

# Three approaches for the design of adaptive time-splitting methods

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# Theme

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

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

**Linear case.** For linear evolution equations, exponential operator splitting methods can be cast into general form (time grid points  $0 = t_0 < \dots < t_N \leq T$  and stepsizes  $\tau_{n-1} = t_n - t_{n-1}$ )

$$u'(t) = Au(t) + Bu(t), \quad t \in (0, T),$$

$$u_n = \prod_{j=1}^s e^{b_j \tau_{n-1} B} e^{a_j \tau_{n-1} A} u_{n-1} \approx u(t_n) = e^{\tau_{n-1}(A+B)} u(t_{n-1}), \quad n \in \{1, \dots, N\}.$$

## 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).

# Outline

**Main theme.** Design and theoretical analysis of local error estimators for exponential operator splitting methods.

## Outline.

- Splitting methods
- Local error estimators
  - Embedded splitting methods
  - Defect-based local error estimators
  - Associated approximations
- Numerical illustrations
  - Nonlinear Schrödinger equations
  - Diffusion-reaction systems

# Exponential operator splitting methods for nonlinear evolution equations

Calculus of Lie-derivatives and Gröbner–Alekseev formula

# 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.

# Calculus of Lie-derivatives

**Formal calculus.** Calculus of Lie-derivatives is suggestive of less involved linear case, see for instance HAIRER, LUBICH, WANNER (2002) and SANZ-SERNA, CALVO (1994).

**Problem.** Consider **nonlinear evolution equation** on Banach space involving (unbounded) nonlinear operator  $F : D(F) \subseteq X \rightarrow X$

$$u'(t) = F(u(t)), \quad t \in (0, T).$$

Employ **formal notation** for exact solution

$$u(t) = \mathcal{E}_F(t, u(0)) = e^{tD_F} u(0), \quad t \in [0, T].$$

**Evolution operator, Lie-derivative.** For (unbounded) nonlinear operator  $G : D(G) \subseteq X \rightarrow X$  define evolution operator and Lie-derivative by

$$e^{tD_F} G v = G(\mathcal{E}_F(t, v)), \quad D_F G v = G'(v) F(v).$$

**Remark.** Definition of Lie-derivative is natural extension of identity  $L = \frac{d}{dt} \Big|_{t=0} e^{tL}$

$$\frac{d}{dt} \Big|_{t=0} e^{tD_F} G v = \frac{d}{dt} \Big|_{t=0} G(\mathcal{E}_F(t, v)) = G'(\mathcal{E}_F(t, v)) F(\mathcal{E}_F(t, v)) \Big|_{t=0} = G'(v) F(v) = D_F G v.$$



# 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  $j \in \{1, \dots, 4\}$ .

# Nonlinear variation-of-constants formula

## Theorem (Gröbner–Aleksseev formula)

*The solutions to the autonomous problem*

$$\begin{cases} \frac{d}{dt} \mathcal{E}_G(t - t_0, v) = G(\mathcal{E}_G(t - t_0, v)), & t_0 \leq t \leq T, \\ \mathcal{E}_G(0, v) = v, \end{cases}$$

*and the related non-autonomous problem*

$$\begin{cases} \frac{d}{dt} \mathcal{E}_{G+R}(t, t_0, v) = G(\mathcal{E}_{G+R}(t, t_0, v)) + R(t), & t_0 \leq t \leq T, \\ \mathcal{E}_{G+R}(t_0, t_0, v) = v, \end{cases}$$

*satisfy the integral relation*

$$\mathcal{E}_{G+R}(t, t_0, v) - \mathcal{E}_G(t - t_0, v) = \int_{t_0}^t \partial_2 \mathcal{E}_G(t - \tau, \mathcal{E}_{G+R}(\tau, t_0, v)) R(\tau) d\tau.$$

