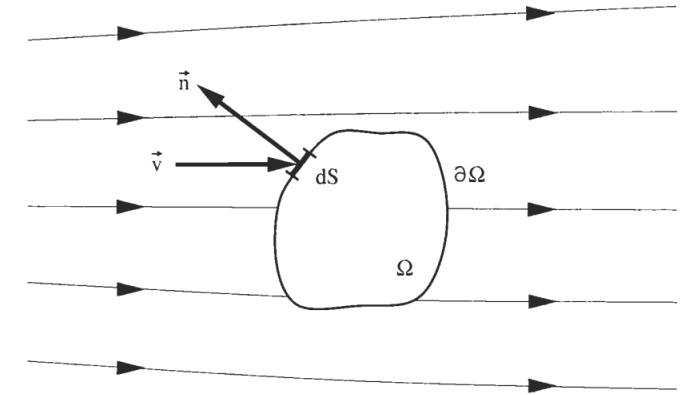


## Part-II: Compressible FVM methods in OF

## The Conservation Equations in Integral Form



$$\frac{\partial}{\partial t} \int_{\Omega} \vec{W} d\Omega + \oint_{\partial\Omega} (\vec{F}_c - \vec{F}_v) dS = \int_{\Omega} \vec{Q} d\Omega.$$

$$\vec{W} = \begin{bmatrix} \rho \\ \rho u \\ \rho v \\ \rho w \\ \rho E \end{bmatrix}.$$

$$\vec{F}_c = \begin{bmatrix} \rho V \\ \rho u V + n_x p \\ \rho v V + n_y p \\ \rho w V + n_z p \\ \rho H V \end{bmatrix}$$

$$V \equiv \vec{v} \cdot \vec{n} = n_x u + n_y v + n_z w$$

$$\vec{F}_v = \begin{bmatrix} 0 \\ n_x \tau_{xx} + n_y \tau_{xy} + n_z \tau_{xz} \\ n_x \tau_{yx} + n_y \tau_{yy} + n_z \tau_{yz} \\ n_x \tau_{zx} + n_y \tau_{zy} + n_z \tau_{zz} \\ n_x \Theta_x + n_y \Theta_y + n_z \Theta_z \end{bmatrix},$$

$$\Theta_x = u \tau_{xx} + v \tau_{xy} + w \tau_{xz} + k \frac{\partial T}{\partial x}$$

$$\Theta_y = u \tau_{yx} + v \tau_{yy} + w \tau_{yz} + k \frac{\partial T}{\partial y}$$

$$\Theta_z = u \tau_{zx} + v \tau_{zy} + w \tau_{zz} + k \frac{\partial T}{\partial z}$$

$$\vec{Q} = \begin{bmatrix} 0 \\ \rho f_{e,x} \\ \rho f_{e,y} \\ \rho f_{e,z} \\ \rho \vec{f}_e \cdot \vec{v} + \dot{q}_h \end{bmatrix}.$$

*Integration of time derivative*

$$\frac{\partial}{\partial t} \int_{\Omega} \vec{W} d\Omega + \oint_{\partial\Omega} (\vec{F}_c - \vec{F}_v) dS = \int_{\Omega} \vec{Q} d\Omega.$$

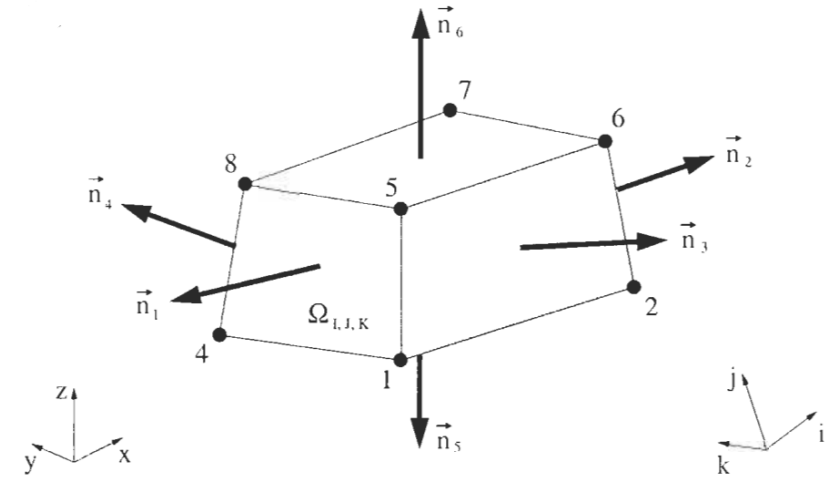
*Step 2*

$$\frac{\partial}{\partial t} \int_{\Omega} \vec{W} d\Omega = \Omega \frac{\partial \vec{W}}{\partial t}.$$

*Step 1*

$$\frac{\partial \vec{W}}{\partial t} = -\frac{1}{\Omega} \left[ \oint_{\partial\Omega} (\vec{F}_c - \vec{F}_v) dS - \int_{\Omega} \vec{Q} d\Omega \right].$$

*Step 3*

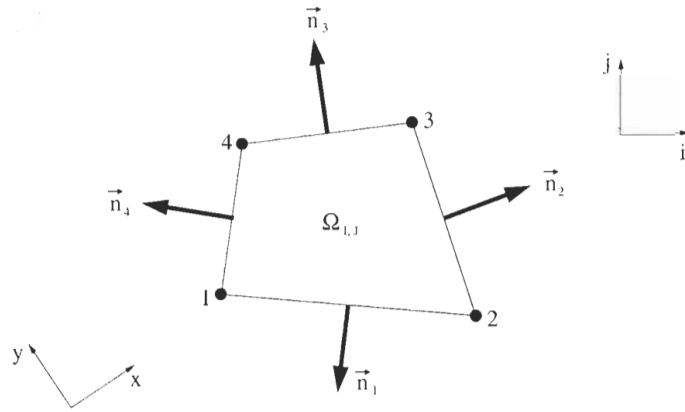


*Step 4*

$$\frac{d\vec{W}_{I,J,K}}{dt} = -\frac{1}{\Omega_{I,J,K}} \left[ \sum_{m=1}^{N_F} (\vec{F}_c - \vec{F}_v)_m \Delta S_m - (\vec{Q}\Omega)_{I,J,K} \right]$$

