

Source Terms

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Seminar Course - Fundamentals of Wave Simulation - Solving Hyperbolic Systems of PDEs

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Our reference equation is

$$q_t + f(q)_x = \psi(q) \quad (1)$$

where

- the homogeneous equation $q_t + f(q)_x = 0$ is **hyperbolic**
- $\psi(q)$ (the **source terms**) don't depend on derivatives of q
 - $\Rightarrow q_t = \psi(q)$ is an independent system of ODEs

A standard example that will be used to illustrate the following numerical methods is the **advection-reaction equation**

$$q_t + \bar{u}q_x = -\beta q. \quad (2)$$

It can be seen as the model for the transport along a flow of a radioactive substance, where

- β is the **decay rate**
- \bar{u} is the (constant) **transport speed**
- $q(x, 0) = \hat{q}(x)$ is the **initial condition**.

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Exact solution

Along the characteristic $\frac{dx}{dt} = \bar{u}$ we have $\frac{dq}{dt} = -\beta q$ and it follows that

$$q(x, t) = e^{-\beta t} \hat{q}(x - \bar{u}t). \quad (3)$$

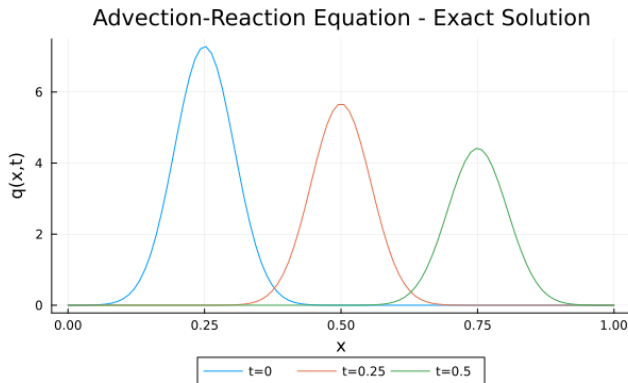


Figure: Evolution of the exact solution of the advection-reaction equation with $\bar{u} = 1$, $\beta = 1$, and $\hat{q} = \text{Gaussian}(0.25, 0.003)$.

For this specific example we can easily compute an **unsplit method**

$$\begin{aligned}q_t &= -\bar{u}q_x - \beta q \\ \frac{Q_i^{n+1} - Q_i^n}{\Delta t} &= -\bar{u} \frac{Q_i^n - Q_{i-1}^n}{\Delta x} - \beta Q_i^n \\ Q_i^{n+1} &= Q_i^n - \bar{u} \frac{\Delta t}{\Delta x} (Q_i^n - Q_{i-1}^n) - \Delta t \beta Q_i^n\end{aligned}$$

which is first-order accurate and stable for $0 < \bar{u} \frac{\Delta t}{\Delta x} \leq 1$.

Note

The full Taylor expansion of (2) can be written formally as

$$\begin{aligned} e^{-\Delta t(\bar{u}\partial_x + \beta)} q(x, t) &:= q(x, t + \Delta t) = \\ &= \sum_{j=0}^{\infty} \frac{(\Delta t)^j}{j!} \partial_t^j q(x, t) = \sum_{j=0}^{\infty} \frac{(\Delta t)^j}{j!} (-\bar{u}\partial_x - \beta)^j q(x, t). \end{aligned} \quad (4)$$

The operator $e^{-\Delta t(\bar{u}\partial_x + \beta)}$ is called **solution operator** for the equation (2) over a time step of length Δt .

In the case of the advection equation, we can split it into two subproblems:

$$\text{Problem A: } q_t + \bar{u}q_x = 0, \quad (5)$$

$$\text{Problem B: } q_t = -\beta q. \quad (6)$$

The idea is to apply the two methods in an alternating manner, using standard solving strategies, e.g.:

$$\text{A-step: } Q_i^* = Q_i^n - \frac{\bar{u}\Delta t}{\Delta x}(Q_i^n - Q_{i-1}^n), \quad (7)$$

$$\text{B-step: } Q_i^{n+1} = Q_i^* - \beta\Delta t Q_i^*. \quad (8)$$

One may think that given that both Q_i^* and Q_i^{n+1} are calculated using Δt , the solution is valid for time $2\Delta t$, but it is not really the case: in fact if we combine the two stages and eliminate Q_i^* , we obtain

$$Q_i^{n+1} = Q_i^n - \frac{\bar{u}\Delta t}{\Delta x}(Q_i^n - Q_{i-1}^n) - \beta\Delta t Q_i^n + \frac{\bar{u}\beta\Delta t^2}{\Delta x}(Q_i^n - Q_{i-1}^n),$$

which differs from the unsplit method for the last term:

$$Q_i^{n+1} = Q_i^n - \bar{u}\frac{\Delta t}{\Delta x}(Q_i^n - Q_{i-1}^n) - \Delta t\beta Q_i^n$$

