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以C3D8R为例介绍如何在VUMAT中得到单元编号

Taking C3D8R as an example to introduce how to obtain element numbers in **VUMAT**



VUMAT不像UMAT直接给出了单元的编号, VUMAT需要借助一个子程序vumatXtrArg 得到,这个方法最早发布 在如下链接中

Unlike UMAT, VUMAT does not directly provide the element number. VUMAT needs to use a subroutine called vumatXtrArg to obtain it, and this method was first published in the following link

https://polymerfem.com/community/constitutive-models/nblock-in-vumat/

相信有不少朋友需要,摘录如下 I believe many friends will need it, here is the excerpt

You are using C3D8R elements which means continuum 3D 8 noded reduced integration. A reduced integration of this kind has a single Gauss pt.

Thats why you have matching GP and element numbers.

Regarding the second, apply the following cover,

```
subroutine vumat (
c Read only -
1 jblock, ndir, nshr, nstatev, nfieldv, nprops, lanneal,
2 stepTime, totalTime, dt, cmname, coordMp, charLength,
3 props, density, strainInc, relSpinInc,
4 tempOld, stretchOld, defgradOld, fieldOld,
5 stressOld, stateOld, enerInternOld, enerInelasOld,
6 tempNew, stretchNew, defgradNew, fieldNew,
c Write only -
7 stressNew, stateNew, enerInternNew, enerInelasNew)
C
include vaba param.inc
```

dimension jblock(*), props(nprops), density(*), coordMp(*),

```
dimension jblock(*), properties(nprops), density(*), coordinateMp(*),
1 charLength(*), strainInc(*),
1 characteristic length(*), strain increment(*),
2 relSpinInc(*), tempOld(*),
3 stretchOld(*), 3 previous stretch(*),
4 defgradOld(*), 4 previous deformation gradient(*),
5 fieldOld(*), stressOld(*),
6 stateOld(*), enerInternOld(*),
7 enerlnelasOld(*), tempNew(*),
8 stretchNew(*),
9 defgradNew(*),
1 fieldNew(*),
2 stressNew(*), stateNew(*),
3 enerInternNew(*), enerInelasNew(*)
C
character*80 cmname
parameter (
1 i umt nblock = 1,
2 i umt npt = 2,
3 i_umt_layer = 3,
4 i umt kspt = 4,
5 i umt noel = 5)
call vumatXtrArg (jblock(i umt nblock),
1 ndir, nshr, nstatev, nfieldv, nprops, lanneal,
2 stepTime, totalTime, dt, cmname, coordMp, charLength,
3 props, density, strainInc, relSpinInc,
4 tempOld, stretchOld, defgradOld, fieldOld,
5 stressOld, stateOld, enerInternOld, enerInelasOld,
6 tempNew, stretchNew, defgradNew, fieldNew,
7 stressNew, stateNew, enerInternNew, enerInelasNew,
stressNew, stateNew, enerInternNew, enerInelasNew,
8 jblock(i umt noel), jblock(i umt npt),
jblock(i_umt_noel), jblock(i_umt_npt),
9 jblock(i umt layer), jblock(i umt kspt))
jblock(i umt layer), jblock(i umt kspt))
return
end
C -----C
subroutine vumatXtrArg (
c read only -
1 nblock, ndir, nshr, nstatev, nfieldv, nprops, lanneal,
```

- 2 stepTime, totalTime, timeinc, cmname, coordMp, charLength,
- 3 props, density, strainInc, relSpinInc,
- 4 tempOld, stretchOld, defgradOld, fieldOld,
- 3 stressOld, stateOld, enerInternOld, enerInelasOld,
- 6 tempNew, stretchNew, defgradNew, fieldNew,
- c write only -
- 5 stressNew, stateNew, enerInternNew, enerInelasNew,
- c read only extra arguments -
- 6 nElement, nMatPoint, nLayer, nSecPoint)

include vaba_param.inc

C

c all arrays dimensioned by (*) are not used in this algorithm

dimension props(nprops), density(nblock),

- 1 strainInc(nblock,ndir+nshr),
- 2 relSpinInc(nblock,nshr), defgradOld(nblock,9),
- 4 stressOld(nblock,ndir+nshr),
- 5 stateOld(nblock,nstatev), enerInternOld(nblock),
- 6 enerlnelasOld(nblock),
- 7 stretchNew(nblock,ndir+nshr), defgradNew(nblock,9),
- 8 stressNew(nblock,ndir+nshr)
- 8 stressNew(nblock, ndir+nshr)

dimension enerInelasNew(nblock), stateNew(nblock, nstatev),

dimension enerlnelasNew(nblock), stateNew(nblock,nstatev),

1 enerInternNew(nblock)

dimension nElement(nblock),nMatPoint(nblock),nLayer(nblock),

dimension nElement(nblock), nMatPoint(nblock), nLayer(nblock),

1 nSecPoint(nblock)

character*80 cmname

c make computations for every material point

C ------

do 1000 nblck = 1,nblock

Nelem=nElement(nblck)! here we go 😀

blah blah blah

enddo

return

end

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