$$\frac{d\vec{W}_{I,J,K}}{dt} = -\frac{1}{\Omega_{I,J,K}} \vec{R}_{I,J,K}$$

## Geometrical Quantities of a Control Volume



*Volume of cell*

$$\Omega_{I,J} = \frac{b}{2} [(x_1 - x_3)(y_2 - y_4) + (x_4 - x_2)(y_1 - y_3)]$$

*Face area of cell*

$$\vec{S}_m = \begin{bmatrix} S_{x,m} \\ S_{y,m} \end{bmatrix} = \vec{n}_m \Delta S_m$$

$$\vec{S}_1 = b \begin{bmatrix} y_2 - y_1 \\ x_1 - x_2 \end{bmatrix}$$

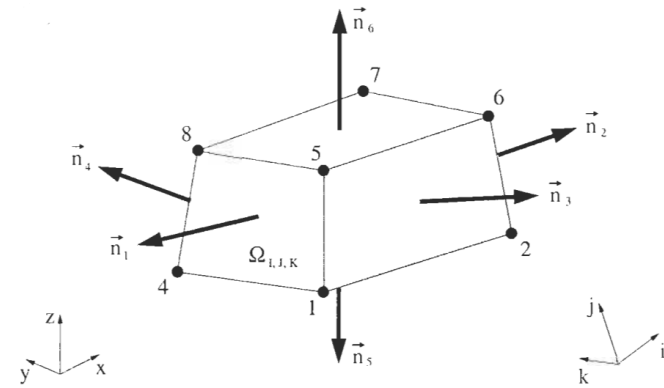
$$\vec{S}_2 = b \begin{bmatrix} y_3 - y_2 \\ x_2 - x_3 \end{bmatrix}$$

$$\vec{S}_3 = b \begin{bmatrix} y_4 - y_3 \\ x_3 - x_4 \end{bmatrix}$$

$$\vec{S}_4 = b \begin{bmatrix} y_1 - y_4 \\ x_4 - x_1 \end{bmatrix}$$

$$\vec{n}_m = \frac{\vec{S}_m}{\Delta S_m}$$

$$\Delta S_m = |\vec{S}_m| = \sqrt{S_{x,m}^2 + S_{y,m}^2}$$



*Face area of cell*

$$\Delta x_A = x_8 - x_1, \quad \Delta x_B = x_5 - x_4,$$

$$\Delta y_A = y_8 - y_1, \quad \Delta y_B = y_5 - y_4,$$

$$\Delta z_A = z_8 - z_1, \quad \Delta z_B = z_5 - z_4$$

$$\vec{S}_1 = \frac{1}{2} \begin{bmatrix} \Delta y_A \Delta z_B - \Delta z_A \Delta y_B \\ \Delta z_A \Delta x_B - \Delta x_A \Delta z_B \\ \Delta x_A \Delta y_B - \Delta y_A \Delta x_B \end{bmatrix}$$

$$\Delta S_m = \sqrt{S_{x,m}^2 + S_{y,m}^2 + S_{z,m}^2}$$

*Volume of cell*

$$\Omega_{I,J,K} = \frac{1}{3} \sum_{m=1}^{m=6} (\vec{r}_{\text{mid}} \cdot \vec{S})_m$$

$$\vec{r}_{\text{mid},1} = \frac{1}{4} (\vec{r}_1 + \vec{r}_5 + \vec{r}_8 + \vec{r}_4)$$

## Discretization of Convective Terms

- *Various methods available for discretization are*

- ☐ *Central,*

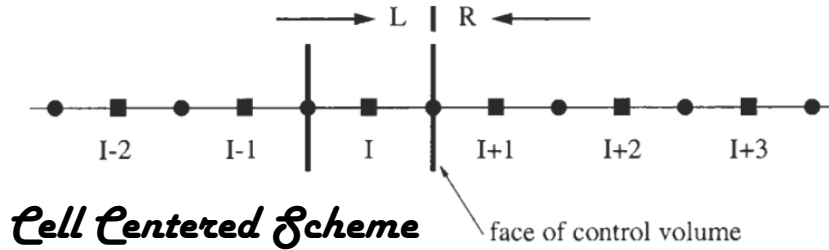
- ☐ *Flux-Vector Splitting*

- ☐ *Flux-difference Splitting*

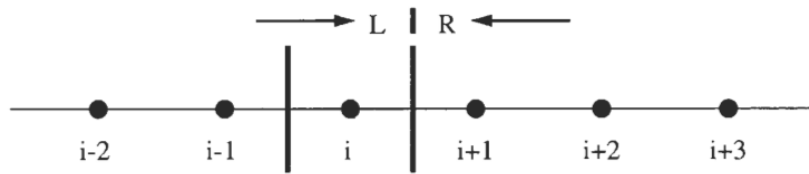
- ☐ *Total Variation Diminishing (TVD) and*

- ☐ *Fluctuation splitting*

## Left & Right States and Stencil



*Cell Centered Scheme*



*Cell Vertex Scheme*

### *Central Scheme*

- Uses same number of cells left and right of interface
- Flux is "centered" by symmetric cell distribution across cell interface

### *Upwinding*

- Uses cells based on information propagation direction based on characteristics
- Asymmetric usage of cells across cell interface

$$U_R = U_{I+1} - \frac{\epsilon}{4} [(1 + \hat{\kappa})\Delta_- + (1 - \hat{\kappa})\Delta_+] U_{I+1}$$

$$U_L = U_I + \frac{\epsilon}{4} [(1 + \hat{\kappa})\Delta_+ + (1 - \hat{\kappa})\Delta_-] U_I.$$

$$\Delta_+ U_I = U_{I+1} - U_I$$

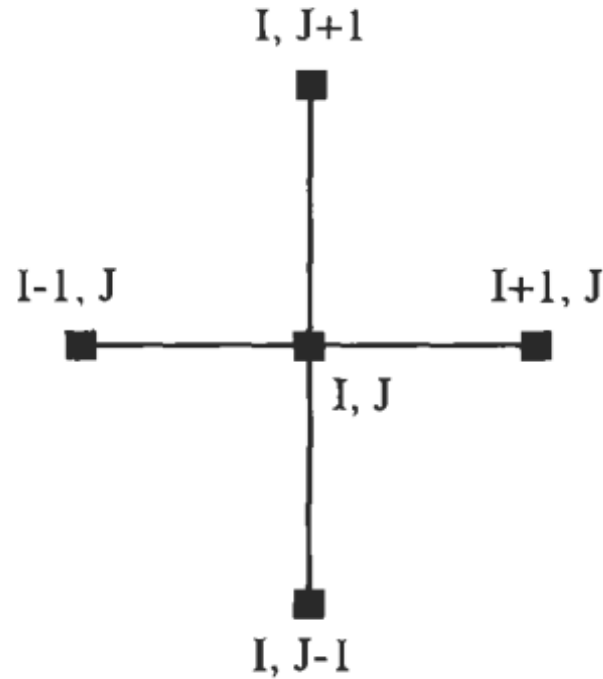
$$\Delta_- U_I = U_I - U_{I-1}.$$

Based on values of  $\epsilon$  and  $\hat{\kappa}$  several upwind schemes can be defined such as

