Computational Methods for the Euler Equations

Before discussing the Euler Equations and computational methods for them, let's look at what we've learned so far:

Assumptions/Flow type 2-D, Incompressible, Irrotational Inviscid
3-D, Incompressible, Irrotational Inviscid, Small disturbance
3-D, Subsonic compressible, Irrotational, Inviscid, Small disturbance

Euler CFD	3-D, Compressible (no M_{∞} limit), Rotational,
	Shocks, Inviscid

The only major effect missing after this week will be viscous-related effects.

2-D Euler Equations in Integral Form

Consider an arbitrary area (i.e. a fixed control volume) through which flows a compressible inviscid flow:



Note: Path around surface is taken so that interior of control volume is on left.

Conservation of Mass

$$\begin{pmatrix} rate \ of \ change \\ of \ mass \ in \ C \end{pmatrix} + \begin{pmatrix} rate \ of \ mass \ flow \\ out \ of \ C \end{pmatrix} = 0$$

$$Mass \ in \ C = \iint_C \ \rho dA \qquad \text{where} \ \ \rho = destiny \ of \ fluid$$

$$\Rightarrow rate \ of \ change = \frac{d}{dt} \iint_C \ \rho dA \qquad of \ mass \ in \ C$$

Now, the rate of mass flowing out of *C* : Mass flow out of $C = \oint_{\mathcal{H}} \rho \vec{u} \bullet \vec{n} dS$

 $\vec{u} =$ velocity vector

$$\Rightarrow \qquad \frac{d}{dt} \iint_{C} \rho dA + \oint_{\partial C} \rho \vec{u} \bullet \vec{n} dS = 0$$

Conservation of x-momentum

Recall that: total rate of change momentum = \sum forces

For x-momentum this gives:

 $\begin{pmatrix} rate of change of \\ x - momentum in C \end{pmatrix} + \begin{pmatrix} rate of x - momflow \\ out of C \end{pmatrix} = \Sigma$ Forces in x-direction $\frac{d}{dt} \iint_{C} \rho u dA + \oint_{\infty} \rho u \bar{u} \bullet \bar{n} dS = \Sigma$ Forces in x-direction Now looking closer at x-forces for an inviscid compressible flow we

Now, looking closer at x-forces, for an inviscid compressible flow we only have pressure (ignoring gravity). Normal to surface

Recall pressure acts normal to the surface





Gives x-direction

$$\frac{d}{dt} \iint_{C} \rho u dA + \oint_{\mathcal{K}} \rho u \vec{u} \bullet \vec{n} dS = -\oint_{\mathcal{K}} p \vec{n} \bullet \vec{i} \, dS$$

Conservation of y-momentum

This follows exactly the same as the x-momentum:

$$\frac{d}{dt} \iint\limits_{C} \rho v dA + \oint\limits_{\mathcal{K}} \rho v \vec{u} \bullet \vec{n} dS = -\oint\limits_{\mathcal{K}} p \vec{n} \bullet \vec{j} dS$$

Conservation of Energy

Recalling your thermodynamics:

$$\begin{pmatrix} total \ rate \ of \ change \\ of \ energy \ in \ C \end{pmatrix} = \begin{pmatrix} work \ done \ on \\ fluid \ in \ C \end{pmatrix} + \begin{pmatrix} heat \ added \\ to \ C \end{pmatrix}$$
For the Euler equations, we ignore the possibility of heat addition.

$$\begin{pmatrix} \text{total rate of change} \\ \text{of energy in } C \end{pmatrix} = \begin{pmatrix} \text{rate of change of} \\ \text{energy in } C \end{pmatrix} + \begin{pmatrix} \text{rate of energy} \\ \text{flow out of } C \end{pmatrix}$$
The total energy of the fluid is:

The total energy of the fluid is:

$$\rho E = \rho e + \frac{1}{2}\rho(u^2 + v^2)$$

Total Internal energy energy Kinetic energy

Note:
$$e = c_v T$$
 where $c_v \equiv$ specific heat at constant volume

Static temperature

So,

$$\begin{pmatrix} total \ rate \ of \ change \\ of \ energy \ in \ C \end{pmatrix} = \frac{d}{dt} \iint_{C} \rho E dA + \oint_{\mathcal{K}} \rho E \vec{u} \bullet \vec{n} dS$$

