

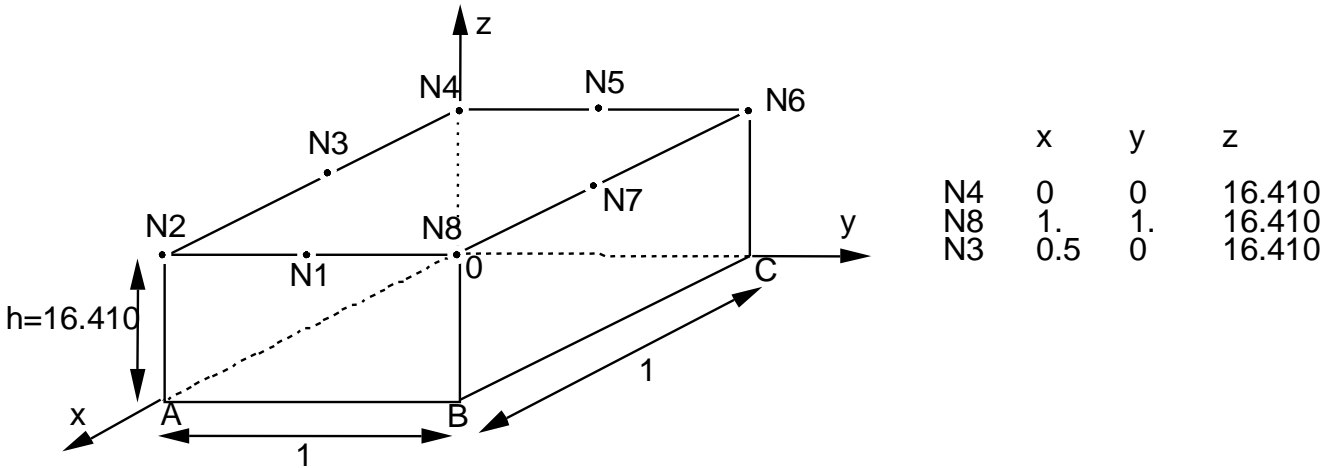
HPLV101 - Homogenization of a Summarized homogeneous

material:

This test tests, in a commonplace situation where the material is homogeneous, the resolution of the thermal and mechanical problems steady, with loadings corresponding to a variation in temperature and an imposed strain, close to those corresponding to the elementary problems of the method of periodic homogenization.

1 Problem of reference

1.1 Geometry



1.2 Material properties

Modelization A

$$E = 1.0 \text{ MPa}$$

$$\nu = 0.3$$

$$k = 1.0 \text{ W l(m. } ^\circ\text{C)}$$

$$C_p = 0 \text{ J l(} ^\circ\text{C.m}^3\text{)}$$

Modelization B

$$E_L = 1.0 \text{ MPa}$$

$$E_T = 0.9 \text{ MPa}$$

$$E_N = 0.8 \text{ MPa}$$

$$\nu_{LT} = 0.1$$

$$\nu_{LN} = 0.25$$

$$\nu_{TN} = 0.3333333$$

$$k = 1.0 \text{ W l(m. } ^\circ\text{C)}$$

$$C_p = 0 \text{ J l(} ^\circ\text{C.m}^3\text{)}$$

1.3 Boundary conditions and loadings

- Mechanics 3D:
 - Plane $z=0$: $dz=0$ for the membrane loading;
 $dx=0, dy=0$ for the loading of cross-bending
 - $y=0$ $y=1$: $dy=0$
 - Planes $x=0$ $x=1$ $dx=0$
 - :
 - The node is outside $dz=0$ (for the only loading of bending)
the field of definition
with a right profile of
the EXCLU type
node: O

Loading:

$$\begin{aligned} \text{déformation membranaire: } E &= \begin{pmatrix} -1 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix} \\ \text{flexion uniforme imposée: } E &= \begin{pmatrix} \frac{z}{z_0} & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix} \end{aligned}$$

$$z_0 = 1 \text{ m}$$

- Mechanics 2D, plane stresses:

Center: $x=0$ $dx=0$ (these conditions do not correspond to the application of the method of homogenization).
The node is outside the field of definition with a right profile of the EXCLU type
node: O $dy=0$

Loading: thermal strain $E = \begin{pmatrix} -1 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix}$ uniforme imposée

- 3D and 2D:

Plane $x=0$ $temp=0$ (this condition does not correspond to the application of the method of homogenization).

