

Stabil

MATLAB toolbox for Structural Mechanics

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Chapter 1

Getting started

1.1 About Stabil

Stabil is a MATLAB toolbox for structural mechanics, based on the finite element method. Classical finite element beam and truss elements, as well as shell elements and elements for 2D and 3D elasticity have been implemented, and the toolbox can be used to solve a variety of static and dynamic structural problems. In addition, Stabil contains a number of postprocessing functions tailored to structural analysis, including plots of member forces and displacements of a structure. The user can interact with Stabil at a low level of abstraction or a high level of abstraction, which is further detailed in this manual. Due to this multi-level approach, the toolbox is suitable for educational purposes and for use in a research environment: the high level functions allow for an easy and efficient implementation of many common problems, while the low level functions facilitate customization and the implementation of novel finite element techniques in a research context.

Stabil is written in standard MATLAB language, so that the source code is available to the user. No additional MATLAB toolboxes are required and the toolbox can be used in a Windows, Linux, or MacOS environment.

1.2 Obtaining and installing Stabil

Stabil can be downloaded from the internet at <https://bwk.kuleuven.be/bwm/stabil>. It is distributed as a ZIP archive, which should be extracted to a directory on the hard disk, e.g. `C:\Users\%username%\Documents\matlab\stabil`. After extraction of the ZIP archive, this directory should contain a number of `m`-files.

The Stabil directory must be subsequently added to the MATLAB path to make the toolbox functions available in MATLAB:

- In MATLAB, click on ‘Set Path...’ in the ‘Home’ tab.
- Click on ‘Add Folder’ and select the Stabil directory.
- Save the path and close the dialog window.

1.3 Terms of use

Stabil is free software: you can redistribute it and/or modify it under the terms of the GNU General Public License as published by the Free Software Foundation, either version 3 of the License, or (at your option) any later version. The toolbox can be used and modified for educational and research purposes. However, it cannot be used for consulting, nor commercialized in any form. Regardless of any provision of the GNU General Public License, Stabil may not be used for commercial purposes without explicit written permission from the authors.

Furthermore, if you use Stabil for research, please include a reference to Stabil in scientific publications of your work (articles, reports, books, etc.).

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1.4 Scope and structure of this document

The present document is the user's guide to Stabil version 4.0. It is a combination of a reference guide, providing an overview of all functions of Stabil, and a tutorial, presenting a set of examples to illustrate the use of Stabil. This document is not meant as a textbook: a theoretical background is only considered where this is necessary to define the functionality of Stabil in a clear way. The reader is referred to various works on Matrix analysis of Structures [5], Finite Elements [1, 3, 4, 6–9], or Dynamics of Structures [2] for a broad theoretical background.

This document is composed of the following chapters:

Chapter 1. Getting started (p. 1)

The aim of this chapter is to get the user started with Stabil and the accompanying user's guide. The installation procedure of Stabil is explained, the terms of use of Stabil are clarified, and the scope and structure of the user's guide are discussed.

Chapter 2. Static analysis of structures (p. 3)

In this chapter, the use of stabil for the static analysis of structures is explained through a number of examples.

Chapter 3. Dynamic analysis of structures (p. 25)

In this chapter, the use of stabil for the dynamic analysis of structures is explained through a number of examples.

Chapter 4. Element guide (p. 37)

This chapter presents an overview of the element types available in Stabil. The conventions (local coordinate system, nodal connectivity) for each element type are given and reference is made to the Stabil functions related to the element implementation.

Chapter 5. Functions — By category (p. 55)

This chapter gives an overview of all functions in Stabil, organized by category.

Chapter 6. Functions — Alphabetical list (p. 57)

This chapter consists of an alphabetical list of the functions in Stabil. The syntax, the input and output arguments, and the use of each function are described in detail. The information provided in this chapter is also accessible at the MATLAB prompt through the `help` command.

Chapter 2

Static analysis of structures

2.1 Example 1.1: static analysis of a frame

The basic concepts of the StabIl toolbox are introduced through the example of a simple frame structure, shown in figure 2.1. The frame has a height and width of 4 m and is clamped at node 1 and pinned at node 5. An internal hinge is present at nodes 2 and 3, and a diagonal brace is present between nodes 1 and 4. At node 4 a point load $F = 5 \text{ kN}$ is applied and the vertical beam on the left is loaded by a distributed load $p = 2 \text{ kN/m}$. The beams and columns have a concrete rectangular cross-section with a width of 0.2 m and a height of 0.4 m. The concrete has a Young's modulus of 30 GPa and a Poisson's ratio of 0.2. The diagonal brace has a circular steel cross-section with a diameter of 8 mm with a Young's modulus of 210 GPa and a Poisson coefficient of 0.3.

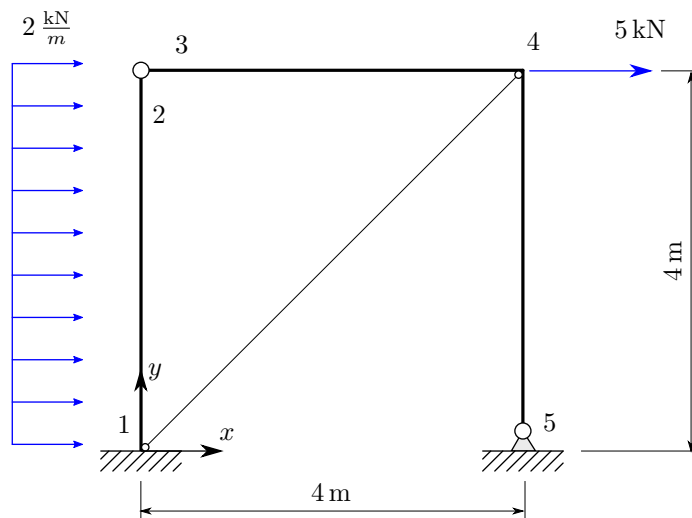


Figure 2.1: Simple frame structure.

The code to compute the deformation and bending moment in the frame structure is listed below:

```
% StaBIL manual
% Example 1.1: static analysis of a frame
% Units: m, kN

% Nodes=[NodID X Y Z]
Nodes= [1    0  0  0;
        2    0  4  0;
        3    0  4  0;
        4    4  4  0;
        5    4  0  0;
        6    1  5  0]; % reference node

% Check the node coordinates as follows:
figure
plotnodes(Nodes);
```

```

% Element types -> {EltTypeID EltName}
Types= {1      'beam';
        2      'truss'};

b=0.10;
h=0.25;
r=0.004;

% Sections=[SecID A      ky  kz  Ixx Iyy Izz      yt  yb  zt  zb]
Sections= [1      b*h      Inf  Inf  0  0  b*h^3/12  h/2  h/2  b/2  b/2;
          2      pi*r^2  NaN  NaN  NaN  NaN  NaN      NaN  NaN  NaN  NaN];

% Materials=[MatID      E nu];
Materials= [1      30e6 0.2;          % concrete
           2      210e6 0.3];        % steel

% Elements=[EltID TypID SecID MatID n1 n2 n3]
Elements= [1      1      1      1      1 2 6;
          2      1      1      1      3 4 6;
          3      1      1      1      5 4 6;
          4      2      2      2      1 4 NaN];

% Check node and element definitions as follows:
hold('on');
plotelem(Nodes,Elements,Types);
title('Nodes and elements');

% Degrees of freedom
% Assemble a column matrix containing all DOFs at which stiffness is
% present in the model:
DOF=getdof(Elements,Types);

% Remove all DOFs equal to zero from the vector:
% - 2D analysis: select only UX,UY,ROTZ
% - clamp node 1
% - hinge at node 5
seldof=[0.03; 0.04; 0.05; 1.00; 5.01; 5.02];
DOF=removedof(DOF,seldof);

% Assembly of stiffness matrix K
K=asmmk(Nodes,Elements,Types,Sections,Materials,DOF);

% Nodal loads: 5 kN horizontally on node 4.
seldof=[4.01];
PLoad= [5];

% Assembly of the load vectors:
P=nodalvalues(DOF,seldof,PLoad);

% Distributed loads are specified in the global coordinate system
% DLoads=[EltID n1globalX n1globalY n1globalZ ...]
DLoads= [1      2 0 0 2 0 0];

P=P+elemloads(DLoads,Nodes,Elements,Types,DOF);

% Constraint equations: Constant=Coef1*DOF1+Coef2*DOF2+ ...
% Constraints=[Constant Coef1 DOF1 Coef2 DOF2 ...]
Constr= [0      1      2.01 -1      3.01;
        0      1      2.02 -1      3.02];

% Add constraint equations
[K,P]=addconstr(Constr,DOF,K,P);

% Solve K * U = P
U=K\P;

```



```

% Plot displacements
figure
plotdisp(Nodes,Elements,Types,DOF,U,DLoads,Sections,Materials)

% The displacements can be displayed as follows:
printdisp(Nodes,DOF,U);

% Compute element forces
Forces=elemforces(Nodes,Elements,Types,Sections,Materials,DOF,U,DLoads);

% The element forces can be displayed in a orderly table:
printforc(Elements,Forces);

% Plot element forces
figure
plotforc('norm',Nodes,Elements,Types,Forces,DLoads)
title('Normal forces')

figure
plotforc('sheary',Nodes,Elements,Types,Forces,DLoads)
title('Shear forces')

figure
plotforc('momz',Nodes,Elements,Types,Forces,DLoads)
title('Bending moments')

% Plot stresses
figure
plotstress('snorm',Nodes,Elements,Types,Sections,Forces,DLoads)
title('Normal stresses due to normal forces')

figure
plotstress('smomzt',Nodes,Elements,Types,Sections,Forces,DLoads)
title('Normal stresses due to bending moments around z: top')

figure
plotstress('smomzb',Nodes,Elements,Types,Sections,Forces,DLoads)
title('Normal stresses due to bending moments around z: bottom')

figure
plotstress('smax',Nodes,Elements,Types,Sections,Forces,DLoads)
title('Maximal normal stresses')

figure
plotstress('smin',Nodes,Elements,Types,Sections,Forces,DLoads)
title('Minimal normal stresses')

```

After the definition of model parameters, the analysis starts by defining the nodes of the model, where the nodes are defined in the (x,y) -plane. The nodes are represented by a matrix that contain node numbers as a first column, followed by the x , y and z coordinates of each node. Next, the element types are defined by a cell array, **Types**, containing type numbers followed by a string (either **beam** or **truss**). The **Sections** matrix defines the section properties (cross section, moments of inertia,...). In a similar way, the **Materials** matrix defines material properties (the Young's modulus and Poisson coefficient). The model definition is then completed by providing an element connectivity table **Elements**, which refers to the previously defined element, material, section and node numbers. In order to uniquely define a local coordinate system for the beam elements, use is made of a reference node (node 6). The local x -axis is directed from the first node of the element to the last node, the local y -axis is perpendicular to the local x -axis with its origin corresponding to the first node of the element, and pointing in the direction of the reference node. The same convention applies in a three-dimensional setting, as indicated in the **beam** element reference sheet on page 38.

Next, the degrees of freedom (DOF's) are specified. The DOF's are defined using a **node.index** approach, where the digits to the left of the decimal point refer to the node number and the digits to the right of the decimal point refer to degree of freedom. By convention, the degrees of freedom 01, 02, and 03 correspond to translations u_x , u_y , and u_z in the global coordinate directions whereas the degrees of freedom 04, 05, and 06

Node	U_x [m]	U_y [m]	U_z [m]	φ_x [rad]	φ_y [rad]	φ_z [rad]
1	0.0000e+0	0.0000e+0	0.0000e+0	0.0000e+0	0.0000e+0	0.0000e+0
2	6.6633e-3	3.2297e-6	0.0000e+0	0.0000e+0	0.0000e+0	-1.8161e-3
3	6.6633e-3	3.2297e-6	0.0000e+0	0.0000e+0	0.0000e+0	4.0356e-4
4	6.6538e-3	-3.6160e-5	0.0000e+0	0.0000e+0	0.0000e+0	-8.3665e-4
5	0.0000e+0	0.0000e+0	0.0000e+0	0.0000e+0	0.0000e+0	-2.0769e-3
6	0.0000e+0	0.0000e+0	0.0000e+0	0.0000e+0	0.0000e+0	0.0000e+0
Element Node	N [kN]	V_y [kN]	V_z [kN]	T [kNm]	M_y [kNm]	M_z [kNm]
1 i	6.0557e-1	6.2201e+0	-0.0000e+0	-0.0000e+0	-0.0000e+0	-8.8804e+00
1 j	6.0557e-1	-1.7799e+0	0.0000e+0	0.0000e+0	0.0000e+0	-4.4409e-16
2 i	-1.7799e+0	6.0557e-1	-0.0000e+0	-0.0000e+0	-0.0000e+0	0.0000e+00
2 j	-1.7799e+0	6.0557e-1	0.0000e+0	0.0000e+0	0.0000e+0	2.4223e+00
3 i	-6.7799e+0	-6.0557e-1	-0.0000e+0	-0.0000e+0	-0.0000e+0	6.6613e-16
3 j	-6.7799e+0	-6.0557e-1	0.0000e+0	0.0000e+0	0.0000e+0	-2.4223e+00
4 i	8.7318e+0	-0.0000e+0	-0.0000e+0	-0.0000e+0	-0.0000e+0	0.0000e+00
4 j	8.7318e+0	0.0000e+0	0.0000e+0	0.0000e+0	0.0000e+0	-0.0000e+00

Table 2.1: Nodal displacements, reaction forces, and member forces for example 1.1.

correspond to rotations φ_x , φ_y , φ_z around the global coordinate axes. For example, DOF 3.02 represents the translation u_y of node 3. This approach of defining the degrees of freedom for the problem at hand is very instructive, since it enforces reasoning on the kinematics of the structure and how boundary conditions are accounted for. This manual input of the vector of degrees of freedom is seen as a low-level functionality of the toolbox. Alternatively, the high-level functions (`getdof`, `selectdof`, `removedof`) could be used to generate and process this DOF vector, allowing for more complex structural models to be analyzed in an efficient way.

The finite element stiffness matrix is next assembled using the `asmkm` function. This function takes the `Nodes`, `Elements`, `Types`, `Sections`, and `Materials` variables that define the finite element model and assembles the sparse global stiffness matrix K corresponding to the `dof` vector. The `asmkm` function is a high-level function that loops over the various elements of the mesh and calls low-level functions that generate element stiffness matrices.

In `Stabil`, both nodal loads as well as distributed loads on elements can be considered, and dedicated functions are available to either generate nodal loads on DOFs (the `nodalvalues` function) or distributed loads on elements (the `elemloads` function). The load vector P is available in the Matlab environment, and has the same size as the `dof`-vector.

In order to account for the internal hinge between nodes 2 and 3, constraint equations are added to the system of equations to couple the horizontal and vertical displacements of both nodes. This is achieved through the function `addconstr` that modifies the stiffness matrix K to account for linear constraint equations that relate various degrees of freedom in the system.

The finite element system of equations ($Ku = f$) is solved using the Matlab \ (left divide, `mldivide`) command. Since the global stiffness matrix K is a sparse matrix, a sparse solver is used by Matlab by default. The resulting displacement vector u has the same size as the `dof` vector and the load vector P and is equally available in the Matlab environment.

In `Stabil`, a number of postprocessing functions is available to easily plot the deformed structure or the resulting member forces. Figure 2.2 shows the resulting displacement of the structure as obtained with the `plotdisp` command. The `plotdisp` command takes the model definition (`Nodes`, `Elements`, `Types`), the displacement solution and corresponding `dof` vector, and (optionally) the distributed load with the section and material definitions (`Sections`, and `Materials`). A key feature of the `Stabil` toolbox is that the deformed shape is plotted in an exact way, accounting for the (cubic) shape functions of the beam element and the deformation of the element due to distributed loads. This allows to demonstrate how distributed loads are reduced to equivalent nodal loads.

Figure 2.3 shows the member forces in the structure as plotted with the `plotforc` command. Like the `plotdisp` command, the `plotforc` command accounts for the effect of distributed loads, as is apparent in the quadratic variation of the bending moment in the left column in figure 2.3.

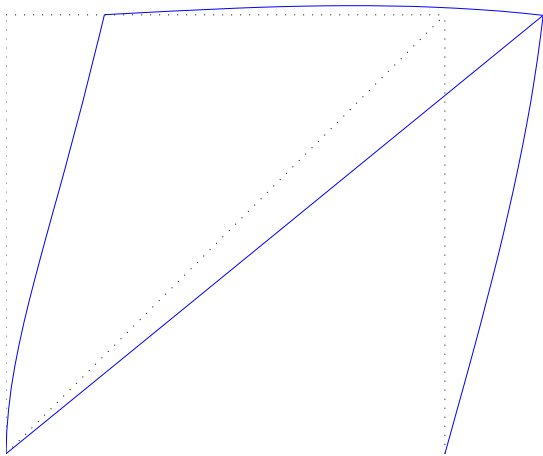


Figure 2.2: Deformed frame structure.

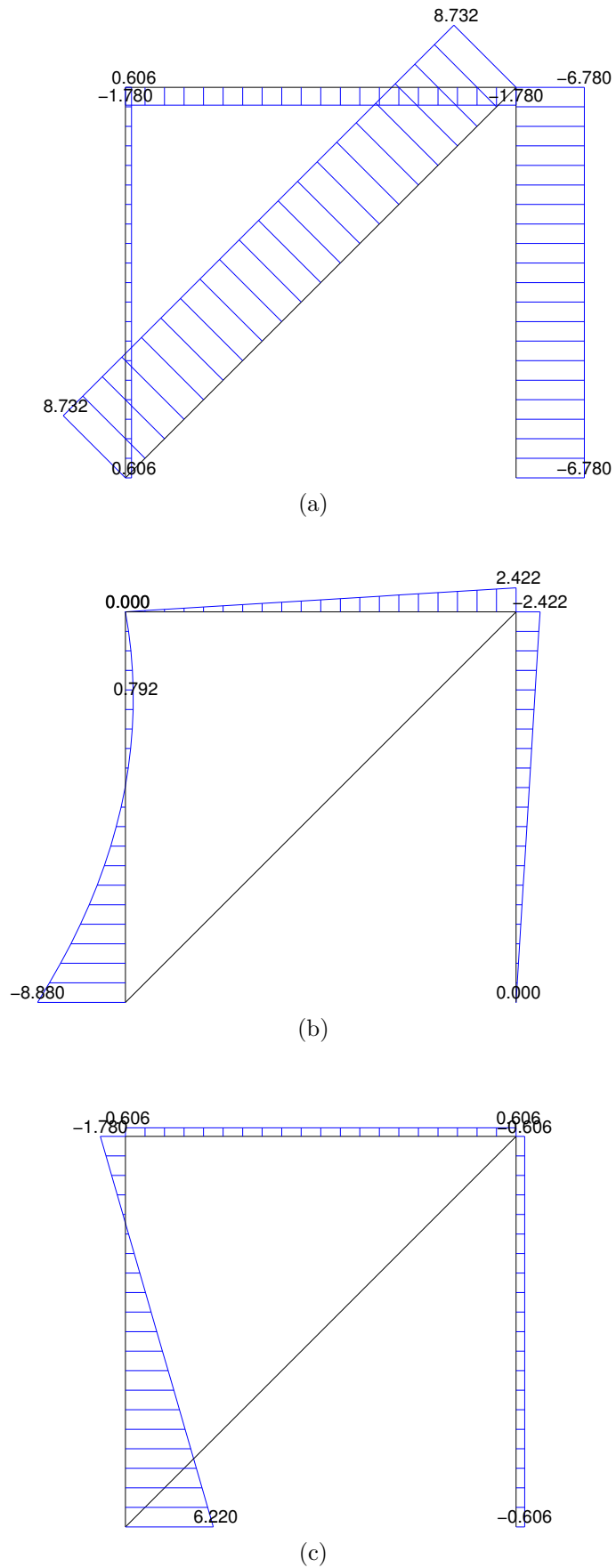


Figure 2.3: (a) Normal forces [kN], (b) bending moments [kNm], and (c) shear forces [kN] in the frame structure.

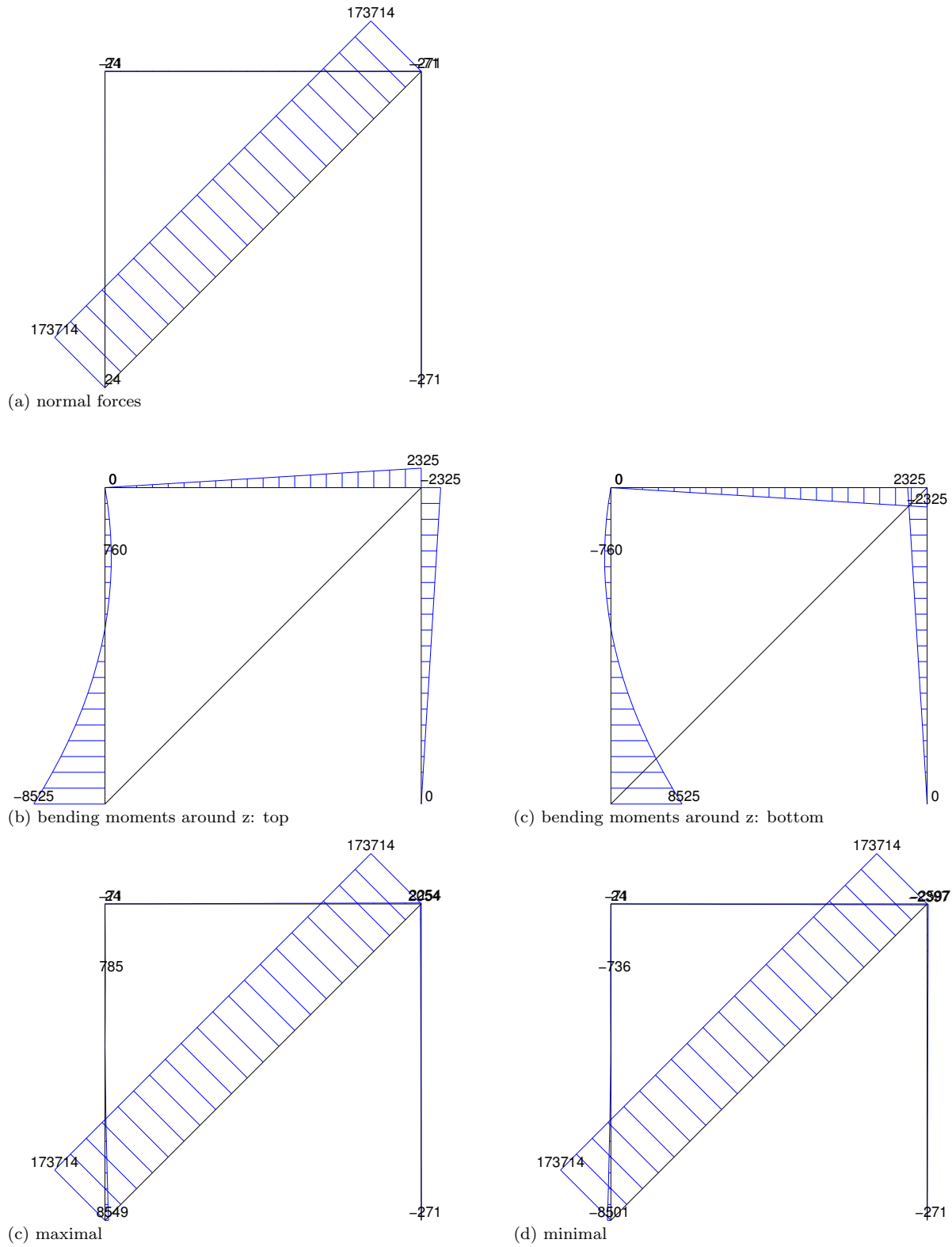


Figure 2.4: Normal stresses.

2.2 Example 1.2: static analysis of a 3D frame

In order to demonstrate the three-dimensional capabilities of StabIl, the following example considers a 3D frame structure (figure 2.5). The frame consists of rectangular concrete beams (Young's modulus $E = 35 \times 10^9 \text{ N/m}^2$ and Poisson coefficient $\nu = 0.2$) with a height of 0.5 m and a width of 0.2 m and concrete columns of 0.3 m by 0.2 m.

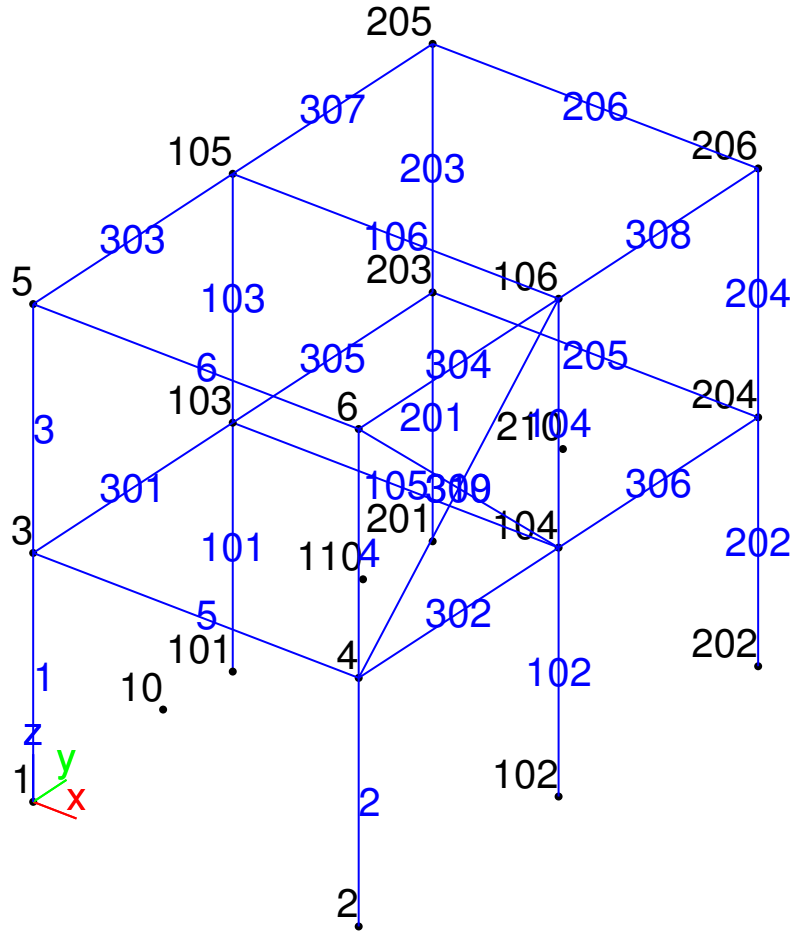


Figure 2.5: Example 1.2: nodes, elements, loads and boundary conditions.

The following input file is structured in a similar way as example 1.1. It starts by defining nodes, element types, sections, and materials. Subsequently, the elements are defined by the use of a reference node. The `repro` command is employed to replicate rows of the `Nodes` and `Elements` matrix. The `getdof` and `seldof` commands are then used to generate and process the DOF vector. After assembling the stiffness matrix and assembling the loads, the system is solved, the forces and reaction forces are computed and results are finally plotted, resulting in figures 2.6 to 2.9 for the different load cases and load combinations.

```
% StaBIL manual
% Example 1.2: static analysis of a 3D frame
% Units: m, kN

% Types={EltTypID EltName}
Types= {1      'beam';
        2      'truss'};

% Sections
bCol=0.2; % Column section width
hCol=0.3; % Column section height
bBeam=0.2; % Beam section width
hBeam=0.5; % Beam section height
Atruss=0.001;
```

```

% Sections=[SecID A ky kz Ixx Iyy Izz yt yb zt zbl]
Sections= [1 hCol*bCol Inf Inf 0.196*bCol^3*hCol bCol^3*hCol/12 ...
           hCol^3*bCol/12 hCol/2 hCol/2 bCol/2 bCol/2; % Columns
          2 hBeam*bBeam Inf Inf 0.249*bBeam^3*hBeam bBeam^3*hBeam/12 ...
           hBeam^3*bBeam/12 hBeam/2 hBeam/2 bBeam/2 bBeam/2; % Beams
          3 Atruss NaN NaN NaN NaN NaN NaN NaN NaN];

% Materials=[MatID E nu rho];
Materials= [1 35e9 0.2 2500; %concrete
            2 210e9 0.3 7850; %steel

L=5;
H=3.5;
B=4;

% Nodes=[NodID X Y Z]
Nodes= [1 0 0 0;
        2 L 0 0]
Nodes=reprow(Nodes,1:2,2,[2 0 0 H])
Nodes=[Nodes;
        10 2 0 2] % reference node
Nodes=reprow(Nodes,1:7,2,[100 0 B 0])

figure
plotnodes(Nodes);

% Elements=[EltID TypID SecID MatID n1 n2 n3]
Elements=[ 1 1 1 1 1 3 10;
           2 1 1 1 2 4 10];
Elements=reprow(Elements,1:2,1,[2 0 0 0 2 2 0])
Elements=[ Elements;
           5 1 2 1 3 4 10;
           6 1 2 1 5 6 10];
Elements=reprow(Elements,1:6,2,[100 0 0 0 100 100 100])
Elements=[ Elements;
           301 2 3 2 3 103 NaN;
           302 2 3 2 4 104 NaN];
Elements=reprow(Elements,19:20,1,[2 0 0 0 2 2 0])
Elements=reprow(Elements,19:22,1,[4 0 0 0 100 100 0])
Elements=[ Elements;
           309 2 3 2 4 106 NaN;
           310 2 3 2 6 104 NaN];

hold('on');
plotelem(Nodes,Elements,Types);
title('Nodes and elements');

% Plot elements in different colors in order to check the section definitions
figure
plotelem(Nodes,Elements(find(Elements(:,3))==1,:),Types,'r');
hold('on');
plotelem(Nodes,Elements(find(Elements(:,3))==2,:),Types,'g');
plotelem(Nodes,Elements(find(Elements(:,3))==3,:),Types,'b');
title('Elements: sections')

% Degrees of freedom
DOF=getdof(Elements,Types);

% Boundary conditions: hinges
seldof=[ 1.01; 1.02; 1.03; 2.01; 2.02; 2.03;
         101.01; 101.02; 101.03; 102.01; 102.02; 102.03;
         201.01; 201.02; 201.03; 202.01; 202.02; 202.03];

DOF=removedof(DOF,seldof);

```

```

% Assembly of stiffness matrix K
K=asmmk(Nodes,Elements,Types,Sections,Materials,DOF);

% Loads

% Own weight
DLoadsOwn=accel([0 0 9.81],Elements,Types,Sections,Materials);

% Wind load

% DLoads=[EltID n1globalX n1globalY n1globalZ ...]
DLoadsWind =[1 0 0    0 0 1500 0;
              2 0 0    0 0 1500 0;
              3 0 1500 0 0 1500 0;
              4 0 1500 0 0 1500 0];

DLoads=multdloads(DLoadsOwn,DLoadsWind);

P=elemloads(DLoads,Nodes,Elements,Types,DOF);

% Solve K * U = P
U=K\P;

figure
plotdisp(Nodes,Elements,Types,DOF,U(:,1),DLoads(:,1),Sections,Materials)
title('Displacements: own weight')

figure
plotdisp(Nodes,Elements,Types,DOF,U(:,2),DLoads(:,2),Sections,Materials)
title('Displacements: wind')

% Compute forces
[ForcesLCS,ForcesGCS]=elemforces(Nodes,Elements,Types,Sections,Materials,DOF,U,DLoads);

% Compute reaction forces for load case 1
Freac=reaction(Elements,ForcesGCS(:,1),[1.03; 2.03; 101.03; 102.03; 201.03; 202.03])

% Plot element forces for load case 1
figure
plotforc('norm',Nodes,Elements,Types,ForcesLCS(:,1),DLoads(:,1))
title('Normal forces: Own weight')
figure
plotforc('sheary',Nodes,Elements,Types,ForcesLCS(:,1),DLoads(:,1))
title('Shear forces along y: Own weight')
figure
plotforc('shearz',Nodes,Elements,Types,ForcesLCS(:,1),DLoads(:,1))
title('Shear forces along z: Own weight')
figure
plotforc('momx',Nodes,Elements,Types,ForcesLCS(:,1),DLoads(:,1))
title('Torsional moments: Own weight')
figure
plotforc('momy',Nodes,Elements,Types,ForcesLCS(:,1),DLoads(:,1))
title('Bending moments around y: Own weight')
figure
plotforc('momz',Nodes,Elements,Types,ForcesLCS(:,1),DLoads(:,1))
title('Bending moments around z: Own weight')

% Plot element forces for load case 2
figure
plotforc('norm',Nodes,Elements,Types,ForcesLCS(:,2),DLoads(:,2))
title('Normal forces: Wind')
figure
plotforc('sheary',Nodes,Elements,Types,ForcesLCS(:,2),DLoads(:,2))
title('Shear forces along y: Wind')
figure
plotforc('shearz',Nodes,Elements,Types,ForcesLCS(:,2),DLoads(:,2))

```



```

title('Shear forces along z: Wind')
figure
plotforc('momx',Nodes,Elements,Types,ForcesLCS(:, :, 2),DLoads(:, :, 2))
title('Torsional moments: Wind')
figure
plotforc('momy',Nodes,Elements,Types,ForcesLCS(:, :, 2),DLoads(:, :, 2))
title('Bending moments around y : Wind')
figure
plotforc('momz',Nodes,Elements,Types,ForcesLCS(:, :, 2),DLoads(:, :, 2))
title('Bending moments around z: Wind')

% Load combinations

% Safety factors
gamma_own=1.35;
gamma_wind=1.5;

% Combination factors
psi_wind=1;

% Load combination (Ultimate Limit State, ULS)
U_ULS=gamma_own*U(:, 1)+gamma_wind*psi_wind*U(:, 2);
Forces_ULS=gamma_own*ForcesLCS(:, :, 1)+gamma_wind*psi_wind*ForcesLCS(:, :, 2);
DLoads_ULS(:, 1)=DLoads(:, 1, 1)
DLoads_ULS(:, 2:7)=gamma_own*DLoads(:, 2:7, 1)+gamma_wind*psi_wind*DLoads(:, 2:7, 2);

figure
plotdisp(Nodes,Elements,Types,DOF,U_ULS,DLoads_ULS,Sections,Materials)

printdisp(Nodes,DOF,U_ULS);

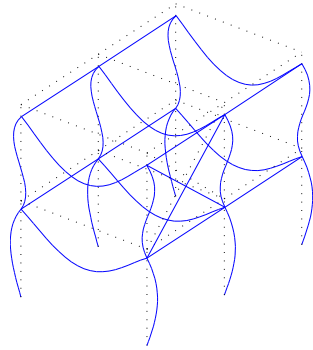
printforc(Elements,Forces_ULS);

% Plot element forces
figure
plotforc('norm',Nodes,Elements,Types,Forces_ULS,DLoads_ULS)
title('Normal forces: ULS')
figure
plotforc('sheary',Nodes,Elements,Types,Forces_ULS,DLoads_ULS)
title('Shear forces along y: ULS')
figure
plotforc('shearz',Nodes,Elements,Types,Forces_ULS,DLoads_ULS)
title('Shear forces along z: ULS')
figure
plotforc('momx',Nodes,Elements,Types,Forces_ULS,DLoads_ULS)
title('Torsional moments: ULS')
figure
plotforc('momy',Nodes,Elements,Types,Forces_ULS,DLoads_ULS)
title('Bending moments around y: ULS')
figure
plotforc('momz',Nodes,Elements,Types,Forces_ULS,DLoads_ULS)
title('Bending moments around z: ULS')

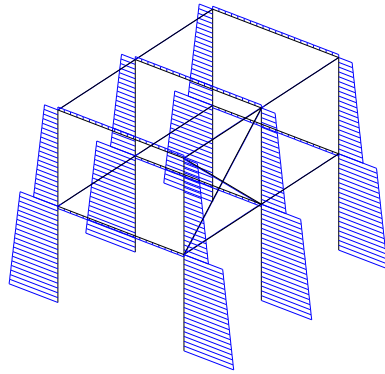
% Plot stresses
figure
plotstress('snorm',Nodes,Elements,Types,Sections,Forces_ULS,DLoads_ULS)
title('Normal stresses due to normal forces')
figure
plotstress('smomyt',Nodes,Elements,Types,Sections,Forces_ULS,DLoads_ULS)
title('Normal stresses due to bending moments around y: top')
figure
plotstress('smomyb',Nodes,Elements,Types,Sections,Forces_ULS,DLoads_ULS)
title('Normal stresses due to bending moments around y: bottom')
figure
plotstress('smomzt',Nodes,Elements,Types,Sections,Forces_ULS,DLoads_ULS)
title('Normal stresses due to bending moments around z: top')

```

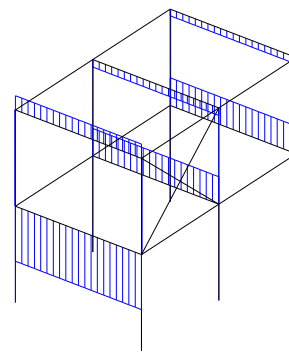
```
figure
plotstress('smomzb',Nodes,Elements,Types,Sections,Forces_ULS,DLoads_ULS)
title('Normal stresses due to bending moments around z: bottom')
figure
plotstress('smax',Nodes,Elements,Types,Sections,Forces_ULS,DLoads_ULS)
title('Maximal normal stresses')
figure
plotstress('smin',Nodes,Elements,Types,Sections,Forces_ULS,DLoads_ULS)
title('Minimal normal stresses')
```



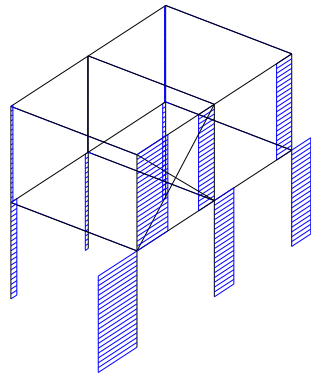
(a) Displacements



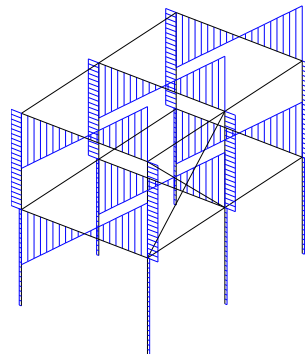
(b) Normal forces



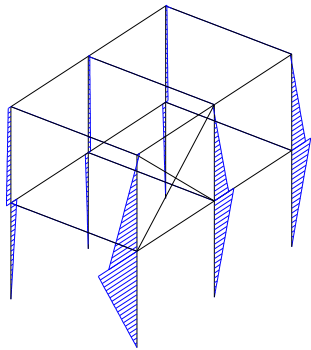
(c) Torsional moments



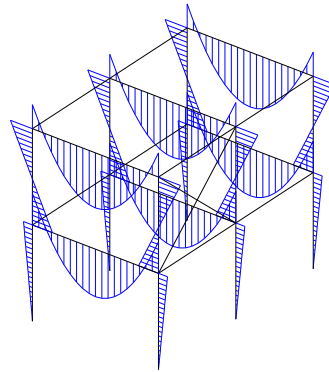
(d) Shear forces along z



(e) Shear forces along y



(f) Bending moments around y



(g) Bending moments around z

Figure 2.6: Results for load case 1 of example 1.2.

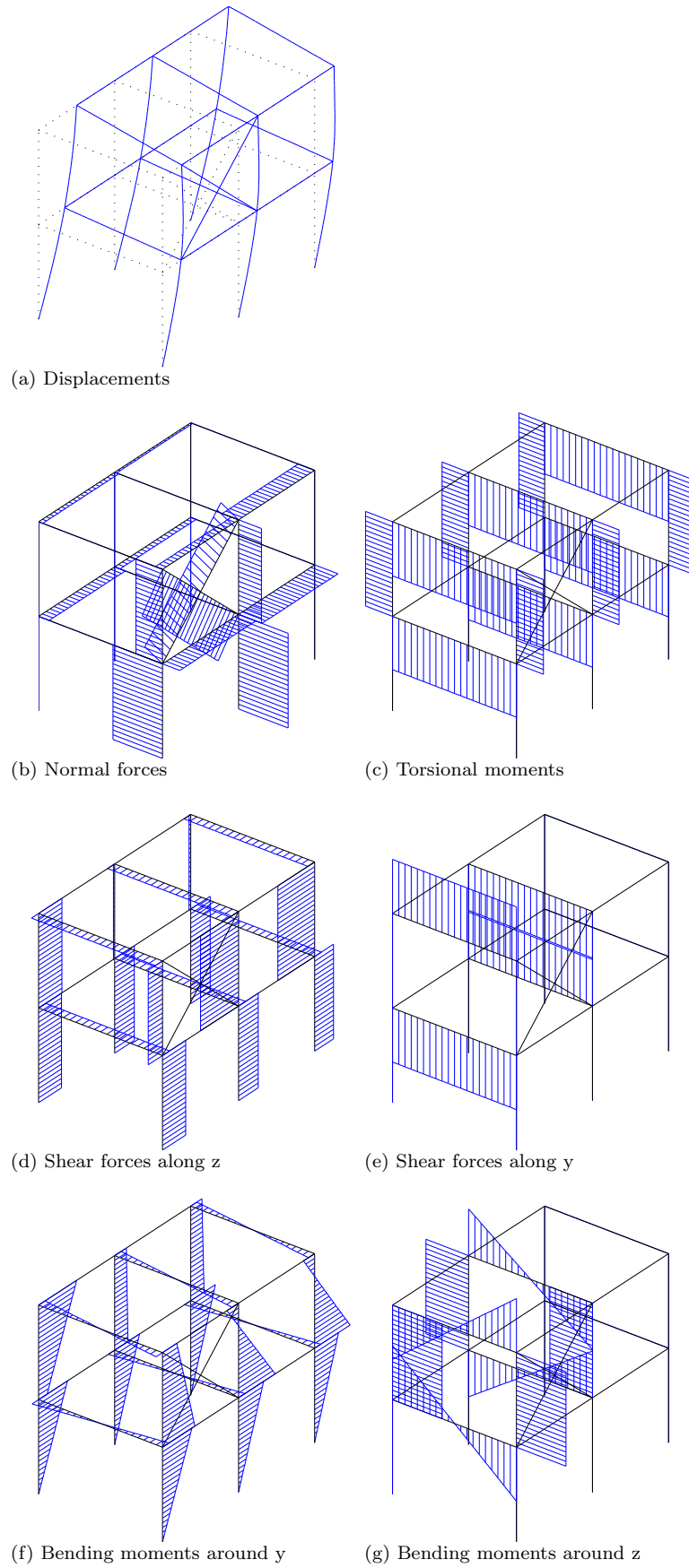


Figure 2.7: Results for load case 2 of example 1.2.

