

From last time

$$\int_V \left[ \hat{T} (c \hat{T} + \nabla \cdot \hat{q} - \hat{Q}) + \hat{q} (k^{-1} \nabla \hat{q} + \nabla T + k \hat{q}) \right] dV$$

$$\int_{d\mathcal{S}} \left[ \hat{T} (c \hat{T} - c T) n_x + \hat{T} (\hat{q}_n - q_n) n_x + (\hat{q} k^{-1} \nabla (\hat{q} - q) n_x + \hat{q} (\hat{T} - T) n_x) \right] d\mathcal{S}$$

weak statement

(WKH)

$$\int_V \left( -\hat{T} c T - \nabla \cdot \hat{T} \hat{q} - \hat{T} \hat{Q} \right) + \underbrace{\varepsilon \hat{q} (k^{-1} \nabla \hat{q} + \nabla T + k \hat{q})}_{\text{const eqn}} dV$$

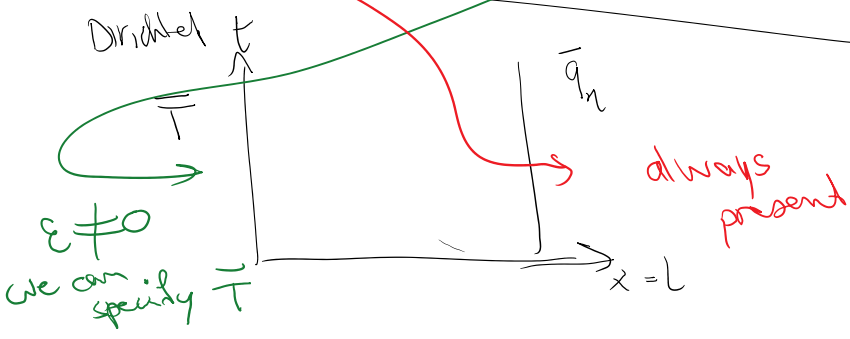
$$\int_{d\mathcal{S}} \left\{ \hat{T} c T n_x + \hat{T} \hat{q}_n n_x + \varepsilon \hat{q} k^{-1} \nabla (\hat{q} - q) n_x + \varepsilon \hat{q} (\hat{T} - T) n_x \right\} d\mathcal{S} = 0$$

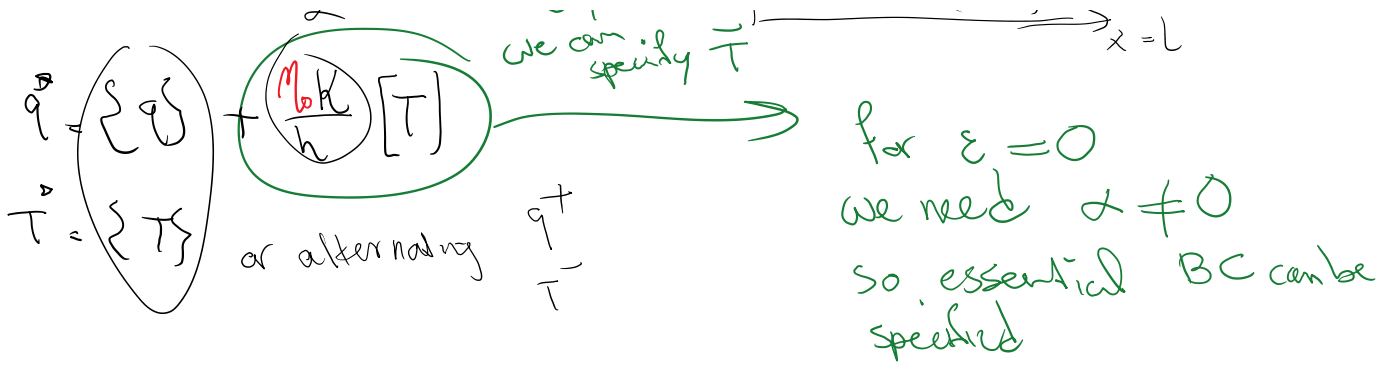
parabolic  $\frac{2F}{1F}$  like above  $q = -k \nabla T$ , constitutive eqn goes away

