



PARA2012 – Elmer Tutorial

Elmer

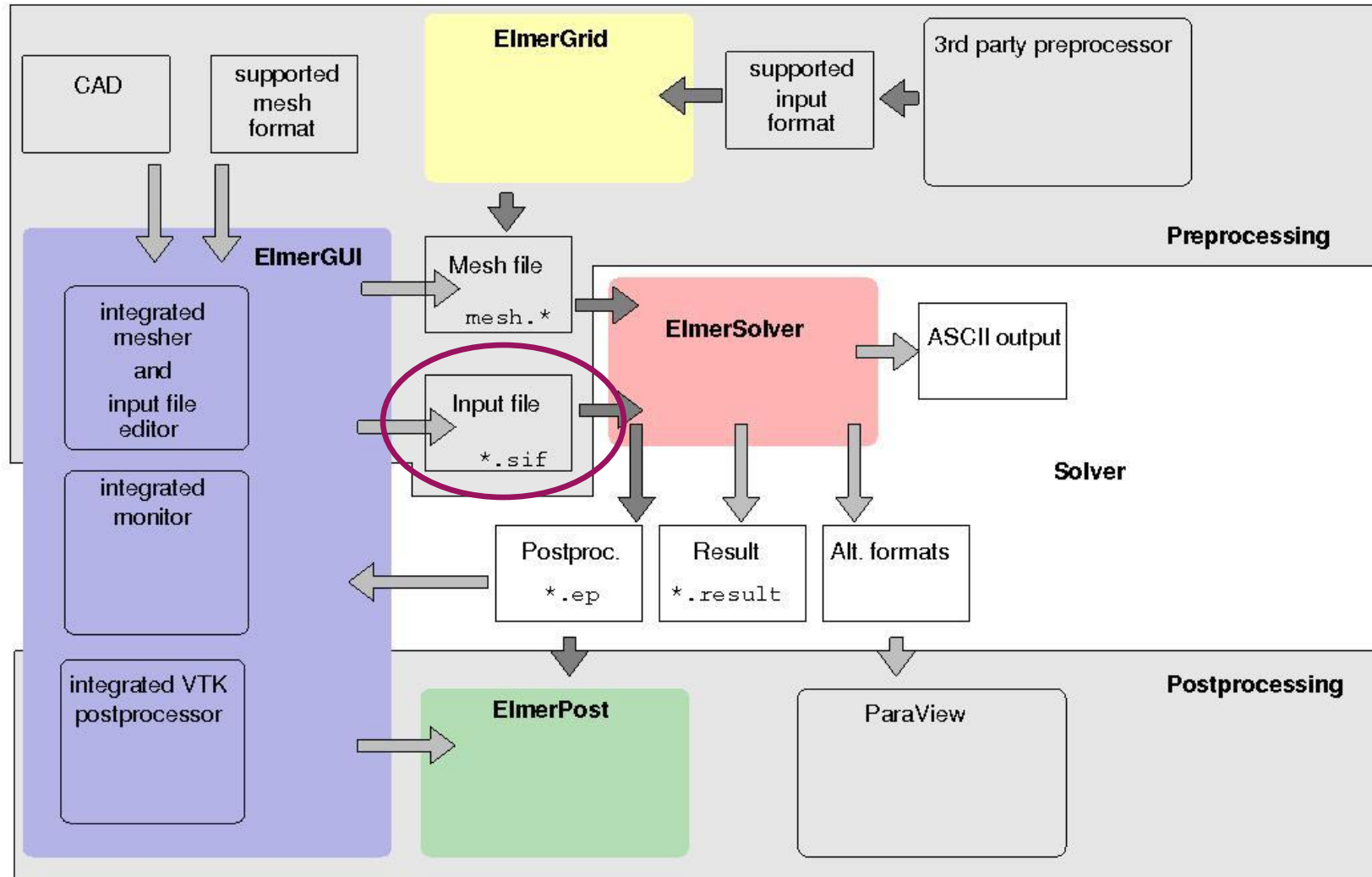
Elmer Solver Input File - Almost all about the text interface to ElmerSolver

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 - Constants
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 - Body
 - Equation
 - Body Force
 - Material
 - Initial Condition
 - Boundary Condition
- Tables and Arrays
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Elmer - Modules