- First order upwind
- Second order upwind
- Upwind biased linear interpolation
- Second order central Scheme

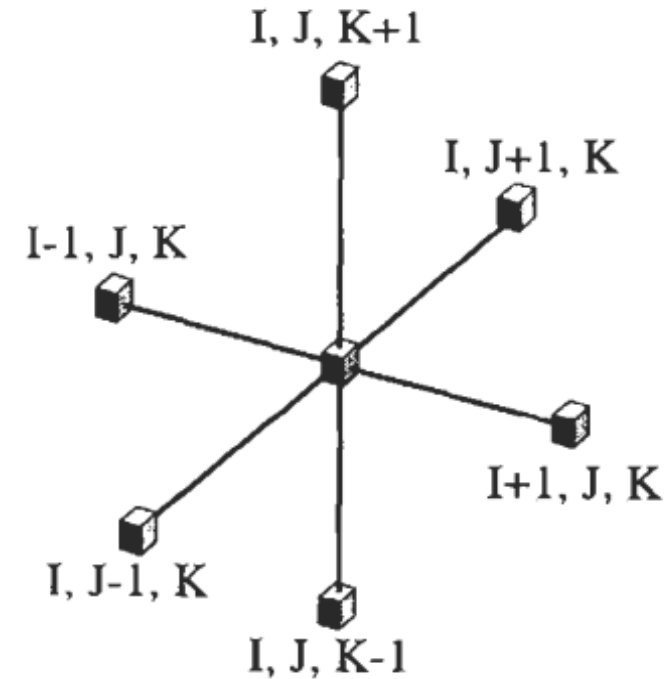
*Union of cells involved in computation of derived quantities is Stencil*

(a)



$2D$

(b)



$3D$

$(I, J) \quad (I + 1, J) \quad (I, J + 1) \quad (I - 1, J) \quad (I, J - 1)$

$(I, J, K) \quad (I + 1, J, K) \quad (I, J + 1, K) \quad (I - 1, J, K)$   
 $(I, J - 1, K) \quad (I, J, K - 1) \quad (I, J, K + 1)$

## Flux Vector Splitting Scheme

### 1. Van Leer's Scheme

$$\vec{F}_c = \vec{F}_c^+ + \vec{F}_c^- ,$$

$$(M_n)_{I+1/2} = \left( \frac{V}{c} \right)_{I+1/2} ,$$

$$(M_n)_{I+1/2} = M_L^+ + M_R^- ,$$

$$M_L^+ = \begin{cases} M_L & \text{if } M_L \geq +1 \\ \frac{1}{4}(M_L + 1)^2 & \text{if } |M_L| < 1 \\ 0 & \text{if } M_L \leq -1 \end{cases} ,$$

$$M_R^- = \begin{cases} 0 & \text{if } M_R \geq +1 \\ \frac{1}{4}(M_R - 1)^2 & \text{if } |M_R| < 1 \\ M_R & \text{if } M_R \leq -1 \end{cases} .$$

$$M_L = \frac{V_L}{c_L} , \quad M_R = \frac{V_R}{c_R} .$$

$$\vec{F}_c^\pm = \begin{bmatrix} f_{\text{mass}}^\pm \\ f_{\text{mass}}^\pm [n_x(-V \pm 2c)/\gamma + u] \\ f_{\text{mass}}^\pm [n_y(-V \pm 2c)/\gamma + v] \\ f_{\text{mass}}^\pm [n_z(-V \pm 2c)/\gamma + w] \\ f_{\text{energy}}^\pm \end{bmatrix} .$$

$$|M_n| < 1$$

$$f_{\text{mass}}^+ = +\rho_L c_L \frac{(M_L + 1)^2}{4}$$

$$f_{\text{mass}}^- = -\rho_R c_R \frac{(M_R - 1)^2}{4}$$

$$f_{\text{energy}}^\pm = f_{\text{mass}}^\pm \left\{ \frac{[(\gamma - 1)V \pm 2c]^2}{2(\gamma^2 - 1)} + \frac{u^2 + v^2 + w^2 - V^2}{2} \right\}_{L/R} .$$

$$|M_n| \geq 1$$

$$\vec{F}_c^+ = \vec{F}_c \quad \vec{F}_c^- = 0 \quad \text{if } M_n \geq +1$$

$$\vec{F}_c^+ = 0 \quad \vec{F}_c^- = \vec{F}_c \quad \text{if } M_n \leq -1 .$$



## 2. Advection Upstream Splitting Method (AUSM)

$$\vec{F}_c = V \begin{bmatrix} \rho \\ \rho u \\ \rho v \\ \rho w \\ \rho H \end{bmatrix} + \begin{bmatrix} 0 \\ n_x p \\ n_y p \\ n_z p \\ 0 \end{bmatrix}.$$

$$(\bullet)_{L/R} = \begin{cases} (\bullet)_L & \text{if } M_{I+1/2} \geq 0 \\ (\bullet)_R & \text{otherwise} \end{cases}$$

*Convective flux splitting same as Van Leer*

$$p_{I+1/2} = p_L^+ + p_R^-$$

$$p_L^+ = \begin{cases} p_L & \text{if } M_L \geq +1 \\ \frac{p_L}{4} (M_L + 1)^2 (2 - M_L) & \text{if } |M_L| < 1 \\ 0 & \text{if } M_L \leq -1 \end{cases},$$

$$p_R^- = \begin{cases} 0 & \text{if } M_R \geq +1 \\ \frac{p_R}{4} (M_R - 1)^2 (2 + M_R) & \text{if } |M_R| < 1 \\ p_R & \text{if } M_R \leq -1 \end{cases}.$$

