \dots \circ \mathcal{E}_B(b_1 \tau, \cdot) \circ \mathcal{E}_A(a_1 \tau, \cdot) \approx \mathcal{E}_F(\tau, \cdot).$$

Compact formulation

Compact formulation. Calculus of Lie-derivatives permits compact formulation and reveals analogies to significantly simpler linear case

$$\begin{aligned} & e^{a_1 \tau D_A} e^{b_1 \tau D_B} \dots e^{a_s \tau D_A} e^{b_s \tau D_B} \\ &= \mathcal{E}_B(b_s \tau, \cdot) \circ \mathcal{E}_A(a_s \tau, \cdot) \circ \dots \circ \mathcal{E}_B(b_1 \tau, \cdot) \circ \mathcal{E}_A(a_1 \tau, \cdot). \end{aligned}$$

Recipe. In order to extend result for linear case to nonlinear case,

- replace operator A, B by Lie-derivatives D_A, D_B and
- reverse order of evolution operators.

Example methods ($p = 1, 2$)

Low-order methods.

- First-order Lie–Trotter splitting method

$$a_1 = 1 = b_1, \quad \mathcal{S}_F(\tau, \cdot) = e^{\tau D_B} e^{\tau D_A}.$$

- Second-order Strang splitting method

$$a_1 = \frac{1}{2} = a_2, \quad b_1 = 1, \quad b_2 = 0,$$
$$\mathcal{S}_F(\tau, \cdot) = e^{\frac{1}{2}\tau D_A} e^{\tau D_B} e^{\frac{1}{2}\tau D_A}.$$

Example methods ($p = 4$)

Higher-order methods.

- Symmetric **fourth-order splitting method** by BLANES, MOAN (2002)

$$\begin{aligned}a_1 &= 0, & a_2 &= 0.245298957184271 = a_7, \\a_3 &= 0.604872665711080 = a_6, & a_4 &= \frac{1}{2} - (a_2 + a_3) = a_5, \\b_1 &= 0.0829844064174052 = b_7, & b_2 &= 0.3963098014983680 = b_6, \\b_3 &= -0.0390563049223486 = b_5, & b_4 &= 1 - 2(b_1 + b_2 + b_3).\end{aligned}$$

Stability ensured for **evolution equations of Schrödinger type**.

- Symmetric **fourth-order splitting method** by YOSHIDA ($s = 4$, complex variant of famous scheme)

$$\begin{aligned}\alpha &= 0.3243964040201711829761560 - 0.1345862724908066967894444 i, \\ \beta &= 0.3512071919596576340476880 + 0.2691725449816133935788885 i, \\ a_1 &= \frac{1}{2} \alpha, & a_2 &= \frac{1}{2} (\alpha + \beta) = a_3, & a_4 &= a_1, \\ b_1 &= \alpha = b_3, & b_2 &= \beta, & b_4 &= 0.\end{aligned}$$

Stability ensured for **evolution equations of parabolic type**, since $\Re(a_j), \Re(b_j) \geq 0$ for $j \in \{1, \dots, 4\}$.

Approaches for design and analysis of local error estimators

Local error estimators

Approaches. Study different approaches for design and theoretical analysis of local error estimators for splitting methods.

- **Embedded splitting methods**

O. KOCH, CH. NEUHAUSER, M. TH. *Embedded exponential operator splitting methods for the time integration of nonlinear evolution equations* (2013).

- **A posteriori local error estimators**

W. AUZINGER, O. KOCH, M. TH. *Defect-based local error estimators for splitting methods, with application to Schrödinger equations. Part I. The linear case* (2012).

W. AUZINGER, O. KOCH, M. TH. *Defect-based local error estimators for splitting methods, with application to Schrödinger equations. Part II. Higher-order methods for linear problems* (2014).

W. AUZINGER, H. HOFSTÄTTER, O. KOCH, M. TH. *Defect-based local error estimators for splitting methods, with application to Schrödinger equations. Part III. The nonlinear case* (2015).

- **Approximations with negligible additional cost** (recent work with SERGIO and FERNANDO)

Simplification. Specify local error estimators for first time step ($\tau > 0$).