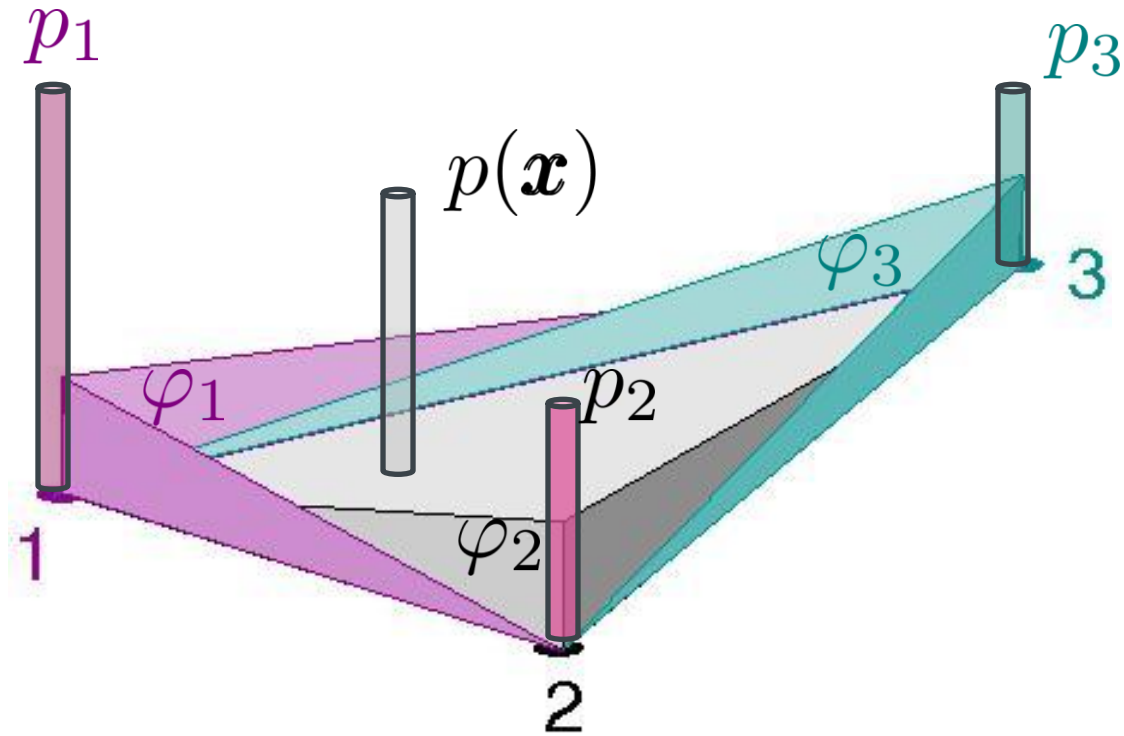
From the PDE to the SIF

- General advection-diffusion problem

$$c\rho \left(\frac{\partial \Psi}{\partial t} + \mathbf{u} \nabla \Psi \right) = \nabla \cdot \underbrace{(-\kappa \nabla \Psi)}_{\mathbf{q}} + \rho \sigma$$

- For instance, heat transfer problem: $\Psi = T$
- Coupled to (Navier-)Stokes via velocity: \mathbf{u}
- Non-linearities via material parameters, e.g.,
 $c = c(\Psi)$

From the PDE to the SIF



$$p(\mathbf{x}) = p_1 \varphi_1 |_{\mathbf{x}} + p_2 \varphi_2 |_{\mathbf{x}} + p_3 \varphi_3 |_{\mathbf{x}}$$

