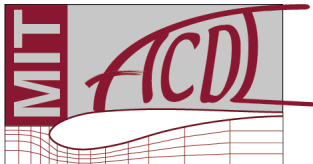
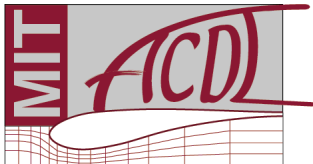

An Introduction to the Discontinuous Galerkin Method

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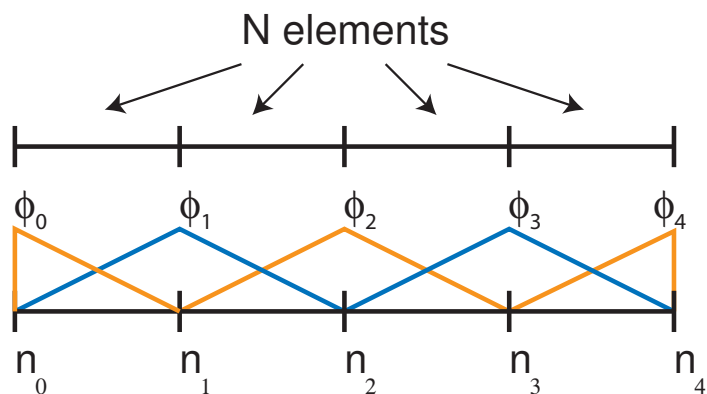
- What is DG? An Example
- Why DG for Computational Fluid Dynamics?
- Project X
- DG vs. Finite Volume
- DG for Elliptic Problems
- p -Multigrid for Higher-order DG Discretizations
- Grid Adaptation
- Conclusions and Ongoing Work



- Examine 1-D convection equation with source term:

$$au_x = \cos(x) \quad \text{on} \quad [0, \pi]$$

- Homogenous Dirichlet BCs \Rightarrow Exact solution: $u(x) = \frac{\sin(x)}{a}$
- Standard FEM procedure:



- Approximate: $u(x) = \sum_j U_j \phi_j(x)$
- ϕ_j are “hat” basis functions
- $N + 1$ basis functions in 1-D

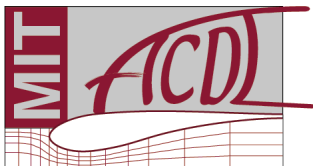
- Goal: solve for the coefficients U_j (discrete vector \mathbf{U})
- Weak form of equations:

$$\int_{\Omega} au_x \phi_i dx = \int_{\Omega} \cos(x) \phi_i dx$$
$$- \int_{\Omega} au \phi_{i,x} dx + \left[(au) \phi_i \right]_{\partial\Omega_L}^{\partial\Omega_R} = \int_{\Omega} \cos(x) \phi_i dx$$

- Substitute $u(x) = \sum_j U_j \phi_j(x)$

$$- \sum_j U_j \int_{\Omega} a \phi_j \phi_{i,x} dx + aU_N \phi_i(\pi) - aU_0 \phi_i(0) = \int_{\Omega} \cos(x) \phi_i dx$$

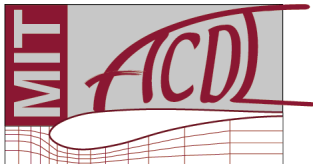
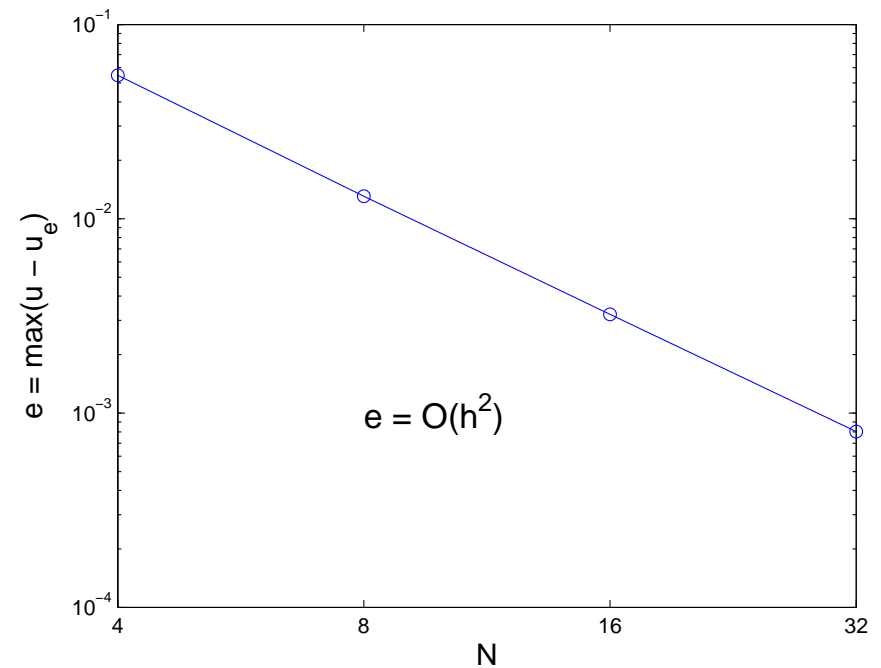
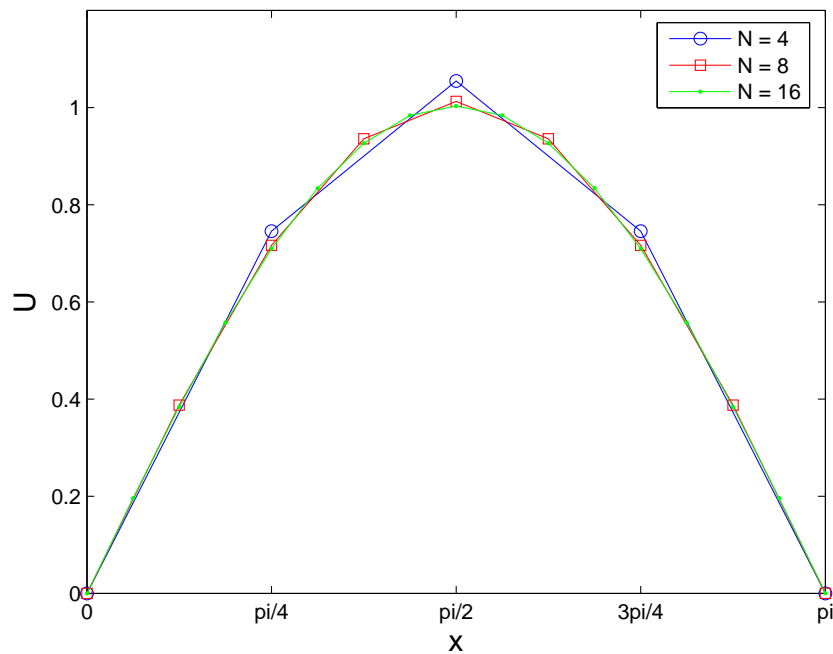
$$\text{Linear System: } \mathbf{AU} = \mathbf{F}$$



Solutions to Example



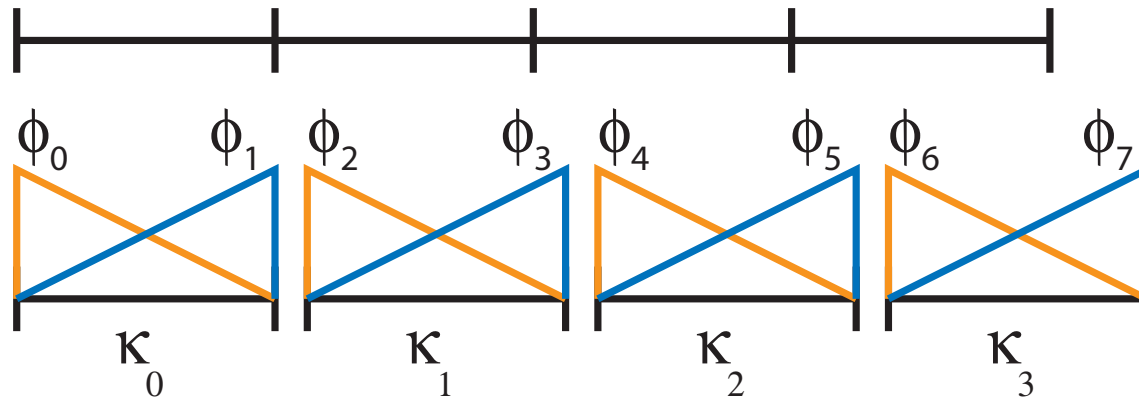
- For $a = 1$, $u_e = \sin(x)$, we obtain:



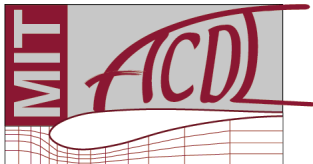
Where Discontinuous Galerkin Differs



- Formulation is the same as standard FEM.
- DG difference is in the choice of basis.
- Specifically, no continuity constraint between elements.



- ϕ_j associated with *elements*.
- $2N$ linear basis functions in 1-D.



$$\int_{\Omega} au_x \phi_i dx = \int_{\Omega} \cos(x) \phi_i dx$$

$$\sum_{\kappa} \int_{\kappa} au_x \phi_i dx = \sum_{\kappa} \int_{\kappa} \cos(x) \phi_i dx$$

$$\sum_{\kappa} \left\{ - \int_{\kappa} au \phi_{i,x} dx + \left[(au) \phi_i \right]_{\kappa-1/2}^{\kappa+1/2} \right\} = \sum_{\kappa} \int_{\kappa} \cos(x) \phi_i dx$$

$$- \int_{\kappa} au \phi_{i,x} dx + \left[(au) \phi_i \right]_{\kappa-1/2}^{\kappa+1/2} = \int_{\kappa} \cos(x) \phi_i dx$$

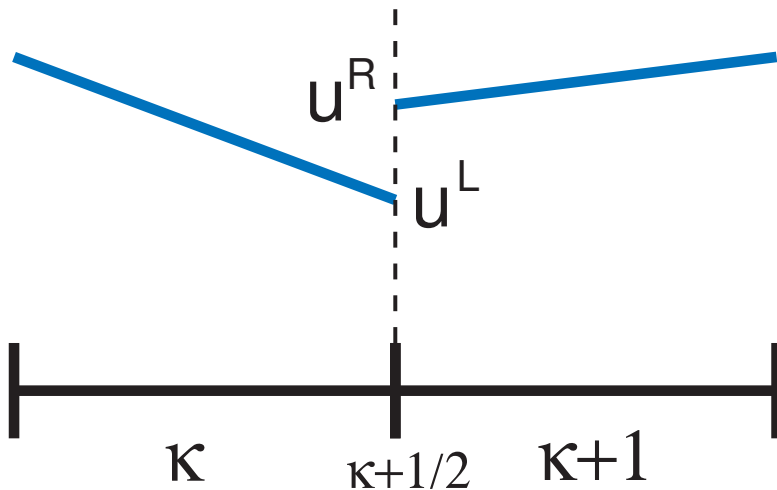
What About Element Interfaces?



- Discontinuity creates problem in the weak form:

$$-\int_{\kappa} au\phi_{i,x}dx + \left[(au)\phi_i \right]_{\kappa-1/2}^{\kappa+1/2} = \int_{\kappa} \cos(x)\phi_i dx$$

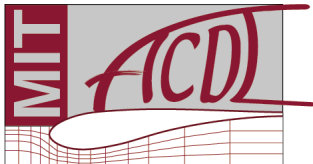
- (au) is multi-valued at element interfaces.



