

Hyperbolic Equations

Hyperbolic Equations

- classic 2nd order Form elliptic in space
2nd derivative in time

$$\frac{\partial^2 u}{\partial t^2} - \nabla \cdot C \nabla u = f(u, \frac{\partial u}{\partial t}, \frac{\partial u}{\partial x}, \dots)$$

- Prototype Equation to study:

$$\frac{\partial^2 u}{\partial t^2} + \tilde{\gamma} \frac{\partial u}{\partial t} - c^2 \frac{\partial^2 u}{\partial x^2} = 0 \quad \text{"Telegraph Egn"}$$

c - "wave speed" ; $\tilde{\gamma}$ - dissipation or "loss factor"

$\tilde{\gamma} = 0 \Rightarrow$ "Wave Equation"

Hyperbolic Equations

- Recall BCs + ICs needed:

$$\begin{array}{ccc} u \text{ or } \frac{\partial u}{\partial x} \text{ or } a \frac{\partial u}{\partial x} + bu & \begin{array}{c} \text{Diagram of a rectangular domain in the } (x, \tau) \text{ plane. The horizontal axis is } x \text{ and the vertical axis is } \tau. \text{ A wavy line is drawn across the top of the rectangle, representing a boundary condition.} \end{array} & u \text{ or } \frac{\partial u}{\partial x} \text{ or } a \frac{\partial u}{\partial x} + bu \\ & u \text{ and } \frac{\partial u}{\partial \tau} & \end{array}$$

Hyperbolic Equations

- can always decompose into coupled set of 1st order PDEs

$$(1) \quad \frac{\partial u}{\partial t} = c_1 \frac{\partial v}{\partial x}$$

$$(2) \quad \frac{\partial v}{\partial t} + \tau v = c_2 \frac{\partial u}{\partial x}$$

$$\Rightarrow \underbrace{\frac{\partial^2 v}{\partial t^2} + \tau \frac{\partial v}{\partial t} - c_2^2 \frac{\partial^2 v}{\partial x^2}}_{\substack{\text{Differentiate (1)} \\ \text{in time and substitute (1)}}} = 0$$

Or : Differentiate (1) in time and substitute (2)

$$\frac{\partial^2 u}{\partial t^2} = c_1 \frac{\partial}{\partial x} \left(\frac{\partial v}{\partial t} \right) = -c_1 \underbrace{\tau \frac{\partial v}{\partial x}}_{\frac{1}{c_1} \frac{\partial u}{\partial t}} + c_2^2 \frac{\partial^2 u}{\partial x^2}$$

$$\text{So: } \frac{\partial^2 u}{\partial t^2} + \tau \frac{\partial u}{\partial t} - c_2^2 \frac{\partial^2 u}{\partial x^2} = 0$$

Hyperbolic Equations

- BCs + ICs for coupled system...



Hyperbolic Equations

Time Stepping Strategies

- consider 2nd order system
- obvious approach ... replace $\frac{\partial^2 u}{\partial t^2}$ w/
centered 2nd order approximation ...
requires 3 levels in time!

a) Explicit:
$$\frac{\partial_t^2 u_i^l}{\Delta t^2} + \frac{\tau}{2\Delta t} \frac{\partial_t u_i^l}{\Delta t} - c^2 \frac{\partial_x^2 u_i^l}{h^2} = 0$$

$$\frac{u_i^{l+1} - 2u_i^l + u_i^{l-1}}{\Delta t^2} + \frac{\tau}{2\Delta t} (u_i^{l+1} - u_i^{l-1}) - \frac{c^2}{h^2} (u_{i-1}^l - 2u_i^l + u_{i+1}^l) = 0$$

Hyperbolic Equations

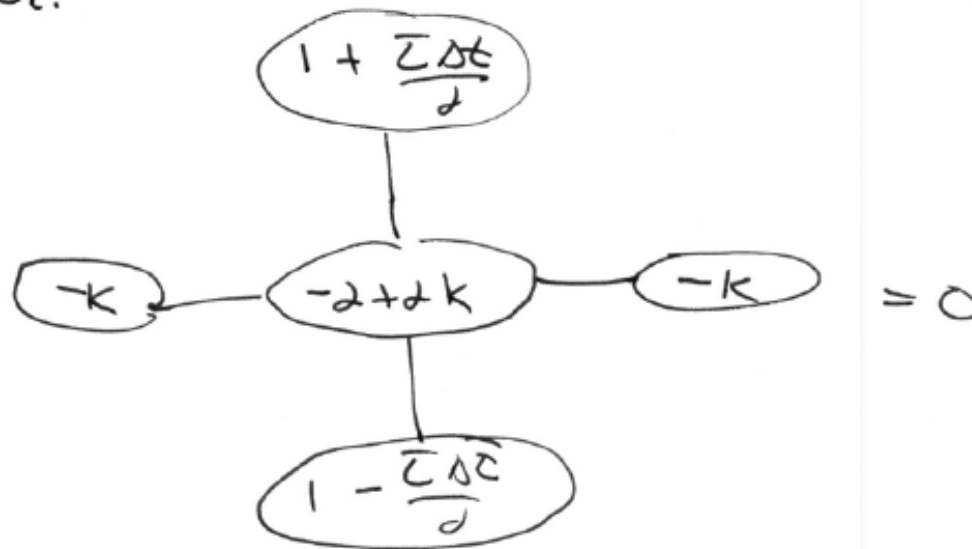
multiply through by Δt^2 :