Commuting vs Non-commuting operators

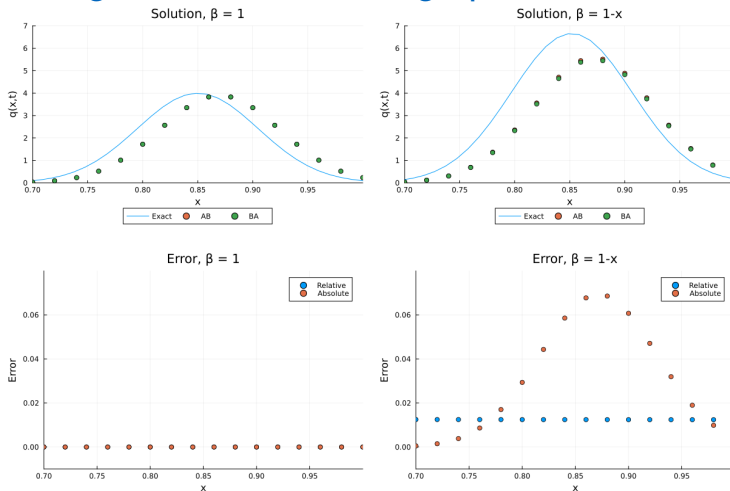


Figure: Comparison between the exact solution of the advection-reaction equation and the split method with the two different orders of steps. The problem has $\bar{u} = 1$, $\hat{q} = \text{Gaussian}(0.25, 0.003)$, $\Delta x = \Delta t = 0.02$, $t = 0.6$.

Consider the more general formulation

$$q_t = (\mathcal{A} + \mathcal{B})q \quad (9)$$

where \mathcal{A} and \mathcal{B} can be differential operators. We assume that they don't explicitly depend on t , so that we can write

$$q_{tt} = (\mathcal{A} + \mathcal{B})q_t = (\mathcal{A} + \mathcal{B})^2 q \quad (10)$$

If we Taylor expand the solution at time t and use the notation defined in (4), we easily get to

$$q(x, \Delta t) = \sum_{j=0}^{\infty} \frac{\Delta t^j}{j!} (\mathcal{A} + \mathcal{B})^j q(x, 0) = e^{\Delta t(\mathcal{A} + \mathcal{B})} q(x, 0). \quad (11)$$

With **Godunov splitting**, we obtain

$$q^*(x, \Delta t) = e^{\Delta t \mathcal{A}} q(x, 0) \quad (12)$$

and

$$q^{**}(x, \Delta t) = e^{\Delta t \mathcal{B}} q^*(x, \Delta t) = e^{\Delta t \mathcal{B}} e^{\Delta t \mathcal{A}} q(x, 0). \quad (13)$$

The splitting error is then

$$q(x, \Delta t) - q^{**}(x, \Delta t) = (e^{\Delta t(\mathcal{A}+\mathcal{B})} - e^{\Delta t \mathcal{B}} e^{\Delta t \mathcal{A}}) q(x, 0). \quad (14)$$

If we Taylor expand qx^{**} , we obtain

$$q^{**}(x, \Delta t) = (I + \Delta t(\mathcal{A} + \mathcal{B}) + \frac{1}{2} \Delta t^2 (\mathcal{A}^2 + 2\mathcal{B}\mathcal{A} + \mathcal{B}^2) + \dots) q(x, 0). \quad (15)$$

With the **Strang splitting** we are approximating $e^{\Delta t(\mathcal{A}+\mathcal{B})}$ by $e^{\frac{1}{2}\Delta t\mathcal{A}}e^{\Delta t\mathcal{B}}e^{\frac{1}{2}\Delta t\mathcal{A}}$. The Taylor expansion shows in fact that

$$e^{\frac{1}{2}\Delta t\mathcal{A}}e^{\Delta t\mathcal{B}}e^{\frac{1}{2}\Delta t\mathcal{A}} = I + \Delta t(\mathcal{A} + \mathcal{B}) + \frac{1}{2}\Delta t^2(\mathcal{A}^2 + \mathcal{A}\mathcal{B} + \mathcal{B}\mathcal{A} + \mathcal{B}^2) + \mathcal{O}(\Delta t^3). \quad (16)$$