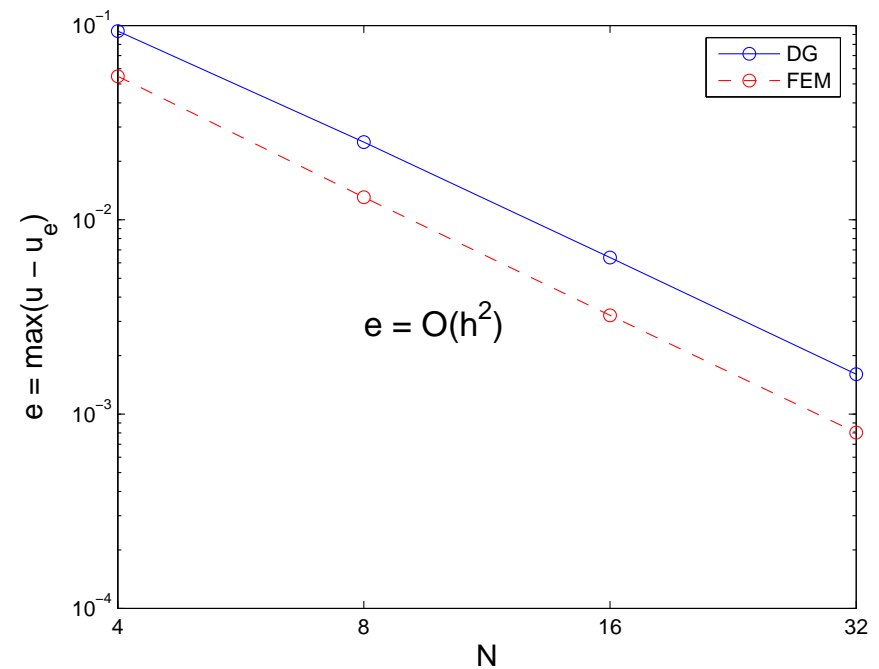
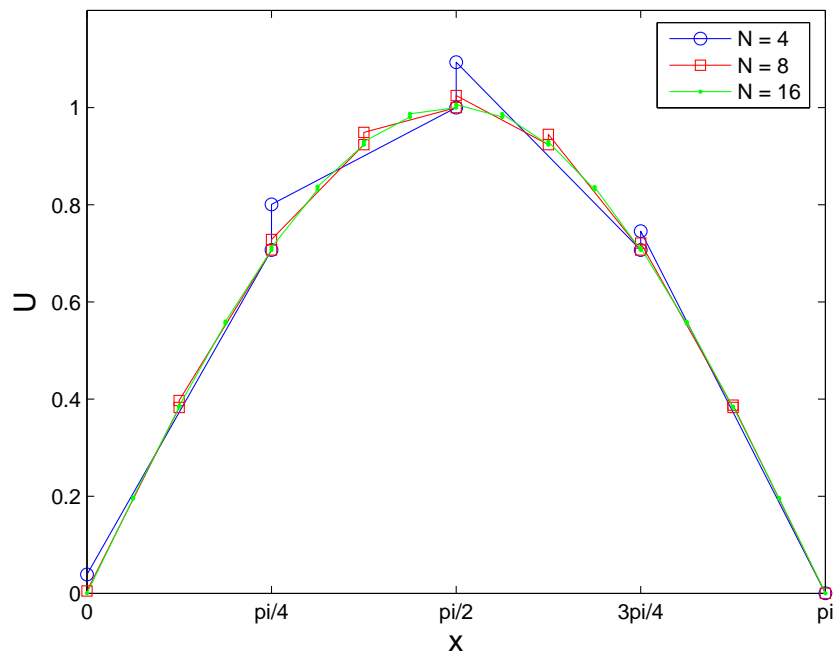
- Fortunately, this problem has been well studied by the Finite Volume community.
- Solution: use numerical flux function

$$(au) \rightarrow \mathcal{H}(u^L, u^R, a).$$

- In this case use pure upwinding.



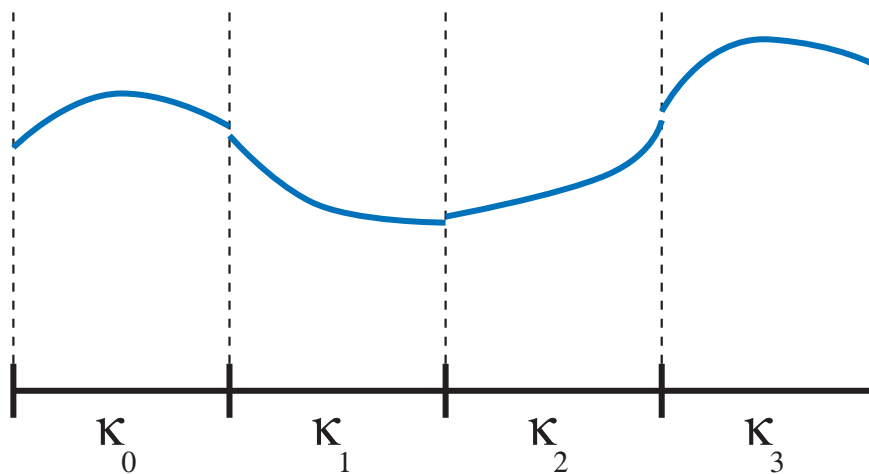
- For the same 1-D problem, $u_e = \sin(x)$, we obtain:



- DG seems to add additional DOFs for no reason.
- Why do we want to use DG over standard FEM?

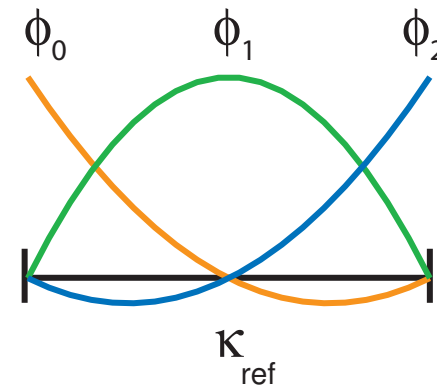
- Main Benefit:

Easy to Introduce Higher Order ($p > 1$) interpolation



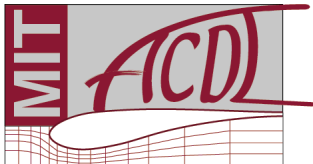
$(p + 1)$ Basis Functions per element

Lagrange Basis
Functions: $p = 2$



- Basic iterative solution schemes (e.g. block Jacobi) are stable for higher order (i.e. no need for multistaging).
- Discretization lends itself to solution via p-multigrid.
- Easy parallelization.
- In the field of CFD for aerodynamics, these benefits seem to outweigh the cost of extra DOFs.

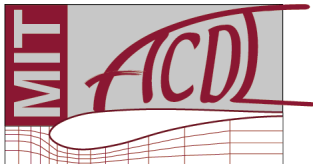
- DG solver package for computational fluid dynamics.
- Started in the Fall of 2002.
- Involved in seven M.S. theses and six ongoing Ph.D. theses.
- Currently:
 - ▶ Approximately 100,000 lines of code.
 - ▶ Capability ranges from grid generation to visualization.
 - ▶ Various side projects: grid adaptation, geometry optimization, turbulence modeling, shock limiting, unsteady studies.
 - ▶ End goal is full Reynolds-Averaged Navier-Stokes in 3-D.



- Project X Team Goal:
 - ▶ To improve the aerothermal design process for complex 3D configurations by significantly reducing the time from geometry to solution at engineering-required accuracy using high-order adaptive methods

Who are we?

- Todd Oliver
- Chris Fidkowski
- Garrett Barter
- Bob Haimes
- Prof. David Darmofal
- Prof. Jaime Peraire
- Tan Bui
- Shannon Cheng
- James Lu
- Pete Whitney
- Doug Quattrochi

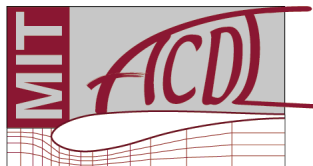
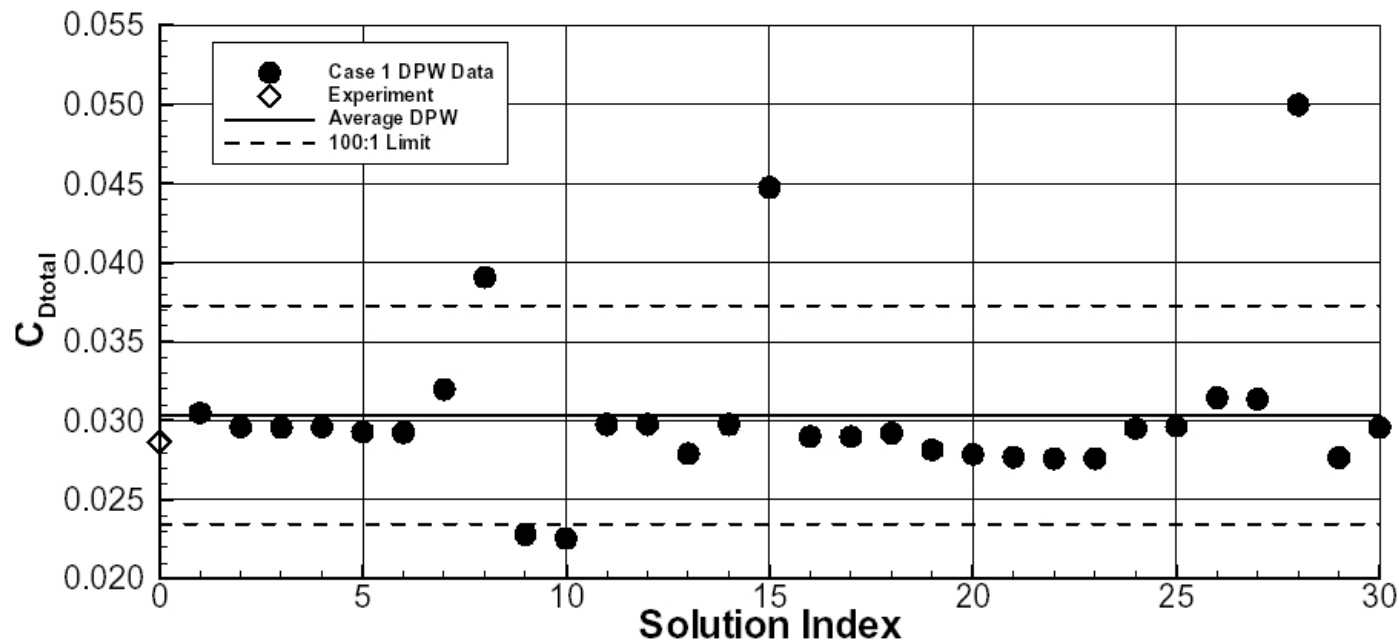


Accuracy of Current CFD Methods



Results of an AIAA Drag Prediction Workshop comparing the performance of industry-standard CFD codes:

AIAA-2002-0841



- State of CFD in applied aerodynamics
 - ▶ Finite-volume with at best second order accuracy
 - ▶ Questions exist whether current discretizations are capable of achieving desired accuracy levels in practical time
- Decrease computational time and gridding requirements by increasing solution order

$$\log T = wd \left(-\frac{1}{p} \log E + \log p \right) - \log F + \text{const}$$