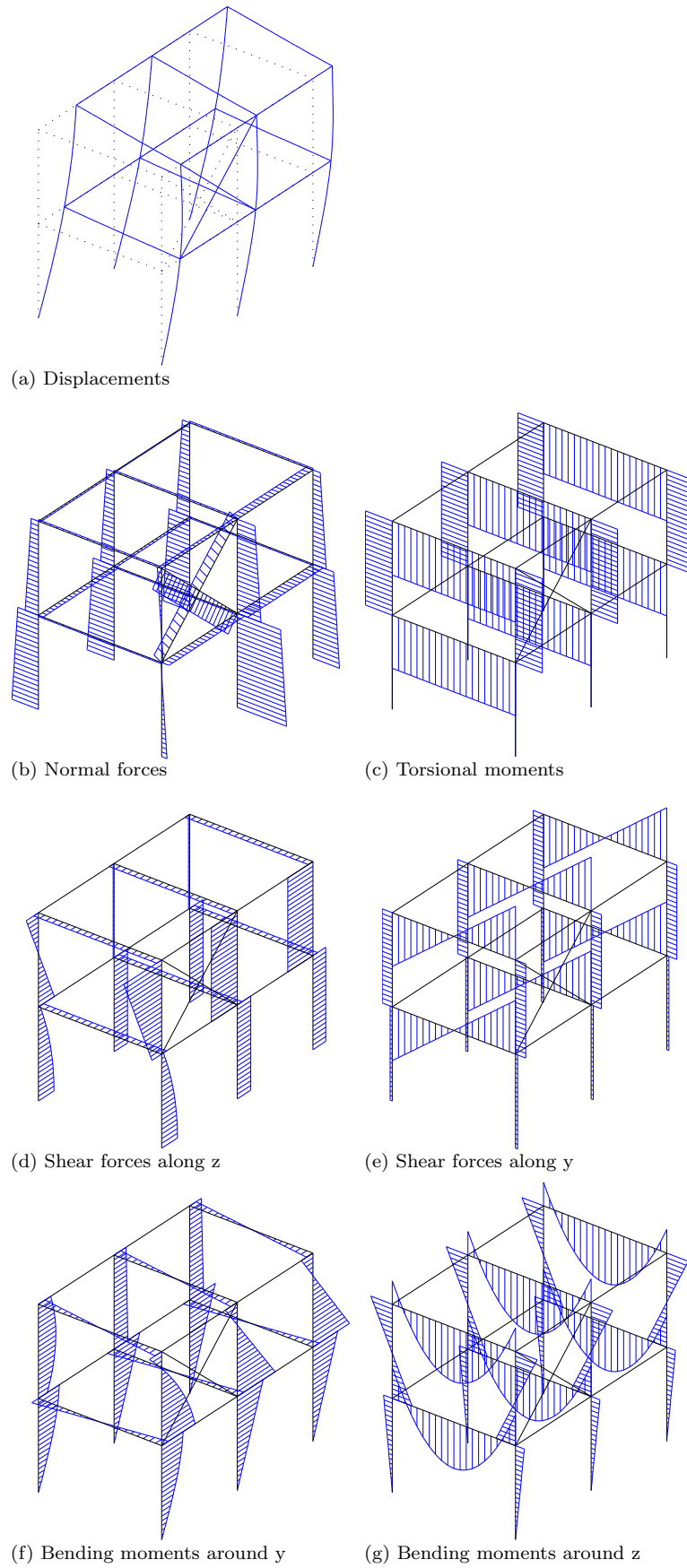


Figure 2.8: Results for load combination ULS of example 1.2.

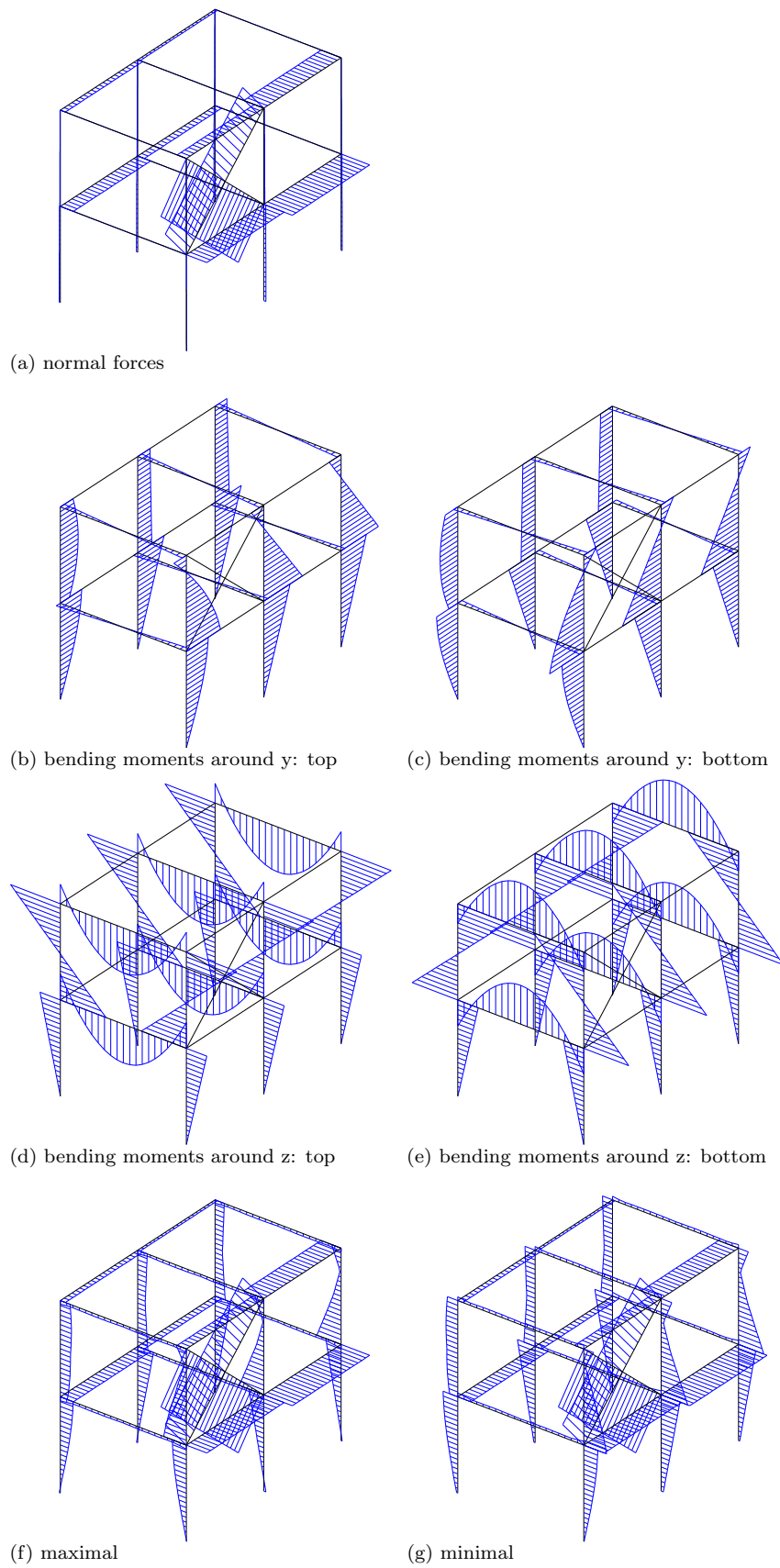


Figure 2.9: Normal stresses for load combination ULS of example 1.2.

2.3 Example 1.3: static analysis of a plate with a circular hole

This example demonstrates the use of StabIl for the solution of 2D plane strain continuum problems. A thin disk with a circular hole is subjected to uniaxial tension. The stress concentration around the hole is examined:

```
% StaBIL manual
% Example 1.3: static analysis of a plate with a circular hole

% Stress concentration around a circular hole in a disk.
r = 1; % Radius of hole [m]
L = 20; % Width of plate [m]
p = 1; % Uniform pressure [N/m^2]
n = 40; % Mesh parameter (nElem = m*n)
m = 20; % Mesh parameter
Types = {1 'shell8'}; % {EltTypeID EltName}
Sections = [1 1]; % [SecID t]
Materials = [1 200000 0]; % [MatID E nu]

% define lines
Line1 = [r 0 0; L/2 0 0];
Line2 = [L/2 0 0; L/2 L/2 0; 0 L/2 0];
Line3 = [0 L/2 0; 0 r 0];
Line4 = [r*sin((0:5).*pi/10) r*cos((0:5).*pi/10) zeros(6,1)];

[Nodes,Elements,Edge1,Edge2,Edge3,Edge4] = makemesh(Line1,Line2,Line3,Line4,...
    m,n,Types(1,:),Sections(1,1),Materials(1,1),'L2method','linear');

% Check mesh:
figure;
plotnodes(Nodes,'numbering','off');
hold('on')
plotelem(Nodes,Elements,Types,'numbering','off');
title('Nodes and elements');

% Select all dof:
DOF = getdof(Elements,Types);
% Apply boundary conditions:
% - Line1 and Line3: symmetry condition
seldof = [0.03;0.04;0.05;0.06;Edge1+0.02;Edge3+0.01];
DOF = removedof(DOF,seldof);

% Assemble K:
K = asmkm(Nodes,Elements,Types,Sections,Materials,DOF);

% Apply load to upper edge (equivalent nodal forces):
C = selectdof(DOF,Edge2(1:(end-1)/2+1)+0.02);
P = C.*[1/6; repmat([2/3; 1/3],((length(Edge2)-1)/2-2)/2,1); 2/3; 1/6]*p*L/m;
U = K\P;

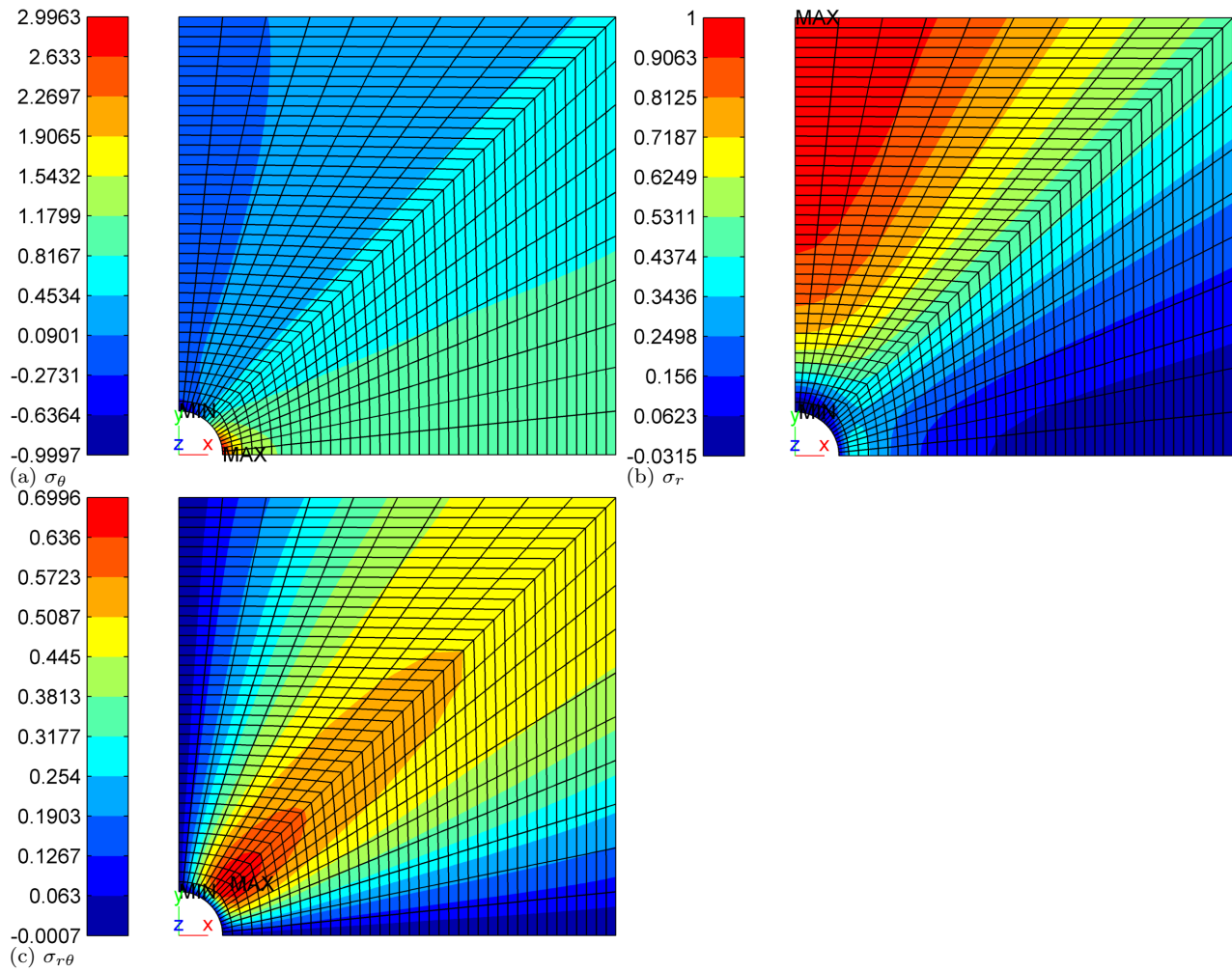
% Calculate stress in cylindrical global coordinate system
[SeGCS]=elemstress(Nodes,Elements,Types,Sections,Materials,DOF,U,'gcs','cyl');

% Get nodal stress from element solution
[SnGCS] = nodalstress(Nodes,Elements,Types,SeGCS);

% plot results
figure;
plotstresscontourf('sx',Nodes,Elements,Types,SnGCS)
title('\sigma_{r}')
figure;
plotstresscontourf('sy',Nodes,Elements,Types,SnGCS)
title('\sigma_{\theta}')
figure;
```

```
plotstresscontourf('sxy',Nodes,Elements,Types,SnGCS)
title('\sigma_{r\theta}')
```

Figure 2.10

**Figure 2.10:** Results for the hole-in-disk problem.

2.4 Example 1.4: static analysis of a barrel vault roof

A barrel vault roof subjected to its self weight is analysed. The curved edges are simply supported and the straight edges are free. Due to symmetry only a quarter of the roof is modelled and symmetry boundary conditions are applied.

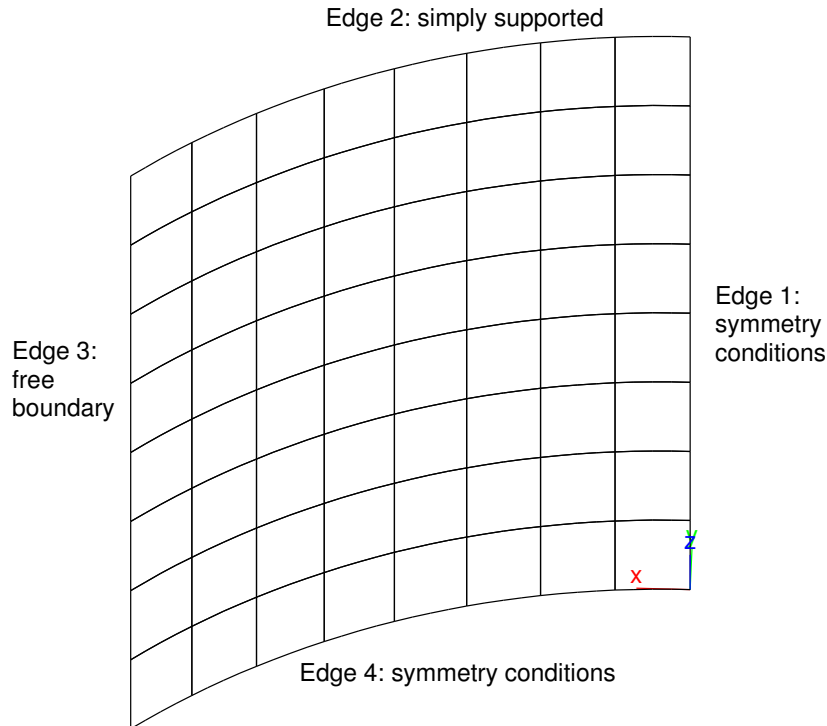


Figure 2.11: 8×8 mesh of a quarter of a cylindrical roof.

```
% StaBIL manual
% Example 1.4: static analysis of a barrel vault roof

% A BARREL VAULT ROOF SUBJECTED TO ITS SELF WEIGHT
% Reference: COOK, CONCEPTS AND APPL OF F.E.A., 2ND ED., 1981, PP. 284-287.

%% Parameters
R=25;           % Radius of cylindrical roof
L=50;           % Length of cylindrical roof
t=0.25;         % Thickness of roof
theta = 40*pi/180; % Angle of cylindrical roof
E = 4.32*10^8;  % Youngs modulus
nu = 0;         % Poisson coefficient
rho = 36.7347;  % Density
N = 8;          % Number of elements

%% Mesh

% Lines = [x1 y1 z1;x2 y2 z2;...]
Line1 = [0 0 0;0 0 L/2];
Line2 = [R*sin(theta*(0:0.1:1).') R*cos(theta*(0:0.1:1).')-R L*ones(11,1)/2];
Line3 = [R*sin(theta) R*cos(theta)-R L/2; R*sin(theta) R*cos(theta)-R 0];
Line4 = [R*sin(theta*(1:-0.1:0).') R*cos(theta*(1:-0.1:0).')-R zeros(11,1)];

% Specify element type for mesh
Materials = [1 E nu rho];
Sections = [1 t];
```

```

Types = {1 'shell8'};

% Mesh the area between lines 1,2,3,4 with N * N elements of type shell8,
% section number 1 and material number 1.
[Nodes,Elements,Edge1,Edge2,Edge3,Edge4] = ...
    makemesh(Line1,Line2,Line3,Line4,N,N,Types(1,:),1,1);

% Check mesh:
figure;
plotnodes(Nodes,'numbering','off');
hold('on')
plotelem(Nodes,Elements,Types,'numbering','off');
title('Nodes and elements');

%% Assemble stiffness matrix

% Select all dof:
DOF = getdof(Elements,Types);

% Apply boundary conditions:
% - Line1 and Line4: symmetry condition
% - Line2: simply supported
sdof = [Edge1+0.01;Edge1+0.06;Edge1+0.05;Edge2+0.02;Edge2+0.01;
        Edge2+0.06;Edge4+0.03;Edge4+0.04;Edge4+0.05];
DOF = removedof(DOF,sdof);

% Assemble K:
K = asmkk(Nodes,Elements,Types,Sections,Materials,DOF);

%% Solution

% Apply gravitational acceleration and determine equivalent nodal forces:
DLoads=accel([0 9.8 0],Elements,Types,Sections,Materials);
P=elemloads(DLoads,Nodes,Elements,Types,DOF);

% Solve K * U = P:
U = K\P;

% Plot displacements:
figure;
plotdisp(Nodes,Elements,Types,DOF,U)
title('Displacements')

% Check target displacement:
TP1 = selectdof(DOF,intersect(Edge3,Edge4)+0.02);
Utp1 = TP1*U
ratio_u = -Utp1/0.3016

%% Stress

% Determine element stress in global and local(element) coordinate system:
[SeGCS,SeLCS,vLCS]=elemstress(Nodes,Elements,Types,Sections,Materials,DOF,U);

% print stress:
printstress(Elements,SeGCS)

% plot stress contour:
figure;
plotstresscontour('sx',Nodes,Elements,Types,SeGCS,'location','bot')
title('sx in gcs (element solution)')

% plot filled contours:
figure;
plotstresscontourf('sx',Nodes,Elements,Types,SeGCS,'location','bot')
title('sx in gcs (element solution)')

```

```

figure;
plotstresscontour('sx',Nodes,Elements,Types,SeLCS,'location','bot')
title('sx in lcs')

% plot lcs for shell elements
figure;
plotlcs(Nodes,Elements,Types,vLCS)
title('local coordinate system')

% Calculate nodal solution
% SnGCS: stress arranged per element
% SnGCS2: stress arranged per node
[SnGCS,SnGCS2]=nodalstress(Nodes,Elements,Types,SeGCS);
[SnLCS,SnLCS2]=nodalstress(Nodes,Elements,Types,SeLCS);
figure;
plotstresscontour('sx',Nodes,Elements,Types,SnGCS,'location','bot');
title('sx in gcs (nodal solution)')

% Stress ratios:

ratio_sz = SnGCS2(intersect(Edge3,Edge4),16)/358420
ratio_st = SnGCS2(intersect(Edge4,Edge1),14)/(-213400)

%% Shell forces

% Shell forces (element solution):
[FeLCS]=elemshellf(Elements,Sections,SeLCS);
figure;
plotshellfcontour('my',Nodes,Elements,Types,FeLCS)
title('my (element solution)')

% Nodal solution:
[FnLCS,FnLCS2]=nodalshellf(Nodes,Elements,Types,FeLCS);
figure;
plotshellfcontour('my',Nodes,Elements,Types,FnLCS)
title('my (nodal solution)')

%% Principal stress

% Principal stresses:
[Spr,Vpr]=principalstress(Elements,SnGCS);
figure;
plotstresscontour('s1',Nodes,Elements,Types,Spr,'location','bot');
title('s1')
% plot principal stresses:
figure;
plotprincstress(Nodes,Elements,Types,Spr,Vpr)
title('principal stresses')

```

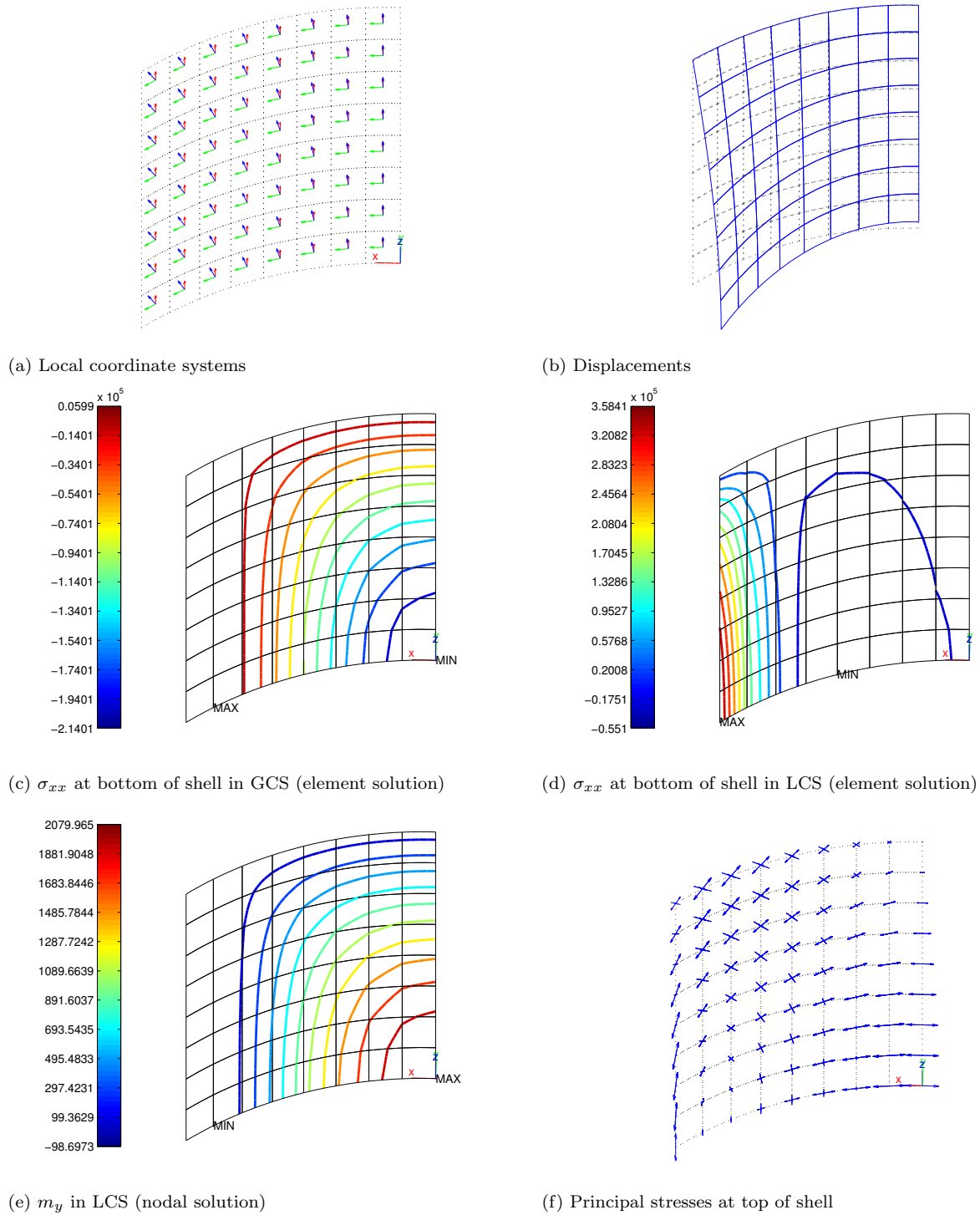


Figure 2.12: Cylindrical roof: results.

Chapter 3

Dynamic analysis of structures

3.1 Example 2.1: dynamic analysis of a frame

The frame that was treated in section 2.1 is reconsidered, computing the eigenmodes and dynamic response of the structure.

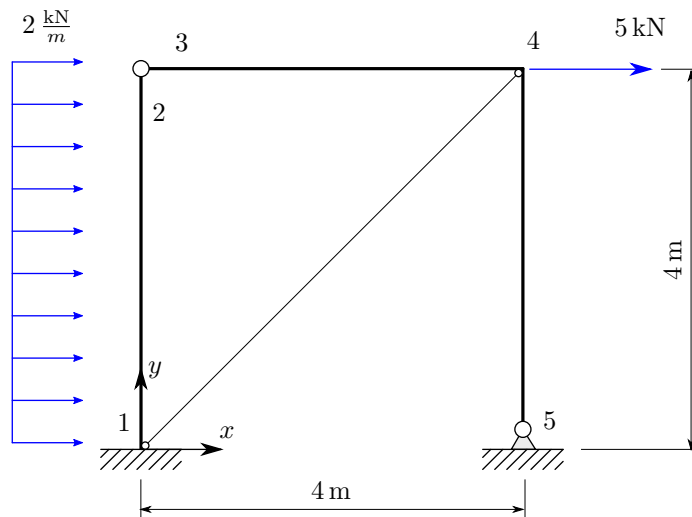


Figure 3.1: Simple frame structure.

3.1.1 Model

```
% StaBIL manual
% Example 2.1: dynamic analysis: model
% Units: m, N

L=4;
H=4;
nElemCable=8;

% Nodes=[NodID X Y Z]
Nodes= [1 0 0 0;
        2 L 0 0];
Nodes=reprow(Nodes,1:2,4,[2 0 H/4 0]);
Nodes= [Nodes;
        11 0 H 0];
Nodes=reprow(Nodes,11,3,[1 L/4 0 0]);
Nodes= [Nodes;
        15 1 5 0];           % reference node
Nodes= [Nodes;
```

```

16 0 0 0];
Nodes=repro(Nodes,16,nElemCable,[1 L/nElemCable H/nElemCable 0]);

% Check the node coordinates as follows:
figure
plotnodes(Nodes);

% Element types -> {EltTypID EltName}
Types= {1 'beam'};

b=0.10;
h=0.25;
r=0.004;

% Sections=[SecID A ky kz Ixx Iyy Izz]
Sections= [1 b*h Inf Inf 0 0 b*h^3/12;
           2 pi*r^2 Inf Inf 0 0 pi*r^4/4];

% Materials=[MatID E nu];
Materials= [1 30e9 0.2 2500; % concrete
            2 210e9 0.3 7850]; % steel

% Elements=[EltID TypID SecID MatID n1 n2 n3]
Elements= [1 1 1 1 1 3 15;
           2 1 1 1 1 2 4 15];
Elements=repro(Elements,1:2,3,[2 0 0 0 2 2 0]);
Elements=[ Elements;
          9 1 1 1 11 12 15;
          10 1 1 1 12 13 15;
          11 1 1 1 13 14 15;
          12 1 1 1 14 10 15;
          13 1 2 2 16 17 15];
Elements=repro(Elements,13,(nElemCable-1),[1 0 0 0 1 1 0]);

% Check node and element definitions as follows:
hold('on');
plotelem(Nodes,Elements,Types);
title('Nodes and elements');

% Degrees of freedom
DOF=getdof(Elements,Types);

% Boundary conditions
seldof=[0.03; 0.04; 0.05; 1.01; 1.02; 1.06; 2.01; 2.02; 16.01; 16.02];

DOF=removedof(DOF,seldof);

% Assembly of stiffness matrix K
[K,M]=asmkm(Nodes,Elements,Types,Sections,Materials,DOF);

% DLoads=[EltID n1globalX n1globalY n1globalZ ...]
DLoads=[1 2000 0 0 2000 0 0;
        3 2000 0 0 2000 0 0;
        5 2000 0 0 2000 0 0;
        7 2000 0 0 2000 0 0];
b=elemloads(DLoads,Nodes,Elements,Types,DOF); % Spatial distribution, nodal (nDOF * 1)

% Constraint equations: Constant=Coef1*DOF1+Coef2*DOF2+ ...
% Constraints=[Constant Coef1 DOF1 Coef2 DOF2 ...]
Constr=[0 1 9.01 -1 11.01;
        0 1 9.02 -1 11.02;
        0 1 10.01 -1 (16.01+nElemCable);
        0 1 10.02 -1 (16.02+nElemCable)];

[K,b,M]=addconstr(Constr,DOF,K,b,M);

```

3.1.2 Eigenmodes

```
% StaBIL manual
% Example 2.1: dynamic analysis: eigenvalue problem
% Units: m, N

% Assembly of M and K
tutorialdyna;

% Eigenvalue problem
nMode=12;
[phi,omega]=eigfem(K,M,nMode);

% Display eigenfrequencies
disp('Lowest eigenfrequencies [Hz]');
disp(omega/2/pi);

% Plot eigenmodes
figure;
plotdisp(Nodes,Elements,Types,DOF,phi(:,1),'DispMax','off')
figure;
plotdisp(Nodes,Elements,Types,DOF,phi(:,2),'DispMax','off')
figure;
plotdisp(Nodes,Elements,Types,DOF,phi(:,5),'DispMax','off')
figure;
plotdisp(Nodes,Elements,Types,DOF,phi(:,8),'DispMax','off')
figure;
plotdisp(Nodes,Elements,Types,DOF,phi(:,11),'DispMax','off')
figure;
plotdisp(Nodes,Elements,Types,DOF,phi(:,12),'DispMax','off')

% Animate eigenmodes
figure;
animdisp(Nodes,Elements,Types,DOF,phi(:,1))
title('Eigenmode 1')

figure;
animdisp(Nodes,Elements,Types,DOF,phi(:,2))
title('Eigenmode 2')

figure;
animdisp(Nodes,Elements,Types,DOF,phi(:,5))
title('Eigenmode 5')

figure;
animdisp(Nodes,Elements,Types,DOF,phi(:,8))
title('Eigenmode 8')

figure;
animdisp(Nodes,Elements,Types,DOF,phi(:,11))
title('Eigenmode 11')

figure;
animdisp(Nodes,Elements,Types,DOF,phi(:,12))
title('Eigenmode 12')
```

3.1.3 Modal superposition: time domain: piecewise exact integration

```
% StaBIL manual
% Example 2.1: dynamic analysis: modal superposition: piecewise exact integration
% Units: m, N

% Assembly of M and K
```

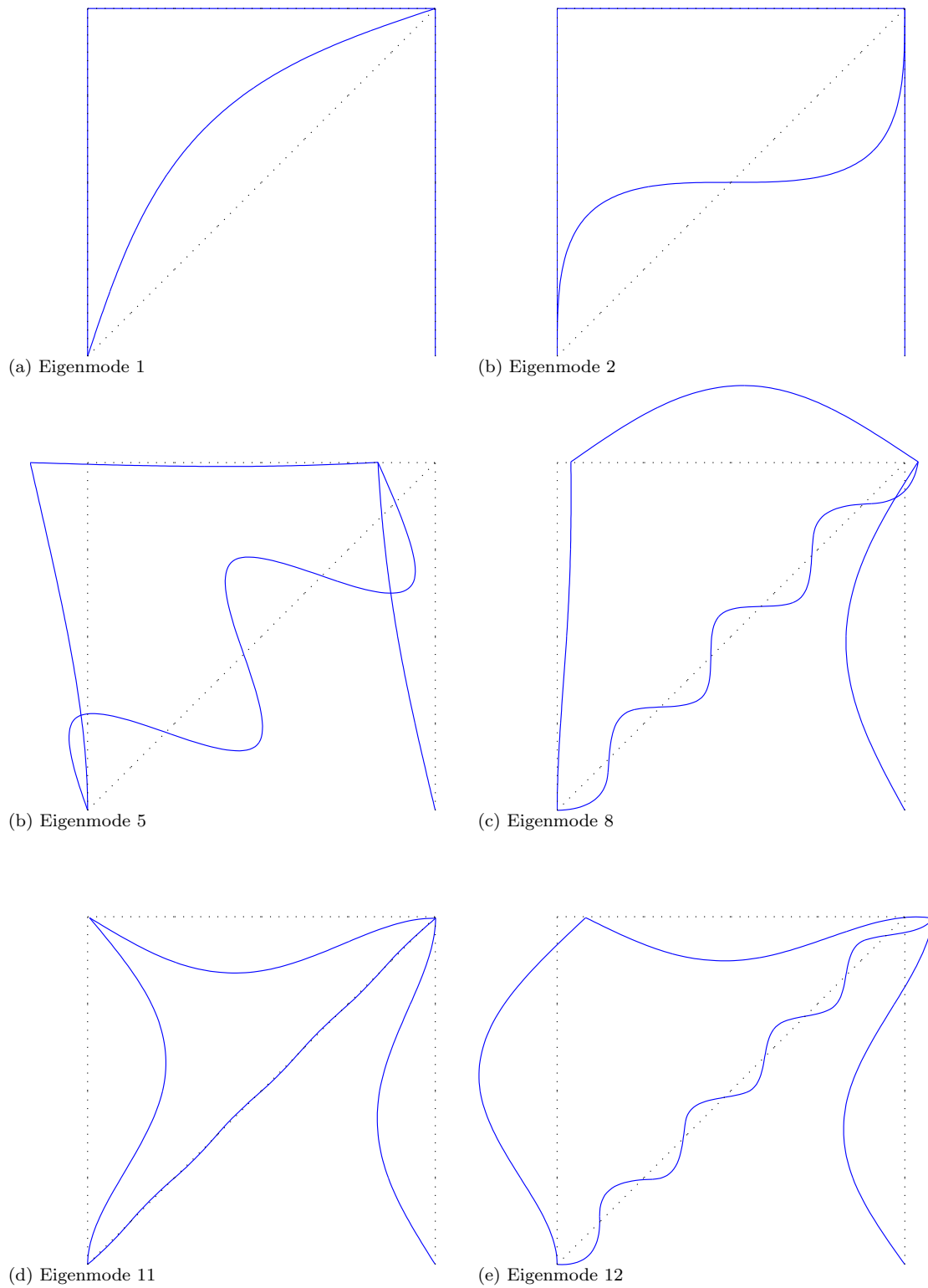


Figure 3.2: Results of example 2.1.

```
tutorialdyna;

% Sampling parameters
T=2.5;           % Time window
dt=0.002;        % Time step
N=T/dt;          % Number of samples
```



```

t=[0:N-1]*dt; % Time axis

% Eigenvalue analysis
nMode=12; % Number of modes to take into account
[phi,omega]=eigfem(K,M,nMode); % Calculate eigenmodes and eigenfrequencies
xi=0.07; % Constant modal damping ratio

% Excitation
bm=phi.'*b; % Spatial distribution, modal (nMode * 1)
q=zeros(1,N); % Time history (1 * N)
q((t>=0.50) & (t<0.60))=1; % Time history (1 * N)
pm=bm*q; % Modal excitation (nMode * N)

% Modal analysis
x=msupt(omega,xi,t,pm,'zoh');

% Modal displacements -> nodal displacements
u=phi*x; % Nodal response (nDOF * N)

% Figures
figure;
plot(t,x);
title('Modal response (piecewise linear exact integration)');
xlabel('Time [s]');
xlim([0 4.1])
ylabel('Displacement [m kg^{0.5}]');
legend([repmat('Mode ',nMode,1) num2str([1:nMode].')]);

figure;
c=selectdof(DOF,[9.01; 13.02; 17.02]);
plot(t,c*u);
title('Nodal response (piecewise linear exact integration)');
xlabel('Time [s]');
xlim([0 4.1])
ylabel('Displacement [m]');
legend('9.01','13.02','17.02');

% Movie
figure;
animdisp(Nodes,Elements,Types,DOF,u);

% Display
disp('Maximum modal response');
disp(max(abs(x),[],2));

disp('Maximum nodal response 9.01 13.02 17.02');
disp(max(abs(c*u),[],2));

```

3.1.4 Modal superposition: transform to frequency domain

```

% StaBIL manual
% Example 2.1: dynamic analysis: direct method: frequency domain
% Units: m, N

% Assembly of M and K
tutorialdyna;

% Sampling parameters
N=2048; % Number of samples
dt=0.002; % Time step
T=N*dt; % Period
F=N/T; % Sampling frequency
df=1/T; % Frequency resolution
t=[0:N-1]*dt; % Time axis

```

```

f=[0:N/2-1]*df;           % Positive frequencies corresponding to FFT [Hz]
Omega=2*pi*f;             % Idem [rad/s]

% Eigenvalue analysis
nMode=12;                 % Number of modes to take into account
[phi,omega]=eigfem(K,M,nMode); % Calculate eigenmodes and eigenfrequencies
xi=0.07;                  % Constant modal damping ratio

% Excitation
bm=phi.*b;                % Spatial distribution, modal (nMode * 1)
q=zeros(1,N);             % Time history (1 * N)
q((t>=0.50) & (t<0.60))=1; % Time history (1 * N)
Q=fft(q);                 % Frequency content (1 * N)
Q=Q(1:N/2);               % Frequency content, positive freq (1 * N/2)
Pm=bm*Q;                  % Modal excitation, positive freq (nMode * N/2)

% Modal analysis
[X,H]=msupf(omega,xi,Omega,Pm); % Modal response, positive freq (nMode * N/2)

% F-dom -> t-dom
X=[X, zeros(nMode,1), conj(X(:,end:-1:2))];
x=ifft(X,[],2);           % Modal response (nMode * N)

% Modal displacements -> nodal displacements
u=phi*x;                  % Nodal response (nDOF * N)

% Figures
figure;
subplot(3,2,1);
plot(t,q,'.-');
xlim([0 4.1]);
ylim([0 1.2]);
title('Excitation time history');
xlabel('Time [s]');
ylabel('Force [N/m]');

subplot(3,2,2);
plot(f,abs(Q)/F,'.-');
title('Excitation frequency content');
xlabel('Frequency [Hz]');
ylabel('Force [N/m/Hz]');

subplot(3,2,4);
plot(f,abs(H),'.-');
title('Modal transfer function');
xlabel('Frequency [Hz]');
ylabel('Displacement [m/N]');
legend([repmat('Mode ',nMode,1) num2str(1:nMode).']);

subplot(3,2,6);
plot(f,abs(X(:,1:N/2))/F,'.-');
title('Modal response');
xlabel('Frequency [Hz]');
ylabel('Displacement [m kg^{0.5}/Hz]');

subplot(3,2,5);
plot(t,x);
title('Modal response (calculation in f-dom)');
xlabel('Time [s]');
xlim([0 4.1]);
ylabel('Displacement [m kg^{0.5}]');

figure;
plot(t,x);
title('Modal response (calculation in f-dom)');
xlabel('Time [s]');

```

```

xlim([0 4.1])
ylabel('Displacement [m kg^{0.5}]');
legend([repmat('Mode ',nMode,1) num2str([1:nMode].')]);

figure;
c=selectdof(DOF,[9.01; 13.02; 17.02]);
plot(t,c*u);
title('Nodal response (calculation in f-dom)');
xlabel('Time [s]');
xlim([0 4.1])
ylabel('Displacement [m]');
legend('9.01','13.02','17.02');

% Movie
figure;
animdisp(Nodes,Elements,Types,DOF,u);

% Display
disp('Maximum modal response');
disp(max(abs(x),[],2));

disp('Maximum nodal response 9.01 13.02 17.02');
disp(max(abs(c*u),[],2));

