Norton behaviour description

file: Norton.mfront
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date: 23 / 11 / 06

This viscoplastic behaviour is fully determined by the evolution of the equivalent viscoplastic strain p as a function f of the Von Mises stress $\sigma_{\rm eq}$:

$$\dot{p} = f\left(\sigma_{\rm eq}\right) = A\,\sigma_{\rm eq}^E$$

where:

- \bullet A is the Norton coefficient .
- \bullet E is the Norton exponent .

A and E are declared as material properties .

Source code

```
OParser IsotropicMisesCreep;
   @Behaviour Norton;
   @Author Helfer Thomas;
   @Date 23/11/06;
   @Description{
     This viscoplastic behaviour is fully determined by the evolution
     of the equivalent viscoplastic strain "\(p\)" as a function "\(f\)"
     of the Von Mises stress "\(\sigmaeq\)":
     "\dot{p}=f\paren{\sigmaeq}=A\,\sigmaeq^{E}"
10
     "\]"
     where:
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13
     - "\(A\)" is the Norton coefficient.
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     - "\(E\)" is the Norton exponent.
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16
     "\(A\)" and "\(E\)" are declared as material properties.
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19
   @UMATFiniteStrainStrategies[umat] {None,LogarithmicStrain1D};
20
21
   //! The Norton coefficient
   OMaterialProperty real A;
```

```
A.setEntryName("NortonCoefficient");
   //! The Norton coefficient
26
   @MaterialProperty real E;
   E.setEntryName("NortonExponent");
28
   @FlowRule{
30
    /*!
      * The return-mapping algorithm used to integrate this behaviour
32
      * requires the definition of (f) and (\langle f \rangle_{sigmaeq}) (see
      * Osimo_computational_1998 and Ohelfer_generateur_2013 for
      * details).
36
      * We introduce an auxiliary variable called `tmp` to
      * limit the number of call to the 'pow' function
     const real tmp = A*pow(seq,E-1);
40
             = tmp*seq;
41
     df_dseq = E*tmp;
43
```