$$u_i^{l+1} - 2u_i^l + u_i^{l-1} + \frac{2\Delta t}{\Delta x} (u_i^{l+1} - u_i^{l-1}) - \frac{c^2 \Delta t^2}{h^2} (u_{i-1}^l - 2u_i^l + u_{i+1}^l)$$

$$K = \frac{c^2 \Delta t^2}{h^2}$$

↪ "Courant #"

MOLECULE:

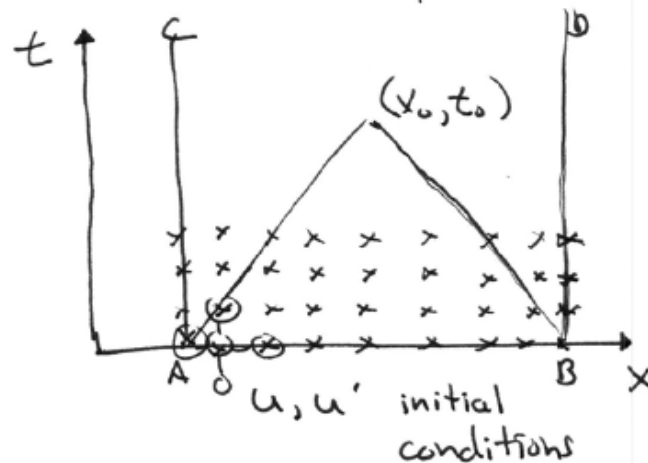


Hyperbolic Equations

- Features

- centered in $x + \epsilon \Rightarrow O(h^2 + \Delta t^2)$
- No matrices
- Pointwise propagation ... experience w/ Parabolic suggests stability constraint

- Recall characteristic Family for wave Equation



Hyperbolic Equations

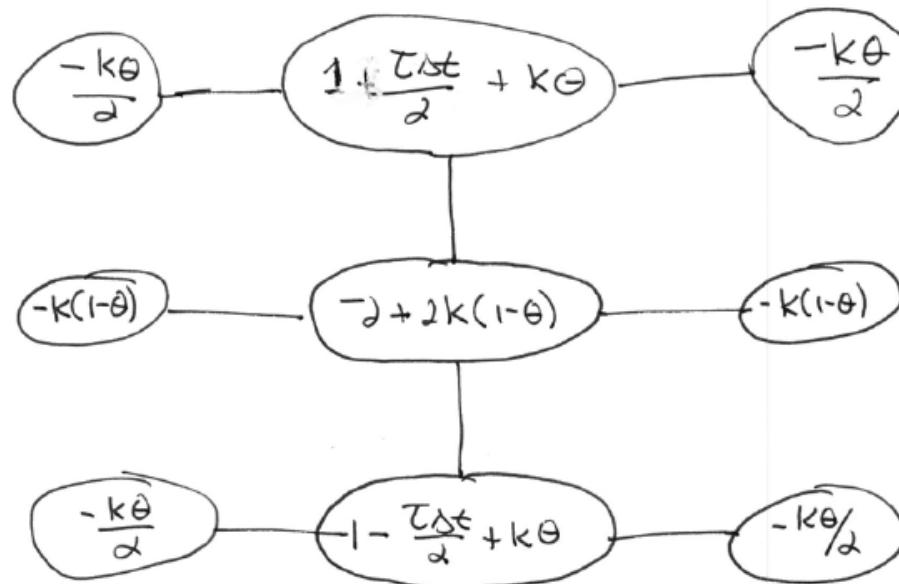
- Knowing initial conditions, can get to (x_0, t_0) without influence of AC or BD....
suspect stability limits similar to previous work!
- Turns out: don't get convergence as $h, \Delta t \rightarrow 0$
for $\frac{\Delta t}{h} > 1/c$

