



Input Data Format Specification

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Contents

1	Introduction	3
2	Running the code	3
3	Syntax and general rules	4
4	Output file record	6
5	Job description record	6
6	Analysis record	6
6.1	Linear static analysis	8
6.2	LinearStability	8
6.3	EigenValueDynamic	8
6.4	NIDEIDynamic	9
6.5	DEIDynamic	9
6.6	DIIDynamic	10
6.7	IncrementalLinearStatic	10
6.8	NonLinearStatic	10
6.8.1	Extended syntax	11
6.8.2	Standard syntax	16
6.9	Adaptive linear static	16
6.10	Adaptive nonlinear static	17
6.11	Stationary transport problem	18
6.12	Transient transport problem - linear case	19
6.13	Transient transport problem - nonlinear case	19
6.14	Transient incompressible flow - CBS Algorithm	20
6.15	Transient incompressible flow SUPG/PSPG Algorithm	21
6.16	Staggered Problem	22
6.17	Sparse linear solver parameters	22
6.18	Eigen value solvers	24
6.19	Dynamic load balancing parameters	24
6.20	Error estimators and indicators	25
6.21	Material interfaces	27
6.22	Mesh generator interfaces	28
6.23	Initialization modules	28
6.24	Export modules	29
7	Domain record(s)	31
7.1	Output manager record	32
7.2	Components size record	33
7.3	Dof manager records	33
7.4	Element records	37
7.5	Cross section records	38
7.6	Material type records	39
7.7	Nonlocal barrier records	39
7.8	Load and boundary conditions	40
7.9	Initial conditions	43

7.10	Time functions records	43
7.11	Xfem manager record and associated records	44
8	Examples	45
8.1	Beam structure	45
8.2	Plane stress example	46
9	Examples - parallel mode	47
9.1	Node-cut example	47
9.2	Element-cut example	48
10	Figures	50

1 Introduction

This report describes in details the format and structure of OOFEM text input file. Input file can be created by any available text editor or can be generated by some conversion program or by some FEM pre-processor.

Some parts of this document are related parallel version of the code (POOFEM). These parts are distinguished by typing the text in sans serif font family. In a parallel mode a set of input files must be provided, each one for particular process and related partition. The input file corresponding to some partition is called partition input file. Its name is composed from two parts. The first part (referred as base name) is user-defined. The second part (called partition name), divided from first part by full-stop, is particular partition rank number. The partitions are numbered from zero. It is assumed, that all partitions input files have the same base name. The partition name is appended automatically to input file name, user provides only base name on input. On the other hand, the partition name is not appended to output file name, as specified in output file record (see section 4).

The parallel version requires the unique global numbering for dof managers (nodes) and elements. The global numbering is necessary to link partitions together. However, the dof managers and optionally elements at interpartition boundaries, should be explicitly marked, in order to distinguish different type of relations between their remote counterparts (see further).

2 Running the code

The program may be executed by typing

```
oofem [option [parameter]] ...
```

on the command line prompt with the following command line options:

- **-f string**
oofem input file name, if not present, program interactively reads this parameter.
- **-r int**
Restarts the analysis from given solution step. The corresponding context file (*.osf) must exist.
- **-rn**
Turns on the profile renumbering. Default is off. It will not work in parallel mode.
- **-ar int**
Restarts the adaptive computation from given solution step. Requires the corresponding context file (*.osf) and domain input file (*.din) to exists. The domain input file describes the new mesh, its syntax is identical to syntax of input file, but it does not contains the output file record, job description record and analysis record.
- **-qo string**
Redirect the standard output stream (stdout) to given file.
- **-qe string**
Redirect standard error stream (stderr) to given file.

- **-context**

Forces the creation of context file for each solution step.

The parallel version uses the MPI (Message Passing Interface) standard for message-passing communication. Thus, to execute POOFEM program, users must know the procedure for beginning MPI jobs on their selected computer system(s). For instance, when using the MPICH implementation of MPI and many others, the following command initiates a program that uses eight processors:

```
mpirun -np 8 poofem program_options
```

3 Syntax and general rules

Input file is created by records. In the current implementation, the particular record is represented by one line in input file. The order of records in file is compulsory, and is following

1. output file record, see section 4,
2. job description record, see section 5,
3. analysis record, see section 6,
4. domain record, see section 7,
5. output manager record, see section 7.1,
6. components size record, see section 7.2,
7. node & element-side record(s), see section 7.3,
8. element record(s), see section 7.4,
9. cross section record(s), see section 7.5,
10. material type record(s), see section 7.6,
11. nonlocal barriers record(s), see section 7.7,
12. load, boundary conditions record(s), see section 7.8,
13. initial conditions record(s), see section 7.9,
14. time functions record(s), see section 7.10.
15. optional xfem manager and associated record(s), see section 7.11

When line begins with '#' character, then it is skipped by parser and provide a way, how to include comments inside input file.

Records contain many fields, each field is characterized by keyword and its associated field value. Some keywords have no field values. Except the first field, the order of remaining fields in record is optional. There are several exceptions, which will be described in particular sections. The type of field value is specific to keyword. General format of field, characterized by keyword "Keyword" with corresponding value (marked as #) of type "type" is **Keyword**_(type) #. If keyword is variable, depending on entity type which is described by particular

record, then the keyword is preceded by star. We call such keyword as entity keyword (For example keyword, which describes element type). The possible substitutions for entity keyword are typed using **Typewriter** font family. Often, many fields are specific to particular entity keyword. Then we describe general format of record without these specific fields and we describe them in entity keyword specific part of record description. Sometimes, there are fields without keywords. These are usually compulsory, with fixed position in record. Their occurrence is described using so called named values. The format of named value is “(**name#**)_(type)”, where a value of required type should be substituted.

In a special case of analysis record, the input record can be followed by optional meta-step input records (see section 6). Then certain fields originally in analysis record should appear in meta-step record instead. This is marked by adding “M” superscript to keyword. Then the field format is **Keyword**^M_(type) #.

As already mentioned, each field value is typed. The possible types are:

- in - integer number;
- rn - real number;
- ch - character (usually for unknowns description ('d' for displacement unknown, 't' for temperature unknown, ...));
- ia - integer array, format of integer array is “size val(1) ... val(size)”, where size, val(1), ..., val(size) are integer numbers. Values are separated by one or more spaces;
- rn - real array, format of real array is “size val(1) ... val(size)”, where size is integer number and val(1), ..., val(size) are real numbers. Values are separated by one or more spaces;
- dc - dictionary. Dictionary consist of pairs, each pair has key (character type) and its associated value (integer type). Format of dictionary is “size key(1) val(1) ... key(size) val(size)”, where size is integer number, key(1), ..., key(size) are single character values, and val(1), ..., val(size) are real numbers. Values are separated by one or more spaces;
- rl - range list. Range list syntax is { number1 .. numberN (start1 end1) (start2 end2)}. The enclosing brackets are compulsory. The range list represent list of integer values. Single values can be specified using single values (number1, ..., NumberN). The range of values (all numbers from startI to endI including startI and endI can be specified using range value in the form (startI endI). The range is described using its start and end values enclosed in parenthesis. Any number of ranges and single values can be used to specify range list.
- et - entity type. For example, it describes the finite element type. Possible type values are mentioned in specific sections.
- s - character string. The string have to be enclosed in quotes (”) following after corresponding keyword.
- expr - function expression. The expression have to be enclosed in quotes (”). The expression is evaluated by internal parser and represent mathematical expressions as a function of certain variables. The variable names

and meaning are described in specific sections. The usual arithmetic operators like $-$, $+$, $*$, $/$ are supported and their evaluation order is taken into account. The evaluation order can be changed using parenthesis. Several built-in functions are supported (sqrt, sin, cos, tan, atan, asin and acos) - these must be typed using lowercase letters and their arguments must be enclosed in parenthesis.

The general format of record is

```
keyword1(type) # [keyword2(type) #] \
... \
[keywordXX(type) #]
```

The keywords and their values are separated by one or more spaces. Please note, that a single record corresponds to one input line in input file. If any keyword or field value is enclosed within angle brackets $\langle \rangle$ then it is related to parallel version of oofem is not available in sequential version. When some field is enclosed in brackets $[]$, then it's use is optional and often overwrites the default behavior or adds additional (but optional) information or property (for example adds a loading to node).

4 Output file record

This record has no keywords and contains character string, which describes the path to output file. If file with the same name exists, it will be overwritten.

5 Job description record

This record has no keywords and contains character string, which describes the job. This description will appear in the output file.

6 Analysis record

This record describes the type of analysis, which should be performed. The general format of this record can be specified in

- “standard-syntax”


```
*AnalysisType nsteps(in) # \
                  [renumber(in) #] \
                  attributes(string) # \
                  [ninitmodules(in) #] \
                  [nmodules(in) #] \
                  [nxfemman(in) #]
```
- “meta step-syntax”


```
*AnalysisType nmsteps(in) # \
                  [ninitmodules(in) #] \
                  [nmodules(in) #] \
                  [nxfemman(in) #]
```

immediately followed by *nmsteps* meta step records with the following

syntax:

```
| nsteps(in) # attributes(string) #
```

The *nsteps* parameter determines the number of “metasteps”. The meta step represent sequence of solution steps with common attributes. There is expected to be *nsteps* subsequent metastep records. The meaning of meta step record parameters (or analysis record parameters in “standard syntax”) is following:

- *nsteps* - determines number of subsequent solution steps within metastep. Nonzero value of optional parameter *renumber* turns on the equation renumbering to optimize the profile of characteristic matrix (uses Sloan algorithm). By default, profile optimization is not performed. It will not work in parallel mode.
- *attributes* - contains the metastep related attributes of analysis (and solver), which are valid for corresponding solution steps within meta step. If used in standard syntax, the attributes are valid for all solution step.
- *ninitmodules* - number of initialization module records for given problem. The initialization modules are specified after meta step section (or after analysis record, if no metasteps are present). Initialization modules allow to initialize the state variables by values previously computed by external software. The available initialization modules are described in section 6.23.
- *nmodules* - number of export module records for given problem. The export modules are specified after initialization modules. Export modules allow to export computed data into external software for postprocessing. The available export modules are described in section 6.24.
- *xfemman* - number of xfem managers. Xfem managers wrap eXtended Finite Element (xfem) data: they manage the list of enrichment items and associated enrichment functions, and their geometry representation. The number of xfem managers required is problem dependent, typically a single xfem manager is used. The syntax of xfem manager record and related records is described in section 7.11.

Not all of analysis types support the metastep syntax, and if not mentioned, the standard-syntax is expected. Currently, supported analysis types are

- Linear static analysis, see section 6.1,
- Eigen value dynamic, see section 6.3,
- Direct explicit nonlinear dynamics, see section 6.4,
- Direct explicit (linear) dynamics, see section 6.5,
- Implicit linear dynamic, see section 6.6,
- Incremental **linear** static problem, see section 6.7,
- Non-linear static analysis, see section 6.8.

6.1 Linear static analysis

LinearStatic	nsteps _(in) #	\
	[renumber _(in) #]	\
	[sparselinsolverparams ₍₎ #]	\
	[sparselinsolverparams ₍₎ #]	

Linear static analysis. Parameter *nsteps* indicates the number of loading cases. Series of loading cases is maintained as sequence of time-steps. For each load case an auxiliary time-step is generated with time equal to load case number. Load vectors for each load case are formed as load vectors at this auxiliary time.

Nonzero value of optional parameter *renumber* turns on the equation renumbering to optimize the profile of characteristic matrix (uses Sloan algorithm). By default, profile optimization is not performed. It will not work in parallel mode.

The *sparselinsolverparams* parameter describes the sparse linear solver attributes and is explained in section 6.17. Can be used in parallel mode. The input record must be the same for all processors. At present, parallel version requires PETSc module.

6.2 LinearStability

LinearStability	nroot _(in) #	\
	[renumber _(in) #]	\
	rtolv _(rn) #	\
	[eigensolverparams ₍₎ #]	

Solves linear stability problem. Only first *nroot* smallest eigenvalues and corresponding eigenvectors will be computed. Relative convergence tolerance is specified using *rtolv* parameter.

Nonzero value of optional parameter *renumber* turns on the equation renumbering to optimize the profile of characteristic matrix (uses Sloan algorithm). By default, profile optimization is not performed. It will not work in parallel mode.

The *eigensolverparams* parameter describes the sparse linear solver attributes and is explained in section 6.18. Can be used in parallel mode. The input record must be the same for all processors. Parallel version requires PETSc and SLEPc modules.

6.3 EigenValueDynamic

EigenValueDynamic	nroot _(in) #	\
	[renumber _(in) #]	\
	rtolv _(rn) #	\
	[eigensolverparams ₍₎ #]	

Represents the eigen value dynamic analysis. Only *nroot* smallest eigenvalues and corresponding eigenvectors will be computed. Relative convergence criteria is governed using *rtolv* parameter. Nonzero value of optional parameter *renumber* turns on the equation renumbering to optimize the profile of characteristic matrix (uses Sloan algorithm). By default, profile optimization is not performed. It will not work in parallel mode.

The *eigensolverparams* parameter describes the sparse linear solver attributes and is explained in section 6.18. Can be used in parallel mode. The input record must be the same for all processors. Parallel version requires PETSc and SLEPc modules.

6.4 NIDEIDynamic

NIDEIDynamic	nsteps _(in) #	\
	[renumber _(in) #]	\
	dumpcoef _(rn) #	\
	[deltaT _(rn) #]	

Represents the direct explicit nonlinear dynamic integration. The central difference method with diagonal mass matrix is used, damping matrix is assumed to be proportional to mass matrix, $\mathbf{C} = \text{dumpcoef} * \mathbf{M}$, where \mathbf{M} is diagonal mass matrix. *deltaT* is time step length used for integration, which may be reduced by program in order to satisfy solution stability conditions. Parameter *nsteps* specifies how many time steps will be analyzed. Nonzero value of optional parameter *renumber* turns on the equation renumbering to optimize the profile of characteristic matrix (uses Sloan algorithm). By default, profile optimization is not performed. It will not work in parallel mode.