Embedded splitting methods

Examples and theoretical basis

Embedded splitting methods

Heuristic approach. Consider splitting method of nonstiff order p

$$u_1 = \prod_{j=1}^s e^{a_{s+1-j}\tau D_A} e^{b_{s+1-j}\tau D_B} u_0.$$

Design related splitting method of nonstiff order \hat{p} such that **certain coefficients coincide**

$$\hat{u}_1 = \prod_{j=1}^{\hat{s}} e^{\hat{a}_{s+1-j}\tau D_A} e^{\hat{b}_{s+1-j}\tau D_B} u_0.$$

Use difference between two approximations as **local error estimator**

$$\text{err}_{\text{local}} = \|u_1 - \hat{u}_1\|_X.$$

Remark. Approach in spirit of *embedded Runge-Kutta methods* (but with higher cost).

Example (Schrödinger equations)

Example. Favourable scheme ($p = 4$, BLANES & MOAN) and embedded scheme ($\hat{p} = 3$, KOCH & TH.).

j	a_j	j	b_j
1	0	1,7	0.0829844064174052
2,7	0.245298957184271	2,6	0.3963098014983680
3,6	0.604872665711080	3,5	-0.0390563049223486
4,5	$1/2 - (a_2 + a_3)$	4	$1 - 2(b_1 + b_2 + b_3)$

j	\hat{a}_j	j	\hat{b}_j
1	a_1	1	b_1
2	a_2	2	b_2
3	a_3	3	b_3
4	a_4	4	b_4
5	0.3752162693236828	5	0.4463374354420499
6	1.4878666594737946	6	-0.0060995324486253
7	-1.3630829287974774	7	0

Example (Parabolic equations)

Example. Complex scheme ($p = 4$, YOSHIDA) and embedded scheme ($\hat{p} = 3$, KOCH & TH.).

j	a_j	j	b_j
1	0	1,4	0.1621982020100856 + 0.0672931362454034i
2,4	0.3243964040201712 + 0.1345862724908067i	2,3	0.3378017979899144 - 0.0672931362454034i
3	0.3512071919596576 - 0.2691725449816134i		

j	\hat{a}_j	j	\hat{b}_j
1	a_1	1	b_1
2	0.4157701540561051 + 0.2129482257474245i	2	0.4052251807333103 + 0.1988642124619028i
3	0.3855092282056243 - 0.1105557092016989i	3	0.4325766172566041 - 0.2661573487073062i
4	0.1987206177382706 - 0.1023925165457255i	4	0

Theoretical justification

Theoretical justification. Consider high-order splitting methods and employ **local error representations** that are suitable for nonlinear evolution equations involving **unbounded operators**.

Theorem (Th. 2008, Th. 2012, Koch & Neuhauser & Th. 2013)

A splitting method of nonstiff order p admits the (formal) expansion

$$\mathcal{L}_F(t, v) = \sum_{k=1}^p \sum_{\substack{\mu \in \mathbb{N}^k \\ |\mu| \leq p-k}} \frac{1}{\mu!} t^{k+|\mu|} C_{k\mu} \prod_{\ell=1}^k ad_{D_A}^{\mu_\ell}(D_B) e^{tD_A} v + R_{p+1}(t, v),$$

$$C_{k\mu} = \sum_{\lambda \in \Lambda_k} \alpha_\lambda \prod_{\ell=1}^k b_{\lambda_\ell} c_{\lambda_\ell}^{\mu_\ell} - \prod_{\ell=1}^k \frac{1}{\mu_\ell + \dots + \mu_k + k - \ell + 1}.$$

Main tools. Calculus of Lie derivatives, Gröbner–Alekseev formula.

Remark. Further considerations show that resulting (redundant) stiff order conditions $C_{k\mu} = 0$ coincide with nonstiff order conditions.

Theoretical justification

Remarks.

- Application of local error representation to different classes of (non)linear evolution equations such as Schrödinger equations or diffusion-reaction systems requires characterisation of domains of iterated Lie-commutators (regularity and consistency requirements).
- In connection with **Schrödinger equations**, it is often justified to assume that exact solution is regular. For **linear equations** versus **nonlinear equations**, the regularity requirements are

$$D = H^p(\Omega), \quad D = H^{2p}(\Omega).$$

- For sufficiently regular solutions (bounded in D), above local error representation implies

$$\mathcal{L}_F(\tau, v) = \mathcal{S}_F(\tau, v) - \mathcal{E}_F(\tau, v) = \mathcal{O}(\tau^{p+1}).$$

Provided that $\hat{p} > p$, this justifies use of local error estimator

$$\text{err}_{\text{local}} = \|u_1 - \hat{u}_1\|_X = \mathcal{O}(\tau^{p+1}).$$

Global error estimate (Full discretisations)

Discretisation. Full discretisation of **nonlinear Schrödinger equations** (GPE) by **high-order variable stepsize time-splitting methods** combined with **pseudo-spectral methods** (Fourier, Sine, Hermite).