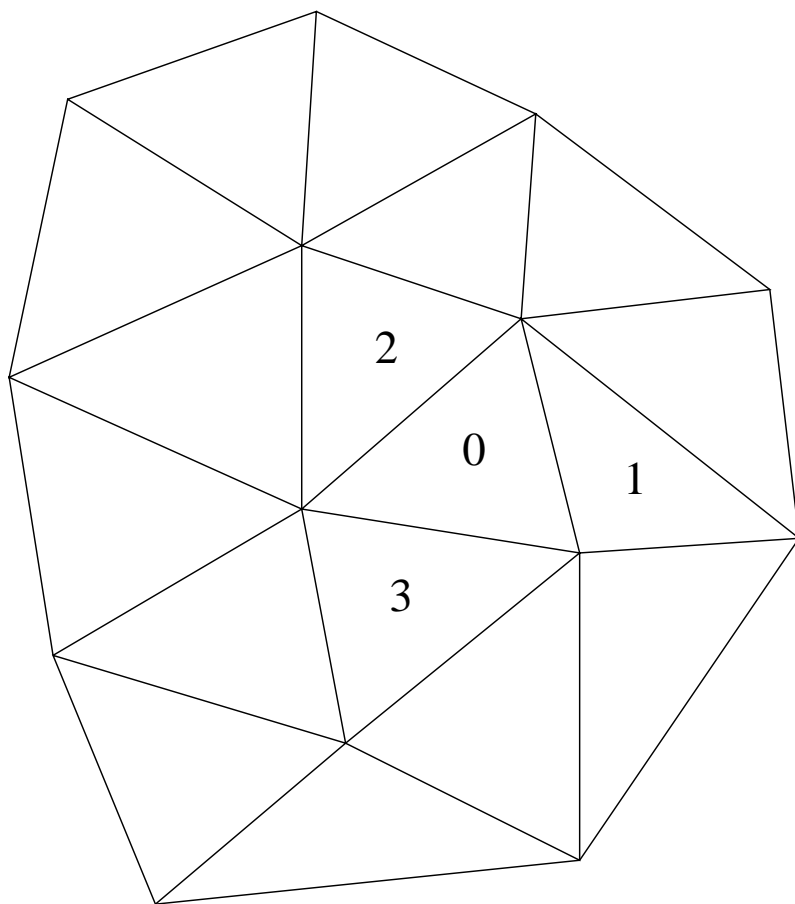
- T = time to solution
- p = discretization order
- E = desired error level ($E \ll 1$)
- w = solution complexity
- d = dimension of problem
- F = computational speed



$$\mathbf{u}_t + \nabla \cdot \mathcal{F}_i(\mathbf{u}) = 0$$

- \mathbf{u} is the state vector and \mathcal{F}_i is the inviscid flux.
- In two dimensions:

$$\mathbf{u} = \begin{bmatrix} \rho \\ \rho u \\ \rho v \\ \rho E \end{bmatrix} \quad \mathcal{F}_i^x = \begin{bmatrix} \rho u \\ \rho u^2 + p \\ \rho uv \\ \rho uH \end{bmatrix}, \quad \mathcal{F}_i^y = \begin{bmatrix} \rho v \\ \rho uv \\ \rho v^2 + p \\ \rho vH \end{bmatrix}.$$

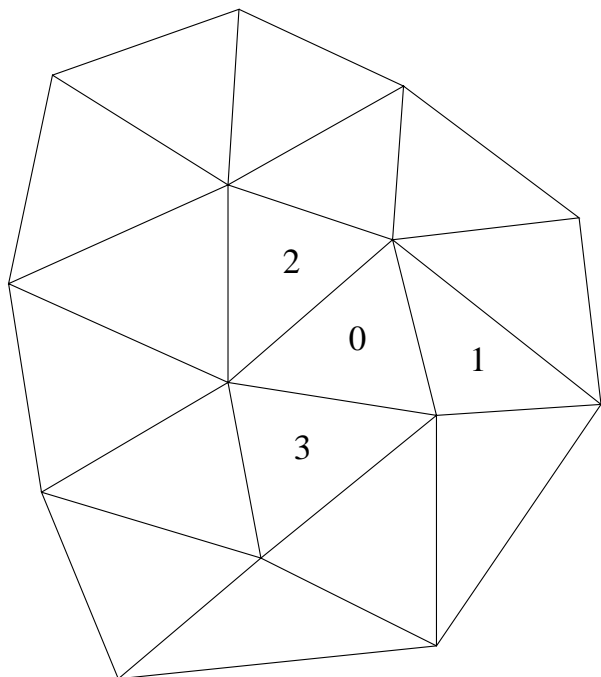


Integrate Euler equations over triangle 0 and use Green's theorem:

$$\frac{d}{dt} \int_{A_0} \mathbf{u} \, d\mathbf{x} + \sum_{k=1}^3 \int_{0k} \mathcal{F}_i(\mathbf{u}) \cdot \hat{\mathbf{n}} \, ds = 0$$

If we assume \mathbf{u} is constant in each triangle ...

\mathbf{u} constant in each triangle:



$$\frac{d\mathbf{u}_0}{dt} A_0 + \sum_{k=1}^3 \int_{0k} \mathcal{H}_i(\mathbf{u}_0, \mathbf{u}_k, \hat{\mathbf{n}}_{0k}) ds = 0$$

$\mathcal{H}_i(\mathbf{u}_L, \mathbf{u}_R, \hat{\mathbf{n}}_{LR})$ is flux function that determines inviscid flux in $\hat{\mathbf{n}}_{LR}$ direction from left and right states, \mathbf{u}_L and \mathbf{u}_R .

Example flux functions: Godunov, Roe, Osher, Van Leer, Lax-Friedrichs, etc.

This discretization has a solution error which is $O(h)$ where h is mesh size.

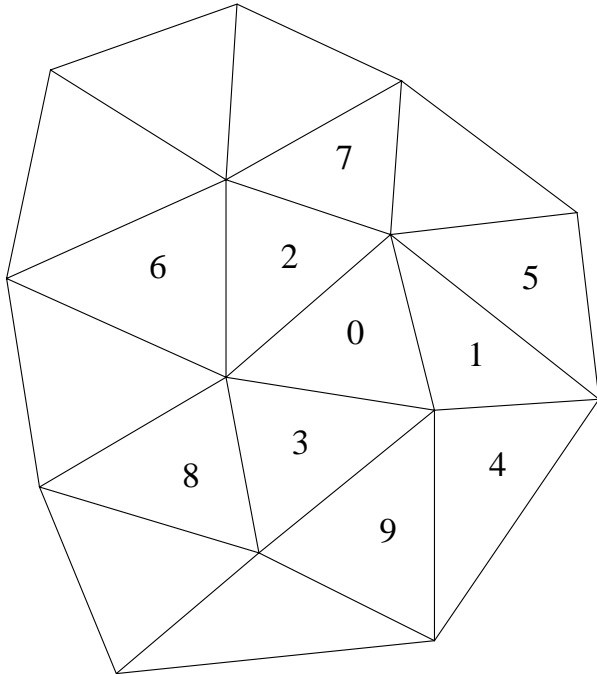
In each triangle, reconstruct a linear solution, $\tilde{\mathbf{u}}$, using neighboring averages:

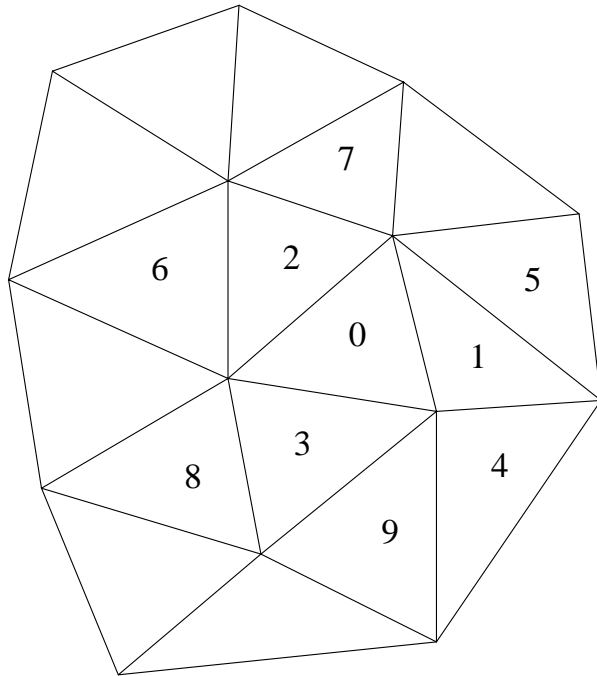
$$\begin{aligned}\tilde{\mathbf{u}}_0 &\equiv \mathbf{u}_0 + (\mathbf{x} - \mathbf{x}_0) \cdot \nabla \mathbf{u}_0, \\ \nabla \mathbf{u}_0 &\equiv \nabla \mathbf{u}_0(\mathbf{u}_0, \mathbf{u}_1, \mathbf{u}_2, \mathbf{u}_3).\end{aligned}$$

Apply conservation law on triangle:

$$\frac{d\mathbf{u}_0}{dt} A_0 + \sum_{k=1}^3 \int_{0k} \mathcal{H}_i(\tilde{\mathbf{u}}_0, \tilde{\mathbf{u}}_k, \hat{\mathbf{n}}_{0k}) ds = 0$$

On smooth meshes and flows, solution error is $O(h^2)$.





- + Increased accuracy on given mesh without additional degrees of freedom
- Difficulty in achieving higher-order on unstructured meshes and near boundaries
- Stabilizing multi-stage methods necessary for local iterative schemes
- Matrix fill-in results in high-memory requirements

Consider steady state problem and define discrete residual for cell j ,

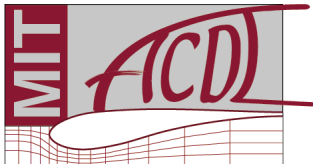
$$\mathbf{R}_j(\mathbf{u}) \equiv \sum_{k=1}^3 \int_{jk} \mathcal{H}_i(\tilde{\mathbf{u}}_j, \tilde{\mathbf{u}}_k, \hat{\mathbf{n}}_{jk}) ds = 0.$$

A Jacobi iterative method to solve this problem is,

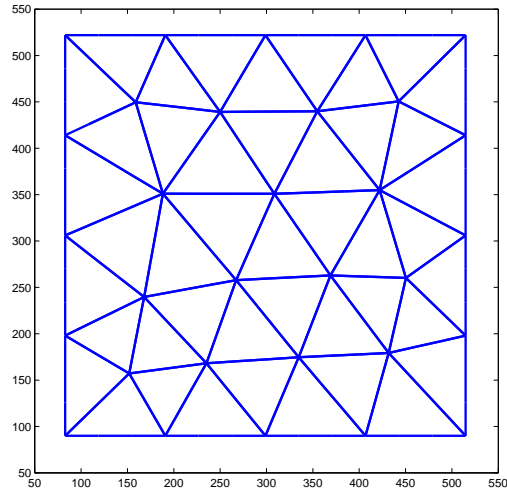
$$\mathbf{u}_j^{n+1} = \mathbf{u}_j^n - \omega (\partial \mathbf{R}_j / \partial \mathbf{u}_j)^{-1} \mathbf{R}_j(\mathbf{u}).$$

For any finite ω , Jacobi is unstable for higher-order. One solution is a multi-stage method,

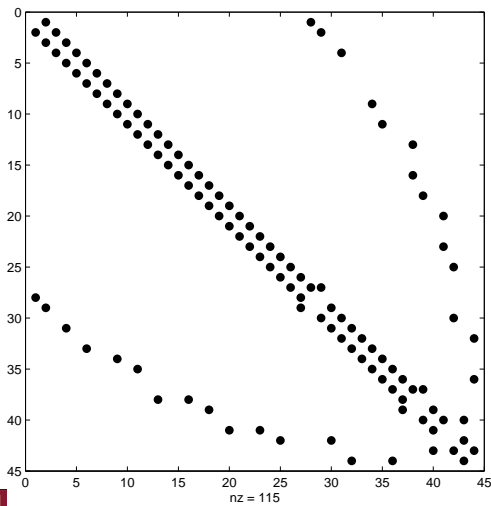
$$\begin{aligned} \hat{\mathbf{u}}_j &= \mathbf{u}_j^n - \hat{\omega} (\partial \mathbf{R}_j / \partial \mathbf{u}_j)^{-1} \mathbf{R}_j(\mathbf{u}^n) \\ \mathbf{u}_j^{n+1} &= \mathbf{u}_j^n - \omega (\partial \mathbf{R}_j / \partial \mathbf{u}_j)^{-1} \mathbf{R}_j(\hat{\mathbf{u}}) \end{aligned} \quad \Leftarrow \text{Requires two residual evaluations.}$$



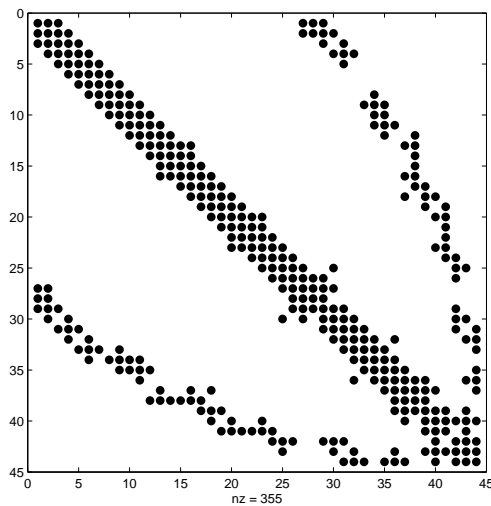
Matrix Fill for Higher-order Finite Volume