After n time steps we obtain

$$Q^n = \underbrace{(e^{\frac{1}{2}\Delta t\mathcal{A}}e^{\Delta t\mathcal{B}}e^{\frac{1}{2}\Delta t\mathcal{A}})(e^{\frac{1}{2}\Delta t\mathcal{A}}e^{\Delta t\mathcal{B}}e^{\frac{1}{2}\Delta t\mathcal{A}})\dots(e^{\frac{1}{2}\Delta t\mathcal{A}}e^{\Delta t\mathcal{B}}e^{\frac{1}{2}\Delta t\mathcal{A}})}_{n \text{ times}} Q^0. \quad (17)$$

Another way of obtaining the same result is

$$\begin{aligned} Q^1 &= e^{\Delta t \mathcal{A}} e^{\Delta t \mathcal{B}} Q^0 \\ Q^2 &= e^{\Delta t \mathcal{B}} e^{\Delta t \mathcal{A}} Q^1 \\ &\vdots \end{aligned} \tag{18}$$

It is essentially the same, but with Δt instead of $\frac{1}{2}\Delta t$.

- **Pros:** computationally cheaper (fewer function evaluations)
- **Cons:** difficult to implement with variable time steps, needs an even number of iterations

If the ODE $q_t = \psi(q)$ is **stiff** (i.e. such that an extremely small time step is required to solve it with an explicit numerical method), then an **implicit** method is needed. The usual choice is the **trapezoidal rule**:

$$Q_i^{n+1} = Q_i^* + \frac{\Delta t}{2} [\psi(Q_i^*) + \psi(Q_i^{n+1})]. \quad (19)$$

Another nice property of the split methods is that they only require the ODE part to be solved implicitly: the hyperbolic part can still be solved with explicit methods.

If the source term is concentrated in space, it is called **stiff** in analogy with the ODEs.

We can say that the solution is evolving on a **slow manifold** in state space and perturbing the solution causes it to produce a rapid transient response followed again by a slow evolution. A very simple example is

$$u'(t) = -\frac{u(t)}{\tau} \quad (20)$$

where of course $u \equiv 0$ is the slow manifold. For some problems the only solution is using an adaptive mesh refinement whereas in other cases using implicit methods may be enough thanks to their good stability properties.

The trapezoidal rule seems to work fine for ODEs, but fails for hyperbolic equations with stiff source terms. If we consider (20), the trapezoidal method yields

$$U^{n+1} = \left(\frac{1 - \frac{1}{2} \frac{\Delta t}{\tau}}{1 + \frac{1}{2} \frac{\Delta t}{\tau}} \right) U^*. \quad (21)$$

If we start on the slow manifold, in this case $U^* = 0$, then $U^{n+1} = 0$ and we remain in the slow manifold. In any other case, if $-\frac{\Delta t}{\tau} \rightarrow -\infty$, then $U^{n+1} = -U^*$ and this means that the coefficient in (21) makes the solution oscillate in time rather than decay as we would expect.

This is due to the fact that the trapezoidal rule is an **A-stable** method: the stability region is the left half plane, but the problem is that it passes through the point at infinity on the Riemann sphere. That's what causes the oscillatory behaviour. We need an **L-stable** method, where the point at infinity is inside the stability region and so the coefficient approaches a value less than 1 in magnitude.

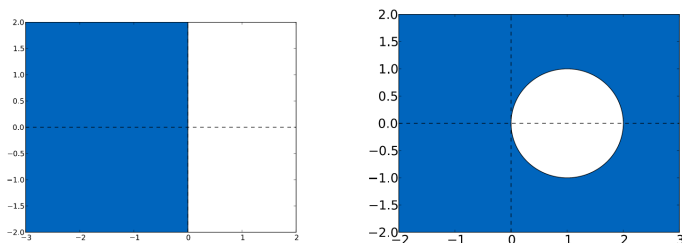


Figure: Stability region of the trapezoidal rule (left) and implicit Euler (right)

The **BDF** (Backward Differentiation Formulas) methods have this characteristic and the simplest one is the backward Euler method. In fact for (20)

$$U^{n+1} = U^* - \frac{\Delta t}{\tau} U^{n+1} \Rightarrow U^{n+1} = \left(\frac{1}{1 + \frac{\Delta t}{\tau}} \right) U^*. \quad (22)$$

This time the coefficient approaches zero for $\frac{\Delta t}{\tau} \rightarrow \infty$. However, the implicit Euler method is only first-order accurate.

- [1] R. J. LeVeque, **Finite Volume Methods for Hyperbolic Problems**. Cambridge: Cambridge University Press, 2002.
- [2] <https://github.com/matilde-t/SeminarCourse-FundamentalsOfWaveSimulation>
- [3] <https://github.com/clawpack/apps/tree/master/fvmbook/chap17>
- [4] https://github.com/clawpack/riemann_book