Theorem (Th. 2012)

Provided that exact solution remains bounded in fractional power space X_β defined by principal linear part for $\beta \geq p$, global error estimate holds

$$\|u_{NM} - u(t_N)\|_X \leq C \left(\|u_0 - u(0)\|_X + \tau_{\max}^p + M^{-q} \right).$$

Extensions.

- Time-dependent Gross–Pitaevskii equations with additional rotation term, see HOFSTÄTTER, KOCH, TH. (2014).
- Multi-revolution composition time-splitting pseudo-spectral methods for highly oscillatory problems (with CHARTIER, MÉHATS).

Defect-based a posteriori local error estimators

Examples and theoretical basis

Alternative approach

Approach. Consider nonlinear evolution equation and deduce **evolution equation of similar form** for splitting operator

$$\begin{aligned}\frac{d}{dt} \mathcal{E}_F(t, \cdot) &= (D_A + D_B) \mathcal{E}_F(t, \cdot), \\ \mathcal{S}_F(t, \cdot) &= \prod_{j=1}^s e^{a_{s+1-j} t D_A} e^{b_{s+1-j} t D_B}.\end{aligned}$$

Employ Gröbner–Aleksiev formula and suitable further expansion.

- For low-order methods, obtain **compact local error representation**

$$\begin{aligned}\mathcal{L}_F(t, \cdot) &= \int_0^t \int_0^{\tau_1} e^{\tau_1 D_A} e^{\tau_2 D_B} [D_A, D_B] e^{(\tau_1 - \tau_2) D_B} e^{(t - \tau_1) D_F} d\tau_2 d\tau_1 \\ &= \int_0^t \int_0^{\tau_1} \partial_2 \mathcal{E}_F(t - \tau_1, \mathcal{S}_F(\tau_1, \cdot)) \partial_2 \mathcal{E}_B(\tau_1 - \tau_2, \mathcal{E}_A(\tau_1, \cdot)) [B, A] (\mathcal{E}_B(\tau_2, \mathcal{E}_A(\tau_1, \cdot))) d\tau_2 d\tau_1.\end{aligned}$$

- As rigorous extension to higher-order splitting methods becomes **highly involved**, study linear case and use formal extension.

Remark. Related approach studied in context of (non)linear Schrödinger equations in **semi-classical regime**, see DESCOMBES, TH, (2010, 2012).

A posteriori local error estimators (Linear case)

Approach. Consider **linear evolution equation**

$$\begin{cases} \mathcal{E}'_F(t) = (A+B) \mathcal{E}_F(t), & t \in (0, T), \\ \mathcal{E}_F(0) = I. \end{cases}$$

For **splitting method**, deduce evolution equation of similar form

$$\begin{cases} \mathcal{S}'_F(t) = (A+B) \mathcal{S}_F(t) + \mathcal{D}_F(t), & t \in (0, T), \\ \mathcal{S}_F(0) = I. \end{cases}$$

Local error $\mathcal{L}_F = \mathcal{S}_F - \mathcal{E}_F$ satisfies evolution equation

$$\begin{cases} \mathcal{L}'_F(t) = (A+B) \mathcal{L}_F(t) + \mathcal{D}_F(t), & t \in (0, T), \\ \mathcal{L}_F(0) = 0. \end{cases}$$

Employ variation-of-constants formula to obtain **integral representation** for **local error** involving **defect**

$$\mathcal{L}_F(t) = \int_0^t e^{(t-\tau)(A+B)} \mathcal{D}_F(\tau) \, d\tau.$$

A posteriori local error estimators (Linear case)

Approach. Recall **integral representation** for local error

$$\mathcal{L}_F(t) = \int_0^t \underbrace{e^{(t-\tau)(A+B)} \mathcal{D}_F(\tau)}_{=f(\tau)} d\tau.$$

Apply **Hermite quadrature approximation** and use that validity of order conditions implies $f(0) = \dots = f^{(p-1)}(0) = 0$

$$\underbrace{\sum_{\ell=0}^{p-1} \omega_{\ell} t^{\ell+1} f^{(\ell)}(0)}_{=0} + \frac{1}{p+1} t f(t) - \int_0^t f(\tau) d\tau = \mathcal{O}(t^{p+2}).$$