## *Some other convective flux discretization schemes continued.....*

Various Computational Schemes for Euler Equations		
Central Schemes	First Order Upwind Schemes	Second Order Upwind Schemes
<b>1. Combined Space-Time Integration</b> <b>(a) Explicit Schemes</b> Lax-Friedrichs – First order (1954) Lax-Wendroff – Second order (1960) <b>(b) Two-Step Explicit Schemes</b> Richtmyer and Morton (1967) MacCormack (1969) LeRat and Peyret (1974) <b>(c) Implicit Schemes</b> MacCormack (1981) Casier, Deconinck, Hirsch (1983) LeRat (1979, 1983) <b>2. Separate Space-Time Integration</b> <b>(a) Implicit Schemes</b> Briley and McDonald (1975) Beam and Warming (1976) <b>(b) Explicit Schemes (Multistage Runge-Kutta)</b> Jameson, Schmidt, Turkel (1981)	<b>1. Flux Vector Splitting</b> Courant, Isaacson, and Reeves (1952) Moretti (1979) Steger and Warming (1981) VanLeer (1982) <b>2. Godunov Methods-Riemann Solvers</b> <b>(a) Exact Riemann Solvers</b> Godunov (1959) – First order VanLeer (1979) – Second order Woodward and Colella (1984) Ben-Artzi and Falcovitz (1984) <b>(b) Approximate Riemann Solvers</b> Roe (1981) Enquist and Osher (1980) Osher (1982) Harten, Lax, Van Leer (1983)	<b>1. Extrapolation</b> <b>(a) Variable Extrapolation (MUSCL)</b> Van Leer (1979) <b>(b) Flux Extrapolation</b> Van Leer (1979) <b>2. Explicit TVD Upwind</b> VanLeer (1974) Harten (1983) Osher (1984) Osher and Chakravarthy (1984) <b>3. Implicit TVD Upwind</b> Yee (1986) <b>4. Central TVD Implicit or Explicit</b> Davis (1984) Roe (1985) Yee (1985) <b>5. Essentially Nonoscillatory Scheme</b> Harten and Osher (1987) <b>6. Flux Corrected Transport</b> Boris and Book (1973)

## Implementation of Central Scheme in OpenFoam: **rhoCentralFoam**

### Features of rhoCentralFoam

- Finite Volume, density based solver.
- Non-staggered/collocated field variables.
- Central/Central-upwind scheme.
- Applicable on polyhedral Mesh.
- Suitable for high speed viscous flows.

## **Some of the Important Solution Methodology for High Speed Flows**

- Notable methods for producing accurate non-oscillatory solutions are (convective terms discretization of governing equations)
- Monotone upstream-centred scheme for conservation laws
- Piece wise parabolic method (PPM)
- Essentially non-oscillatory scheme (ENO)
- Weighted ENO (WENO)

The above methods are complicated as they involve Riemann solvers, characteristic decomposition and Jacobian evaluation.

They are complex to implement !!

- To avoid the difficulty of implementation and to reduce cost of computation, **Central schemes** were devised which was first proposed by Nessyahu and Tadmor as second order accurate version of Lax-Friedrichs scheme.
- The more advanced versions of central scheme widely used to solve industrial problems are **Kurganov-Tadmor** (KT) and **Kurganov-Noelle-Petrova** (KNP) schemes
- These schemes are implemented in **OpenFOAM** package, with “**rhoCentralFoam**” as the solver name

## Governing Equations : Differential Form

*Continuity Equation*

$$\frac{\partial \rho}{\partial t} + \nabla \cdot [\mathbf{u}\rho] = 0$$

*Momentum Equation*

$$\frac{\partial(\rho \mathbf{u})}{\partial t} + \nabla \cdot [\mathbf{u}(\rho \mathbf{u})] + \nabla p + \nabla \cdot \mathbf{T} = 0$$

*Conservation of Total Energy*

$$\frac{\partial(\rho E)}{\partial t} + \nabla \cdot [\mathbf{u}(\rho E)] + \nabla \cdot (\mathbf{u}p) + \nabla \cdot (\mathbf{T}\mathbf{u}) + \nabla \cdot \mathbf{j} = 0$$



- Here  $\rho$  is the mass density  $\mathbf{u}$  is the fluid velocity,  $p$  is the pressure,  
Total energy density  $E = e + 0.5 |\mathbf{u}|^2$  with 'e' as specific internal energy  
 $\mathbf{j}$  as the diffusion heat flux vector,  $\mathbf{T}$  is the viscous shear stress tensor (defined positive in compression)

$$\mathbf{T} = -2\mu \text{dev}(\mathbf{D})$$

$\mu$  is the dynamic viscosity (Sutherland's law or constant)

$$\mathbf{D} = \frac{1}{2} [\nabla \mathbf{u} + \nabla \mathbf{u}^T] \text{ (Deformation gradient tensor)}$$

$$\text{Deviatoric component } \text{dev}(\mathbf{D}) = \mathbf{D} - \left(\frac{1}{3}\right) \text{tr}(\mathbf{D}) \mathbf{I} \quad \mathbf{I} \text{ is the unit tensor}$$