```

3.1.5 Direct time integration

```

% StaBIL manual
% Example 2.1: dynamic analysis: direct time integration: trapezium rule
% Units: m, N

% Assembly of M, K and C
tutorialdyna;
[phi,omega]=eigfem(K,M);           % Calculate eigenmodes and eigenfrequencies
xi=0.07;                           % Damping ratio
nModes=length(K)-size(Constr,1);
C=M.*phi(:,1:nModes)*diag(2*xi*omega(1:nModes))*phi(:,1:nModes).'*M;
                                   % Modal -> full damping matrix C

% Sampling parameters
T=2.5;                             % Time window
dt=0.002;                          % Time step
N=T/dt;                            % Number of samples
t=[0:N-1]*dt;                     % Time axis

% Excitation
q=zeros(1,N);                      % Time history (1 * N)
q((t>=0.50) & (t<0.60))=1;        % Time history (1 * N)
p=b*q;                             % Nodal excitation (nDOF * N)

% Direct time integration - trapezium rule
alpha=1/4;
delta=1/2;
theta=1;
u=newmark(M,C,K,dt,p,[alpha delta theta]);

% Figures
figure;
c=selectdof(DOF,[9.01; 13.02; 17.02]);
plot(t,c*u);
title(['Nodal response (direct time integration)']);
xlabel('Time [s]');
xlim([0 4.1])
ylabel('Nodal displacements [m]');
legend('9.01','13.02','17.02');

```

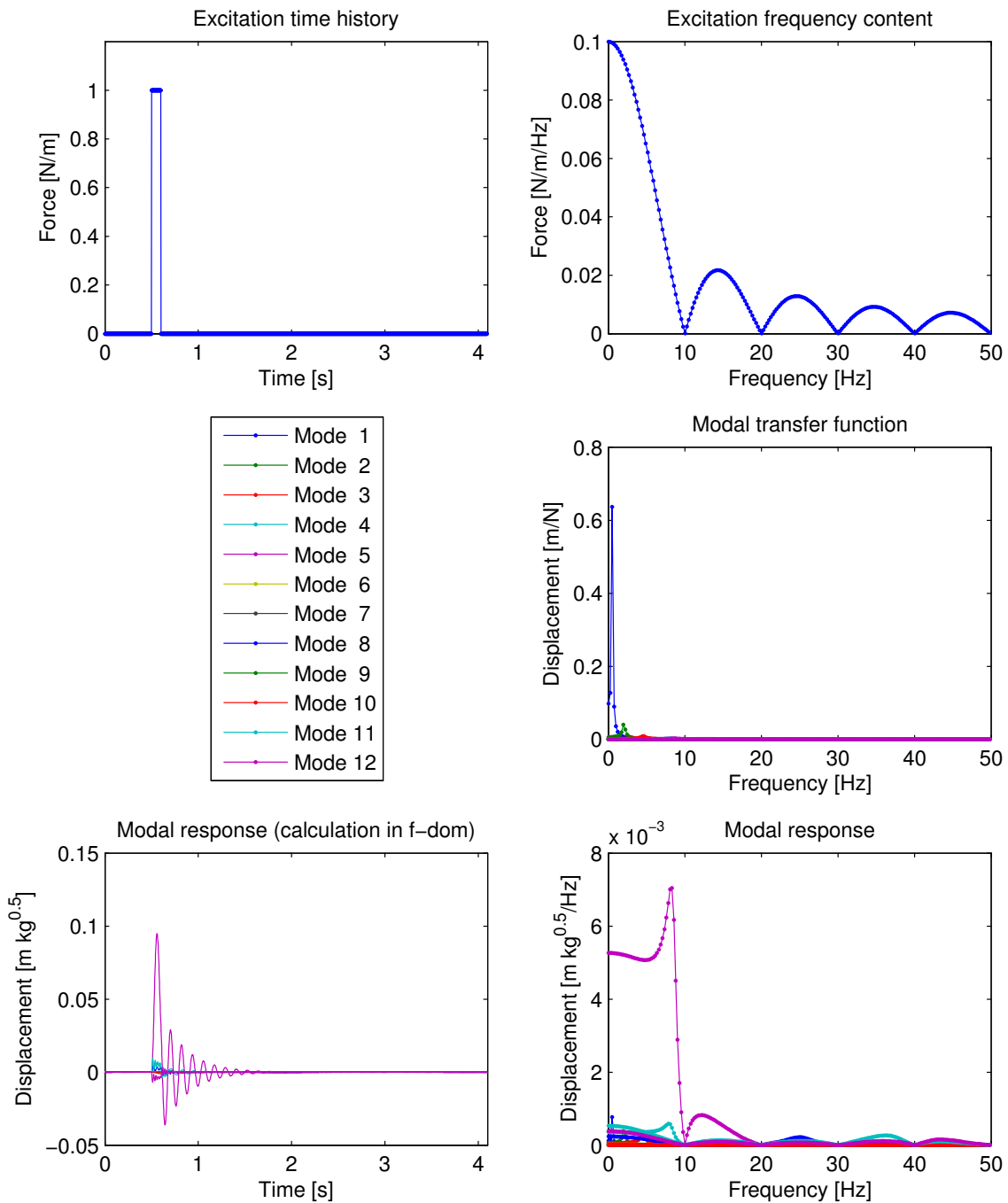


Figure 3.3: Modal superposition in frequency domain.

```
% Movie
figure;
animdisp(Nodes,Elements,Types,DOF,u);

% Display
disp('Maximum nodal response 9.01 13.02 17.02');
disp(max(abs(c*u), [], 2));
```

3.1.6 Direct solution in the frequency domain

```
% StaBIL manual
% Example 2.1: dynamic analysis: modal superposition: transform to f-dom
```

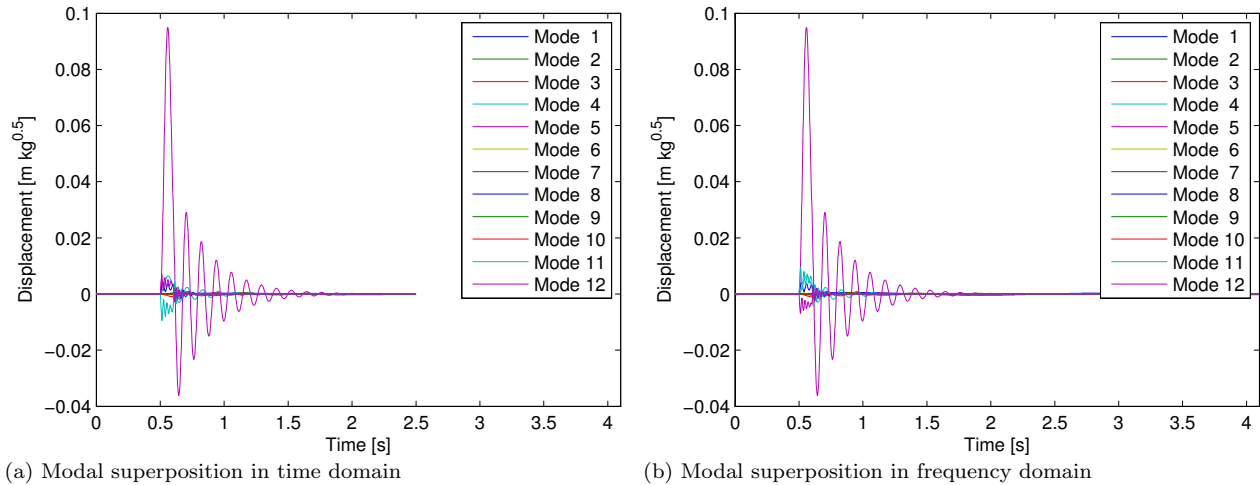


Figure 3.4: Modal response: comparison.

```
% Units: m, N

% Assembly of M, K and C
tutorialdyna;
[phi,omega]=eigfem(K,M);           % Calculate eigenmodes and eigenfrequencies
xi=0.07;                           % Damping ratio
nModes=length(K)-size(Constr,1);
C=M.*phi(:,1:nModes)*diag(2*xi*omega(1:nModes))*phi(:,1:nModes).'*M;
                                   % Modal -> full damping matrix C

% Sampling parameters
N=2048;                            % Number of samples
dt=0.002;                          % Time step
T=N*dt;                            % Period
F=N/T;                             % Sampling frequency
df=1/T;                            % Frequency resolution
t=[0:N-1]*dt;                     % Time axis
f=[0:N/2-1]*df;                   % Positive frequencies corresponding to FFT [Hz]
Omega=2*pi*f;                     % Idem [rad/s]

% Excitation
q=zeros(1,N);                     % Time history (1 * N)
q((t>=0.50) & (t<0.60))=1;       % Time history (1 * N)
Q=fft(q);                          % Frequency content (1 * N)
Q=Q(1:N/2);                       % Frequency content, positive freq (1 * N/2)
Pd=b*Q;                           % Nodal excitation, positive freq (nDOF * N/2)

% Solution for each frequency
Ud=zeros(size(Pd));
for k=1:N/2
    Kd=-Omega(k)^2*M+Omega(k)*i*C+K;
    Ud(:,k)=Kd\Pd(:,k);
end

% F-dom -> t-dom
Ud=[Ud, zeros(length(K),1), conj(Ud(:,end:-1:2))];
u=ifft(Ud,[],2);                  % Nodal response (nDOF * N)

% Figures
figure;
c=selectdof(DOF,[9.01; 13.02; 17.02]);
plot(t,c*u);
title('Nodal response (direct method in f-dom)');
xlabel('Time [s]');
xlim([0 4.1])
```

```

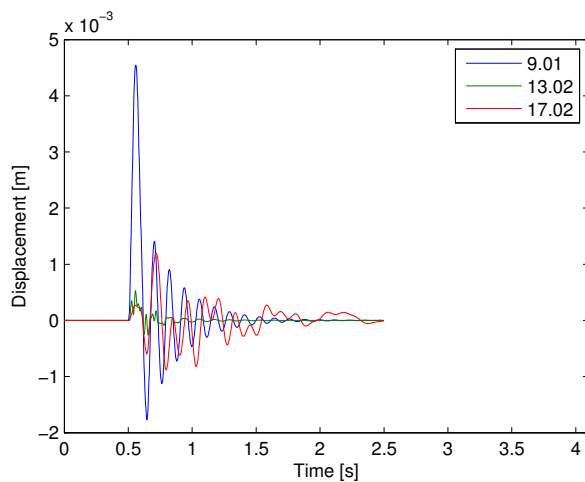
ylabel('Displacement [m]');
legend('9.01','13.02','17.02');

% Movie
figure;
animdisp(Nodes,Elements,Types,DOF,u);

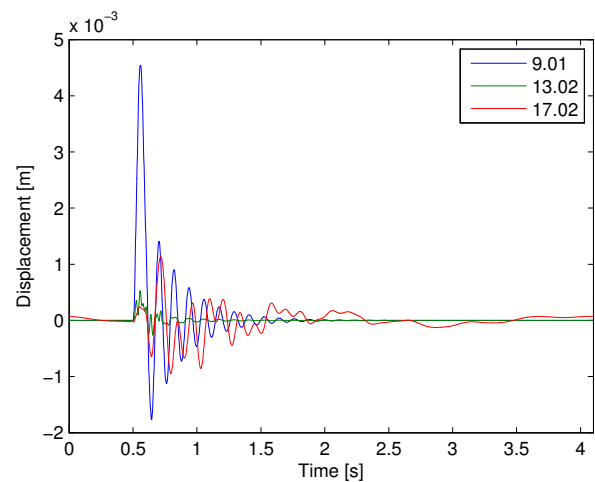
% Display
disp('Maximum nodal response 9.01 13.02 17.02');
disp(max(abs(c*u),[],2));

```

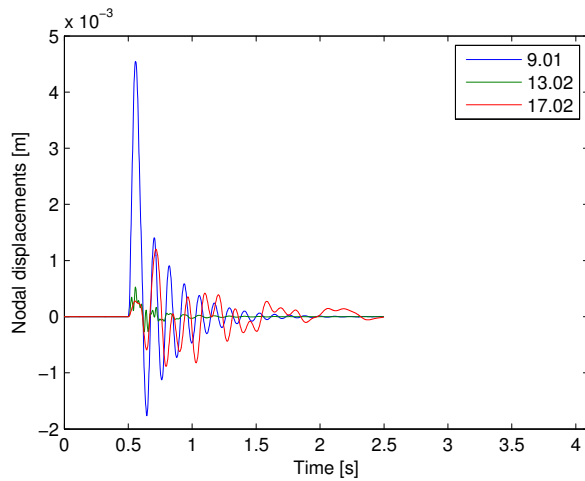
3.1.7 Comparison



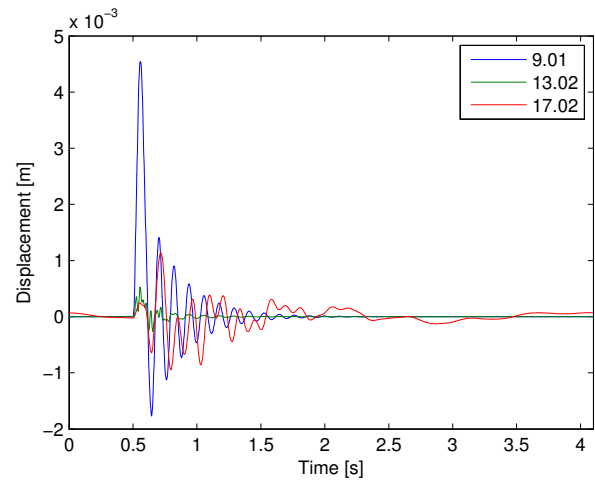
(a) Modal superposition in time domain



(b) Modal superposition in frequency domain



(b) Direct time integration



(c) Direct method in frequency domain

Figure 3.5: Nodal response: comparison.

Computational cost in seconds

	Time domain	Freq domain
Modal superposition	0.167	0.147
Direct method	0.213	0.789

3.2 Example 2.2: dynamic analysis of a plate

In this example the eigenfrequencies of a simply supported rectangular plate are calculated using shell4 and compared with the theoretical solution.

```
% StaBIL manual
% Example 2.2: dynamic analysis of a plate

% dynamic plate problem (dkt element)

% parameters

Lx = 10;           % length x-direction
Ly = 10;           % length y-direction
t = 1;             % thickness plate
E = 200000;        % E-modulus
nu = 0.3;          % Poisson-coeff.
rho = 7000;        % mass density
m = 20;            % number of elements in x-direction
n = 20;            % number of elements in y-direction
Types = {1 'shell4'}; % {EltTypID EltName}
Sections = [1 t];   % [SecID t]
Materials = [1 E nu rho]; % [MatID E nu]

% mesh

Line1 = [0 0 0;Lx 0 0];
Line2 = [Lx 0 0;Lx Ly 0];
Line3 = [Lx Ly 0;0 Ly 0];
Line4 = [0 Ly 0;0 0 0];

[Nodes,Elements,Edge1,Edge2,Edge3,Edge4] = makemesh(Line1,Line2,Line3,Line4,n,m,Types(1,:),1,1);
figure;
plotnodes(Nodes);
figure;
plotelem(Nodes,Elements,Types);

% DOFs (simply supported plate)

DOF = getdof(Elements,Types);
sdof = [0.01;0.02;0.06;[Edge1;Edge2;Edge3;Edge4]+0.03;[Edge1;Edge3]+0.05;[Edge2;Edge4]+0.04];
DOF = removedof(DOF,sdof);

% K & M

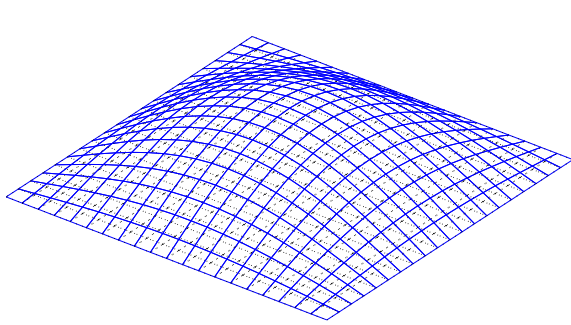
[K,M] = asmkm(Nodes,Elements,Types,Sections,Materials,DOF);

% eigenmodes

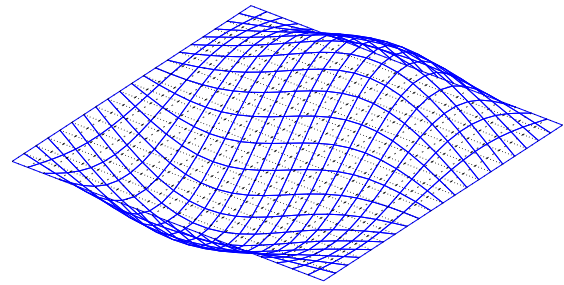
nMode = 10;
[phi,omega] = eigfem(K,M,nMode);
figure;
animdisp(Nodes,Elements,Types,DOF,phi(:,1));

% analytical solution

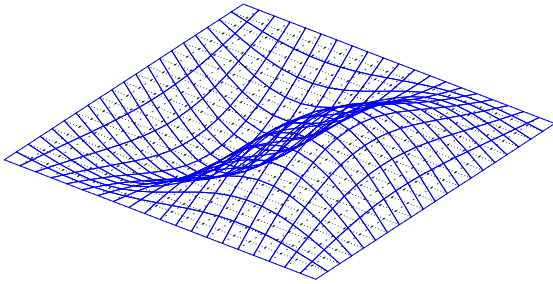
[mm,nn] = meshgrid((1:m),(1:n));
aomega = sqrt(E*t^2/(12*(1-nu^2)*rho))*((mm*pi/Lx).^2+(nn*pi/Ly).^2);
aomega = reshape(aomega,numel(aomega),1);
aomega = sort(aomega);
ratio = omega./aomega(1:nMode)
```



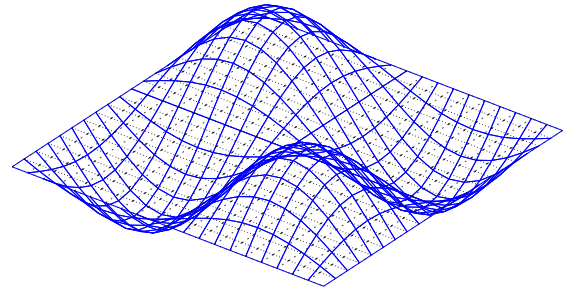
(a) 0.1268 Hz



(b) 0.3170 Hz



(c) 0.3170 Hz



(d) 0.5050 Hz

Figure 3.6: The first four eigenmodes of a thin plate.

$f_{\text{FEM}}[\text{Hz}]$	$f_{\text{analytical}}[\text{Hz}]$	$f_{\text{FEM}}/f_{\text{analytical}}$
0.1268	0.1270	0.9983
0.3170	0.3176	0.9982
0.3170	0.3176	0.9982
0.5050	0.5082	0.9939
0.6355	0.6352	1.0004
0.6355	0.6352	1.0004
0.8202	0.8258	0.9932
0.8202	0.8258	0.9932
1.0848	1.0799	1.0046
1.0848	1.0799	1.0046

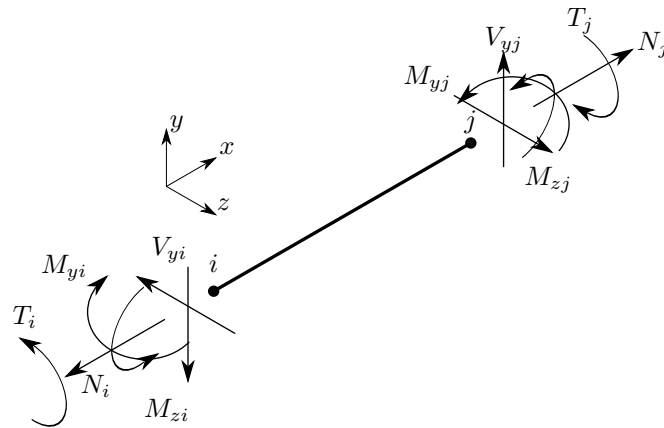
Table 3.1: Eigenfrequencies of a thin plate.

Chapter 4

Element guide

This chapter presents an overview of the element types available in Stabil. The conventions (local coordinate system, nodal connectivity) for each element type are given and reference is made to the Stabil functions related to the element implementation.

beam



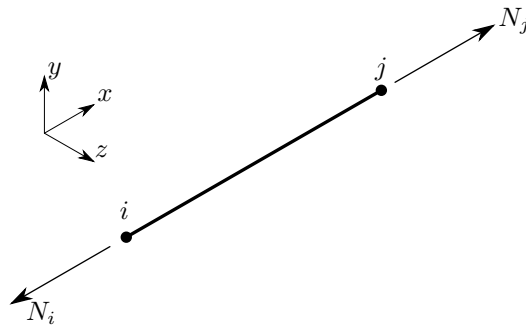
Description

Euler-Bernoulli beam element with a cubic interpolation of the beam deflection.

Functions

dof_beam	Element degrees of freedom for a beam element.	p.117
ke_beam	Beam element stiffness and mass matrix in global coordinate system.	p.168
kelcs_beam	Beam element stiffness and mass matrix in local coordinate system.	p.164
trans_beam	Transform coordinate system for a beam element.	p.291
loads_beam	Equivalent nodal forces for a beam element in the GCS.	p.192
loadslcs_beam	Equivalent nodal forces for a beam element in the LCS.	p.189
accel_beam	Compute the distributed loads for a beam due to an acceleration.	p.59
forces_beam	Compute the element forces for a beam element.	p.153
forceslcs_beam	Compute the element forces for a beam element in the LCS.	p.151
nelcs_beam	Shape functions for a beam element.	p.206
nedloadlcs_beam	Shape functions for a distributed load on a beam element.	p.205
coord_beam	Coordinates of the beam elements for plotting.	p.78
disp_beam	Return matrices to compute the displacements of the deformed beams.	p.99
dispgcs2lcs_beam	Transform the element displacements to the LCS for a beam.	p.97
fdiagrgcs_beam	Return matrices to plot the forces in a beam element.	p.144
fdiagrlcs_beam	Force diagram for a beam element in LCS.	p.148
sdiagrgcs_beam	Return matrices to plot the stresses in a beam element.	p.253
sdiagrlcs_beam	Stress diagram for a beam element in LCS.	p.256

truss



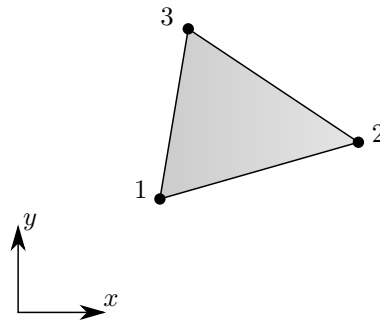
Description

Linear truss element.

Functions

<code>dof_truss</code>	Element degrees of freedom for a truss element.	p.134
<code>ke_truss</code>	Truss element stiffness and mass matrix in global coordinate system.	p.186
<code>kelcs_truss</code>	Truss element stiffness and mass matrix in local coordinate system.	p.167
<code>trans_truss</code>	Transform coordinate system for a truss element.	p.297
<code>loads_truss</code>	Equivalent nodal forces for a truss element in the GCS.	p.199
<code>accel_truss</code>	Compute the distributed loads for a truss due to an acceleration.	p.66
<code>forces_truss</code>	Compute the element forces for a truss element.	p.155
<code>forceslcs_truss</code>	Compute the element forces for a truss element in the LCS.	p.152
<code>coord_truss</code>	Coordinates of the truss elements for plotting.	p.96
<code>disp_truss</code>	Return matrices to compute the displacements of the deformed trusses.	p.115
<code>dispgcs2lcs_truss</code>	Transform the element displacements to the LCS for a truss.	p.98
<code>fdiagrgcs_truss</code>	Return matrices to plot the forces in a truss element.	p.146
<code>sdiagrgcs_truss</code>	Return matrices to plot the stresses in a truss element.	p.255

plane3



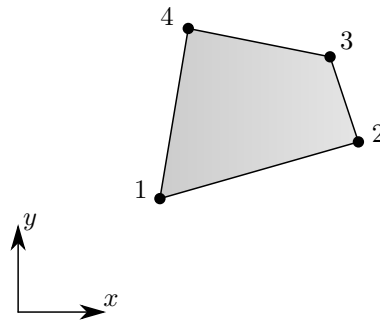
Description

The `plane3` element is a 3-node linear isoparametric 2D triangular plane element, commonly referred to as the “Constant Strain Triangle” (CST).

Functions

<code>dof_plane3</code>	Element degrees of freedom for a <code>plane3</code> element.	p.121
<code>ke_plane3</code>	Element stiffness and mass matrix in global coordinate system.	p.173
<code>coord_plane3</code>	Coordinates of the <code>plane3</code> element for plotting.	p.83
<code>disp_plane3</code>	Return matrices to compute the displacements of the deformed element.	p.103
<code>se_plane3</code>	Compute the element stresses for a <code>plane3</code> element in the GCS.	p.264
<code>selcs_plane3</code>	Compute the element stresses for a <code>plane3</code> element in the LCS.	p.257
<code>patch_plane3</code>	Patch information of the <code>plane3</code> elements for plotting.	p.213

plane4



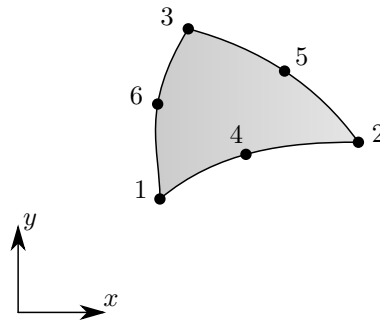
Description

The `plane4` element is a 4-node linear isoparametric 2D quadrilateral plane element.

Functions

<code>dof_plane4</code>	Element degrees of freedom for a <code>plane4</code> element.	p.122
<code>ke_plane4</code>	Element stiffness and mass matrix in global coordinate system.	p.174
<code>coord_plane4</code>	Coordinates of the <code>plane3</code> element for plotting.	p.84
<code>disp_plane4</code>	Return matrices to compute the displacements of the deformed element.	p.104
<code>se_plane4</code>	Compute the element stresses for a <code>plane4</code> element in the GCS.	p.265
<code>selcs_plane4</code>	Compute the element stresses for a <code>plane4</code> element in the LCS.	p.258
<code>patch_plane4</code>	Patch information of the <code>plane4</code> elements for plotting.	p.214

plane6



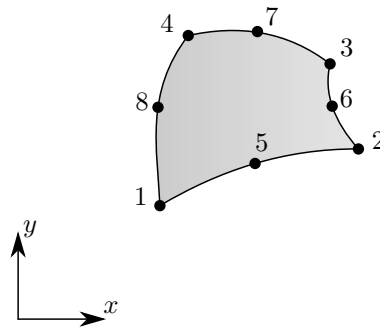
Description

The `plane6` element is a 6-node quadratic isoparametric 2D triangular plane element.

Functions

<code>dof_plane6</code>	Element degrees of freedom for a <code>plane6</code> element.	p.123
<code>ke_plane6</code>	Element stiffness and mass matrix in global coordinate system.	p.175
<code>coord_plane6</code>	Coordinates of the <code>plane3</code> element for plotting.	p.85
<code>disp_plane6</code>	Return matrices to compute the displacements of the deformed element.	p.105
<code>se_plane6</code>	Compute the element stresses for a <code>plane6</code> element in the GCS.	p.266
<code>selcs_plane6</code>	Compute the element stresses for a <code>plane6</code> element in the LCS.	p.259
<code>patch_plane6</code>	Patch information of the <code>plane6</code> elements for plotting.	p.215

plane8



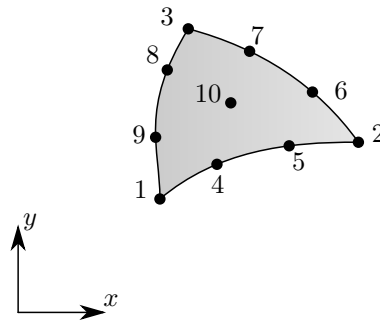
Description

The `plane8` element is a 8-node quadratic isoparametric 2D rectangular plane element.

Functions

<code>dof_plane8</code>	Element degrees of freedom for a <code>plane8</code> element.	p.124
<code>ke_plane8</code>	Element stiffness and mass matrix in global coordinate system.	p.176
<code>coord_plane8</code>	Coordinates of the <code>plane3</code> element for plotting.	p.86
<code>disp_plane8</code>	Return matrices to compute the displacements of the deformed element.	p.106
<code>se_plane8</code>	Compute the element stresses for a <code>plane8</code> element in the GCS.	p.267
<code>selcs_plane8</code>	Compute the element stresses for a <code>plane8</code> element in the LCS.	p.260
<code>patch_plane8</code>	Patch information of the <code>plane8</code> elements for plotting.	p.216

plane10



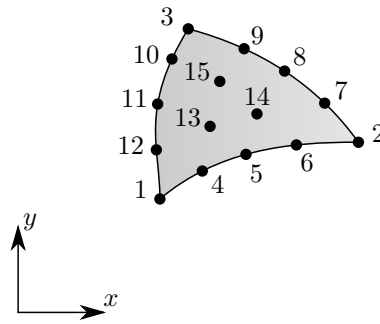
Description

The plane10 element is a 10-node cubic isoparametric 2D triangular plane element.

Functions

dof_plane10	Element degrees of freedom for a plane10 element.	p.119
ke_plane10	Element stiffness and mass matrix in global coordinate system.	p.171
coord_plane10	Coordinates of the plane3 element for plotting.	p.81
disp_plane10	Return matrices to compute the displacements of the deformed element.	p.101

plane15



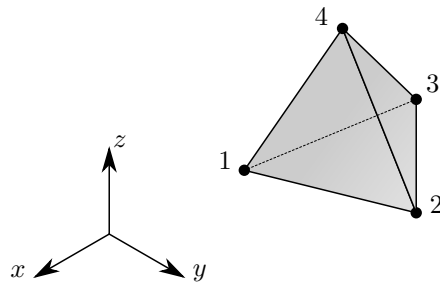
Description

The plane15 element is a 15-node quartic isoparametric 2D triangular plane element.

Functions

dof_plane15	Element degrees of freedom for a plane8 element.	p.120
ke_plane15	Element stiffness and mass matrix in global coordinate system.	p.172
coord_plane15	Coordinates of the plane3 element for plotting.	p.82
disp_plane15	Return matrices to compute the displacements of the deformed element.	p.102

solid4



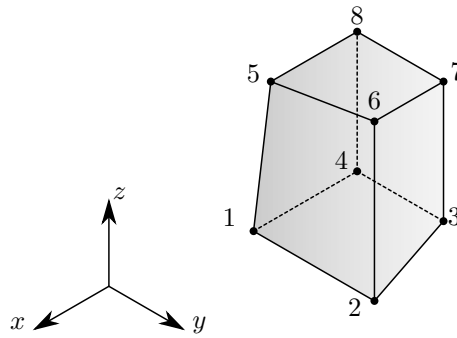
Description

The `solid4` element is a 4-node linear isoparametric 3D solid tetrahedral element.

Functions

<code>dof_solid4</code>	Element degrees of freedom for a <code>solid4</code> element.	p.132
<code>ke_solid4</code>	Element stiffness and mass matrix in global coordinate system.	p.184
<code>coord_solid4</code>	Coordinates of the <code>plane3</code> element for plotting.	p.94
<code>patch_solid4</code>	Patch information of the <code>solid4</code> elements for plotting.	p.222

solid8



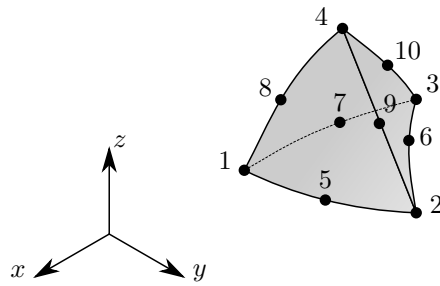
Description

The `solid8` element is a 8-node linear isoparametric 3D solid element.

Functions

<code>dof_solid8</code>	Element degrees of freedom for a solid4 element.	p.133
<code>ke_solid8</code>	Element stiffness and mass matrix in global coordinate system.	p.185
<code>coord_solid8</code>	Coordinates of the plane3 element for plotting.	p.95
<code>patch_solid8</code>	Patch information of the solid8 elements for plotting.	p.223

solid10



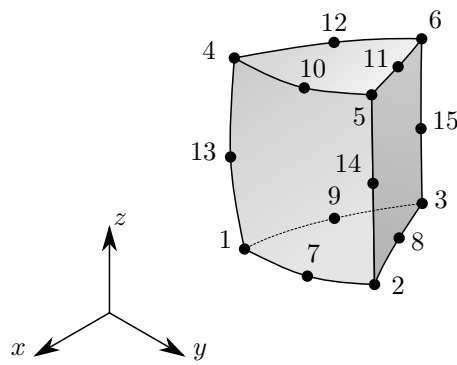
Description

The `solid10` element is a 10-node quadratic isoparametric 3D solid tetrahedral element.

Functions

<code>dof_solid10</code>	Element degrees of freedom for a <code>solid10</code> element.	p.129
<code>ke_solid10</code>	Element stiffness and mass matrix in global coordinate system.	p.181
<code>coord_solid10</code>	Coordinates of the <code>plane3</code> element for plotting.	p.91
<code>disp_solid10</code>	Return matrices to compute the displacements of the deformed element.	p.111
<code>patch_solid10</code>	Patch information of the <code>solid10</code> elements for plotting.	p.220

solid15

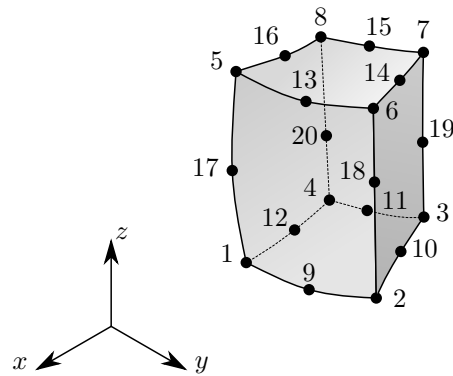


Description

The `solid15` element is a 15-node quadratic isoparametric 3D solid tetrahedral element.

<code>dof_solid15</code>	Element degrees of freedom for a <code>solid15</code> element.	p.130
<code>ke_solid15</code>	Element stiffness and mass matrix in global coordinate system.	p.182
<code>coord_solid15</code>	Coordinates of the <code>plane3</code> element for plotting.	p.92
<code>disp_solid15</code>	Return matrices to compute the displacements of the deformed element.	p.112

solid20



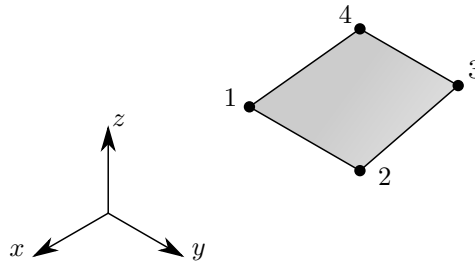
Description

The `solid20` element is a 20-node quadratic isoparametric 3D solid tetrahedral element.

Functions

<code>dof_solid20</code>	Element degrees of freedom for a <code>solid20</code> element.	p.131
<code>ke_solid20</code>	Element stiffness and mass matrix in global coordinate system.	p.183
<code>coord_solid20</code>	Coordinates of the <code>plane3</code> element for plotting.	p.93
<code>disp_solid20</code>	Return matrices to compute the displacements of the deformed element.	p.113
<code>patch_solid20</code>	Patch information of the <code>solid10</code> elements for plotting.	p.221

shell4



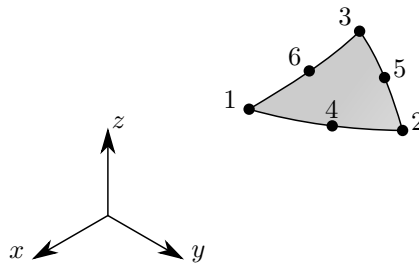
Description

Shell element consisting of a bilinear membrane element and four overlaid DKT triangles for the bending stiffness.

Functions

dof_shell4	Element degrees of freedom for a shell4 element.	p.126
ke_shell4	Shell4 element stiffness and mass matrix in global coordinate system.	p.178
kelcs_shell4	Shell4 element stiffness and mass matrix in local coordinate system.	p.166
trans_shell4	Transform coordinate system for a shell4 element.	p.293
ke_dkt	DKT plate element stiffness and mass matrix.	p.169
q_dkt	Q matrix for a DKT element.	p.247
sh_qs4	Shape functions for a quadrilateral serendipity element with 4 nodes.	p.274
sh_t	Shape functions for a triangular plate element.	p.276
se_shell4	Compute the element stresses for a shell4 element.	p.268
selcs_shell4	Compute the element stresses for a shell4 element.	p.261
accel_shell4	Compute the distributed loads for a shell due to an acceleration.	p.61
loads_shell4	Equivalent nodal forces for a shell4 element in the GCS.	p.194
loadslcs_shell4	Equivalent nodal forces for a shell4 element in the LCS.	p.191
pressure_shell4	Equivalent nodal forces for a shell4 element in the GCS due to a pressure.	p.239
coord_shell4	Coordinates of the shell4 elements for plotting.	p.88
disp_shell4	Matrices to compute the displacements of the deformed shell.	p.108
scontour_shell4	Matrix to plot contours in a shell4 element.	p.251
patch_shell4	Patch information of the shell4 elements for plotting.	p.217
grid_shell4	Grid in natural coordinates for mapped meshing.	p.162

shell6

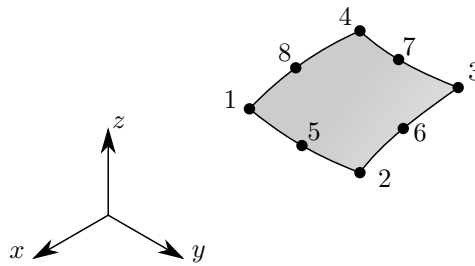


Description

This element is based on chapter 15 of Zienkiewicz [9].

dof_shell6	Element degrees of freedom for a shell4 element.	p.127
ke_shell6	Shell6 element stiffness and mass matrix in global coordinate system.	p.179
ke_dkt	DKT plate element stiffness and mass matrix.	p.169
q_dkt	Q matrix for a DKT element.	p.247
se_shell6	Compute the element stresses for a shell4 element.	p.269
accel_shell6	Compute the distributed loads for a shell due to an acceleration.	p.62
loads_shell6	Equivalent nodal forces for a shell6 element in the GCS.	p.195
pressure_shell6	Equivalent nodal forces for a shell6 element in the GCS due to a pressure.	p.240
coord_shell6	Coordinates of the shell6 elements for plotting.	p.89
disp_shell6	Matrices to compute the displacements of the deformed shell.	p.109
patch_shell6	Patch information of the shell6 elements for plotting.	p.218

shell8



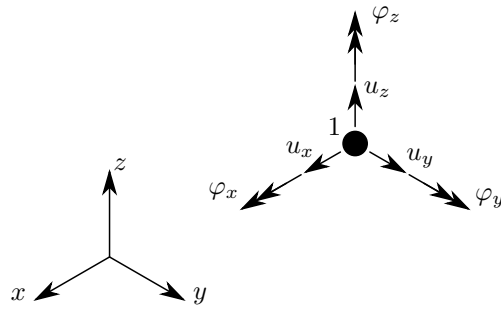
Description

This element is based on chapter 15 of Zienkiewicz [9].

Functions

<code>dof_shell8</code>	Element degrees of freedom for a shell8 element.	p.128
<code>ke_shell8</code>	shell8 element stiffness and mass matrix in global coordinate system.	p.180
<code>sh_qs8</code>	Shape functions for a quadrilateral serendipity element with 8 nodes.	p.275
<code>b_shell8</code>	B matrix for a shell8 element in global coordinate system.	p.73
<code>se_shell8</code>	Compute the element stresses for a shell8 element.	p.270
<code>accel_shell8</code>	Compute the distributed loads for a shell due to an acceleration.	p.63
<code>loads_shell8</code>	Equivalent nodal forces for a shell8 element in the GCS.	p.196
<code>pressure_shell8</code>	Equivalent nodal forces for a shell8 element in the GCS due to a pressure.	p.241
<code>coord_shell8</code>	Coordinates of the shell8 elements for plotting.	p.90
<code>disp_shell8</code>	Matrices to compute the displacements of the deformed shell.	p.110
<code>scontour_shell8</code>	Matrix to plot contours in a shell8 element.	p.252
<code>patch_shell8</code>	Patch information of the shell8 elements for plotting.	p.219
<code>grid_shell8</code>	Grid in natural coordinates for mapped meshing.	p.163

mass



Description

Single point concentrated mass element with 6 degrees of freedom.