<PNIDEIDynamic>	<nsteps _(in) #	\
	<dumpcoef _(rn) #	\
	<[deltaT _(rn) #]	\
	<*commode>	\
	<[nonlocalext ₍₎ #]	

Represents the parallel direct explicit nonlinear dynamic integration. The central difference method with diagonal mass matrix is used, damping matrix is assumed to be proportional to mass matrix, $\mathbf{C} = \text{dumpcoef} * \mathbf{M}$, where \mathbf{M} is diagonal mass matrix. *deltaT* is time step length used for integration, which may be reduced by program in order to satisfy solution stability conditions. Parameter *nsteps* specifies how many time steps will be analyzed. The **<*commode>** keyword can be one from following:

- **nodecutmode** - indicates the node cut partitioning scheme is used. The cut is led along element edges or surfaces (see figures 5, 6, and 7).
- **elementcutmode** - the element cut partitioning is used. The partitioning cut led across the element edges or surfaces (see figures 8 and 9).

The *nonlocalext* turns on the nonlocal constitutive extension. The extension considers a band of remote elements involved in computation of nonlocal variables (see fig. 7 illustrating this approach for node-cut partitioning).

6.5 DEIDynamic

DEIDynamic	nsteps _(in) #	\
	dumpcoef _(rn) #	\
	[deltaT _(rn) #]	

Represent the **linear** explicit integration scheme for dynamic problem solution. The central difference method with diagonal mass matrix is used, damping matrix is assumed to be proportional to mass matrix, $\mathbf{C} = \text{dumpcoef} * \mathbf{M}$, where \mathbf{M} is diagonal mass matrix. deltaT is time step length used for integration, which may be reduced by program in order to satisfy solution stability conditions. Parameter $nsteps$ specifies how many time steps will be analyzed.

6.6 DIIDynamic

DIIDynamic	nsteps _(in) #	\
	[renumber _(in) #]	\
	deltaT _(rn) #	\
	alpha _(rn) #	\
	beta _(rn) #	\
	Psi _(rn) #	\

Represents direct implicit integration of linear dynamic problems. Damping is modeled as Rayleigh damping ($\mathbf{c} = \alpha * \mathbf{M} + \beta * \mathbf{K}$). Parameter Psi determines integration method used, for $Psi = 1$ the Newmark and for $Psi \geq 1.37$ the Wilson method will be used. Parameter deltaT is required time integration step length. Nonzero value of optional parameter $renumber$ turns on the equation renumbering to optimize the profile of characteristic matrix (uses Sloan algorithm). By default, profile optimization is not performed. It will not work in parallel mode.

6.7 IncrementalLinearStatic

IncrLinearStatic	endOfTimeOfInterest _(rn) #	\
	prescribedTimes _(ra) #	\
	[renumber _(in) #]	

Represents incremental **linear** static problem. The problem is solved as series of linear solutions and is intended to be used for solving linear creep problems or incremental perfect plasticity. Nonzero value of optional parameter $renumber$ turns on the equation renumbering to optimize the profile of characteristic matrix (uses Sloan algorithm). By default, profile optimization is not performed. It will not work in parallel mode.

Supports the changes of static scheme (applying, removing and changing boundary conditions) during the analysis.

Response is computed in times defined by *prescribedTimes* array. These times should include times, when generally the boundary conditions are changing, and in other times of interest. (For linear creep analysis, the values should be uniformly distributed on log-time scale, if no change in loading or boundary conditions). The time at the end of interested is specified using *endOfTimeOfInterest* parameter.

6.8 NonLinearStatic

NonLinearStatic

Non-linear static analysis. The problem can be solved under direct load or displacement control, indirect control, or by their arbitrary combination. Can be

used in parallel mode. The input record must be the same for all processors. At present, parallel version requires PETSc module. By default all material nonlinearities will be included, geometrical not. To include geometrically nonlinear effect one must specify level of non-linearity in element records. There are two different ways, how to specify the parameters - the extended and standard syntax.

6.8.1 Extended syntax

The extended syntax uses the “metastep” concept and has the following format:

```

NonLinearStatic  [nmsteps(in) #] \
                  nsteps(in) # \
                  [renumber(in) #] \
                  [contextOutputStep(in) #] \
                  [sparselinsolverparams(string) #] \
                  [nonlinform(in) #] \
                  <[nonlocstiff(in) #]> \
                  <[nonlocalext() #]> \
                  <[loadbalancing() #]>

```

This record is immediately followed by metastep records with the format described below. The analysis parameters have following meaning

- *nmsteps* - determines the number of “metasteps”, default is 1.
- *nsteps* - determines number of solution steps.
- Nonzero value of optional parameter *renumber* turns on the equation renumbering to optimize the profile of characteristic matrix (uses Sloan algorithm). By default, profile optimization is not performed. It will not work in parallel mode.
- *contextOutputStep* - causes the context file to be created for every contextOutputStep-th step and when needed. Useful for postprocessing.
- The *sparselinsolverparams* parameter describes the sparse linear solver attributes and is explained in section 6.17.
- *nonlinform* - formulation of non-linear problem. If == 1 (default), total Lagrangian formulation in undeformed original shape is used (first-order theory). If == 2, the equilibrated displacements are added to original ones and updated in each time step (second-order theory).
- *nonlocstiff* - determines whether the tangent stiffness extension for nonlocal models is activated. If == 0 (default) this option is not active. If == 1 the support for nonlocal tangent stiffness is activated.
- The *nonlocalext* turns on the nonlocal constitutive extension. The extension considers a band of remote elements involved in computation of nonlocal variables (see fig. 7 illustrating this approach for node-cut partitioning).
- The *loadbalancing* parameter describes the dynamic load balancing attributes and is explained in section 6.19.

The metastep record has following general syntax:

nsteps _(in) #	\
[controllmode _(in) #]	\
[deltat _(rn) #]	\
[stiffmode _(in) #]	\
[refloadmode _(in) #]	\
solverParams ₍₎ #	\
[sparselinsolverparams _(string) #]	\
[donotfixload ₍₎ #]	\

where

- *controllmode* - determines the type of solution control used for corresponding meta step. if == 0 then indirect control will be used to control solution process (arc-length method, default). if == 1 then direct displacement or load control will be used (Newton-Raphson solver). In the later mode, one can apply the prescribed load increments as well as control displacements.
- *deltaT* - is time step length. If not specified, it is set equal to 1,0. Each solution step has associated the corresponding intrinsic time, at which the loading is generated. The *deltaT* determines the spacing between solution steps on time scale.
- *stiffMode* - If == 0 (default) then tangent stiffness will be used at new step beginning and whenever numerical method will ask for stiffness update. If == 1 the use of secant tangent will be forced. The secant stiffness will be used at new step beginning and whenever numerical method will ask for stiffness update. If == 2 then original elastic stiffness will be used during the whole solution process.
- The *refloadmode* parameter determines how the reference force load vector is obtained from given totalLoadVector and initialLoadVector. The initialLoadVector describes the part of loading which does not scale. Works only for force loading, other non-force components (temperature, prescribed displacements should always given in total values). If refLoadInputMode is 0 (rlm_total, default) then the reference incremental load vector is defined as totalLoadVector assembled at given time. If refLoadInputMode is 1 (rlm_incremental) then the reference load vector is obtained as incremental load vector at given time.
- *solverParams* - parameters of solver. The solver type is determined using *controllmode*.
- The *sparselinsolverparams* parameter describes the sparse linear solver attributes and is explained in section 6.17.
- By default, reached load at the end of metastep will be maintained in subsequent steps as fixed, non scaling load and load level will be reset to zero. This can be changed using keyword *donotfixload*, which if present, causes the loading to continue, not resetting the load level. For the indirect control the reached loading will not be fixed, however, the new reference loading vector will be assembled for the new metastep.

The direct solver corresponds to *controllmode*=1 and the Newton-Raphson solver is used. Under the direct control, the total load vector assembled for specific solution step represents the load level, where equilibrium is searched. The implementation supports also displacement control - it is possible to prescribe one or more displacements by applying “quasi prescribed” boundary condition(s)¹ The load level then represents the time, where the equilibrium has been found. The Newton-Raphson solver parameters (*solverParams*) for load-control are:

```

maxiter(in) # \
[minsteplength(in) #] \
[minIter(in) #] \
[manrmsteps(in) #] \
[ddm(ia) #] [ddv(ra) #] [ddltf(in) #] \
[linesearch(in) #] [lsearchamp(rn) #] \
[lsearchmaxeta(rn) #] [lsearchtol(rn) #] \
[nccdg(in) # ccdgl(ia) # ... ccdgn(ia) # ] \
rtolv() # [rtolf() # rtold() # ]

```

where

- *maxiter* determines the maximum number of iterations allowed to reach equilibrium. If equilibrium is not reached, the step length (corresponding to time) is reduced.
- *minsteplength* parameter is the minimum step length allowed.
- *minIter* - minimum number of iterations which always proceed during the iterative solution.
- If *manrmsteps* parameter is nonzero, then the modified N-R scheme is used, with the stiffness updated after *manrmsteps* steps.
- *ddm* is array specifying the degrees of freedom, which displacements are controlled. Let the number of these DOFs is N. The format of *ddm* array is 2*N dofman1 idof1 dofman2 idof2 ... dofmanN idofN, where the dofmani is the number of i-th dof manager and idofi is the corresponding DOF number.
- *ddv* is array of relative weights of controlled displacements, the size should be equal to N. The actual value of prescribed dofs is defined as a product of its weight and the value of load time function specified using *ddltf* parameter (see below).
- *ddltf* number of load time function, which is used to evaluate the actual displacements of controlled dofs.
- *linesearch* nonzero value turns on line search algorithm. The *lsearchtol* defines tolerance (default value is 0.8), amplification factor can be specified using *lsearchamp* parameter (should be in interval (1,10)), and parameter *lsearchmaxeta* defines maximum limit on the length of iterative step (allowed range is (1.5,15)).

¹However, the problem does not supports the changes of static system. But it is possible to apply direct displacement control without requiring BC applied (see nrsolver documentation). Therefore it is possible to combine direct displacement control with direct load control or indirect control.

- *nccdg* allows to define one or more DOF groups, that are used for evaluation of convergence criteria. Each DOF is checked if it is a member of particular group and in this case its contribution is taken into account when evaluating the convergence criteria for that group. By default, if *nccdg* is not specified, one group containing all DOF types is created. The value of *nccdg* parameter defines the number of DOF type groups. For each group, the corresponding DOF types need to be specified using *ccdgi* parameter, where '#' should be replaced by group number (numbering starts from 1). This array contains the DofIDItem values, that identify the physical meaning of DOFs in the group. The values and their physical meaning is defined by DofIDItem enum type (see src/oofemlib/dofiditem.h for reference).
- *rtolv* determines relative convergence norm (both for displacement iterative change vector and for residual unbalanced force vector). Optionally, the *rtolf* and *rtold* parameters can be used to define independent relative convergence criteria for unbalanced forces and displacement iterative change. If the default convergence criteria is used, the parameters *rtolv*, *rtolf*, and *rtold* are real values. If the convergence criteria DOF groups are used (see below the description of *nccdg* parameter) then they should be specified as real valued arrays of *nccdg* size, and individual values define relative convergence criteria for each individual dof group.

The indirect solver corresponds to *controllmode*=0 and the CALM solver is used. The value of reference load vector is determined by *refloadmode* parameter mentioned above at the first step of each metastep. *However, the user must ensure, that the same value of reference load vector could be obtained for all solution steps of particular metastep (this is necessary for restart and adaptivity to work).* The corresponding meta step solver parameters (*solverParams*) are:

```

Psi(rn) # \
MaxIter(in) # \
stepLength(rn) # \
[minStepLength(in) #] \
[initialStepLength(rn) #] \
[forcedInitialStepLength(rn) #] \
[reqIterations(in) #] \
[minIter(in) #] \
[manrmsteps(in) #] \
[hpcmode(in) #] [hpc(ia) #] [hpcw(ia) #] \
[lineSearch(in) #] [lsearchamp(rn) #] \
[lsearchmaxeta(rn) #] [lsearchtol(rn) #] \
[nccdg(in) # ccdg1(ia) # ... ccdgN(ia) #] \
rtolv() # [rtolf() # rtold() # ]

```

where

- *Psi* - CALM Ψ control parameter. For $\Psi = 0$ displacement control is applied. For nonzero values the load control applies together with displacement control (ALM). For large Ψ load control apply.
- *MaxIter* - determines the maximum number of iteration allowed to reach equilibrium state. If this limit is reached, restart follows with smaller step length.

- *stepLength* - determines the maximum value of arc-length (step length).
- *minStepLength* - minimum step length. The step length will never be smaller. If convergence problems are encountered and step length cannot be decreased, computation terminates.
- *initialsteplength* - determines the initial step length (the arc-length). If not provided, the maximum step length (determined by *stepLength* parameter) will be used as the value of initial step length.
- *forcedInitialStepLength* - When simulation is restarted, the last predicted step length is used. Use *forcedInitialStepLength* parameter to override the value of step length. This parameter will also override the value of initial step length set by *initialsteplength* parameter.
- *reqIterations* - approximate number of iterations controlled by changing the step length.
- *minIter* - minimum number of iterations which always proceed during the iterative solution. *reqIterations* are set to be the same, *MaxIter* are increased if lower.
- *manrmsteps* - Forces the use of accelerated Newton Raphson method, where stiffness is updated after *manrmsteps* steps. By default, the modified NR method is used (no stiffness update).
- *hpcmode* Parameter determining the alm mode. Possible values are: 0 - (default) full ALM with quadratic constrain and all dofs, 1 - (default, if *hpc* parameter used) full ALM with quadratic constrain, taking into account only selected dofs (see *hpc* param), 2 - linearized constrain in displacements only, taking into account only selected dofs with given weight (see *hpc* and *hpcw* parameters).
- *hpc* - Special parameter for Hyper-plane control, when only selected DOFS are taken account in ALM step length condition. Important mainly for material nonlinear problems with strong localization. This array selects the degrees of freedom, which displacements are controlled. Let the number of these DOFs is N. The format of *ddm* array is 2*N dofman1 idof1 dofman2 idof2 ... dofmanN idofN, where the dofmani is the number of i-th dof manager and idofi is the corresponding DOF number.
- *hpcw* - Array of dof weights in linear constraint. The dof ordering is determined by *hpc* param, size of array should be N.
- *linsearch* nonzero value turns on line search algorithm. The *lsearchtol* defines tolerance, amplification factor can be specified using *lsearchamp* parameter (should be in interval (1,10)), and parameter *lsearchmaxeta* defines maximum limit on the length of iterative step (allowed range is (1.5,15)).
- *nccdg* allows to define one or more DOF groups, that are used for evaluation of convergence criteria. Each DOF is checked if it is a member of particular group and in this case its contribution is taken into account when evaluating the convergence criteria for that group. By default, if *nccdg* is

not specified, one group containing all DOF types is created. The value of *nccdg* parameter defines the number of DOF type groups. For each group, the corresponding DOF types need to be specified using *ccdg#* parameter, where '#' should be replaced by group number (numbering starts from 1). This array contains the DofIDItem values, that identify the physical meaning of DOFs in the group. The values and their physical meaning is defined by DofIDItem enum type (see src/oofemlib/dofiditem.h for reference).