## Proof

The fundamental identity

$$\begin{aligned}\partial_2 \mathcal{E}_G(t, v) G(v) &= \partial_2 \mathcal{E}_G(t, \mathcal{E}_G(0, v)) G(\mathcal{E}_G(0, v)) = \left. \frac{d}{ds} \right|_{s=0} \mathcal{E}_G(t, \mathcal{E}_G(s, v)) \\ &= \left. \frac{d}{ds} \right|_{s=0} \mathcal{E}_G(t+s, v) = G(\mathcal{E}_G(t, v))\end{aligned}$$

implies  $\partial_2 \mathcal{E}_G(t-\tau, w) G(w) = G(\mathcal{E}_G(t-\tau, w))$  for  $w = \mathcal{E}_{G+R}(\tau, t_0, v)$ . A brief calculation shows the stated relation

$$\begin{aligned}\mathcal{E}_{G+R}(t, t_0, v) - \mathcal{E}_G(t-t_0, v) &= \mathcal{E}_G(t-\tau, \mathcal{E}_{G+R}(\tau, t_0, v)) \Big|_{\tau=t_0}^t \\ &= \int_{t_0}^t \frac{d}{d\tau} \mathcal{E}_G(t-\tau, \mathcal{E}_{G+R}(\tau, t_0, v)) d\tau \\ &= \int_{t_0}^t \left( \partial_2 \mathcal{E}_G(t-\tau, \mathcal{E}_{G+R}(\tau, t_0, v)) (G(\mathcal{E}_{G+R}(\tau, t_0, v)) + R(\tau)) \right. \\ &\quad \left. - G(\mathcal{E}_G(t-\tau, \mathcal{E}_{G+R}(\tau, t_0, v))) \right) d\tau \\ &= \int_{t_0}^t \partial_2 \mathcal{E}_G(t-\tau, \mathcal{E}_{G+R}(\tau, t_0, v)) R(\tau) d\tau.\end{aligned}$$



# 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  $X_\beta$  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 in present situation 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  $\mathcal{D}_F(t, v) = D_A e^{tD_A} e^{tD_B} v - e^{tD_A} e^{tD_B} D_A v$  and

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

**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

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

# 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 ( $d = 1$ ,  $\omega = 2$ ,  $\vartheta = 1$ )

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

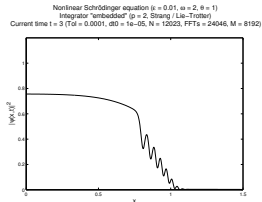
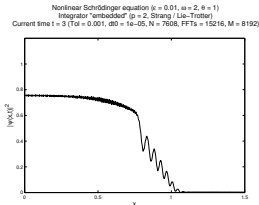
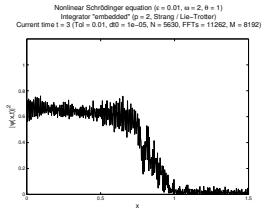
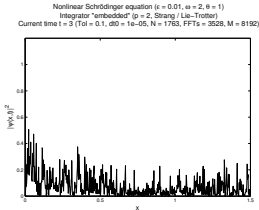
**Small value of (semi-classical) parameter**  $\varepsilon > 0$  causes high oscillations in initial condition and solution

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

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

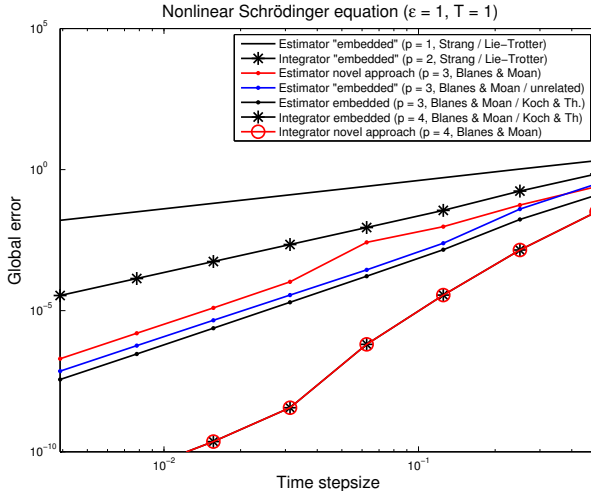
# Illustration (Solution behaviour for $\varepsilon = 10^{-2}$ )

**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.

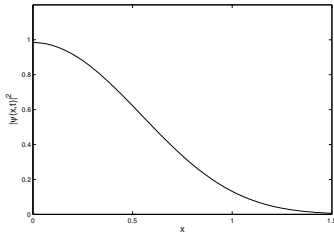


# Illustration ( $\varepsilon = 1$ )

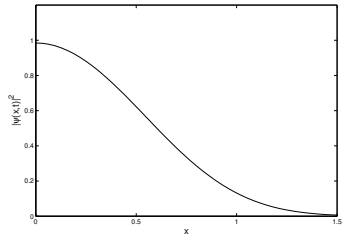
**Comparison.** Compare approach based on embedded splitting methods with novel approach. Obtain expected results for  $\varepsilon = 1$ ,  $T = 10$ ,  $\text{Tol} = 10^{-4}$ .