parabolic 1F

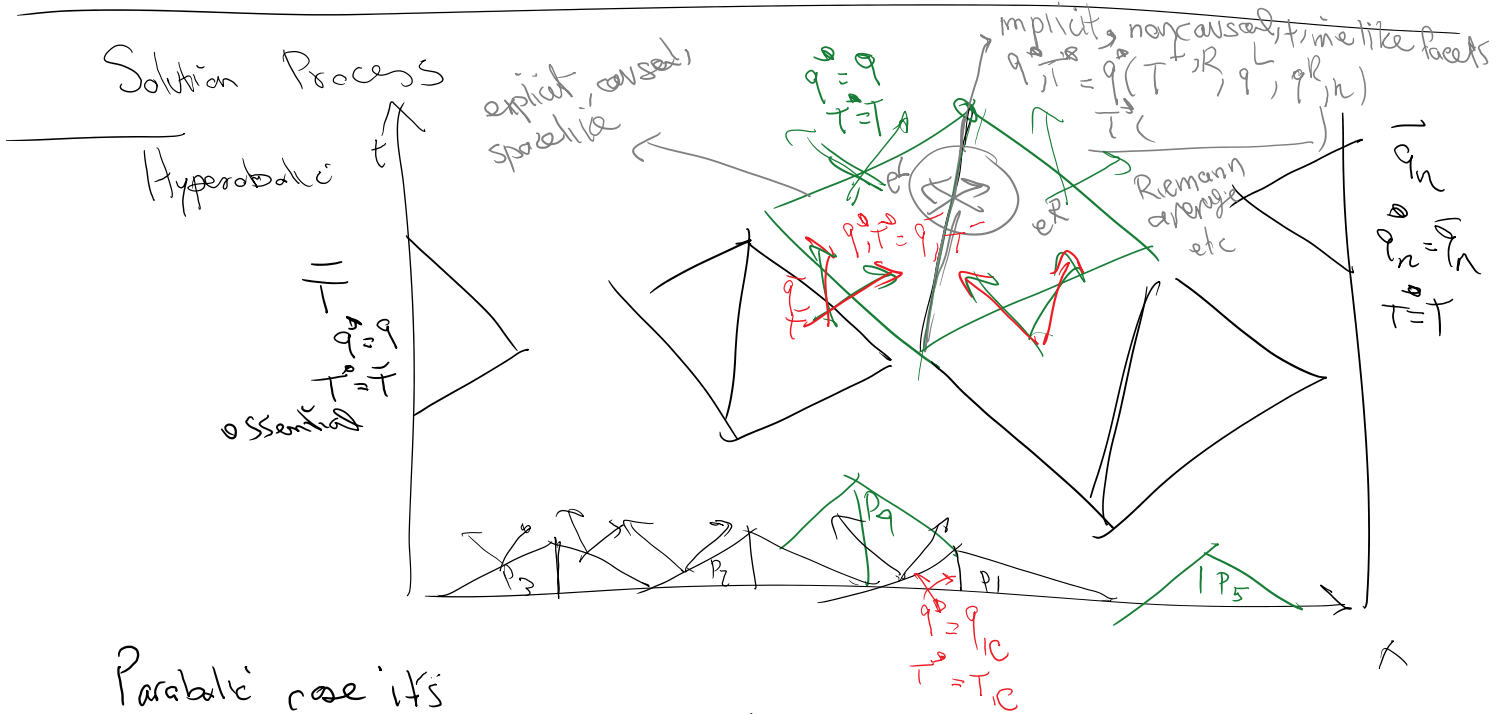
$$\int_V (-\hat{T} c T - \nabla \cdot \hat{T} \hat{q} - \hat{T} \hat{Q}) dV + \int_{d\mathcal{S}} \left\{ \hat{T} c T n_x + \hat{T} \hat{q}_n n_x + \varepsilon \hat{q} k^{-1} \nabla (\hat{q} - q) n_x + \varepsilon \hat{q} (\hat{T} - T) n_x \right\} d\mathcal{S} = 0$$

For 1F parabolic case:  
 Natural Boundary condition (red term) can always be enforced  
 Essential BC:  
 - We need  $\varepsilon \neq 0$   
 - Or for  $\varepsilon = 0$  Refer to fluxes for parabolic problem