For any **splitting method of order p** , obtain **asymptotically correct defect-based a posteriori local error estimator**

$$\mathcal{P}_F(t) = \frac{1}{p+1} t \mathcal{D}_F(t), \quad \mathcal{P}_F(t) - \mathcal{L}_F(t) = \mathcal{O}(t^{p+2}).$$

A posteriori local error estimators

Result. Asymptotically correct a posteriori local error estimator associated with splitting method of order p given by

$$\mathcal{S}_k^m(t) = \prod_{j=k}^m e^{b_j t B} e^{a_j t A}, \quad \mathcal{S}_F(t) = \mathcal{S}_1^s(t),$$

$$\mathcal{S}_F(t) - \mathcal{E}_F(t) = \mathcal{O}(t^{p+1}),$$

$$\mathcal{D}_F = \sum_{k=1}^s \mathcal{S}_k^s a_k A \mathcal{S}_1^{k-1} + \sum_{k=1}^{s-1} \mathcal{S}_{k+1}^s b_k B \mathcal{S}_1^k - (A + (1 - b_s) B) \mathcal{S}_F(t),$$

$$\mathcal{D}_F = \frac{1}{p+1} t \mathcal{D}_F, \quad \mathcal{D}_F(t) - \mathcal{L}_F(t) = \mathcal{O}(t^{p+2}).$$

Extension to **nonlinear evolution equations** by calculus of Lie-derivatives.

Theoretical analysis. In context of linear Schrödinger equations, rigorous analysis given in AUZINGER, KOCH, TH. (2012, 2014). Corresponding result for nonlinear case deduced in AUZINGER, HOFSTÄTTER, KOCH, TH. (2015) for second-order Strang splitting method.

Special case (Lie–Trotter splitting method)

Special case. A posteriori local error estimator for **Lie–Trotter splitting method** applied to **linear evolution equation** given by

$$\mathcal{P}_F(t, v) = \frac{1}{2} t \mathcal{D}_F(t, v), \quad \mathcal{D}_F(t, v) = (e^{tB} e^{tA} A - A e^{tB} e^{tA}) v.$$

Extension to **nonlinear case** yields

$$\mathcal{D}_F(t, v) = \partial_2 \mathcal{E}_B(t, \mathcal{E}_A(t, v)) \partial_2 \mathcal{E}_A(t, v) A v - A \mathcal{E}_B(t, \mathcal{E}_A(t, v)).$$

Explanation. Extension by formal calculus of Lie-derivatives implies $\mathcal{D}_F(t, v) = D_A e^{tDA} e^{tDB} v - e^{tDA} e^{tDB} D_A v$ and

$$G(v) = e^{tDA} e^{tDB} v = \mathcal{E}_B(t, \mathcal{E}_A(t, v)), \quad G'(v) = \partial_2 \mathcal{E}_B(t, \mathcal{E}_A(t, v)) \partial_2 \mathcal{E}_A(t, v), \\ e^{tDA} e^{tDB} D_A v = A \mathcal{E}_B(t, \mathcal{E}_A(t, v)), \quad D_A e^{tDA} e^{tDB} v = G'(v) A v = \partial_2 \mathcal{E}_B(t, \mathcal{E}_A(t, v)) \partial_2 \mathcal{E}_A(t, v) A v.$$

Remark. Improved approximation $\mathcal{S}_F(t, \cdot) - \mathcal{P}_F(t, \cdot) = \mathcal{E}_F(t, \cdot) + \mathcal{O}(t^{p+2})$.

Realisation and computational effort. Realisation for nonlinear Schrödinger equations (**Gross–Pitaevskii equation**) straightforward. Computational effort comparable with splitting pair Lie/Strang (two additional applications of A required, FFT)

$$\mathcal{P}(t, v) = e^{-it(U+\theta|w|^2)} \left(A w - i\theta t (A w |w|^2 + \overline{A w} w^2) \right) - A e^{-it(U+\theta|w|^2)} w, \quad w = e^{tA} v.$$