- 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)



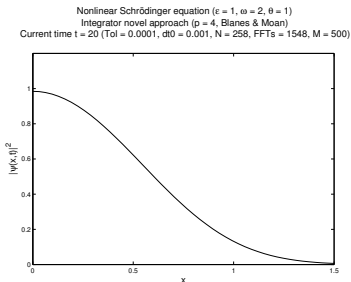
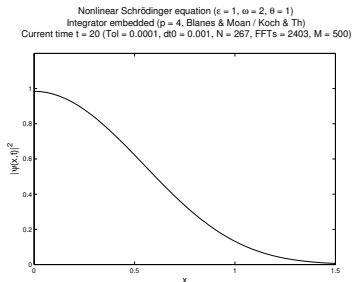
Nonlinear Schrödinger equation ( $\varepsilon = 1$ ,  $\omega = 2$ ,  $\theta = 1$ )  
Integrator novel approach ( $p = 4$ , Blanes & Moan)  
Current time  $t = 20$  (Tol = 0.0001, dt0 = 0.001, N = 258, FFTs = 1548, M = 500)



# Illustration ( $\varepsilon = 1$ )

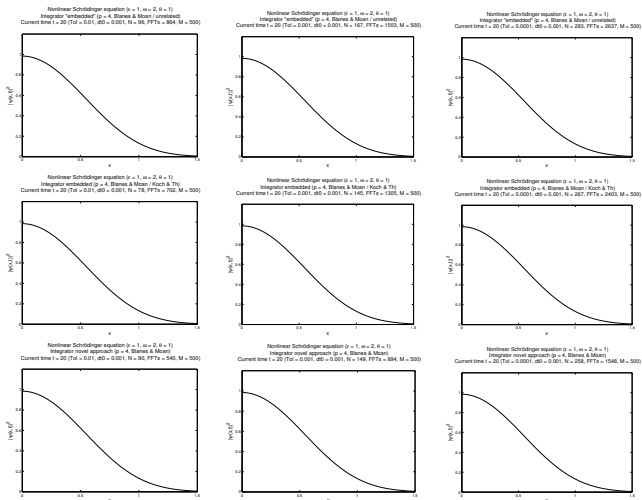
**Comparison.** Compare approach based on embedded splitting methods with novel approach. Obtain expected results for  $\varepsilon = 1$ ,  $T = 10$ ,  $\text{Tol} = 10^{-4}$ .

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



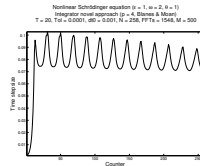
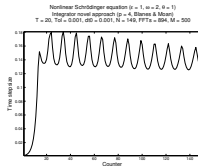
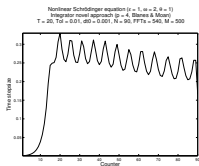
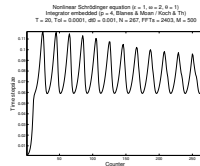
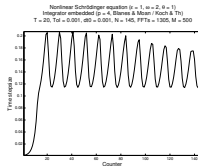
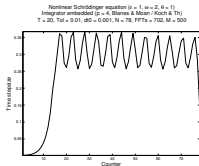
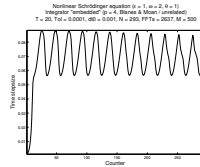
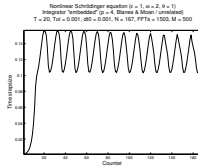
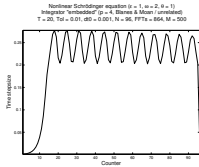
# Illustration ( $\varepsilon = 1$ )

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



# Illustration ( $\varepsilon = 1$ )

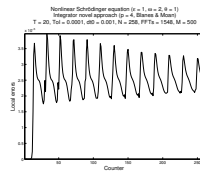
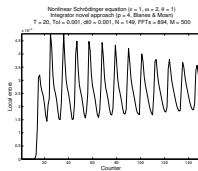
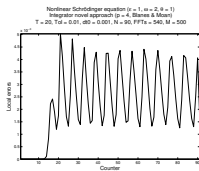
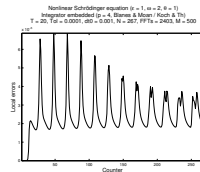
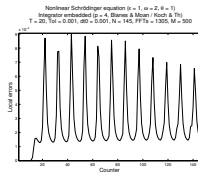
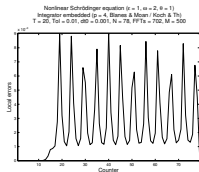
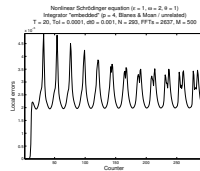
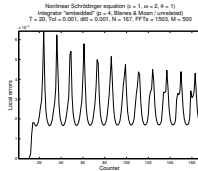
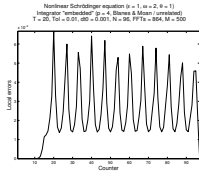
## Time stepsizes. Compare sequences of time stepsizes.