Loading: imposed $G = (-1,0,0)$ uniform gradient.

2 Reference solution

2.1 Method of calculating used for the reference solution

- In thermal: the steady thermal problem is solved:

$$\int_{\Omega} \nabla T \cdot \mathbf{K} \cdot \nabla \theta = \int_{\Omega} \mathbf{G} \cdot \mathbf{K} \cdot \nabla \theta, \quad \forall \theta \in V, \quad \text{avec } \mathbf{G} = \begin{pmatrix} -1 \\ 0 \\ 0 \end{pmatrix}$$

Note:

The boundary conditions chosen here are not those necessary to the method of homogenization: one would find indeed $T = 0$ everywhere.

The solution is then (checking the conditions defined in [§1.3]): $T(x, y, z) = -x$

Potential energy is then with the equilibrium: $W^{th} = -\frac{1}{2} \int_{\Omega} \nabla T \cdot \mathbf{K} \cdot \nabla T = -\frac{1}{2} |\Omega|$ ici

- In mechanics: one solves the problem of elastostatic:

$$\int_{\Omega} \boldsymbol{\varepsilon}(\mathbf{u}) \cdot \mathbf{A} \cdot \boldsymbol{\varepsilon}(\mathbf{v}) = \int_{\Omega} \mathbf{E} \cdot \mathbf{A} \cdot \boldsymbol{\varepsilon}(\mathbf{v}), \quad \forall \mathbf{v} \in W,$$

for the cases:

loading 3D membrane	loading 3D of bending	loading 2D plane stresses
$\mathbf{E} = \begin{pmatrix} -1 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix}$	$\mathbf{E} = \begin{pmatrix} z & 0 & 0 \\ z_0 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix}$	$\mathbf{E} = \begin{pmatrix} -1 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix}$

the solutions are:

- in 3D, membrane **loading** and isotropic elasticity: $\mathbf{u}(x, y, z) = \left(0, 0, -\frac{\nu z}{(1-\nu)} \right)$

potential energy with the equilibrium is:

$$W^{pot} = -\frac{1}{2} \int_{\Omega} \boldsymbol{\varepsilon}(\mathbf{u}) \cdot \mathbf{A} \cdot \boldsymbol{\varepsilon}(\mathbf{u}) = -\left(\frac{\nu}{1-\nu} \right)^2 \cdot \frac{|\Omega|}{2} (\lambda + 2\mu)$$

- in 3D, membrane **loading** and orthotropic elasticity: $\mathbf{u}(x, y, z) = (0, 0, -\beta z)$ with $\beta = \frac{C_{13}}{C_{33}}$ is

$$\beta = \frac{\frac{E_N}{E_L} (\nu_{LN} + \nu_{LT} \nu_{TN})}{1 - \frac{E_N}{E_T} \nu_{TN}^2}$$

because the local coordinate system is not confused with the total

reference (nautical angles being worth all 90°).

$$W^{pot} = -\frac{1}{2} \int_{\Omega} (\mathbf{u}) \cdot \mathbf{A} \cdot \boldsymbol{\varepsilon}(\mathbf{u}) = -\beta^2 \cdot \frac{|\Omega|}{2} C_{33}$$

$$\text{where } C_{33} = \frac{1 - \nu_{NT}\nu_{TN}}{E_T E_N \Delta} \text{ and } \Delta = \frac{1 - \nu_{LT}\nu_{TL} - \nu_{LN}\nu_{NL} - \nu_{TN}\nu_{NT} - 2\nu_{LT}\nu_{TN}\nu_{NL}}{E_L E_T E_N}$$