- *rtolv* determines relative convergence norm (both for displacement iterative change vector and for residual unbalanced force vector). Optionally, the *rtolf* and *rtold* parameters can be used to define independent relative convergence criteria for unbalanced forces and displacement iterative change. If the default convergence criteria is used, the parameters *rtolv*, *rtolf*, and *rtold* are real values. If the convergence criteria DOF groups are used (see below the description of *nccdg* parameter) then they should be specified as real valued arrays of *nccdg* size, and individual values define relative convergence criteria for each individual dof group.

6.8.2 Standard syntax

In this case, all parameters (for analysis as well as for the solver) are supplied in analysis record. The default meta step is created for all solution steps required. Then the meta step attributes are specified within analysis record. The format of analysis record is then following

```
NonLinearStatic  nsteps(in) # \
                 [nonlocstiff(in) #] \
                 [contextOutputStep(in) #] \
                 [controllmode(in) #] \
                 [deltat(rn) #] \
                 rtolv(rn) # \
                 [stiffmode(in) #] \
                 lstype(in) # \
                 smtype(in) # \
                 solverParams() # \
                 [nonlinform(in) #] \
                 <[nonlocstiff(in) #]> \
                 <[nonlocalext() #]> \
                 <[loadbalancing() #]>
```

The meaning of parameters is the same as for extended syntax.

6.9 Adaptive linear static

```
Adaptlinearstatic  nsteps(in) # \
                  [renumber(in) #] \
                  [sparselinsolverparams() #] \
                  [meshpackage(in) #] \
                  errorestimatorparams() #
```

Adaptive linear static analysis. Multiple loading cases are not supported. Due to linearity of a problem, the complete reanalysis from the beginning is done after

adaptive remeshing. After first step the error is estimated, information about required density is generated (using mesher interface) and solution terminates. If the error criteria is not satisfied, then the new mesh and corresponding input file is generated and new analysis should be performed until the error is acceptable. Currently, the available error estimator for linear problems is Zienkiewicz-Zhu. Please note, that adaptive framework requires specific functionality provided by elements and material models. For details, see element and material model manuals.

- Parameter *nsteps* indicates the number of loading cases. Should be set to 1.
- Nonzero value of optional parameter *renumber* turns on the equation renumbering to optimize the profile of characteristic matrix (uses Sloan algorithm). By default, profile optimization is not performed. It will not work in parallel mode.
- The *sparselinsolverparams* parameter describes the sparse linear solver attributes and is explained in section 6.17.
- The *meshpackage* parameter selects the mesh package interface, which is used to generate information about required mesh density for new remeshing. The supported interfaces are explained in section 6.22. By default, the T3d interface is used.
- The *errorrestimatorparams* parameter contains the parameters of Zienkiewicz Zhu Error Estimator. These are described in section 6.20.

6.10 Adaptive nonlinear static

Adaptnlinearstatic	Nonlinearstaticparams ₍₎ #	\
	[equilmc _(in) #]	\
	[renumber _(in) #]	\
	[meshpackage _(in) #]	\
	[eetype ₍₎ #]	\
	errorestimatorparams ₍₎ #	

Represents Adaptive Non-LinearStatic problem. Solution is performed as a series of increments (loading or displacement). The error is estimated at the end of each load increment (after equilibrium is reached), and based on reached error, the computation continues, or the new mesh densities are generated and solution stops. Then the new discretization should be generated. The truly adaptive approach is supported, so the computation can be restarted from the last step (see section 2), solution is mapped to new mesh (separate solution step) and new load increment is applied. Of course, one can start the analysis from the very beginning using new mesh. Currently, the available estimators/indicators include only linear Zienkiewicz-Zhu estimator and scalar error indicator. Please note, that adaptive framework requires specific functionality provided by elements and material models. For details, see element and material model manuals.

- Set of parameters *Nonlinearstaticparams* are related to nonlinear analysis. They are described in section 6.8.

- Parameter *equilmc* determines, whether after mapping of primary and internal variables to new mesh the equilibrium is restored or not before new load increment is applied. The possible values are: 0 (default), when no equilibrium is restored, and 1 forcing the equilibrium to be restored before applying new step.
- The *meshpackage* parameter selects the mesh package interface, which is used to generate information about required mesh density for new remeshing. The supported interfaces are explained in section 6.22. By default, the T3d interface is used.
- Parameter *eetype* determines the type of error estimator/indicator to be used. The parameters *errorestimatorparams* represent set of parameters corresponding to selected error estimator. For description, follow to section 6.20.

6.11 Stationary transport problem

StationaryProblem	nsteps _(in) #	\
	[renumber _(in) #]	\
	[sparselinsolverparams ₍₎ #]	\
	[exportfields ₍₎ #]	\
	[atomicfields _(in) #]	\

Stationary transport problem. Series of loading cases is maintained as sequence of time-steps. For each load case an auxiliary time-step is generated with time equal to load case number. Load vectors for each load case are formed as load vectors at this auxiliary time. The *sparselinsolverparams* parameter describes the sparse linear solver attributes and is explained in section 6.17. Nonzero value of optional parameter *renumber* turns on the equation renumbering to optimize the profile of characteristic matrix (uses Sloan algorithm). By default, profile optimization is not performed. It will not work in parallel mode.

If the present problem is used within the context of staggered-like analysis, the temperature field obtained by the solution can be exported and made available to any subsequent analyses. For example, temperature field obtained by present analysis can be taken into account in subsequent mechanical analysis. To allow this, the temperate must be “exported”. This can be done by adding *exportfields* parameter, followed by *atomicfields* array. This array contains the atomic field identifiers, which tell the problem to register its primary unknowns under given atomic identifiers. Currently available atomic ids are: Temperature (corresponding value is 1). Then the subsequent analyses can get access to exported fields (via atomic id) and take them into account, if they support such feature.

6.12 Transient transport problem - linear case

NonStationaryProblem	nsteps _(in) #	\
	[renumber _(in) #]	\
	deltaT _(rn) #	\
	alpha _(rn) #	\
	[lumpedcapa ₍₎ #]	\
	[sparselinsolverparams ₍₎ #]	\
	[exportfields ₍₎ #]	\
	[atomicfields _(in) #]	\

Linear implicit integration scheme for transient transport problems. The generalized midpoint rule (sometimes called α -method) is used for time discretization, with alpha parameter, which has limits $0 \leq \alpha \leq 1$. For $\alpha = 0$ explicit Euler forward method is obtained, for $\alpha = 0.5$ implicit trapezoidal rule is recovered, which is unconditionally stable, second-order accurate in Δt , and $\alpha = 1.0$ yields implicit Euler backward method, which is unconditionally stable, and first-order accurate in Δt . *deltaT* is time step length used for integration, *nsteps* parameter specifies number of time steps to be solved. Nonzero value of optional parameter *renumber* turns on the equation renumbering to optimize the profile of characteristic matrix (uses Sloan algorithm). By default, profile optimization is not performed. It will not work in parallel mode. If *lumpedcapa* is set, then the stabilization of numerical algorithm using lumped capacity matrix will be used, reducing the initial oscillations. See section 6.17 for explanation of *exportfields* and *atomicfields* parameters. Note: This problem type **requires transport module** and it can be used only when this module is included in your oofem configuration.

Note: This problem type **requires transport module** and it can be used only when this module is included in your oofem configuration.

6.13 Transient transport problem - nonlinear case

NITransientTransportProblem	nsteps _(in) #	\
	[renumber _(in) #]	\
	deltaT _(rn) #	\
	alpha _(rn) #	\
	[lumpedcapa ₍₎ #]	\
	[nsmax _(in) #]	\
	rtol _(rn) #	\
	[manrmsteps _(in) #]	\
	[sparselinsolverparams ₍₎ #]	\
	[exportfields ₍₎ #]	\
	[atomicfields _(in) #]	\

Implicit integration scheme for transient transport problems. The generalized midpoint rule (sometimes called α -method) is used for time discretization, with alpha parameter, which has limits $0 \leq \alpha \leq 1$. For $\alpha = 0$ explicit Euler forward method is obtained, for $\alpha = 0.5$ implicit trapezoidal rule is recovered, which is unconditionally stable, second-order accurate in Δt , and $\alpha = 1.0$ yields implicit Euler backward method, which is unconditionally stable, and first-order accurate in Δt . *deltaT* is time step length used for integration, *nsteps* parame-

ter specifies number of time steps to be solved. Parameter *maxiter* determines the maximum number of iterations allowed to reach equilibrium (default is 30). The convergence is reached, when norms of both residual fluxes and iterative change of solution vector is less than the value given by *rtol*. If *manrmsteps* parameter is nonzero, then the modified N-R scheme is used, with the left-hand side matrix updated after *manrmsteps* steps. If *lumpedcapa* is set, then the stabilization of numerical algorithm using lumped capacity matrix will be used, reducing the initial oscillations. Nonzero value of optional parameter *renumber* turns on the equation renumbering to optimize the profile of characteristic matrix (uses Sloan algorithm). By default, profile optimization is not performed. It will not work in parallel mode. See section 6.17 for explanation of *exportfields* and *atomicfields* parameters.

Note: This problem type **requires transport module** and it can be used only when this module is included in your oofem configuration.

6.14 Transient incompressible flow - CBS Algorithm

CBS	nsteps _(in) # [renumber _(in) #] deltaT ₀ # [theta1 _(in) #] [theta2 _(in) #] [cmflag _(in) #] [scaleflag _(in) # lscale _(in) # uscale _(in) # dscale _(in) #] [lstype _(in) #] [smttype _(in) #]	\ \ \ \ \ \
-----	---	----------------------------

Solves the transient incompressible flow using algorithm based on Characteristics Based Split (CBS, for reference see O.C.Zienkiewics and R.L.Taylor: The Finite Element Method, 3rd volume, Butterworth-Heinemann, 2000). At present, only semi-implicit form of the algorithm is available and energy equation, yielding the temperature field, is not solved. Parameter *nsteps* determines number of solution steps. Parameter *deltaT* is time step length used for integration. This time step will be automatically adjusted to satisfy integration stability limits $\Delta t \leq \frac{h}{|\mathbf{u}|}$ and $\Delta t \leq \frac{h^2}{2\nu}$, if necessary. Parameters *theta1* and *theta2* are integration constants, $\theta_1, \theta_2 \in < \frac{1}{2}, 1 >$. If *cmflag* is given a nonzero value, then consistent mass matrix will be used instead of (default) lumped one.

The characteristic equations can be solved in non-dimensional form. To enable this, the *scaleflag* should have a nonzero value, and the following parameters should be provided: *lscale*, *uscale*, and *dscale* representing typical length, velocity, and density scales.

Parameter *lstype* allows to select solver for linear system of equations. Parameter *smttype* allows to select sparse matrix storage scheme. The scheme should be compatible with solver type. See section 6.17 for further details.

Nonzero value of optional parameter *renumber* turns on the equation renumbering to optimize the profile of characteristic matrix (uses Sloan algorithm). By default, profile optimization is not performed. It will not work in parallel mode.

6.15 Transient incompressible flow SUPG/PSPG Algorithm

SUPG	nsteps _(in) #	\
	[renumber _(in) #]	\
	deltaT _(rn) #	\
	rtolv _(rn) #	\
	[atolv _(rn) #]	\
	[stopmaxiter _(in) #]	\
	[alpha _(rn) #]	\
	[cmflag _(in) #]	\
	[deltatlft _(in) #]	\
	[miflag _(in) #]	\
	[scaleflag _(in) # lscale _(in) # uscale _(in) # dscale _(in) #]	\
	[lstype _(in) #] [smtpe _(in) #]	\

Solves the transient incompressible flow using stabilized formulation based on SUPG and PSPG stabilization terms. The stabilization provides stability and accuracy in the solution of advection-dominated problems and permits usage of equal-order interpolation functions for velocity and pressure. Furthermore, stabilized formulation significantly improves convergence rate in iterative solution of large nonlinear systems of equations.

By changing the value α , different methods from “Generalized mid-point family” can be chosen, i.e., Forward Euler ($\alpha = 0$), Midpoint rule ($\alpha = 0.5$), Galerkin ($\alpha = 2/3$), and Backward Euler ($\alpha = 1$). Except the first one, all the methods are implicit and require matrix inversion for solution. Some results from an energy method analysis suggest unconditional stability for $\alpha \geq 0.5$ for the generalized mid-point family. As far as accuracy is concerned, the midpoint rule is to be generally preferred.

Parameter *nsteps* determines number of solution steps. Parameter *deltaT* is time step length used for integration. Alternatively, the load time function can be used to determine time step length for particular solution step. The load time function number is determined by parameter *deltatlft* and its value evaluated for solution step number should yield the step length.