# Illustration ( $\varepsilon = 1$ )

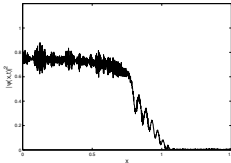
## Local errors. Compare local errors.



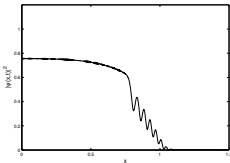
# Illustration ( $\varepsilon = 10^{-2}$ )

**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?

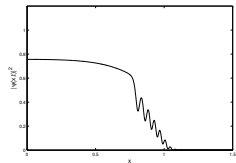
Nonlinear Schrödinger equation ( $\kappa = 0.01, \omega = 2, \theta = 1$ )  
Integrator embedded ( $p = 4$ , Blanes & Moan / Koch & Th)  
Current time  $t = 3$  (Tol = 0.01, ddt =  $1e-05$ , N = 2370, FFTs = 21384, M = 8192)



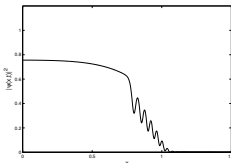
Nonlinear Schrödinger equation ( $\kappa = 0.01, \omega = 2, \theta = 1$ )  
Integrator embedded ( $p = 4$ , Blanes & Moan / Koch & Th)  
Current time  $t = 3$  (Tol = 0.001, ddt =  $1e-05$ , N = 2468, FFTs = 23239, M = 8192)



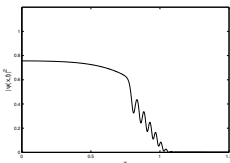
Nonlinear Schrödinger equation ( $\kappa = 0.01, \omega = 2, \theta = 1$ )  
Integrator embedded ( $p = 4$ , Blanes & Moan / Koch & Th)  
Current time  $t = 3$  (Tol = 0.0001, ddt =  $1e-05$ , N = 2287, FFTs = 20601, M = 8192)



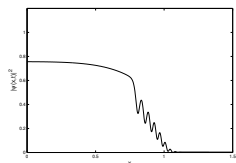
Nonlinear Schrödinger equation ( $\kappa = 0.01, \omega = 2, \theta = 1$ )  
Integrator novel approach ( $p = 4$ , Blanes & Moan)  
Current time  $t = 3$  (Tol = 0.01, ddt =  $1e-05$ , N = 5765, FFTs = 34752, M = 8192)



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)



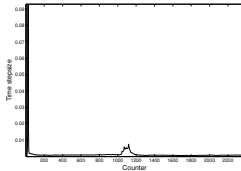
Nonlinear Schrödinger equation ( $\kappa = 0.01, \omega = 2, \theta = 1$ )  
Integrator novel approach ( $p = 4$ , Blanes & Moan)  
Current time  $t = 3$  (Tol = 0.0001, ddt =  $1e-05$ , N = 5825, FFTs = 52950, M = 8192)



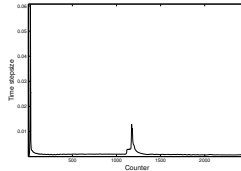
# Illustration ( $\varepsilon = 10^{-2}$ )

**Time stepsizes.** Compare sequences of time stepsizes.

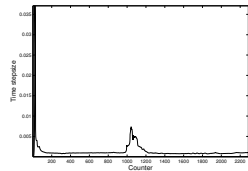
Nonlinear Schrödinger equation ( $\varepsilon = 0.01, \omega = 2, \theta = 1$ )  
Integrator embedded ( $p = 4$ , Blanes & Moan / Koch & Th)  
 $T = 3, \text{Tol} = 0.01, \text{d0} = 1\text{e-}05, N = 2370, \text{FFTs} = 21384, M = 8192$