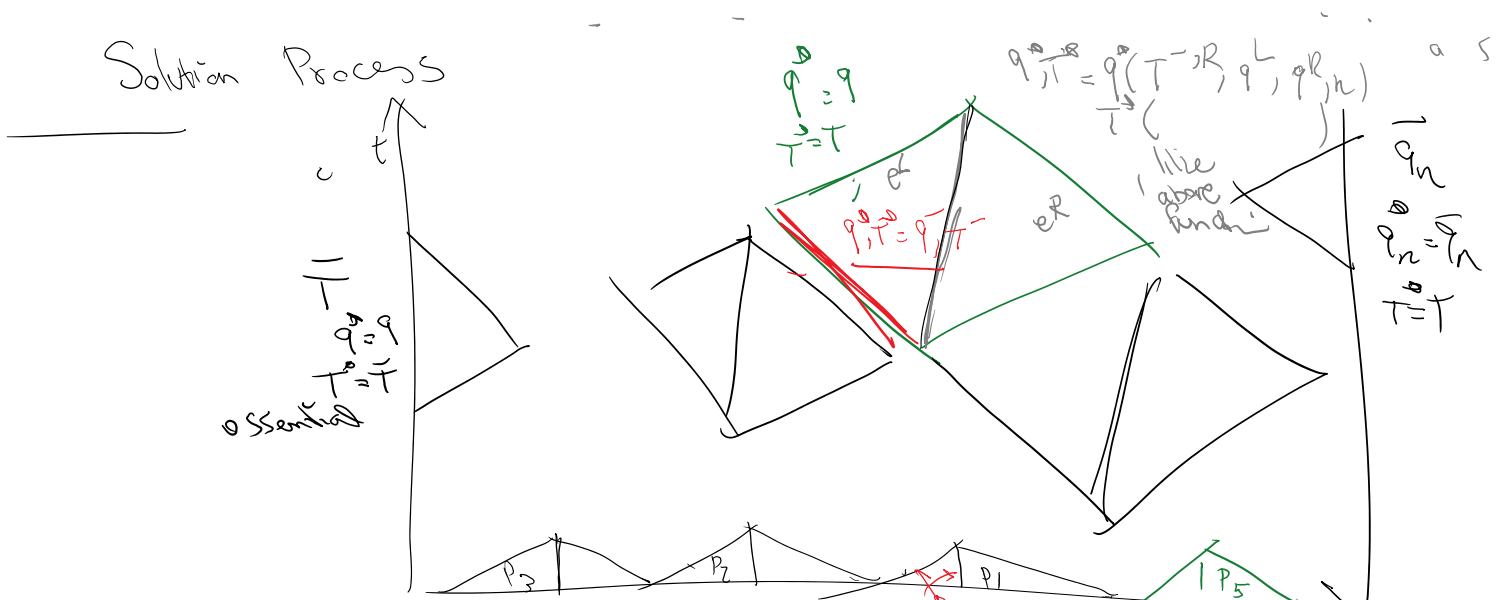


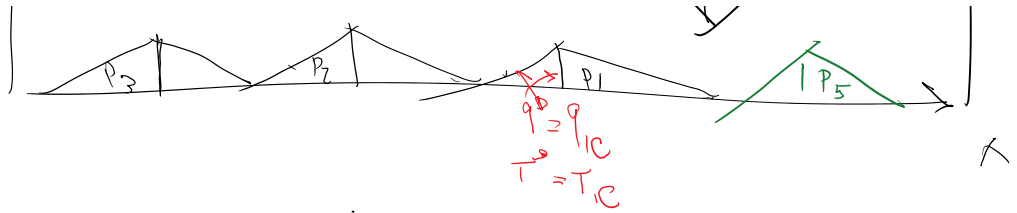


- Unlike elliptic case, the alpha term is not needed for stability. We just need this for enforcing essential BC when  $\varepsilon = 0$



Parabolic case it's similar but there is no characteristics





- For any nonhorizontal facet, the star value really depends on the two sides values.
- For inflow (red) we can use this because the inflow solution is available and the time outflow side is being solved. -> **We make the approximation to let the star equal to the earlier solution.**
- For **outflow** (green) we cannot have a local solution process if the star value is similar to interior facets (i.e. depending on the two sides). We **make an approximation by letting the star equal to the interior trace.**

The parabolic case is stable if the time advance is limited by element size<sup>2</sup>.