- in 3D, loading of **bending** : $\mathbf{u}(x, y, z) = \left(0, 0, \frac{+ \nu z^2}{2z_0(1-\nu)} \right)$;

$$\bar{W}^{pot} = - \left(\frac{\nu}{1-\nu} \right)^2 \cdot \frac{|\Omega|}{2} (\lambda + 2\mu) \cdot \frac{h^2}{3z_0^2}$$

- in 3D, loading of **bending** and orthotropic elasticity: $\mathbf{u}(x, y, z) = \left(0, 0, -\beta \frac{z}{L} \right)$ with $\beta = \frac{C_{13}}{C_{33}}$ is

$$\beta = \frac{\frac{E_N}{E_L} (\nu_{LN} + \nu_{LT}\nu_{TN})}{1 - \frac{E_N}{E_T} \nu_{TN}^2} \text{ because the local coordinate system is not confused with the total}$$

reference (nautical angles being worth all 90°).

$$\bar{W}^{pot} = - \frac{1}{2} \int_{\Omega} \boldsymbol{\varepsilon}(\mathbf{u}) \cdot \mathbf{A} \cdot \boldsymbol{\varepsilon}(\mathbf{u}) = -\beta^2 \cdot \frac{|\Omega|}{2} C_{33} \frac{h^2}{3z_0^2}$$

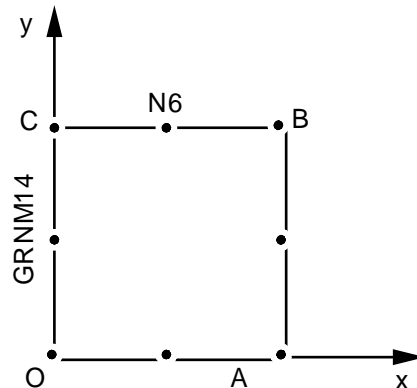
$$\text{where } C_{33} = \frac{1 - \nu_{NT}\nu_{TN}}{E_T E_N \Delta} \text{ and } \Delta = \frac{1 - \nu_{LT}\nu_{TL} - \nu_{LN}\nu_{NL} - \nu_{TN}\nu_{NT} - 2\nu_{LT}\nu_{TN}\nu_{NL}}{E_L E_T E_N}$$

- in 2D, plane loading: $\mathbf{u}(x, y) = (-x, 0)$;

$$\bar{W}^{pot} = \frac{-|\Omega|}{2(1-\nu^2)}$$

3 Modelization A

3.1 Characteristic of the modelization



Boundary conditions and loading:

Thermal: GROUP_NO: GRNM14: TEMP: 0.0
 PRE_GRAD_TEMP: FLUX_X: - 1.0

Mechanics: GROUP_NO: GRNM14: DX: 0.0
(plane stresses)