The diffusion heat flux vector defined by

$$\mathbf{j} = -k \nabla T \text{ (Fourier's Law)}$$

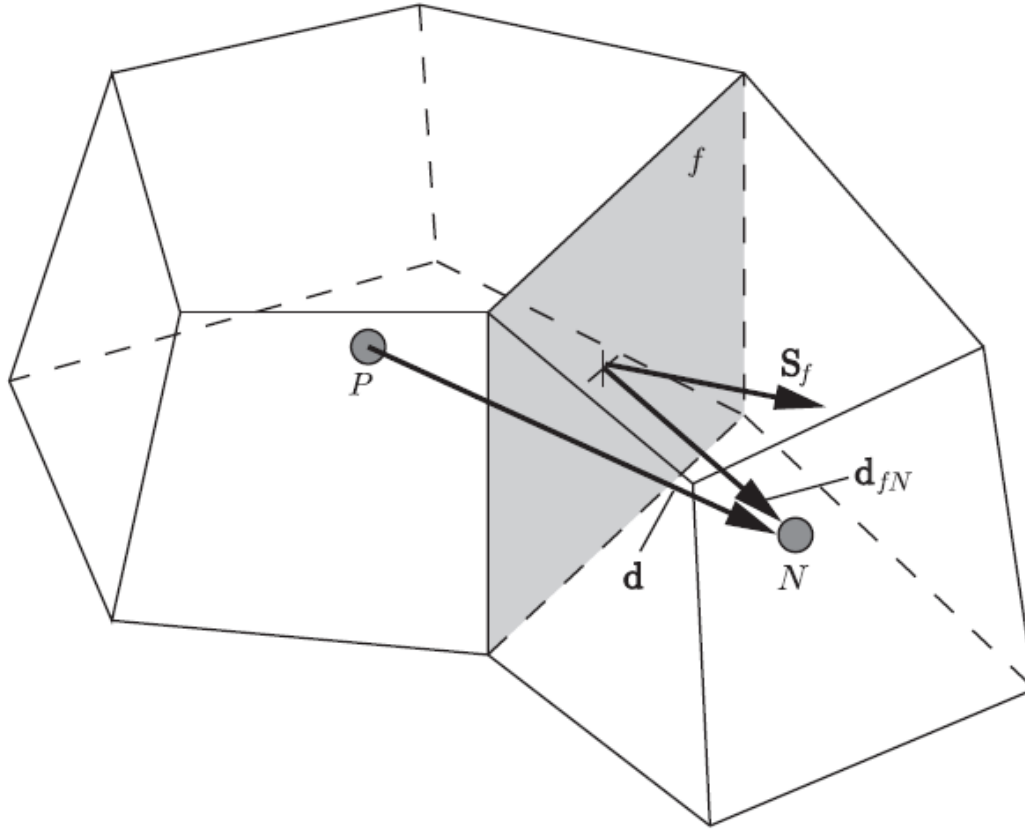
$k$  = thermal conductivity

$T$  = temperature

For calorically perfect gas,  $p = \rho R T$  and  $e = C_v T = (\gamma - 1) R T$

$\gamma = C_p / C_v$ ,  $R$  is characteristic gas constant

## Computational Method



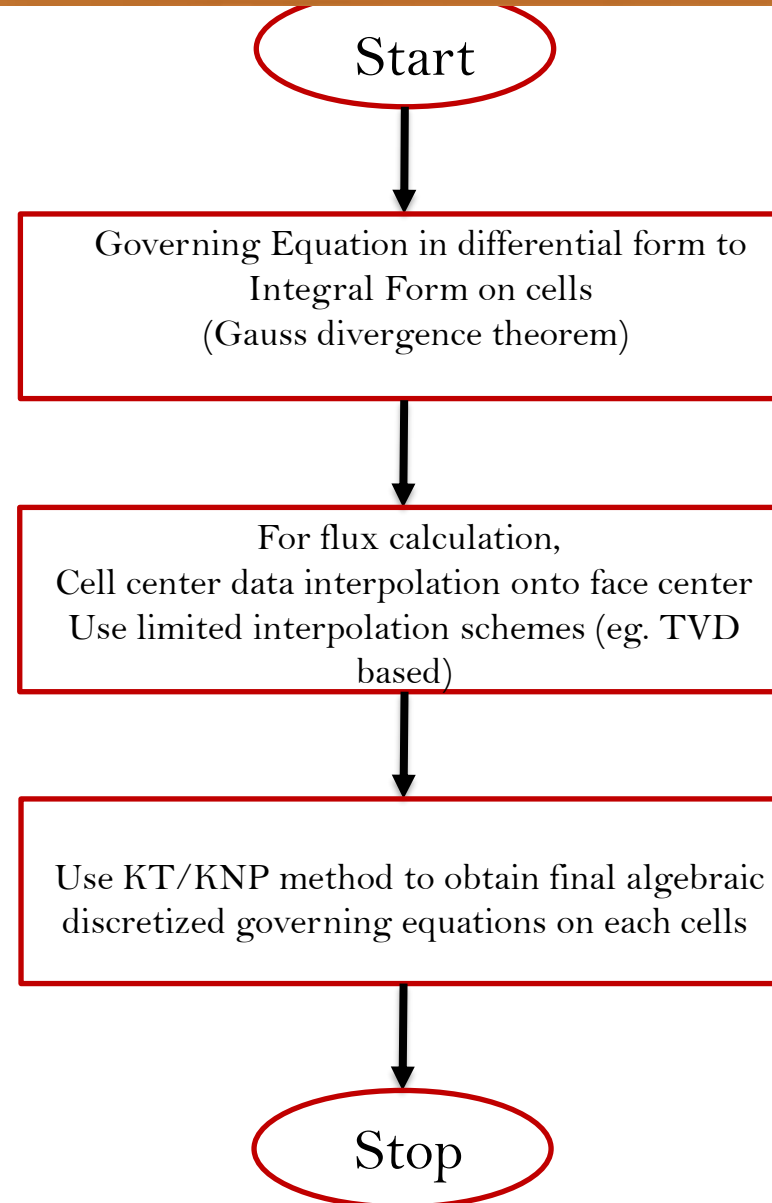
$d_{fN}$  vector connecting face centre and cell centre of neighbouring cell

$d$  Vector connecting cell centre  $\mathbf{P}$  and cell centre  $\mathbf{N}$

$\mathbf{S}_f$  face area vector (outwards is positive)

All variables are stored at cell centres ( $\mathbf{P}, \mathbf{N}, \dots$ )





*The KT/KNP convective flux discretization method shall be explained in upcoming slides.....*

***Procedure to obtain discretized algebraic governing equation on collocated finite volume grid***

## The Algebraic Discretized Equation

- The general variable  $\psi$  (tensor of any rank) can be convected in the flow field, the general term associated with convection in GE is:

$$\nabla \cdot (\mathbf{u}\psi)$$

Application of Gauss divergence theorem

$$\iiint \nabla \cdot (\mathbf{u}\psi) dV = \iint d\mathbf{S} \cdot [\mathbf{u}\psi] = \sum_f S_f \cdot [\mathbf{u}_f \psi_f] = \sum_f \phi_f \psi_f$$

Here,  $\phi_f$  is the volumetric flux

## Convective Terms

- The general variable  $\psi$  (tensor of any rank) can be convected in the flow field, the general term associated with convection in GE is:

$$\nabla \cdot (\mathbf{u}\psi)$$

Application of Gauss divergence theorem

$$\iiint \nabla \cdot (\mathbf{u}\psi) dV = \iint d\mathbf{S} \cdot [\mathbf{u}\psi] = \sum_f S_f \cdot [\mathbf{u}_f \psi_f] = \sum_f \phi_f \psi_f$$

Here,  $\phi_f$  is the volumetric flux

$$\alpha = \begin{cases} \frac{1}{2} & \text{for } KT \text{ method} \\ \frac{\lambda_{f+}}{\lambda_{f+} + \lambda_{f-}} & \text{for } KNP \text{ method} \end{cases}$$

$$\lambda_{f+} = \max(c_{f+}|\mathbf{S}_f| + \phi_{f+}, c_{f-}|\mathbf{S}_f| + \phi_{f-}, 0)$$

$$\lambda_{f-} = \max(c_{f+}|\mathbf{S}_f| - \phi_{f+}, c_{f-}|\mathbf{S}_f| - \phi_{f-}, 0)$$

$$\omega_f = \begin{cases} \alpha \max(\lambda_{f+}, \lambda_{f-}) & \text{for } KT \text{ method} \\ \alpha(1 - \alpha)(\lambda_{f+} + \lambda_{f-}) & \text{for } KNP \text{ method} \end{cases}$$