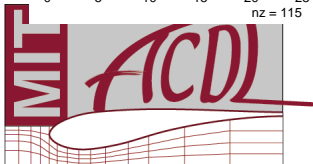
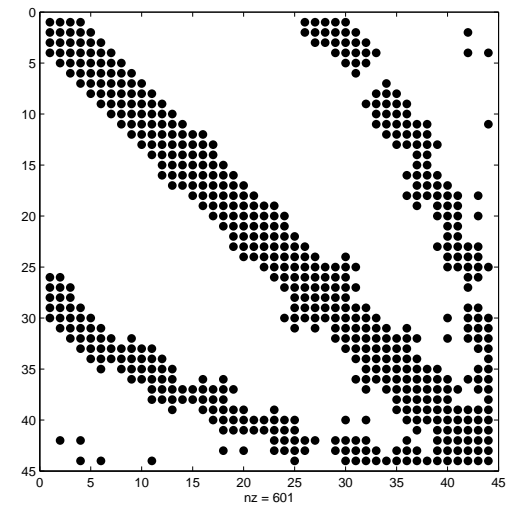
First-order



Second-order



Third-order



- Start from strong form of governing equations:

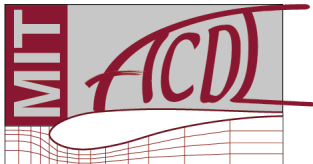
$$\mathbf{u}_t + \nabla \cdot \mathcal{F}_i(\mathbf{u}) = 0.$$

- Look for a solution $\mathbf{u}_h \in \mathcal{V}_h^p$.
- Multiply governing equation by weight function $\mathbf{v}_h \in \mathcal{V}_h^p$ and integrate over element $\kappa \in T_h$:

$$\int_{\kappa} \mathbf{v}_h^T [(\mathbf{u}_h)_t + \nabla \cdot \mathcal{F}_i] d\mathbf{x} = 0.$$

- Integrate second term by parts (assume interior element):

$$\int_{\kappa} \mathbf{v}_h^T (\mathbf{u}_h)_t d\mathbf{x} - \int_{\kappa} \nabla \mathbf{v}_h^T \cdot \mathcal{F}_i d\mathbf{x} + \int_{\partial\kappa} \mathbf{v}_h^{+T} \mathcal{H}_i(\mathbf{u}_h^+, \mathbf{u}_h^-, \hat{\mathbf{n}}) ds = 0.$$



- DG weighted residual:

$$\int_{\kappa} \mathbf{v}_h^T (\mathbf{u}_h)_t d\mathbf{x} - \int_{\kappa} \nabla \mathbf{v}_h^T \cdot \mathcal{F}_i d\mathbf{x} + \int_{\partial\kappa} \mathbf{v}_h^{+T} \mathcal{H}_i(\mathbf{u}_h^+, \mathbf{u}_h^-, \hat{\mathbf{n}}) ds = 0.$$

- For $p = 0$ solution, this reduces to:

$$(\mathbf{u}_{\kappa})_t A_{\kappa} + \int_{\partial\kappa} \mathcal{H}_i(\mathbf{u}_h^+, \mathbf{u}_h^-, \hat{\mathbf{n}}) ds = 0.$$

- Thus, $p = 0$ DG is identical to first-order finite volume.

- Find $\mathbf{u}_h \in \mathcal{V}_h^p$ such that $\forall \mathbf{v}_h \in \mathcal{V}_h^p$,

$$\sum_{\kappa \in T_h} \left\{ \int_{\kappa} \mathbf{v}_h^T(\mathbf{u}_h)_t d\mathbf{x} - \int_{\kappa} \nabla \mathbf{v}_h^T \cdot \mathcal{F}_i d\mathbf{x} \right\} + \int_{\Gamma_i} \mathbf{v}_h^{+T} \mathcal{H}_i(\mathbf{u}_h^+, \mathbf{u}_h^-, \hat{\mathbf{n}}) ds + \int_{\partial\Omega} \mathbf{v}_h^{+T} \mathcal{H}_i^b(\mathbf{u}_h^+, \mathbf{u}_h^b, \hat{\mathbf{n}}) ds = 0.$$

- Boundary conditions enforced weakly through $\mathcal{H}_i^b(\mathbf{u}_h^+, \mathbf{u}_h^b, \hat{\mathbf{n}})$ where \mathbf{u}_h^b is determined from desired boundary conditions and outgoing characteristics.
- For smooth problems, the error of this scheme is expected to be $O(h^{p+1})$.

- Increased accuracy on given mesh requires additional degrees of freedom
- + Higher-order accuracy not hampered on unstructured meshes nor near boundaries
- + Local iterative methods are stable
- + Matrix fill-in maintains block sparsity of $p = 0$

$$\int_{\kappa} \mathbf{v}_h^T (\mathbf{u}_h)_t d\mathbf{x} - \int_{\kappa} \nabla \mathbf{v}_h^T \cdot \mathcal{F}_i d\mathbf{x} + \int_{\partial\kappa} \mathbf{v}_h^{+T} \mathcal{H}_i(\mathbf{u}_h^+, \mathbf{u}_h^-, \hat{\mathbf{n}}) ds = 0.$$

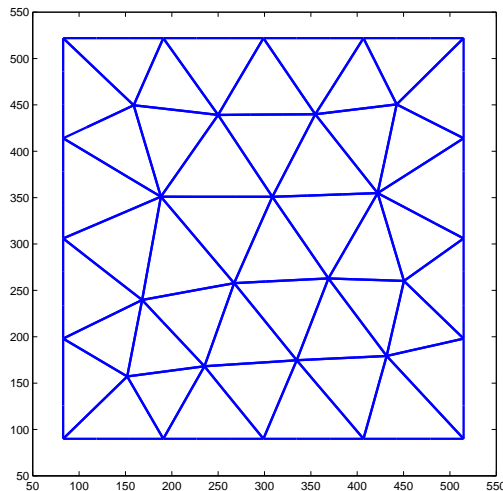
An elemental block Jacobi iterative method to solve this problem is,

$$\mathbf{u}_j^{n+1} = \mathbf{u}_j^n - \omega (\partial \mathbf{R}_j / \partial \mathbf{u}_j)^{-1} \mathbf{R}_j(\mathbf{u}).$$

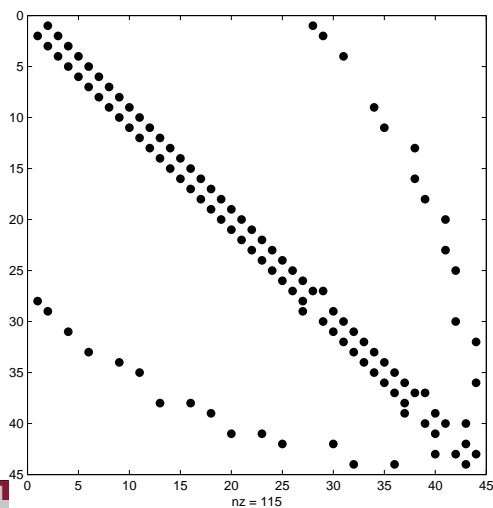
where $\partial \mathbf{R}_j / \partial \mathbf{u}_j$ is the diagonal block for the element j .

For $0 < \omega < 1$, elemental block Jacobi is stable independent of p .

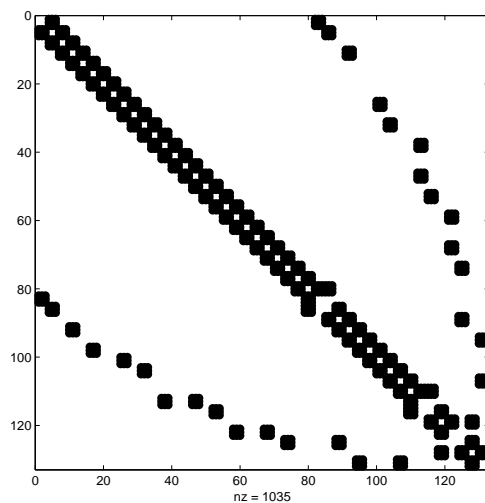
Matrix Fill for Higher-order DG



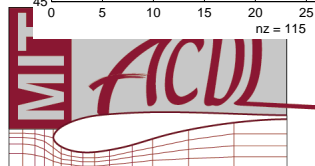
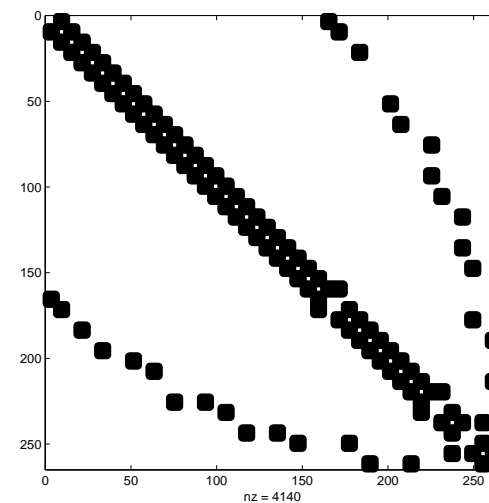
First-order ($p = 0$)



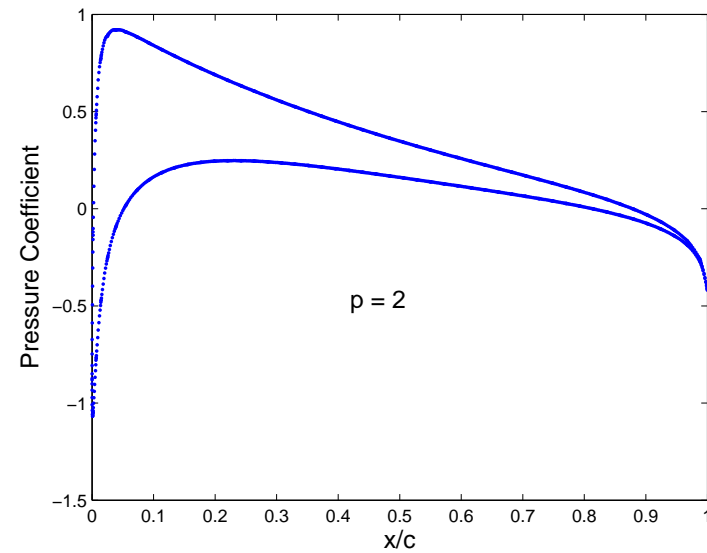
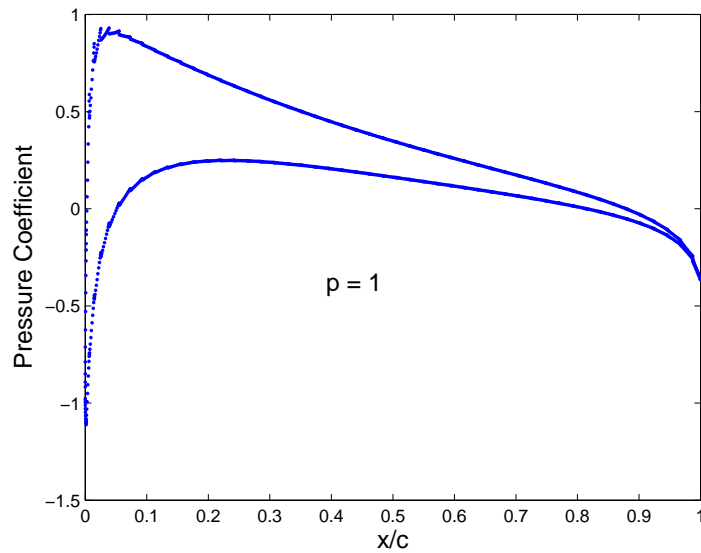
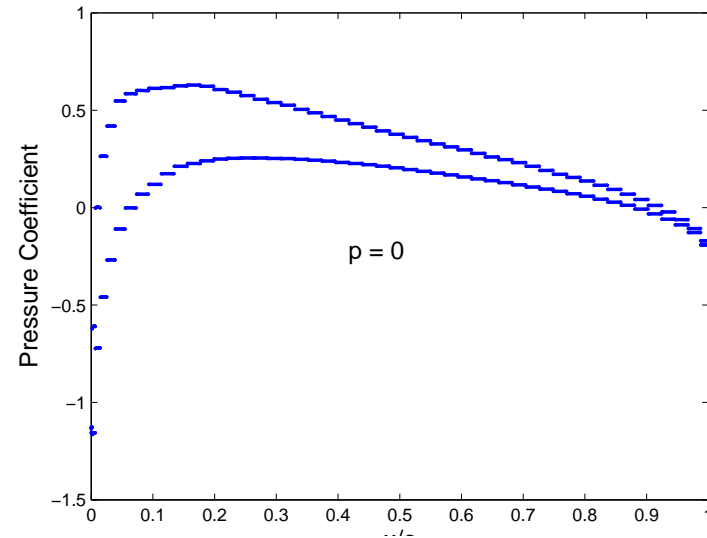
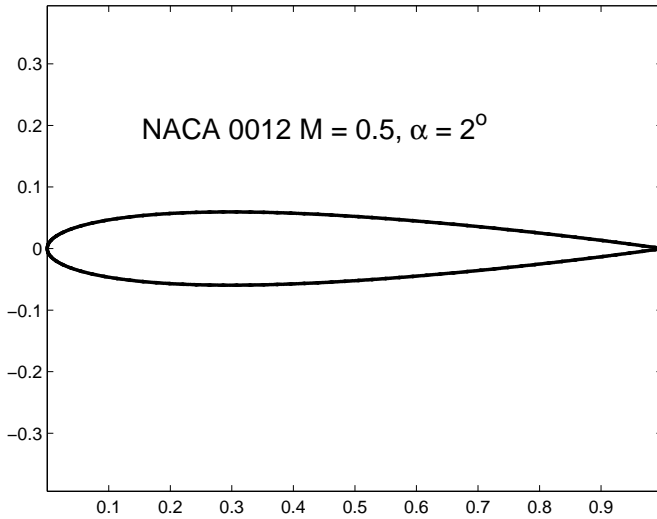
Second-order ($p = 1$)