Hyperbolic Equations

2nd order Egn (cont)

b) Implicit

$$(u_i^{l+1} - 2u_i^l + u_i^{l-1}) + \frac{\tau \Delta t}{\alpha} (u_i^{l+1} - u_i^{l-1}) - k \delta_x^2 \left[\theta \left(\frac{u_i^{l+1} + u_i^{l-1}}{2} \right) + (1-\theta) u_i^l \right] = 0$$



Hyperbolic Equations

- solution of system of Equations per-time step
- centered in $(x, t) \Rightarrow O(h^2 + \Delta t^2)$
- stability: $\Theta \geq 1/2$ unconditional

$$k < 1/(1-2\Theta) \quad \text{otherwise}$$

- Shortest waves propagate + damp (phase distortions possible)

Hyperbolic Equations

- Courant, Friedrichs and Lewy Condition

↳ convergence requires numerical characteristics (i.e. domain of dependence of FD) to have slopes less than $1/c$ (i.e. must include domain of dependence of PDE)

$$\therefore \frac{\Delta t}{h} < 1/c \Rightarrow \underbrace{\frac{c \Delta t}{h}} < 1$$

This is essentially
a stability requirement

Hyperbolic Equations

- Lets look at formal stability analysis

$$u_i^{l+1} = \gamma_0 u_i^l ; u_{i+1}^l = u_i^l e^{i\alpha h}$$

$$u_i^{l+1} - 2u_i^l + u_i^{l-1} + \frac{\tau \Delta t}{2} (u_i^{l+1} - u_i^l) + K (u_{i-1}^l - 2u_i^l + u_{i+1}^l) = 0$$

$$\gamma_0 - 2 + \frac{1}{\gamma_0} + \frac{\tau \Delta t}{2} (\gamma_0 - \frac{1}{\gamma_0}) - K (2 \cos \alpha h - 2) = 0$$

$$\gamma_0' \left(1 + \frac{\tau \Delta t}{2} \right) + \gamma_0 (2K(1 - \cos \alpha h) - 2) + \left(1 - \frac{\tau \Delta t}{2} \right) = 0$$

$$\text{Need } |\gamma_0| \leq 1; \text{ requires } \frac{c}{a} \leq 1; |b| \leq a+c$$

Hyperbolic Equations

$$\rightarrow \frac{c}{a} \leq 1 \Rightarrow \frac{1 - \frac{2\sigma t}{j}}{1 + \frac{2\sigma t}{j}} \leq 1 \quad \underline{\text{always!}}$$

$$\rightarrow |b| \leq a + c$$

$$|2k(1 - \cosh) - 2| \leq 2 \Rightarrow |k(1 - \cosh) - 1| \leq 1$$

$$\text{i.e.} \quad -1 \leq k(1 - \cosh) - 1 \leq 1$$

$$\text{consider left side: } -1 \leq k(1 - \cosh) - 1$$

always true since $k > 0$

$$1 - \cosh > 0$$

Hyperbolic Equations

right side: $K(1 - \cos \phi h) \leq 1$

$$K \leq \frac{1}{1 - \cos \phi h}$$

most restrictive limit is stability restriction!
as usual ... short waves are potentially most
unstable i.e. $\phi h = \pi$

CONCLUDE:

$$K \leq 1 ; \frac{C^2 \Delta t^2}{h^2} \leq 1 \text{ OR } \frac{C \Delta t}{h} \leq 1$$

"Courant" condition

Hyperbolic Equations

- So what kind of errors do we expect w/ explicit scheme...

- Experience tells us "short waves" are the most difficult to handle ... look at θ_0 ($\theta h = \pi$)

Recall:
$$\gamma_0^2 \left(1 + \frac{\tau \Delta t^2}{2}\right) + \gamma_0 (2k(1 - \cos \theta h) - 2) + \left(1 - \frac{\tau \Delta t^2}{2}\right) = 0$$

at $\theta h = \pi$

$$\gamma_0^2 \left(1 + \frac{\tau \Delta t^2}{2}\right) + \gamma_0 (4k - 2) + \left(1 - \frac{\tau \Delta t^2}{2}\right) = 0$$

Hyperbolic Equations

$$\gamma_0 = \frac{2 - 4K \pm \sqrt{(4K-2)^2 - 4\left(1 - \left(\frac{\Delta t \tau}{2}\right)^2\right)}}{2(1 + \tau \Delta t/2)}$$

$$= \frac{2 - 4K \pm j \left(4\left(1 - \left(\frac{\Delta t \tau}{2}\right)^2\right) - (4K-2)^2 \right)^{1/2}}{2(1 + \tau \Delta t/2)}$$

generally get decay (Good News), but also phase distortion!!
(BAD NEWS)