Explanation. With $G(v) = e^{-it(U+\theta|w|^2)} w$, $G'(v) = e^{-it(U+\theta|w|^2)} \left(e^{tA(\cdot)} - i\theta t (\overline{w} e^{tA(\cdot)} + w \overline{e^{tA(\cdot)}}) \right) w$ obtain

$$e^{tDA} e^{tDB} D_A v = A e^{-it(U+\theta|w|^2)} w, \quad D_A e^{tDA} e^{tDB} v = e^{-it(U+\theta|w|^2)} \left(A w - i\theta t (A w |w|^2 + \overline{A w} w^2) \right).$$

Associated approximations with negligible additional computational cost

Examples and theoretical basis

Novel approach

Approach. Consider nonlinear evolution equation

$$\begin{cases} u'(t) = A(u(t)) + B(u(t)), & t \in (0, T), \\ u(0) = u_0. \end{cases}$$

Realise higher-order splitting method in straightforward manner

$u = u_n$

for $j = 1 : s$

$u = \mathcal{E}_A(a_j \tau_n, u)$ Solution of subproblem $u'(t) = A(u(t))$

$u = \mathcal{E}_B(b_j \tau_n, u)$ Solution of subproblem $u'(t) = B(u(t))$

end

$u_{n+1} = u$

Use suitable **linear combination of intermediate values** to compute associated approximation that serves as **local error estimator**.

Novel approach

Schrödinger equations. Consider splitting method by BLANES, MOAN

$$p = 4, \quad s = 7.$$

Associated **third-order approximation** obtained by certain linear combination of intermediate values yields local error estimator

$$u = u_n$$

$$u_{\text{Estimator}} = \alpha_0 u$$

for $j = 1 : s$

$$u = \mathcal{E}_A(a_j \tau_n, u)$$

$$u_{\text{Estimator}} = u_{\text{Estimator}} + \alpha_{2j-1} u$$

$$u = \mathcal{E}_B(b_j \tau_n, u)$$

$$u_{\text{Estimator}} = u_{\text{Estimator}} + \alpha_{2j} u$$

end

$$u_{n+1} = u$$

$$\text{Local error estimator} = u - u_{\text{Estimator}}$$

Parabolic equations. Consider instead splitting method by YOSHIDA with complex coefficients and melt two subsequent time steps ($p = 4, s = 7$).

Novel approach

Benefit. Compared to approaches based on embedded splitting methods or defect-based local error estimators, novel approach leads to local error estimators with **negligible additional computational cost**.

Open questions.

- Provide coefficients for favourable higher-order splitting methods.
- Numerical tests confirm stability of associated approximations.
Rigorous argument?

Numerical examples

Local error control

Local error control. Use of **local error control** to adjust time stepsize

$$\tau_{\text{optimal}} = \tau_{\text{current}} \cdot \min \left(\alpha_{\max}, \max \left(\alpha_{\min}, \sqrt[p+1]{\alpha \cdot \frac{\text{tol}}{\text{err}_{\text{local}}}} \right) \right),$$
$$\alpha_{\max} = 1.5, \quad \alpha_{\min} = 0.2, \quad \alpha = 0.25,$$

enhances **reliability** and **efficiency** of time integration.

Illustration (Schrödinger equation)

Test equation (see BAO ET AL.). Consider **nonlinear Schrödinger equation** under harmonic potential ($d = 1$, $\omega = 2$, $\vartheta = 1$)

$$i \partial_t \psi(x, t) = \left(-\frac{1}{2} \varepsilon \Delta + \frac{1}{\varepsilon} U(x) + \frac{1}{\varepsilon} \vartheta |\psi(x, t)|^2 \right) \psi(x, t).$$

Small value of (semi-classical) parameter $\varepsilon > 0$ causes high oscillations in initial condition and solution

$$\psi(x, 0) = \varrho_0(x) e^{i \frac{1}{\varepsilon} \sigma_0(x)}, \quad \varrho_0(x) = e^{-x^2}, \quad \sigma_0(x) = -\ln(e^x + e^{-x}).$$

Use **Fourier spectral space discretisation** combined with **fourth-order time-splitting method** by BLANES & MOAN ($x \in [-8, 8]$, $M = 8192$, $t \in [0, 3]$).

Illustration (Solution behaviour for $\varepsilon = 10^{-2}$)

First observation. Even a simple local error control for second-order Strang splitting method based on first-order Lie–Trotter splitting method is useful to **enhance reliability!** See Movie.