Third-order ($p = 2$)



Inviscid Flow Example



- Navier-Stokes Equations: $\mathbf{u}_t + \nabla \cdot \mathcal{F}_i(\mathbf{u}) - \nabla \cdot \mathcal{F}_v(\mathbf{u}, \nabla \mathbf{u}) = 0$
- $\mathcal{F}_v = \mathcal{A}_v \nabla \mathbf{u} = (\mathbf{F}_v^x, \mathbf{F}_v^y)$ is the viscous flux vector

$$\mathbf{F}_v^x = \begin{pmatrix} 0 \\ \frac{2}{3}\mu(2\frac{\partial u}{\partial x} - \frac{\partial v}{\partial y}) \\ \mu(\frac{\partial u}{\partial y} + \frac{\partial v}{\partial x}) \\ \frac{2}{3}\mu(2\frac{\partial u}{\partial x} - \frac{\partial v}{\partial y})u + \mu(\frac{\partial u}{\partial y} + \frac{\partial v}{\partial x})v + \kappa\frac{\partial T}{\partial x} \end{pmatrix},$$

$$\mathbf{F}_v^y = \begin{pmatrix} 0 \\ \mu(\frac{\partial u}{\partial y} + \frac{\partial v}{\partial x}) \\ \frac{2}{3}\mu(2\frac{\partial v}{\partial y} - \frac{\partial u}{\partial x}) \\ \frac{2}{3}\mu(2\frac{\partial v}{\partial y} - \frac{\partial u}{\partial x})v + \mu(\frac{\partial u}{\partial y} + \frac{\partial v}{\partial x})u + \kappa\frac{\partial T}{\partial y} \end{pmatrix}$$

- Model problem for viscous terms of N-S: 1-D, scalar Poisson's equation

$$-u_{xx} = f \quad \text{on} \quad [-1, 1]$$

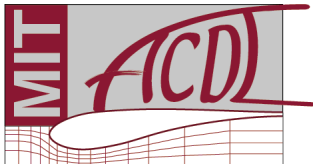
- Proceed as for Euler:

- ▶ Triangulate domain into non-overlapping elements $\kappa \in T_h$
- ▶ Define solution and test function space \mathcal{V}_h^p

- Discrete formulation: Find $u_h \in \mathcal{V}_h^p$ such that $\forall v_h \in \mathcal{V}_h^p$,

$$\sum_{\kappa \in T_h} \left\{ - [v_h \widehat{u}_x]_{x_{\kappa-1/2}}^{x_{\kappa+1/2}} + \int_{\kappa} (v_h)_x (u_h)_x dx \right\} = \sum_{\kappa \in T_h} \left\{ \int_{\kappa} v_h f dx \right\}$$

- Need to define \widehat{u}_x



- No upwinding mechanism \Rightarrow choose central flux

$$\widehat{u}_x = \frac{1}{2} \left((u_h)_x^L + (u_h)_x^R \right)$$

- Discrete formulation becomes: Find $u_h \in \mathcal{V}_h^p$ such that $\forall v_h \in \mathcal{V}_h^p$,

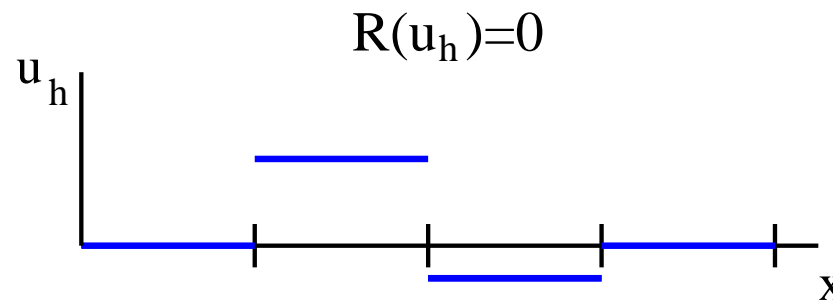
$$\sum_{\kappa \in T_h} \left\{ - \left[\frac{1}{2} v_h \left((u_h)_x^L + (u_h)_x^R \right) \right]_{x_{\kappa-1/2}}^{x_{\kappa+1/2}} + \int_{\kappa} (v_h)_x (u_h)_x dx \right\} = \sum_{\kappa \in T_h} \left\{ \int_{\kappa} v_h f dx \right\}$$

- PROBLEM: Scheme is inconsistent!

- Examine Laplace's equation with homogeneous Dirichlet BCs

$$\begin{aligned} -u_{xx} &= 0 \quad \text{on } [-1, 1] \\ u(-1) &= u(1) = 0 \end{aligned}$$

- Exact solution: $u(x) = 0$



- If $(u_h)_x = 0$ everywhere, discrete equations satisfied exactly regardless of magnitude of u_h

- Introduce new variable, $q = u_x$, such that

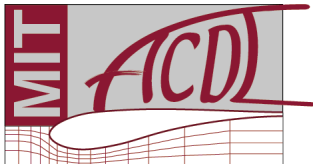
$$\begin{aligned} -q_x &= f \\ q - u_x &= 0 \end{aligned}$$

- Discrete formulation: Find $u_h \in \mathcal{V}_h^p$ and $q_h \in \mathcal{V}_h^p$ such that $\forall v_h \in \mathcal{V}_h^p$ and $\forall \tau_h \in \mathcal{V}_h^p$,

$$\sum_{\kappa \in T_h} \left\{ - \left[v_h \hat{q} \right]_{x_{\kappa-1/2}}^{x_{\kappa+1/2}} + \int_{\kappa} (v_h)_x q_h dx \right\} - \sum_{\kappa \in T_h} \left\{ \int_{\kappa} v_h f dx \right\} = 0$$

$$\sum_{\kappa \in T_h} \left\{ \int_{\kappa} \tau_h q_h dx + \int_{\kappa} (\tau_h)_x u_h dx - \left[\tau_h \hat{u} \right]_{x_{\kappa-1/2}}^{x_{\kappa+1/2}} \right\} = 0$$

- Need to choose \hat{q} and \hat{u}



- Scheme 1: no upwinding mechanism \Rightarrow choose central fluxes

$$\hat{u} = \frac{1}{2}(u_h^L + u_h^R); \quad \hat{q} = \frac{1}{2}(q_h^L + q_h^R)$$

- ▶ Sub-optimal order of accuracy for odd p
- ▶ Stencil no longer compact

- Scheme 2: With $\llbracket s \rrbracket = s^L - s^R$ and $\{s\} = 0.5(s^L + s^R)$,

$$\hat{u} = \{u_h\}; \quad \hat{q} = \{(u_h)_x\} - \eta_f \{\delta_f\}$$

- ▶ New variable, $\delta_f(\llbracket u \rrbracket)$, is a jump penalty term.
- ▶ Optimal accuracy and compact stencil.

- Nonlinear discrete equations can be written

$$\mathbf{R}(\mathbf{u}_h) = 0$$

- Use a preconditioned iterative scheme

$$\mathbf{u}_h^{n+1} = \mathbf{u}_h^n - \mathbf{P}^{-1} \mathbf{R}(\mathbf{u}_h^n)$$

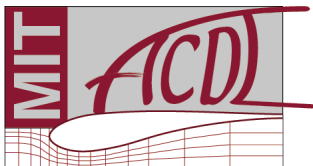
- Preconditioner

- ▶ Block-element smoothing

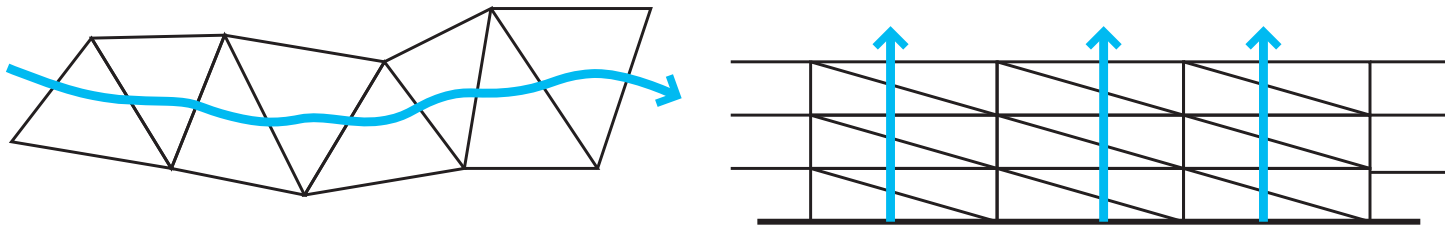
- ◆ $\mathbf{P} = \mathbf{M}_{block} \Rightarrow$ Block diagonal of the Jacobian

- ▶ Line-element smoothing

- ◆ $\mathbf{P} = \mathbf{M}_{line} \Rightarrow$ Block tridiagonal systems from Jacobian



- Motivation: Transport of information in Navier-Stokes equations characterized by convection-diffusion like phenomena
 - ▶ Inviscid regions: characteristic directions set by convection
 - ▶ Viscous regions: diffusion effects can be stronger