Functions

<code>dof_mass</code>	Element degrees of freedom for a mass element.	p.118
<code>ke_mass</code>	Element stiffness and mass matrix in global coordinate system.	p.170
<code>coord_mass</code>	Coordinates of the mass element for plotting.	p.80
<code>disp_mass</code>	Return matrices to compute the displacements of the deformed element.	p.100
<code>patch_mass</code>	Patch information of the mass elements for plotting.	p.212

Chapter 5

Functions — By category

5.1 General functions

<code>getdof</code>	Get the vector with the degrees of freedom of the model.	p.159
<code>asmkm</code>	Assemble stiffness and mass matrix.	p.70
<code>removedof</code>	Remove DOF with Dirichlet boundary conditions equal to zero.	p.249
<code>addconstr</code>	Add constraint equations to the stiffness matrix and load vector.	p.67
<code>tconstr</code>	Return matrices to apply constraint equations.	p.288
<code>nodalvalues</code>	Construct a vector with the values at the selected DOF.	p.210
<code>elemloads</code>	Equivalent nodal forces.	p.138
<code>accel</code>	Compute the distributed loads due to an acceleration.	p.58
<code>elemforces</code>	Compute the element forces.	p.137
<code>elemdisp</code>	Select the element displacements from the global displacement vector.	p.136
<code>selectdof</code>	Select degrees of freedom.	p.262
<code>unselectdof</code>	Unselect degrees of freedom.	p.299
<code>selectnode</code>	Select nodes by location.	p.263
<code>repro</code>	Replicate rows from a matrix.	p.250
<code>multdloads</code>	Combine distributed loads.	p.204

5.2 Postprocessing

<code>plotnodes</code>	Plot the nodes.	p.232
<code>plotelem</code>	Plot the elements.	p.227
<code>plotdisp</code>	Plot the displacements.	p.225
<code>plotforc</code>	Plot the forces.	p.229
<code>plotlcs</code>	Plot the local element coordinate systems.	p.231
<code>plotstress</code>	Plot the stresses.	p.236
<code>plotstresscontour</code>	Plot stress contours.	p.237
<code>plotstresscontourf</code>	Plot filled contours of stresses.	p.238
<code>animdisp</code>	Animate the displacements.	p.68
<code>getmovie</code>	Get the movie from a figure where an animation has been played.	p.160
<code>printdisp</code>	Display the displacements in the command window.	p.243
<code>printforc</code>	Display the forces in the command window.	p.244

5.3 Dynamics

<code>eigfem</code>	Compute eigenmodes and eigenfrequencies.	p.135
<code>msupt</code>	Modal superposition in the time domain.	p.203
<code>msupf</code>	Modal superposition in the frequency domain.	p.202
<code>cdiff</code>	Direct time integration for dynamic systems - central diff. method.	p.74
<code>newmark</code>	Direct time integration for dynamic systems - Newmark method	p.207
<code>wilson</code>	Direct time integration for dynamic systems - Wilson-theta method	p.302

5.4 General shell functions

<code>elempressure</code>	Equivalent nodal forces for pressure \perp shell surface.	p.139
<code>gaussq</code>	Gauss points for 2D numerical integration.	p.156
<code>nodalshellf</code>	Compute the nodal shell forces/moments from element solution.	p.208
<code>nodalstress</code>	Compute the nodal stresses from element solution.	p.209
<code>plotprincstress</code>	Plot the principal stresses.	p.233
<code>plotshellfcontour</code>	Plot contour lines of forces/moments per unit length.	p.234
<code>plotstresscontour</code>	Plot stress contour lines.	p.237
<code>plotshellfcontourf</code>	Plot filled contours of shell forces.	p.235
<code>principalstress</code>	Compute the principal stresses and directions.	p.242
<code>printshellf</code>	Display forces/moments in command window (shell elements).	p.245
<code>printstress</code>	Display stresses in command window (shell elements).	p.246

Chapter 6

Functions — Alphabetical list

accel

ACCEL Compute the distributed loads due to an acceleration.

```
DLoads=accel(Accelxyz,Nodes,Elements,Types,Sections,Materials)
```

computes the distributed loads due to an acceleration.

In order to simulate gravity, accelerate the structure in the direction opposite to gravity.

Accelxyz	Acceleration	[Ax Ay Az] (1 * 3)
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Types	Element type definitions	{TypID EltName Option1 ... }
Sections	Section definitions	[SecID SecProp1 SecProp2 ...]
Materials	Material definitions	[MatID MatProp1 MatProp2 ...]
DLoads	Distributed loads	[EltID n1globalX n1globalY n1globalZ ...]

See also ELEMLOADS, ACCEL_BEAM, ACCEL_TRUSS.

accel_beam

ACCEL_BEAM Compute the distributed loads for a beam due to an acceleration.

DLoads=accel_beam(Accelxyz,[],Elements,Sections,Materials,Options)
computes the distributed loads for a beam due to an acceleration.
In order to simulate gravity, accelerate the structure in the direction
opposite to gravity.

Accelxyz	Acceleration	[Ax Ay Az] (1 * 3)
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Sections	Section definitions	[SecID SecProp1 SecProp2 ...]
Materials	Material definitions	[MatID MatProp1 MatProp2 ...]
Options	Element options	{Option1 Option2 ...}
DLoads	Distributed loads	[EltID n1globalX n1globalY n1globalZ ...]

See also ACCEL, ACCEL_TRUSS.

accel_shell2

ACCEL_SHELL2 Compute the distributed loads for a SHELL2 element due to an acceleration.

`DLoads=accel_shell2(Accelxyz,Nodes,Elements,Sections,Materials,Options)`
 computes the distributed loads for a SHELL2 element due to an acceleration.
 In order to simulate gravity, accelerate the structure in the direction
 opposite to gravity.

<code>Accelxyz</code>	Acceleration	<code>[Ax Ay Az] (1 * 3)</code>
<code>Nodes</code>	Node definitions	<code>[NodeID x y z]</code>
<code>Elements</code>	Element definitions	<code>[EltID TypID SecID MatID n1 n2 ...]</code>
<code>Sections</code>	Section definitions	<code>[SecID SecProp1 SecProp2 ...]</code>
<code>Materials</code>	Material definitions	<code>[MatID MatProp1 MatProp2 ...]</code>
<code>Options</code>	Element options	<code>{Option1 Option2 ...}</code>
<code>DLoads</code>	Distributed loads	<code>[EltID n1globalX n1globalY n1globalZ ...]</code>

accel_shell4

ACCEL_SHELL4 Compute the distributed loads for shell4 elements due to an acceleration.

`DLoads = accel_shell4(Accelxyz,[],Elements,Sections,Materials,Options)`
computes the distributed loads for shell4 elements due to an acceleration.
In order to simulate gravity, accelerate the structure in the direction opposite to gravity.

<code>Accelxyz</code>	Acceleration	<code>[Ax Ay Az] (1 * 3)</code>
<code>Elements</code>	Element definitions	<code>[EltID TypID SecID MatID n1 n2 ...]</code>
<code>Sections</code>	Section definitions	<code>[SecID SecProp1 SecProp2 ...]</code>
<code>Materials</code>	Material definitions	<code>[MatID MatProp1 MatProp2 ...]</code>
<code>Options</code>	Element options	<code>{Option1 Option2 ...}</code>
<code>DLoads</code>	Distributed loads	<code>[EltID n1globalX n1globalY n1globalZ ...]</code>

See also `ACCEL`, `ACCEL_TRUSS`.

accel_shell6

ACCEL_SHELL6 Compute the distributed loads for shell6 elements due to an acceleration.

```
DLoads = accel_shell6(Accelxyz,[],Elements,Sections,Materials,Options)
computes the distributed loads for shell8 elements due to an acceleration.
In order to simulate gravity, accelerate the structure in the direction
opposite to gravity.
```

Accelxyz	Acceleration	[Ax Ay Az] (1 * 3)
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Sections	Section definitions	[SecID SecProp1 SecProp2 ...]
Materials	Material definitions	[MatID MatProp1 MatProp2 ...]
Options	Element options struct. Fields:	
	-MatType: 'isotropic' (default) or 'orthotropic'	
DLoads	Distributed loads	[EltID n1globalX n1globalY n1globalZ ...]

See also ACCEL, ACCEL_TRUSS.

accel_shell8

ACCEL_SHELL8 Compute the distributed loads for shell8 elements due to an acceleration.

```
DLoads = accel_shell8(Accelxyz,[],Elements,Sections,Materials,Options)
computes the distributed loads for shell8 elements due to an acceleration.
In order to simulate gravity, accelerate the structure in the direction
opposite to gravity.
```

Accelxyz	Acceleration	[Ax Ay Az] (1 * 3)
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Sections	Section definitions	[SecID SecProp1 SecProp2 ...]
Materials	Material definitions	[MatID MatProp1 MatProp2 ...]
Options	Element options struct. Fields:	
	-MatType: 'isotropic' (default) or 'orthotropic'	
DLoads	Distributed loads	[EltID n1globalX n1globalY n1globalZ ...]

See also ACCEL, ACCEL_TRUSS.

accel_solid20

ACCEL_SOLID20 Compute the distributed loads for solid20 elements due to an acceleration.

`DLoads = accel_solid20(Accelxyz,[],Elements,Sections,Materials,Options)`
 computes the distributed loads for solid20 elements due to an acceleration.
 In order to simulate gravity, accelerate the structure in the direction
 opposite to gravity.

<code>Accelxyz</code>	Acceleration	<code>[Ax Ay Az] (1 * 3)</code>
<code>Elements</code>	Element definitions	<code>[EltID TypID SecID MatID n1 n2 ...]</code>
<code>Sections</code>	Section definitions	<code>[SecID SecProp1 SecProp2 ...]</code>
<code>Materials</code>	Material definitions	<code>[MatID MatProp1 MatProp2 ...]</code>
<code>Options</code>	Element options	<code>{Option1 Option2 ...}</code>
<code>DLoads</code>	Distributed loads	<code>[EltID n1globalX n1globalY n1globalZ ...]</code>

See also `ACCEL`, `ACCEL_TRUSS`.

accel_solid8

ACCEL_SOLID8 Compute the distributed loads for solid8 elements due to an acceleration.

`DLoads = accel_solid8(Accelxyz,[],Elements,Sections,Materials,Options)`
computes the distributed loads for solid8 elements due to an acceleration.
In order to simulate gravity, accelerate the structure in the direction opposite to gravity.

<code>Accelxyz</code>	Acceleration	<code>[Ax Ay Az] (1 * 3)</code>
<code>Elements</code>	Element definitions	<code>[EltID TypID SecID MatID n1 n2 ...]</code>
<code>Sections</code>	Section definitions	<code>[SecID SecProp1 SecProp2 ...]</code>
<code>Materials</code>	Material definitions	<code>[MatID MatProp1 MatProp2 ...]</code>
<code>Options</code>	Element options	<code>{Option1 Option2 ...}</code>
<code>DLoads</code>	Distributed loads	<code>[EltID n1globalX n1globalY n1globalZ ...]</code>

See also `ACCEL`, `ACCEL_TRUSS`.

accel_truss

ACCEL_TRUSS Compute the distributed loads for a truss due to an acceleration.

`DLoads=accel_truss(Accelxyz,[],Elements,Sections,Materials,Options)`
 computes the distributed loads for a truss due to an acceleration.
 In order to simulate gravity, accelerate the structure in the direction
 opposite to gravity.

<code>Accelxyz</code>	Acceleration	<code>[Ax Ay Az] (1 * 3)</code>
<code>Elements</code>	Element definitions	<code>[EltID TypID SecID MatID n1 n2 ...]</code>
<code>Sections</code>	Section definitions	<code>[SecID SecProp1 SecProp2 ...]</code>
<code>Materials</code>	Material definitions	<code>[MatID MatProp1 MatProp2 ...]</code>
<code>Options</code>	Element options	<code>{Option1 Option2 ...}</code>
<code>DLoads</code>	Distributed loads	<code>[EltID n1globalX n1globalY n1globalZ ...]</code>

See also `ACCEL`, `ACCEL_BEAM`.

addconstr

ADDCONSTR Add constraint equations to the stiffness matrix and load vector.

```
[K,F]=addconstr(Constr,DOF,K,F)
[K,F,M]=addconstr(Constr,DOF,K,[],M)
[K,F,M]=addconstr(Constr,DOF,K,F,M)
```

modifies the stiffness matrix, the mass matrix and the load vector according to the applied constraint equations. The dimensions of the stiffness matrix, the mass matrix and the load vector are kept the same. The resulting stiffness and mass matrix are not symmetric anymore. This function can be used as well to apply imposed displacements.

Constr	Constraint equation: Constant=CoefS*SlaveDOF+CoefM1*MasterDOF1+CoefM2*MasterDOF2+... [Constant CoefS SlaveDOF CoefM1 MasterDOF1 CoefM2 MasterDOF2 ...]
DOF	Degrees of freedom (nDOF * 1)
K	Stiffness matrix (nDOF * nDOF)
F	Load vector (nDOF * nSteps)
M	Mass matrix (nDOF * nDOF)

animdisp

ANIMDISP Animate the displacements.

```
DispScal=animdisp(Nodes,Elements,Types,DOF,U)
animates the displacements.
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Types	Element type definitions	{TypID EltName Option1 ... }
DOF	Degrees of freedom	(nDOF * 1)
U	Displacements	(nDOF * nSteps)
DispScal	Displacement scaling	

ANIMDISP(...,ParamName,ParamValue) sets the value of the specified parameters. The following parameters can be specified:

'DispScal'	Displacement scaling. Default: 'auto'.
'Handle'	Plots in the axis with this handle. Default: current axis.
'Fps'	Frames per second. Default: 12.
'CreateMovie'	Saves the movie in the userdata of the axis of the figure. Use getmovie to get the movie from the axis. Default: 'off'.
'Counter'	Displays the number of the frame for transient displacements. Default: 'on'.

Additional parameters are redirected to the PLOTDISP function which plots the individual frames of the movie.

See also GETMOVIE, PLOTDISP.

argdimchk

ARGDIMCHK Validate input argument dimensions.

`msg = ARGDIMCHK(arg1,size1,arg2,size2,...)` returns an appropriate error message if the dimensions of `arg1,arg2,...` do not comply with `size1,size2,...` respectively. If they do comply, an empty matrix is returned. `size1,size2,...` are cell arrays consisting of at least 2 cells. Each cell corresponds to a dimension of `arg1,arg2,...` and contains a number (to constrain the dimension explicitly) or a string (to constrain the dimension implicitly). E.g. the expression:

```
ERROR(ARGDIMCHK( ...
    omega,{ 'nMode',    1      }, ...
    Phi,  { 'nDof',    'nMode' }, ...
    xi,   { 'nMode',    1      }, ...
    b,    { 'nDof',    1      }, ...
    q,    { 1,          'nOmega'}, ...
    Omega,{ 1,          'nOmega'}, ...
    c,    { 'nSelDof', 'nDof'  }));
```

checks if `omega,Phi,xi,b,q,Omega,c` are all 2-dimensional variables and if

```
SIZE(omega,2)==1      SIZE(Phi,2)==SIZE(omega,1)
SIZE(xi,1)==SIZE(omega,1)  SIZE(xi,2)==1
SIZE(b,1)==SIZE(Phi,1)    SIZE(b,2)==1
SIZE(q,1)==1            SIZE(Omega,1)==1
SIZE(Omega,2)==SIZE(q,2)  SIZE(c,2)==SIZE(Phi,1)
```

If not, an appropriate error message is shown.

asmkm

ASMKM Assemble stiffness and mass matrix.

```
[K,M] = ASMKM(Nodes,Elements,Types,Sections,Materials,DOF)
K      = ASMKM(Nodes,Elements,Types,Sections,Materials,DOF)
K      = ASMKM(Nodes,Elements,Types,Sections,Materials)
```

assembles the stiffness and the mass matrix using the finite element method.

```
[K,~,dKdx] = ASMKM(Nodes,Elements,Types,Sections,Materials,DOF,dNodesdx,dSectionsdx)
assembles the stiffness matrix using the finite element method and
additionally computes the derivatives of the stiffness matrix with
respect to the design variables x. The derivatives of the mass matrix
have not yet been implemented.
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Types	Element type definitions	{TypID EltName Option1 ...}
Sections	Section definitions	[SecID SecProp1 SecProp2 ...]
Materials	Material definitions	[MatID MatProp1 MatProp2 ...]
DOF	Degrees of freedom	(nDOF * 1)
dNodesdx	Node definitions derivatives	(SIZE(Nodes) * nVar)
dSectionsdx	Section definitions derivatives	(SIZE(Sections) * nVar)
K	Stiffness matrix	(nDOF * nDOF)
M	Mass matrix	(nDOF * nDOF)
dKdx	Stiffness matrix derivatives	(CELL(nVar,1))

See also KE_TRUSS, KE_BEAM.

be_plane3

e_plane3 is a function.

```
[BeGCS] = be_plane3(Node, Section, Material, UeGCS, Options, gcs)
```

b_shell6

B_SHELL6 b matrix for a shell6 element in global coordinate system.

[Bg,J] = b_shell6(Ni,dN_dxi,dN_deta,zetar,Node,h,v1i,v2i,v3i) returns the element b matrix in the global coordinate system and the Jacobian of the parametric transformation. Both are evaluated in the natural coordinates (xi,eta and zetar) which were used to calculate Ni,dN_dxi,dN_deta and zetar.

Node	Node definitions	[x y z] (6 * 3)
	Nodes should have the following order:	
	3	
	\	
	6 5	
	\	
	1--4--2	
Ni	Shape functions for quadratic serendipity element	(6 * 1)
dN_dxi	first derivatives of shape functions Ni	(6 * 1)
dN_deta	first derivatives of shape functions Ni	(6 * 1)
h	scalar or vector containing thickness	scalar or (6 * 1)
v(1,2,3)i	components of the local coordinate system in node i	(6 * 3)
d	Nodal offset from shell mid plane	scalar (default = 0)
Bg	b matrix of shell8 element	(6 * 36)
J	Jacobian of the parametric transformation	(3 * 3)

See also SE_SHELL8, KE_SHELL8.

b_shell8

B_SHELL8 b matrix for a shell8 element in global coordinate system.

[Bg,J] = b_shell8(Ni,dN_dxi,dN_deta,zetar,Node,h,v1i,v2i,v3i) returns the element b matrix in the global coordinate system and the Jacobian of the parametric transformation. Both are evaluated in the natural coordinates (xi,eta and zetar) which were used to calculate Ni,dN_dxi,dN_deta and zetar.

Node	Node definitions	[x y z] (8 * 3)
	Nodes should have the following order: 4-----7-----3 8 6 1-----5-----2	
Ni	Shape functions for quadratic serendipity element	(8 * 1)
dN_dxi	first derivatives of shape functions Ni	(8 * 1)
dN_deta	first derivatives of shape functions Ni	(8 * 1)
h	scalar or vector containing thickness	scalar or (8 * 1)
v(1,2,3)i	components of the local coordinate system in node i	(8 * 3)
d	Nodal offset from shell mid plane	scalar (default = 0)
Bg	b matrix of shell8 element	(6 * 48)
J	Jacobian of the parametric transformation	(3 * 3)

See also SE_SHELL8, KE_SHELL8.

cdiff

CDIFF Direct time integration for dynamic systems - central diff. method.
[u,t] = CDIFF(M,C,K,dt,p,u0,u1) applies the central difference method for the calculation of the nodal displacements u of the dynamic system with the system matrices M, C and K due to the excitation p.

M Mass matrix (nDof * nDof)
C Damping matrix (nDof * nDof)
K Stiffness matrix (nDof * nDof)
dt Time step of the integration scheme (1 * 1). Should be small enough to ensure the stability and the precision of the integration scheme.
p Excitation (nDof * N). p(:,k) corresponds to time point t(k).
u0 Displacements at time point t(1)-dt (nDof * 1). Defaults to zero.
u1 Displacements at time point t(1) (nDof * 1). Defaults to zero.
u Displacements (nDof * N). u(:,k) corresponds to time point t(k).
t Time axis (1 * N), defined as t = [0:N-1] * dt.

checkunique

CHECKUNIQUE Check if vector contains unique elements.

```
checkunique(p,name)  
find non-unique elements in a vector and print error
```

```
p      Vector with elements to be checked  
name   Name of elements in p
```

cmat_isotropic

CMAT_ISOTROPIC Constitutive matrix for isotropic materials.

```
[C,C_lambda,C_mu]==cmat_isotropic(problem,Section,Material)
```

computes the constitutive matrix for isotropic materials.

problem

Section Section definition

Material Material definition

C Constitutive matrix (nStress * nStrain)

C_lambda Contribution of lambda to C (nStress * nStrain)

C_mu Contribution of mu to C (nStress * nStrain)

cmat_shell8

CMAT_SHELL8 Constitutive matrix for shell8 element

```
C = cmat_shell8(MatType,Material,k)
C = cmat_shell8(MatType,Material)
returns the constitutive matrix for shell8 elements.
```

```
MatType    Material type: 'isotropic' or 'orthotropic'
Material    Material definition
            isotropic: [E nu rho]
            orthotropic: [Exx Eyy nuxy muxy muyz muzx theta rho]
k           Geometric coefficient for non-uniform shear stress (default = 1.2)
C           Constitutive matrix (5 * 5)
```

See also KE_SHELL8

coord_beam

COORD_BEAM Coordinates of the beam elements for plotting.

```
[X,Y,Z]=coord_beam(Nodes,NodeNum)
```

returns the coordinates of the beam elements for plotting.

Nodes	Node definitions	[NodID x y z] (nNodes * 4)
NodeNum	Node numbers	[NodID1 NodID2 NodID3] (nElem * 3)
X	X coordinates	(2 * nElem)
Y	Y coordinates	(2 * nElem)
Z	Z coordinates	(2 * nElem)

See also COORD_TRUSS, PLOTELEM.

coord_kbeam

COORD_BEAM Coordinates of the beam elements for plotting.

```
coord_beam(Nodes,NodeNum)
```

returns the coordinates of the beam elements for plotting.

Nodes	Node definitions	[NodID x y z] (nNodes * 4)
NodeNum	Node numbers	[NodID1 NodID2 NodID3] (nElem * 3)
X	X coordinates	(2 * nElem)
Y	Y coordinates	(2 * nElem)
Z	Z coordinates	(2 * nElem)

See also COORD_TRUSS, PLOTELEM.

coord_mass

COORD_MASS Coordinates of the mass element for plotting.

```
[X,Y,Z]=coord_mass(Nodes,NodeNum)
```

returns the coordinates of the mass element for plotting.

Nodes	Node definitions	[NodID x y z] (nNode * 4)
NodeNum	Node numbers	[NodID1] (nElem * 1)
X	X coordinates	(1 * nElem)
Y	Y coordinates	(1 * nElem)
Z	Z coordinates	(1 * nElem)

See also COORD_BEAM, PLOTELEM.

coord_plane10

COORD_PLANE10 Coordinates of the plane elements for plotting.

```
[X,Y,Z] = coord_plane10(Nodes,NodeNum)
```

returns the coordinates of the plane10 elements for plotting.

Nodes	Node definitions	[NodID x y z] (nNodes * 4)
NodeNum	Node numbers	[NodID1 NodID2 ...] (nElem * 10)
X	X coordinates	(15 * nElem)
Y	Y coordinates	(15 * nElem)
Z	Z coordinates	(15 * nElem)

See also COORD_TRUSS, PLOTELEM.

coord_plane15

COORD_PLANE15 Coordinates of the plane elements for plotting.

```
[X,Y,Z] = coord_plane15(Nodes,NodeNum)
```

returns the coordinates of the plane15 elements for plotting.

Nodes	Node definitions	[NodID x y z] (nNodes * 4)
NodeNum	Node numbers	[NodID1 NodID2 ...] (nElem * 15)
X	X coordinates	(15 * nElem)
Y	Y coordinates	(15 * nElem)
Z	Z coordinates	(15 * nElem)

See also COORD_TRUSS, PLOTELEM.

coord_plane3

COORD_PLANE3 Coordinates of PLANE3 element sides for plotting.

```
[X,Y,Z]=coord_plane3(Nodes,Nodenumbers)
```

returns the coordinates of PLANE3 element sides for plotting.

Nodes	Node definitions [NodID x y z] (nNodes * 3)
Nodenumbers	Node numbers [NodID1 NodID2 NodID3] (nElem * 3)
X	X coordinates (3 * nElem)
Y	Y coordinates (3 * nElem)
Z	Z coordinates (3 * nElem)

coord_plane4

COORD_PLANE4 Coordinates of the plane elements for plotting.

```
[X,Y,Z] = coord_plane4(Nodes,NodeNum)
```

returns the coordinates of the plane4 elements for plotting.

Nodes	Node definitions	[NodID x y z] (nNodes * 4)
NodeNum	Node numbers	[NodID1 NodID2 NodID3 NodID4] (nElem * 4)
X	X coordinates	(4 * nElem)
Y	Y coordinates	(4 * nElem)
Z	Z coordinates	(4 * nElem)

See also COORD_TRUSS, PLOTELEM.

coord_plane6

COORD_PLANE6 Coordinates of the plane elements for plotting.

```
[X,Y,Z] = coord_plane6(Nodes,NodeNum)
```

returns the coordinates of the plane6 elements for plotting.

Nodes	Node definitions	[NodID x y z] (nNodes * 4)
NodeNum	Node numbers	[NodID1 NodID2 ...] (nElem * 6)
X	X coordinates	(6 * nElem)
Y	Y coordinates	(6 * nElem)
Z	Z coordinates	(6 * nElem)

See also COORD_TRUSS, PLOTELEM.

coord_plane8

COORD_PLANE8 Coordinates of the shell8 elements for plotting.

```
coord_plane8(Nodes,NodeNum)
```

returns the coordinates of the plane8 elements for plotting.

Nodes	Node definitions	[NodID x y z] (nNodes * 4)
NodeNum	Node numbers	[NodID1 NodID2 NodID3 NodID4] (nElem * 8)
X	X coordinates	(20 * nElem)
Y	Y coordinates	(20 * nElem)
Z	Z coordinates	(20 * nElem)

See also COORD_TRUSS, PLOTELEM.

coord_shell2

COORD_SHELL2 Coordinates of SHELL2 elements for plotting.

[X,Y,Z]=coord_shell2(Nodes,NodeNum)
returns the coordinates of the SHELL2 elements for plotting.

Nodes	Node definitions	[NodID x y z] (nNodes * 4)
NodeNum	Node numbers	[NodID1 NodID2] (nElem * 2)
X	X coordinates	(2 * nElem)
Y	Y coordinates	(2 * nElem)
Z	Z coordinates	(2 * nElem)

coord_shell4

COORD_SHELL4 Coordinates of the shell elements for plotting.

```
[X,Y,Z] = coord_shell4(Nodes,NodeNum)
```

returns the coordinates of the shell4 elements for plotting.

Nodes	Node definitions	[NodID x y z] (nNodes * 4)
NodeNum	Node numbers	[NodID1 NodID2 NodID3 NodID4] (nElem * 4)
X	X coordinates	(4 * nElem)
Y	Y coordinates	(4 * nElem)
Z	Z coordinates	(4 * nElem)

See also COORD_TRUSS, PLOTELEM.

coord_shell6

COORD_SHELL6 Coordinates of the shell6 elements for plotting.

```
[X,Y,Z] = coord_shell6(Nodes,NodeNum)
```

returns the coordinates of the shell6 elements for plotting.

Nodes	Node definitions	[NodID x y z] (nNodes * 4)
NodeNum	Node numbers	[NodID1 NodID2 NodID3 NodID4] (nElem * 6)
X	X coordinates	(15 * nElem)
Y	Y coordinates	(15 * nElem)
Z	Z coordinates	(15 * nElem)

See also COORD_TRUSS, PLOTELEM.

coord_shell8

COORD_SHELL8 Coordinates of the shell8 elements for plotting.

```
[X,Y,Z] = coord_shell8(Nodes,NodeNum)
```

returns the coordinates of the shell8 elements for plotting.

Nodes	Node definitions	[NodID x y z] (nNodes * 4)
NodeNum	Node numbers	[NodID1 NodID2 NodID3 NodID4] (nElem * 8)
X	X coordinates	(20 * nElem)
Y	Y coordinates	(20 * nElem)
Z	Z coordinates	(20 * nElem)

See also COORD_TRUSS, PLOTELEM.

coord_solid10

COORD_SOLID10 Coordinates of SOLID10 element sides for plotting.

```
[X,Y,Z]=coord_solid10(Nodes,Nodenumbers)
```

returns the coordinates of SOLID10 element sides for plotting.

Nodes Node definitions [NodID x y z] (nNodes * 4)

Nodenumbers Node numbers [NodID1 NodID2 NodID3 NodID4] (nElem * 4)

X X coordinates (4 * nElem)

Y Y coordinates (4 * nElem)

Z Z coordinates (4 * nElem)

coord_solid15

COORD_SOLID8 Coordinates of SOLID15 element sides for plotting.

```
[X,Y,Z]=coord_solid15(Nodes,Nodenumbers)
```

returns the coordinates of SOLID15 element sides for plotting.

Nodes Node definitions [NodID x y z] (nNodes * 4)

Nodenumbers Node numbers [NodID1 NodID2 NodID3 NodID4] (nElem * 4)

X X coordinates (4 * nElem)

Y Y coordinates (4 * nElem)

Z Z coordinates (4 * nElem)

coord_solid20

COORD_SOLID20 Coordinates of SOLID20 element sides for plotting.

```
[X,Y,Z]=coord_solid20(Nodes,Nodenumbers)
```

returns the coordinates of SOLID20 element sides for plotting.

Nodes Node definitions [NodID x y z] (nNodes * 4)

Nodenumbers Node numbers [NodID1 NodID2 NodID3 NodID4] (nElem * 4)

X X coordinates (4 * nElem)

Y Y coordinates (4 * nElem)

Z Z coordinates (4 * nElem)

coord_solid4

COORD_SOLID4 Coordinates of SOLID4 element sides for plotting.

```
[X,Y,Z]=coord_rshell(Nodes,Nodenumbers)
```

returns the coordinates of RSHELL element sides for plotting.

Nodes Node definitions [NodID x y z] (nNodes * 4)

Nodenumbers Node numbers [NodID1 NodID2 NodID3 NodID4] (nElem * 4)

X X coordinates (4 * nElem)

Y Y coordinates (4 * nElem)

Z Z coordinates (4 * nElem)

coord_solid8

COORD_SOLID8 Coordinates of SOLID8 element sides for plotting.

```
[X,Y,Z]=coord_solid8(Nodes,Nodenumbers)
```

returns the coordinates of the SOLID8 element sides for plotting.

Nodes Node definitions [NodID x y z] (nNodes * 4)

Nodenumbers Node numbers [NodID1 NodID2 NodID3 NodID4] (nElem * 4)

X X coordinates (4 * nElem)

Y Y coordinates (4 * nElem)

Z Z coordinates (4 * nElem)

coord_truss

COORD_TRUSS Coordinates of the truss elements for plotting.

```
[X,Y,Z]=coord_truss(Nodes,NodeNum)
```

returns the coordinates of the truss elements for plotting.

Nodes	Node definitions	[NodID x y z] (nNodes * 4)
NodeNum	Node numbers	[NodID1 NodID2] (nElem * 2)
X	X coordinates	(2 * nElem)
Y	Y coordinates	(2 * nElem)
Z	Z coordinates	(2 * nElem)

See also COORD_BEAM, PLOTELEM.

dispgcs2lcs_beam

DISPGCS2LCS_BEAM Transform the element displacements to the LCS for a beam.

```
UeLCS=dispgcs2lcs_beam(UeGCS,Node)
```

transforms the element displacements from the GCS to the LCS for a beam element.

Node	Node definitions	[x y z] (3 * 3)
------	------------------	-----------------

UeGCS	Displacements in the GCS (12 * 1)
-------	-----------------------------------

UeLCS	Displacements in the LCS (12 * 1)
-------	-----------------------------------

See also DISPGCS2LCS_TRUSS.

dispgcs2lcs_truss

DISPGCS2LCS_TRUSS Transform the element displacements to the LCS for a truss.

```
UeLCS=dispgcs2lcs_truss(UeGCS,Node)
transforms the element displacements from the GCS to the LCS for a truss
element.
```

Node	Node definitions	[x y z] (2 * 3)
UeGCS	Displacements in the GCS	(6 * 1)
UeLCS	Displacements in the LCS	(6 * 1)

See also DISPGCS2LCS_BEAM.

disp_beam

DISP_BEAM Return matrices to compute the displacements of the deformed beams.

```
[Ax,Ay,Az,B,Cx,Cy,Cz] = DISP_BEAM(Nodes,Elements,DOF,DLoads,Sections,Materials,Points)
[Ax,Ay,Az,B,Cx,Cy,Cz] = DISP_BEAM(Nodes,Elements,DOF,DLoads,Sections,Materials)
[Ax,Ay,Az,B] = DISP_BEAM(Nodes,Elements,DOF,[],Sections,Materials)
[Ax,Ay,Az,B] = DISP_BEAM(Nodes,Elements,DOF)
    returns the matrices to compute the displacements of the deformed
    beams. The coordinates of the specified points along the deformed
    beam elements are computed using  $X=Ax*U+Cx*DLoad+B(:,1)$ ;
     $Y=Ay*U+Cy*DLoad+B(:,2)$  and  $Z=Az*U+Cz*DLoad+B(:,3)$ . The matrices
    Cx,Cy and Cz superimpose the displacements that occur due to the
    distributed loads if all nodes are fixed.
```

```
[Ax,Ay,Az,B,Cx,Cy,Cz,dAxdx,dAydx,dAzdxd,Cxdx,dCydx,dCzdx]
    = DISP_BEAM(Nodes,Elements,DOF,DLoads,Sections,Materials,Points,dNodesdx,
    dDLoadsdx,dSectionsdx)
    additionally computes the derivatives of the displacements with
    respect to the design variables x.
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
DOF	Degrees of freedom	(nDOF * 1)
DLoads	Distributed loads	[EltID n1globalX n1globalY n1globalZ ...]
		(use an empty array [] when shear deformation is considered but no DLoads are present)
Sections	Section definitions	[SecID SecProp1 SecProp2 ...]
Materials	Material definitions	[MatID MatProp1 MatProp2 ...]
Points	Points in the local coordinate system	(1 * nPoints)
dNodesdx	Node definitions derivatives	(SIZE(Node) * nVar)
dDLoadsdx	Distributed loads derivatives	(SIZE(DLoad) * nVar)
Ax	Matrix to compute the x-coordinates of the deformations	
Ay	Matrix to compute the y-coordinates of the deformations	
Az	Matrix to compute the z-coordinates of the deformations	
B	Matrix which contains the x-, y- and z-coordinates of the undeformed structure	
Cx	Matrix to compute the x-coordinates of the deformations	
Cy	Matrix to compute the y-coordinates of the deformations	
Cz	Matrix to compute the z-coordinates of the deformations	
dAxdx, dAydx, dAzdxd, dCxdx, dCydx, dCzdx	Derivatives of the matrices to compute the coordinates of the interpolation points after deformation	

See also DISP_TRUSS, PLOTDISP, NELCS_BEAM, NEDLOADLCS_BEAM.

disp_mass

DISP_MASS Matrices to compute the displacements of the deformed mass element

```
[Ax,Ay,Az,B,Cx,Cy,Cz]
    =disp_beam(Nodes,Elements,DOF,EltIDDLoad,Sections,Materials,Points)
[Ax,Ay,Az,B,Cx,Cy,Cz]
    =disp_beam(Nodes,Elements,DOF,EltIDDLoad,Sections,Materials)
[Ax,Ay,Az,B]
    =disp_beam(Nodes,Elements,DOF,[],Sections,Materials)
[Ax,Ay,Az,B]
    =disp_beam(Nodes,Elements,DOF)
```

returns the matrices to compute the displacements of the deformed mass.
 The coordinates of the node of the mass element are
 computed using $X=Ax*U+B(:,1)$; $Y=Ay*U+B(:,2)$ and
 $Z=Az*U+B(:,3)$. The matrices Cx,Cy and Cz superimpose the
 displacements that occur due to the distributed loads if all nodes are fixed.

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
DOF	Degrees of freedom	(nDOF * 1)
Ax	Matrix to compute the x-coordinates of the deformations	
Ay	Matrix to compute the y-coordinates of the deformations	
Az	Matrix to compute the z-coordinates of the deformations	
B	Matrix which contains the x-, y- and z-coordinates of the undeformed structure	
Cx	Matrix to compute the x-coordinates of the deformations	
Cy	Matrix to compute the y-coordinates of the deformations	
Cz	Matrix to compute the z-coordinates of the deformations	

See also DISP_TRUSS, PLOTDISP, DISP_SHELL8.

disp_plane10

DISP_PLANE10 Matrices to compute the displacements of the deformed plane.

```
[Ax,Ay,Az,B] = disp_plane10(Nodes,Elements,DOF,U)
returns the matrices to compute the displacements of the deformed plane.
The coordinates of the nodes of the plane10 element are
computed using X=Ax*U+B(:,1); Y=Ay*U+B(:,2) and
Z=Az*U+B(:,3).
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
DOF	Degrees of freedom	(nDOF * 1)
Ax	Matrix to compute the x-coordinates of the deformations	
Ay	Matrix to compute the y-coordinates of the deformations	
Az	Matrix to compute the z-coordinates of the deformations	
B	Matrix which contains the x-, y- and z-coordinates of the undeformed structure	

See also DISP_TRUSS, PLOTDISP, DISP_SHELL4.

disp_plane15

DISP_PLANE15 Matrices to compute the displacements of the deformed plane.

```
[Ax,Ay,Az,B] = disp_plane15(Nodes,Elements,DOF,U)
returns the matrices to compute the displacements of the deformed plane.
The coordinates of the nodes of the plane15 element are
computed using X=Ax*U+B(:,1); Y=Ay*U+B(:,2) and
Z=Az*U+B(:,3).
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
DOF	Degrees of freedom	(nDOF * 1)
Ax	Matrix to compute the x-coordinates of the deformations	
Ay	Matrix to compute the y-coordinates of the deformations	
Az	Matrix to compute the z-coordinates of the deformations	
B	Matrix which contains the x-, y- and z-coordinates of the undeformed structure	

See also DISP_TRUSS, PLOTDISP, DISP_SHELL4.

disp_plane3

DISP_PLANE3 Matrices to compute the displacements of the deformed plane3.

```
[Ax,Ay,Az,B] = disp_plane3(Nodes,Elements,DOF,U)
returns the matrices to compute the displacements of the deformed plane3 element.
The coordinates of the nodes of the plane3 element are
computed using X=Ax*U+B(:,1); Y=Ay*U+B(:,2) and
Z=Az*U+B(:,3).
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
DOF	Degrees of freedom	(nDOF * 1)
Ax	Matrix to compute the x-coordinates of the deformations	
Ay	Matrix to compute the y-coordinates of the deformations	
Az	Matrix to compute the z-coordinates of the deformations	
B	Matrix which contains the x-, y- and z-coordinates of the undeformed structure	

See also DISP_TRUSS, PLOTDISP, DISP_SHELL8.

disp_plane4

DISP_PLANE4 Matrices to compute the displacements of the deformed plane4.

```
[Ax,Ay,Az,B] = disp_plane4(Nodes,Elements,DOF,U)
returns the matrices to compute the displacements of the deformed plane4.
The coordinates of the nodes of the plane4 element are
computed using X=Ax*U+B(:,1); Y=Ay*U+B(:,2) and
Z=Az*U+B(:,3).
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
DOF	Degrees of freedom	(nDOF * 1)
Ax	Matrix to compute the x-coordinates of the deformations	
Ay	Matrix to compute the y-coordinates of the deformations	
Az	Matrix to compute the z-coordinates of the deformations	
B	Matrix which contains the x-, y- and z-coordinates of the undeformed structure	

See also DISP_TRUSS, PLOTDISP, DISP_SHELL8.

disp_plane6

DISP_PLANE6 Matrices to compute the displacements of the deformed plane.

```
[Ax,Ay,Az,B] = disp_plane6(Nodes,Elements,DOF,U)
returns the matrices to compute the displacements of the deformed plane.
The coordinates of the nodes of the plane6 element are
computed using X=Ax*U+B(:,1); Y=Ay*U+B(:,2) and
Z=Az*U+B(:,3).
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
DOF	Degrees of freedom	(nDOF * 1)
Ax	Matrix to compute the x-coordinates of the deformations	
Ay	Matrix to compute the y-coordinates of the deformations	
Az	Matrix to compute the z-coordinates of the deformations	
B	Matrix which contains the x-, y- and z-coordinates of the undeformed structure	

See also DISP_TRUSS, PLOTDISP, DISP_SHELL4.

disp_plane8

DISP_PLANE8 Return matrices to compute the displacements of the deformed elements.

```
[Ax,Ay,Az,B]=disp_plane8(Nodes,Elements,DOF)
```

returns the matrices to compute the displacements of the deformed elements.

The coordinates of the specified points along the deformed beams element are computed using

```
X=Ax*U+B(:,1)
```

```
Y=Ay*U+B(:,2)
```

```
Z=Az*U+B(:,3)
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
DOF	Degrees of freedom	(nDOF * 1)
Points	Points in local coordinate system	(1 * nPoints)
Ax	Matrix to compute the x-coordinates of the deformations	
Ay	Matrix to compute the y-coordinates of the deformations	
Az	Matrix to compute the z-coordinates of the deformations	
B	Matrix which contains the x-, y- and z-coordinates of the undeformed structure	

disp_shell2

DISP_SHELL2 Return matrices to compute the displacements of deformed SHELL2 elements.

