

Non-linear analysis – Advanced topics



Code_Aster, Salome-Meca course material

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Outline

- ▶ The `DEFI_LIST_INST` command
- ▶ Multi-step analysis (reuse and initial state)
- ▶ Advanced algorithms controls
- ▶ Real-time monitoring

The `DEFI_LIST_INST` command:

More control on non-linear analysis

More control on non-linear analysis

▶ When to use `DEFI_LIST_INST` ?

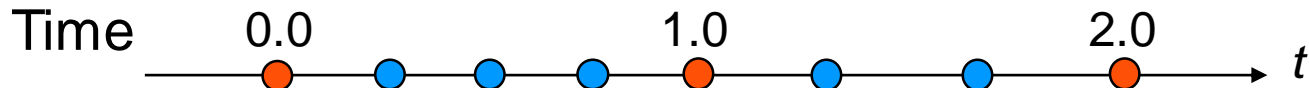
- ▶ For automatic time-step control
- ▶ Event-driven system

The `DEFI_LIST_INST` command:

Advanced time-step control

Advanced time-step control

- ▶ Non-linear problem is parametrized by time – Basic



```
L_REAL=DEFI_LIST_REEL( DEBUT=0.0,  
                       INTERVALLE=(  
                           _F(JUSQU_A=1.0,NOMBRE=4, ),  
                           _F(JUSQU_A=2.0,NOMBRE=3, ),  
                       ), )
```

```
RESUN = STAT_NON_LINE(...  
                      INCREMENT=_F(LIST_INST=L_REAL)  
                      ...)
```



Deprecated in V13.3 => use DEFI_LIST_INST instead !

Advanced time-step control

- ▶ Non-linear problem is parametrized by time – Advanced



Initial time list defined by `DEFI_LIST_REEL`

```
L_REAL = DEFI_LIST_REEL(  
    DEBUT=0.0,  
    INTERVALLE=( _F(JUSQU_A=2.0,NOMBRE=1,) , ) , )
```

```
DEFLIST = DEFI_LIST_INST(  
    DEFI_LIST = _F(LIST_INST = L_REAL,) , )
```

```
RESUN = STAT_NON_LINE(...  
    INCREMENT=_F(LIST_INST= DEFLIST)  
    ...)
```

Advanced time-step control

▶ Time-step control – Manual mode:

- From user's defined list of values: `DEFI_LIST_REEL`

```
DEFLIST = DEFI_LIST_INST (  
    DEFI_LIST = _F (MODE = 'MANUEL' ,  
        LIST_INST = L_REAL ,) ,)
```

- From user's defined list of values: Python's list in `VALE`

```
DEFLIST = DEFI_LIST_INST (  
    DEFI_LIST = _F (MODE = 'MANUEL' ,  
        VALE = (1. , 2. , 3. ,) ,)
```


Advanced time-step control

▶ Time-step control – Manual mode

▶ From previous computation:

- First computation with automatic mode for instance => results RESU1
- Second computation with previous list, divided by two for instance (SUBD_PAS=2)

```
DEFLIST = DEFI_LIST_INST(  
    DEFI_LIST = _F(MODE = 'MANUEL' ,  
                  RESULTAT = RESU1 ,  
                  SUBD_PAS = 2) ,)
```

Advanced time-step control

▶ Time-step control – Automatic mode

- ▶ Definition of global parameters: minimum and maximum step, maximum number of steps

```
DEFLIST = DEFI_LIST_INST (  
    DEFI_LIST = _F (MODE = 'AUTO' ,  
        PAS_MINI = 1.E-3 ,  
        PAS_MAXI = 0.1  
        NB_PAS_MAXI = 1000.) ,)
```



- ▶ PAS_MINI reached => stop
- ▶ PAS_MAXI reached => time step is capped

Advanced time-step control

▶ Time-step control – Automatic mode

- ▶ Definition of automatic adaptation schemes in `ADAPTATION` keyword: how to compute next time step ?