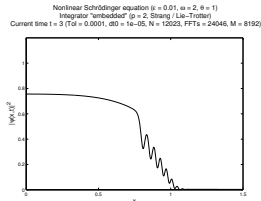
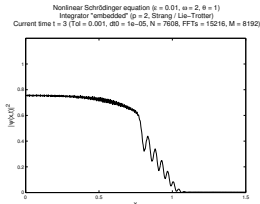
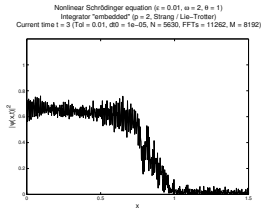
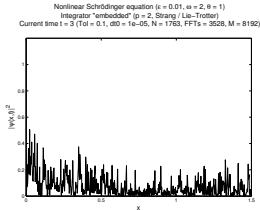


Illustration (Global error)

Expectation. Use of higher-order methods will enhance efficiency.

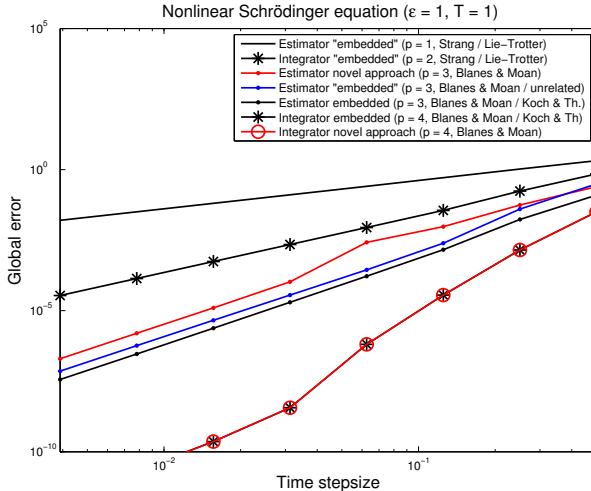
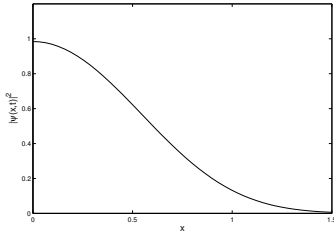


Illustration ($\varepsilon = 1$)

Comparison. Compare approach based on embedded splitting methods with novel approach. Obtain expected results for $\varepsilon = 1$, $T = 10$, $\text{Tol} = 10^{-4}$.

- **Higher-order method superior** to low-order method (e.g. with respect to number of FFT transforms).

Nonlinear Schrödinger equation ($\varepsilon = 1$, $\omega = 2$, $\theta = 1$)
Integrator "embedded" ($p = 2$, Strang / Lie-Trotter)
Current time $t = 20$ (Tol = 0.0001, dt0 = 0.001, N = 4830, FFTs = 9660, M = 500)



Nonlinear Schrödinger equation ($\varepsilon = 1$, $\omega = 2$, $\theta = 1$)
Integrator novel approach ($p = 4$, Blanes & Moan)
Current time $t = 20$ (Tol = 0.0001, dt0 = 0.001, N = 258, FFTs = 1548, M = 500)

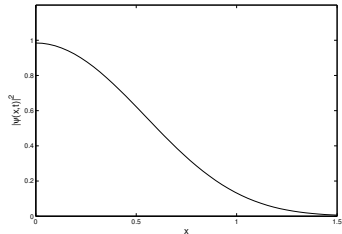


Illustration ($\varepsilon = 1$)

Comparison. Compare approach based on embedded splitting methods with novel approach. Obtain expected results for $\varepsilon = 1$, $T = 10$, $\text{Tol} = 10^{-4}$.

- **Good performance of novel approach** in comparison with embedded method (e.g. with respect to number of FFT transforms).

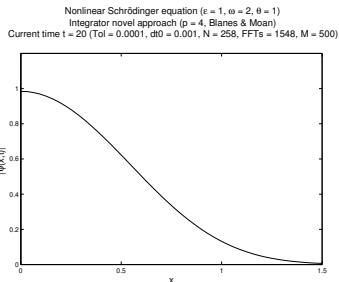
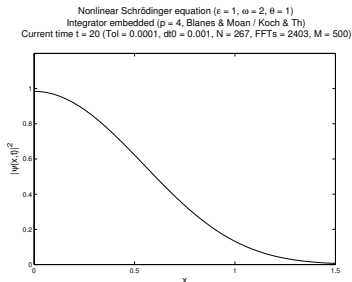
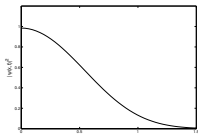


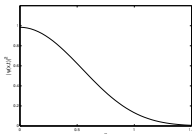
Illustration ($\varepsilon = 1$)

Best performance. Observe best performance for novel approach (FFTs).