Nonzero value of optional parameter *renumber* turns on the equation renumbering to optimize the profile of characteristic matrix (uses Sloan algorithm). By default, profile optimization is not performed. It will not work in parallel mode.

Parameters *rtolv* and *atolv* allow to specify relative and absolute errors norms for residual vector. The equilibrium iteration process will be stopped when both error limits are satisfied or when the number of iteration exceeds the value given by parameter *stopmaxiter*.

If *cmflag* is given a nonzero value, then consistent mass matrix will be used instead of (default) lumped one.

The algorithm allows to solve the flow of two immiscible fluids in fixed spatial domain (currently only in 2d). This can be also used for solving free surface problems, where one of the fluids should represent air. To enable multi-fluid analysis, user should set parameter *miflag*. The supported values are described in section 6.21. Please note, that the initial distribution of reference fluid volume should be provided as well as constitutive models for both fluids.

The characteristic equations can be solved in non-dimensional form. To enable this, the *scaleflag* should have a nonzero value, and the following parameters should be provided: *lscale*, *uscale*, and *dscale* representing typical length, velocity, and density scales.

Parameter *lstype* allows to select solver for linear system of equations. Parameter *smttype* allows to select sparse matrix storage scheme. Please note, that the present algorithm leads to a non-symmetrical system matrix. The scheme should be compatible with solver type. See section 6.17 for further details.

6.16 Staggered Problem

StaggeredProblem	nsteps _(in) #	\
	deltaT ₍₎ #	\
	prob1 _(s) # prob2 _(s) #	

Represent so-called staggered analysis. This can be described as an sequence of sub-problems, where the result of some sub-problem in the sequence can depend on results of previous sub-problems in sequence. Typical example is heat transfer analysis followed by mechanical analysis taking into account the temperature field generated by the heat transfer analysis. The actual implementation supports only sequence of two sub-problems. The sub-problems are described using sub-problem input files. The syntax of sub-problem input file is the same as for standalone problem. The only addition is that sub-problems should export their solution fields so that they became available for subsequent sub-problems. The subproblem input files are described using *prob1* and *prob2* parameters, which are strings containing a path to sub-problem input files, the *prob1* contains input file path of the first sub-problem, which runs first for each solution step, the *prob2* contains input file path of the second sub-problem. The solution steps are generated by the master-problem (represented by Staggered-Problem). Therefore any sub-problem time-stepping parameters are ignored (even if they are required by sub-problem input syntax) and only master problem parameters are relevant. *deltaT* is time step length used for integration, *nsteps* parameter specifies number of time steps to be solved.

Note: This problem type **is included in transport module** and it can be used only when this module is configured. Note: All material models derived from StructuralMaterial base will take into account the external registered temperature field, if provided.

6.17 Sparse linear solver parameters

The sparselinsolverparams field has the following general syntax:

[lstype ^M _(in) #]		\
[smttype _(in) #]		\
solverParams ^M _(string) #		

where parameter *lstype* allows to select solver for linear system of equations. Currently supported values are 0 (default) for direct solver (ST_Direct), 1 for Iterative Method Library (IML) solver (ST_IML), 2 for Spooles direct solver, 3 for Petsc library family of solvers, and 4 for DirectSparseSolver (ST_DSS). Parameter *smttype* allows to select sparse matrix storage scheme. The scheme should be compatible with solver type. Currently supported values (marked as

“id”) are summarized in table (1). The 0 value is default and selects the symmetric skyline (SMT_Skyline). Ther possible storage formats include unsymmetric skyline (SMT_SkylineU), compressed column (SMT_CompCol), dynamically growing compressed column (SMT_DynCompCol), symmetric compressed column (SMT_SymCompCol), spooles library storage format (SMT_SpoolesMtrx), PETSc library matrix representation (SMT_PetscMtrx, a sparse serial/parallel matrix in AIJ format), and DSS compatible matrix representations (SMT_DSS_*). The allowed *lstype* and *smttype* combinations are summarized in the table (1), together with solver parameters related to specific solver.

Storage format	id	Sparse solver				
		Direct	IML	Spooles	Petsc	DSS
SMT_Skyline	0	+	+			
SMT_SkylineU	1	+	+			
SMT_CompCol	2		+			
SMT_DynCompCol	3		+			
SMT_SymCompCol	4		+			
SMT_DynCompRow	5		+			
SMT_SpoolesMtrx	6			+		
SMT_PetscMtrx	7				+	
SMT_DSS_sym_LDL	8					+
SMT_DSS_sym_LL	9					+
SMT_DSS_unsym_LU	10					+

Table 1: Solver and stoorage scheme compatibility.

The solver parameters in *solverParams* depend on the solver type and are summarized in table (2).

Solver type	id	Solver parameters
ST_Direct	0	
ST_IML	1	[stype _(in) #] [lstol _(rn) #] [lster _(in) #] [lsprecond _(in) #] [precondattributes _(string) #]
ST_Spooles	2	[msglvl _(in) #] [msgfile _(s) #]
ST_Petsc	3	see Petsc manual, for details ²
ST_DSS	4	

Table 2: Solver parameters.

The *stype* allows to select particular iterative solver from IML library, currently supported values are 0 (default) for Conjugate-Gradient solver, 1 for GMRES solver. Parameter *lstol* represents the maximum value of residual after the final iteration and the *lster* is maximum number of iteration for iterative solver. The *precondattributes* parameters contains the optional preconditioner parameters. The *lsprecond* parameter determines the type of preconditioner

²User can set several run-time options, e.g., -ksp_type [cg, gmres, bicg, bcgs] -pc_type [jacobi, bjacobi, none, ilu, ...] -ksp_monitor -ksp_rtol <rtol> -ksp_view -ksp_converged_reason. These options will override those that are default (PETSC KSPSetFromOptions() routine is called after any other customization routines).

to be used. The possible values of *lsprecond* together with supported storage schemes and their descriptions are summarized in table (3).

Precond type	id	Compatible storage	Description and parameters
IML_VoidPrec	0	all	No preconditioning
IML_DiagPrec	1	all	Diagonal preconditioning
IML_ILUPrec	2	SMT_CompCol SMT_DynCompCol	Incomplete LU Decomposition with no fill up
IML_ILUPrec	3	SMT_DynCompRow	Incomplete LU (ILUT) with fillup. The <i>precondattributes</i> are: [droptol _(rn) #] [partfill _(in) #]. <i>droptol</i> dropping tolerance <i>partfill</i> level of fill-up
IML_ICPrec	4	SMT_SymCompCol SMT_CompCol	Incomplete Cholesky with no fill up

Table 3: Preconditioning summary.

6.18 Eigen value solvers

The eigensolverparams field has the following general syntax:

$$\left[\begin{array}{l} \text{stype}^{M_{(in)}} \# \\ \text{smttype}_{(in)} \# \\ \text{solverParams}^{M_{(string)}} \# \end{array} \right. \quad \backslash$$

where parameter *stype* allows to select solver type. Parameter *smttype* allows to select sparse matrix storage scheme. The scheme should be compatible with solver type. Currently supported values of *stype* are summarized in table 4.

Solver type	stype id	solver parameters
Subspace Iteration	0 (default)	
Inverse Iteration	1	
SLEPc solver	2	requires “smttype 7” see also SLEPc manual

Table 4: Eigen Solver parameters.

6.19 Dynamic load balancing parameters

There are in general two basic factors causing load imbalance between individual subdomains: (i) one coming from application nature, such as switching from linear to nonlinear response in certain regions or local adaptive refinement, and (ii) external factors, caused by resource reallocation, typical for nondedicated cluster environments, where individual processors are shared by different applications and users, leading to time variation in allocated processing power. The load balance recovery is achieved by repartitioning of the problem domain and transferring the work (represented typically by finite elements) from one

subdomain to another. This section describes the structure and syntax of parameters related to dynamic load balancing. The corresponding part of analysis record has the following general syntax:

[lbflag ^M _(in) #]	\
[forcelb1 _(in) #]	\
[wtp _(ia) #]	\
[lbstep _(in) #]	\
[relwct _(rn) #]	\
[abswct _(rn) #]	\
[minwct _(rn) #]	\

where the parameters have following meaning:

- *lbflag*, when set to nonzero value activates the dynamic load balancing. Default value is zero.
- *forcelb1* forces the load rebalancing after the first solution step, when set to nonzero value.
- *wtp* allows to activate optional load balancing plugins. At present, the only supported value is 1, that activates nonlocal plugin, necessary for nonlocal averaging to work properly when dynamic load balancing is active.
- *lbstep* rebalancing, if needed, is performed only every *lbstep* solution step. Default value is 1 (recover balance after every step, if necessary).
- *relwcr* sets relative wall-clock imbalance treshold. When achieved relative imbalance between wall clock solution time of individual processors is greater than provided treshold, the rebalancing procedure will be activated.
- *abswct* sets absolute wall-clock imbalance treshold. When achieved absolute imbalance between wall clock solution time of individual processors is greater than provided treshold, the rebalancing procedure will be activated.
- *minwct* minimum absolute imbalance to perform relative imbalance check using *relwcr* parameter, otherwise only absolute check is done. Default value is 0.

At present, the load balancing support requires ParMETIS module to be configured and compiled.

6.20 Error estimators and indicators

The currently supported values of *eetype* are in table 5.

- EET_SEI - Represents scalar error indicator. It indicates element error based on the value of some suitable scalar value (for example damage level, plastic strain level) obtained from the element integration points and corresponding material model.
- EET_ZZEE - The implementation of Zienkiewicz Zhu Error Estimator. It requires the special element algorithms, which may not be available for all element types.

- EET_CZZSI - The implementation of combined criteria: Zienkiewicz Zhu Error Estimator for elastic regime and scalar error indicator in non-linear regime.

Error estimator/indicator	<i>eetype</i>
EET_SEI	0
EET_ZZEE	1
EET_CZZSI	2

Table 5: Supported error estimators and indicators.

The sets of parameters (*errorestimatorparams* field) used to configure each error estimator are different

- EET_SEI


```

[regionskipmap(ia) #]
vartype(in) #
minlim(rn) # maxlim(rn) #
mindens(rn) # maxdens(rn) # defdens(rn) #
[remeshingdensityratio(rn) #]
```

```

\
\
\
\
```

 - *regionskipmap* parameter allows to skip some regions. The error is not evaluated in these regions and default mesh density is used. The size of this array should be equal to number of regions and nonzero entry indicates region to skip.
 - *vartype* parameter determines the type of internal variable to be used as error indicator. Currently supported value is 1, representing damage based indicator.
 - If the indicator value is in range given by parameters (*minlim*, *maxlim*) then the proposed mesh density is linearly interpolated within range given by parameters (*mindens*, *maxdens*). If indicator value is less than value of *minlim* parameter then value of *defdens* parameter is used as required density, if it is larger than *maxlim* then *maxdens* is used as required density.
 - *remeshingdensityratio* parameter determines the allowed ratio between proposed density and actual density. The remeshing is forced, whenever the actual ratio is smaller than this value. Default value is equal to 0.80.
- EET_ZZEE


```

[regionskipmap(ia) #]
normtype(in) #
requirederror(rn) #
minelemsize(rn) #
```

```

\
\
\
```

 - *regionskipmap* parameter allows to skip some regions. The error is not evaluated in these regions and default mesh density is used. The size of this array should be equal to number of regions and nonzero entry indicates region to skip.

- *normtype* Allows select the type of norm used in evaluation of error. Default value is to use L2 norm (equal to 0), value equal to 1 uses the energy norm.
- *requirederror* parameter determines the required error to obtain (in percents/100).
- *minelemsize* parameter allows to set minimum limit on element size.
- EET_CZZSI - combination of parameters for EET_SEI and EET_ZZEE; the in elastic regions are driven using EET_SEI, the elastic are driven by EET_ZZEE.

6.21 Material interfaces

The material interfaces are used to represent and track the position of various interfaces on fixed grids. Typical examples include free surface, evolving interface between two materials, etc. Available representations include:

MI	miflag	Compatibility
LEPlic	0	2D triangular
LevelSet	1	2D triangular

- LEPlic- representation based on Volume-Of-Fluid approach; the initial distribution of VOF fractions should be specified for each element (see element manual)
 - | **[refvol_(rn) #]**
- parameter *refvol* allows to set initial volume of reference fluid, then the reference volume is computed in each step and printed, so the accuracy and mass conservation can be monitored.

LevelSet- level set based representation

| **[levelset_(ra) # OR refmatpolyx_(ra) # refmatpolyy_(ra) #]** \

| **[lsra_(in) #] [rdt_(rn) #] [rerr_(rn) #]**

- *levelset* allows to specify the initial level set values for all nodes directly. The size should be equal to total number of nodes within the domain.
- Parameters *refmatpolyx* and *refmatpolyy* allow to initialize level set by specifying interface geometry as 2d polygon. Then polygon describes the initial zero level set, and level set values are then defined as signed distance from this polygon. Positive values are on the left side when walking along polygon. The parameter *refmatpolyx* specifies the x-coordinates of polygon vertices, parameter *refmatpolyy* y-coordinates. Please note, that level set must be initialized, either using *levelset* parameter or using *refmatpolyx* and *refmatpolyy*.
- Parameter *lsra* allows to select level set reinitialization algorithm. Currently supported values are 0 (re-initializes the level set representation by solving $d_\tau = S(\phi)(1 - |\nabla d|)$ to steady state, default), 1 (uses fast marching method to build signed distance level set representation).

- Parameters *rdt* *rerr* are used to control reinitialization algorithm for *lsra* = 0. *rdt* allows to change time step of integration algorithm and parameter *rerr* allows to change default error limit used to detect steady state.

6.22 Mesh generator interfaces

The mesh generator interface is responsible to provide a link to specific mesh generator. The supported values of *meshpackage* parameter are

- MPT_T3D: *meshpackage* = 0. T3d mesh interface. Default. Supports both 1d, 2d (triangles) and 3d (tetrahedras) meshes. Reliable.
- MPT_TARGE2: *meshpackage* = 1. Interface to Targe2 2D mesh generator.