From the PDE to the SIF



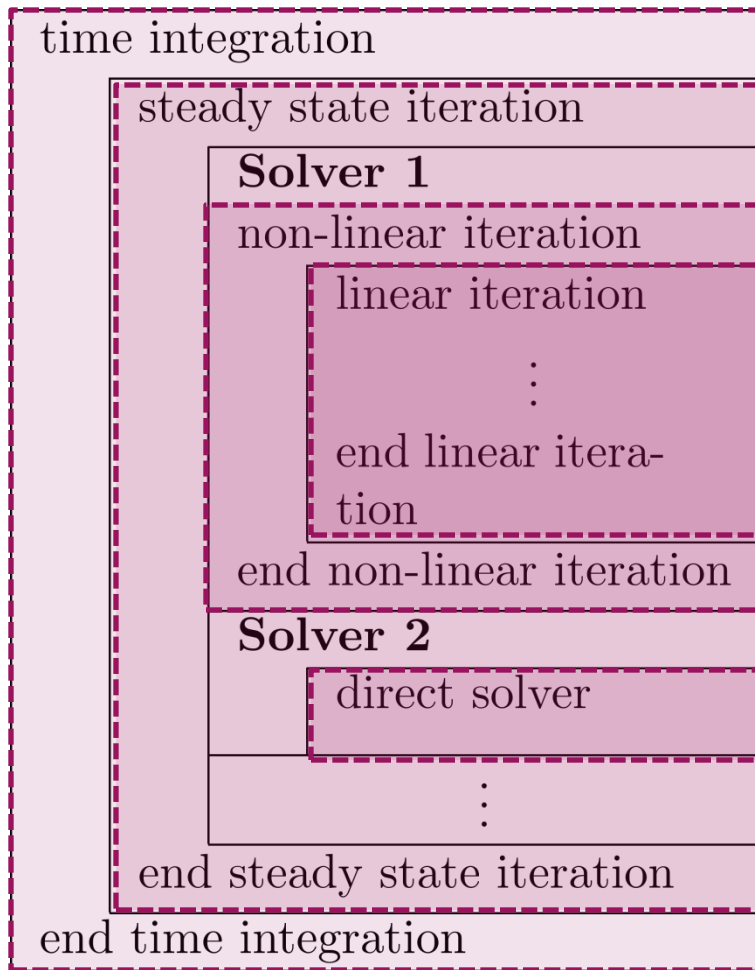
➤ Weak formulation:

$$\underbrace{\Psi_\beta a(\Delta t) \int_{\Omega} \rho c \varphi_\beta \varphi_\alpha d\Omega}_{\text{M}} + \underbrace{\Psi_\beta \int_{\Omega} [\rho c \mathbf{u} \cdot \nabla \varphi_\beta \varphi_\alpha + \kappa \nabla \varphi_\beta \cdot \nabla \varphi_\alpha] d\Omega}_{\text{S}} =$$

$$\underbrace{\int_{\partial\Omega} (\mathbf{q} \varphi_\alpha) \cdot \mathbf{n} d\Omega}_{\text{nat. BC}} + \underbrace{\int_{\Omega} \rho \sigma \varphi_\alpha d\Omega}_{\text{f}}$$

- Time integration
- Steady State: dependence on other variables
- Non-linear iteration: internal dependence

From the PDE to the SIF



1. Timestep Intervals
2. Steady State Max Iterations
3. Nonlinear Max Iterations
4. Linear System Max Iterations
4. Linear System Convergence Tolerance
3. Nonlinear System Convergence Tolerance
4. Linear System Convergence Tolerance
2. Steady State Convergence Tolerance
- 1.

Syntax of SIF

- Do not use non-printable characters!
 - No Tabulators, etc.
- A comment is preceded by a !
- Parameters in general have to be casted by their type:
 - **Real, Integer, Logical, String, File**
 - Exception: entry in Keyword DB
- Arrays have to be declared with the name:
Array (4) = Real 1.0 2.0 3.0 4.0

Sections of SIF

- The SIF is structured into sections
 - Header
 - Constants
 - Simulation
 - Solver
 - Body
 - Equation
 - Body Force
 - Material
 - Initial Condition
 - Boundary Condition

The contents of each section is between the keyword above and an **End**-statement

Sections of SIF: Header

- Declares search paths for mesh

Header

```
Mesh DB "." "dirname"
```

End

- preceding path + directory name of mesh database
- Replace path and *dirname* to fit your case

Sections of SIF: Constants

- Declares simulation-wide constants

Constants

```
Gas Constant = Real 8.314E00
```

```
Gravity (4) = 0 -1 0 9.81
```

```
End
```

- a casted scalar constant
- Gravity vector, an array with a registered name

Sections of SIF: Simulation



- Declares details of the simulation:

Simulation

```
Coordinate System = "Cartesian 2D"
```

- choices:
Cartesian{1D,2D,3D},
Polar{2D,3D},
Cylindric, Cylindric
Symmetric, Axi
Symmetric

```
Coordinate Mapping(3) = Integer 1 2 3
```

- Permute, if you want to interchange directions in mesh

```
Simulation Type ="Transient"
```