$$\Delta t^{i+1} = c. \Delta t^i ?$$

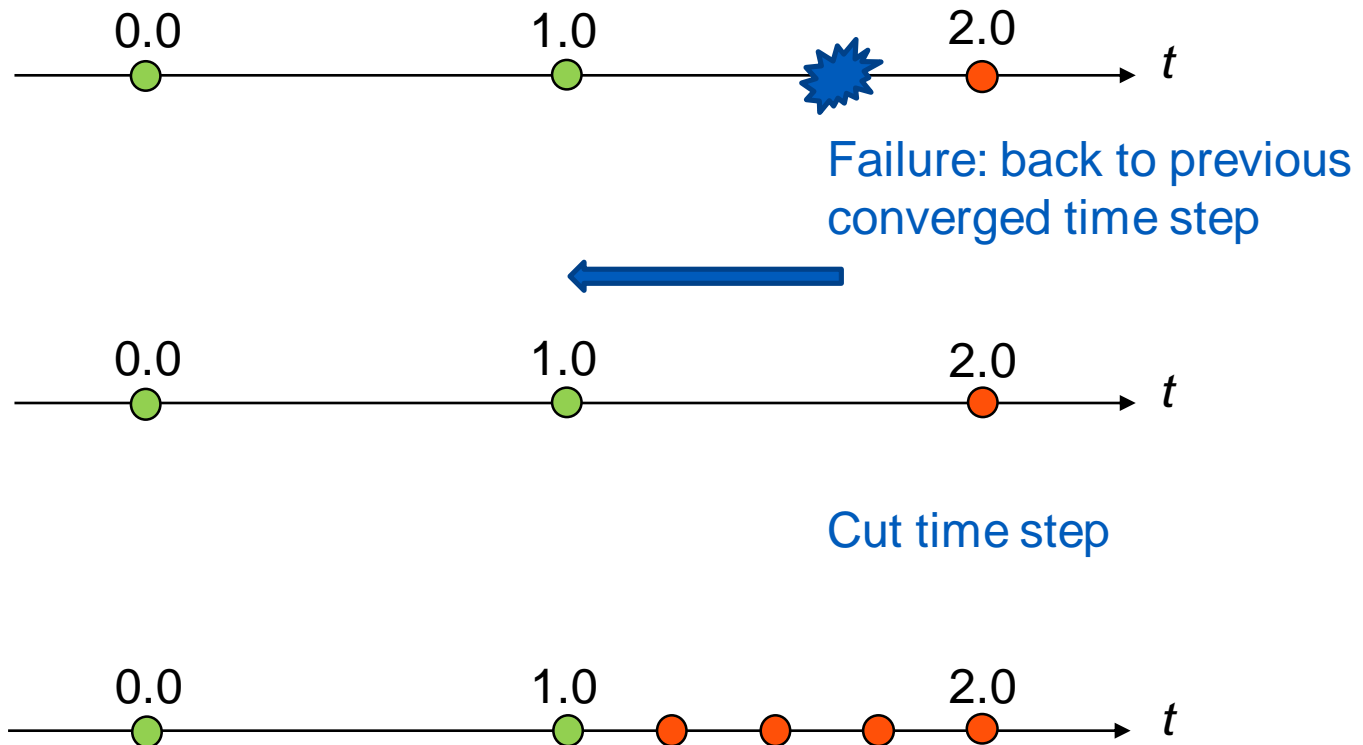
- ▶ **Event-driven system: `EVENEMENT+MODE_CALCUL_TPLUS`**

```
DEFLIST = DEFI_LIST_INST (  
    DEFI_LIST = _F (MODE = 'AUTO' , ...) ,  
    ADAPTATION = _F (EVENEMENT = 'SEUIL' ,  
        MODE_CALCUL_TPLUS = 'FIXE' ,  
        PCENT_AUGM = 50. , ) , )
```

Advanced time-step control

▶ Time-step control – Manage step time

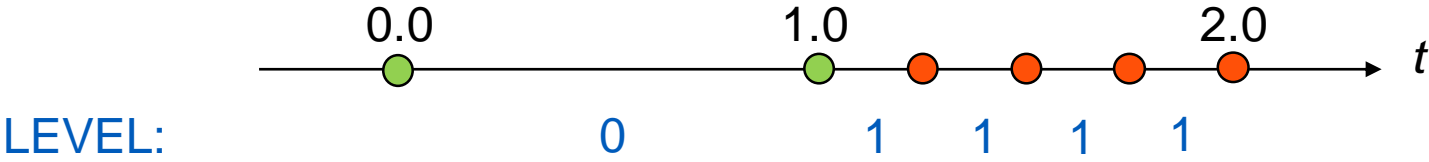
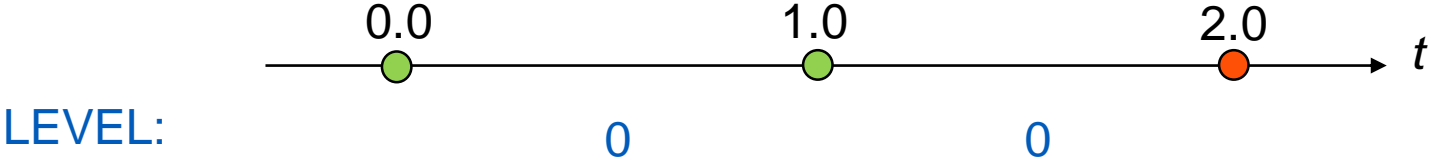
- ▶ **Event-driven system: EVENEMENT= 'ERREUR' +ACTION= 'DECOUPE**




Advanced time-step control

▶ Time-step control – The LEVEL

▶ Level of time step: initial from `DEFI_LIST_REEL` is **LEVEL 0**



 LEVEL = NIVEAU

LEVEL is recursive ...