6.23 Initialization modules

Initialization modules allow to initialize the state variables using data previously computed by external software. The number of initialization module records is specified in analysis record using *ninitmodules* parameter (see the initial part of section 6). The general format is the following:

```
| *EntType initfile(string) #
```

The file name following the keyword “initfile” specifies the path to the file that contains the initialization data and should be given without quotes.

Currently, the only supported initialization module is

- Gauss point initialization module

```
| GPInitModule initfile(string) #
```

- Each Gauss point is represented by one line in the initialization file.
- The Gauss points should be given in a specific order, based on the element number and the Gauss point number, in agreement with the mesh specified in later sections.

- Each line referring to a Gauss point should contain the following data:

```
| elnum(in) # gpnum(in) # coords(ra) # ng(in) #
| var_1_id(in) # values_1(ra) #
| ...
| var_ng_id(in) # values_ng(ra) #
```

\\
\\
\\

- *elnum* is the element number
- *gpnum* is the Gauss point number
- *coords* are the coordinates of the Gauss point
- *ng* is the number of groups of variables that will follow
- *var_1_id* is the identification number of variable group number 1 (according to the definitions in *internalstatetype.h*)
- *values_1* are the values of variables in group number 1
- *var_ng_id* is the identification number of variable group number ng

- *values_ng* are the values of variables in group number *ng*
- Example:
“37 4 3 0.02 0.04 0.05 3 52 1 0.23 62 1 0.049 1 6 0 -2.08e+07 0 0 0 0”
means that Gauss point number 4 of element number 37 has coordinates $x = 0.02$, $y = 0.04$ and $z = 0.05$ and the initial values are specified for 3 groups of variables;
the first group (variable ID 52) is of type `IST_DamageScalar` (see `internalstatetype.h`) and contains 1 variable (since it is a scalar) with value 0.23;
the second group (ID 62) is of type `IST_CumPlasticStrain` and contains 1 variable with value 0.049;
the third group is of type `IST_StressTensor` and contains 6 variables (stress components σ_x , σ_y , etc.) with values 0, -2.08e+07, 0, 0, 0, 0

6.24 Export modules

Export modules allow to export computed data into external software for post-processing. The number of export module records is specified in analysis record using *nmodules* parameter (see the initial part of section 6). The general format is the following:

*EntType	[tstep_all]	\
	[tstep_step _(in) #]	\
	[tsteps_out _(r1) #]	\
	[domain_all]	\
	[domain_mask _(in) #]	

To select all solution steps, in which output will be performed, use *tstep_all*. To select each *tstep_step*-nth step, use *tstep_step* parameter. In order to select only specific solution steps, the *tsteps_out* list can be specified, supplying solution step number list in which output will be done. To select output for all domain of the problem the *domain_all* keyword can be used. To select only specific domains, *domain_mask* array can be used, where the values of the array specify the domain numbers to be exported. Currently, the supported export modules are following

- VTK export

vtk	[vars _(ia) #]	\
	[primvars _(ia) #]	\
	[cellvars _(ia) #]	\
	[stype _(in) #]	\
	[regionstoskip _(ia) #]	

- The array *vars* contains identifiers for those internal variables which are to be exported. The id values are defined by `InternalStateType` enumeration, which is defined in include file “src/oofemlib/internal-statetype.h”.
- The array *primvars* contains identifiers of primary variables to be exported. The possible values correspond to the values of enumerated type `UnknownType`, which is again defined in “src/oofemlib/unknown-type.h”. Please note, that the values corresponding to enumerated

type values start from zero, if not specified directly and that not all values are supported by particular material model or analysis type.

- The array *cellvars* contains identifiers of constant variables defined on an element (cell), e.g. a material number. Identifier numbers are specified in “src/oofemlib/internalstatetype.h”.
- The parameter *stype* allows to select smoothing procedure for internal variables, which is used to compute nodal values from values in integration points.

When vtk is configured to operate in region-by-region mode (the default), then the export is done on region by region basis. In this case, the smoothing is performed only over particular region, where only elements in this region participate. Then the *regionstoskip* parameter allows to skip given regions when generating output, so only remaining regions (and corresponding elements and nodes) will be exported.

- Homogenization of stresses and strains in the global coordinate system. Corresponding stress and strain components are summed and averaged over the volume. It is possible to select material numbers from which the averaging occurs. The averaging works for 3D domains with an extension to trusses. A truss is considered as a volume element with oriented stress and strain components along the truss axis. The transformation to global components occurs before averaging.

hom	[scale _(rn) #]	\
	[MatNum _(ia) #]	

- The parameter *scale* multiplies all averaged stresses and strains. *scale*=1 by default.
- An integer array *MatNum* specifies which material numbers are taken into account. All material numbers are averaged by default.
- Gauss point export is useful if one needs to plot a certain variable (such as damage) as a function of a spatial coordinate using tools like gnuplot. It generates files with data organized in columns, each row representing one Gauss point. In this way, one can plot e.g. the damage distribution along a one-dimensional bar.

gpexportmodule	[vars _(ia) #]	\
	[ncoords _(in) #]	

- The array *vars* contains identifiers for those internal variables which are to be exported. The id values are defined by InternalStateType enumeration, which is defined in include file “src/oofemlib/internal-statetype.h”.
- Parameter *ncoords* specifies the number of spatial coordinates to be exported at each Gauss point. Depending on the spatial dimension of the domain, the points can have one, two or three coordinates. If *ncoords* is set to -1, only those coordinates that are actually used are exported. If *ncoords* is set to 0, no coordinates are exported. If *ncoords* is set to a positive integer, exactly *ncoords* coordinates are

exported. If *ncoords* exceeds the actual number of coordinates, the actual coordinates are supplemented by zeros. For instance, if we deal with a 2D problem, the actual number of coordinates is 2. For *ncoords*=3, the two actual coordinates followed by 0 will be exported. For *ncoords*=1, only the first coordinate will be exported.

The Gauss point export module creates a file with extension “gp” after each step for which the output is performed. This file contains a header with lines starting by the symbol #, followed by the actual data section. Each data line corresponds to one Gauss point and contains the following data:

1. element number,
2. material number,
3. Gauss point number,
4. contributing volume around Gauss point,
5. Gauss point global coordinates (written as a real array of length *ncoords*),
6. internal variables according to the specification in *vars* (each written as a real array of the corresponding length).

Example:

```
“GPExportModule 1 tstep_step 100 domain_all ncoords 2 vars 5 4 13 31
64 65”
```

means that the *.gp file will be written after each 100 steps and will contain for each of the Gauss points in the entire domain its 2 coordinates and also internal variables of type 4, 13, 31, 64 and 65, which are the strain tensor, damage tensor, maximum equivalent strain level, stress work density and dissipated work density. Of course, the material model must be able to deliver such variables. The size of the strain tensor depends on the spatial dimension, and the size of the damage tensor depends on the spatial dimension and type of model (e.g., for a simple isotropic damage model it will have just 1 component while for an anisotropic damage model it may have more). The other variables in this example are scalars, but they will be written as arrays of length 1, so the actual value will always be preceded by “1” as the length of the array. Since certain internal variables have the meaning of densities (per unit volume or area, again depending on the spatial dimension), it is useful to have access to the contributing volume of the Gauss point. The product of this contributing volume and the density gives an additive contribution to the total value of the corresponding variable. This can be exploited e.g. to evaluate the total dissipated energy over the entire domain.

7 Domain record(s)

This set of records describes the whole domain and its type. Depending on the type of problem, there may be one or several domain records. If not indicated, one domain record is default for all problem types.

The domain type is used to resolve the default number of DOFs in node and their physical meaning. Format is following

| **domain** ***domainType**

The ***domainType** can be one from the following

- The **2dPlaneStress** and **2d-Truss** modes declare two default dofs per node (u-displacement, v-displacement),
- The **3d** mode declares three default dofs per node (u-displacement, v-displacement, w-displacement),
- The **2dMindlinPlate** mode declares three default dofs per node (w-displacement, u-rotation, v-rotation).
- The **3dShell** mode declares six default dofs per node (displacement and rotation along each axis).
- The **2dBeam** mode declares three default dofs per node (u-displacement, w-displacement, v-rotation).
- The **2dIncompFlow** mode declares three default dofs per node (u-velocity, v-velocity, and pressure). The default number of dofs per node as well as their physical meaning can be overloaded in particular dof manager record (see section 7.3).

The further records describe particular domain components - OutputManagers, DofManagers, Elements, CrossSection models, Material Models, Boundary and Initial Conditions and Load time functions.

7.1 Output manager record

The output manager controls output. It can filter output to specific solution steps, and within these selected steps allows also to filter output only to specific dof managers and elements. The format of output manager record is

OutputManager	[tstep_all]	\
	[tstep_step _(in) #]	\
	[tsteps_out _(r1) #]	\
	[dofman_all]	\
	[dofman_output _(r1) #]	\
	[dofman_except _(r1) #]	\
	[element_all]	\
	[element_output _(r1) #]	\
	[element_except _(r1) #]	\

To select all solution steps, in which output will be performed, use *tstep_all*. To select each *tstep_step*-nth step, use *tstep_step* parameter. In order to select only specific solution steps, the *tsteps_out* list can be specified, supplying solution step number list in which output will be done. The combination of *tstep_step* and *tsteps_out* parameters is allowed.

Output manager allows also to filter output to only specific dof managers and elements. If these specific members are selected, the output happens only in selected solution steps. The *dofman_all* and *element_all* parameters select all dof managers or elements respectively. Parameter arrays *dofman_output* and

element_output allow to select only specific members. Numbers of selected members are then contained in *dofman_output* or *element_output* lists respectively. The previously selected members can be explicitly de-selected by specifying their component numbers in *dofman_except* or *element_except* lists.

7.2 Components size record

This record describes the number of components in related domain. The particular records will follow immediately in input file. The general format is:

```

ndofman(in) # \
nelem(in) # \
ncrosssect(in) # \
nmat(in) # \
nbc(in) # \
nic(in) # \
nltf(in) # \
[nbarrier(in) #]

```

where *ndofman* represents number of dof managers (e.g. nodes) and their associated records, *nelem* represents number of elements and their associated records, *ncrosssect* is number of cross sections and their records, *nmat* is number of material models and their records, *nbc* represents number of boundary conditions (including loads) and their records, *nic* parameter determines the number of initial conditions, and *nltf* represents number of time functions and their associated records. The optional parameter *nbarrier* represents the number of nonlocal barriers and their records. If not specified, no barriers are assumed.

7.3 Dof manager records

These records describe individual DofManager records (i.e. nodes or element sides (if they manage some DOFs)). The general format is following:

```

*DofManagerType (num#)(in) \
                  [load(ra) #] \
                  [ndofs(in) #] \
                  DofIDMask(ia) #] \
                  [bc(ia) #] \
                  [ic(ia) #] \
                  [doftype(ia) #] \
                  masterMask(ia) #] \
                  <[shared() #]> <[remote() #]> <[null() #]> \
                  <[partitions(ia) #]>

```

The order of particular records is optional, the dof manager number is determined by (**num#**)₍₎ parameter. The numbering of individual dof managers is arbitrary, it could be even non-continuous. In this context, one could think of dof manager number as a label that is assigned to individual dof manager and by which the dof manager is referenced. In parallel mode, the label represents a global id across all partitions.

The applied primary (Dirichlet) boundary conditions are specified using "bc" record, while natural boundary conditions using "load" parameter.

- The size of "bc" array (primary bc) should be equal to number of DOFs in dof manager and i-th value relates to i-th DOF - the ordering and physical

meaning of DOFs is determined by domain record and can be optionally specified for each dof manager individually (see next paragraph). The values of this array are corresponding boundary condition record numbers or zero, if no primary bc is applied to corresponding DOF. The compatible boundary condition type are required: primary conditions require "BoundaryCondition" records.

- The load "array" contains record numbers of natural boundary conditions that are applied. The required record type for natural condition is "Nodal-Load". The actual value is the summation of all contributions, if more than one natural bc is applied. See section on boundary conditions for the syntax. Please note, that the values of natural bc for individual DOFs are specified in its record, not in dofmanager record.

By default, if "bc" and/or "load" parameters are omitted, no primary and/or natural bc are applied. Analogously, initial conditions are represented using *ic* array. The size of *ic* array should be equal to number of DOFs in dof manager. The values of this array are corresponding initial condition record numbers or zero, if no initial condition is applied to corresponding DOF (in this case zero value is assumed as value of initial condition).

By default, the number of DOFs per node or side and their physical meanings are determined by domain record (***domainType** keyword). If it is necessary to have different number of DOFs, then the *ndofs* field determines number of DOFs in DofManager, and their physical meaning is determined by *DofIDMask* array. Size of this array should be equal to *ndofs* value. Each item of *DofIDMask* array describes the physical meaning of corresponding DOF in dof manager. Currently the following values are supported: {u-displacement=1, v-displacement=2, w-displacement=3, u-rotation=4, v-rotation=5, w-rotation=6, u-velocity=7, v-velocity=8, w-velocity=, temperature=7}. **It is not allowed to have two DOFs with the same physical meaning in the same DofManager.**

Parameters *dofType* and *masterMask* allows to connect some dof manager's dofs (so-called "slave" dofs) to corresponding dof (according to their physical meaning) of another dof manager (so-called "master" dof). The master slave principle allows for example simple modeling of structure hinges, where multiple elements are connected by introducing multiple nodes (with same coordinates) sharing the same displacement dofs and each one possessing their own rotational dofs. Parameter *dofType* determines the type of (slave) dof to create. Currently supported values are 0 for master DOF, 1 for simpleSlave DOF (linked to another single master DOF), and 2 for general slave dof, that can depend on different DOFs belonging to different dof managers. If *dofType* is not specified, then by default all DOFs are created as master DOFs. If provided, *masterMask* is also required. The meaning of *masterMask* parameter is depending on type of particular dofManager, and will be described in corresponding sections.

The *shared* indicates, that dofmanager is shared by neighboring partitions. The contributions from all contributing domains are summed. Typical for node cut algorithm (see figures 6 and 7).