Hyperbolic Equations

e.g. $K=1/4$:
$$r_0 = \frac{1 \pm j \sqrt{4(1 - (\frac{\Delta t}{2})^2) - 1}}{2(1 + Z\Delta t/2)}$$

let $Z = .015$
 $h = 5$
 $\Delta t = .25$ }
$$r_0 = \frac{1 \pm j \sqrt{4(1 - (1.875e-3)^2) - 1}}{2(1 + 1.875e-3)}$$

$$= \frac{1 \pm 1.732j}{2.00375} \Rightarrow |r_0| < 1$$
$$\text{Arg}(r_0) \approx \pm \pi/3$$

Amplitude damped ; expect phase errors

FD Conservation

FD Conservation

- Conservation Law + Constitutive Relation \Rightarrow PDE

$$\nabla \cdot \mathbf{q} = \sigma$$

Flux \nearrow source/volume

$$\mathbf{q} = -k \nabla u \Rightarrow \nabla \cdot (-k \nabla u) = \sigma$$

\nwarrow scalar surrogate for q

e.g.	<u>q</u>	<u>Conserved</u>	<u>u</u>	<u>σ</u>
	Heat Flux	Thermal Energy	T	Heating rate
	Diffusion Flux	Molecules of species	C	Reaction rate
	Mass Flux	Fluid mass	ρ	Evaporation rate
	\vdots	\vdots	\vdots	\vdots

FD Conservation

Note: Θ need not be constant

- radioactive decay: $\Theta = -kC$

storage: $\Theta = -\frac{\partial C}{\partial t}$

$$\nabla \cdot (-D \nabla c) = \underbrace{-\frac{\partial C}{\partial t} - kC}_{\sigma}$$

(Paraboliz
-more later)

FD Conservation

- PDE is local conservation statement

- Global Conservation:

$$\int (\text{PDE}) dV \Rightarrow \int \nabla \cdot \mathbf{q} dV = \int \phi dV$$

Divergence Theorem: $\oint \mathbf{q} \cdot \hat{\mathbf{n}} ds = \int \phi dV$

Rate of escape = sum of sources

or $\oint -K \nabla u \cdot \hat{\mathbf{n}} ds = \int \phi dV$

$$\oint -K \frac{\partial u}{\partial n} ds = \int \phi dV$$

FD Conservation

• Numerical Conservation $\Rightarrow \int () dV = \sum ()_i \Delta V_i$

e.g. 1-D $\sum_i \int_{i-1/2}^{i+1/2} ()_i dx$



$\frac{\partial}{\partial x} \ll \frac{\partial u}{\partial x} = -\theta$ For 1 box:

FD Conservation

$$\int_{i-1/2}^{i+1/2} \left(\frac{\partial}{\partial x} k \frac{\partial u}{\partial x} \right)_i dx = - \int_{i-1/2}^{i+1/2} \sigma dx$$

$$\underbrace{k \frac{\partial u}{\partial x} \Big|_{i+1/2}}_{-q_{i+1/2}} - \underbrace{k \frac{\partial u}{\partial x} \Big|_{i-1/2}}_{+q_{i-1/2}} = -\sigma_i h$$



so σ_i represents all sources in the box associated w/ Node i

$$\sigma_i = \int_{x_{i-1/2}}^{x_{i+1/2}} \sigma dx / (x_{i+1/2} - x_{i-1/2})$$

FD Conservation

Now:

$$\begin{aligned} -q_{i+1/2} + q_{i-1/2} &= K_{i+1/2} \left(\frac{u_{i+1} - u_i}{h} \right) - K_{i-1/2} \left(\frac{u_i - u_{i-1}}{h} \right) \\ &= -\theta_i h \end{aligned}$$

Or

$$\frac{K_{i+1/2} \left(\frac{u_{i+1} - u_i}{h} \right) - K_{i-1/2} \left(\frac{u_i - u_{i-1}}{h} \right)}{h} = -\theta_i$$

FD Conservation

But left-hand side is exactly our FD approximation at node i ... so to conserve we must view θ_i as average of sources in box

$$\text{i.e. } \theta_i = \int_{x_{i-1/2}}^{x_{i+1/2}} \theta dx / (x_{i+1/2} - x_{i-1/2})$$

Then $\sum_{i=1}^N$ (FD equation $\neq i$):

$$-q_{N+1/2} + q_{1/2} = - \sum_{i=1}^N \theta_i h$$

Internal $q_{i+1/2}$ cancel

FD Conservation

Then $\sum_{i=1}^N$ (FD equation $\neq i$):

$$-q_{N+1/2} + q_{1/2} = - \sum_{i=1}^N \theta_i h \quad \text{Internal } \underline{q_{i+1/2}} \text{ cancel}$$



- Type I BCs at 0, N+1 \Rightarrow can solve ...