THE NODE IS OUTSIDE THE FIELD OF DEFINITION WITH A
RIGHT PROFILE OF THE EXCLU TYPE NODE: O
DY: 0.0
PRE_EPSI: EPXX: - 1.0

3.2 Characteristics of the mesh

Many nodes: 8

Number of meshes and types: 1 QUAD8

3.3 Values tested

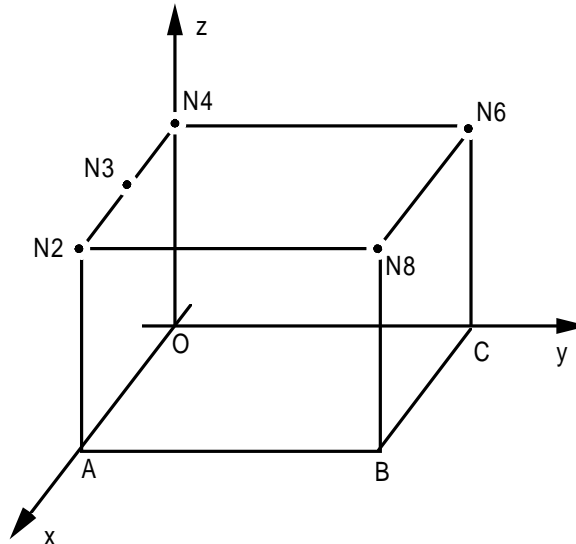
Not	Quantity	Reference
<i>CMP</i>		
<i>A</i>	<i>TEMP</i>	- 1.0000
<i>A</i>	<i>DX</i>	- 1.0000
<i>N6</i>	<i>DX</i>	- 0.5000
Mesh	Potential energy	equilibrium
with the Thermal		
<i>MI</i>	Reference	- 0.500000000
<i>MI</i>	Mechanics	- 0.549450550

3.4 Remarks

Code_Aster provides the value of strain energy, equal contrary to potential energy to the equilibrium (elastic case).

4 Modelization B

4.1 Characteristic of the modelization



Name of meshes of the sides:	ZEGALO	YEGALO	YEGAL1	XEGALO	XEGAL1
Summits:	BCOA	OAN2N4	BCN6N8	CON4N6	ABN8N2

Boundary conditions:

```

ZERO: DEFI_CONSTANTE (VALE: 0.0) ;
FCT1:  DEFI_FONCTION  (Nom_para:  "Z",  VALE:  (0.0
0.0.1.0.1.0));
Thermal:
GROUP_NO: XEGAL0: TEMP: 0.0
PRE_GRAD_TEMP: FLUX_X: -1.0
Mechanics:
GROUP_NO:  YEGALO: DY = 0.0
           XEGAL1: DX = 0.0
           YEGAL1: DY = 0.0
           XEGALO: Membrane DZ =
0.0 Case:
GROUP_NO:  ZEGALO: DZ = 0.0
PRE_EPSI: EPXX: -1.0
Case bending:
GROUP_NO:  ZEGALO: DX = ZERO,  DY = ZERO
THE NODE IS OUTSIDE THE FIELD OF DEFINITION WITH A RIGHT
PROFILE OF THE EXCLU TYPE NODE: 0
DZ = ZERO
PRE_EPSI:                                     EPXX: FCT1
    
```

4.2 Characteristic of the mesh

Many nodes: 20
Number of meshes and types: 1 HEXA20

4.3 Values tested

In isotropic elasticity

Case	Thermal	Quantit y	Not
Reference	<i>temp</i>	<i>N8</i>	- 1.000000
	<i>temp</i>	<i>N3</i>	- 0.500000
Mechanics membrane	<i>dz</i>	<i>N4</i>	- 7.03285714
	<i>dz</i>	<i>N8</i>	- 7.03285714
Mechanics bending	<i>dz</i>	<i>N4</i>	57.70459285
	<i>dz</i>	<i>N8</i>	57.70459285

Nets	Potential energy	equilibrium
with the Thermal		
<i>MI</i>	Reference	- 8.20500
<i>MI</i>	Mechanical	
	Membrane	- 2.0287088
	Bending	- 1.8210238 102

In orthotropic elasticity

Case	Thermal	Quantit y	Not
Reference	<i>temp</i>	<i>N8</i>	- 1.000000
	<i>temp</i>	<i>N3</i>	- 0.500000
Mechanics membrane	<i>dz</i>	<i>N4</i>	- 6.63044894
	<i>dz</i>	<i>N8</i>	- 6.63044894
Mechanics bending	<i>dz</i>	<i>N4</i>	54.40283358
	<i>dz</i>	<i>N8</i>	54.40283358

Nets	Potential energy	equilibrium
with the Thermal		
<i>MI</i>	Reference	- 8.20500

5 Summary of the results

the results are exact with errors rounding close, since the sought solutions are part of space of the finite elements selected for the modelization.