Remote DofManager is indicated by *remote* parameter. Then DofManager in active domain is only mirror of some remote DofManager and it is necessary to copy remote values into local ones. Typical for element cut (see fig. 9). The *null* parameter indicates so-called null DofManager. The null DofManager should be shared only by remote elements (these are only introduced for nonlocal constitutive

model to allow effective local averaging, so only local material value to be averaged are transferred for these remote elements). Null nodes are therefore used only for computing real integration point coordinates of remote elements and there is no reason to maintain their unknowns (they have no equation number assigned, see fig. 7). They do not contribute to local partition governing equation. Only one of the *null remote shared* parameters can be used for particular DofManagers. If no one is used, the DofManager is maintained as local for particular partition.

The list of remote partitions sharing corresponding DofManager or list containing remote partition containing remote DofManager counterpart is specified using *partitions* parameter. The local partition should not be included in the list. The slaves are allowed, but masters have to be in the same partition. The masters can be again remote copies.

Supported DofManagerType keywords are

- Node record

Node	coords _(ra) #	\
	[lcs _(ra) #]	

Represent an abstraction for finite element node. The node coordinates in space (given by global coordinate system) are described using *coords* field. This array contains x, y and possibly z (depends on problem under consideration) coordinate of node. By default, the coordinate system in node is global coordinate system. User defined local coordinate system in node is described using *lcs* array. This array contains six numbers, where the first three numbers represent a directional vector of the local x-axis, and the next three numbers represent a directional vector of the local y-axis. The local z-axis is determined using a vector product. A right-hand coordinate system is assumed. If user defined local coordinate system in node is specified, then the boundary conditions and applied loading are specified in this local coordinate system. The reactions and displacements are also in *lcs* system at the output.

The node can create only master DOFs and SimpleSlave DOFs, so the allowable values of *dofType* array are in range 0,1. For the Node dof manager, the *masterMask* is the array of size equal to number of DOFs, and the i-th value determines the master dof manager, to which i-th dof is directly linked (the dof with same physical meaning are linked together). The local coordinate system in node with same linked dofs is supported, but it should be exactly the same as on master.

- Rigid arm record

RigidArmNode	coords _(ra) #	\
	master _(in) #	\
	[masterMask _(ia) #]	

Represent node connected to other node (called master) using rigid arm. Rigid arm node possesses no degrees of freedom - all dofs are mapped to master dofs. The introduction of rigid arm connected nodes allows to avoid very stiff elements used for modelling the rigid-arm connection. The rigid arm node maps its dofs to master dofs using simple transformations (small rotations are assumed). Therefore, the contribution to rigid arm node are localized directly to master related equations. *The rigid arm node can not have its own boundary or initial conditions, they are determined*

completely from master dof conditions. Currently it is possible to map only certain dofs - see `dofType`. Linked DOFs should have `dofType` value equal to 2, non-linked (primary) DOFs 0.

Rigid arm node can be loaded independently of master. The node coordinates in space (given by global coordinate system) are described using `coords` field. This array contains x, y and possibly z (depends on problem under consideration) coordinate of node. The `master` parameter is the master node number, to which rigid arm node dofs are mapped. The current implementation allows chaining of rigid arm nodes. The optional parameter `masterMask` allows to specify how particular mapped DOF depends on master DOFs. The size of `masterMask` array should be equal to number of DOFs. For all linked DOFs (with corresponding `dofType` value equal to 2) the corresponding value of `masterMask` array should be 1.

- Hanging node

HangingNode	coords _(ra)	#	\
type _(in)		#	\
masters _(ia)		#	\
[ksi _(rn)		#]	\
[eta _(rn)		#]	\
[dzeta _(rn)		#]	\
[weights _(ra)		#]	\

Hanging node connected to other nodes (masters) using generalized interpolation. Hanging node possesses no degrees of freedom (except unlined dofs) - all values are interpolated from corresponding master nodes and their DOFs. The introduction of hanging nodes allows, for example, to include reinforcing bar elements inside arbitrary fe mesh of concrete specimen or facilitates the local refinement of fe-mesh. The hanging nodes can be in chain. The contributions of hanging node are localized directly to master related equations. The hanging node can have its own boundary or initial conditions, but only for primary unlinked DOFs, for linked DOFs, these conditions are determined completely from master dof conditions. The local coordinate system should be same for all master nodes. On the other hand, hanging node can be loaded independently of master.

Values of array `dofType` can have following values: 0-primary DOF, 2-linked DOF.

If parameter `type` equals to zero, then user have to specify for each master its weight (determined by `weights` array). Currently, there is no possibility to have different weights for different DOFs - all DOFs share the same weight).

If parameter `type` is not equal to zero value, then the value determines the type of interpolation used to evaluate receiver unknowns and position and `weights` array is not required. The value is defined as 100*(number of master nodes) + 10*(order of polynomial approximation) + dimension. For example: 211 - linear truss, 312 - linear triangle, 412 - linear rectangle, or 413 - linear tetrahedron. On input, the hanging node coordinates (specified using `coord` array) as well as its natural coordinates (parameters `ksi`, `eta`, `dzeta` determine first, second, and third natural coordinate) should be provided. Natural coordinates are defined as local coordinate that yield

node global coordinates for selected interpolation type (isoparametric interpolation (same order for geometry and unknowns) is assumed).

- Element side

| **ElementSide**

Represents an abstraction for element side, which holds some unknowns.

7.4 Element records

These records specify a description of particular elements. The general format is following:

```
*ElementType (num#)(in) \
               mat(in) # crossSect(in) # nodes(ia) # \
               [bodyLoads(ia) #] [boundaryLoads(ia) #] \
               [activityl(in) #] [lcs(ra) #] \
               <*ElementType> <[partitions(ia) #]> <[remote() #]>
```

The order of element records is optional, the element number is determined by **(num#)**₀ parameter. The numbering of individual elements is arbitrary, it could be even non-continuous. In this context, one could think of element number as a label that is assigned to individual elements and by which the element is referenced. In parallel mode, the label represents a global id across all partitions.

Element material is described by parameter *mat*, which contains corresponding material record number. Element cross section is determined by cross section with *crossSect* record number. Element dof managers (nodes, sides, etc.) defining element geometry are specified using *nodes* array.

Body load acting on element is specified using *bodyLoads* array. Components of this array are corresponding load record numbers. The loads should have the proper type (body load type), otherwise error will be generated.

Boundary load acting on element boundary is specified using *boundaryLoads* array. The format of this array is

$$2 \cdot size \quad lnum(1) \ id(1) \ \dots \ lnum(size) \ id(size),$$

where *size* is total number of loadings applied to element, *lnum(i)* is the applied load number, and *id(i)* is the corresponding entity number, to which the load is applied (for example a side or a surface number). The entity numbering is element dependent and is described in element specific sections. The applied loads must be of proper type (boundary load type), otherwise error is generated.

The support for element insertion and removal during the analysis is provided. One can specify optional time function (identified by its id using *activityl_{tf}* parameter). The nonzero value of this time function indicates, whether the element is active (nonzero value, the default) or inactive (zero value) at particular solution step. Currently available only for structural elements.

Orientation of local coordinates can be specified using *lcs* array. This array contains six numbers, where the first three numbers represent a directional vector of local x-axis, and the next three numbers represent a directional vector of local y-axis. The local z-axis is determined using the vector product. The *lcs* array on the element is particularly useful for modeling of orthotropic materials which follow the element orientation. On a beam or truss element, the *lcs*

array has no effect and the 1D element orientation is aligned with the global xx component.

The *remote* forces the element to be remote element. Remote element does not contribute to local partition governing equation. They are introduced in order to implement band of elements involved in computation of nonlocal variables (see fig. 7 illustrating this approach for node-cut partitioning). They role is to provide local mirror of corresponding remote partition element integration point values which undergo nonlocal averaging on local partition. If not used, element is assumed to be local partition element. When *remote* is used, the *partitions* parameter should contain remote partition number, where corresponding element is local (this array should have size equal to one).

Available material models, their outline and corresponding parameters are described in separate **Element Library Manual**.

7.5 Cross section records

These records specify a cross section model descriptions. The general format is following:

```
| *CrossSectType (num#)(in)
```

The order of particular cross section records is optional, cross section model number is determined by (num#)₀ parameter. The numbering should start from one and should end at n, where n is the number of records.

The crossSectType keyword can be one from following possibilities

- | **SimpleCS** [thick_(rn) #] [width_(rn) #]

Represents integral type of cross section model. In current implementation, such cross section is described using cross section thick (*thickVal*) and width (*widthVal*). For some problems (for example 3d), the corresponding volume and cross section dimensions are determined using element geometry, and then you can omit *thickVal* and *widthVal* parameters.

- | **LayeredCS** nLayers_(in) # \

 LayerMaterials_(ia) # \

 Thicks_(ra) # Widths_(ra) # \

 midSurf_(rn) #

Represents the layered cross section model, based on geometrical hypothesis, that cross sections remain planar after deformation. Number of layers is determined by *nLayers* parameter. Materials for each layer are specified by *LayerMaterials* array. For each layer is necessary to input geometrical characteristic, thick - using *Thicks* array, and width - using *Widths* array. Position of mid surface is determined by its distance from bottom of cross section using *midSurf* parameter (normal and momentum forces are then computed with regard to it's position). Elements using this cross section model must implement layered cross section extension. For information see element library manual.

- | **FiberedCS** nfibers_(in) # fibermaterials_(ia) # \

 thicks_(ra) # widths_(ra) # thick_(rn) # width_(rn) # \

 fiberycentrecoords_(ra) # fiberzcentrecoords_(ra) #

Cross section represented as a set of rectangular fibers. It is based on geometrical hypothesis, that cross sections remain planar after deformation

(3d generalization of layered approach for beams). Parameter *nfibers* determines the number of fibers that together form the overall cross section. The model requires to specify a material model corresponding to particular fiber using *fibermaterials* array. This array should contain for each fibre corresponding material model number (the material model specified on element level has no meaning in this particular case). **The geometry of cross section is determined from fiber dimensions and fiber positions, all input in local coordinate system of the beam (yz plane).** The thick and width of each fiber are determined using *thicks* and *widths* arrays. The overall thick and width are specified using parameters *thick* and *width*. Positions of particular fibers are specified by providing coordinates of center of each fiber using *fiberycentrecoords* array for y-coordinates and *fiberzcentrecoords* array for z-coordinates.

7.6 Material type records

These records specify a material model description. The general format is following:

```
| *MaterialType (num#)(in) d(rn) #
```

The order of particular material records is optional, the material number is determined by (num#)₀ parameter. The numbering should start from one and should end at n, where n is the number of records. Material density is compulsory parameter and it's value is given by *d* parameter.

Available material models, their outline and corresponding parameters are described in separate **Material Library Manual**.

7.7 Nonlocal barrier records

Nonlocal material models of integral type are based on replacement of certain suitable local quantity in local constitutive law by their nonlocal counterparts, that are obtained as weighted average over some characteristic volume. The weighted average is computed as a sum of a remote value multiplied by weight function value. The weight function typically depend on a distance between remote and receiver points and decreases with increasing distance. In some cases, it is necessary to disregard mutual interaction between some points (for example if they are on the opposite sides of a thin notch, which prevents the nonlocal interactions to take place). The barriers are the way how to introduce these constrains. The barrier represent a curve (in 2D) or surface (in 3D). When the line connecting receiver and remote point intersects a barrier, the barriers is activated and the corresponding interaction is not taken into account.

Currently, the supported barrier types are following:

- Polyline barrier

```
| polylinebarrier (num#)(in) vertexnodes(ia) # \
| [xcoordindx(in) #] [ycoordindx(in) #]
```

This represents a polyline barrier for 2D problems. Barrier is a polyline, defined as a sequence of nodes representing vertices. The vertices are specified using parameter *vertexnodes* array, which contains the node

numbers. The optional parameters *xcoordindx* and *ycoordindx* allow to select the plane (xy, yz, or xz), where the barrier is defined. The *xcoordindx* is the first coordinate index, *ycoordindx* is the second. The default values are 1 for *xcoordindx* and 2 for *ycoordindx*, representing barrier in xy plane.

- Symmetry barrier

```
| symmetrybarrier (num#)(in) origin(ra) # \
| normals(ra) # activemask(ia) #
```

Implementation of symmetry barrier, that allows to specify up to three planes (orthogonal ones) of symmetry. This barrier allows to model the symmetry of the averaged field on the boundary without the need of modeling the other part of structure across the plane of symmetry. It is based on modifying the integration weights of source points to take into account the symmetry. The potential symmetry planes are determined by specifying orthogonal right-handed coordinate system, where axes represent the normals of corresponding symmetry planes. Parameter *origin* determines the origin of the coordinate system, the *normals* array contains three components of x-axis direction vector, followed by three components of y-axis direction vector (expressed in global coordinate system). The z-axis is determined from the orthogonality conditions. Parameter *activemask* allows to specify active symmetry planes; i-th nonzero value activates the symmetry barrier for plane with normal determined by corresponding coordinate axis (x=1, y=2, z=3).

7.8 Load and boundary conditions

These records specify description of boundary conditions. The general format is following:

```
| *EntType (num#)(in) \
| loadTimeFunction(in) # [valType(in) #]
```

The order of particular records is optional, boundary condition number is determined by (num#)₀ parameter. The numbering should start from one and should end at n, where n is the number of records. Time function value (given by *loadTimeFunction* parameter) is a multiplier, using which each component (value of loading or value of boundary condition) describes its time variation. The optional parameter *valType* allows to determine the physical meaning of bc value, which is sometimes required. Supported values are (1 - temperature, 2 - force/traction, 3 - pressure, 4 - humidity, 5 - velocity, 6 - displacement). Currently, EntType keyword can be one from

- Dirichlet boundary condition

```
| BoundaryCondition prescribedvalue(rn) # \
| [d(rn) #] \
| [isImposedTimeFunction(in) #]
```

Represents boundary condition. Prescribed value is specified using *prescribedvalue* parameter. The physical meaning of value is fully determined by corresponding DOF. Optionally, the prescribed value can be specified using *d* parameter. It is introduced for compatibility reasons. If *prescribedvalue* is specified, then *d* is ignored. The nonzero value of *isImposedTime-*

Function time function indicates that given boundary condition is active, zero value indicates not active boundary condition in given time (the bc does not exist). By default, the boundary condition apply at any time.