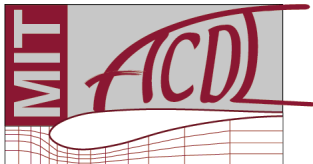
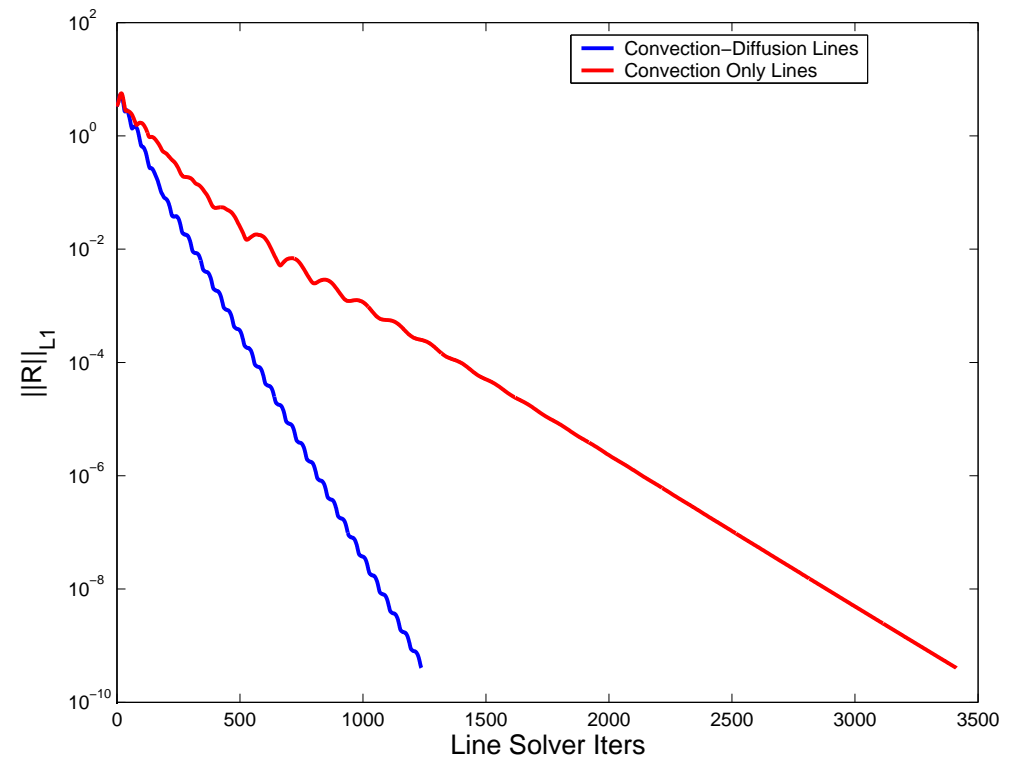
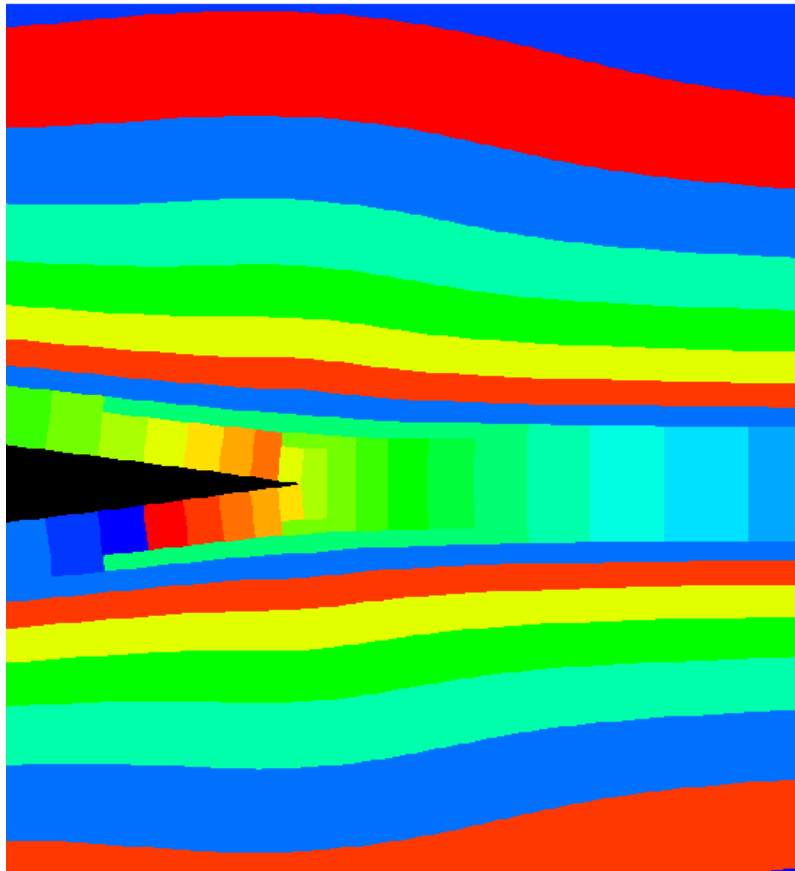


- Procedure:
 - ▶ Construct lines of elements based on measure of influence
 - ▶ Build and invert M_{line} , which is a set of block tridiagonal systems from the full Jacobian

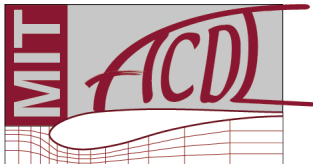
Example Lines and Performance



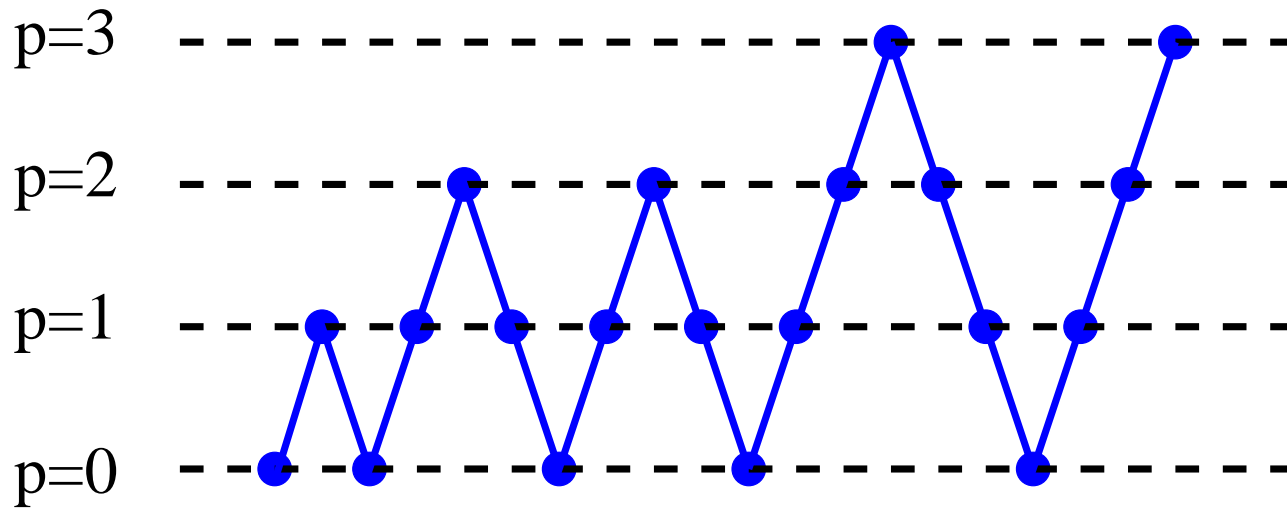
Trailing edge of NACA 0012



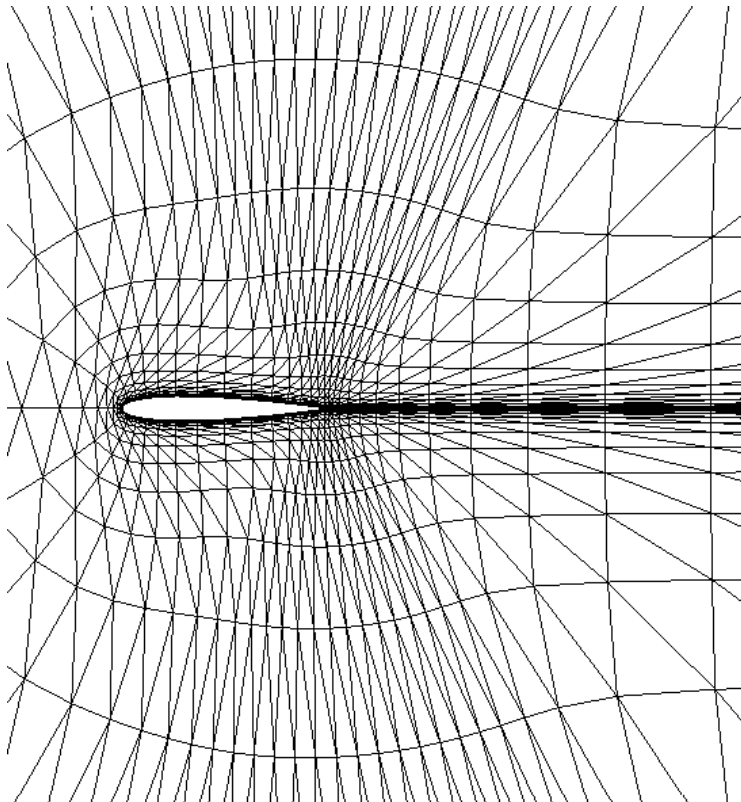
- Observation: Smoothers are inefficient at eliminating low frequency error modes on fine level
- h -Multigrid
 - ▶ Spatially coarse grid used to correct solution on fine grid
 - ▶ Grid coarsening is complex on unstructured meshes
- p -Multigrid (Ronquist & Patera, Helenbrook et al., Fidkowski & Darmofal)
 - ▶ Low order ($p - 1$) approximation used to correct high order (p) solution
 - ▶ Natural implementation in DG discretization on unstructured meshes



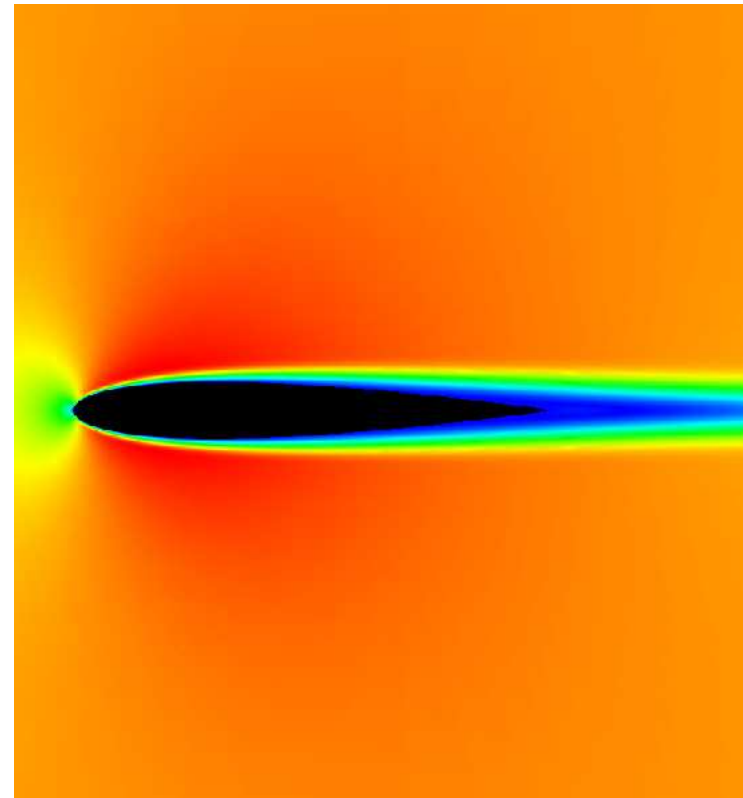
- Iterate on coarsest levels first and use these solutions as initial guesses for fine level solutions.
- Line solver used as smoother
- Full Approximation Scheme (FAS) used to deal with non-linearity