Similar background:

**An explicit discontinuous Galerkin scheme with local time-stepping for general unsteady diffusion equations**

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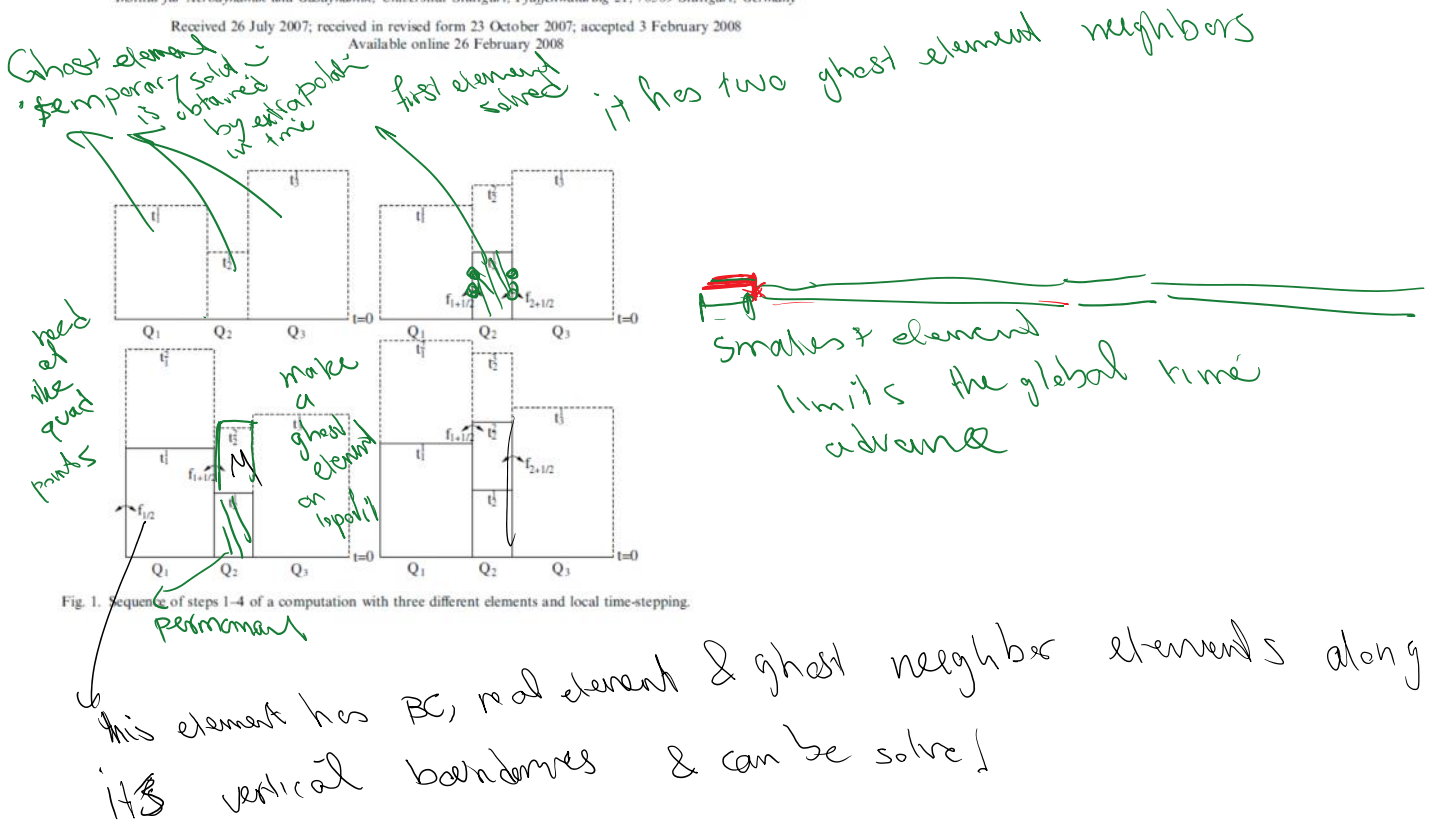


Fig. 1. Sequence of steps 1-4 of a computation with three different elements and local time-stepping.