The work done on the fluid is through pressure forces and is equal to the pressure forces multiplied by (i.e. acting in) the velocity direction:

$$(work) = \oint_{\infty} (-p\bar{n}) \bullet \bar{u}dS$$

Pressure force

$$\Rightarrow \quad \frac{d}{dt} \iint_{C} \rho E dA + \oint_{\mathcal{K}} \rho E \vec{u} \bullet \vec{n} dS = -\oint_{\mathcal{K}} p \vec{n} \bullet \vec{u} dS$$

Summary of 2-D Euler Equations

$$\frac{d}{dt} \iint_{C} \rho dA + \oint_{\mathcal{X}} \rho \vec{u} \cdot \vec{n} dS = 0$$

$$\frac{d}{dt} \iint_{C} \rho u dA + \oint_{\mathcal{X}} \rho u \vec{u} \cdot \vec{n} dS = -\oint_{\mathcal{X}} p \vec{n} \cdot \vec{i} dS$$

$$\frac{d}{dt} \iint_{C} \rho v dA + \oint_{\mathcal{X}} \rho v \vec{u} \cdot \vec{n} dS = -\oint_{\mathcal{X}} p \vec{n} \cdot \vec{j} dS$$

$$\frac{d}{dt} \iint_{C} \rho E dA + \oint_{\mathcal{X}} \rho E \vec{u} \cdot \vec{n} dS = -\oint_{\mathcal{X}} p \vec{n} \cdot \vec{n} dS$$

These are often written very compactly as:

$$\frac{d}{dt}\iint_{c} U dA + \oint_{sc} \left(F\vec{i} + G\vec{j} \right) \bullet \vec{n} ds = 0$$



$$G \equiv \begin{bmatrix} \rho v \\ \rho v u \\ \rho v^{2} + p \\ \rho v H \end{bmatrix}$$

Conservative state vector

Flux vector for x-direction

Flux vector for y-direction

$$H \equiv total \ enthalpy \equiv E + \frac{p}{\rho}$$

Ideal gas: $p = \rho RT = (\gamma - 1) \left[\rho E - \frac{1}{2} \rho (u^2 + v^2) \right]$
A Finite Volume Scheme for the 2-D Euler Eqns.

Here's the basic idea:

(1) Divide up (i.e. discretize) the domain into simple geometric shapes (triangles and quads)



Cell 0 is surrounded by cells 1, 2, & 3. i.e. cell 0 has 3 neighbors: cell 1, 2, & 3.

Nearest neighbors

- (2) Decide how to place the unknowns in the grid.
 - (a) <u>Cell-centered</u>: cell-average values of the conservative state vector are stored for each cell.
 - (b) <u>Node-based</u>: point values of the conservative state vector are stored at each node.

The debate still rages about which of these options is best. We will look at cellcentered schemes because these are easiest (although not necessarily the best). Also, they are very widely used in the aerospace industry.

(3) Approximate the 2-D integral Euler equation on the grid to determine the chosen unknowns.





Refined Mesh - 7506 Nodes





Refined Mesh - 7506 Nodes

Let's look in detail at step (3):



Cells: 0,1, 2, 3 Nodes: a, b, c, d, e, f

Cell-average unknowns:

$$U_{0} = \begin{bmatrix} \rho_{0} \\ (\rho u)_{0} \\ (\rho v)_{0} \\ (\rho E)_{0} \end{bmatrix} \qquad \qquad U_{1} = \begin{bmatrix} \rho_{1} \\ (\rho u)_{1} \\ (\rho v)_{1} \\ (\rho E)_{1} \end{bmatrix} \qquad \qquad U_{2} = \dots \dots \dots \dots$$

Specifically, we define U_0 as:

 $U_{0} \equiv \frac{1}{A_{0}} \iint_{C_{0}} U dA \qquad \text{where } \begin{cases} C_{0} \equiv cell \ 0 \\ A_{0} \equiv area \ of \ cell \ 0 \end{cases}$

Now, we apply conservation eqns:

$$\frac{d}{dt} \iint_{\mathcal{C}_0} U dA + \oint_{\mathcal{K}_0} \left(F \vec{i} + G \vec{j} \right) \bullet \vec{n} dS = 0$$