- \hookrightarrow Conservation boundaries not at 0, N+1
- \hookrightarrow Conservation independent of $\theta_0 + \theta_{N+1}$

- Type II + Type III ... bring PDE molecule to boundary

FD Conservation

e.g. $-K \frac{\partial u}{\partial x} = f_0$ specified ... then

FD molecule at node 0 :

$$K_{1/2} \left(\frac{u_{-1} - 2u_0 + u_1}{h} \right) = -f_0 h \quad (K_{1/2} = K_{-1/2})$$

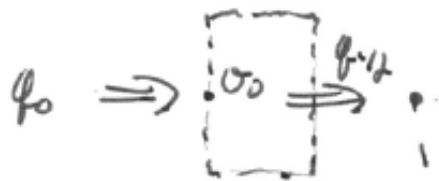
But BC says : $k_0 \frac{u_{-1} - u_1}{2h} = f_0$

$$\Rightarrow u_{-1} = \frac{2hf_0}{k_0} + u_1 = \frac{2hf_0}{K_{1/2}} + u_1 \quad \left(k_0 = \frac{K_{1/2} + K_{3/2}}{2} \right)$$

FD Conservation

so we get $2q_0 + 2\left(\frac{u_1 - u_0}{h}\right)k_{1/2} = -\sigma_0 h$

or $q_0 + \underbrace{k_{1/2}\left(\frac{u_1 - u_0}{h}\right)}_{-q_{1/2}} = -\sigma_0 h/2$



This provides the missing half box!

Similarly at node $N+1$: $-q_{N+1} + q_{N+1/2} = -\sigma h/2$

FD Conservation

So For Global conservation add in the two boundary molecules ...

$$\begin{aligned} -q_{N+1} + q_0 &= -\sum_{i=1}^N \phi_i h - \frac{\phi_0 h}{2} - \frac{\phi_{N+1} h}{2} \\ &= -\sum_{i=0}^N (\phi_i + \phi_{i+1}) h/2 \end{aligned}$$

Trapezoidal Rule Integration

This is the FD conservation statement!

FD Conservation

- At Type I boundaries need to use all molecules to conserve... use the unused (in solution) molecules at boundary to compute fluxes

$$q_0 = -K_{1/2} \left(\frac{u_1 - u_0}{h} \right) - \frac{\sigma_0 h}{2}$$

This is the
intuitive FD
expression...
looks $\mathcal{O}(h)$

correction term to
account for half box
makes everything $\mathcal{O}(h^2)$

similar at node $N+1$:

$$-q_{N+1} = K_{N+1/2} \left(\frac{u_{N+1} - u_N}{h} \right) - \sigma_{N+1} h/2$$

FD Conservation

In 2-D:

$$\frac{1}{h} \left[k_{i+1/2} \left(\frac{u_{i+1,j} - u_{i,j}}{h} \right) - k_{i-1/2} \left(\frac{u_{i,j} - u_{i-1,j}}{h} \right) \right] \\ + \frac{1}{h} \left[k_{j+1/2} \left(\frac{u_{i,j+1} - u_{i,j}}{h} \right) - k_{j-1/2} \left(\frac{u_{i,j} - u_{i,j-1}}{h} \right) \right]$$

$$= -\Theta_{i,j}$$

$$-h q_{i+1/2}^x + h q_{i-1/2}^x - h q_{j+1/2}^y + h q_{j-1/2}^y = -\Theta_{i,j} h^2$$

FD Conservation



$\bar{\sigma}_{ij}$ is 2D average

i.e.

$$\bar{\sigma}_{ij} = \frac{\int_{y_{j-1/2}}^{y_{j+1/2}} \int_{x_{i-1/2}}^{x_{i+1/2}} \sigma dx dy}{(x_{i+1/2} - x_{i-1/2})(y_{j+1/2} - y_{j-1/2})}$$

- Note: if $\sigma = Q \delta(x-x_0, y-y_0)$ then for
 (i, j) corresponding to (x_0, y_0)

$$\bar{\sigma}_{ij} = \frac{\iint Q \delta(x-x_0, y-y_0) dx dy}{h^2} = \underbrace{Q/h^2}_{\text{source strength divided by area of 1 box!!}}$$