```
[Ax,Ay,Az,B,Cx,Cy,Cz]
    =disp_shell2(Nodes,Elements,DOF,EltIDDLoad,Sections,Materials,Points)
[Ax,Ay,Az,B,Cx,Cy,Cz]
    =disp_shell2(Nodes,Elements,DOF,EltIDDLoad,Sections,Materials)
[Ax,Ay,Az,B]
    =disp_shell2(Nodes,Elements,DOF,[],Sections,Materials)
[Ax,Ay,Az,B]
    =disp_shell2(Nodes,Elements,DOF)
```

returns the matrices to compute the displacements of deformed SHELL2 elements. The coordinates of the specified points along the deformed SHELL2 element are computed using $X=A_x*U+C_x*DLoad+B(:,1)$; $Y=A_y*U+C_y*DLoad+B(:,2)$ and $Z=A_z*U+C_z*DLoad+B(:,3)$. The matrices C_x, C_y and C_z superimpose the displacements that occur due to the distributed loads if all nodes are fixed. In the current implementation, $C_x=C_y=C_z=0$, i.e. the local effect of the distributed loads is ignored.

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
DOF	Degrees of freedom	(nDOF * 1)
EltIDDLoad	Elements with distributed loads	[EltID] (use empty array [] when shear deformation is considered but no DLoads are present)
Sections	Section definitions	[SecID SecProp1 SecProp2 ...]
Materials	Material definitions	[MatID MatProp1 MatProp2 ...]
Points	Points in the local coordinate system	(1 * nPoints)
Ax	Matrix to compute the x-coordinates of the deformations	
Ay	Matrix to compute the y-coordinates of the deformations	
Az	Matrix to compute the z-coordinates of the deformations	
B	Matrix which contains the x-, y- and z-coordinates of the undeformed structure	
Cx	Matrix to compute the x-coordinates of the deformations	
Cy	Matrix to compute the y-coordinates of the deformations	
Cz	Matrix to compute the z-coordinates of the deformations	

disp_shell4

DISP_SHELL4 Matrices to compute the displacements of the deformed shell4.

```
[Ax,Ay,Az,B] = disp_shell4(Nodes,Elements,DOF,U)
returns the matrices to compute the displacements of the deformed shell4.
The coordinates of the nodes of the shell4 element are
computed using X=Ax*U+B(:,1); Y=Ay*U+B(:,2) and
Z=Az*U+B(:,3).
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
DOF	Degrees of freedom	(nDOF * 1)
Ax	Matrix to compute the x-coordinates of the deformations	
Ay	Matrix to compute the y-coordinates of the deformations	
Az	Matrix to compute the z-coordinates of the deformations	
B	Matrix which contains the x-, y- and z-coordinates of the undeformed structure	

See also DISP_TRUSS, PLOTDISP, DISP_SHELL8.

disp_shell6

DISP_SHELL6 Matrices to compute the displacements of the deformed shell.

```
[Ax,Ay,Az,B] = disp_shell6(Nodes,Elements,DOF,U)
returns the matrices to compute the displacements of the deformed shell.
The coordinates of the nodes of the shell6 element are
computed using X=Ax*U+B(:,1); Y=Ay*U+B(:,2) and
Z=Az*U+B(:,3).
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
DOF	Degrees of freedom	(nDOF * 1)
Ax	Matrix to compute the x-coordinates of the deformations	
Ay	Matrix to compute the y-coordinates of the deformations	
Az	Matrix to compute the z-coordinates of the deformations	
B	Matrix which contains the x-, y- and z-coordinates of the undeformed structure	

See also DISP_TRUSS, PLOTDISP, DISP_SHELL4.

disp_shell8

DISP_SHELL8 Matrices to compute the displacements of the deformed shell.

```
[Ax,Ay,Az,B] = disp_shell8(Nodes,Elements,DOF,U)
returns the matrices to compute the displacements of the deformed shell.
The coordinates of the nodes of the shell8 element are
computed using X=Ax*U+B(:,1); Y=Ay*U+B(:,2) and
Z=Az*U+B(:,3).
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
DOF	Degrees of freedom	(nDOF * 1)
Ax	Matrix to compute the x-coordinates of the deformations	
Ay	Matrix to compute the y-coordinates of the deformations	
Az	Matrix to compute the z-coordinates of the deformations	
B	Matrix which contains the x-, y- and z-coordinates of the undeformed structure	

See also DISP_TRUSS, PLOTDISP, DISP_SHELL4.

disp_solid10

DISP_SOLID10 Return matrices to compute the displacements of the deformed elements.

```
[Ax,Ay,Az,B]=disp_solid10(Nodes,Elements,DOF,Points)
[Ax,Ay,Az,B]=disp_solid10(Nodes,Elements,DOF)
returns the matrices to compute the displacements of the deformed elements.
The coordinates of the specified points along the deformed beams element are
computed using  $X=Ax*U+B(:,1)$ ;  $Y=Ay*U+B(:,2)$  and  $Z=Az*U+B(:,3)$ .
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
DOF	Degrees of freedom	(nDOF * 1)
Points	Points in local coordinate system	(1 * nPoints)
Ax	Matrix to compute the x-coordinates of the deformations	
Ay	Matrix to compute the y-coordinates of the deformations	
Az	Matrix to compute the z-coordinates of the deformations	
B	Matrix which contains the x-, y- and z-coordinates of the undeformed structure	

See also DISP_TRUSS, PLOTDISP.

disp_solid15

DISP_SOLID15 Return matrices to compute the displacements of the deformed elements.

```
[Ax,Ay,Az,B]=disp_solid15(Nodes,Elements,DOF,Points)
[Ax,Ay,Az,B]=disp_solid15(Nodes,Elements,DOF)
returns the matrices to compute the displacements of the deformed elements.
The coordinates of the specified points along the deformed beams element are
computed using  $X=Ax*U+B(:,1)$ ;  $Y=Ay*U+B(:,2)$  and  $Z=Az*U+B(:,3)$ .
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
DOF	Degrees of freedom	(nDOF * 1)
Points	Points in local coordinate system	(1 * nPoints)
Ax	Matrix to compute the x-coordinates of the deformations	
Ay	Matrix to compute the y-coordinates of the deformations	
Az	Matrix to compute the z-coordinates of the deformations	
B	Matrix which contains the x-, y- and z-coordinates of the undeformed structure	

See also DISP_TRUSS, PLOTDISP.

disp_solid20

DISP_SOLID20 Return matrices to compute the displacements of the deformed elements.

```
[Ax,Ay,Az,B]=disp_solid20(Nodes,Elements,DOF,Points)
[Ax,Ay,Az,B]=disp_solid20(Nodes,Elements,DOF)
returns the matrices to compute the displacements of the deformed elements.
The coordinates of the specified points along the deformed beams element are
computed using  $X=Ax*U+B(:,1)$ ;  $Y=Ay*U+B(:,2)$  and  $Z=Az*U+B(:,3)$ .
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
DOF	Degrees of freedom	(nDOF * 1)
Points	Points in local coordinate system	(1 * nPoints)
Ax	Matrix to compute the x-coordinates of the deformations	
Ay	Matrix to compute the y-coordinates of the deformations	
Az	Matrix to compute the z-coordinates of the deformations	
B	Matrix which contains the x-, y- and z-coordinates of the undeformed structure	

See also DISP_TRUSS, PLOTDISP.

disp_solid8

DISP_SOLID8 Return matrices to compute the displacements of the deformed elements.

```
[Ax,Ay,Az,B]=disp_solid8(Nodes,Elements,DOF,Points)
```

```
[Ax,Ay,Az,B]=disp_solid8(Nodes,Elements,DOF)
```

returns the matrices to compute the displacements of the deformed elements.

The coordinates of the specified points along the deformed beams element are computed using $X=Ax*U+B(:,1)$; $Y=Ay*U+B(:,2)$ and $Z=Az*U+B(:,3)$.

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
DOF	Degrees of freedom	(nDOF * 1)
Points	Points in local coordinate system	(1 * nPoints)
Ax	Matrix to compute the x-coordinates of the deformations	
Ay	Matrix to compute the y-coordinates of the deformations	
Az	Matrix to compute the z-coordinates of the deformations	
B	Matrix which contains the x-, y- and z-coordinates of the undeformed structure	

See also DISP_TRUSS, PLOTDISP.

disp_truss

DISP_TRUSS Return matrices to compute the displacements of the deformed trusses.

```
[Ax,Ay,Az,B]=disp_truss(Nodes,Elements,DOF,[],[],[],Points)
[Ax,Ay,Az,B]=disp_truss(Nodes,Elements,DOF)
    returns the matrices to compute the displacements of the deformed
    trusses. The coordinates of the specified points along the deformed
    truss elements are computed using X=Ax*U+B(:,1); Y=Ay*U+B(:,2)
    and Z=Az*U+B(:,3).
```

```
[Ax,Ay,Az,B,Cx,Cy,Cz,dAxdx,dAydx,dAzdx,dCxdx,dCydx,dCzdx]
    = DISP_TRUSS(Nodes,Elements,DOF,EltIDDLoad,Sections,Materials,Points,dNodesdx,
    dDLoadsdx(,dSectionsdx))
    additionally computes the derivatives of the displacements with
    respect to the design variables x.
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
DOF	Degrees of freedom	(nDOF * 1)
Points	Points in the local coordinate system	(1 * nPoints)
dNodesdx	Node definitions derivatives	(SIZE(Node) * nVar)
dDLoads	Distributed loads derivatives	(SIZE(DLoad) * nVar)
Ax	Matrix to compute the x-coordinates of the deformations	
Ay	Matrix to compute the y-coordinates of the deformations	
Az	Matrix to compute the z-coordinates of the deformations	
B	Matrix which contains the x-, y- and z-coordinates of the undeformed structure	

```
dAxdx, dAydx, dAzdx, dCxdx, dCydx, dCzdx
    Derivatives of the matrices to compute the coordinates of the
    interpolation points after deformation
```

See also DISP_BEAM, PLOTDISP.

dloadgcs2lcs

DLOADGCS2LCS Distributed load transformation from GCS to LCS.

```
DLoadLCS = DLOADGCS2LCS(T,DLoad)
    transforms the distributed load definitions in the global
    coordinate system (algebraic convention) to the local coordinate
    system (beam convention).
```

```
[DLoadLCS,dDLoadLCSdx] = DLOADGCS2LCS(T,DLoad,dTdx,dDLoaddx)
    transforms the distributed load definitions in the global
    coordinate system (algebraic convention) to the local coordinate
    system (beam convention), and additionally computes the derivatives
    of the distributed load information with respect to the design
    variables x.
```

T	Element transformation matrix	(6 * 6)
DLoad	Distributed loads in GCS (6/8 * nLC * nDLoads)	[n1globalX; n1globalY; n1globalZ; ...]
dTdx	Transformation matrix derivatives	(6 * 6 * nVar)
dDLoaddx	Distributed loads derivatives (GCS)	(SIZE(DLoad) * nVar)
DLoadLCS	Distributed loads in LCS (6/8 * nLC * nDLoads)	[n1localX; n1localY; n1localZ; ...]
dDLoadLCSdx	Distributed loads derivatives (LCS)	(SIZE(DLoadLCS) * nVar)

dof_beam

DOF_BEAM Element degrees of freedom for a beam element.

DOF = dof_beam(NodeNum) builds the vector with the
builds the vector with the degrees of freedom for the beam element.

NodeNum	Node definitions	[NodID1 NodID2]	(1 * 2)
DOF	Degrees of freedom		(12 * 1)

See also GETDOF.

dof_mass

DOF_MASS Element degrees of freedom for a mass element.

dof = dof_truss(NodeNum) builds the vector with the labels of the degrees of freedom for which stiffness is present in the mass element.

NodeNum	Node numbers	[NodID1]	(1)
dof	Degrees of freedom	(6 * 1)	

See also GETDOF.

dof_plane10

DOF_PLANE10 Element degrees of freedom for a plane10 element.

dof = dof_plane10(NodeNum) builds the vector with the labels of the degrees of freedom for which stiffness is present in the plane10 element.

NodeNum	Node definitions	[NodID1 NodID2 ... NodIDn]	(1 * 10)
dof	Degrees of freedom		(20 * 1)

See also GETDOF.

dof_plane15

DOF_PLANE15 Element degrees of freedom for a plane15 element.

dof = dof_plane15(NodeNum) builds the vector with the labels of the degrees of freedom for which stiffness is present in the plane15 element.

NodeNum	Node definitions	[NodID1 NodID2 ... NodIDn]	(1 * 15)
dof	Degrees of freedom		(30 * 1)

See also GETDOF.

dof_plane3

DOF_PLANE3 Element degrees of freedom for a plane3 element.

dof = dof_plane3(NodeNum) builds the vector with the labels of the degrees of freedom for which stiffness is present in the plane3 element.

NodeNum	Node definitions	[NodID1 NodID2 NodID3]	(1 * 3)
dof	Degrees of freedom		(6 * 1)

See also GETDOF.

dof_plane4

DOF_PLANE4 Element degrees of freedom for a plane4 element.

dof = dof_plane4(NodeNum) builds the vector with the labels of the degrees of freedom for which stiffness is present in the plane4 element.

NodeNum	Node definitions	[NodID1 NodID2 ... NodIDn]	(1 * 4)
dof	Degrees of freedom		(8 * 1)

See also GETDOF.

dof_plane6

DOF_PLANE6 Element degrees of freedom for a plane6 element.

dof = dof_plane6(NodeNum) builds the vector with the labels of the degrees of freedom for which stiffness is present in the plane6 element.

NodeNum	Node definitions	[NodID1 NodID2 ... NodIDn]	(1 * 6)
dof	Degrees of freedom		(12 * 1)

See also GETDOF.

dof_plane8

of_plane8 is a function.

```
dof = dof_plane8(Nodenumbers)
```


dof_shell2

DOF_SHELL2 Element degrees of freedom for a SHELL2 element.

DOF = dof_shell2(NodeNum) returns a vector with the degrees of freedom for a SHELL2 element.

NodeNum	Node definitions	[NodID1 NodID2]	(1 * 2)
DOF	Degrees of freedom		(12 * 1)

dof_shell4

DOF_SHELL4 Element degrees of freedom for a shell4 element.

dof = dof_shell4(NodeNum) builds the vector with the labels of the degrees of freedom for which stiffness is present in the shell4 element.

NodeNum	Node definitions	[NodID1 NodID2 ... NodIDn]	(1 * 4)
dof	Degrees of freedom		(24 * 1)

See also GETDOF.

dof_shell6

DOF_SHELL6 Element degrees of freedom for a shell6 element.

dof = dof_shell6(NodeNum) builds the vector with the labels of the degrees of freedom for which stiffness is present in the shell6 element.

NodeNum	Node definitions	[NodID1 NodID2 ... NodIDn]	(1 * 6)
dof	Degrees of freedom		(36 * 1)

See also GETDOF.

dof_shell8

DOF_SHELL8 Element degrees of freedom for a shell8 element.

dof = dof_shell8(NodeNum) builds the vector with the labels of the degrees of freedom for which stiffness is present in the shell8 element.

NodeNum	Node definitions	[NodID1 NodID2 ... NodIDn]	(1 * 8)
dof	Degrees of freedom		(48 * 1)

See also GETDOF.

dof_solid10

DOF_SOLID10 Element degrees of freedom for a solid10 element.

dof = dof_solid10(NodeNum) builds the vector with the labels of the degrees of freedom for which stiffness is present in the solid10 element.

NodeNum	Node definitions	[NodID1 NodID2 ... NodIDn]	(1 * 4)
dof	Degrees of freedom		(30 * 1)

See also GETDOF.

dof_solid15

DOF_SOLID15 Element degrees of freedom for a solid10 element.

dof = dof_solid15(NodeNum) builds the vector with the labels of the degrees of freedom for which stiffness is present in the solid15 element.

NodeNum	Node definitions	[NodID1 NodID2 ... NodIDn]	(1 * 4)
dof	Degrees of freedom		(45 * 1)

See also GETDOF.

dof_solid20

DOF_SOLID20 Element degrees of freedom for a solid20 element.

dof = dof_solid20(NodeNum) builds the vector with the labels of the degrees of freedom for which stiffness is present in the solid20 element.

NodeNum	Node definitions	[NodID1 NodID2 ... NodIDn]	(1 * 4)
dof	Degrees of freedom		(60 * 1)

See also GETDOF.

dof_solid4

DOF_SOLID4 Element degrees of freedom for a solid4 element.

dof = dof_solid4(NodeNum) builds the vector with the labels of the degrees of freedom for which stiffness is present in the solid4 element.

NodeNum	Node definitions	[NodID1 NodID2 NodID3 NodID4]	(1 * 4).
dof	Degrees of freedom		(12 * 1).

See also GETDOF.

dof_solid8

DOF_SOLID8 Element degrees of freedom for a solid8 element.

dof = dof_solid8(NodeNum) builds the vector with the labels of the degrees of freedom for which stiffness is present in the solid8 element.

NodeNum	Node definitions	[NodID1 NodID2 ... NodIDn]	(1 * 4)
dof	Degrees of freedom		(24 * 1)

See also GETDOF.

dof_truss

DOF_TRUSS Element degrees of freedom for a truss element.

dof = dof_truss(NodeNum) builds the vector with the degrees of freedom for the truss element.

NodeNum	Node numbers	[NodID1 NodID2]	(1 * 2)
dof	Degrees of freedom		(6 * 1)

See also GETDOF.

eigfem

EIGFEM Compute the eigenmodes and eigenfrequencies of the finite element model.

```
[phi,omega]=eigfem(K,M,nMode)
```

```
[phi,omega]=eigfem(K,M)
```

computes the eigenmodes and eigenfrequencies of the finite element model.

K Stiffness matrix (nDOF * nDOF)

M Mass matrix (nDOF * nDOF)

nMode Number of eigenmodes and eigenfrequencies (default: all)

phi Eigenmodes (in columns) (nDOF * nMode)

omega Eigenfrequencies [rad/s] (nMode * 1)

elemdisp

ELEMDISP Select the element displacements from the global displacement vector.

UeGCS = elemdisp(Type,NodeNum,DOF,U) selects the element displacements from the global displacement vector.

Type	Element type e.g. 'beam','truss', ...
NodeNum	Node numbers (1 * nNodes)
DOF	Degrees of freedom (nDOF * 1)
U	Displacements (nDOF * nLC)
UeGCS	Element displacements

See also ELEMFORCES, DOF_BEAM, DOF_TRUSS.

elemforces

ELEMFORCES Compute the element forces.

```
[ForcesLCS,ForcesGCS] = ELEMFORCES(Nodes,Elements,Types,Sections,Materials,DOF,U,DLoads,TLoads)
[ForcesLCS,ForcesGCS] = ELEMFORCES(Nodes,Elements,Types,Sections,Materials,DOF,U)
    computes the element forces in the local (beam convention) and the
    global (algebraic convention) coordinate system.
```

```
[ForcesLCS,ForcesGCS,dForcesLCSdx,dForcesGCSdx]
    = ELEMFORCES(Nodes,Elements,Types,Sections,Materials,DOF,U,DLoads,TLoads,
                  dNodesdx,dSectionsdx,dUdx,dDLoadsdx)
    additionally computes the derivatives of the element forces with
    respect to the design variables x.
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Types	Element type definitions	{TypID EltName Option1 ...}
Sections	Section definitions	[SecID SecProp1 SecProp2 ...]
Materials	Material definitions	[MatID MatProp1 MatProp2 ...]
DOF	Degrees of freedom	(nDOF * 1)
U	Displacements	(nDOF * nLC)
DLoads	Distributed loads	[EltID n1globalX n1globalY n1globalZ ...]
TLoads	Temperature loads	[EltID Tyt Tyb Tzt Tzb]
dNodesdx	Node definitions derivatives	(SIZE(Nodes) * nVar)
dSectionsdx	Section definitions derivatives	(SIZE(Sections) * nVar)
dUdx	Displacements derivatives	(nDOF * nLC * nVar)
dDLoadsdx	Distributed loads derivative	(SIZE(DLoads) * nVar)
ForcesLCS	Element forces in LCS (beam convention)	[N Vy Vz T My Mz] (nElem * 12 * nLC)
ForcesGCS	Element forces in GCS (algebraic convention)	(nElem * 12 * nLC)
dForcesLCSdx	Element forces derivatives in LCS	(nElem * 12 * nLC * nVar)
dForcesGCSdx	Element forces derivatives in GCS	(nElem * 12 * nLC * nVar)

See also FORCES_TRUSS, FORCES_BEAM.

elemloads

ELEMLOADS Equivalent nodal forces.

```
F = ELEMLOADS(DLoads,Nodes,Elements,Types,DOF)
computes the equivalent nodal forces of a distributed load
(in the global coordinate system).
```

```
[F,dFdx] = ELEMLOADS(DLoads,Nodes,Elements,Types,DOF,dDLoadsdx,dNodesdx)
additionally computes the derivatives of the equivalent nodal forces
with respect to the design variables x.
```

DLoads	Distributed loads	[EltID n1globalX n1globalY n1globalZ ...]
Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Types	Element type definitions	{TypID EltName Option1 ... }
DOF	Degrees of freedom	(nDOF * 1)
dDLoadsdx	Distributed loads derivatives	(SIZE(DLoads) * nVar)
dNodesdx	Node definitions derivatives	(SIZE(Nodes) * nVar)
F	Load vector	(nDOF * nLC)
dFdx	Load vector derivatives	(nDOF * nLC * nVar)

See also LOADS_TRUSS, LOADS_BEAM, NODALVALUES.

elempressure

ELEMPRESSURE Equivalent nodal forces for a pressure load on a shell element.

```
F = elempressure(Pressures,Nodes,Elements,Types,DOF)
computes the equivalent nodal forces of a distributed pressure
(in the local coordinate system xyz, with z perpendicular to the surface).
```

Pressures	Pressure on surface or edge	[EltID n1localZ n2localZ n3localZ ...]
Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Types	Element type definitions	{TypID EltName Option1 ... }
DOF	Degrees of freedom	(nDOF * 1)
F	Load vector	(nDOF * 1)

See also PRESSURE_SHELL8, PRESSURE_SHELL4, NODALVALUES.

elemsize

ELEMSIZE Compute element length/area/volume.

```
S = ELEMSIZE(Nodes,Elements,Types)
```

computes the size of all elements, depending on element type. For a 1D line element it computes length, for a 2D plate element it computes area, and for a 3D solid element it computes volume.

```
[S,dSdx] = ELEMSIZE(Nodes,Elements,Types,dNodesdx)
```

additionally computes the derivatives of the size with respect to the design variables x.

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Types	Element type definitions	{TypID EltName Option1 ... }
dNodesdx	Node definitions derivatives	(SIZE(Nodes) * nVar)
S	Element sizes	
dSdx	Element sizes derivatives	

See also `SIZE_BEAM`, `SIZE_TRUSS`, `ELEMVOLUMES`.

ELEMSTRESS Compute the element stresses.

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Types	Element type definitions	{TypID EltName Option1 ... }
Sections	Section definitions	[SecID SecProp1 SecProp2 ...]
Materials	Material definitions	[MatID MatProp1 MatProp2 ...]
DOF	Degrees of freedom (nDOF * 1)	
U	Displacements (nDOF * 1)	

```
'GCS'      Sets the gcs in which SeGCS is computed.
            Default: 'cart'. Type of values:
            'cart': cartesian coordinate system
            'cyl'  : cylindrical coordinate system
            'sph'  : spherical coordinate system
```

See also SE_SHELL8, SE_SHELL4.

elemtloads

ELEMTLOADS Equivalent nodal forces for temperature loading.

`F=elemtloads(TLoads,Nodes,Elements,Types,Sections,Materials,DOF)`
 computes the equivalent nodal forces of a temperature gradient
 (in the global coordinate system).

TLoads	Temperature gradient	[EltID Tyt Tyb Tzt Tzb]
	Tkt and Tkb correspond to the temperatures at the top and the bottom of the profile when the k-axis (LCS) points up (k=y,z)	
Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Types	Type definitions	{TypID EltName Option1 ... }
Sections	Section definitions	[SecID A ky kz Ixx Iyy Izz yt yb zt zb]
Materials	Material definitions	[MatID E nu rho alpha]
DOF	Degrees of freedom	(nDOF * 1)
F	Load vector	(nDOF * 1)

See also TLOADS_TRUSS, TLOADS_BEAM.

elemvolumes

ELEMVOLUMES Compute element volumes.

```
V = ELEMVOLUMES(Nodes,Elements,Types,Sections)
computes the volume for all elements.
```

```
[V,dVdx] = ELEMVOLUMES(Nodes,Elements,Types,Sections,dNodesdx,dSectionsdx)
computes the volume for all elements, as well as the derivatives of the
volume with respect to the design variables x.
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Types	Element type definitions	{TypID EltName Option1 ... }
Sections	Section definitions	[SecID SecProp1 SecProp2 ...]
dNodesdx	Node definitions derivatives	(SIZE(Nodes) * nVar)
dSectionsdx	Section definitions derivatives	(SIZE(Sections) * nVar)
V	Element volumes	(nElem * 1)
dVdx	Element volumes derivatives	(nElem * nVar)

See also VOLUME_BEAM, VOLUME_TRUSS, ELEMSIZES.

fdiargcs_beam

FDIARGCS_BEAM Return matrices to plot the forces in a beam element.

```
[ElemGCS,FdiagrGCS,ElemExtGCS,ExtremaGCS,Extrema]
    = fdiargcs_beam(ftype,Forces,Node,[],[],DLoad,Points)
[ElemGCS,FdiagrGCS,ElemExtGCS,ExtremaGCS,Extrema]
    = fdiargcs_beam(ftype,Forces,Node,[],[],DLoad)
```

returns the coordinates of the points along the beam in the global coordinate system and the coordinates of the forces with respect to the beam in the global coordinate system. These can be added in order to plot the forces: ElemGCS+FdiagrGCS. The coordinates of the points with extreme values and the coordinates of the extreme values with respect to the beam are given as well and can be similarly added: ElemExtGCS+ExtremaGCS. Extrema is the list with the correspondig extreme values.

ftype	'norm'	Normal force (in the local x-direction)
	'sheary'	Shear force in the local y-direction
	'shearz'	Shear force in the local z-direction
	'momx'	Torsional moment (around the local x-direction)
	'momy'	Bending moment around the local y-direction
	'momz'	Bending moment around the local z-direction
Forces	Element forces in LCS (beam convention) [N; Vy; Vz; T; My; Mz] (12 * 1)	
Node	Node definitions	[x y z] (3 * 3)
DLoad	Distributed loads	[n1globalX; n1globalY; n1globalZ; ...] (6 * 1)
Points	Points in the local coordinate system (1 * nPoints)	
ElemGCS	Coordinates of the points along the beam in GCS (nPoints * 3)	
FdiagrGCS	Coordinates of the force with respect to the beam in GCS (nValues * 3)	
ElemExtGCS	Coordinates of the points with extreme values in GCS (nValues * 3)	
ExtremaGCS	Coordinates of the extreme values with respect to the beam in GCS (nValues * 3)	
Extrema	Extreme values (nValues * 1)	

See also PLOTFORC, FDIAGRLCS_BEAM, FDIARGCS_TRUSS.

fdiargcs_shell2

FDIARGCS_SHELL2 Return matrices to plot the forces in a SHELL2 element.

```
[ElemGCS,FdiagrGCS,ElemExtGCS,ExtremaGCS,Extrema]
    = fdiargcs_shell2(ftype,Forces,Node,Section,Material,DLoad,Points)
[ElemGCS,FdiagrGCS,ElemExtGCS,ExtremaGCS,Extrema]
    = fdiargcs_shell2(ftype,Forces,Node,Section,Material,DLoad)
```

returns the coordinates of the points along the SHELL2 in the global coordinate system and the coordinates of the forces with respect to the element in the global coordinate system. These can be added in order to plot the forces: ElemGCS+FdiagrGCS. The coordinates of the points with extreme values and the coordinates of the extreme values with respect to the element are given as well and can be similarly added: ElemExtGCS+ExtremaGCS. Extrema is the list with the correspondig extreme values.

ftype	'Nphi'	Normal force (per unit length) in meridional direction
	'Qphi'	Transverse force (per unit length) in meridional direction
	'Mphi'	Bending moment (per unit length) in meridional direction
	'Ntheta'	Normal force (per unit length) in circumferential direction
	'Mtheta'	Bending moment (per unit length) in circumferential direction
Forces	Element forces in LCS (beam convention) [N; Vy; 0; 0; 0; Mz] (12 * 1)	
Node	Node definitions [x y z] (3 * 3)	
DLoad	Distributed loads [n1globalX; n1globalY; n1globalZ; ...] (6 * 1)	
Points	Points in the local coordinate system (1 * nPoints)	
ElemGCS	Coordinates of the points along the element in GCS (nPoints * 3)	
FdiagrGCS	Coordinates of the force with respect to the element in GCS (nValues * 3)	
ElemExtGCS	Coordinates of the points with extreme values in GCS (nValues * 3)	
ExtremaGCS	Coordinates of the extreme values with respect to the element in GCS (nValues * 3)	
Extrema	Extreme values (nValues * 1)	

fdiargcs_truss

FDIARGCS_TRUSS Return matrices to plot the forces in a truss element.

```
[ElemGCS,FdiagrGCS,ElemExtGCS,ExtremaGCS,Extrema]
    = fdiargcs_truss(ftype,Forces,Node,[],[],[],Points)
[ElemGCS,FdiagrGCS,ElemExtGCS,ExtremaGCS,Extrema]
    = fdiargcs_truss(ftype,Forces,Node)
```

returns the coordinates of the points along the truss in the global coordinate system and the coordinates of the forces with respect to the truss in the global coordinate system. These can be added in order to plot the forces: ElemGCS+FdiagrGCS. The coordinates of the points with extreme values and the coordinates of the extreme values with respect to the truss are given as well and can be similarly added: ElemExtGCS+ExtremaGCS. Extrema is the list with the correspondig extreme values.

```
ftype      'norm'      Normal force (in the local x-direction)
Forces      Element forces in LCS [N; 0; 0; 0; 0; 0] (12 * 1)
Node        Node definitions      [x y z] (3 * 3)
Points      Points in the local coordinate system (1 * nPoints)
ElemGCS      Coordinates of the points along the truss in GCS (nPoints * 3)
FdiagrGCS    Coordinates of the force with respect to the truss in GCS
                                     (nValues * 3)
ElemExtGCS   Coordinates of the points with extreme values in GCS (nValues * 3)
ExtremaGCS   Coordinates of the extreme values with respect to the truss in GCS
                                     (nValues * 3)

Extrema      Extreme values (nValues * 1)
```

See also PLOTFORC, FDIARGCS_BEAM.

fdiagrlcs

FDIAGR LCS Return force diagrams in LCS.

```
FdiagrLCS = FDIAGR LCS(ftype,Nodes,Elements,Types,Forces,DLoads,Points)
FdiagrLCS = FDIAGR LCS(ftype,Nodes,Elements,Types,Forces,DLoads)
FdiagrLCS = FDIAGR LCS(ftype,Nodes,Elements,Types,Forces)
    computes element force values in all interpolation points (in beam
    convention) for beam and truss elements.
```

```
[FdiagrLCS,dFdiagrLCSdx]
    = FDIAGR LCS(ftype,Nodes,Elements,Types,Forces,DLoads,...
                  Points,dNodesdx,dForcesdx,dDLoadsdx)
    additionally computes the derivatives of the force values with
    respect to the design variables x.
```

ftype	'norm'	Normal force (in the local x-direction)
	'sheary'	Shear force in the local y-direction
	'shearz'	Shear force in the local z-direction
	'momx'	Torsional moment (around the local x-direction)
	'momy'	Bending moment around the local y-direction
	'momz'	Bending moment around the local z-direction
Nodes	Node definitions [NodID x y z]	
Elements	Element definitions [EltID TypID SecID MatID n1 n2 ...]	
Types	Element type definitions {TypID EltName Option1 ... }	
Forces	Element forces in LCS (beam convention) [N Vy Vz T My Mz] (nElem * 12)	
DLoads	Distributed loads [EltID n1globalX n1globalY n1globalZ ...]	
Points	Points in the local coordinate system (1 * nPoints)	
dNodesdx	Node definitions derivatives (SIZE(Nodes) * nVar)	
dForcesdx	Element forces in LCS derivatives (SIZE(Forces) * nVar)	
dDLoadsdx	Distributed loads derivatives (SIZE(DLoads) * nVar)	
FdiagrLCS	Element force values at the points (nElem * nPoints * nLC)	
dFdiagrLCSdx	Element force values derivatives (nElem * nPoints * nLC * nVar)	

See also PLOTFORC, FDIAGR GCS_BEAM, FDIAGR GCS_TRUSS.

fdiagrlcs_beam

FDIAGRLCS_BEAM Force diagram for a beam element in LCS.

[FdiagrLCS,loc,Extrema] = FDIAGRLCS_BEAM(ftype,Forces,DLoadLCS,L,Points)
 computes the elements forces at the specified points. The extreme values for an element with a single DLoad are analytically determined. The extreme values for an element with multiple DLoads are calculated in the interpolation points only.

[FdiagrLCS,loc,Extrema,dFdiagrLCSdx]
 = FDIAGRLCS_BEAM(ftype,Forces,DLoadLCS,L,Points,dForcesdx,dDLoadLCSdx,dLdx)
 additionally computes the derivatives of the element force values with respect to the design variables x.

ftype	'norm'	Normal force (in the local x-direction)
	'sheary'	Shear force in the local y-direction
	'shearz'	Shear force in the local z-direction
	'momx'	Torsional moment (around the local x-direction)
	'momy'	Bending moment around the local y-direction
	'momz'	Bending moment around the local z-direction
Forces	Element forces in LCS (beam convention) [N; Vy; Vz; T; My; Mz] (12 * nLC)	
DLoadLCS	Distributed loads in LCS [n1localX; n1localY; n1localZ; ...] (6 * nLC)	
L	Beam length	
Points	Points in the local coordinate system (1 * nPoints)	
dForcesdx	Element forces in LCS derivatives	(SIZE(Forces) * nVar)
dDLoadLCSdx	Distributed loads derivatives	(SIZE(DLoadLCS) * nVar)
dLdx	Beam length derivatives	(1 * nVar)
FdiagrLCS	Element forces at the points	(1 * nPoints * nLC)
loc	Locations of the extreme values	(nValues * nLC)
Extrema	Extreme values	(nValues * nLC)
	loc and Extrema are only calculated when nLC = 1 (for plotting).	
	If this is not the case their calculation is omitted for efficiency.	
dFdiagrLCSdx	Element force value derivatives	(1 * nPoints * nLC * nVar)

See also FDIAGRGCS_BEAM.

fdiagrlcs_shell2

FDIAGRCS_SHELL2 Force diagram for a SHELL2 element in LCS.

```
[FdiagrLCS,loc,Extrema] = fdiagrlcs_shell2(ftype,Forces,DLoadLCS,L,Points)
computes the elements forces at the specified points. The extreme values are
obtained by enumeration.
```

ftype	'Nphi'	Normal force (per unit length) in meridional direction
	'Qphi'	Transverse force (per unit length) in meridional direction
	'Mphi'	Bending moment (per unit length) in meridional direction
	'Ntheta'	Normal force (per unit length) in circumferential direction
	'Mtheta'	Bending moment (per unit length) in circumferential direction
Forces	Element forces in LCS (beam convention) [N; Vy; 0; 0; 0; Mz] (12 * 1)	
DLoadLCS	Distributed loads in LCS [nlocalX; nlocalY; nlocalZ; ...] (6 * 1) or (12 * 1)	
Points	Points in the local coordinate system (1 * nPoints)	
FdiagrLCS	Element forces at the points (1 * nPoints)	
loc	Locations of the extreme values (nValues * 1)	
Extrema	Extreme values (nValues * 1)	

fdiagrlcs_truss

FDIAGRLCS_TRUSS Force diagram for a truss element in LCS.

```
[FdiagrLCS,loc,Extrema] = FDIAGRLCS_TRUSS(ftype,Forces,DLoadLCS,L,Points)
computes the element forces at the specified points.
```

```
[FdiagrLCS,loc,Extrema,dFdiagrLCSdx]
    = FDIAGRLCS_TRUSS(ftype,Forces,DLoadLCS,L,Points,dForcesdx,dDLoadLCSdx,dLdx)
additionally computes the derivatives of the element force values with
respect to the design variables x.
```

ftype	'norm'	Normal force (in the local x-direction)
Forces	Element forces in LCS	[N; 0; 0; 0; 0; 0] (12 * 1)
Node	Node definitions	[x y z] (3 * 3)
Points	Points in the local coordinate system	(1 * nPoints)
dForcesdx	Element forces in LCS derivatives	(SIZE(Forces) * nVar)
FdiagrLCS	Element forces at the points	(1 * nPoints * nLC)
loc	Locations of the extreme values	(nValues * nLC)
Extrema	Extreme values	(nValues * nLC)
dFdiagrLCSdx	Element force value derivatives	(1 * nPoints * nLC * nVar)

See also FDIAGRGCS_TRUSS.

forceslcs_beam

FORCESLCS_BEAM Compute the element forces for a beam element in the LCS.

```
Forces = FORCESLCS_BEAM(KeLCS,UeLCS,DLoadLCS,L,TLoadLCS,A,E,alpha,Iyy,Izz,hy,hz)
Forces = FORCESLCS_BEAM(KeLCS,UeLCS,DLoadLCS,L)
Forces = FORCESLCS_BEAM(KeLCS,UeLCS)
```

computes the element forces for the beam element in the local coordinate system (algebraic convention).

```
[Forces,dForcesdx] = FORCESLCS_BEAM(KeLCS,UeLCS,DLoadLCS,L,[],[],[],...
                                     [],[],[],[],[],dKeLCSdx,dUeLCSdx,dDLoadLCSdx,dLdx)
```

```
[Forces,dForcesdx] = FORCESLCS_BEAM(KeLCS,UeLCS,[],[],[],[],[],[],[],...
                                     [],[],[],dKeLCSdx,dUeLCSdx,dDLoadLCSdx,dLdx)
```

additionally computes the derivatives of the element forces with respect to the design variables x.

KeLCS	Element stiffness matrix	(12 * 12)
UeLCS	Displacements	(12 * nLC)
DLoadLCS	Distributed loads	[n1localX; n1localY; n1localZ; ...]
L	Beam length	
DLoadLCS	Distributed loads	[n1localX; n1localY; n1localZ; ...] (6 * 1)
TLoadLCS	Temperature load	[dTm; dTy; dTz] (3 * 1)
A	Cross-sectional area	
E	Young's modulus	
alpha	Linear thermal expansion coefficient	
Iyy	Area moment of inertia for bending around local y-axis	
Izz	Area moment of inertia for bending around local z-axis	
hy	Profile height (in local y-direction)	
hz	Profile height (in local z-direction)	
dKeLCSdx	Element stiffness matrix derivatives	(CELL(nVar,1))
dUeLCSdx	Displacements derivatives	(SIZE(UeLCS) * nVar)
dDLoadLCSdx	Distributed loads derivatives	(SIZE(DLoadLCS) * nVar)
dLdx	Beam length derivatives	(1 * nVar)
Forces	Element forces	[N; Vy; Vz; T; My; Mz] (12 * nLC)
dForcesdx	Element forces derivatives	(12 * nLC * nVar)

See also **FORCES_BEAM**, **FORCESLCS_TRUSS**, **ELEMFORCES**

forceslcs_truss

FORCESLCS_TRUSS Compute the element forces for a truss element in the LCS.

```
Forces = FORCESLCS_TRUSS(KeLCS,UeLCS,dTm,A,E,alpha)
Forces = FORCESLCS_TRUSS(KeLCS,UeLCS)
    computes the element forces for the truss element in the local coordinate
    system (algebraic convention).

[Forces,dForcesdx] = FORCESLCS_TRUSS(KeLCS,UeLCS,[],[],[],[],dKeLCSdx,dUeLCSdx)
    additionally computes the derivatives of the element forces with
    respect to the design variables x.
```

KeLCS	Element stiffness matrix	(6 * 6)	
UeLCS	Displacements	(6 * nLC)	
dKeLCSdx	Element stiffness matrix derivatives		(CELL(nVar,1))
dUeLCSdx	Displacements derivatives		(SIZE(UeLCS) * nVar)
Forces	Element forces		[N; 0; 0] (6 * nLC)
dForcesdx	Element forces derivatives		(6 * nLC * nVar)

See also **FORCES_TRUSS**, **FORCESLCS_BEAM**, **ELEMFORCES**

forces_beam

FORCES_BEAM Compute the element forces for a beam element.

```
[ForcesLCS,ForcesGCS] = FORCES_BEAM(Node,Section,Material,UeGCS,DLoad,TLoad,Options)
[ForcesLCS,ForcesGCS] = FORCES_BEAM(Node,Section,Material,UeGCS,DLoad)
[ForcesLCS,ForcesGCS] = FORCES_BEAM(Node,Section,Material,UeGCS)
    computes the element forces for the beam element in the local and
    the global coordinate system (algebraic convention).
```

```
[ForcesLCS,ForcesGCS,dForcesLCSdx,dForcesGCSdx]
    = FORCES_BEAM(Node,Section,Material,UeGCS,DLoad,TLoad,Options,dNodedx,...
                  dSectiondx,dUeGCSdx,dDLoaddx)
    = FORCES_BEAM(Node,Section,Material,UeGCS,[],[],Options,dNodedx,...
                  dSectiondx,dUeGCSdx)
    additionally computes the derivatives of the element forces with
    respect to the design variables x.
```

Node	Node definitions	[x y z] (3 * 3)
Section	Section definition	[A ky kz Ixx Iyy Izz]
Material	Material definition	[E nu]
UeGCS	Displacements	(12 * nLC)
DLoad	Distributed loads	[n1globalX; n1globalY; n1globalZ; ...] (6 * 1)
TLoad	TLoad	[dTm; dTy; dTz] (3 * 1)
Options	Element options	{Option1 Option2 ...}
dNodedx	Node definitions derivatives	(SIZE(Node) * nVar)
dSectiondx	Section definitions derivatives	(SIZE(Section) * nVar)
dUeGCSdx	Displacements derivatives	(SIZE(UeGCS) * nVar)
dDLoaddx	Distributed loads derivatives	(SIZE(DLoad) * nVar)
ForcesLCS	Element forces in the LCS	(12 * nLC)
ForcesGCS	Element forces in the GCS	(12 * nLC)
dForcesLCSdx	Element forces derivatives in LCS	(12 * nLC * nVar)
dForcesGCSdx	Element forces derivatives in GCS	(12 * nLC * nVar)

See also **FORCESLCS_BEAM**, **ELEMFORCES**.

forces_shell2

FORCES_BEAM Compute the element forces for a SHELL2 element.

```
[ForcesLCS,ForcesGCS]=forces_shell2(Node,Section,Material,UeGCS,DLoad,TLoad,Options)
[ForcesLCS,ForcesGCS]=forces_shell2(Node,Section,Material,UeGCS,DLoad)
[ForcesLCS,ForcesGCS]=forces_shell2(Node,Section,Material,UeGCS)
computes the element forces for the SHELL2 element in the local and the
global coordinate system (algebraic convention).
```

Node	Node definitions	[x y z] (3 * 3)
Section	Section definition	[h]
Material	Material definition	[E nu]
UeGCS	Displacements (12 * 1)	
DLoad	Distributed loads	[n1globalX; n1globalY; n1globalZ; ...] (6 * 1)
TLoad	TLoad	[dTm; dTy; dTz] (3 * 1)
Options	Element options	{Option1 Option2 ...}
ForcesLCS	Element forces in the LCS	(12 * 1)
ForcesGCS	Element forces in the GCS	(12 * 1)

forces_truss

FORCES_TRUSS Compute the element forces for a truss element.

```
[ForcesLCS,ForcesGCS] = FORCES_TRUSS(Node,Section,Material,UeGCS,[],TLoad)
[ForcesLCS,ForcesGCS] = FORCES_TRUSS(Node,Section,Material,UeGCS)
    computes the element forces for the truss element in the local and the
    global coordinate system (algebraic convention).

[ForcesLCS,ForcesGCS,dForcesLCSdx,dForcesGCSdx]
    = FORCES_TRUSS(Node,Section,Material,UeGCS,[],TLoad,[],dNodedx,...
    dSectiondx,dUeGCSdx)
    additionally computes the derivatives of the element forces with
    respect to the design variables x.
```

Node	Node definitions	[x y z] (2 * 3)
Section	Section definition	[A]
Material	Material definition	[E]
UeGCS	Displacements	(6 * nLC)
TLoad	Temperature load	[dTm]
Options	Element options	{Option1 Option2 ...}
dNodedx	Node definitions derivatives	(SIZE(Node) * nVar)
dSectiondx	Section definitions derivatives	(SIZE(Section) * nVar)
dUeGCSdx	Displacements derivatives	(SIZE(UeGCS) * nVar)
dDLoaddx	Distributed loads derivatives	(SIZE(DLoad) * nVar)
ForcesLCS	Element forces in the LCS	(12 * nLC)
ForcesGCS	Element forces in the GCS	(12 * nLC)
dForcesLCSdx	Element forces derivatives in LCS	(12 * nLC * nVar)
dForcesGCSdx	Element forces derivatives in GCS	(12 * nLC * nVar)

See also **FORCESLCS_TRUSS**, **ELEMFORCES**.

gaussq

GAUSSQ Gauss points for 2D numerical integration.