## Gradient Terms

$$\iiint \nabla(\psi) dV = \iint d\mathbf{S} \psi = \sum_f \mathbf{s}_f \psi_f$$

$$\sum_f \mathbf{s}_f \psi_f = \sum_f [\alpha \mathbf{s}_f \psi_{f+} + (1 - \alpha) \mathbf{s}_f \psi_{f-}]$$

## Laplacian Terms

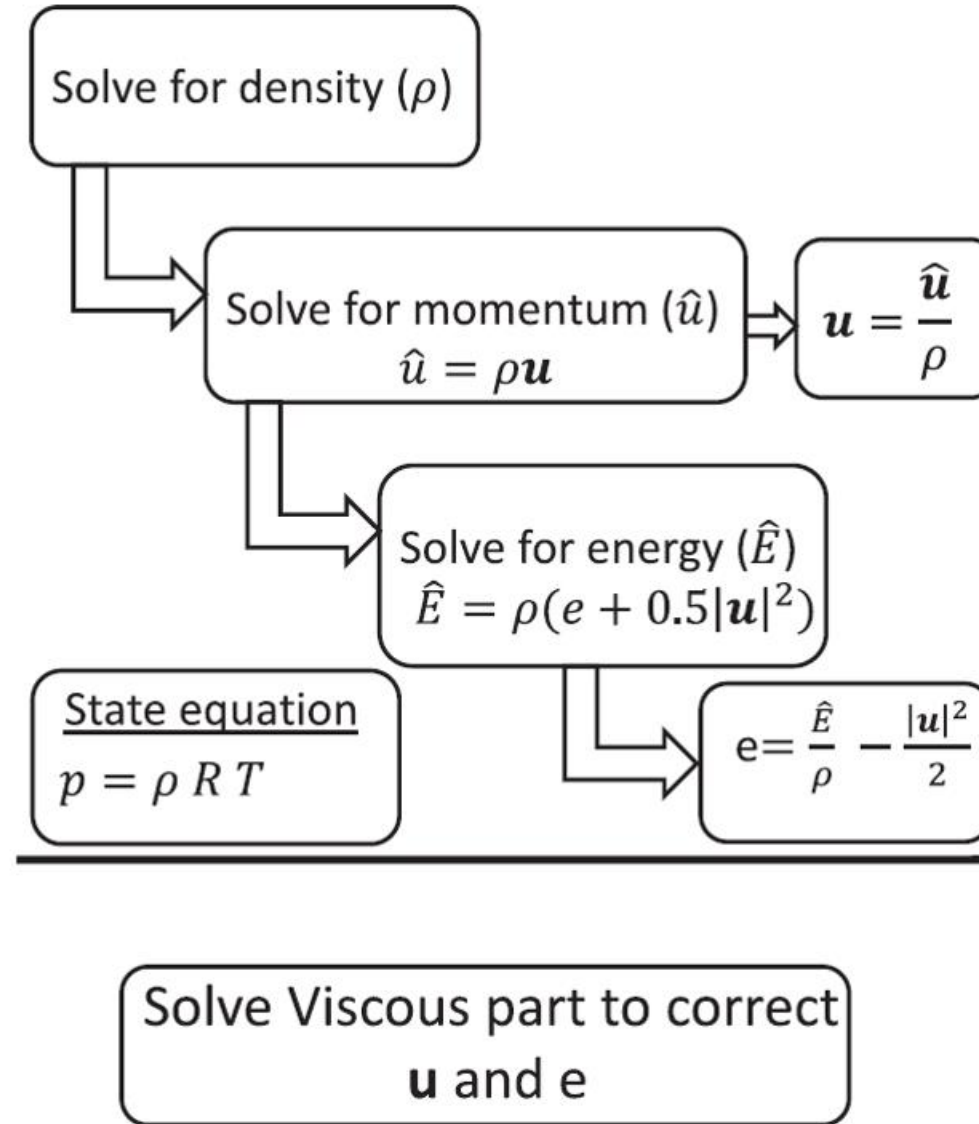
$$\iiint \nabla \cdot (\Gamma \nabla \psi) dV = \iint dS (\Gamma \nabla \psi) = \sum_f \Gamma_f \mathbf{S}_f \cdot (\nabla \psi)_f$$

For Non-orthogonal Mesh,

$$\mathbf{S}_f \cdot (\nabla \psi)_f = A(\psi_N - \psi_P) + \mathbf{a} \cdot (\nabla \psi)_f$$

$$A = |\mathbf{S}_f|^2 / \mathbf{S}_f \cdot \mathbf{d}$$

$$\mathbf{a} = \mathbf{S}_f - Ad$$



*Solution Algorithm of rhoCentralFoam*

## OpenFoam File : Implementation

```
surfaceScalarField rho_pos(interpolate(rho, pos));
surfaceScalarField rho_neg(interpolate(rho, neg));

surfaceVectorField rhoU_pos(interpolate(rhoU, pos, U.name()));
surfaceVectorField rhoU_neg(interpolate(rhoU, neg, U.name()));

volScalarField rPsi("rPsi", 1.0/psi);
surfaceScalarField rPsi_pos(interpolate(rPsi, pos, T.name()));
surfaceScalarField rPsi_neg(interpolate(rPsi, neg, T.name()));

surfaceScalarField e_pos(interpolate(e, pos, T.name()));
surfaceScalarField e_neg(interpolate(e, neg, T.name()));

surfaceVectorField U_pos("U_pos", rhoU_pos/rho_pos);
surfaceVectorField U_neg("U_neg", rhoU_neg/rho_neg);

surfaceScalarField p_pos("p_pos", rho_pos*rPsi_pos);
surfaceScalarField p_neg("p_neg", rho_neg*rPsi_neg);

surfaceScalarField phiv_pos("phiv_pos", U_pos & mesh.Sf());
// Note: extracted out the orientation so becomes unoriented
phiv_pos.setOriented(false);
surfaceScalarField phiv_neg("phiv_neg", U_neg & mesh.Sf());
phiv_neg.setOriented(false);
```

```
volScalarField c("c", sqrt(thermo.Cp()/thermo.Cv()*rPsi));
surfaceScalarField cSf_pos
(
    "cSf_pos",
    interpolate(c, pos, T.name())*mesh.magSf()
);

surfaceScalarField cSf_neg
(
    "cSf_neg",
    interpolate(c, neg, T.name())*mesh.magSf()
);
```