$M = 0.5$, $Re = 5000$, $\alpha = 0$
Grids are from Swanson at NASA Langley

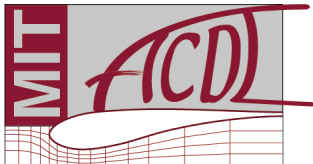
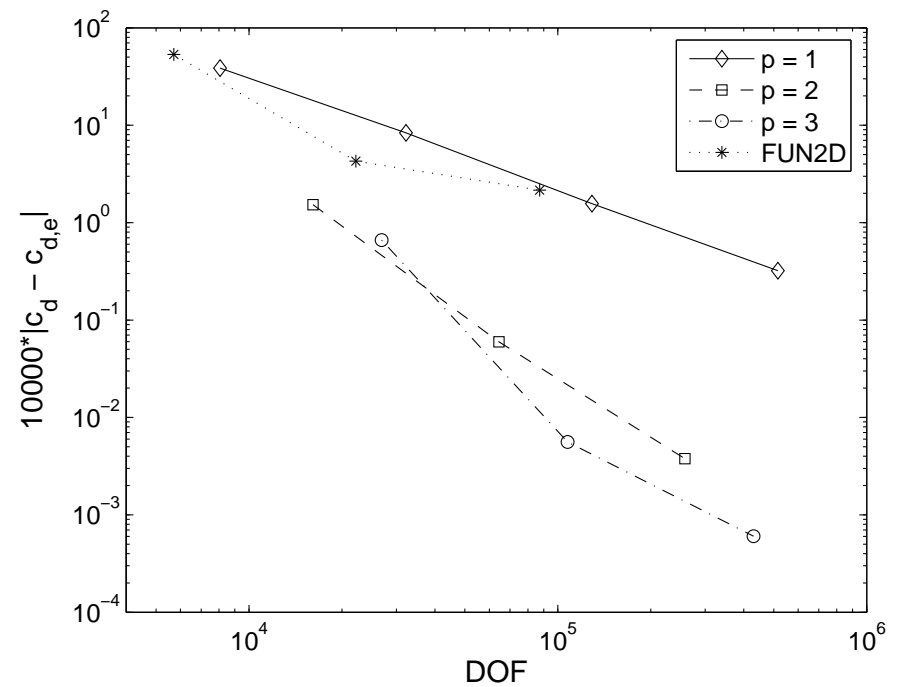
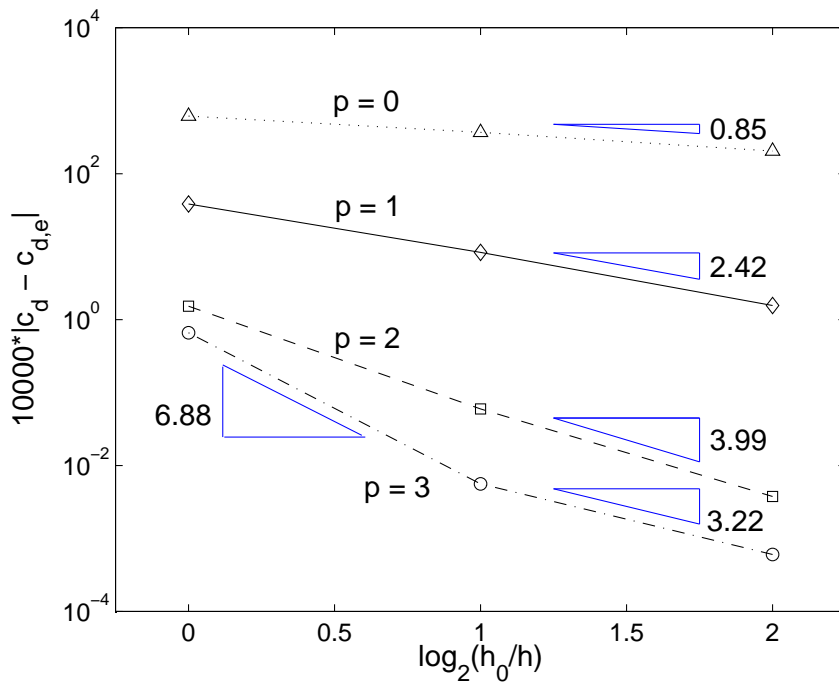


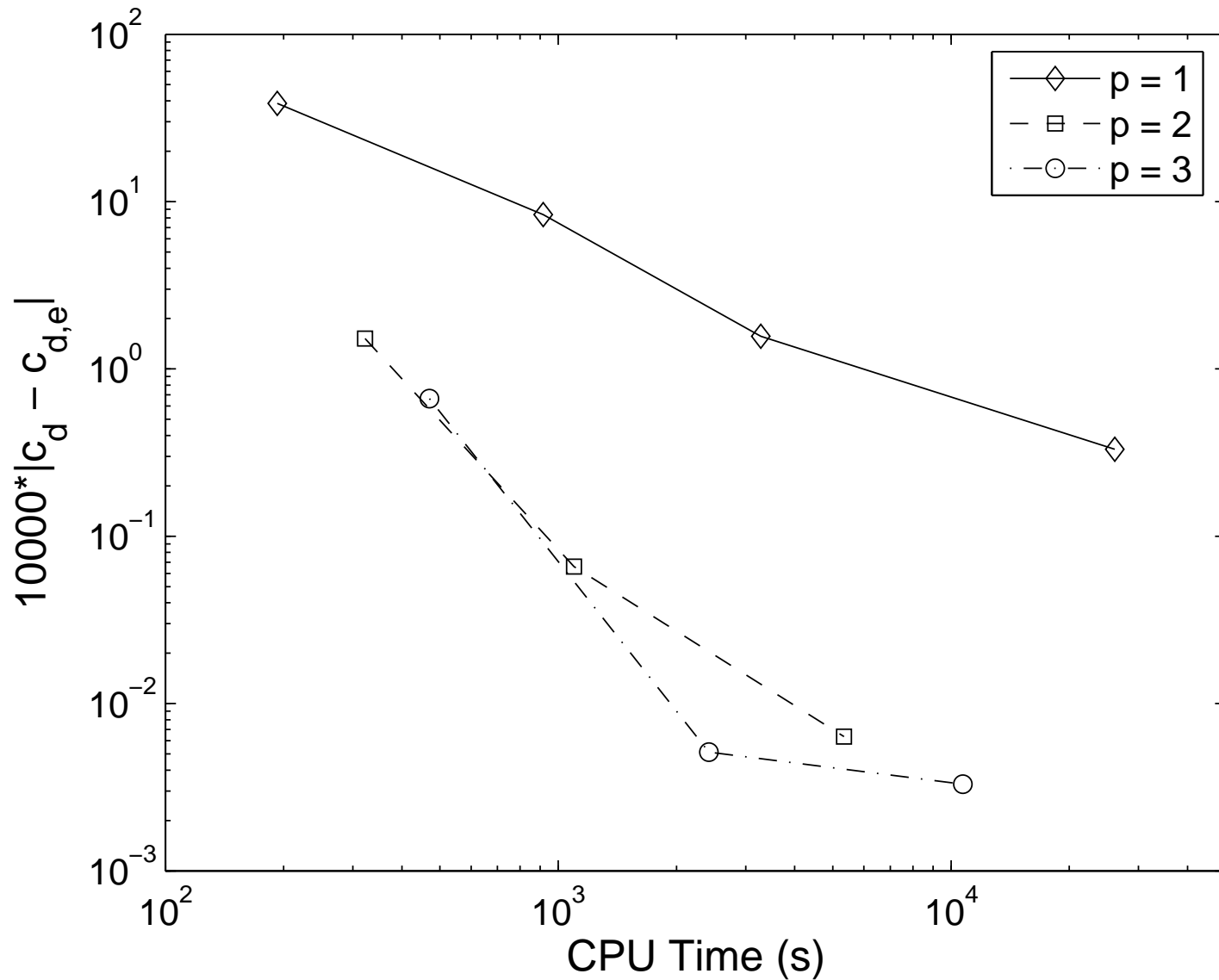
2112 element mesh



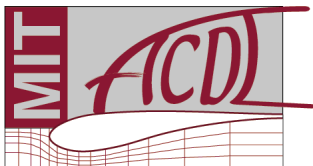
Mach contours

Drag Error Convergence





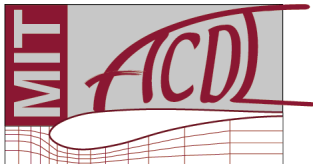
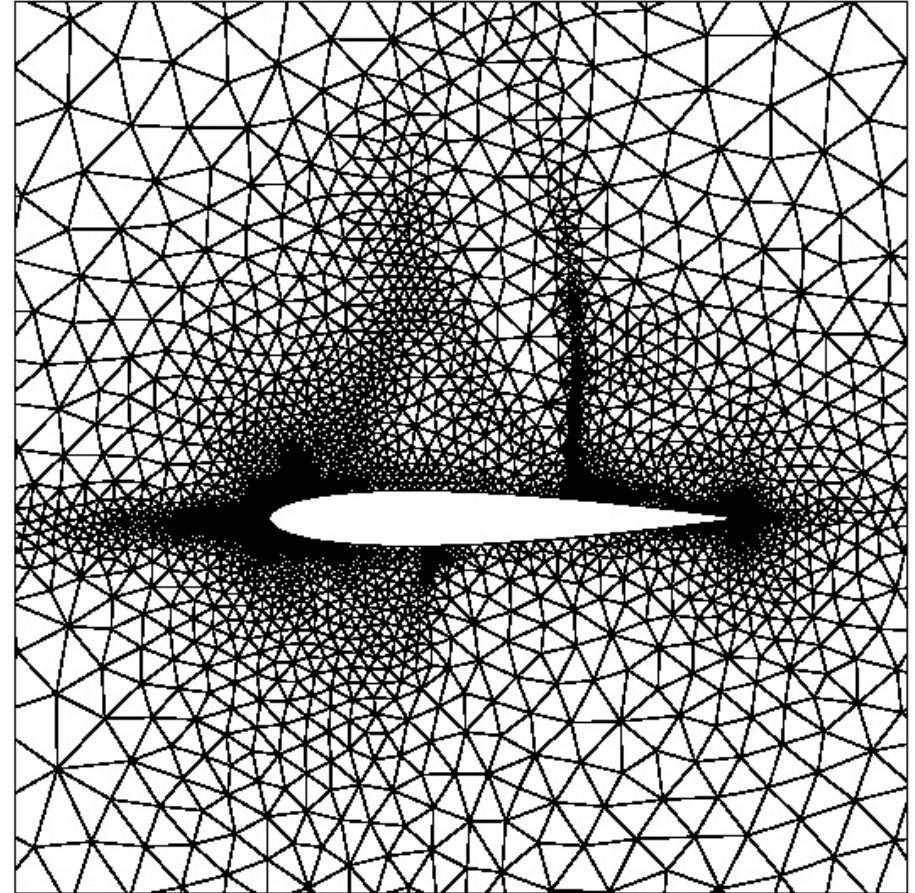
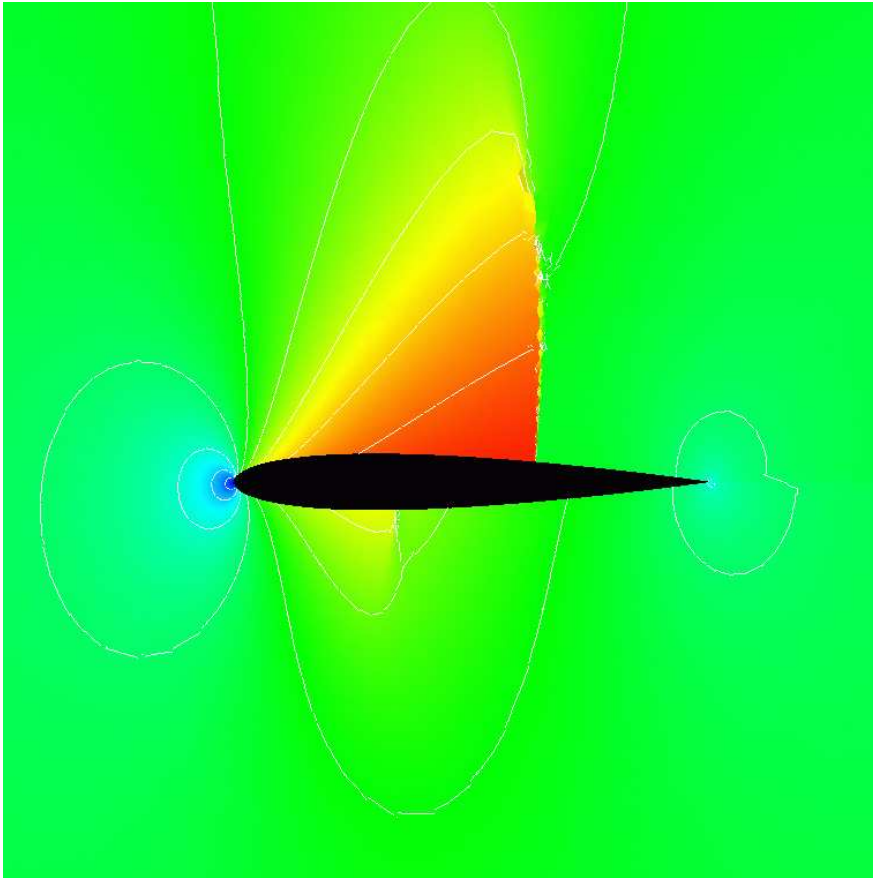
- For smooth flows, large high-order elements are ideal for accuracy and efficiency
- When singularities are present, h-adaptation is necessary to maintain accuracy
- Where to adapt? One option is output-based adaptation (Venditti & Darmofal)
 - ▶ Pick output of interest (lift, drag) and a desired error tolerance
 - ▶ Solve for output adjoint in addition to the flow solution
 - ▶ Use adjoint and flow solutions to estimate which elements contribute most/least to output error
 - ▶ Refine/coarsen mesh until the output error tolerance is met



