







Principles of CFD

- Our aim is the approximate solution of the governing PDE-s via numerical methods.
 - Four major methods are in use:
 - finite difference method,
 finite element method,
 - spectral methods,
 - finite volume method.
- · In the field of CFD finite volume methods prevail.
- The domain is subdivided into smaller volumes (cells).
- This process is called: the grid generation or meshing. One single value of every field variable is stored in each cell center. Within the cell the field variables are approximated by linear functions.

00000 Finite volume method 0000

- The integral form of the above conservation equations are discretized. Divergence terms are converted into surface integrals over the facets enclosing the cells. The numeric approximation of the flux integral for one facet depend only on the unknown ϕ values stored in the centers of the two neighboring cells adjacent to the facet. As a result of this so called discretization process, every transport equation provides one (non-finear) adjebraic equation per cell, e.g. if we have 5 transport equations and 1 000 000 cells then we obtain a system of 5000 000 non-linear adjebraic equations. In the case of time dependent problems we have to solve this system of equation in every time sten.

- case of time dependent problems we have to solve this system of equation in every time step. Each algebraic equations contain unknown e values from one particular cell and from all neighboring cells. This is e.g. 5 unknowns per equations to tetrahedral grids. Due to the high number of unknowns and the non-linearity of the system of equations **iterative** solution methods have to be used. The solution is first **initialized**, and than, iteratively refined thus **converging** towards the final solution. Integral of fluxes over the boundary facets meed to be defined in consistence with the physical characteristics of the region outside of the boundary by imposing additional mathematical conditions: **boundary** constituents. Surface integrals are numerically evaluated for every small facets connecting two neighboring cells. These values express the flow rates of conserved quantities (mass, momentum, energy). When we numerically calculate the integral of such conserved quantities for the whole domain, the surface integrals for internal facets are canceled, therefore the conservation equations for the whole domain are exactly fulfilled. E.g. the mass flow-rate through the intel's tequal to that of the outlet. This is called as the **conservative behavior** of the finite volume method.