`[x,H] = gaussq(n)` returns the coordinates and weights for a 2D gauss-legendre quadrature.

<code>n</code>	number of points in one direction = 2 or 3	
<code>x</code>	coordinates of gauss-points	$(n^2 * 2)$
<code>H</code>	weights used in summation	$(1 * n^2)$

See also KE_SHELL8, KELCS_SHELL4

gaussqtet

GAUSSQTET Gauss points for 3D numerical integration on a tetrahedron.

`[x,H] = gaussqtet(n)` returns the coordinates and weights for a 3D gauss-legendre quadrature on a tetrahedron. The coordinates are returned in natural coordinates (i.e. no volume coordinates)

<code>n</code>	number of integration points	
<code>x</code>	coordinates of the integration points	(<code>n * 3</code>)
<code>H</code>	weights used in summation	(<code>1 * n</code>)

See also GAUSSQ, GAUSSQTRI

gaussqtri

GAUSSQTRI Gauss points for 2D numerical integration on a triangle.

`[x,H] = gaussqtri(n)` returns the coordinates and weights for a 2D triangular gauss-legendre quadrature.

<code>n</code>	number of integration points	
<code>x</code>	coordinates of the integration points	<code>(n * 2)</code>
<code>H</code>	weights used in summation	<code>(1 * n)</code>

See also **GAUSSQ**

getdof

GETDOF Get the vector with the degrees of freedom of the model.

DOF=getdof(Elements,Types) builds the vector with the labels of the degrees of freedom for which stiffness is present in the finite element model.

Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Types	Element type definitions	{TypID EltName Option1 ... }
DOF	Degrees of freedom	(nDOF * 1)

See also DOF_TRUSS, DOF_BEAM, GETDOF.

getmovie

GETMOVIE Get the movie from a figure where an animation has been played.

```
mov=getmovie(h)  
gets the movie from the userdata of the axis of a figure where an animation  
has been played. In order to save the movie in the userdata the animation  
should have been played using animdisp(...,'CreateMovie','on'). This  
function blocks the command prompt until the movie has become available.
```

h Axis handle.
mov Structured array with movie frames.

The movie can be played with `movie(gcf,mov)`.

See also ANIMDISP, MOVIE.

grid_plane4

GRID_PLANE4 Grid in natural coordinates for mapped meshing.

```
[s,t,NodeNum,Elements] = grid_plane4(m,n,Type,Section,Material)
```

returns matrices of a grid in the natural coordinate system, which can be used for mapped meshing.

s	s-coordinate of nodes (1 * nNodes)
t	t-coordinate of nodes (1 * nNodes)
NodeNum	Node numbers order on grid ((m+1) * (n+1))
Elements	Node numbers are saved per element here (nElem * 8)
Type	Type ID of meshed elements
Section	Section ID of meshed elements
Material	Material ID of meshed elements

See also MAKEMESH, GRID_SHELL8.

grid_shell4

GRID_SHELL4 Grid in natural coordinates for mapped meshing.

```
[s,t,NodeNum,Elements] = grid_shell4(m,n,Type,Section,Material)
```

returns matrices of a grid in the natural coordinate system, which can be used for mapped meshing.

s	s-coordinate of nodes (1 * nNodes)
t	t-coordinate of nodes (1 * nNodes)
NodeNum	Node numbers order on grid ((m+1) * (n+1))
Elements	Node numbers are saved per element here (nElem * 8)
Type	Type ID of meshed elements
Section	Section ID of meshed elements
Material	Material ID of meshed elements

See also MAKEMESH, GRID_SHELL8.

grid_shell8

GRID_SHELL8 Grid in natural coordinates for mapped meshing.

```
[s,t,NodeNum,Elements] = grid_shell8(m,n,Type,Section,Material)
```

returns matrices of a grid in the natural coordinate system, which can be used for mapped meshing.

s	s-coordinate of nodes (1 * nNodes)
t	t-coordinate of nodes (1 * nNodes)
NodeNum	Node numbers order on grid ((m+1) * (n+1))
Elements	Node numbers are saved per element here (nElem * 12)
Type	Type ID of meshed elements
Section	Section ID of meshed elements
Material	Material ID of meshed elements

See also MAKEMESH, GRID_SHELL4.

kelcs_beam

KELCS_BEAM Beam element stiffness and mass matrix in local coordinate system.

```
[KeLCS,MeLCS] = KELCS_BEAM(L,A,ky,kz,Ixx,Iyy,Izz,E,nu,rho,Options)
[KeLCS,MeLCS] = KELCS_BEAM(L,A,ky,kz,Ixx,Iyy,Izz,E,nu,rho)
KeLCS          = KELCS_BEAM(L,A,ky,kz,Ixx,Iyy,Izz,E,nu)
returns the element stiffness and mass matrix in the local coordinate system
for a two node beam element (isotropic material).
```

```
[KeLCS,~,dKeLCSdx] = KELCS_BEAM(L,A,ky,kz,Ixx,Iyy,Izz,E,nu,rho,Options,dLdx,dAdx,...
                                dkydx,dkzdx,dIxxdx,dIyydx,dIzzdx)
```

returns the element stiffness matrix in the local coordinate system for a two node beam element (isotropic material), and additionally computes the derivatives of the stiffness matrix with respect to the design variables x. The derivatives of the mass matrix have not yet been implemented.

L	Beam length	
A	Beam cross section area	
ky	Shear deflection factor $A_{sy} = ky * A$	
kz	Shear deflection factor $A_{sz} = kz * A$	
Ixx	Moment of inertia	
Iyy	Moment of inertia	
Izz	Moment of inertia	
E	Young's modulus	
nu	Poisson coefficient	
rho	Mass density	
Options	Options for the mass matrix: {'lumped'}, {'norotatoryinertia'}	
dLdx	Beam length derivatives	(1 * nVar)
dAdx	Beam cross section area derivatives	(1 * nVar)
dkydx	Shear deflection factor derivatives	(1 * nVar)
dkzdx	Shear deflection factor derivatives	(1 * nVar)
dIxxdx	Moment of inertia derivatives	(1 * nVar)
dIyydx	Moment of inertia derivatives	(1 * nVar)
dIzzdx	Moment of inertia derivatives	(1 * nVar)
KeLCS	Element stiffness matrix	(12 * 12)
MeLCS	Element mass matrix	(12 * 12)
dKeLCSdx	Element stiffness matrix derivatives	(CELL(nVar,1))

See also KE_BEAM, KELCS_TRUSS.

kelcs_shell2

KELCS_SHELL2 SHELL2 element stiffness matrix in local coordinate system.

KeLCS = kelcs_shell2(r1,phi,L,h,E,nu) returns the element stiffness matrix in the local coordinate system for a two-node axisymmetric shell element. The global y-axis is assumed to be the axis of symmetry.

r1	Radial coordinate of node 1
phi	Slope with respect to the xz-plane
h	Shell thickness
E	Young's modulus
nu	Poisson coefficient
KeLCS	Element stiffness matrix (6 * 6)

kelcs_shell4

KELCS_SHELL4 shell element stiffness and mass matrix in element coordinate system.

```
[Ke,Me] = kelcs_shell4(Node_lc,h,E,nu,rho)
```

```
Ke       = kelcs_shell4(Node_lc,h,E,nu)
```

returns the element stiffness and mass matrix in the element coordinate system for a four node shell element (isotropic material).

Node Node definitions [x y z] (4 * 3)

Nodes should have the following order:

```
4-----3
```

```
|           |  
|           |  
|           |
```

```
1-----2
```

h Shell thickness

E Young's modulus

nu Poisson coefficient

rho Mass density

Options Element options [NOT available yet] {Option1 Option2 ...}

KeLCS Element stiffness matrix (24 * 24)

MeLCS Element mass matrix (24 * 24)

This element is a flat shell element that consists of a bilinear membrane element and four overlaid DKT triangles for the bending stiffness.

See also KE_SHELL8, ASMKM, KE_DKT.

kelcs_truss

KELCS_TRUSS Truss element stiffness and mass matrix in local coordinate system.

```
[KeLCS,MeLCS] = KELCS_TRUSS(L,A,E,rho,Options)
```

```
[KeLCS,MeLCS] = KELCS_TRUSS(L,A,E,rho)
```

```
KeLCS = KELCS_TRUSS(L,A,E)
```

returns the element stiffness and mass matrix in the local coordinate system for a two node truss element (isotropic material).

```
[KeLCS,~,dKeLCSdx] = KELCS_TRUSS(L,A,E,rho,Options,dLdx,dAdx)
```

returns the element stiffness matrix in the local coordinate system for a two node truss element (isotropic material), and additionally computes the derivatives of the stiffness matrix with respect to the design variables x . The derivatives of the mass matrix have not yet been implemented.

L Truss length

A Truss cross section area

E Young's modulus

rho Mass density

Options Options for the mass matrix: {'lumped'}

dLdx Truss length derivatives (1 * nVar)

dAdx Truss cross section area derivatives (1 * nVar)

KeLCS Element stiffness matrix (6 * 6)

MeLCS Element mass matrix (6 * 6)

dKeLCSdx Element stiffness matrix derivatives (CELL(nVar,1))

See also KE_TRUSS, KELCS_BEAM.

ke_beam

KE_BEAM Beam element stiffness and mass matrix in global coordinate system.

`[Ke,Me] = KE_BEAM(Node,Section,Material,Options)` returns the element stiffness and mass matrix in the global coordinate system for a two node beam element (isotropic material).

`[Ke,~,dKedx] = KE_BEAM(Node,Section,Material,Options,dNodedx,dSectiondx)` returns the element stiffness matrix in the global coordinate system for a two node beam element (isotropic material), and additionally computes the derivatives of the stiffness matrix with respect to the design variables `x`. The derivatives of the mass matrix have not yet been implemented.

<code>Node</code>	Node definitions	<code>[x y z] (3 * 3)</code>
<code>Section</code>	Section definition	<code>[A ky kz Ixx Iyy Izz]</code>
<code>Material</code>	Material definition	<code>[E nu rho]</code>
<code>Options</code>	Struct containing element options. Fields:	
	<code>.lumped</code>	Construct lumped mass matrix: {true false (default)}
	<code>.rotationalInertia</code>	Include rotational inertia: {true (default) false}
<code>dNodedx</code>	Node definitions derivatives	<code>(SIZE(Node) * nVar)</code>
<code>dSectiondx</code>	Section definitions derivatives	<code>(SIZE(Section) * nVar)</code>
<code>Ke</code>	Element stiffness matrix	<code>(12 * 12)</code>
<code>Me</code>	Element mass matrix	<code>(12 * 12)</code>
<code>dKedx</code>	Element stiffness matrix derivatives	<code>(CELL(nVar,1))</code>

See also `KELCS_BEAM`, `TRANS_BEAM`, `ASMKM`, `KE_TRUSS`.

ke_dkt

KE_DKT DKT plate element stiffness and mass matrix.

```
[Ke,Me] = ke_dkt(Node,h,E,nu,rho)
Ke       = ke_dkt(Node,h,E,nu)
```

returns the element stiffness and mass matrix in the global coordinate system for a three node plate element (isotropic material) in the xy-plane.

Node	Node definitions	[x y] (3 * 2)
h	Plate thickness (uniform or defined in nodes)	[h]/[h1 h2 h3]
E	Young's modulus	
nu	Poisson coefficient	
rho	Mass density	
Ke	Element stiffness matrix (24 * 24)	
Me	Element mass matrix (24 * 24)	

This element is a Discrete Kirchhoff Triangle plate element. This code is a MATLAB code based on the E-2 element FORTRAN coding, described in: Construction of new efficient three-node triangular thin plate bending elements, C. Jeyachandrabose and J. Kirkhope, Computers & Structures Vol. 23, No. 5, pp. 587-603, 1986.

See also Q_DKT, SH_T, KE_DKT4.

ke_mass

KE_MASS mass element system matrices in the global coordinate system.

[Ke,Me] = ke_mass(Node,Section,Material,Options) returns the element stiffness and mass matrix in the global coordinate system for a concentrated mass element

Node	Node definitions	[x1 y1 z1] (1 * 3)
Section	Section definition	[m]
Material	Material definition	[]
Options	Element options	{Option1 Option2 ...}
Ke	Element stiffness matrix	(6 * 6)
Me	Element mass matrix	(6 * 6)

ke_plane10

KE_PLANE10 plane element stiffness and mass matrix in global coordinate system.

`[Ke,Me] = ke_plane10(Node,Section,Material,Options)` returns the element stiffness and mass matrix in the global coordinate system for a 10-node plane triangular element. Plane10 only operates in the 2D xy-plane so that z-coordinates should be equal to zero.

Node Node definitions [x y z] (10 * 3)
Nodes should have the following order:

```

  3
  | \
 8  7
  |  \
 9 10 6
  |    \
 1--4--5--2

```

Section Section definition [h] (only used in plane stress)

Material Material definition [E nu rho]

Options Struct containing optional parameters. Fields:

.problem Plane stress or plane strain

{'2dstress' (default) | '2dstrain'}

.nXi Number of Gauss integration points

{1 | 3 | 4 | 6 | 7 (default) | 12 | 13 | 16 | 19 | 28 | 33 | 37}

Ke Element stiffness matrix (20 * 20)

Me Element mass matrix (20 * 20)

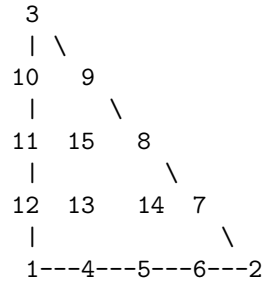
See also KE_BEAM, ASMKM, KE_TRUSS.

ke_plane15

KE_PLANE15 plane element stiffness and mass matrix in global coordinate system.

`[Ke,Me] = ke_plane15(Node,Section,Material,Options)` returns the element stiffness and mass matrix in the global coordinate system for a 15-node plane triangular element. Plane15 only operates in the 2D xy-plane so that z-coordinates should be equal to zero.

Node Node definitions [x y z] (15 * 3)
Nodes should have the following order:



Section Section definition [h] (only used in plane stress)

Material Material definition [E nu rho]

Options Struct containing optional parameters. Fields:

.problem Plane stress or plane strain

{'2dstress' (default) | '2dstrain'}

.nXi Number of Gauss integration points

{1 | 3 | 4 | 6 | 7 | 12 | 13 (default) | 16 | 19 | 28 | 33 | 37}

Ke Element stiffness matrix (30 * 30)

Me Element mass matrix (30 * 30)

See also KE_BEAM, ASMKM, KE_TRUSS.

ke_plane3

KE_PLANE3 plane element stiffness and mass matrix in global coordinate system.

[Ke,Me] = ke_plane3(Node,Section,Material,Options) returns the element stiffness and mass matrix in the global coordinate system for a 3-node CST element. Plane3 only operates in the 2D xy-plane so that z-coordinates should be equal to zero.

Node	Node definitions	[x y z] (3 * 3)
Section	Section definitions	[h] (only used in plane stress)
Material	Material definition	[E nu rho]
Options	Struct containing optional parameters. Fields:	
	.problem Plane stress or plane strain	
	{'2dstress' (default) '2dstrain'}	
Ke	Element stiffness matrix (6 * 6)	
Me	Element mass matrix (6 * 6)	

ke_plane4

KE_PLANE4 plane element stiffness and mass matrix in global coordinate system.

`[Ke,Me] = ke_plane4(Node,Section,Material,Options)` returns the element stiffness and mass matrix in the global coordinate system for a 4-node plane element. Plane4 only operates in the 2D xy-plane so that z-coordinates should be equal to zero.

Node Node definitions [x y z] (4 * 3)
Nodes should have the following order:

```

      4---3
      |   |
      1---2

```

Section Section definition [h] (only used in plane stress)

Material Material definition [E nu rho]

Options Struct containing optional parameters. Fields:

 .problem Plane stress, plane strain or axisymmetrical
 {'2dstress' (default) | '2dstrain' | 'axisym'}

 .bendingmodes Include (non-conforming) bending modes
 {true (default) | false}

 .integration Full (2x2) or reduced (1x1) integration
 {'full' (default) | 'reduced'}

Ke Element stiffness matrix (8 * 8)

Me Element mass matrix (8 * 8)

See also KE_BEAM, ASMKM, KE_TRUSS.

ke_plane6

KE_PLANE6 plane element stiffness and mass matrix in global coordinate system.

`[Ke,Me] = ke_plane6(Node,Section,Material,Options)` returns the element stiffness and mass matrix in the global coordinate system for a 6-node plane triangular element. Plane6 only operates in the 2D xy-plane so that z-coordinates should be equal to zero.

Node Node definitions [x y z] (6 * 3)
Nodes should have the following order:

```

3
| \
6 5
|  \
1--4--2

```

Section	Section definition	[h] (only used in plane stress)
Material	Material definition	[E nu rho]
Options	Struct containing optional parameters. Fields:	
	.problem	Plane stress or plane strain {'2dstress' (default) '2dstrain'}
	.nXi	Number of Gauss integration points {1 3 4 6 7 (default) 12 13 16 19 28 33 37}
Ke	Element stiffness matrix (12 * 12)	
Me	Element mass matrix (12 * 12)	

See also KE_BEAM, ASMKM, KE_TRUSS.

ke_plane8

KE_PLANE8 plane element stiffness and mass matrix in global coordinate system.

`[Ke,Me] = ke_plane8(Node,Section,Material,Options)` returns the element stiffness and mass matrix in the global coordinate system for a 8-node plane element. Plane8 only operates in the 2D xy-plane so that z-coordinates should be equal to zero.

Node	Node definitions	[x y z] (8 * 3)
Section	Section definition	[h] (only used in plane stress)
Material	Material definition	[E nu rho]
Options	Struct containing optional parameters. Fields:	
	.problem Plane stress or plane strain	
	{'2dstress' (default) '2dstrain' 'axisym'}	
Ke	Element stiffness matrix (16 * 16)	
Me	Element mass matrix (16 * 16)	

ke_shell2

KE_SHELL2 SHELL2 element stiffness matrix in global coordinate system.

Ke = ke_shell2(Node,Section,Material) returns the element stiffness matrix in the global coordinate system for a two-node axisymmetric shell element. The global y-axis is assumed to be the axis of symmetry.

Node	Node definitions	[x y z] (3 * 3)
Section	Section definition	[h]
Material	Material definition	[E nu]
Ke	Element stiffness matrix	(6 * 6)

ke_shell4

KE_SHELL4 shell element stiffness and mass matrix in global coordinate system.

```
[Ke,Me] = ke_shell4(Node,Section,Material,Options)
Ke       = ke_shell4(Node,Section,Material,Options)
returns the element stiffness and mass matrix in the global coordinate system
for a four node shell element (isotropic material).
```

Node Node definitions [x y z] (4 * 3)
 Nodes should have the following order:
 4-----3
 | |
 | |
 | |
 1-----2

Section	Section definition	[h]
Material	Material definition	[E nu rho]
Options	Element options	{Option1 Option2 ...}
Ke	Element stiffness matrix	(24 * 24)
Me	Element mass matrix	(24 * 24)

This shell element consists of a bilinear membrane element and four overlaid DKT triangles for the bending stiffness.

See also KE_BEAM, ASMKM, KE_TRUSS.

ke_shell6

KE_SHELL6 shell element stiffness and mass matrix in global coordinate system.

```
[Ke,Me] = ke_shell6(Node,Section,Material,Options)
Ke       = ke_shell6(Node,Section,Material,Options)
returns the element stiffness and mass matrix in the global coordinate system
for an eight node shell element.
```

Node	Node definitions [x y z] (6 * 3) Nodes should have the following order: 3 \ 6 5 \ 1--4--2
Section	Section definition [h] or [h1 h2 h3] (uniform thickness or defined in corner nodes(1,2,3))
Material	Material definition [E nu rho] or [Exx Eyy nuxy muxy muyz muzx theta rho]
Options	Element options struct. Fields: -LCSType: determine the reference local element coordinate system. Values: 'element' (default) or 'global' -MatType: 'isotropic' (default) or 'orthotropic' -Offset: nodal offset from shell midplane. Values: 'top', 'mid' (default), 'bot' or numerical value
Ke	Element stiffness matrix (36 * 36)
Me	Element mass matrix (36 * 36)

This element is based on chapter 15 of
The Finite Element Method: for Solid and Structural Mechanics,
Zienkiewicz (2005).

See also KE_BEAM, ASMKM, KE_TRUSS.

ke_shell8

KE_SHELL8 Shell element stiffness and mass matrix in global coordinate system.

```
[Ke,Me] = ke_shell8(Node,Section,Material,Options)
```

```
Ke = ke_shell8(Node,Section,Material,Options)
```

returns the element stiffness and mass matrix in the global coordinate system for an eight node shell element.

Node Node definitions [x y z] (8 * 3)

Nodes should have the following order:

```
4-----7-----3
```

```
|           |
```

```
8           6
```

```
|           |
```

```
1-----5-----2
```

Section Section definition [h] or [h1 h2 h3 h4]

(uniform thickness or defined in corner nodes(1,2,3,4))

Material Material definition [E nu rho] or

[Exx Eyy nuxy muxy muyz muzx theta rho]

Options Element options struct. Fields:

-LCSType: determine the reference local element coordinate system. Values:

'element' (default) or 'global'

-MatType: 'isotropic' (default) or 'orthotropic'

-Offset: nodal offset from shell midplane. Values:

'top', 'mid' (default), 'bot' or numerical value

-RotaryInertia: 0 (default) or 1

Ke Element stiffness matrix (48 * 48)

Me Element mass matrix (48 * 48)

This element is based on chapter 15 of

The Finite Element Method: for Solid and Structural Mechanics,
Zienkiewicz (2005).

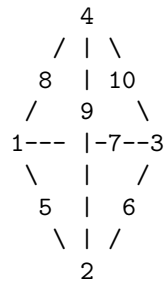
See also KE_BEAM, ASMKM, KE_TRUSS.

ke_solid10

KE_SOLID10 Compute the element stiffness and mass matrix for a solid10 element.

`[Ke,Me] = ke_solid10(Node,Section,Material,Options)` computes element stiffness and mass matrix in the global coordinate system for a solid10 element.

Node Node definitions [x y z] (10 * 3)
Nodes should have the following order:



Section Section definition []
Material Material definition [E nu rho]
Options Struct containing optional parameters. Fields:
 .nXi Number of Gauss integration points
 {1 | 4 (default) | 5}
Ke Element stiffness matrix (30 * 30)
Me Element mass matrix (30 * 30)

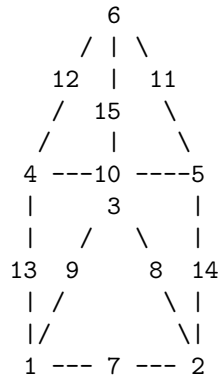
See also KE_BEAM, ASMKM, KE_TRUSS.

ke_solid15

KE_SOLID15 Compute the element stiffness and mass matrix for a solid15 element.

`[Ke,Me] = ke_solid15(Node,Section,Material,Options)` computes element stiffness and mass matrix in the global coordinate system for a solid15 prismatic element.

Node Node definitions [x y z] (15 * 3)
Nodes should have the following order:



Section Section definition []
Material Material definition [E nu rho]
Options Element options struct. Fields: []
Ke Element stiffness matrix (45 * 45)
Me Element mass matrix (45 * 45)

ke_solid20

KE_SOLID20 Compute the element stiffness and mass matrix for a solid20 element.

`[Ke,Me] = ke_solid20(Node,Section,Material,Options)` computes element stiffness and mass matrix in the global coordinate system for a solid20 element.

Node Node definitions [x y z] (20 * 3)

Nodes should have the following order:

```

      8---15-----7
      /|          /|
    16 |          14 |
      / 20        / 19
      / |          / |
    5---13-----6   |
    |   4--11|----3
    | /       | /
  17 12      18 10
    | /       | /
    | /       | /
    1-----9-----2

```

Section Section definition []

Material Material definition [E nu rho]

Options Element options struct. Fields: []

Ke Element stiffness matrix (60 * 60)

Me Element mass matrix (60 * 60)

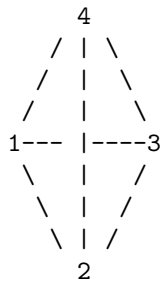
See also **KE_BEAM**, **ASMKM**, **KE_TRUSS**.

ke_solid4

KE_SOLID4 Compute the element stiffness and mass matrix for a solid4 element.

`[Ke,Me] = ke_solid4(Node,Section,Material,Options)` computes element stiffness and mass matrix in the global coordinate system for a solid4 element.

Node Node definitions `[x y z]` (4 * 3)
Nodes should have the following order:



Section Section definition `[]`
Material Material definition `[E nu rho]`
Options Element options struct. Fields: `[]`
Ke Element stiffness matrix (12 * 12)
Me Element mass matrix (12 * 12)

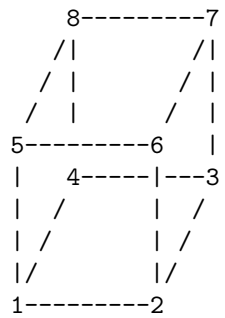
See also **KE_BEAM**, **ASMKM**, **KE_TRUSS**.

ke_solid8

KE_SOLID8 Compute the element stiffness and mass matrix for a solid8 element.

```
[Ke,Me] = ke_solid8(Node,Section,Material,Options)
[Ke,Me] = ke_solid8(Node,Section,Material,Options)
computes element stiffness and mass matrix in the global coordinate system
                                                    for a solid8 element.
```

Node Node definitions [x y z] (8 * 3)
Nodes should have the following order:



Section Section definition []
Material Material definition [E nu rho]
Options Element options struct. Fields: []
Ke Element stiffness matrix (24 * 24)
Me Element mass matrix (24 * 24)

See also KE_BEAM, ASMKM, KE_TRUSS.

ke_truss

KE_TRUSS Truss element stiffness and mass matrix in global coordinate system.

`[Ke,Me] = KE_TRUSS(Node,Section,Material,Options)` returns the element stiffness and mass matrix in the global coordinate system for a two node truss element (isotropic material).

`[Ke,~,dKedx] = KE_TRUSS(Node,Section,Material,Options,dNodedx,dSectiondx)` returns the element stiffness matrix in the global coordinate system for a two node truss element (isotropic material), and additionally computes the derivatives of the stiffness matrix with respect to the design variables `x`. The derivatives of the mass matrix have not yet been implemented.

Node	Node definitions	[x1 y1 z1; x2 y2 z2] (2 * 3)
Section	Section definition	[A]
Material	Material definition	[E nu rho]
Options	Struct containing element options. Fields:	
	.lumped Construct lumped mass matrix:	
	{true false (default)}	
dNodedx	Node definitions derivatives	(SIZE(Node) * nVar)
dSectiondx	Section definitions derivatives	(SIZE(Section) * nVar)
Ke	Element stiffness matrix	(6 * 6)
Me	Element mass matrix	(6 * 6)
dKedx	Element stiffness matrix	(CELL(nVar,1))

See also `KELCS_TRUSS`, `TRANS_TRUSS`, `ASMKM`, `KE_BEAM`.

lconstr

LCONSTR Add linear constraint equations to the stiffness matrix and load vector
using Lagrange multipliers.

```
[L,K,F]=lconstr(Constr,DOF,K,F)
[L,K,F,M]=lconstr(Constr,DOF,K,[],M)
[L,K,F,M]=lconstr(Constr,DOF,K,F,M)
```

modifies the stiffness matrix, the mass matrix and the load vector according to the applied constraint equations. The dimensions of the stiffness matrix, the mass matrix and the load vector increase with the number of constraints. The resulting stiffness and mass matrix are symmetric.

Constr	Constraint equation: Constant=CoefS*SlaveDOF+CoefM1*MasterDOF1+CoefM2*MasterDOF2+... [Constant CoefS SlaveDOF CoefM1 MasterDOF1 CoefM2 MasterDOF2 ...]
DOF	Degrees of freedom (nDOF * 1)
L	Selection matrix for displacements: $U = L*(K \setminus F)$; (nDOF * (nDOF+nConstr))
K	Stiffness matrix (nDOF * nDOF)
F	Load vector (nDOF * nSteps)
M	Mass matrix (nDOF * nDOF)

linkandcheck

LINKANDCHECK Link two matrices and check for presence.

```
p3=linkandcheck(p1,p2,name)
```

Link two matrices and check for presence

p1 Vector with elements to be link and checked

p2 Vector with elements to be link and checked

name Name of elements in p2

p3 Lowest absolute index in p2 for each element in p1

loadslcs_beam

`LOADSLCS_BEAM` Equivalent nodal forces for a beam element in the LCS.

```
FLCS = LOADSLCS_BEAM(DLoadLCS,L)
```

computes the equivalent nodal forces of a distributed load
(in the local coordinate system).

```
[FLCS,dFLCSdx] = LOADSLCS_BEAM(DLoadLCS,L,dDLoadLCSdx,dLdx)
```

additionally computes the derivatives of the equivalent nodal forces
with respect to the design variables `x`.

<code>DLoadLCS</code>	Distributed loads	<code>[n1localX; n1localY; n1localZ; ...]</code>	$(6/8 * nLC)$
<code>L</code>	Beam length		
<code>dDLoadLCSdx</code>	Distributed loads derivatives		$(6/8 * nLC * nVar)$
<code>dLdx</code>	Beam length derivatives		$(1 * nVar)$
<code>FLCS</code>	Load vector		$(12 * nLC)$
<code>dFLCSdx</code>	Load vector derivatives		$(12 * nLC * nVar)$

See also `LOADS_BEAM`.

loadslcs_shell2

LOADSLCS_SHELL2 Equivalent nodal forces for a SHELL2 element in the LCS.

```
FLCS = loadslcs_shell2(DLoadLCS,L)
computes the equivalent nodal forces of a distributed load
(in the local coordinate system).
```

DLoadLCS	Distributed loads	[n1localX; n1localY; n1localZ; ...]
		(6 * 1)

L	Element length
---	----------------

FLCS	Load vector (12 * 1)
------	----------------------

loadslcs_shell4

LOADSLCS_SHELL4 Equivalent nodal forces for a shell4 element in the LCS.

```
F = loadslcs_shell4(DLoadLCS,Node)
computes the equivalent nodal forces of a distributed load
(in the local coordinate system).
```

DLoadLCS	Distributed loads	[n1localX; n1localY; n1localZ; ...]
	in corner Nodes	(12 * 1)
Node	Node definitions	[x y z] (4 * 3)
FLCS	Load vector	(24 * 1)

See also LOADSLCS_BEAM, ELEMLOADS, LOADS_TRUSS.

loads_beam

LOADS_BEAM Equivalent nodal forces for a beam element in the GCS.

```
F = LOADS_BEAM(DLoad,Node)
```

computes the equivalent nodal forces of a distributed load
(in the global coordinate system).

```
[F,dFdx] = LOADS_BEAM(DLoad,Node,dDLoaddx,dNodedx)
```

additionally computes the derivatives of the equivalent nodal forces
with respect to the design variables x.

DLoad	Distributed loads	[n1globalX; n1globalY; n1globalZ; ...] (6/8 * nLC * nDLoads)
Node	Node definitions	[x y z] (3 * 3)
dDLoaddx	Distributed loads derivatives	(6/8 * nLC * nVar * nDLoads)
dNodedx	Node definitions derivatives	(SIZE(Node) * nVar)
F	Load vector	(12 * nLC)
dFdx	Load vector derivatives	(12 * nLC * nVar)

See also **LOADSLCS_BEAM**, **ELEMLoads**, **LOADS_TRUSS**.

loads_shell2

LOADS_SHELL2 Equivalent nodal forces for a SHELL2 element in the GCS.

F = loads_shell2(DLoad,Node)
computes the equivalent nodal forces of a distributed load
(in the global coordinate system).

DLoad	Distributed loads	[n1globalX; n1globalY; n1globalZ; ...] (6 * 1)
Node	Node definitions	[x y z] (3 * 3)
F	Load vector	(6 * 1)

loads_shell4

`LOADS_SHELL4` Equivalent nodal forces for a shell4 element in the GCS.

```
F = loads_shell4(DLoad,Node)
computes the equivalent nodal forces of a distributed load
(in the global coordinate system).
```

<code>DLoad</code>	Distributed loads in corner Nodes	<code>[n1globalX; n1globalY; n1globalZ; ...]</code> <code>(12 * 1)</code>
<code>Node</code>	Node definitions	<code>[x y z] (4 * 3)</code>
<code>F</code>	Load vector	<code>(24 * 1)</code>

See also `LOADSLCS_BEAM`, `ELEMLOADS`, `LOADS_TRUSS`.

loads_shell6

LOADS_SHELL6 Equivalent nodal forces for a shell6 element in the GCS.

```
F = loads_shell6(DLoad,Node)
computes the equivalent nodal forces of a distributed load
(in the global coordinate system).
```

DLoad	Distributed loads	[n1globalX; n1globalY; n1globalZ; ...]
	in corner Nodes	(9 * 1)
Node	Node definitions	[x y z] (6 * 3)
F	Load vector	(36 * 1)

See also **LOADSLCS_BEAM**, **ELEMLOADS**, **LOADS_TRUSS**.

loads_shell8

LOADS_SHELL8 Equivalent nodal forces for a shell8 element in the GCS.

```
F = loads_shell8(DLoad,Node)
computes the equivalent nodal forces of a distributed load
(in the global coordinate system).
```

DLoad	Distributed loads	[n1globalX; n1globalY; n1globalZ; ...]
	in corner Nodes	(12 * 1)
Node	Node definitions	[x y z] (8 * 3)
F	Load vector	(48 * 1)

See also **LOADSLCS_BEAM**, **ELEMLOADS**, **LOADS_TRUSS**.

loads_solid20

LOADS_SOLID20 Equivalent nodal forces for a solid20 element.

```
F = loads_solid20(DLoad,Node)
computes the equivalent nodal forces of a distributed load
(in the global coordinate system).
```

DLoad	Distributed loads	[n1globalX; n1globalY; n1globalZ; ...]
	in corner Nodes	(24 * 1)
Node	Node definitions	[x y z] (8 * 3)
F	Load vector	(24 * 1)

See also LOADSLCS_BEAM, ELEMLOADS, LOADS_TRUSS.

loads_solid8

LOADS_SOLID8 Equivalent nodal forces for a solid8 element.

```
F = loads_solid8(DLoad,Node)
computes the equivalent nodal forces of a distributed load
(in the global coordinate system).
```

DLoad	Distributed loads	[n1globalX; n1globalY; n1globalZ; ...]
	in Nodes	(24 * 1)
Node	Node definitions	[x y z] (8 * 3)
F	Load vector	(24 * 1)

See also LOADSLCS_BEAM, ELEMLOADS, LOADS_TRUSS.

loads_truss

LOADS_TRUSS Equivalent nodal forces for a truss element in the GCS.

```
F = LOADS_TRUSS(DLoad,Node)
```

computes the equivalent nodal forces of a distributed load
(in the global coordinate system).

```
[F,dFdx] = LOADS_TRUSS(DLoad,Node,dDLoaddx,dNodedx)
```

additionally computes the derivatives of the equivalent nodal forces
with respect to the design variables x.

DLoad	Distributed load	[n1globalX n1globalY n1globalZ ...]
	(6/8 * nLC)	

Node	Node definitions	[x y z] (2 * 3)
------	------------------	-----------------

dDLoaddx	Distributed load derivatives	(6/8 * nLC * nVar)
----------	------------------------------	--------------------

dNodedx	Node definitions derivatives	(SIZE(Node) * nVar)
---------	------------------------------	---------------------

F	Load vector	(6 * nLC)
---	-------------	-----------

dFdx	Load vector derivatives	(6 * nLC * nVar)
------	-------------------------	------------------

See also **ELEMLOADS**, **LOADS_BEAM**.

Isolver

LSOLVER Solves linear system of equations.

```
[U,fK] = lsolver(K,P,save_factor,check_crpoint,fK)
solves the linear system  $K*U=P$ .
```

```
K      Coefficient matrix [n * n]
P      Right-hand side      [n * 1]
save_factor Save factorization {true | false}
check_crpoint Check critical points {true | false}
fK     Struct containing factorization of K
```

meshcat

eshcat is a function.

```
[Nodes, Elements] = meshcat(varargin)
```

msupf

MSUPF Modal superposition in the frequency domain.

`[X,H] = MSUPF(omega,xi,Omega,Pm)` calculates the modal transfer functions H and the modal displacements $x(t) = X * \text{EXP}(i * \text{Omega} * t)$ of a dynamic system with the eigenfrequencies ω and the modal damping ratios ξ due to the modal excitation $p_m(t) = P_m * \text{EXP}(i * \text{Omega} * t)$.

`omega` Eigenfrequencies [rad/s] ($n_{\text{Mode}} * 1$)
`xi` Modal damping ratios [-] ($n_{\text{Mode}} * 1$) or ($1 * 1$), constant modal damping is assumed in the ($1 * 1$) case.
`Omega` Excitation frequencies [rad/s] ($1 * N$).
`Pm` Complex amplitude of the modal excitation ($n_{\text{Mode}} * N$).
`X` Complex amplitude of the modal displacements ($n_{\text{Mode}} * N$).
`H` Modal transfer functions ($n_{\text{Mode}} * N$).

msupt

MSUPT Modal superposition in the time domain.

`x = MSUPT(omega,xi,t,pm,x1,y1,interp)` calculates the modal displacements $x(t)$ of a dynamic system with the eigenfrequencies ω and the modal damping ratios ξ due to the modal excitation $p(t)$, given the initial conditions x_1 and y_1 .

The solution of the modal differential equations is performed by means of the piecewise exact method. `interp` can be 'foh' (first order hold) or 'zoh' (zero order hold). In the first case the excitation is assumed to vary linearly within each time step while in the second case it is assumed to be constant: if $t(k) \leq t < t(k+1)$ then $p(t) = p(k)$.

<code>omega</code>	Eigenfrequencies [rad/s] ($n_{\text{Mode}} * 1$)
<code>xi</code>	Modal damping ratios [-] ($n_{\text{Mode}} * 1$) or ($1 * 1$), constant modal damping is assumed in the ($1 * 1$) case.
<code>t</code>	Time points ($1 * N$) of the sampling of p , x and t .
<code>pm</code>	Modal excitation ($n_{\text{Mode}} * N$).
<code>x1</code>	Modal displacements at time point $t(1)$ ($n_{\text{Mode}} * 1$), defaults to zero.
<code>y1</code>	Modal velocities at time point $t(1)$ ($n_{\text{Mode}} * 1$), defaults to zero.
<code>interp</code>	Interpolation scheme: 'foh' or 'zoh', defaults to 'foh'.
<code>x</code>	Modal displacements ($n_{\text{Mode}} * N$).

multdloads

MULTDLOADS Combine distributed loads.