Introduction to the Method of Weighted Residuals

Method of Weighted Residuals

- Consider $\nabla^2 \mathbf{u} + \mathbf{f} \mathbf{u} = \mathbf{g} \rightarrow \underbrace{(\nabla^2 + \mathbf{f})}_{\mathbf{L}} \mathbf{u} = \mathbf{g}$

- Finite Difference Approach

- Approximate \mathbf{L} w/ $\mathbf{L}_{ij} \rightarrow \underbrace{(\delta^2 + \mathbf{f})}_{\mathbf{L}_{ij}} \mathbf{u}_{ij} = \mathbf{g}_{ij}$

i.e. replace “differential” operator w/ “difference” operator \rightarrow get “exact solution” to “approximate operator”

- Limitations:

- cumbersome on irregular meshes
- curved boundaries difficult to handle
- \mathbf{u} only found (i,j) points – need interpolation strategy

Method of Weighted Residuals

- Weighted Residuals

- Approximate \mathbf{u} as $\hat{u} = \sum_{j=1}^N c_j \Phi_j(x, y, z)$
 - c_j : coefficients
 - Φ_j : known function
 - “basis” func.
 - “trial” func.
 - “expansion” func.

- Define “Residual” $\rightarrow R(\hat{u}) = L\hat{u} - g$
- For exact solution: $R(\hat{u}) = 0$ everywhere then
- Want $R(\hat{u}) = 0$ to vanish in average way, one way is in weighted integral sense

$$\iiint R(\hat{u}) W(x, y, z) dx dy dz = 0$$

any function
of position

- “ $R(\hat{u})$ orthogonal to all $W(x, y, z)$ ”

Method of Weighted Residuals

- So for $\hat{u} : R(\hat{u}) \neq 0 \dots$ choose 'N' c_j 's such

that $\langle R(\hat{u}), W_i \rangle = 0$ for $i = 1, 2, 3, \dots, N$

“Inner Product” $\Rightarrow \langle \mathbf{a}, \mathbf{b} \rangle \equiv \iiint \mathbf{a} \bullet \mathbf{b} \, dx dy dz$

- W_i 's set of “weighting” functions \rightarrow finite!
local “testing” functions

- Use ‘N’ independent W_i 's \rightarrow generate ‘N’ equations in ‘N’ unknown c_j 's

$$\langle R(\hat{u}), W_i \rangle = \sum_{j=1}^N c_j \langle L(\phi_j), W_i \rangle = \langle g, W_i \rangle \text{ for each } W_i(x, y) \text{ } i = 1, 2, \dots, N$$

- Necessary, but not sufficient for $\hat{u} = u$

Method of Weighted Residuals

- Continuous function must be zero if it is orthogonal to every member of a complete set \therefore WRM can be thought of as a technique which enforces orthogonality between basis and weighting function sets
 - General idea is that the basis is a subset of a complete set that can represent any function, i.e. the true solution
 - As number of coefficients goes up, the approximation approaches the complete set

Method of Weighted Residuals

- Take home interpretation:
 - PDE is determined from first principles
 - Introduce the idea of a weighting function which is a known function of space – on its introduction, the physics behavior is prescribed to a region
 - Integration formulation
 - Introduction of basis as an approximation to the solution – on its introduction, it provides a description of how the solution will be treated behaviorally within in a local region

Method of Weighted Residuals

- Weighted Residual Methods Summary
 - L is typically a differential operator
 - W_i not complete in practice (N finite), but make “ $R(u)$ orthogonal to 1st N members of a complete set”
 - “Approximate solution” which exactly satisfies “differential relations” in PDE
 - Is an “integral” formulation

Method of Weighted Residuals

- What's involved numerically?

- Expand unknown solution as $\hat{u} = \sum_{j=1}^N c_j \phi_j(\mathbf{x}, \mathbf{y}, \mathbf{z})$
 - Finite sum
 - $\phi_j(\mathbf{x}, \mathbf{y}, \mathbf{z})$ known function of (x, y, z)
 - c_j 's unknown coefficients to determine
- Generate system of equations in unknowns
$$R(\hat{u}) = \langle (L(\hat{u}) - g), W_i(\mathbf{x}, \mathbf{y}, \mathbf{z}) \rangle = 0 \text{ for } i = 1, 2, 3, \dots, N$$
$$\sum_{j=1}^N c_j \langle L(\phi_j), W_i(\mathbf{x}, \mathbf{y}, \mathbf{z}) \rangle = \langle g, W_i(\mathbf{x}, \mathbf{y}, \mathbf{z}) \rangle \text{ for } i = 1, 2, 3, \dots, N$$
 - $W_i(x, y, z)$ known functions of (x, y, z)
 - For each i , generate algebraic equation in c_j 's
 - Creates 'N' equations in 'N' unknowns