- Steady State or
Transient

```
Output Intervals(2) = 10 1
```

- Interval of results being written to disk

Sections of SIF: Simulation



- Declares details of the simulation:

```
Steady State Max Iterations = 10
```

```
Steady State Min Iterations = 2
```

```
Timestepping Method = "BDF"
```

```
Timestep Intervals(2) = 10 100
```

```
Timestep Sizes(2) = 0.1 1.0
```

```
Output File = "name.result"
```

```
Post File = "name.ep"
```

- How many min/max rounds on one timelevel/in a steady state simulation
- Choices: **BDF** or **Crank-Nicholson**
- Has to match array dimension of **Timestep Sizes**
- The length of one time step
- Contains data for restarting
- Contains ElmerPost data

Sections of SIF: Simulation



- Declares details of the simulation:

```
Restart File = "previous.result"
```

```
Restart Position = 10
```

```
Restart Time = 100
```

```
Initialize Dirichlet Condition =  
False
```

```
Restart Before Initial Conditions =  
True
```

```
Max Output Level = 3
```

```
End
```

- Restart from this file at file-entry (not necessarily timestep!) no. 10 and set time to 100 time-units
- Default is True. If false, Dirichlet conditions are called at Solver execution and not at beginning
- Default is False. If True, then Initial Condition can overwrite previous results
- Level of verbosity. 1 = basics, 9 = all and everything

Sections of SIF: Solver



- Declares a physical model to be solved

`Solver 3`

`Equation = "Navier-Stokes"`

`Exec Solver = "Always"`

`Linear System Solver = "Iterative"`

`Linear System Iterative Method =
"BiCGStab"`

`Linear System Convergence Tolerance
= 1.0E-06`

`Linear System Abort Not Converged =
True`

`Linear System Preconditioning =
"ILU2"`

- Numbering from 1 (priority)
- The name of the equation
- **Before/After Simulation/Timestep**
- Choices: **Iterative** or **Direct**
- Lots of choices here
- Convergence criterion
- If not True (default) continues simulation in any case
- Lots of choices

Sections of SIF: Solver

➤ Solving the **Linear(ized) Problem**

$$\mathbf{A} \Psi = \mathbf{f}$$

➤ Keyword:

Linear System Solver

➤ 3 ways to do that in Elmer:

1. **Direct** methods (= inversion of \mathbf{A})
2. **Iterative** methods (=working with approximations to \mathbf{A})
3. Multi-grid methods (not discussed here)

Sections of SIF: Solver

➤ Direct linear system solver

➤ Keyword:

Linear System Direct Method

- Banded (default) LAPack
 - UMFPACK Unsymmetric MultiFrontal method (only serial)
 - MUMPS Unsymmetric MultiFrontal method (only parallel)
- Sometimes the only way to go (if bad conditioned)
 - Costly: Elimination takes $\sim N^3$ operations and needs to store N^2 unknowns in memory

Sections of SIF: Solver

Iterative solvers:

- Krylov subspace: $\mathbf{x}_n \in \text{span}(K_n)$

$$K_n = [\mathbf{f}, \mathbf{A}\mathbf{f}, \mathbf{A}^2\mathbf{f}, \mathbf{A}^3\mathbf{f}, \dots, \mathbf{A}^{n-1}\mathbf{f}]$$

$$\mathbf{R} = (\mathbf{f} - \mathbf{A}\mathbf{x}_n) \rightarrow \min.$$

– Linear System Iterative Method

- GMRES Generalized Minimal Residual Method
- CG, CGS, BiCGStab Conjugate Gradient
- TFQMR Transpose-free quasi-minimal residual
- GCR Generalized Conjugate Residual