`DLoads = MULTDLOADS(DLoads_1,DLoads_2,...,DLoads_k)`
 combines the distributed loads of multiple load cases into one 3D array.
 Each 3D-plane corresponds to a single load case. In the presence of
 partial distributed loads, every row will have identical starting
 and ending points in the 3rd dimension to allow for accurate combination
 of load cases. A distributed load with a starting and/or ending point
 value of 'NaN' will be considered as a load on the entire element
 Calculations will be more efficient compared to a partial distributed
 load with starting point equal to zero and ending point equal to the
 length of the element.

<code>DLoads_k</code>	Distributed loads	<code>[EltID n1globalX n1globalY n1globalZ ...]</code> without partial DLoads: <code>(nElem_k * 7)</code> with partial DLoads: <code>(nElem_k * 9)</code>
<code>DLoads</code>	Distributed loads	<code>[EltID n1globalX n1globalY n1globalZ ...]</code> <code>(maxnElem * 7 * k)</code> <code>(maxnElem * 9 * k)</code>

See also ELEMLOADS.

nedloadlcs_beam

NEDLOADLCS_BEAM Interpolation functions for a distributed load on a beam element.

```
NeLCS = NEDLOADLCS_BEAM(Points)
NeLCS = NEDLOADLCS_BEAM(Points,phi_y,phi_z)
determines the values of the interpolation functions for a distributed
load in the specified points (LCS). These are used to compute the
displacements that occur due to the distributed loads if all nodes are
fixed.
```

```
NeLCS = NEDLOADLCS_BEAM(Points,[],[],a,b,L)
NeLCS = NEDLOADLCS_BEAM(Points,phi_y,phi_z,a,b,L)
determines the values of the interpolation functions for a partial
distributed load in the specified points (LCS). The load starts at a
distance 'a' and ends at distance 'b' from the first node of the element:
1) from 0 to a: no DLoad,
2) from a to b: DLoad,
3) from b to L: no DLoad.
```

```
[NeLCS,dNeLCSdx] = NEDLOADLCS_BEAM(Points,[],[],a,b,L,dadx,dbdx,dLdx)
[NeLCS,dNeLCSdx] = NEDLOADLCS_BEAM(Points,phi_y,phi_z,a,b,L,dadx,dbdx,dLdx)
additionally computes the derivatives of the interpolation functions of
a partial distributed load with respect to the design variables x.
Note: the derivatives of the interpolation functions for a distributed
load on the complete element are zero.
```

Points	Points in the local coordinate system	(1 * nPoints)
phi_y	Shear deformation constant in y dir	(scalar)
phi_z	Shear deformation constant in z dir	(scalar)
a	Local starting point for the DLoad	(scalar)
b	Local ending point for the DLoad	(scalar)
L	Element length	(scalar)
dadx	Local starting point derivatives	(1 * nVar)
dbdx	Local ending point derivatives	(1 * nVar)
dLdx	Element length derivatives	(1 * nVar)
NeDLoad	Interpolated values	(nPoints * 6)
dNeDLoaddx	Interpolated values derivatives	(nPoints * 6 * nVar)

See also DISP_BEAM, NELCS_BEAM.

nelcs_beam

NELCS_BEAM Shape functions for a beam element.

```
NeLCS = nelcs_beam(Points)
NeLCS = nelcs_beam(Points,phi_y,phi_z)
determines the values of the shape functions in the specified points.
```

Points	Points in the local coordinate system (1 * nPoints).
phi_y	Shear deformation constant in the y-direction (1 * 1).
phi_z	Shear deformation constant in the z-direction (1 * 1).
NeLCS	Values (nPoints * 12).

See also DISP_BEAM.

newmark

NEWMARK Direct time integration for dynamic systems - Newmark method

`[u,v,a,t] = NEWMARK(M,C,K,dt,p,u0,v0,a0,[alpha delta])` applies the Newmark method for the calculation of the nodal displacements `u`, velocities `v` and accelerations `a` of the dynamic system with the system matrices `M`, `C` and `K` due to the excitation `p`.

`M` Mass matrix (`nDof * nDof`)

`C` Damping matrix (`nDof * nDof`)

`K` Stiffness matrix (`nDof * nDof`)

`dt` Time step of the integration scheme (`1 * 1`). Should be small enough to ensure the stability and the precision of the integration scheme.

`p` Excitation (`nDof * N`). `p(:,k)` corresponds to time point `t(k)`.

`u0` Displacements at time point `t(1)-dt` (`nDof * 1`). Defaults to zero.

`v0` Velocities at time point `t(1)-dt` (`nDof * 1`). Defaults to zero.

`a0` Accelerations at time point `t(1)-dt` (`nDof * 1`). Defaults to zero.

`u` Displacements (`nDof * N`). `u(:,k)` corresponds to time point `t(k)`.

`t` Time axis (`1 * N`), defined as `t = [0:N-1] * dt`.

nodalshellf

NODALSHELLF Compute the nodal shell forces/moments per unit length
from the element solution.

[FnLCS,FnLCS2] = nodalshellf(Nodes,Elements,Types,FeLCS) computes the
nodal forces from the element solution.

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Types	Element type definitions	{TypID EltName Option1 ... }
Sections	Section definitions	[SecID SecProp1 SecProp2 ...]
FeLCS	Element forces/moments per unit length in corner nodes IJKL (nElem * 32)	[Nx Ny Nxy Mx My Mxy Vx Vy]
FnLCS	Nodal forces/moments per unit length in corner nodes IJKL (nElem * 32)	[Nx Ny Nxy Mx My Mxy Vx Vy]
FnLCS2	Nodal forces per node (nNodes * 9) (9 = NodID + 8 fcomp.)	[NodID Nx Ny Nxy Mx My Mxy Vx Vy]

See also ELEMShellF.

nodalstress

NODALSTRESS Compute the nodal stresses from the element solution.

```
[Sn,Sn2] = nodalstress(Nodes,Elements,Types,Se)
computes the nodal stresses from the element solution.
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Types	Element type definitions	{TypID EltName Option1 ... }
Sections	Section definitions	[SecID SecProp1 SecProp2 ...]
Se	Element stresses in corner nodes IJKL and at top/mid/bot of shell (nElem * 72) 72 = 6 stresscomp. * 4 nodes * 3 locations	[sxx syy szz sxy syz sxz ...]
Sn	Nodal stress in corner nodes IJKL and at top/mid/bot of shell (nElem * 72) 72 = 6 stresscomp. * 4 nodes * 3 locations	[sxx syy szz sxy syz sxz ...]
Sn2	Nodal stresses per node at top/mid/bot of shell (nNodes * 19) (19 = NodID + 3 locations * 6 scomp.)	[NodID sxx syy szz sxy syz sxz ...]

See also ELEMSTRESS.

nodalvalues

NODALVALUES Construct a vector with the values at the selected DOF.

```
F=nodalvalues(DOF,seldof,values)
```

constructs a vector with the values at the selected DOF. This function can be used to obtain a load vector, initial displacements, velocities or accelerations.

DOF Degrees of freedom (nDOF * 1)

seldof Selected degrees of freedom [NodID.dof] (nValues * 1)

values Corresponding values [Value] (nValues * nSteps)

F Load vector (nDOF * nSteps)

See also ELEMLOADS.

patch_beam

PATCH_BEAM Patch information of the beam elements for plotting.

[pxyz,pind,pvalue] = patch_beam(Nodes,NodeNum,Values) returns matrices to plot patches of beam elements.

Nodes	Node definitions [NodID x y z].
NodeNum	Node numbers [NodID1 NodID2 NodID3] (nElem * 3).
Values	Values assigned to nodes used for coloring (nElem * 3).
pxyz	Coordinates of Nodes (3*nElem * 3).
pind	Indices of Nodes (nElem * 3).
pvalue	Values arranged per Node (3*nElem * 1).

patch_mass

PATCH_MASS Patch information of the mass elements for plotting.

[pxyz,pind,pvalue] = patch_mass(Nodes,NodeNum,Values) returns matrices to plot patches of mass elements.

Nodes	Node definitions [NodID x y z].
NodeNum	Node numbers [NodID1 NodID2 NodID3] (nElem).
Values	Values assigned to nodes used for coloring (nElem).
pxyz	Coordinates of Nodes (nElem * 3).
pind	Indices of Nodes (nElem).
pvalue	Values arranged per Node (nElem).

patch_plane3

PATCH_PLANE4 Patch information of the plane4 elements for plotting.

[pxyz,pind,pvalue] = patch_plane4(Nodes,NodeNum,Values) returns matrices to plot patches of plane4 elements.

Nodes	Node definitions	[NodID x y z]
NodeNum	Node numbers	[NodID1 NodID2 NodID3 NodID4] (nElem * 4)
Values	Values assigned to nodes used for coloring	(nElem * 4)
pxyz	Coordinates of Nodes	(4*nElem * 3)
pind	indices of Nodes	(nElem * 4)
pvalue	Values arranged per Node	(4*nElem * 1)

See also PLOTSTRESSCONTOURF, PLOTSHELLFCONTOURF.

patch_plane4

PATCH_PLANE4 Patch information of the plane4 elements for plotting.

[pxyz,pind,pvalue] = patch_plane4(Nodes,NodeNum,Values) returns matrices to plot patches of plane4 elements.

Nodes	Node definitions	[NodID x y z]
NodeNum	Node numbers	[NodID1 NodID2 NodID3 NodID4] (nElem * 4)
Values	Values assigned to nodes used for coloring	(nElem * 4)
pxyz	Coordinates of Nodes	(4*nElem * 3)
pind	Indices of Nodes	(nElem * 4)
pvalue	Values arranged per Node	(4*nElem * 1)

See also PLOTSTRESSCONTOURF, PLOTSHELLFCONTOURF.

patch_plane6

PATCH_PLANE4 Patch information of the plane4 elements for plotting.

[pxyz,pind,pvalue] = patch_plane4(Nodes,NodeNum,Values) returns matrices to plot patches of plane4 elements.

Nodes	Node definitions	[NodID x y z]
NodeNum	Node numbers	[NodID1 NodID2 NodID3 NodID4] (nElem * 4)
Values	Values assigned to nodes used for coloring	(nElem * 4)
pxyz	Coordinates of Nodes	(4*nElem * 3)
pind	indices of Nodes	(nElem * 4)
pvalue	Values arranged per Node	(4*nElem * 1)

See also PLOTSTRESSCONTOURF, PLOTSHELLFCONTOURF.

patch_plane8

PATCH_PLANE4 Patch information of the plane4 elements for plotting.

[pxyz,pind,pvalue] = patch_plane4(Nodes,NodeNum,Values) returns matrices to plot patches of plane4 elements.

Nodes	Node definitions	[NodID x y z]
NodeNum	Node numbers	[NodID1 NodID2 NodID3 NodID4] (nElem * 4)
Values	Values assigned to nodes used for coloring	(nElem * 4)
pxyz	Coordinates of Nodes	(4*nElem * 3)
pind	indices of Nodes	(nElem * 4)
pvalue	Values arranged per Node	(4*nElem * 1)

See also PLOTSTRESSCONTOURF, PLOTSHELLFCONTOURF.

patch_shell4

PATCH_SHELL4 Patch information of the shell4 elements for plotting.

[pxyz,pind,pvalue] = patch_shell4(Nodes,NodeNum,Values) returns matrices to plot patches of shell4 elements.

Nodes	Node definitions	[NodID x y z]
NodeNum	Node numbers	[NodID1 NodID2 NodID3 NodID4] (nElem * 4)
Values	Values assigned to nodes used for coloring	(nElem * 4)
pxyz	Coordinates of Nodes	(4*nElem * 3)
pind	indices of Nodes	(nElem * 4)
pvalue	Values arranged per Node	(4*nElem * 1)

See also PLOTSTRESSCONTOURF, PLOTSHELLFCONTOURF.

patch_shell6

PATCH_SHELL6 Patch information of the shell6 elements for plotting.

[pxyz,pind,pvalue] = patch_shell6(Nodes,NodeNum,Values) returns matrices to plot patches of shell6 elements.

Nodes	Node definitions	[NodID x y z]
NodeNum	Node numbers	[NodID1 NodID2 NodID3 NodID4] (nElem * 6)
Values	Values assigned to nodes used for coloring	(nElem * 6)
pxyz	Coordinates of Nodes	(6*nElem * 3)
pind	indices of Nodes	(nElem * 8)
pvalue	Values arranged per Node	(6*nElem * 1)

See also PLOTSTRESSCONTOURF, PLOTSHELLFCONTOURF.

patch_shell8

PATCH_SHELL8 Patch information of the shell8 elements for plotting.

```
[pxyz,pind,pvalue] = patch_shell8(Nodes,NodeNum,Values)
returns matrices to plot patches of shell8 elements.
```

Nodes	Node definitions	[NodID x y z]	
NodeNum	Node numbers	[NodID1 NodID2 NodID3 NodID4]	(nElem * 8)
Values	Values assigned to nodes used for coloring		(nElem * 8)
pxyz	Coordinates of Nodes		(8*nElem * 3)
pind	indices of Nodes		(nElem * 8)
pvalue	Values arranged per Node		(8*nElem * 1)

See also PLOTSTRESSCONTOURF, PLOTSHELLFCONTOURF.

patch_solid10

PATCH_SOLID20 Patch information of the solid10 elements for plotting.

```
[pxyz,pind,pvalue] = patch_solid20(Nodes,NodeNum,Values)
returns matrices to plot patches of solid10 elements.
```

Nodes	Node definitions	[NodID x y z]
NodeNum	Node numbers	[NodID1 NodID2 NodID3 NodID4] (nElem * 10)
Values	Values assigned to nodes used for coloring	(nElem * 4)
pxyz	Coordinates of Nodes	(4*nElem * 3)
pind	indices of Nodes	(nElem * 4)
pvalue	Values arranged per Node	(4*nElem * 1)

See also PATCH_SHELL20.

patch_solid20

PATCH_SOLID20 Patch information of the solid20 elements for plotting.

```
[pxyz,pind,pvalue] = patch_solid20(Nodes,NodeNum,Values)
returns matrices to plot patches of solid20 elements.
```

Nodes	Node definitions	[NodID x y z]
NodeNum	Node numbers	[NodID1 NodID2 NodID3 NodID4] (nElem * 20)
Values	Values assigned to nodes used for coloring	(nElem * 8)
pxyz	Coordinates of Nodes	(8*nElem * 3)
pind	indices of Nodes	(nElem * 8)
pvalue	Values arranged per Node	(8*nElem * 1)

See also PATCH_SHELL8.

patch_solid4

PATCH_SOLID4 Patch information of the solid4 elements for plotting.

```
[pxyz,pind,pvalue] = patch_solid20(Nodes,NodeNum,Values)
returns matrices to plot patches of solid4 elements.
```

Nodes	Node definitions	[NodID x y z]
NodeNum	Node numbers	[NodID1 NodID2 NodID3 NodID4] (nElem * 10)
Values	Values assigned to nodes used for coloring	(nElem * 4)
pxyz	Coordinates of Nodes	(4*nElem * 3)
pind	indices of Nodes	(nElem * 4)
pvalue	Values arranged per Node	(4*nElem * 1)

See also PATCH_SHELL20.

patch_solid8

PATCH_SOLID8 Patch information of the solid8 elements for plotting.

```
[pxyz,pind,pvalue] = patch_solid8(Nodes,NodeNum,Values)  
returns matrices to plot patches of solid8 elements.
```

Nodes	Node definitions	[NodID x y z]
NodeNum	Node numbers	[NodID1 NodID2 NodID3 NodID4] (nElem * 8)
Values	Values assigned to nodes used for coloring	(nElem * 8)
pxyz	Coordinates of Nodes	(8*nElem * 3)
pind	indices of Nodes	(nElem * 8)
pvalue	Values arranged per Node	(8*nElem * 1)

See also PATCH_SHELL8.

patch_truss

PATCH_TRUSS Patch information of the truss elements for plotting.

[pxyz,pind,pvalue] = patch_truss(Nodes,NodeNum,Values) returns matrices to plot patches of truss elements.

Nodes	Node definitions [NodID x y z].
NodeNum	Node numbers [NodID1 NodID2 NodID3] (nElem * 2).
Values	Values assigned to nodes used for coloring (nElem * 2).
pxyz	Coordinates of Nodes (2*nElem * 3).
pind	Indices of Nodes (nElem * 2).
pvalue	Values arranged per Node (2*nElem * 1).

plotdisp

PLOTDISP Plot the displacements.

```
plotdisp(Nodes,Elements,Types,DOF,U,DLoads,Sections,Materials)
plotdisp(Nodes,Elements,Types,DOF,U,[],Sections,Materials)
plotdisp(Nodes,Elements,Types,DOF,U)
DispScal=plotdisp(Nodes,Elements,Types,DOF,U,DLoads,Sections,Materials)
plots the displacements. If DLoads, Sections and Materials are supplied, the
displacements that occur due to the distributed loads if all nodes are fixed,
are superimposed.
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Types	Element type definitions	{TypID EltName Option1 ... }
DOF	Degrees of freedom	(nDOF * 1)
U	Displacements	(nDOF * 1)
DLoads	Distributed loads	[EltID n1globalX n1globalY n1globalZ ...]
	(use empty array [] when shear deformation (in beam element)	
	is considered but no DLoads are present)	
Sections	Section definitions	[SecID SecProp1 SecProp2 ...]
Materials	Material definitions	[MatID MatProp1 MatProp2 ...]
DispScal	Displacement scaling	

plotdisp(...,ParamName,ParamValue) sets the value of the specified parameters. The following parameters can be specified:

'DispScal'	Displacement scaling. Default: 'auto'.
'DispMax'	Mention maximal displacement. Default: 'on'.
'Undeformed'	Plots the undeformed mesh. Default: 'k:'.
'Handle'	Plots in the axis with this handle. Default: current axis.

Additional parameters are redirected to the PLOT3 function which plots the deformations.

[DispScal,h] = plotdisp(...) returns a struct h with handles to all the objects in the plot.

See also DISP_TRUSS, DISP_BEAM, PLOTELEM.

plotdispx

PLOTDISPX Plot element quantity on deformed elements.

```
plotdispx(Nodes,Elements,Types,DOF,U,x);
[h,DispScal] = plotdispx(Nodes,Elements,Types,DOF,U,x,varargin)
plots element quantity on deformed elements.
```

Nodes Node definitions [NodID x y z]
Elements Element definitions [EltID TypID SecID MatID n1 n2 ...]
Types Element type definitions {TypID EltName Option1 ... }
DOF Degrees of freedom (nDOF * 1)
U Displacements (nDOF * 1)
x Elements quantity to plot (nElem * 1)

plotdispx(...,ParamName,ParamValue) sets the value of the specified parameters. The following parameters can be specified:

'GCS' Plot the GCS. Default: 'on'.
'minmax' Add location of min and max stress. Default: 'off'.
'colorbar' Add colorbar. Default: 'off'.
'Undeformed' Plots the undeformed mesh. {'on' | 'off' (default)}
'ncolor' Number of colors in colormap. Default: 10.
'Handle' Plots in the axis with this handle. Default: current axis.

Additional parameters are redirected to the PATCH function which plots the elements.

plotelem

PLOTELEM Plot the elements.

```
plotelem(Nodes,Elements,Types)
plots the elements.
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Types	Element type definitions	{TypID EltName Option1 ... }

`plotelem(...,ParamName,ParamValue)` sets the value of the specified parameters. The following parameters can be specified:

- 'Numbering' Plots the element numbers. Default: 'on'.
- 'GCS' Plots the global coordinate system. Default: 'on'.
- 'Handle' Plots in the axis with this handle. Default: current axis.

Additional parameters are redirected to the `PLOT3` function which plots the elements.

`h = PLOTELEM(...)` returns a struct `h` with handles to all the objects in the plot.

See also `COORD_TRUSS`, `COORD_BEAM`, `PLOTDISP`.

plotelemx

PLOTELEMX Plot element quantity on elements.

`plotelemx(Nodes,Elements,Types,x)` plots element quantity on elements.

Nodes Node definitions [NodID x y z]
Elements Element definitions [EltID TypID SecID MatID n1 n2 ...]
Types Element type definitions {TypID EltName Option1 ... }
x Elements quantity to plot (nElem * 1)
`plotelemx(...,ParamName,ParamValue)` sets the value of the specified parameters. The following parameters can be specified:
 'GCS' Plot the GCS. Default: 'on'.
 'minmax' Add location of min and max stress. Default: 'off'.
 'colorbar' Add colorbar. Default: 'off'.
 'Undeformed' Plots the undeformed mesh. Default: 'k-'.
 'ncolor' Number of colors in colormap. Default: 10.
 'Handle' Plots in the axis with this handle. Default: current axis.
 Additional parameters are redirected to the PATCH function which plots the elements.

See also ELEMSTRESS.

plotforc

PLOTFORC Plot element member forces.

```
plotforc(ftype,Nodes,Elements,Types,Sections,Materials,Forces,DLoads)
plotforc(ftype,Nodes,Elements,Types,Sections,Materials,Forces)
plots the element member forces (in accordance to the beam convention).
```

```
ftype      'norm'      Normal force (in the local x-direction)
(for truss 'sheary'    Shear force in the local y-direction
and beam  'shearz'    Shear force in the local z-direction
elements) 'momx'      Torsional moment (around the local x-direction)
          'momx'      Bending moment around the local y-direction
          'momz'      Bending moment around the local z-direction
ftype      'Nphi'      Normal force (per unit length) in meridional direction
(for       'Qphi'      Transverse force (per unit length) in meridional direction
shell2    'Mphi'      Bending moment (per unit length) in meridional direction
elements) 'Ntheta'     Normal force (per unit length) in circumferential direction
          'Mtheta'     Bending moment (per unit length) in circumferential direction
Nodes      Node definitions      [NodID x y z]
Elements   Element definitions   [EltID TypID SecID MatID n1 n2 ...]
Types      Element type definitions {TypID EltName Option1 ... }
Sections   Section definitions     [SecID SecProp1 SecProp2 ...]
Materials  Material definitions    [MatID MatProp1 MatProp2 ... ]
Forces      Element forces in LCS (beam convention) [N Vy Vz T My Mz]
                                                    (nElem * 12)
DLoads     Distributed loads      [EltID n1globalX n1globalY n1globalZ ...]
```

plotforc(...,ParamName,ParamValue) sets the value of the specified parameters. The following parameters can be specified:

```
'ForcScal' Force scaling. Default: 'auto'.
'MinMax'   Display minimum and maximum value. Default: 'off'.
'Values'   Force values. Default: 'on'.
'Undeformed' Plots the undeformed mesh. Default: 'k-'.
'Handle'   Plots in the axis with this handle. Default: current axis.
```

Additional parameters are redirected to the PLOT3 function which plots the forces.

[ForcScal,h] = plotforc(...) returns a struct h with handles to all the objects in the plot.

See also FDIAGRGCS_BEAM, FDIAGRGCS_TRUSS.

plotgcs

lotgcs is a function.

```
h = plotgcs(lref, h)
```

plotlcs

PLOTLCS Plot the local element coordinate systems.

```
[h,vLCS] = plotlcs(Nodes,Elements,Types)
[h,vLCS] = plotlcs(Nodes,Elements,Types,[],varargin)
plotlcs(Nodes,Elements,Types,vLCS,varargin)
```

Nodes Node definitions [NodID x y z]
 Elements Element definitions [EltID TypID SecID MatID n1 n2 ...]
 Types Element type definitions {TypID EltName Option1 ... }
 vLCS Element coordinate systems (nElem * 9)

plotlcs(...,ParamName,ParamValue) sets the value of the specified parameters. The following parameters can be specified:

'GCS' Plot the GCS. Default: 'on'.
 'Undeformed' Plots the undeformed mesh. Default: 'k-'.
 'Handle' Plots in the axis with this handle. Default: current axis.

See also ELEMSTRESS

plotnodes

PLOTNODES Plot the nodes.

```
plotnodes(Nodes)
plots the nodes.
```

Nodes Node definitions [NodID x y z]

plotnodes(...,ParamName,ParamValue) sets the value of the specified parameters. The following parameters can be specified:

- 'Numbering' Plots the node numbers. Default: 'on'.
- 'GCS' Plots the global coordinate system. Default: 'on'.
- 'Handle' Plots in the axis with this handle. Default: current axis.

Additional parameters are redirected to the PLOT3 function which plots the nodes.

h = PLOTNODES(...) returns a struct h with handles to all the objects in the plot.

See also PLOTELEM, PLOTDISP.

plotprincstress

PLOTPRINCSTRESS Plot the principal stresses in shell elements.

```
plotprincstress(Nodes,Elements,Types,Spr,Vpr)
plots the principal stresses in shell elements with a vector plot.
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Types	Element type definitions	{TypID EltName Option1 ... }
Spr	Principal stresses (nElem * 72)	[s3 s2 s1 0 0 0 ...]
Vpr	Principal dir. matrix (output from principalstress)	{nElem * 12}

plotprincdir(...,ParamName,ParamValue) sets the value of the specified parameters. The following parameters can be specified:

'location'	Location (top,mid,bot). Default: 'top'.
'GCS'	Plot the GCS. Default: 'on'.
'VectScal'	Vector scaling. Default: 'auto'.
'Undeformed'	Plots the undeformed mesh. Default: 'k-'.
'Handle'	Plots in the axis with this handle. Default: current axis.

See also PRINCIPALSTRESS.

plotshellfcontour

PLOTSHELLFCONTOUR Plot forces/moments per unit length in shell elements.

```
plotshellfcontour(ftype,Nodes,Elements,Types,F)
plots force contours (output from ELEMShellF/NODALShellF).
```

ftype	'Nx'	Membrane forces
	'Ny'	
	'Nxy'	
	'Mx'	Bending moments
	'My'	
	'Mxy'	
	'Vx'	Shear forces
	'Vy'	

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Types	Element type definitions	{TypID EltName Option1 ... }
F	Forces/moments per unit length	
	(nElem * 32)	[Nx Ny Nxy Mx My Mxy Vx Vy]

plotshellfcontour(...,ParamName,ParamValue) sets the value of the specified parameters. The following parameters can be specified:

'Ncontour'	Number of contours. Default: '10'.
'GCS'	Plot the GCS. Default: 'on'.
'Undeformed'	Plots the undeformed mesh. Default: 'k-'.
'Handle'	Plots in the axis with this handle. Default: current axis.

See also ELEMShellF, SCONTOUR_SHELL8, SCONTOUR_SHELL4.

plotshellfcontourf

PLOTSHELLFCONTOURF Plot filled contours of shell forces in shell elements.

```
plotshellfcontourf(stype,Nodes,Elements,Types,F)
plotshellfcontourf(stype,Nodes,Elements,Types,F,DOF,U)
plots force contours (output from ELEMSTRESS).
```

fctype	'Nx'	Membrane forces
	'Ny'	
	'Nxy'	
	'Mx'	Bending moments
	'My'	
	'Mxy'	
	'Vx'	Shear forces
	'Vy'	
Nodes	Node definitions [NodID x y z]	
Elements	Element definitions [EltID TypID SecID MatID n1 n2 ...]	
Types	Element type definitions {TypID EltName Option1 ... }	
F	Forces/moments per unit length (nElem * 32) [Nx Ny Nxy Mx My Mxy Vx Vy]	
DOF	Degrees of freedom (nDOF * 1)	
U	Displacements (nDOF * 1)	

plotshellfcontourf(...,ParamName,ParamValue) sets the value of the specified parameters. The following parameters can be specified:

'GCS'	Plot the GCS. Default: 'on'.
'ncolor'	Number of colors in colormap. Default: 10.
'Handle'	Plots in the axis with this handle. Default: current axis.

See also ELEMShellF, PATCH_SHELL8, PATCH_SHELL4.

plotstress

PLOTSTRESS Plot the stresses.

```
plotstress(stype,Nodes,Elements,Types,Sections,Materials,Forces,Dloads)
plotstress(stype,Nodes,Elements,Types,Sections,Materials,Forces)
plots the stresses.
```

stype	'snorm'	Normal stress due to normal force
(for truss	'smomyt'	Normal stress due to bending moment
and beam		around the local y-direction at the top
elements)	'smomyb'	Normal stress due to bending moment
		around the local y-direction at the bottom
	'smomzt'	Normal stress due to bending moment
		around the local z-direction at the top
	'smomzb'	Normal stress due to bending moment
		around the local z-direction at the bottom
	'smax'	Maximal normal stress (normal force and bending moment)
	'smin'	Minimal normal stress (normal force and bending moment)
stype	'sNphi'	Stress due to normal force in meridional direction
(for	'sMphiT'	Stress at the top due to bending moment in
		meridional direction
shell2	'sMphiB'	Stress at the bottom due to bending moment in
		meridional direction
elements)	'sNtheta'	Stress due to normal force in circumferential direction
	'sMthetaT'	Stress at the top due to bending moment in
		circumferential direction
	'sMthetaB'	Stress at the bottom due to bending moment in
		circumferential direction
Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Types	Element type definitions	{TypID EltName Option1 ... }
Sections	Section definitions	[A ky kz Ixx Iyy Izz yt yb zt zb]
Materials	Material definitions	[MatID MatProp1 MatProp2 ...]
Forces	Element forces in LCS (beam convention)	[N Vy Vz T My Mz]
		(nElem * 12)
Dloads	Distributed loads	[EltID n1globalX n1globalY n1globalZ ...]

plotstress(...,ParamName,ParamValue) sets the value of the specified parameters. The following parameters can be specified:

'StressScal'	Stress scaling. Default: 'auto'.
'MinMax'	Display minimum and maximum value. Default: 'off'.
'Values'	Stress values. Default: 'on'.
'Undeformed'	Plots the undeformed mesh. Default: 'k-'.
'Handle'	Plots in the axis with this handle. Default: current axis.

Additional parameters are redirected to the PLOT3 function which plots the stresses.

[StressScal,h] = plotstress(...) returns a struct h with handles to all the objects in the plot.

See also SDIAGRGCS_BEAM, SDIAGRGCS_TRUSS.

plotstresscontour

PLOTSTRESSCONTOUR Plot stress contour lines in shell elements.

```
plotstresscontour(stype,Nodes,Elements,Types,S)
plots stress contours (output from ELEMSTRESS/NODALSTRESS).
```

```
stype      'sx'      Normal stress (in the global/local x-direction)
           'sy'
           'sz'
           'sxy'     Shear stress
           'syz'
           'sxz'
Nodes      Node definitions      [NodID x y z]
Elements   Element definitions   [EltID TypID SecID MatID n1 n2 ...]
Types      Element type definitions {TypID EltName Option1 ... }
S          Element stresses in GCS/LCS [sxx syy szz sxy syz sxz]
                                           (nElem * 72)
```

plotstresscontour(...,ParamName,ParamValue) sets the value of the specified parameters. The following parameters can be specified:

```
'location'   Location (top,mid,bot). Default: 'top'.
'Ncontour'   Number of contours. Default: '10'.
'GCS'        Plot the GCS. Default: 'on'.
'Undeformed' Plots the undeformed mesh. Default: 'k-'.
'Handle'     Plots in the axis with this handle. Default: current axis.
```

See also ELEMSTRESS, SCONTOUR_SHELL8, SCONTOUR_SHELL4.

plotstresscontourf

PLOTSTRESSCONTOURF Plot filled contours of stresses.

```
plotstresscontourf(stype,Nodes,Elements,Types,S)
plotstresscontourf(stype,Nodes,Elements,Types,S,DOF,U)
plots stress contours (output from ELEMSTRESS).
```

stype	'sx'	Normal stress (in the global/local x-direction)
	'sy'	
	'sz'	
	'sxy'	Shear stress
	'syz'	
	'sxz'	
Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Types	Element type definitions	{TypID EltName Option1 ... }
S	Element stresses in GCS/LCS	[sxx syy szz sxy syz sxz] (nElem * 72)

plotstresscontourf(...,ParamName,ParamValue) sets the value of the specified parameters. The following parameters can be specified:

'location'	Location (top,mid,bot). Default: 'top'.
'GCS'	Plot the GCS. Default: 'on'.
'minmax'	Add location of min and max stress. Default: 'on'.
'colorbar'	Add colorbar. Default: 'on'.
'ncolor'	Number of colors in colormap. Default: 10.
'Handle'	Plots in the axis with this handle. Default: current axis.

See also ELEMSTRESS.

pressure_shell4

PRESSURE_SHELL4 Equivalent nodal forces for a shell4 element in the GCS due to a pressure normal to the element surface.

`F = pressure_shell4(Pressure,Node)`
computes the equivalent nodal forces of a pressure load normal to the elements surface.

Pressure	Distributed pressure in corner Nodes	[p1lobalZ; ...] (4 * 1)
Node	Node definitions	[x y z] (4 * 3)
F	Load vector	(24 * 1)

See also `ELEMPRESSURE`, `PRESSURE_SHELL8`.

pressure_shell6

PRESSURE_SHELL6 Equivalent nodal forces for a shell6 element in the GCS due to a pressure normal to the element surface.

`F = pressure_shell(Pressure,Node)`
 computes the equivalent nodal forces of a pressure load normal to the elements surface.

Pressure	Distributed loads in corner Nodes	<code>[p1localZ; ...]</code>	$(3 * 1)$
Node	Node definitions	<code>[x y z]</code>	$(6 * 3)$
F	Load vector		$(36 * 1)$

See also `ELEMPRESSURE`, `PRESSURE_SHELL4`.

pressure_shell8

PRESSURE_SHELL8 Equivalent nodal forces for a shell8 element in the GCS
due to a pressure normal to the element surface.

`F = pressure_shell(Pressure,Node)`
computes the equivalent nodal forces of a pressure load normal to
the elements surface.

Pressure	Distributed loads in corner Nodes	<code>[p1localZ; ...]</code>	$(4 * 1)$
Node	Node definitions	<code>[x y z]</code>	$(8 * 3)$
F	Load vector		$(48 * 1)$

See also `ELEMPRESSURE`, `PRESSURE_SHELL4`.

principalstress

PRINCIPALSTRESS Compute the principal stresses and directions in shell elements.

```
[Spr,Vpr] = principalstress(Elements,SeGCS)
computes the principal stresses and directions.
```

Elements	Element definitions	[EltID TypID SecID MatID n1 n2...]
SeGCS	Element stresses in GCS in corner nodes IJKL and at top/mid/bot of shell (nElem * 72) 72 = 6 stresscomp. * 4 nodes * 3 locations	
		[sxx syy szz sxy syz sxz ...]
Spr	Principal stress (nElem * 72)	
		[s1 s2 s3 0 0 0 ...]
Vpr	Principal stress directions {nElem * 12}	

See also **ELEMSTRESS**, **PLOTPRINCDIR**.

printdisp

PRINTDISP Display the displacements in the command window.

```
printdisp(Nodes,DOF,U)
displays the displacements in the command window.
```

Nodes	Node definitions	[NodID x y z]
DOF	Degrees of freedom	(nDOF * 1)
U	Displacements	(nDOF * 1)

See also PRINTFORC, PLOTDISP.

printforc

PRINTFORC Display the forces in the command window.

```
printforc(Elements,Forces)
displays the forces in the command window.
```

Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Forces	Element forces	[N Vy Vz T My Mz] (nElem * 12)

See also PRINTDISP.

printshellf

PRINTSHELLF Display forces/moments in command window (shell elements).

```
printshellf(Elements,F)
displays shell forces in command window.
```

Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
F	Shellf matrix	[Nx Ny Nxy Mx My Mxy Vx Vy] (nElem * 32)

See also PRINTSTRESS.

printstress

PRINTSTRESS Display stress in command window (shell elements).

`printstress(Elements,S,location)` displays stress in command window.

Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
S	Stress matrix	[sx sy sz sxy syz szx] (nElem * 72)

See also PRINTSHELLF.

q_dkt

`Q_DKT` `Q` matrix for a DKT element.

`Q = Q_DKT(b,c,det)` returns the `Q` matrix of the DKT element,
which is used to compute the curvatures of the plate element

<code>b</code>	Geometrical property of the triangle (see <code>ke_dkt</code>)	(3 * 1)
<code>c</code>	Geometrical property of the triangle	(3 * 1)
<code>det</code>	Determinant of the parametric transformation (=2*area of triangle)	

See also `KE_DKT`, `KELCS_SHELL4`, `SE_SHELL4`.

reaction

REACTION Compute the reaction forces for a degree of freedom.

```
Freac=reaction(Elements,ForcesGCS,seldof)
computes the reaction forces for a degree of freedom.
```

```
Elements    Element definitions            [EltID TypID SecID MatID n1 n2 ...]
ForcesGCS   Element forces in GCS (nElem * 12)
seldof      Selected DOF labels (kDOF * 1)
Freac       Reaction force
```

See also ELEMFORCES.

removedof

REMOVEDOF Remove DOF with Dirichlet boundary conditions equal to zero.

`DOF=removedof(DOF,seldof)` removes the specified Dirichlet boundary conditions from the degrees of freedom vector.

DOF Degrees of freedom (`nDOF * 1`)

seldof Dirichlet boundary conditions equal to zero [`NodID.dof`]

See also `GETDOF`.

reproW

REPROW Replicate rows from a matrix.

```
Matrix=reproW(Matrix,RowSel,nTime,RowInc)
```

replicates the selected rows from a matrix a number of times and adds them below the existing rows. The k-th time the increment RowInc is added k times to the copied rows.

Matrix	Matrix (nRow * nCol)
RowSel	Rows to be copied (1 * kRow)
nTime	Number of times
RowInc	Increments that are added to the different columns of the copied rows (1 * nCol)

scontour_shell4

SCONTOUR_SHELL4 Matrix to plot contours in a shell4 element.

Scontour = scontour_shell4(Node,Stress,Svalues) returns a matrix used for plotting contours in a shell4 element.

Node	Node definitions	[x y z] (4 * 3)
Stress	Stress in nodes	(4 * 1)
Svalues	Values of contours	(nContour * 1)
Scontour	Coordinates of contours in GCS	

See also PLOTSTRESSCONTOUR, PLOTSHELLFCONTOUR.

scontour_shell8

SCONTOUR_SHELL8 Matrix to plot contours in a shell8 element.

Scontour = scontour_shell8(Node,Stress,Svalues) returns a matrix used for plotting contours in a shell4 element.