Method of Weighted Residuals

- But analytic methods...
 - Require special knowledge of how to choose basis and weighting functions
 - Different choices are needed for different problems
 - Usually need an infinite # of them
 - Can't find them for many practical problems
- Numerically want ...
 - Basis and Weighting functions to be simple ... easy to integrate
 - Single choice suitable for many problems
 - Can only use finite #, but want convergence as number used increases

Method of Weighted Residuals

- WRM Function

- Subregion
- Collocation
- Least squares
- Monte Carlo
- Galerkin

- Weighted Residual Method

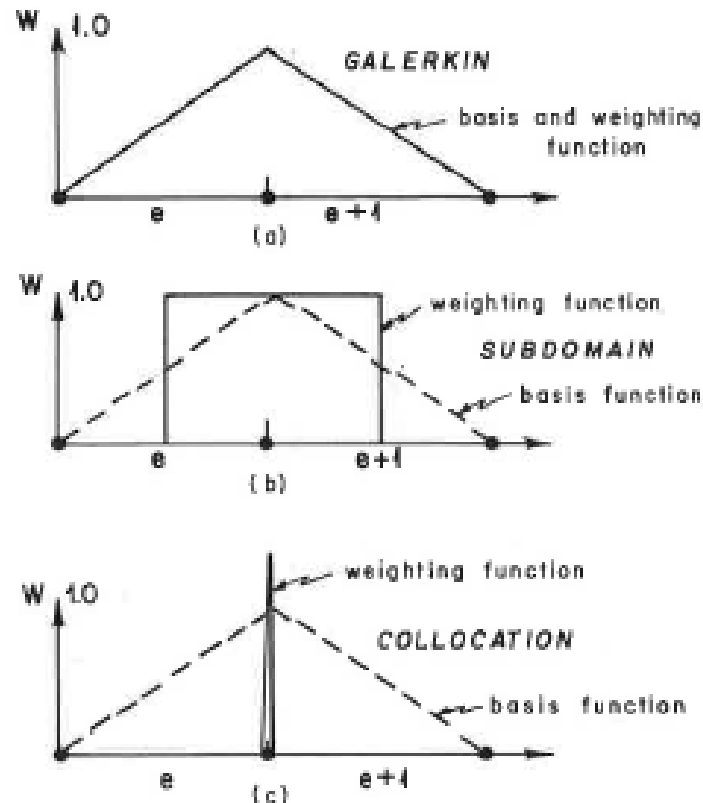


Figure 2.4. Schematic representation of the one-dimensional weighting functions for the Galerkin, subdomain and collocation methods. (It is assumed here that the chapeau function is used as a basis for all methods.)

weighting

N' small
equal to

- residual

weighting

nction
ction are

residual

Galerkin Weighted Residual Method

- Key Feature:
 - Basis function and weighting function are selected to be the same function form
- So what do we choose?
 - Polynomials may be a nice choice
 - Other's exist

Polynomial Basis/Weighting Functions

- Why?
 - Easy to differentiate
 - Represent a complete set for continuous functions ... e.g. Taylor Polynomial
- Lagrange Polynomials
 - Nth order polynomial for N+1 pts
 - Easily automated
 - Handy on uneven grids
 - $\phi_i(\mathbf{x}) = 0$ at $\mathbf{x} = \mathbf{x}_j$ for $i \neq j$
 - $\phi_j(\mathbf{x}) = 1$ at $\mathbf{x} = \mathbf{x}_i$ for $i = j$

$$\phi_i = \prod_{j \neq i} \frac{(\mathbf{x} - \mathbf{x}_j)}{(\mathbf{x}_i - \mathbf{x}_j)}$$

Polynomial Basis/Weighting Functions

- Lagrange Polynomials

- $\phi_j(\mathbf{x}_j) = 1 \Rightarrow \mathbf{u}(\mathbf{x}) = \sum_{j=1}^N \mathbf{c}_j \phi_j(\mathbf{x}, \mathbf{y}) \Rightarrow \mathbf{u}(\mathbf{x}_i) = \mathbf{c}_i$

- $\therefore \mathbf{u}(\mathbf{x}) = \sum_{j=1}^N \mathbf{u}_j \phi_j(\mathbf{x}, \mathbf{y})$

Coefficients are solution at nodes ... similar to FD but have functional form specified in between