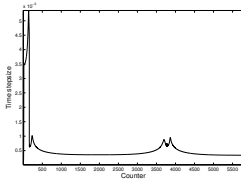
Nonlinear Schrödinger equation ( $\varepsilon = 0.01, \omega = 2, \theta = 1$ )  
Integrator embedded ( $p = 4$ , Blanes & Moan / Koch & Th)  
 $T = 3, \text{Tol} = 0.001, \text{d0} = 1\text{e-}05, N = 2468, \text{FFTs} = 22239, M = 8192$



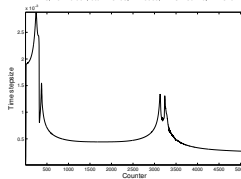
Nonlinear Schrödinger equation ( $\varepsilon = 0.01, \omega = 2, \theta = 1$ )  
Integrator embedded ( $p = 4$ , Blanes & Moan / Koch & Th)  
 $T = 3, \text{Tol} = 0.0001, \text{d0} = 1\text{e-}05, N = 2287, \text{FFTs} = 20601, M = 8192$



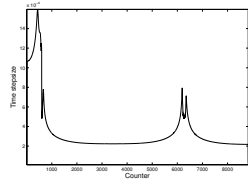
Nonlinear Schrödinger equation ( $\varepsilon = 0.01, \omega = 2, \theta = 1$ )  
Integrator novel approach ( $p = 4$ , Blanes & Moan)  
 $T = 3, \text{Tol} = 0.01, \text{d0} = 1\text{e-}05, N = 5788, \text{FFTs} = 34752, M = 8192$



Nonlinear Schrödinger equation ( $\varepsilon = 0.01, \omega = 2, \theta = 1$ )  
Integrator novel approach ( $p = 4$ , Blanes & Moan)  
 $T = 3, \text{Tol} = 0.001, \text{d0} = 1\text{e-}05, N = 5069, \text{FFTs} = 30420, M = 8192$

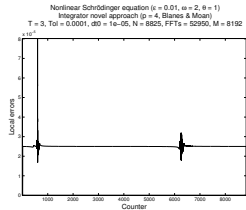
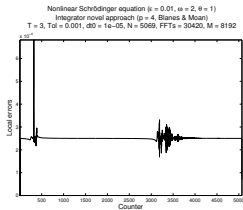
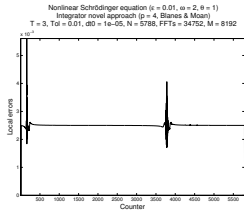
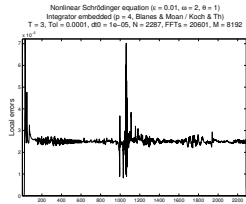
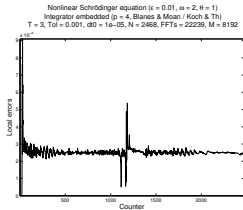
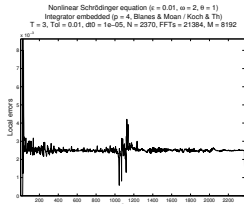


Nonlinear Schrödinger equation ( $\varepsilon = 0.01, \omega = 2, \theta = 1$ )  
Integrator novel approach ( $p = 4$ , Blanes & Moan)  
 $T = 3, \text{Tol} = 0.0001, \text{d0} = 1\text{e-}05, N = 8825, \text{FFTs} = 52950, M = 8192$



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

## Local errors. Compare local errors.



# Illustration (Gray–Scott equations)

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

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

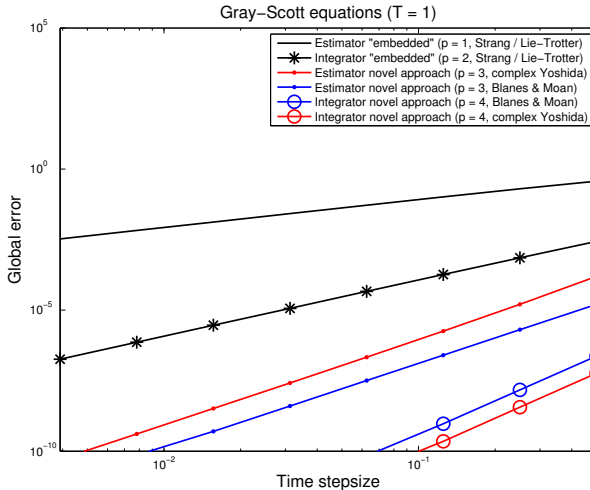
**Special choice of parameters** causes formation of patterns over long times

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

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!**