ADER DG is a very powerful method: Like cSDG it's

- Arbitrarily high order in space and time
- Asynchronous

But here are some concerns:

- For nonlinear PDEs (like fluids the extrapolation in time is very complicated)
- For both hyperbolic and parabolic PDEs the time step is drastically influenced by polynomial order. In SDG, the time advance is not influenced by polynomial order in hyperbolic case.

$$\Delta t \leq \beta(N) \frac{\Delta x^2}{(2N+1)^2 \mu \sqrt{d}}$$
 (3.42)

*correct factor*  
*element size*  
*d: space dimension*  
*space and time*  
*ignore  $\beta$*   
*polynomial order*  
 $\tau \sim \frac{\mu \Delta T}{L^2}$   
 error  $\propto (\Delta t)^{N+1}$  or  $(\Delta t)^N$   
 $N=1 \quad \frac{1}{12} \beta$   
 $N=5 \quad \frac{1}{112} \beta \quad \sim 100\times \text{smaller time step}$

Table 2  
Stability numbers of the optimized STE-DG scheme

N	1	2	3	4	5	6	7
$\beta_{\max}$	1.46	0.8	0.54	0.355	0.28	0.21	0.16
$\beta_{\min}$	1.0	0.33	0.20	0.14	0.10	0.08	0.06
$\beta_{\max}$	2.8	0.86	0.40	0.24	0.16	0.12	0.09

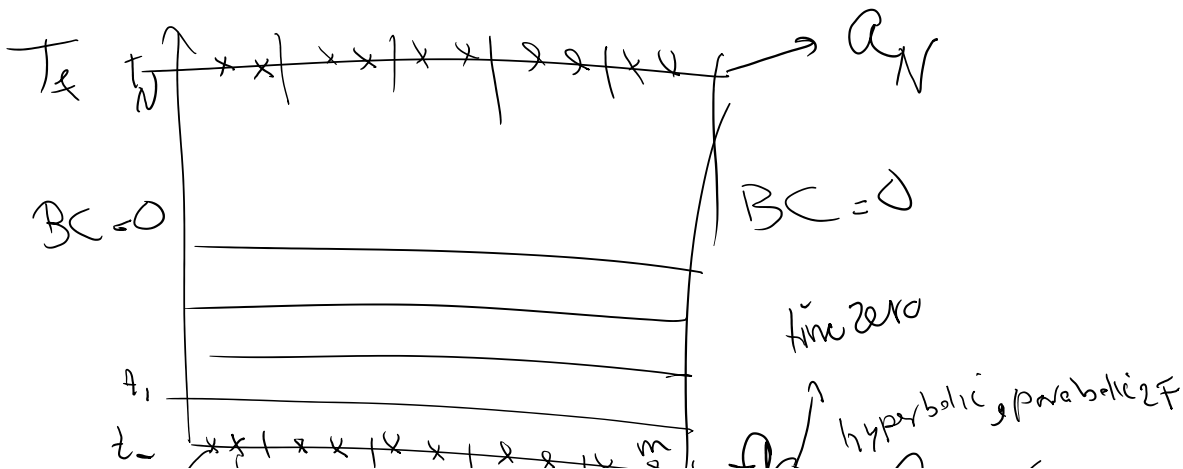
### Global stability analysis of linear PDEs

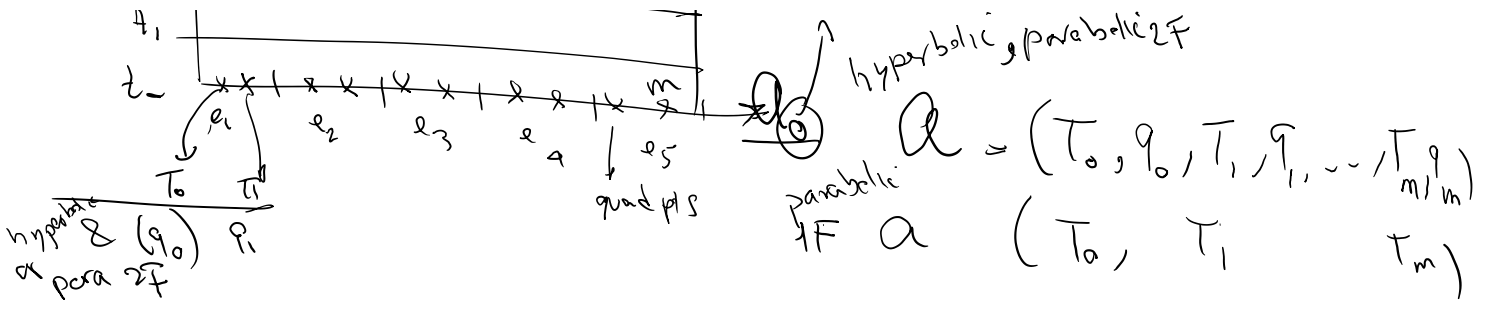
To analyze the stability of the scheme, a periodic problem with a given spatial discretization is considered. As the problem (3.40) is linear, one can construct a matrix  $W$  such that