Node	Node definitions	[x y z] (4 * 3)
Stress	Stress in nodes	(4 * 1) or (8 * 1)
Svalues	Values of contours	(nContour * 1)
Scontour	Coordinates of contours in GCS	

See also PLOTSTRESSCONTOUR, PLOTSHELLFCONTOUR.

sdiagrgcs_beam

SDIAGRGCS_BEAM Return matrices to plot the stresses in a beam element.

```
[ElemGCS,SdiagrGCS,ElemExtGCS,ExtremaGCS,Extrema]
    = sdiagrgcs_beam(stype,Forces,Node,Section,[],DLoad,Points)
[ElemGCS,SdiagrGCS,ElemExtGCS,ExtremaGCS,Extrema]
    = sdiagrgcs_beam(Forces,Node,Section,[],DLoad)
```

returns the coordinates of the points along the beam in the global coordinate system and the coordinates of the stresses with respect to the beam in the global coordinate system. These can be added in order to plot the stresses: ElemGCS+SdiagrGCS. The coordinates of the points with extreme values and the coordinates of the extreme values with respect to the beam are given as well and can be similarly added: ElemExtGCS+ExtremaGCS. Extrema is the list with the corresponding extreme values.

stype	'snorm'	Normal stress due to normal force
	'smomyt'	Normal stress due to bending moment around the local y-direction at the top
	'smomyb'	Normal stress due to bending moment around the local y-direction at the bottom
	'smomzt'	Normal stress due to bending moment around the local z-direction at the top
	'smomzb'	Normal stress due to bending moment around the local z-direction at the bottom
	'smax'	Maximal normal stress (normal force and bending moment)
	'smin'	Minimal normal stress (normal force and bending moment)
Forces	Element forces in LCS (beam convention) [N; Vy; Vz; T; My; Mz] (12 * 1)	
Node	Node definitions	[x y z] (3 * 3)
Section	Section definition	[A ky kz Ixx Iyy Izz yt yb zt zb]
DLoad	Distributed loads	[n1globalX; n1globalY; n1globalZ; ...] (6 * 1)
Points	Points in the local coordinate system (1 * nPoints)	
ElemGCS	Coordinates of the points along the beam in GCS (nPoints * 3)	
SdiagrGCS	Coordinates of the stress with respect to the beam in GCS (nValues * 3)	
ElemExtGCS	Coordinates of the points with extreme values in GCS (nValues * 3)	
ExtremaGCS	Coordinates of the extreme values with respect to the beam in GCS (nValues * 3)	
Extrema	Extreme values (nValues * 1)	

See also PLOTSTRESS, SDIAGRLCS_BEAM, SDIAGRGCS_TRUSS.

sdiagrgcs_shell2

SDIAGRGCS_SHELL2 Return matrices to plot the stresses in a SHELL2 element.

```
[ElemGCS,Sdiagrgcs,ElemExtGCS,ExtremaGCS,Extrema]
    = sdiagrgcs_shell2(ftype,Forces,Node,Section,Material,DLoad,Points)
[ElemGCS,Sdiagrgcs,ElemExtGCS,ExtremaGCS,Extrema]
    = sdiagrgcs_shell2(ftype,Forces,Node,Section,Material,DLoad)
```

returns the coordinates of the points along the SHELL2 in the global coordinate system and the coordinates of the stresses with respect to the element in the global coordinate system. These can be added in order to plot the stresses: ElemGCS+Sdiagrgcs. The coordinates of the points with extreme values and the coordinates of the extreme values with respect to the element are given as well and can be similarly added: ElemExtGCS+ExtremaGCS. Extrema is the list with the correspondig extreme values.

ftype	'sNphi'	Stress due to normal force in meridional direction
	'sMphiT'	Stress at the top due to bending moment in meridional direction
	'sMphiB'	Stress at the bottom due to bending moment in meridional direction
	'sNtheta'	Stress due to normal force in circumferential direction
	'sMthetaT'	Stress at the top due to bending moment in circumferential direction
	'sMthetaB'	Stress at the bottom due to bending moment in circumferential direction
Forces	Element forces in LCS (beam convention) [N; Vy; 0; 0; 0; Mz] (12 * 1)	
Node	Node definitions [x y z] (3 * 3)	
DLoad	Distributed loads [n1globalX; n1globalY; n1globalZ; ...] (6 * 1)	
Points	Points in the local coordinate system (1 * nPoints)	
ElemGCS	Coordinates of the points along the element in GCS (nPoints * 3)	
Fdiagrgcs	Coordinates of the force with respect to the element in GCS (nValues * 3)	
ElemExtGCS	Coordinates of the points with extreme values in GCS (nValues * 3)	
ExtremaGCS	Coordinates of the extreme values with respect to the element in GCS (nValues * 3)	
Extrema	Extreme values (nValues * 1)	

sdiagrgcs_truss

SDIAGRGCS_TRUSS Return matrices to plot the stresses in a truss element.

```
[ElemGCS,SdiagrGCS,ElemExtGCS,ExtremaGCS,Extrema]
    = sdiagrgcs_truss(stype,Forces,Node,Section,[],[],Points)
[ElemGCS,SdiagrGCS,ElemExtGCS,ExtremaGCS,Extrema]
    = sdiagrgcs_truss(stype,Forces,Node,Section)
```

returns the coordinates of the points along the truss in the global coordinate system and the coordinates of the stresses with respect to the truss in the global coordinate system. These can be added in order to plot the stresses: ElemGCS+SdiagrGCS. The coordinates of the points with extreme values and the coordinates of the extreme values with respect to the truss are given as well and can be similarly added: ElemExtGCS+ExtremaGCS. Extrema is the list with the corresponding extreme values.

```
stype      'snorm'      Normal stress due to normal force
Forces      Element forces in LCS [N; 0; 0; 0; 0; 0](12 * 1)
Node        Node definitions      [x y z] (3 * 3)
Section     Section definition     [A ky kz Ixx Iyy Izz yt yb zt zb]
Points      Points in the local coordinate system (1 * nPoints)
ElemGCS     Coordinates of the points along the truss in GCS (nPoints * 3)
SdiagrGCS   Coordinates of the force with respect to the truss in GCS
                                                    (nValues * 3)
ElemExtGCS  Coordinates of the points with extreme values in GCS (nValues * 3)
ExtremaGCS  Coordinates of the extreme values with respect to the truss in GCS
                                                    (nValues * 3)

Extrema     Extreme values (nValues * 1)
```

See also PLOTSTRESS, SDIAGRGCS_BEAM.

sdiagrlcs_beam

SDIAGRCS_BEAM Stress diagram for a beam element in LCS.

[SdiagrLCS,loc,Extrema] = sdiagrlcs_beam(ftype,Forces,DLoadLCS,L,Points)
computes the stresses at the specified points. The extreme values are analytically determined.

stype	'snorm'	Normal stress due to normal force
	'smomyt'	Normal stress due to bending moment around the local y-direction at the top
	'smomyb'	Normal stress due to bending moment around the local y-direction at the bottom
	'smomzt'	Normal stress due to bending moment around the local z-direction at the top
	'smomzb'	Normal stress due to bending moment around the local z-direction at the bottom
	'smax'	Maximal normal stress (normal force and bending moment)
	'smin'	Minimal normal stress (normal force and bending moment)
Forces	Element forces in LCS (beam convention) [N; Vy; Vz; T; My; Mz](12 * 1)	
DLoadLCS	Distributed loads in LCS [n1localX; n1localY; n1localZ; ...](6 * 1)	
Points	Points in the local coordinate system (1 * nPoints)	
Section	Section definition [A ky kz Ixx Iyy Izz yt yb zt zb]	
SdiagrLCS	Stresses at the points (1 * nPoints)	
loc	Locations of the extreme values (nValues * 1)	
Extrema	Extreme values (nValues * 1)	

See also SDIAGRCS_BEAM.

selcs_plane3

SELCS_PLANE3 Compute the element stresses for a plane3 element.

```
[SeLCS] = selcs_plane3(Node,Section,Material,UeGCS,Options)
computes the element stresses in the
local coordinate system for the plane3 element.
```

Node	Node definitions	[x y z] (4 * 3)
Section	Section definition	[h] (only used in plane stress)
Material	Material definition	[E nu rho]
UeLCS	Displacements	(6 * nSteps)
Options	Element options	{Option1 Option2 ...}
SeLCS	Element stresses in LCS in corner nodes IJKL	(9 * nTimeSteps) [sxx syy sxy]

See also **ELEMSTRESS**, **SE_PLANE3**.

selcs_plane4

SELCS_PLANE4 Compute the element stresses for a plane4 element.

```
[SeLCS] = selcs_plane4(Node,Section,Material,UeGCS,Options)
computes the element stresses in the
local coordinate system for the plane4 element.
```

Node	Node definitions	[x y z] (4 * 3)
Section	Section definition	[h] (only used in plane stress)
Material	Material definition	[E nu rho]
UeLCS	Displacements	(8 * nSteps)
Options	Element options	{Option1 Option2 ...}
SeLCS	Element stresses in LCS in corner nodes IJKL	
	(12 * nTimeSteps) [sxx syy sxy] for plane stress / plane strain problems	
	(16 * nTimeSteps) [sxx stheta syy sxy] for axisymmetric problems	

See also **ELEMSTRESS**, **SE_PLANE4**.

selcs_plane6

SELCS_PLANE6 Compute the element stresses for a plane6 element.

```
[SeLCS] = selcs_plane6(Node,Section,Material,UeGCS,Options)
computes the element stresses in the
local coordinate system for the plane6 element.
```

Node	Node definitions	[x y z] (6 * 3)
	Nodes should have the following order:	
	3	
	\	
	6 5	
	\	
	1--4--2	
Section	Section definition	[h] (only used in plane stress)
Material	Material definition	[E nu rho]
UeLCS	Displacements	(6 * nSteps)
Options	Element options	{Option1 Option2 ...}
SeLCS	Element stresses in LCS in corner nodes IJKL	(9 * nTimeSteps) [sxx syy sxy]

See also **ELEMSTRESS**, **SE_PLANE6**.

selcs_plane8

SELCS_PLANE8 Compute the element stresses for a plane4 element.

```
[SeLCS] = selcs_plane8(Node,Section,Material,UeGCS,Options)
computes the element stresses in the
local coordinate system for the plane8 element.
```

Node	Node definitions	[x y z] (4 * 3)
Section	Section definition	[h] (only used in plane stress)
Material	Material definition	[E nu rho]
UeLCS	Displacements	(8 * nSteps)
Options	Element options	{Option1 Option2 ...}
SeLCS	Element stresses in LCS in corner nodes IJKL	(12 * nTimeSteps) [sxx syy sxy]

See also ELEMSTRESS, SE_PLANE8.

selcs_shell4

SELCS_SHELL4 Compute the element stresses for a shell4 element.

```
[SeLCS] = selcs_shell4(Node,Section,Material,UeGCS,Options)
computes the element stresses in the global and the
local coordinate system for the shell4 element.
```

Node	Node definitions	[x y z] (4 * 3)
h	Shell thickness	
E	Young's modulus	
nu	Poisson coefficient	
rho	Mass density	
UeLCS	Displacements (24 * nTimeSteps)	
Options	Element options	{Option1 Option2 ...}
SeLCS	Element stresses in LCS in corner nodes IJKL and at top/mid/bot of shell (72 * nTimeSteps)	[sxx syy szz sxy syz sxz]

See also **ELEMSTRESS**, **SE_SHELL4**.

selectdof

SELECTDOF Select degrees of freedom.

```
L=selectdof(DOF,seldof)
[L,I]=selectdof(DOF,seldof)
L=selectdof(DOF,seldof,'Ordering',ordering)
creates the matrix to extract degrees of freedom from the global degrees of
freedom by matrix multiplication.
```

DOF Degrees of freedom (nDOF * 1)
seldof Selected DOF labels (kDOF * 1)
L Selection matrix (kDOF * nDOF)
I Index vector (kDOF * 1)
ordering 'seldof','DOF' or 'sorted' Default: 'seldof'
 Ordering of L and I similar as seldof, DOF or sorted

selectnode

SELECTNODE Select nodes by location.

```
Nodesel=selectnode(Nodes,x,y,z)
Nodesel=selectnode(Nodes,xmin,ymin,zmin,xmax,ymax,zmax)
selects nodes by location.
```

Nodes	Node definitions	[NodID x y z]
Nodesel	Node definitions of the selected nodes	

se_plane3

SE_PLANE3 Compute the element stresses for a plane3 element.

```
[SeGCS,SeLCS,vLCS] = se_plane3(Node,Section,Material,UeGCS,Options,GCS)
[SeGCS,SeLCS]       = se_plane3(Node,Section,Material,UeGCS,Options,GCS)
SeGCS               = se_plane3(Node,Section,Material,UeGCS,Options,GCS)
```

computes the element stresses in the global and the local coordinate system for the shell3 element.

Node	Node definitions	[x y z] (3 * 3)
	Nodes should have the following order:	
	3	
	\	
	1--2	
Section	Section definition	[h] (only used in plane stress)
Material	Material definition	[E nu rho]
UeGCS	Displacements (8 * nTimeSteps)	
Options	Element options	{Option1 Option2 ...}
GCS	Global coordinate system in which stresses are returned	
	'cart' 'cyl'	
SeGCS	Element stresses in GCS in corner nodes IJKL	
	18 = 6 stress comp. * 3 nodes (18 * nTimeSteps)	
	[sxx syy szz sxy syz sxz]	
SeLCS	Element stresses in LCS in corner nodes IJKL	
	18 = 6 stress comp. * 3 nodes (18 * nTimeSteps)	
	[sxx syy szz sxy syz sxz]	
vLCS	Unit vectors of LCS (1 * 9)	

See also ELEMSTRESS, SE_SHELL8.

se_plane4

SE_PLANE4 Compute the element stresses for a plane4 element.

```
[SeGCS,SeLCS,vLCS] = se_plane4(Node,Section,Material,UeGCS,Options,GCS)
[SeGCS,SeLCS]       = se_plane4(Node,Section,Material,UeGCS,Options,GCS)
SeGCS               = se_plane4(Node,Section,Material,UeGCS,Options,GCS)
```

computes the element stresses in the global and the local coordinate system for the shell4 element.

Node	Node definitions	[x y z] (4 * 3)
	Nodes should have the following order:	
	4---3	
	1---2	
Section	Section definition	[h] (only used in plane stress)
Material	Material definition	[E nu rho]
UeGCS	Displacements (8 * nTimeSteps)	
Options	Element options	{Option1 Option2 ...}
GCS	Global coordinate system in which stresses are returned	'cart' 'cyl'
SeGCS	Element stresses in GCS in corner nodes IJKL	
	24 = 6 stress comp. * 4 nodes (24 * nTimeSteps)	
		[sxx syy szz sxy syz sxz]
SeLCS	Element stresses in LCS in corner nodes IJKL	
	24 = 6 stress comp. * 4 nodes (24 * nTimeSteps)	
		[sxx syy szz sxy syz sxz]
vLCS	Unit vectors of LCS (1 * 9)	

See also ELEMSTRESS, SE_SHELL8.

se_plane6

SE_PLANE6 Compute the element stresses for a plane6 element.

```
[SeGCS,SeLCS,vLCS] = se_plane6(Node,Section,Material,UeGCS,Options,GCS)
[SeGCS,SeLCS]       = se_plane6(Node,Section,Material,UeGCS,Options,GCS)
SeGCS               = se_plane6(Node,Section,Material,UeGCS,Options,GCS)
```

computes the element stresses in the global and the local coordinate system for the plane6 element.

Node	Node definitions	[x y z] (8 * 3)
	Nodes should have the following order:	
	<pre> 3 \ 6 5 \ 1--4--2</pre>	
Section	Section definition	[h] (only used in plane stress)
Material	Material definition	[E nu rho]
UeGCS	Displacements (8 * nTimeSteps)	
Options	Element options	{Option1 Option2 ...}
GCS	Global coordinate system in which stresses are returned	'cart' 'cyl'
SeGCS	Element stresses in GCS in corner nodes IJKL	
	48 = 6 stress comp. * 8 nodes (24 * nTimeSteps)	
		[sxx syy szz sxy syz sxz]
SeLCS	Element stresses in LCS in corner nodes IJKL	
	48 = 6 stress comp. * 8 nodes (24 * nTimeSteps)	
		[sxx syy szz sxy syz sxz]
vLCS	Unit vectors of LCS (1 * 9)	

See also ELEMSTRESS, SE_PLANE8.

se_plane8

SE_PLANE8 Compute the element stresses for a plane8 element.

```
[SeGCS,SeLCS,vLCS] = se_plane8(Node,Section,Material,UeGCS,Options,GCS)
[SeGCS,SeLCS]       = se_plane8(Node,Section,Material,UeGCS,Options,GCS)
SeGCS               = se_plane8(Node,Section,Material,UeGCS,Options,GCS)
```

computes the element stresses in the global and the local coordinate system for the shell8 element.

Node	Node definitions	[x y z] (8 * 3)
	Nodes should have the following order:	
	4-----7-----3	
	8	6
	1-----5-----2	
Section	Section definition	[h] (only used in plane stress)
Material	Material definition	[E nu rho]
UeGCS	Displacements (8 * nTimeSteps)	
Options	Element options	{Option1 Option2 ...}
GCS	Global coordinate system in which stresses are returned	
	'cart' 'cyl'	
SeGCS	Element stresses in GCS in corner nodes IJKL	
	48 = 6 stress comp. * 8 nodes (24 * nTimeSteps)	
	[sxx syy szz sxy syz sxz]	
SeLCS	Element stresses in LCS in corner nodes IJKL	
	48 = 6 stress comp. * 8 nodes (24 * nTimeSteps)	
	[sxx syy szz sxy syz sxz]	
vLCS	Unit vectors of LCS (1 * 9)	

See also ELEMSTRESS, SE_SHELL8.

se_shell4

SE_SHELL4 Compute the element stresses for a shell4 element.

```
[SeGCS,SeLCS,vLCS] = se_shell4(Node,Section,Material,UeGCS,Options,GCS)
[SeGCS,SeLCS]       = se_shell4(Node,Section,Material,UeGCS,Options,GCS)
SeGCS               = se_shell4(Node,Section,Material,UeGCS,Options,GCS)
```

computes the element stresses in the global and the local coordinate system for the shell4 element.

Node	Node definitions	[x y z] (4 * 3)
	Nodes should have the following order:	
Section	Section definition	[h]
Material	Material definition	[E nu rho]
UeGCS	Displacements (24 * nTimeSteps)	
Options	Element options	{Option1 Option2 ...}
GCS	Global coordinate system in which stresses are returned	'cart' 'cyl' 'sph'
SeGCS	Element stresses in GCS in corner nodes IJKL and at top/mid/bot of shell (72 * nTimeSteps)	
	72 = 6 stress comp. * 4 nodes * 3 locations	[sxx syy szz sxy syz sxz]
SeLCS	Element stresses in LCS in corner nodes IJKL and at top/mid/bot of shell (72 * nTimeSteps)	
		[sxx syy szz sxy syz sxz]
vLCS	Unit vectors of LCS (1 * 9)	

See also ELEMSTRESS, SE_SHELL8.

se_shell6

SE_SHELL6 Compute the element stresses for a shell6 element.

```
[SeGCS,SeLCS,vLCS] = se_shell6(Node,Section,Material,UeGCS,Options,gcs)
[SeGCS,SeLCS]       = se_shell6(Node,Section,Material,UeGCS,Options,gcs)
SeGCS               = se_shell6(Node,Section,Material,UeGCS,Options,gcs)
computes the element stresses in the global and the
local coordinate system for the shell6 element.
```

Node	Node definitions [x y z] (6 * 3) Nodes should have the following order: 3 \ 6 5 \ 1--4--2
Section	Section definition [h] or [h1 h2 h3] (uniform thickness or defined in corner nodes(1,2,3))
Material	Material definition [E nu rho] or [Exx Eyy nuxy muxy muyz muzx theta rho]
UeGCS	Displacements (36 * nTimeSteps)
Options	Element options struct. Fields: -LCSType: determine the reference local element coordinate system. Values: 'element' (default) or 'global' -MatType: 'isotropic' (default) or 'orthotropic' -Offset: nodal offset from shell midplane. Values: 'top', 'mid' (default), 'bot' or numerical value
GCS	Global coordinate system in which stresses are returned 'cart' 'cyl' 'sph'
SeGCS	Element stresses in GCS in corner nodes IJKL and at top/mid/bot of shell (72 * nTimeSteps) 72 = 6 stress comp. * 4 nodes * 3 locations (in a triangular element the fourth node is zero) [sxx syy szz sxy syz sxz]
SeLCS	Element stresses in LCS in corner nodes IJKL and at top/mid/bot of shell (54 * nTimeSteps) [sxx syy szz sxy syz sxz]
vLCS	Unit vectors of LCS (3 * 3)

See also ELEMSTRESS, SE_SHELL4.

se_shell8

SE_SHELL8 Compute the element stresses for a shell8 element.

```
[SeGCS,SeLCS,vLCS] = se_shell8(Node,Section,Material,UeGCS,Options,gcs)
[SeGCS,SeLCS]       = se_shell8(Node,Section,Material,UeGCS,Options,gcs)
SeGCS               = se_shell8(Node,Section,Material,UeGCS,Options,gcs)
computes the element stresses in the global and the
local coordinate system for the shell8 element.
```

Node	Node definitions	[x y z] (8 * 3)
	Nodes should have the following order:	
	<pre> 4-----7-----3 8 6 1-----5-----2 </pre>	
Section	Section definition	[h] or [h1 h2 h3 h4] (uniform thickness or defined in corner nodes(1,2,3,4))
Material	Material definition	[E nu rho] or [Exx Eyy nuxy muxy muyz muzx theta rho]
UeGCS	Displacements (48 * nTimeSteps)	
Options	Element options struct. Fields:	
	-LCSType: determine the reference local element coordinate system. Values: 'element' (default) or 'global'	
	-MatType: 'isotropic' (default) or 'orthotropic'	
	-Offset: nodal offset from shell midplane. Values: 'top', 'mid' (default), 'bot' or numerical value	
GCS	Global coordinate system in which stresses are returned 'cart' 'cyl' 'sph'	
SeGCS	Element stresses in GCS in corner nodes IJKL and at top/mid/bot of shell (72 * nTimeSteps) 72 = 6 stress comp. * 4 nodes * 3 locations [sxx syy szz sxy syz sxz]	
SeLCS	Element stresses in LCS in corner nodes IJKL and at top/mid/bot of shell (72 * nTimeSteps) [sxx syy szz sxy syz sxz]	
vLCS	Unit vectors of LCS (3 * 3)	

See also ELEMSTRESS, SE_SHELL4.

se_solid20

SE_SOLID20 Compute the element stresses for a solid20 element.

```
[SeGCS,SeLCS,vLCS] = se_solid20(Node,Section,Material,UeGCS,Options,gcs)
[SeGCS,SeLCS]       = se_solid20(Node,Section,Material,UeGCS,Options,gcs)
SeGCS               = se_solid20(Node,Section,Material,UeGCS,Options,gcs)
```

computes the element stresses in the global and the local coordinate system for the solid20 element.

Node Node definitions [x y z] (20 * 3)

Nodes should have the following order:

```
      8---15----7
      /|        /|
    16 |        14 |
      / 20      / 19
      / |        / |
    5---13----6   |
    |   4--11|----3
    | /      | /
  17 12      18 10
    | /      | /
    | /      | /
    1----9----2
```

se_solid8

SE_SOLID8 Compute the element stresses for a solid8 element.

```
[SeGCS,SeLCS,vLCS] = se_solid8(Node,Section,Material,UeGCS,Options,gcs)
[SeGCS,SeLCS]       = se_solid8(Node,Section,Material,UeGCS,Options,gcs)
SeGCS               = se_solid8(Node,Section,Material,UeGCS,Options,gcs)
computes the element stresses in the global and the
local coordinate system for the solid8 element.
```

Node Node definitions [x y z] (8 * 3)
Nodes should have the following order:

```

      8-----7
      /|      /|
     / |     / |
    /  |    /  |
5-----6   |
|   4-----|---3
|  /       | /
| /        | /
|/         |/
1-----2
```

Section	Section definition	[]
Material	Material definition	[E nu rho]
UeGCS	Displacements (48 * nTimeSteps)	
Options	Element options struct. Fields:	[]
GCS	Global coordinate system in which stresses are returned	'cart' 'cyl' 'sph'
SeGCS	Element stresses in GCS in nodes (48 * nTimeSteps)	
	48 = 6 stress comp. * 8 nodes	[sxx syy szz sxy syz sxz]
SeGCS	Element stresses in GCS in nodes (48 * nTimeSteps)	
	48 = 6 stress comp. * 8 nodes	[sxx syy szz sxy syz sxz]
vLCS	Unit vectors of LCS (3 * 3)	

See also **ELEMSTRESS**, **SE_SHELL4**.

se_truss

`e_truss` is a function.

```
[SeGCS, SeLCS, vLCS] = se_truss(Node, Section, Material, UeGCS, Options, gcs)
```

sh_qs4

SH_QS4 Shape functions for a quadrilateral serendipity element with 4 nodes.

`[Ni,dN_dxi,dN_deta] = sh_qs4(xi,eta)` returns the shape functions and its derivatives to the natural coordinates in the point `(xi,eta)`.

<code>xi</code>	Natural coordinate	(scalar)
<code>eta</code>	Natural coordinate	(scalar)
<code>Ni</code>	Shape functions in point <code>(xi,eta)</code>	(4 * 1)
<code>dN_dxi</code>	Derivative of <code>Ni</code> to <code>xi</code>	(4 * 1)
<code>dN_deta</code>	Derivative of <code>Ni</code> to <code>eta</code>	(4 * 1)

See also KELCS_SHELL4.

sh_qs8

SH_QS8 Shape functions for an 8 node quadrilateral serendipity element.

[Ni,dN_dxi,dN_deta] = sh_qs8(xi,eta) returns the shape functions and its derivatives with respect to the natural coordinates in the point (xi,eta).

xi	natural coordinate	(scalar)
eta	natural coordinate	(scalar)
Ni	shape functions in point (xi,eta)	(8 * 1)
dN_dxi	derivative of Ni to xi	(8 * 1)
dN_deta	derivative of Ni to eta	(8 * 1)

see also KE_SHELL8.

sh_t

SH_T Shape functions for a triangular plate element.

[Ni] = sh_t(L,b,c) returns the shape functions and
its derivatives in point L.

L	Area coordinates [L1,L2,L3]	(3 * 1)
b	Geometrical property of the triangle (see ke_dkt)	(3 * 1)
c	Geometrical property of the triangle	(3 * 1)
Ni	Shape functions and derivatives in point L	(9 * 1)

These shape functions are used to determine the mass matrix of a
triangular plate element.

See also KE_DKT.

sh_t10

SH_T10 Shape functions for a 10 node triangular element.

`[Ni,dN_dxi,dN_deta] = sh_t10(xi,eta)` returns the shape functions and its derivatives with respect to the natural coordinates in the point (xi,eta)

<code>xi</code>	natural coordinate	(scalar)
<code>eta</code>	natural coordinate	(scalar)
<code>Ni</code>	shape functions in point (xi,eta)	(10 * 1)
<code>dN_dxi</code>	derivative of Ni to xi	(10 * 1)
<code>dN_deta</code>	derivative of Ni to eta	(10 * 1)

The nodes have the following order:

see also SH_T3, SH_T6, SH_T15.

sh_t15

SH_T15 Shape functions for a 15 node triangular element.

`[Ni,dN_dxi,dN_deta] = sh_t15(xi,eta)` returns the shape functions and its derivatives with respect to the natural coordinates in the point `(xi,eta)*`

<code>xi</code>	natural coordinate	(scalar)
<code>eta</code>	natural coordinate	(scalar)
<code>Ni</code>	shape functions in point <code>(xi,eta)</code>	(15 * 1)
<code>dN_dxi</code>	derivative of <code>Ni</code> to <code>xi</code>	(15 * 1)
<code>dN_deta</code>	derivative of <code>Ni</code> to <code>eta</code>	(15 * 1)

The nodes have the following order:

see also SH_T3, SH_T6, SH_T15.

sh_t3

SH_T3 Shape functions for a 3 node triangular element.

[Ni,dN_dxi,dN_deta] = sh_t3(xi,eta) returns the shape functions and its derivatives with respect to the natural coordinates in the point (xi,eta).

xi	natural coordinate	(scalar)
eta	natural coordinate	(scalar)
Ni	shape functions in point (xi,eta)	(3 * 1)
dN_dxi	derivative of Ni to xi	(3 * 1)
dN_deta	derivative of Ni to eta	(3 * 1)

sh_t6

SH_T6 Shape functions for an 6 node triangular element.

[Ni,dN_dxi,dN_deta] = sh_t6(xi,eta) returns the shape functions and its derivatives with respect to the natural coordinates in the point (xi,eta).

xi	natural coordinate	(scalar)
eta	natural coordinate	(scalar)
Ni	shape functions in point (xi,eta)	(6 * 1)
dN_dxi	derivative of Ni to xi	(6 * 1)
dN_deta	derivative of Ni to eta	(6 * 1)

see also KE_SHELL6.

sh_vs20

SH_VS20 Shape functions for a volume serendipity element with 20 nodes.

```
[Ni,dNi_dxi,dNi_deta,dNi_dzeta] = sh_vs20(xi,eta,zeta)
returns the shape functions and its derivatives to the natural coordinates
in the point (xi,eta,zeta).
```

xi	Natural coordinate	(scalar)
eta	Natural coordinate	(scalar)
zeta	Natural coordinate	(scalar)
Ni	Shape functions in point (xi,eta)	(20 * 1)
dN_dxi	Derivative of Ni to xi	(20 * 1)
dN_deta	Derivative of Ni to eta	(20 * 1)
dN_dzeta	Derivative of Ni to zeta	(20 * 1)

See also KE_SOLID20.

sh_vs8

SH_VS8 Shape functions for a volume serendipity element with 8 nodes.

```
[Ni,dNi_dxi,dNi_deta,dNi_dzeta] = sh_vs8(xi,eta,zeta)
returns the shape functions and its derivatives to the natural coordinates
in the point (xi,eta,zeta).
```

xi	Natural coordinate	(scalar)
eta	Natural coordinate	(scalar)
zeta	Natural coordinate	(scalar)
Ni	Shape functions in point (xi,eta)	(8 * 1)
dN_dxi	Derivative of Ni to xi	(8 * 1)
dN_deta	Derivative of Ni to eta	(8 * 1)
dN_dzeta	Derivative of Ni to zeta	(8 * 1)

See also KE_SOLID8.

size_beam

SIZE_BEAM Compute beam element size (length).

`S = SIZE_BEAM(Node)` computes the element size (length) of a two node beam element.

`[S,dSdx] = SIZE_BEAM(Node,dNodedx)` additionally computes the derivatives of the element size with respect to the design variables `x`.

<code>Node</code>	Node definitions	<code>[x y z] (3 * 3)</code>
<code>dNodedx</code>	Node definitions derivatives	<code>(SIZE(Node) * nVar)</code>
<code>S</code>	Element size	
<code>dSdx</code>	Element size derivatives	

See also `ELEMSIZES`, `ELEMVOLUMES`, `SIZE_TRUSS`.

size_plane6

`size_plane6` Compute plane6 element size (area).

`s = size_plane6(NodeNum)` computes the size (area) of a plane6 element.

Node	Node definitions	[x y z] (6 * 3)
S	Element size	

size_solid10

size_solid10 Compute solid10 element size (volume).

 S = size_solid10(NodeNum) computes the size of a solid10 element.

Node	Node definitions	[x y z] (10 * 3)
S	Element size	

size_solid4

`size_solid4` Compute solid4 element size (volume).

`s = size_solid4(NodeNum)` computes the size (volume) of a solid4 element.

Node	Node definitions	[x y z] (4 * 3)
S	Element size	

size_truss

`SIZE_TRUSS` Compute truss element size (length).

`S = SIZE_TRUSS(Node)` computes the element size (length) of a two node truss element.

`[S,dSdx] = SIZE_TRUSS(Node,dNodedx)` additionally computes the derivatives of the element size with respect to the design variables `x`.

<code>Node</code>	Node definitions	<code>[x y z] (3 * 3)</code>
<code>dNodedx</code>	Node definitions derivatives	<code>(SIZE(Node) * nVar)</code>
<code>S</code>	Element size	
<code>dSdx</code>	Element size derivatives	

See also `ELEMSIZES`, `ELEMVOLUMES`, `SIZE_BEAM`.

tconstr

TCONSTR Return matrices to apply constraint equations.

```
[T,Q0,MasterDOF]=tconstr(Constr,DOF)
```

returns matrices to apply constraint equations to the stiffness and mass matrix and the load vector: $K_r = T.' * K * T$, $M_r = T.' * M * T$ and $F_r = T.' * (F - K * Q_0)$.

The original displacement vector is computed using $U = T * U_r + Q_0$.

Constr Constraint equation:

 Constant=CoefS*SlaveDOF+CoefM1*MasterDOF1+CoefM2*MasterDOF2+...

 [Constant CoefS SlaveDOF CoefM1 MasterDOF1 CoefM2 MasterDOF2 ...]

DOF Degrees of freedom (nDOF * 1)

tloads_beam

TLOADS_BEAM Equivalent nodal forces for a beam element in the GCS.

```
F = tloads_beam(TLoad,Node,Section,Material)
computes the equivalent nodal forces of a temperature load
(in the global coordinate system).
```

TLoad	Temperature loads	[dTm; dTy; dTz] (3 * 1)
Node	Node definitions	[x y z] (2 * 3)
Section	Section definitions	[A ky kz Ixx Iyy Izz yt yb zt zb]
Material	Material definitions	[E nu rho alpha]
F	Load vector	(12 * 1)

See also **ELEMTLOADS**, **TLOADS_TRUSS**.

tloads_truss

TLOADS_TRUSS Equivalent nodal forces for a truss element in the GCS.

```
F = tloads_truss(TLoad,Node,Section,Material)
computes the equivalent nodal forces of a temperature load
(in the global coordinate system).
```

TLoad	Temperature loads	[n1globalX; n1globalY; n1globalZ; ...] (6 * 1)
Node	Node definition	[x y z] (2 * 3)
Section	Section definition	[A ...]
Material	Material definition	[E nu rho alpha]
F	Load vector	(6 * 1)

See also **ELEMTLOADS**, **TLOADS_BEAM**.

trans_beam

TRANS_BEAM Transform coordinate system for a beam element.

```
t = TRANS_BEAM(Node)
    computes the transformation matrix between the local and the global
    coordinate system for the beam element.
[t,dtdx] = TRANS_BEAM(Node,dNodedx)
    additionally computes the derivatives of the transformation matrix
    with respect to the design variables x.
```

Node	Node definitions	[x y z] (3 * 3)
dNodedx	Node definitions derivatives	(SIZE(Node) * nVar)
t	Transformation matrix	(3 * 3)
dtdx	Transformation matrix derivatives	(3 * 3 * nVar)

See also KE_BEAM, TRANS_TRUSS.

trans_shell2

TRANS_BEAM Transform coordinate system for a beam element.

```
t = trans_beam(Node)
computes the transformation matrix between the local and the global
coordinate system for the SHELL2 element.
```

Node	Node definitions	[x y z] (3 * 3)
t	Transformation matrix	(3 * 3)

See also KE_BEAM, TRANS_TRUSS.

trans_shell4

TRANS_SHELL4 Transform coordinate system for a shell4 element.

```
[t,Node_lc,W] = trans_shell4(Node)
[t,Node_lc]   = trans_shell4(Node)
t             = trans_shell4(Node)
```

computes the transformation matrix between the local and the global coordinate system and the correction matrix for non-coplanar nodes for the shell4 element.

Node	Node definitions	[x y z] (4 * 3)
t	Transformation matrix	(3 * 3)
Node_lc	Nodes in LCS	[x y z] (4 * 3)
W	Correction matrix for warped elements	(24 * 24)

See also KE_BEAM, TRANS_TRUSS.

trans_shell8

TRANS_SHELL8 Transform coordinate system of a shell8 element.

```
t = trans_shell8(Node)
t = trans_shell8(Node,Options)
computes the transformation matrix between the local and the global
coordinate system for stress computations.
```

```
Node      Node definitions      [x y z] (8 * 3)
t          Transformation matrix (3 * 3)
Options    Element options struct. Fields:
            -LCSType: determine the reference local element
                  coordinate system. Values:
                      'element' (default) or 'global'
```

See also SE_SHELL8, TRANS_TRUSS.

trans_solid20

TRANS_SOLID20 Transform coordinate system of a solid8 element.

```
t = trans_solid20(Node)
t = trans_solid20(Node,Options)
computes the transformation matrix between the local and the global
coordinate system for stress computations.
```

Node	Node definitions	[x y z]	(20 * 3)
t	Transformation matrix		(3 * 3)

See also SE_SOLID20, TRANS_TRUSS.

trans_solid8

TRANS_SOLID8 Transform coordinate system of a solid8 element.

```
t = trans_solid8(Node)
t = trans_solid8(Node,Options)
computes the transformation matrix between the local and the global
coordinate system for stress computations.
```

Node	Node definitions	[x y z]	(8 * 3)
t	Transformation matrix		(3 * 3)

See also SE_SOLID8, TRANS_TRUSS.

trans_truss

TRANS_TRUSS Transform coordinate system for a truss element.

```
t = TRANS_TRUSS(Node)
    computes the transformation matrix between the local and the global
    coordinate system for the truss element.
[t,dtdx] = TRANS_TRUSS(Node,dNodedx)
    additionally computes the derivatives of the transformation matrix
    with respect to the design variables x.
```

Node	Node definitions	[x y z] (2 * 3)
dNodedx	Node definitions derivatives	(SIZE(Node) * nVar)
t	Transformation matrix	(3 * 3)
dtdx	Transformation matrix derivatives	(3 * 3 * nVar)

See also KE_TRUSS, TRANS_BEAM.

udiargcs

UDIARGCS Return displacement diagrams in GCS

```
[UxdiagrGCS,UydiagrGCS,UzdiagrGCS] = UDIARGCS(Nodes,Elements,Types,DOF,U,DLoads,
                                                Sections,Materials,Points)
[UxdiagrGCS,UydiagrGCS,UzdiagrGCS] = UDIARGCS(Nodes,Elements,Types,DOF,U,DLoads,
                                                Sections,Materials)
[UxdiagrGCS,UydiagrGCS,UzdiagrGCS] = UDIARGCS(Nodes,Elements,Types,DOF,U,[],
                                                Sections,Materials)
[UxdiagrGCS,UydiagrGCS,UzdiagrGCS] = UDIARGCS(Nodes,Elements,Types,DOF,U)
    computes the displacements of the interpolation points after
    deformation in the global (algebraic) coordinate system. If DLoads,
    Sections and Materials are supplied, the displacements that occur
    due to distributed loads if all nodes are fixed, are superimposed.
```

```
[UxdiagrGCS,UydiagrGCS,UzdiagrGCS,dUxdiagrGCSdx,dUydiagrGCSdx,dUzdiagrGCSdx]
    = UDIARGCS(Nodes,Elements,Types,DOF,U,DLoads,Sections,Materials,Points,
               dNodesdx,dUdx,dDLoadsdx,dSectionsdx)
    additionally computes the derivatives of the displacement values
    with respect to the design variables x.
```

Nodes	Node definitions	[NodID x y z]
Elements	Element definitions	[EltID TypID SecID MatID n1 n2 ...]
Types	Element type definitions	{TypID EltName Option1 ... }
DOF	Degrees of freedom	(nDOF * 1)
U	Displacements	(nDOF * 1)
DLoads	Distributed loads	[EltID n1globalX n1globalY n1globalZ ...] (use empty array [] when shear deformation (in beam element) is considered but no DLoads are present)
Sections	Section definitions	[SecID SecProp1 SecProp2 ...]
Materials	Material definitions	[MatID MatProp1 MatProp2 ...]
Points	Points in the local coordinate system	(1 * nPoints)
dNodesdx	Node definitions derivatives	(SIZE(Nodes) * nVar)
dUdx	Displacements derivatives	(SIZE(U) * nVar)
dDLoadsdx	Distributed loads derivatives	(SIZE(DLoads) * nVar)
UxdiagrGCS	x-direction displacement values at the points	(nElem * nPoints * nLC)
UydiagrGCS	y-direction displacement values at the points	(nElem * nPoints * nLC)
UzdiagrGCS	z-direction displacement values at the points	(nElem * nPoints * nLC)
dUxdiagrGCSdx	x-direction displacement values derivatives	(nElem * nPoints * nLC * nVar)
dUydiagrGCSdx	y-direction displacement values derivatives	(nElem * nPoints * nLC * nVar)
dUzdiagrGCSdx	z-direction displacement values derivatives	(nElem * nPoints * nLC * nVar)

See also PLOTDISP, DISP_TRUSS, DISP_BEAM.

unselectdof

UNSELECTDOF Unselect degrees of freedom.

```
L=unselectdof(DOF,seldof)
[L,I]=unselectdof(DOF,seldof)
creates the matrix to unselect degrees of freedom from the global degrees of
freedom.
```

DOF	Degrees of freedom (nDOF * 1)
seldof	Unselected dof labels (ndof * 1)
L	Selection matrix ((nDOF-ndof) * nDOF)
I	Index vector ((nDOF-ndof) * 1)

See also SELECTDOF.

clearpage

volume_beam

VOLUME_BEAM Compute the volume of a beam element.

V = VOLUME_BEAM(Node,Section) computes the volume of a two-node beam element.

[V,dVdx] = VOLUME_BEAM(Node,Section,dNodedx,dSectiondx) computes the volume of a two node beam element, as well as the derivatives of the volume with respect to the design variables x.

Node	Node definitions	[x y z] (3 * 3)
Sections	Section definitions	[SecID SecProp1 SecProp2 ...]
dNodedx	Node definitions derivatives	(SIZE(Node) * nVar)
dSectionsdx	Section definitions derivatives	(SIZE(Section) * nVar)
V	Element volume	(1 * 1)
dVdx	Element volume derivatives	(nVar * 1)

See also ELEMVOLUMES, VOLUME_TRUSS, ELEMSIZES, SIZE_BEAM.

clearpage

volume_truss

VOLUME_TRUSS Compute the volume of a truss element.

V = VOLUME_TRUSS(Node,Section) computes the volume of a two-node truss element.

[V,dVdx] = VOLUME_TRUSS(Node,Section,dNodedx,dSectiondx) computes the volume of a two node truss element, as well as the derivatives of the volume with respect to the design variables x.

Node	Node definitions	[x y z] (3 * 3)
------	------------------	-----------------

Sections	Section definitions	[SecID SecProp1 SecProp2 ...]
dNodedx	Node definitions derivatives	(SIZE(Node) * nVar)
dSectionsdx	Section definitions derivatives	(SIZE(Section) * nVar)
V	Element volume	(1 * 1)
dVdx	Element volume derivatives	(nVar * 1)

See also ELEMVOLUMES, VOLUME_BEAM, ELEMSIZES, SIZE_TRUSS.

vtrans_solid

VTRANS_SOLID Transformation matrix for stress and strain components in matrix (Voigt) notation.

```
[theta] = vtrans_solid(t)
[theta] = vtrans_solid(t,vtype)
computes the transformation matrix between the local and the global
coordinate system for t stress or strain vector in matrix notation.
```

t	Transformation matrix	(3 * 3)
vtype	Vector type 'stress' (default) 'strain'	
theta	Stress transformation matrix	(6 * 6)

See also TRANS_SOLID8, TRANS_SOLID20.

wilson

WILSON Direct time integration for dynamic systems - Wilson-theta method

`[u,v,a,t] = WILSON(M,C,K,dt,p,u1,v1,[alpha delta theta])` applies the Wilson-theta method for the calculation of the nodal displacements `u`, velocities `v` and accelerations `a` of the dynamic system with the system matrices `M`, `C` and `K` due to the excitation `p`.

`M` Mass matrix (`nDof * nDof`)
`C` Damping matrix (`nDof * nDof`)
`K` Stiffness matrix (`nDof * nDof`)
`dt` Time step of the integration scheme (`1 * 1`). Should be small enough to ensure the stability and the precision of the integration scheme.
`p` Excitation (`nDof * N`). `p(:,k)` corresponds to time point `t(k)`.
`u1` Displacements at time point `t(1)` (`nDof * 1`). Defaults to zero.
`v1` Velocities at time point `t(1)` (`nDof * 1`). Defaults to zero.
`u` Displacements (`nDof * N`). `u(:,k)` corresponds to time point `t(k)`.
`t` Time axis (`1 * N`), defined as `t = [0:N-1] * dt`.

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