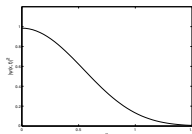
Nonlinear Schrödinger equation ($\nu = 1, \alpha = 2, \theta = 1$)
Integrator "embedded" ($p = 4$, Blanes & Moan / unrefined)
Current time $t = 20$ (Tol = 0.01, $\delta_0 = 0.001$, $N = 98$, FFTs = 864, $M = 500$)



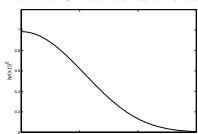
Nonlinear Schrödinger equation ($\nu = 1, \alpha = 2, \theta = 1$)
Integrator "embedded" ($p = 4$, Blanes & Moan / unrefined)
Current time $t = 20$ (Tol = 0.001, $\delta_0 = 0.001$, $N = 167$, FFTs = 1503, $M = 500$)



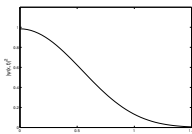
Nonlinear Schrödinger equation ($\nu = 1, \alpha = 2, \theta = 1$)
Integrator "embedded" ($p = 4$, Blanes & Moan / unrefined)
Current time $t = 20$ (Tol = 0.0001, $\delta_0 = 0.001$, $N = 255$, FFTs = 3037, $M = 500$)



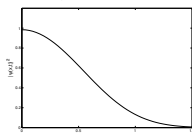
Nonlinear Schrödinger equation ($\nu = 1, \alpha = 2, \theta = 1$)
Integrator embedded ($p = 4$, Blanes & Moan / Koch & Th)



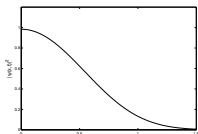
Nonlinear Schrödinger equation ($\nu = 1, \alpha = 2, \theta = 1$)
Integrator embedded ($p = 4$, Blanes & Moan / Koch & Th)
Current time $t = 20$ (Tol = 0.001, $\delta_0 = 0.001$, $N = 145$, FFTs = 1305, $M = 500$)



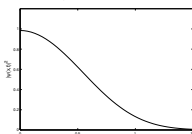
Nonlinear Schrödinger equation ($\nu = 1, \alpha = 2, \theta = 1$)
Integrator embedded ($p = 4$, Blanes & Moan / Koch & Th)
Current time $t = 20$ (Tol = 0.0001, $\delta_0 = 0.001$, $N = 267$, FFTs = 2403, $M = 500$)



Nonlinear Schrödinger equation ($\nu = 1, \alpha = 2, \theta = 1$)
Integrator novel approach ($p = 4$, Blanes & Moan)



Nonlinear Schrödinger equation ($\nu = 1, \alpha = 2, \theta = 1$)
Integrator novel approach ($p = 4$, Blanes & Moan)
Current time $t = 20$ (Tol = 0.001, $\delta_0 = 0.001$, $N = 146$, FFTs = 854, $M = 500$)



Nonlinear Schrödinger equation ($\nu = 1, \alpha = 2, \theta = 1$)
Integrator novel approach ($p = 4$, Blanes & Moan)
Current time $t = 20$ (Tol = 0.0001, $\delta_0 = 0.001$, $N = 258$, FFTs = 1548, $M = 500$)

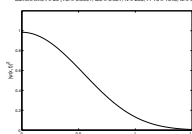
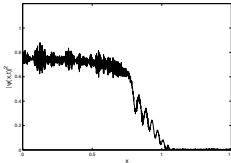


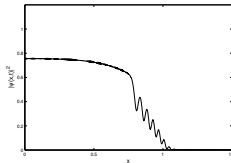
Illustration ($\varepsilon = 10^{-2}$)

Observation. Time integration of test equation for $\varepsilon = 10^{-2}$ is **delicate** task. Number of time steps does not necessarily increase for smaller **tolerances**. Influence of choice of initial time stepsize? Improvement of local error control in this situation?