Section of SIF: Solver

➤ But, before we solve, we usually apply a

pre-conditioner

– Find $\tilde{\mathbf{P}}\mathbf{A}$, but much easier to invert

– $\tilde{\mathbf{P}}\mathbf{A} \sim \mathbf{I}$ has favourable condition number

$$\tilde{\mathbf{P}}^{-1}\mathbf{A}\Psi = \tilde{\mathbf{P}}^{-1}\mathbf{f}$$

Linear System Preconditioning

- None
- Diagonal
- $ILUn$ ($n=0,1,2,\dots$) and $ILUT$

Sections of SIF: Solver



- Declares a physical model to be solved

Nonlinear System Convergence

Tolerance = 1.0E-05

Nonlinear System Max Iterations = 20

Nonlinear System Min Iterations = 1

**Nonlinear System Newton After
Iterations = 10**

**Nonlinear System Newton After
Tolerance = 1.0e-03**

**Steady State Convergence Tolerance =
1.0E-03**

Stabilization Method = Stabilized

End

- Convergence criterion for non-linear problem
- The maximum rounds
- The minimum rounds
- Switch from Picard to Newton scheme after 10 iterations ...
- ... or after this criterion (NV.: has to be smaller than convergence criterion of hit)
- The convergence on the time-level
- Convection needs stabilization. Alternatives: **Bubbles** or **P2/P1**

Sections of SIF: Body



- Declares a physical model to be solved

Body 2

Name = "pipe"

Equation = 2

Material = 2

Body Force = 1

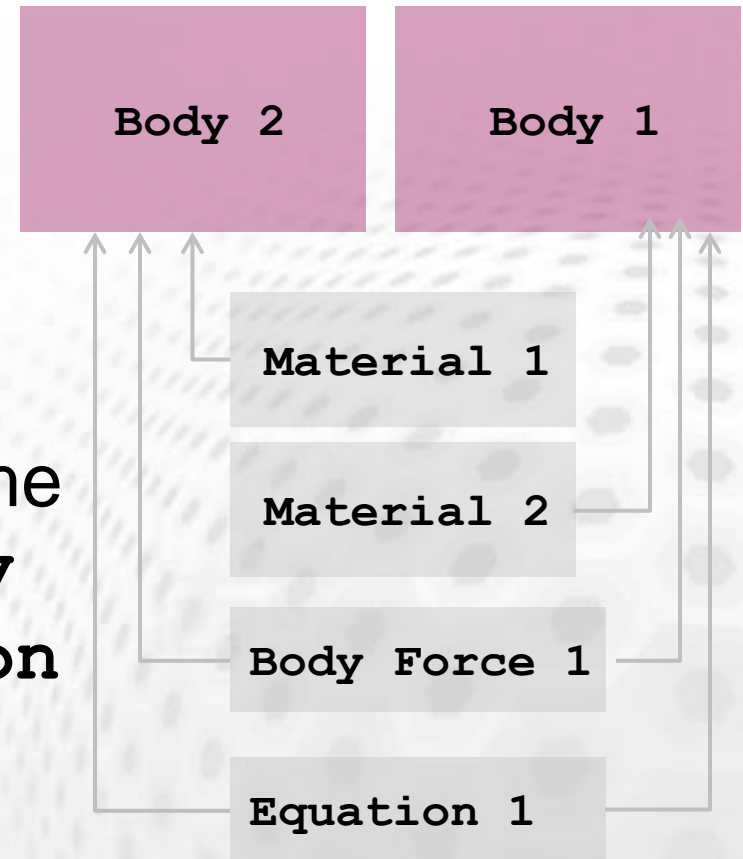
Initial Condition = 2

End

- Numbering from 1 to number of bodies
- Identifier of the body
- The assigned set of equations
- The assigned material section
- The assigned body force
- The assigned initial condition

Sections of SIF: Body

- Each **Body** has to have an **Equation** and **Material** assigned
 - **Body Force**, **Initial Condition** optional
- Two bodies can have the same **Material/Equation/Body Force/Initial Condition** section assigned



Sections of SIF: Equation



- Declares set of solvers for a body

```
Equation 2
```

```
Active Solvers(2) = 1 3
```

```
Convection = Computed
```

```
NS Convect = False
```

```
End
```

- Numbering from 1 to number of equation sets
- Declares the solvers (according to their numbers) to be solved within this set
- Important switch to account for convection term. Alternatives: None and Constant (needs Convection Velocity to be declared in the Material section)
- Sets no convection for Navier-Stokes (=Stokes) alternative:
Flow Model = Stokes
in the Solver section of Navier-Stokes

Sections of SIF: Body Force



- Declares body forces and bulk and execution conditions for a body

```
Body Force 3
```

```
Flow Body Force 1 = 0.0
```

```
Flow Body Force 2 = -9.81
```

```
MyVariable = Real 0.0
```

```
Flow Solution Passive = 1.0
```

```
End
```