*Interpolation of variables from left and right side of interface of control volume*



```
surfaceScalarField ap
(
    "ap",
    max(max(phiv_pos + cSf_pos, phiv_neg + cSf_neg), v_zero)
);

surfaceScalarField am
(
    "am",
    min(min(phiv_pos - cSf_pos, phiv_neg - cSf_neg), v_zero)
);

surfaceScalarField a_pos("a_pos", ap/(ap - am));

surfaceScalarField amaxSf("amaxSf", max(mag(am), mag(ap)));

surfaceScalarField aSf("aSf", am*a_pos);

if (fluxScheme == "Tadmor")
{
    aSf = -0.5*amaxSf;
    a_pos = 0.5;
}

surfaceScalarField a_neg("a_neg", 1.0 - a_pos);

phiv_pos *= a_pos;
phiv_neg *= a_neg;

surfaceScalarField aphiv_pos("aphiv_pos", phiv_pos - aSf);
surfaceScalarField aphiv_neg("aphiv_neg", phiv_neg + aSf);
```

*Calculations for scalar dissipation coefficient  
For  $K_J$  or  $K_{NP}$  scheme based on  
user input flag (fluxScheme=="Tadmor")*

```
phi = aphiv_pos*rho_pos + aphiv_neg*rho_neg;

surfaceVectorField phiU(aphiv_pos*rhoU_pos + aphiv_neg*rhoU_neg);
// Note: reassembled orientation from the pos and neg parts so becomes
// oriented
phiU.setOriented(true);

surfaceVectorField phiUp(phiU + (a_pos*p_pos + a_neg*p_neg)*mesh.Sf());

surfaceScalarField phiEp
(
    "phiEp",
    aphiv_pos*(rho_pos*(e_pos + 0.5*magSqr(U_pos)) + p_pos)
  + aphiv_neg*(rho_neg*(e_neg + 0.5*magSqr(U_neg)) + p_neg)
  + aSf*p_pos - aSf*p_neg
);
```

*Calculation of Convective Fluxes of Governing Equations*

# Compressible FVM: CFD



```
volScalarField muEff("muEff", turbulence->muEff());
volTensorField tauMC("tauMC", muEff*dev2(Foam::T(fvc::grad(U))));
```

Calculation of viscosity and shear stress tensor

```
// --- Solve density
```

```
solve(fvm::ddt(rho) + fvc::div(phi));
```

 Solves density equation

```
// --- Solve momentum
```

```
solve(fvm::ddt(rhoU) + fvc::div(phiUp));
```

Predictor step of momentum equation

```
U.ref() =
```

```
rhoU()
```

```
/rho();
```

Update Boundary Condition on U

```
U.correctBoundaryConditions();
```

```
rhoU.boundaryFieldRef() == rho.boundaryField()*U.boundaryField();
```

```
if (!inviscid)
```

```
{
```

Viscous Corrector  
step of momentum equation

```
solve
```

```
(
```

```
fvm::ddt(rho, U) - fvc::ddt(rho, U)
```

```
- fvm::laplacian(muEff, U)
```

```
- fvc::div(tauMC)
```

```
);
```

```
rhoU = rho*U;
```

```
}
```

```
// --- Solve energy
```

```
surfaceScalarField sigmaDotU
```

```
(
```

```
"sigmaDotU",
```

Calculation of work done by viscous forces

```
(
```

```
fvc::interpolate(muEff)*mesh.magSf()*fvc::snGrad(U)
```

```
+ fvc::dotInterpolate(mesh.Sf(), tauMC)
```

```
)
```

```
& (a_pos*U_pos + a_neg*U_neg)
```

```
solve
```

```
(
```

```
fvm::ddt(rhoE)
```

```
+ fvc::div(phiEp)
```

```
- fvc::div(sigmaDotU)
```

```
);
```

Predictor step of energy equation

```
e = rhoE/rho - 0.5*magSqr(U);
```

```
e.correctBoundaryConditions();
```

Update Boundary Condition on e  
Extract temperature from e

```
thermo.correct();
```

```
rhoE.boundaryFieldRef() ==
```

```
rho.boundaryField()*
```

Update boundary field of rhoE

```
(
```

```
e.boundaryField() + 0.5*magSqr(U.boundaryField())
```

```
);
```

```
if (!inviscid)
```

```
{
```

Viscous Corrector  
step of energy equation

```
solve
```

```
(
```

```
fvm::ddt(rho, e) - fvc::ddt(rho, e)
```

```
- fvm::laplacian(turbulence->alphaEff(), e)
```

```
);
```

```
thermo.correct();
```

```
rhoE = rho*(e + 0.5*magSqr(U));
```

```
}
```



```
p.ref() =  
    rho()  
    /psi();  
p.correctBoundaryConditions();  
rho.boundaryFieldRef() == psi.boundaryField()*p.boundaryField();  
  
turbulence->correct();  
  
runTime.write();  
  
runTime.printExecutionTime(Info);
```