The time-derivative term can be simplified a little:

$$\frac{d}{dt} \iint_{C_0} U dA = A_0 \frac{dU_0}{dt}$$

The surface flux integral can also be simplified a little:

$$\begin{split} \oint_{\partial C_0} \left(F\vec{i} + G\vec{j} \right) \bullet \vec{n} ds &= \int_a^b \left(F\vec{i} + G\vec{j} \right) \bullet \vec{n} dS \\ &+ \int_b^c \left(F\vec{i} + G\vec{j} \right) \bullet \vec{n} dS \\ &+ \int_c^a \left(F\vec{i} + G\vec{j} \right) \bullet \vec{n} dS \end{split}$$

Combining these expressions:

$$A_{0} \frac{dU_{0}}{dt} + \int_{a}^{b} \left(F\vec{i} + G\vec{j}\right) \bullet \vec{n}dS + \int_{b}^{c} \left(F\vec{i} + G\vec{j}\right) \bullet \vec{n}dS$$

+ $\int_{c}^{a} \left(F\vec{i} + G\vec{j}\right) \bullet \vec{n}dS = 0$ No approximations so far!

Now, we make some approximations. Let's look at the surface integral from $a \rightarrow b$: b



The normal can easily be calculated since the <u>face</u> is a straight line between nodes a & b. Recall, the unknowns are stored at all centers. So, what would be a logical approximation for :

$$\int_{a}^{b} (F\vec{i} + G\vec{j}) \bullet \vec{n}_{ab} dS = ???$$
Option #1=
Option #2=
Note: Option #1 \neq option #2 in general.

There is very little difference in practice between these options. Let's stick with:

$$\begin{aligned} \mathfrak{Z}_{ab} &= \int_{a}^{b} \left(F\vec{i} + G\vec{j} \right) \bullet \vec{n}_{ab} dS = \left[\frac{1}{2} (F_{0} + F_{1})\vec{i} + \frac{1}{2} (G_{0} + G_{1})\vec{j} \right] \bullet \vec{n}_{ab} \Delta S_{ab} \\ \mathfrak{Z}_{bc} &= \int_{b}^{c} \left(F\vec{i} + G\vec{j} \right) \bullet \vec{n}_{bc} dS = \left[\frac{1}{2} (F_{0} + F_{2})\vec{i} + \frac{1}{2} (G_{0} + G_{2})\vec{j} \right] \bullet \vec{n}_{bc} \Delta S_{bc} \\ \mathfrak{Z}_{ca} &= \int_{c}^{a} \left(F\vec{i} + G\vec{j} \right) \bullet n_{ca} dS = \left[\frac{1}{2} (F_{0} + F_{3})\vec{i} + \frac{1}{2} (G_{0} + G_{3})\vec{j} \right] \bullet \vec{n}_{ca} \Delta s_{ca} \end{aligned}$$

Where

$$F_0 \equiv F(U_0) \qquad G_0 \equiv G(U_0)$$

$$F_1 \equiv F(U_1) \qquad G_1 \equiv G(U_1)$$

$$F_2 \equiv F(U_2) \qquad G_2 \equiv G(U_2)$$

$$F_3 \equiv F(U_3) \qquad G_3 \equiv G(U_3)$$

Finally, we have to approximate $A_0 \frac{dU_0}{dt}$ somehow. The simplest approach is forward Euler:

$$A_0 \frac{dU_0}{dt} + \mathfrak{I}_{ab} + \mathfrak{I}_{bc} + \mathfrak{I}_{ca} = 0$$

$$A_0 \frac{U_0^{n+1} - U_0^n}{\Delta t} + \mathfrak{I}_{ab}^n + \mathfrak{I}_{bc}^n + \mathfrak{I}_{ca}^n = 0$$

Where $U_o^n \equiv U_o(t^n)$ and $t^n \equiv n\Delta t$, $n \equiv iteration$

And \mathfrak{I}_{ab}^n etc. are defined as:

$$\begin{split} \mathfrak{T}_{ab}^{n} &= \left[\frac{1}{2} \Big(F_{0}^{n} + F_{1}^{n} \Big) \overline{i} + \frac{1}{2} \Big(G_{0}^{n} + G_{1}^{n} \Big) \overline{j} \right] \bullet \overline{n}_{ab} \Delta S_{ab} \\ F_{0}^{n} &= F \Big(U_{0}^{n} \Big) \qquad etc. \\ F_{1}^{n} &= F \Big(U_{1}^{n} \Big) \end{split}$$

For steady solution, basic procedure is to make a guess of U at t = 0 and then iterate until the solution no longer changes. This is called time marching.

Question

What assumptions have we made in developing our 2-D Euler Equation Finite Volume Method?