- Numbering from 1 to number of body forces
- Gravity pointing in negative x-direction applied to Navier-Stokes solver
- A Dirichlet condition for a variable set within the body
- Suspends execution of Navier-Stokes solver (-1.0 = default = execution)

Body forces can be functions of other variables

Sections of SIF: Material



- Declares set of material parameters for body

```
Material 1

Density = 1000.0

Heat Conductivity(3,3) = 1 0 0\
                        0 1 0\
                        0 0 2

Viscosity = Variable Temperature
Real MATC "viscosity(tx)"

MyMaterialParameter = Real 0.0

End
```

- Numbering from 1 to number of material
- Always declare a density (mandatory)
- Parameters can be arrays
- Or functions of other variables
- Non-keyword DB parameters have to be casted

Sections of SIF: Initial Condition



Declares initial conditions for a body

By default restart values are used

```
Initial Condition 2
```

```
Velocity 1 = Variable Coordinate 2
```

```
Real MATC "42.0*(1.0 - tx/100.0)"
```

```
Velocity 2 = 0.0
```

```
MyVariable = Real 20.0
```

```
End
```

- Numbering from 1 to number of IC's
- Initial condition as a function of a variable ...
- ... and as a constant
- Non-keyword DB parameters have to be casted

Sections of SIF: Boundary Condition



- Declares conditions at certain boundaries

```
Boundary Condition 3
```

```
Target Boundaries(2) = 1 4
```

```
Velocity 1 = Variable Coordinate 2
```

```
Real MATC "42.0*(1.0 - tx/100.0)"
```

```
Velocity 2 = 0.0
```

```
Normal-Tangential Velocity = Logical True
```

```
End
```

- Numbering from 1 to number of BC's
- The boundaries of the mesh the BC is assigned to
- Variable as a function and ...
... as a constant
- Set velocities in normal-tangential system

Tables and Arrays



- Tables (piecewise linear or cubic):

```
Density = Variable Temperature
Real cubic
    0 900
    273 1000
    300 1020
    400 1000
End
```

- Arrays:

```
Target Boundaries(3) = 5 7 10

MyParamterArray(3,2) = Real 1 2\
                             3 4\
                             5 6
```

- Expresions:

```
OneThird = Real $1.0/3.0
```

MATC

- Syntax close to C
- Even if-conditions and loops
- Can be use for on-the-fly functions inside the SIF
- Documentation on [web-pages](#)
- Do not use with simple numeric expressions:
is much faster than

```
OneThird = Real $1.0/3.0
```

```
OneThird = Real MATC "1.0/3.0"
```

MATC

- Use directly in section:

```
Heat Capacity = Variable Temperature  
Real MATC "2.1275D03 + 7.253D00*(tx - 273.16)"
```

- Even with more than one dependency:

```
Temp = Variable Latitude, Coordinate 3  
Real MATC "49.13 + 273.16 - 0.7576*tx(0) - 7.992E-03*tx"
```

- Or declare functions (somewhere in SIF, outside a section)

```
$ function stemp(X) {\  
  stemp = 49.13 + 273.16 - 0.7576*X(0) - 7.992E-03*X(1) }  
}
```

– being called:

```
Temp = Variable Latitude, Coordinate 3  
Real MATC "_stemp(tx)"
```

User Defined Functions (UDF)

- Written in Fortran 90
 - Dynamically linked to Elmer
 - Faster, if more complicated computations involved
 - Compilation command **elmerf90**
- ```
elmerf90 myUDF.f90 -o myUDF.f90
```
- Call from within section:

```
MyVariable = Variable Temperature
Real Procedure "myUDF" "myRoutine"
```



# User Defined Functions (UDF)

➤ Example:  $\rho(T[K]) = 1000.0 \cdot [1 - 1 \times 10^{-4} \cdot (T - 273.15)]$

```
FUNCTION getdensity(Model, N, T) RESULT(dens)
USE DefUtils !important definitions
IMPLICIT None
TYPE(Model_t) :: Model
INTEGER :: N
REAL(KIND=dp) :: T, dens
dens = 1000.0_dp*(1.0_dp - 1.0d-04*(T - 273.0_dp))
END FUNCTION getdensity
```

- Definitions loaded from **DefUtils**
- Header: **Model** access-point to all ElmerSolver inside data; Node number **N**; input value **T**