- Nodal fluxes (loads)

| **NodalLoadcomponents**_(ra) # \

| **[cstype**_(in) #]

Concentrated nodal load. The components of nodal load vector are given by *components* parameter. The size of this vector corresponds to a total number of nodal DOFs, and i-th value corresponds to i-th DOF in associated dof manager. The load can be defined in global coordinate system (*cstype* = 0) or in entity - specific local coordinate system (*cstype* = 1, default).

-

| **PrescribedTractionPressureBC**[**isImposedTimeFunction**_(in) #]

Represents pressure boundary condition (of Dirichlet type) due to prescribed tractions. In CBS algorithm formulation the prescribed traction boundary condition leads indirectly to pressure boundary condition in corresponding nodes. This boundary condition implements this pressure bc. The value of bc is determined from applied tractions, that should be specified on element edges/surfaces using suitable boundary loads. The nonzero value of *isImposedTimeFunction* time function indicates that given boundary condition is active, zero value indicates not active boundary condition in given time (the bc does not exist). By default, the boundary condition apply at any time.

Body loads

- Volume flux (load)

| **DeadWeight Components**_(ra) #

Represents dead weight loading applied on element volume (for structural elements). For transport problems, it represents the internal source, i.e. the rate of (heat) generated per unit volume. The magnitude of load for specific i-th DOF is computed as product of material density, corresponding volume and i-th member of *Components* array.

- Structural temperature load

| **StructTemperatureLoad Components**_(ra) #

Represents temperature loading imposed to some elements. The members of *Components* array represent the change of temperature (or change of temperature gradient) corresponding to specific element strain components. See element library manual for details.

- Structural eigenstrain load

| **StructEigenstrainLoad Components**_(ra) #

Prescribes eigenstrain (or stress-free strain) to a structural element. The

array of *Components* is defined in the global coordinate system. The number of components corresponds to a material mode, e.g. plane stress has three components and 3D six. Periodic boundary conditions can be imposed using eigenstrains and master-slave nodes. Consider decomposition of strain into average and fluctuating part

$$\boldsymbol{\varepsilon}(\mathbf{x}) = \langle \boldsymbol{\varepsilon} \rangle + \boldsymbol{\varepsilon}^*(\mathbf{x}) \quad (1)$$

where $\langle \boldsymbol{\varepsilon} \rangle$ can be imposed as eigenstrain over the domain and the solution gives the fluctuating part $\boldsymbol{\varepsilon}^*(\mathbf{x})$. Master-slave nodes have to interconnect opposing boundary nodes of a unit cell.

Boundary loads

- Constant edge fluxes (load)

	ConstantEdgeLoad	ndofs _(in) #	\
		loadType _(in) #	\
		Components _(ra) #	\
		[dofexcludemask _(ia) #]	\
		[csType _(in) #]	\
	[properties _(dc) #]		

- Constant surface fluxes (load)

	ConstantSurfaceLoad	ndofs _(in) #	\
		loadType _(in) #	\
		Components _(ra) #	\
		[dofexcludemask _(ia) #]	\
		[csType _(in) #]	\
	[properties _(dc) #]		

Represent constant edge/surface loads or boundary conditions. The *ndofs* parameter must be set equal to element's number of unknowns. Parameter *loadType* distinguishes the type of boundary condition (supported values are: 2 - prescribed flux input (Neumann boundary condition), 3 - convection bc (Newton boundary condition)). If the boundary condition corresponds to distributed force load, the *Components* array contains components of distributed load corresponding to element unknowns. The load is specified for all DOFs of object to which is associated. For some types of boundary conditions the zero value of load does not mean that the load is not applied (Newton's type of bc, for example). Then some mask, which allows to exclude specific dofs is necessary. The *dofexcludemask* parameter is introduced to allow this. It should have the same size as *Components* array, and by default is filled with zeroes. If some value of *dofExcludeMask* is set to nonzero, then the corresponding componentArray is set to zero and load is not applied for this DOF. If the boundary condition corresponds to prescribed flux input, then the *Components* array contains the components of prescribed input flux corresponding to element unknowns. If the boundary condition is a convection boundary condition (*loadType* = 6) then *Components* array contains the environmental values (temperature of the environment) corresponding to element unknowns, and *properties*

dictionary should contain value of convection coefficient (assumed to be constant) under 'a' key.

The load can be defined in global coordinate system ($csType = 0$, default) or in entity - specific local coordinate system ($csType = 1$).

- Linear edge flux (load)

LinearEdgeLoad	ndofs _(in) #	\
	loadType _(in) #	\
	Components _(ra) #	\
	[dofexcludemask _(ia) #]	\
	[csType _(in) #]	

Represents linear edge load. The meanings of parameters *ndofs*, *csType*, and *loadType* are the same as for **ConstantEdgeLoad**. In *Components* array are stored load components for corresponding unknowns at the beginning of edge (*ndofs* values), followed by values valid for end of edge (*ndofs* values). The load can be defined in global coordinate system ($csType = 0$, default) or in entity - specific local coordinate system ($csType = 1$).

7.9 Initial conditions

These records specify description of initial conditions. The general format is following:

InitialCondition	(num#) _(in)	\
	conditions _(dc) #	

The order of particular records is optional, load, boundary or initial condition number is determined by (**num#**)_o parameter. The numbering should start from one and should end at n, where n is the number of records. Initial parameters are listed in *conditions* dictionary using keys followed by their initial values. Now 'v' key represents velocity and 'a' key represents acceleration.

7.10 Time functions records

These records specify description of time functions, which generally describe time variation of components during solution. The general format is following:

* TimeFuncType	(num#) _(in)	\
	[initialValue _(rn) #]	

The order of these records is optional, time function number is determined by (**num#**)_o parameter. The *initialValue* parameter allows to control the way, how increment of receiver is evaluated for the first solution step. This first solution step increment is evaluated as the difference of value of receiver at this first step and given initial value (which is by default set to zero). The increments of receiver in subsequent steps are computed as a difference between receiver evaluated at given solution step and in previous step.

The numbering should start from one and should end at n, where n is the number of records.

Currently, TimeFuncType keyword can be one from

- Constant function

ConstantFunctionf(t)_(rn) #

Represents the constant time function, with value $f(t)$.

- Peak function

PeakFunction	$\mathbf{t}_{(\text{rn})} \#$	\backslash
	$\mathbf{f}(\mathbf{t})_{(\text{rn})} \#$	

Represents peak time function. If time is equal to t value, then the value of time function is given by $f(t)$ value, otherwise zero value is returned.

- Piecewise function

PiecewiseLinFunction	nPoints _(in) #	\
	t _(ra) # f(t) _(ra) #	

Represents the piecewise time function. The particular time values in t array should be sorted according to time scale. Corresponding time function values are in $f(t)$ array. Value for time, which is not present in t is computed using linear interpolation scheme. Number of time-value pairs is in $nPoints$ parameter.

- Heaviside-like time function

HeavisideLTF	origin _(rn) #	\
	value _(rn) #	

Up to time, given by parameter *origin*, the value of time function is zero. If time greater than *origin* parameter, the value is equal to parameter *value* value.

- User defined

UsrDefLTF	f(t)_(expr) #	\
	[dfdt(t) _(expr) #]	\
	[d2fdt2(t) _(expr) #]	

Represents user defined time function. The expressions can depend on “t” parameter, for which actual time will be substituted and expression evaluated. The function is defined using $f(t)$ parameter, and optionally, its first and second time derivatives using $dfdt(t)$ and $d2fdt2(t)$ parameters. The first and second derivatives may be required, this depend on type of analysis.

Very general, but relatively slow.

7.11 Xfem manager record and associated records

This record creates corresponding xfm manager and specifies the number of its components, in terms of number of enrichment functions, enrichment items, and geometry items. The global shape functions are represented by individual enrichment functions. The enrichment items are abstractions of entities, which are included in the FE model using one or more global functions. Such entity may represent crack, material interface, slipping interface, etc. The particular records will follow immediately in input file. The general format is:

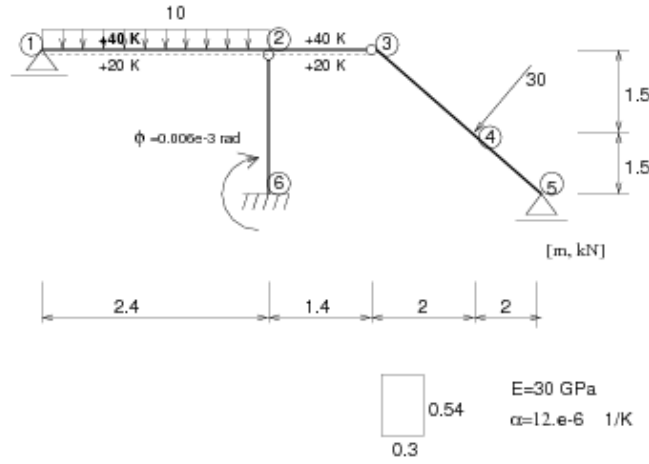


Figure 1: Example 1

```
XfemManager  numberofenrichmentfunctions(in) # \
              numberofenrichmentitems(in) # \
              numberofgeometryitems(in) #
```

where *numberofenrichmentfunctions* represents number of enrichment functions and their associated records, *numberofenrichmentitems* represents number of enrichment items and corresponding records, and *numberofgeometryitems* represents number of geometry items, that are used to define the geometry of individual enrichment items.

The enrichment function records describe individual enrichment functions. The general format is following:

```
*EntType  (num#)(in)
```

The geometry item records describe individual geometries of enrichment items. The general format is following:

```
*EntType  (num#)(in)
```

The enrichment item records describe individual enrichment items. The general format is following:

```
*EntType  (num#)(in) \
          (geometryitem#)(in) \
          (enrichmentfunction#)(in)
```

8 Examples

8.1 Beam structure

The example of input file for simple beam structure is presented. Structure geometry and its constitutive and geometrical properties are shown in fig. (1). The linear static analysis is required, the influence of shear is neglected. Please note that one input line which is too long to fit to page, is separated into two lines using backslash '\' character, but OOFEM requires that it must be typed as one single line.

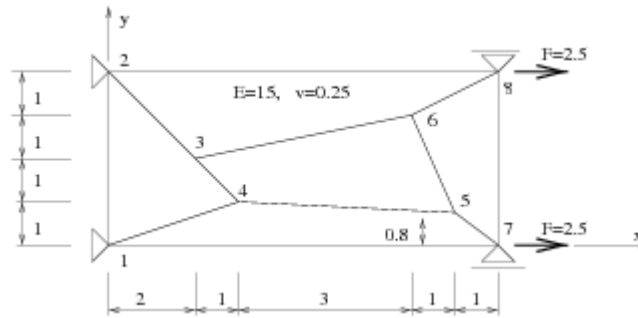


Figure 2: Example 2

```

test41.out
Simple Beam Structure - linear analysis
#only momentum influence to the displacements is taken into account
#beamShearCoeff is artificially enlarged.
LinearStatic 1 nsteps 3
domain 2dBeam
OutputManager timestep_all dofman_all element_all
ndofman 6 nelem 5 ncrosssect 1 nmat 1 nbc 5 nic 0 nltf 3
node 1 coords 3 0. 0. 0. bc 3 0 1 0
node 2 coords 3 2.4 0. 0. bc 3 0 0 0
node 3 coords 3 3.8 0. 0. bc 3 0 0 1
node 4 coords 3 5.8 0. 1.5 bc 3 0 0 0 load 1 4
node 5 coords 3 7.8 0. 3.0 bc 3 0 1 0
node 6 coords 3 2.4 0. 3.0 bc 3 1 1 2
Beam2d 1 nodes 2 1 2 mat 1 crossSect 1 boundaryLoads 2 3 1 bodyLoads 1 5
Beam2d 2 nodes 2 2 3 mat 1 crossSect 1 DofsToCondense 1 6 bodyLoads 1 5
Beam2d 3 nodes 2 3 4 mat 1 crossSect 1 DofsToCondense 1 3
Beam2d 4 nodes 2 4 5 mat 1 crossSect 1
Beam2d 5 nodes 2 6 2 mat 1 crossSect 1 DofsToCondense 1 6
SimpleCS 1 area 0.162 Iy 0.0039366 beamShearCoeff 1.e18 thick 0.54
IsoLE 1 d 1. E 30.e6 n 0.2 tAlpha 1.2e-5
BoundaryCondition 1 loadTimeFunction 1 prescribedvalue 0.0
BoundaryCondition 2 loadTimeFunction 2 prescribedvalue -0.006e-3
ConstantEdgeLoad 3 loadTimeFunction 1 Components 3 0. 10. 0.0\
loadType 3 ndofs 3
NodalLoad 4 loadTimeFunction 1 Components 3 -18.0 24.0 0.0
StructTemperatureLoad 5 loadTimeFunction 3 Components 2 30.0 -20.0
PeakFunction 1 t 1.0 f(t) 1.
PeakFunction 2 t 2.0 f(t) 1.
PeakFunction 3 t 3.0 f(t) 1.