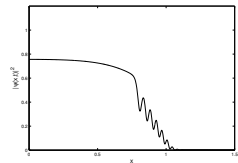
Nonlinear Schrödinger equation ($\kappa = 0.01, \omega = 2, \theta = 1$)
Integrator embedded ($p = 4$, Blanes & Moan / Koch & Th)
Current time $t = 3$ (Tol = 0.01, ddt = $1e-05$, N = 2370, FFTs = 21384, M = 8192)



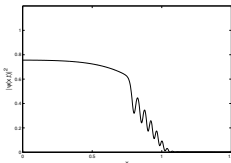
Nonlinear Schrödinger equation ($\kappa = 0.01, \omega = 2, \theta = 1$)
Integrator embedded ($p = 4$, Blanes & Moan / Koch & Th)
Current time $t = 3$ (Tol = 0.001, ddt = $1e-05$, N = 2468, FFTs = 22239, M = 8192)



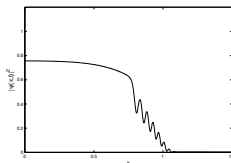
Nonlinear Schrödinger equation ($\kappa = 0.01, \omega = 2, \theta = 1$)
Integrator embedded ($p = 4$, Blanes & Moan / Koch & Th)
Current time $t = 3$ (Tol = 0.0001, ddt = $1e-05$, N = 2287, FFTs = 20601, M = 8192)



Nonlinear Schrödinger equation ($\kappa = 0.01, \omega = 2, \theta = 1$)
Integrator novel approach ($p = 4$, Blanes & Moan)
Current time $t = 3$ (Tol = 0.01, ddt = $1e-05$, N = 5765, FFTs = 34752, M = 8192)



Nonlinear Schrödinger equation ($\kappa = 0.01, \omega = 2, \theta = 1$)
Integrator novel approach ($p = 4$, Blanes & Moan)
Current time $t = 3$ (Tol = 0.001, ddt = $1e-05$, N = 5069, FFTs = 30420, M = 8192)



Nonlinear Schrödinger equation ($\kappa = 0.01, \omega = 2, \theta = 1$)
Integrator novel approach ($p = 4$, Blanes & Moan)
Current time $t = 3$ (Tol = 0.0001, ddt = $1e-05$, N = 5825, FFTs = 52950, M = 8192)

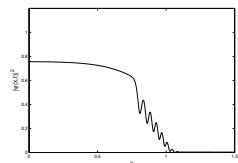


Illustration (Gray–Scott equations)

Test equation . Consider system of **diffusion–reaction equations** ($d = 2$)

$$\begin{cases} \partial_t u(x, t) = (D_u \Delta - \alpha) u(x, t) - u(x, t) (v(x, t))^2 + \alpha, \\ \partial_t v(x, t) = (D_v \Delta - \beta) v(x, t) + u(x, t) (v(x, t))^2. \end{cases}$$

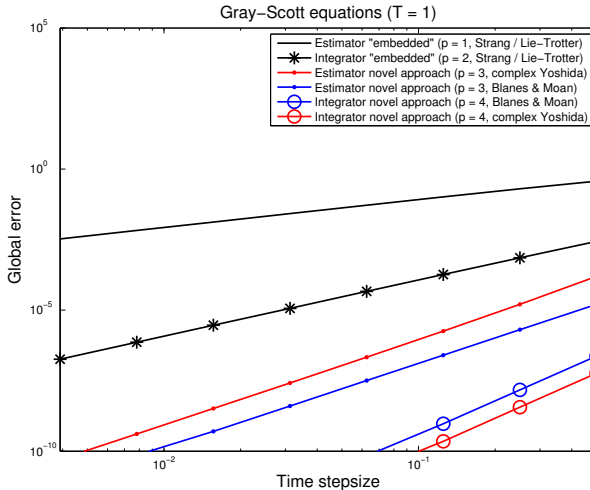
Special choice of parameters causes formation of patterns over long times

$$D_u = 2D_v = 0.16, \quad f = 0.035, \quad k = 0.06, \quad \alpha = f, \quad \beta = f + k,$$

see also NICOLAS P. ROUGIER. Use **Fourier spectral space discretisation** combined with **fourth-order complex time-splitting method** by YOSHIDA ($x \in [-75, 75]$, $M = 150 \times 150$).

Movie
Formation of patterns

Illustration (Global error)



Conclusions and future work

Conclusions.

- Adaptivity in time essential for reliable and efficient numerical simulations.
- Novel approach for time-splitting methods provides local error estimators with negligible additional cost.

Open questions.

- Theoretical understanding of novel local error estimators (favourable stability behaviour). Design of higher-order schemes.
- Detailed study of local error control for semi-classical Schrödinger equation to understand unexpected behaviour.

Thank you!