- everywhere $\sum_{j=1}^N \phi_j(\mathbf{x}, \mathbf{y}) = 1$

- $\sum_{j=1}^N \frac{\partial \phi_j(\mathbf{x}, \mathbf{y})}{\partial \mathbf{x}} = \frac{\partial}{\partial \mathbf{x}} \sum_{j=1}^N \phi_j(\mathbf{x}, \mathbf{y}) = 0$

Polynomial Basis/Weighting Functions

- Global Polynomials

- Potential for disaster ... “polynomial wiggle”
- Can have N zeros
- Sensitive to all u_j - “Global Support”
- Can have large variation between points

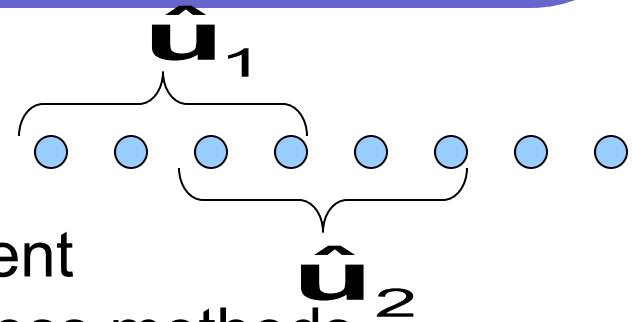
- Local Interpolation

- Use subset of nodes to represent solution over localized areas – so-called finite elements
- Possibilities : Overlapping Domains, Non-overlapping Domains

Polynomial Basis/Weighting Functions

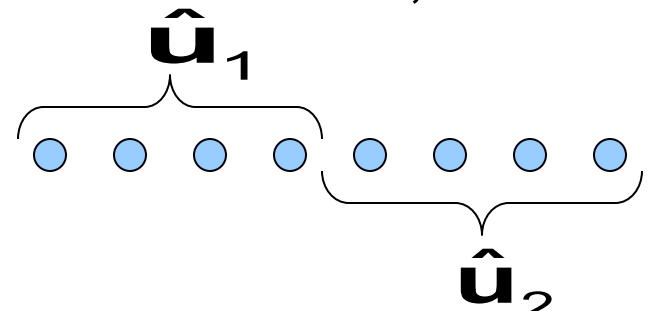
- Overlapping Domains

- Nonunique
- Somewhat difficult to implement
- Further study of this in meshless methods²



- Non-overlapping Domains

- Idea of finite “element” as collection of nodes, i.e. unit of local interpolation
- Unequal spacing of nodes, i.e. conformity of shape
- Basic building block of FEM
- Typically use same type of element throughout for programming ease (unless reason not to do so)



Polynomial Basis/Weighting Functions

- Continuity of $\hat{\mathbf{u}}$ In 1D, if $N+1$ nodes/element
 - $\hat{\mathbf{u}}$ is locally N^{th} order polynomial
 - On element **interior** ... 1st N derivatives continuous
 - ... but, at element boundaries only $\hat{\mathbf{u}}$ is continuous; $\frac{\partial \hat{\mathbf{u}}}{\partial \mathbf{x}}$ changes abruptly
 - “ C^0 ” continuity \rightarrow continuous in 0th derivative
- Higher order continuity possible
 - e.g. “ C^1 ” continuity ... need Hermite polynomial
 - Simplest local unit: Hermite cubic
 - $\frac{\partial \hat{\mathbf{u}}}{\partial \mathbf{x}}$ becomes “nodal parameter”



A little bit of review

Topics Covered

- Difference equations and error
- Polynomials, Taylor, Lagrange, **Splines**, and basis functions
- Finite differences, truncation error, round-off error
- Numerical integration, quadrature, Gauss pts, Gaussian quadrature, truncation error, round-off error

Topics Covered

- Boundary integrals, Div, Grad, Curl
- PDE types, Anisotropy, Divergence Theorem
- Finite Difference Method for solving PDEs
- Internal FDM equations, and boundary conditions (Type 1, 2 , 3)
- Recovering flux from unused molecules

Topics Covered

- Modeling physics with FDM, incorporating heterogeneity, understanding matrix properties
- Iterative solution methods (Jacobi, Gauss seidel, SOR, **Line Methods**)
- Convergence behavior of iterative methods
- Cross derivatives in FDM

Topics Covered

- Advection-diffusive problems
- Parabolic (diffusion equation)
 - Explicit, Semi-implicit, Fully implicit
 - Potential use of higher order methods
- Stability analysis (Fourier, and Matrix Methods)
- Hyperbolic models (telegraph equation)
- FD Conservation