```

8.2 Plane stress example

```

patch100.out
Patch test of PlaneStress2d elements -> pure compression
LinearStatic nsteps 1
domain 2dPlaneStress
OutputManager timestep_all dofman_all element_all
ndofman 8 nelem 5 ncrosssect 1 nmat 1 nbc 2 nic 0 nltf 1

```

```

node 1 coords 3 0.0 0.0 0.0 bc 2 1 1
node 2 coords 3 0.0 4.0 0.0 bc 2 1 1
node 3 coords 3 2.0 2.0 0.0 bc 2 0 0
node 4 coords 3 3.0 1.0 0.0 bc 2 0 0
node 5 coords 3 8.0 0.8 0.0 bc 2 0 0
node 6 coords 3 7.0 3.0 0.0 bc 2 0 0
node 7 coords 3 9.0 0.0 0.0 bc 2 0 1 load 1 2
node 8 coords 3 9.0 4.0 0.0 bc 2 0 1 load 1 2
PlaneStress2d 1 nodes 4 1 4 3 2 crossSect 1 mat 1 NIP 1
PlaneStress2d 2 nodes 4 1 7 5 4 crossSect 1 mat 1 NIP 1
PlaneStress2d 3 nodes 4 4 5 6 3 crossSect 1 mat 1 NIP 1
PlaneStress2d 4 nodes 4 3 6 8 2 crossSect 1 mat 1 NIP 1
PlaneStress2d 5 nodes 4 5 7 8 6 crossSect 1 mat 1 NIP 1
SimpleCS 1 thick 1.0 width 1.0
IsoLE 1 d 0. E 15.0 n 0.25 talpha 1.0
BoundaryCondition 1 loadTimeFunction 1 prescribedvalue 0.0
NodalLoad 2 loadTimeFunction 1 Components 2 2.5 0.0
ConstantFunction 1 f(t) 1.0

```

9 Examples - parallel mode

9.1 Node-cut example

The example shows explicit direct integration analysis of simple structure with two DOFs. The geometry and partitioning is sketched in fig. 3.

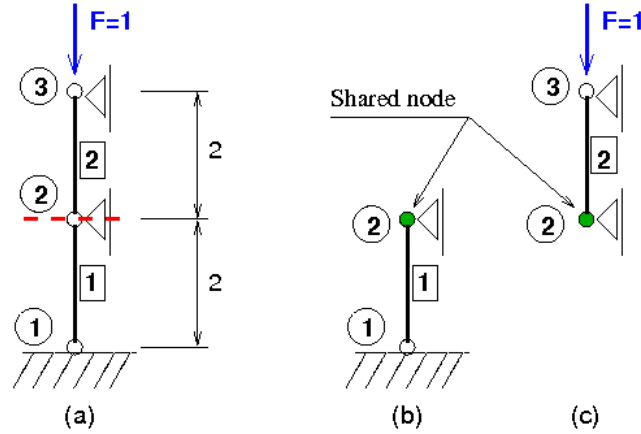


Figure 3: Node-cut partitioning example: (a) whole geometry, (b) partition 0, (c) partition 1.

```

#
# partition 0
#
partest.out.0
Parallel test of explicit oofem computation
#
PN1DEIDynamic nsteps 3 dumpcoef 0.0 deltaT 1.0 NodeCutMode

```



```

domain 2dTruss
#
OutputManager tstep_all dofman_all element_all
ndofman 2 nelem 1 ncrosssect 1 nmat 1 nbc 2 nic 0 nltf 1
#
Node 1 coords 3 0. 0. 0. bc 2 1 1
Node 2 coords 3 0. 0. 2. bc 2 1 0 Shared partitions 1 1
Truss2d 1 nodes 2 1 2 mat 1 crossSect 1
SimpleCS 1 thick 0.1 width 10.0
IsoLE 1 tAlpha 0.000012 d 10.0 E 1.0 n 0.2
BoundaryCondition 1 loadTimeFunction 1 conditions 1 d 0.0
NodalLoad 2 loadTimeFunction 1 Components 2 0. 1.0
ConstantFunction 1 f(t) 1.0

#
# partition 1
#
partest.out.1
Parallel test of explicit oofem computation
#
PNlDEIDynamic nsteps 3 dumpcoef 0.0 deltaT 1.0 NodeCutMode
domain 2dTruss
#
OutputManager tstep_all dofman_all element_all
ndofman 2 nelem 1 ncrosssect 1 nmat 1 nbc 2 nic 0 nltf 1
#
Node 2 coords 3 0. 0. 2. bc 2 1 0 Shared partitions 1 0
Node 3 coords 3 0. 0. 4. bc 2 1 0 load 1 2
Truss2d 2 nodes 2 2 3 mat 1 crossSect 1
SimpleCS 1 thick 0.1 width 10.0
IsoLE 1 tAlpha 0.000012 d 10.0 E 1.0 n 0.2
BoundaryCondition 1 loadTimeFunction 1 conditions 1 d 0.0
NodalLoad 2 loadTimeFunction 1 Components 2 0. 1.0
ConstantFunction 1 f(t) 1.0

```

9.2 Element-cut example

The example shows explicit direct integration analysis of simple structure with two DOFs. The geometry and partitioning is sketched in fig. 3.

```

#
# partition 0
#
partest2.out.0
Parallel test of explicit oofem computation
#
PNlDEIDynamic nsteps 5 dumpcoef 0.0 deltaT 1.0 ElementCutMode
domain 2dTruss
#
OutputManager tstep_all dofman_all element_all

```

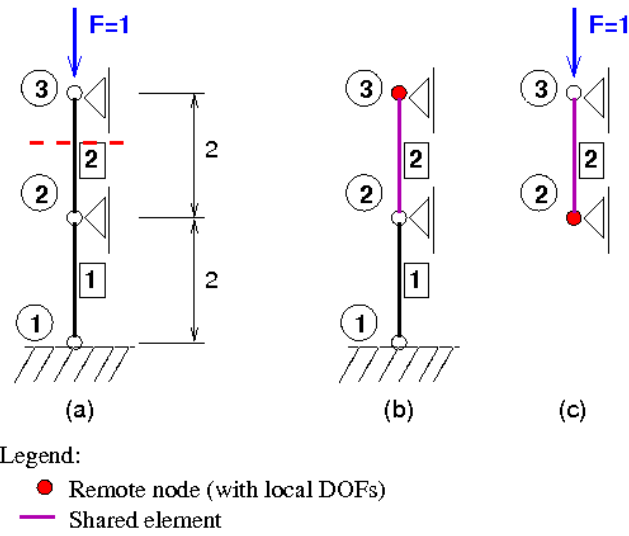


Figure 4: Element-cut partitioning example: (a) whole geometry, (b) partition 0, (c) partition 1.

```
ndofman 3 nelem 2 ncrosssect 1 nmat 1 nbc 2 nic 0 nltf 1
#
Node 1 coords 3 0. 0. 0. bc 2 1 1
Node 2 coords 3 0. 0. 2. bc 2 1 0
Node 3 coords 3 0. 0. 4. bc 2 1 0 \
    Remote partitions 1 1 load 1 2
Truss2d 1 nodes 2 1 2 mat 1 crossSect 1
Truss2d 2 nodes 2 2 3 mat 1 crossSect 1
SimpleCS 1 thick 0.1 width 10.0
IsoLE 1 tAlpha 0.000012 d 10.0 E 1.0 n 0.2
BoundaryCondition 1 loadTimeFunction 1 conditions 1 d 0.0
NodalLoad 2 loadTimeFunction 1 Components 2 0. 1.0
ConstantFunction 1 f(t) 1.0

#
# partition 1
#
partest2.out.1
Parallel test of explicit oofem computation
#
PNlDEIDynamic nsteps 5 dumpcoef 0.0 deltaT 1.0 ElementCutMode
domain 2dTruss
#
OutputManager tstep_all dofman_all element_all
ndofman 2 nelem 1 ncrosssect 1 nmat 1 nbc 2 nic 0 nltf 1
#
Node 2 coords 3 0. 0. 2. bc 2 1 0 Remote partitions 1 0
```

```

Node 3 coords 3 0. 0. 4. bc 2 1 0 load 1 2
Truss2d 2 nodes 2 2 3 mat 1 crossSect 1
SimpleCS 1 thick 0.1 width 10.0
IsoLE 1 tAlpha 0.000012 d 10.0 E 1.0 n 0.2
BoundaryCondition 1 loadTimeFunction 1 conditions 1 d 0.0
NodalLoad 2 loadTimeFunction 1 Components 2 0. 1.0
ConstantFunction 1 f(t) 1.0

```

10 Figures

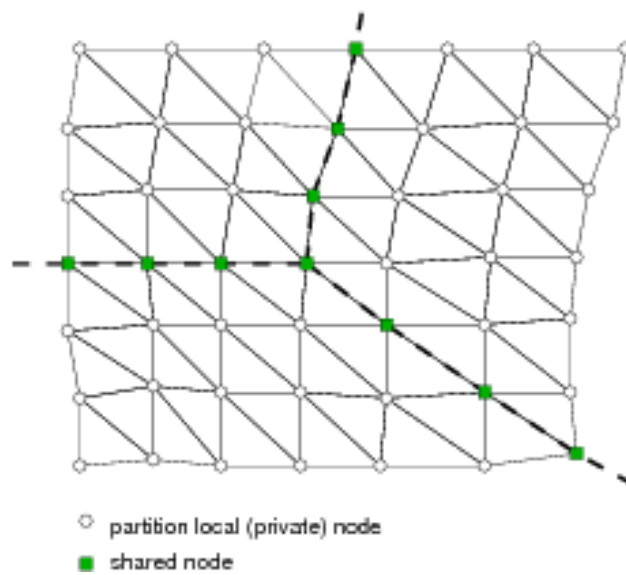


Figure 5: Node-cut partitioning.

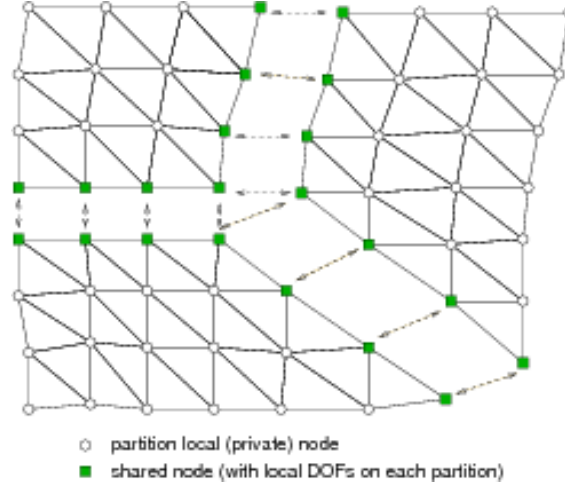


Figure 6: Node-cut partitioning - local constitutive mode.

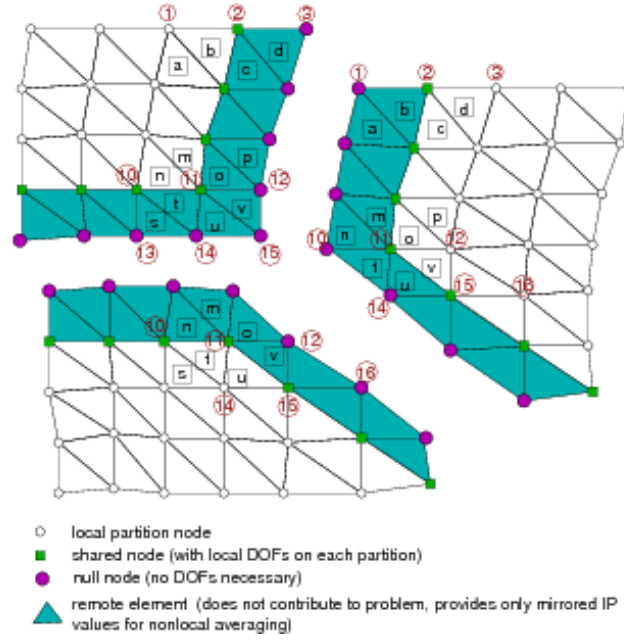


Figure 7: Node-cut partitioning - nonlocal constitutive mode.

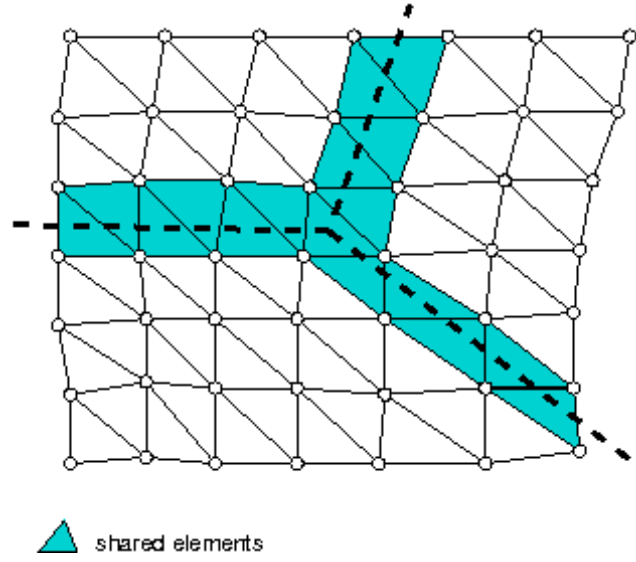


Figure 8: Element-cut partitioning.

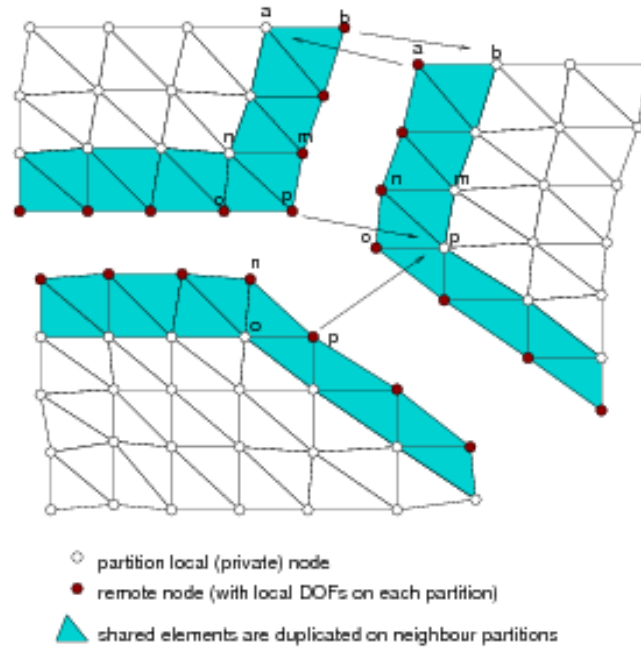


Figure 9: Element-cut partitioning, local constitutive mode.