$$\hat{u}^{\text{new}} = W \hat{u}^{\text{old}}, \quad (3.41)$$

where  $\hat{u}^{\text{old}}$  denote DOF at a common time level  $t^{\text{old}}$  and  $\hat{u}^{\text{new}}$  denote DOF at a common time level  $t^{\text{new}}$  with  $t^{\text{new}} > t^{\text{old}}$ .  $W = W(\Delta t_i)$  depends on the time steps  $\Delta t_i$  of each element  $Q_i$ . Following the matrix method of stability analysis described in [14], the scheme is stable, if the spectral radius  $\rho(W)$  is lower or equal to 1.

We first consider uniform grid spacing and uniform polynomial order, thus,  $\Delta x = \text{const}$  and  $\Delta t = \text{const}$  in the whole computational domain. For an explicit DG scheme discretizing equation (3.40), a stability restriction for the time step has the form

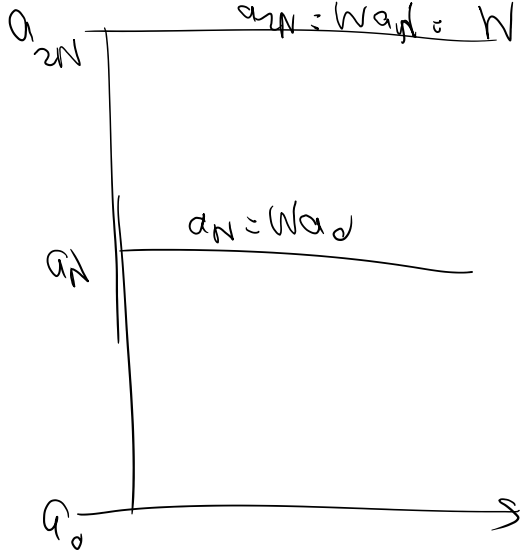




$$a_{qN} = f(a_0)$$

$$a_N = W a_0$$

$$a_{qN} = W a_{q0} = W^2 a_0$$



$$a_{qN} = W^q a_0$$

Eigen decomposition (if it exists)

$$W U = U \Lambda$$

$$U = [u_1 | u_2 | \dots | u_m]$$

right eigenvectors

$$\Lambda = \begin{bmatrix} \lambda_1 & & \\ & \lambda_2 & \\ & & \ddots \\ & & & \lambda_m \end{bmatrix}$$

$$W = U \Lambda U^{-1}$$

$$a_{qN} = W^q a_0$$

$$a_{qN} = U \Lambda^q U^{-1} a_0$$

for all  $\lambda_i$  :  $|\lambda_i| < 1$

$$\|a_{qN}\| < \|a_0\|$$

for all  $\lambda_i$  :  $\underline{| \lambda_i | < 1}$   $\|a_{qn}\| \leq \|a_n\|$

In general  $W$  may not be diagonalizable

→ We can always reduce it to Jordan form

$$M_{n \times n} = \begin{bmatrix} \lambda_1 & & & \\ & \lambda_2 & & \\ & & \ddots & \\ & & & \lambda_n \end{bmatrix}$$

$\rho(W)$   
 spectral radius =  $\text{Min}_{i=1 \dots n} (|\lambda_i|)$

For stability

$\rho(W) < 1$	in general
$\rho(W) \leq 1$	$W$ is diagonalizable

$\rho(W) = 1$  & nondiagonalizable → weak unstable

(readily blows up)

$\rho(W) > 1$  → strong unstable / exponentially blows up