*Update value of pressure field and continue to the next time loop*

- **The Sod's Shock Tube Problem**

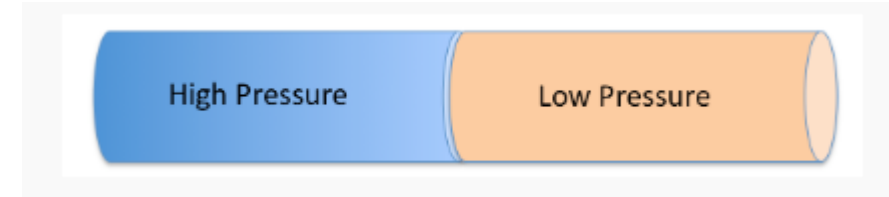
Initial Conditions

The state on the left side of the diaphragm is denoted by L

$$\rho_L = 1.0 \text{ kg/m}^3, p_L = 100000 \text{ Pa}, T_L = 348.4 \text{ K}$$

The state on the right side of the diaphragm is denoted by R

$$\rho_R = 0.125 \text{ kg/m}^3, p_R = 10000 \text{ Pa}, T_R = 278.7 \text{ K}$$



# Demo examples in OF: CFD



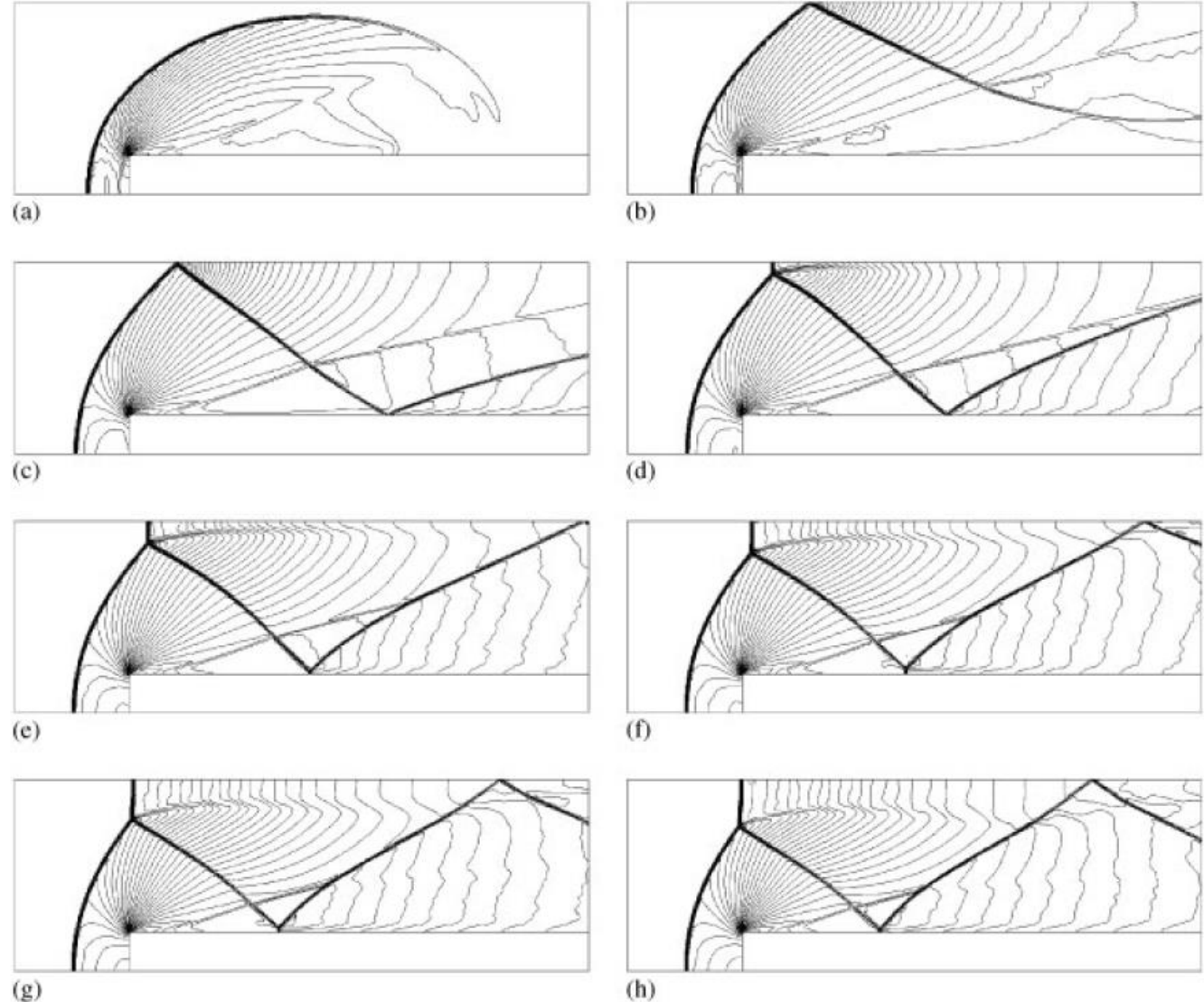
- Forward Step (Woodward and Collela)



Unsteady flow over the forward step  
placed in supersonic stream of air

Mach 3 flow

A standard benchmark test case to access the  
Performance of various numerical schemes

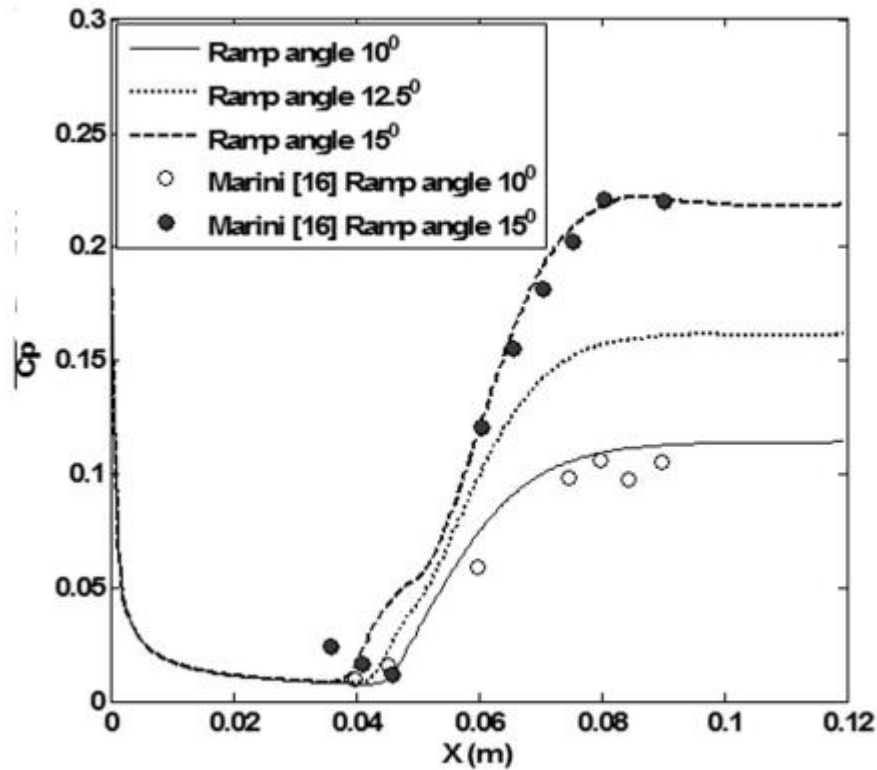




# Demo examples in OF: CFD



- Shock wave boundary layer interaction on compression corner



$U=1380\text{m/s}$   
 $P=191.4345\text{ Pa}$   
 $T=131.7$   
 $T_{\text{wall}} = 300\text{ K}$