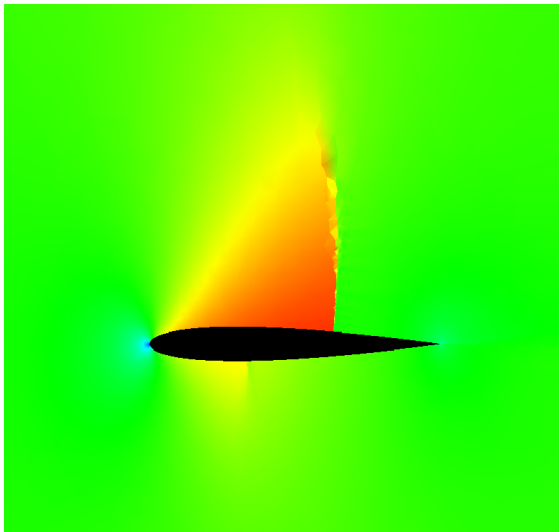
h-Adaptation Example



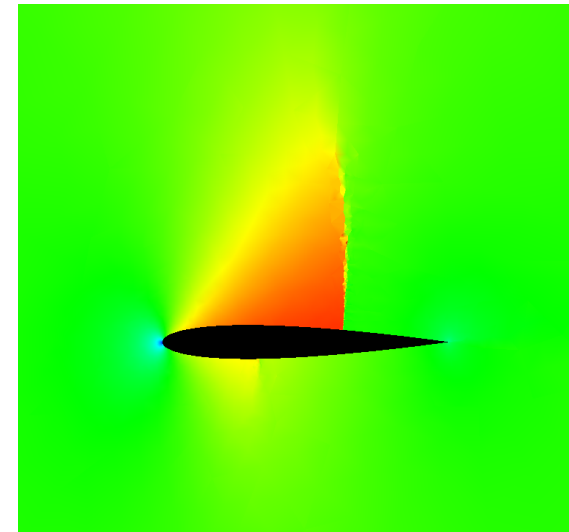
NACA 0012 Transonic test case: $M = 0.8$, $\alpha = 1.25^\circ$



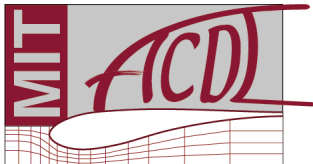
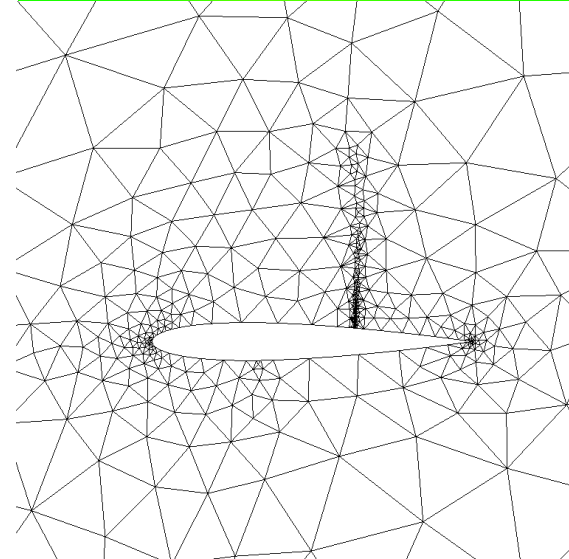
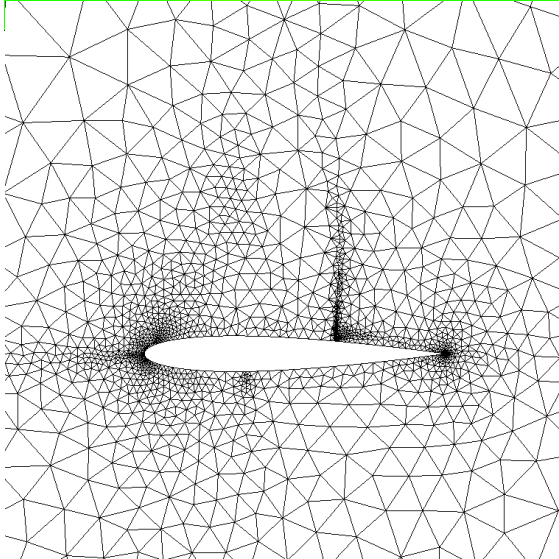
NACA 0012: $M = 0.8, \alpha = 1.25^\circ$



$$\Leftarrow \text{P1}$$
$$49,836 \text{ DOF}$$
$$|C_d - C_{d,\text{exact}}| = 4.0 * 10^{-5}$$

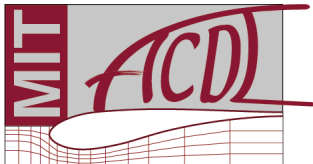


$$\text{P2} \Rightarrow$$
$$30,288 \text{ DOF}$$
$$|C_d - C_{d,\text{exact}}| = 1.2 * 10^{-5}$$



- Turbulence modeling (Todd Oliver)
- Shocks (Garrett Barter)
- Adaptation (Mike Park and Chris Fidkowski)
- Optimization (James Lu)
- Unsteady + Frequency Domain Solver (Tan Bui)
- Axisymmetric Solver + Plasma Physics (Shannon Cheng)
- Explicit solver (Pete Whitney)
- Hypersonic Flow (Doug Quattrochi)

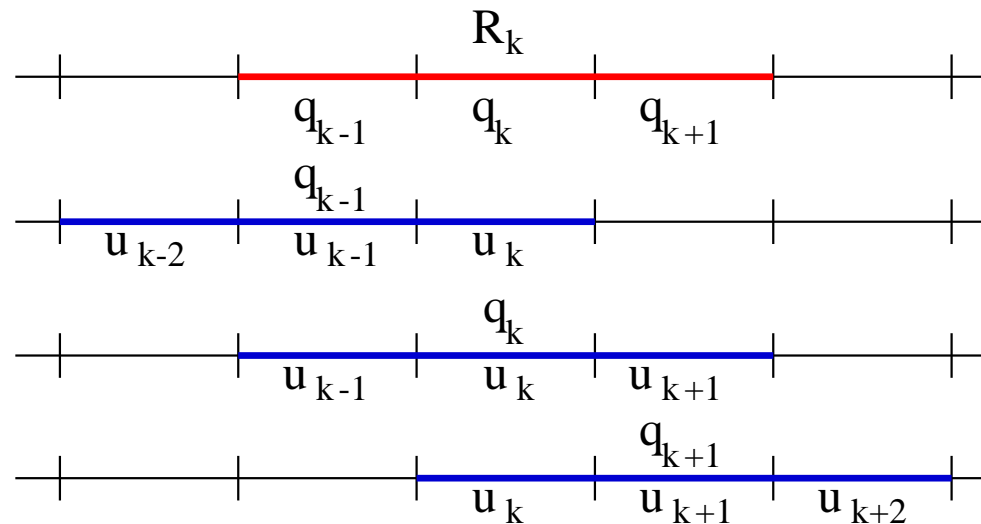
Questions?



- No upwinding mechanism \Rightarrow choose central fluxes

$$\hat{u} = \frac{1}{2}(u_h^L + u_h^R); \quad \hat{q} = \frac{1}{2}(q_h^L + q_h^R)$$

- Sub-optimal order of accuracy for odd p
- Stencil no longer compact



- Define jump, $[[\cdot]]$, and average, $\{\cdot\}$, operators:

$$[[s]] = s^L - s^R \quad \text{and} \quad \{s\} = \frac{1}{2}(s^L + s^R)$$

- Central fluxes become

$$\hat{u} = \{u_h\}; \quad \hat{q} = \{(u_h)_x\} - \{\delta\}$$

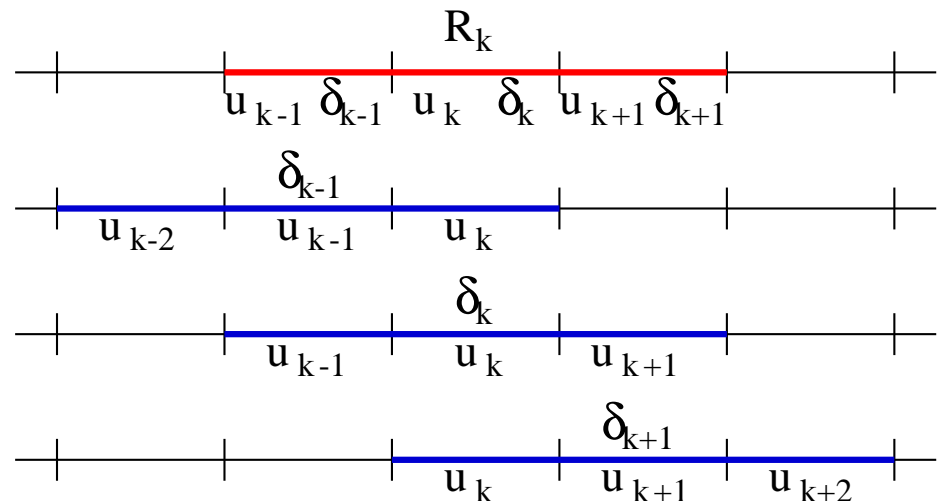
- δ given by following problem: Find $\delta \in \mathcal{V}_h^p$ such that $\forall \tau_h \in \mathcal{V}_h^p$,

$$\sum_{\kappa \in T_h} \int_{\kappa} \tau_h \delta dx = \sum_n \left[[[u_h]] \{\tau_h\} \right]$$

- BR1 becomes: Find $u_h \in \mathcal{V}_h^p$ and such that $\forall v_h \in \mathcal{V}_h^p$,

$$\sum_{\kappa \in T_h} \int_{\kappa} (v_h)_x (u_h)_x dx - \sum_n \left[\llbracket u_h \rrbracket \{ (v_h)_x \} + \llbracket v_h \rrbracket (\{ (u_h)_x \} - \{ \delta \}) \right] = \sum_{\kappa \in T_h} \int_{\kappa} v_h f dx$$

- Stencil extended by δ dependence on u_h



- Goal: Eliminate extended stencil
- Approach: Modify auxiliary variable, δ , previously defined by:

$$\sum_{\kappa \in T_h} \int_{\kappa} \tau_h \delta dx = \sum_n \left[\llbracket u_h \rrbracket \{ \tau_h \} \right]$$

- New variable, δ_f , given by: Find $\delta_f \in \mathcal{V}_h^p$ such that $\forall \tau_h \in \mathcal{V}_h^p$,

$$\int_{\kappa^{L/R}} \tau_h \delta_f^{L/R} dx = \left[\llbracket u_h \rrbracket \{ \tau_h \}^{L/R} \right]_{n_f}$$

- New fluxes have same form as before

$$\hat{u} = \{ u_h \}; \quad \hat{q} = \{ (u_h)_x \} - \eta_f \{ \delta_f \}$$

- Replacing $\{\delta\}$ in BR1 by $\eta_f\{\delta_f\}$ gives BR2
- For proper choice of η_f , can prove optimal order of accuracy
- Stencil is compact