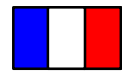
The `DEFI_LIST_INST` command:

Events management

Events management

▶ General event mechanism

- ▶ Event-driven is how to control algorithm: stop computation, manage step time, other actions
- ▶ Direct event-driven: keywords **EVENEMENT** and **ACTION**
- ▶ Time step adaptation event-driven: keywords **ADAPTATION** and **MODE_CALCUL_TPLUS**



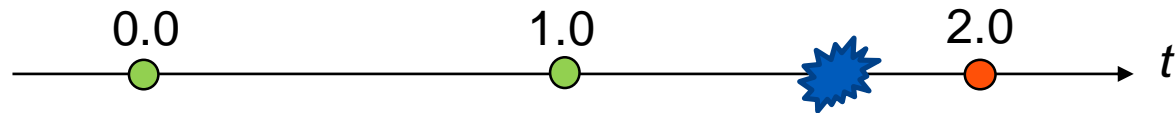
EVENT = **EVENEMENT**

ACTION = **ACTION**

EVENT+ACTION = **ECHEC**

Events management

► The **EVENT** and the **ACTION**: an example for step time management



EVENT: ERROR (ex.: failure in constitutive equation solving, too many Newton iterations, etc)



ACTION: Cut time step

Events management

◆ List of the EVENEMENT for general event management

- ◆ ERREUR: error during constitutive equation solving, too many Newton iterations, contact error, singular matrix
- ◆ DIVE_RESI: residual has increased twice during Newton iterations
- ◆ DELTA_GRANDEUR: the increase of a given quantity has exceeded a value defined by the user (ex: cumulative plastic strain)
- ◆ COLLISION: contact occurs for the first time
- ◆ INTERPENETRATION: contact interpenetration (for penalty methods)
- ◆ INSTABILITE: detection of an instability (CRIT_STAB operator)

Some of these events have parameters (see U4.34.04). Example:

```
_F (EVENEMENT= 'DELTA_GRANDEUR',  
    VALE_REF = 0.1e-2,  
    NOM_CHAM = 'VARI_ELGA',  
    NOM_CMP  = 'V1'),
```

For
VMIS_ISOT_LINE
 $\Delta \varepsilon_{eq}^p > 0.1\%$

Events management

◆ List of ACTION for general event management

- ◆ ARRET: stop computation (previous converged steps are saved)
- ◆ DECOUPE: cut time step
- ◆ ITER_SUPPL: allow more Newton iterations (more than ITER_GLOB_MAXI parameter)
- ◆ ADAPT_COEF_PENA: adaptation of penalty parameter for contact
- ◆ CONTINUE: continue computation

Some of these actions have parameters (see U4.34.04).

Events management

◆ List of the EVENEMENT for AUTO time stepping

- ◆ SEUIL: defining the trigger for adapting the time step in automatic mode
- ◆ TOUT_INST: adaptation occurs at every step time

Some of these events have parameters (see U4.34.04). Example:

```
_F (EVENEMENT= 'SEUIL',  
    NB_INCR_SEUIL=2,  
    NOM_PARA='NB_ITER_NEWTON',  
    CRIT_COMP='LE',  
    VALE_I=5),
```

■ ■ If two successive time steps use less than 5 Newton iterations

Events management

◆ List of the `MODE_CALCUL_TPLUS` for AUTO time stepping

What is c in $\Delta t = c \cdot \Delta t^i$?

- ◆ `FIXE`: c is constant (`PCENT_AUGM` parameter)
- ◆ `DELTA_GRANDEUR`: c is defined by the user from quantities computed (displacement, stress...)
- ◆ `IMPLEX` : c is automatically defined for `IMPLEX`
- ◆ `ITER_NEWTON` : c is defined by the variation of the number of Newton iterations during time stepping

Some of these actions have parameters (see U4.34.04).

The `DEFI_LIST_INST` command:

Summary

The `DEFI_LIST_INST` command – Summary

- ◆ By default: at least one event (`=ERREUR`) with its action (`=DECOUPE`)
- ◆ You can define more than one `ECHEC` (event and its action)
- ◆ You can define more than one `ADAPTATION` (the smallest time step time is selected)
- ◆ Advanced users may define special heuristics for time adaptation or event-driven => these may then be integrated in the official version of Code_Aster

Multi-step analysis

Multi-step analysis

◆ What can I change between two non-linear computations ?

- ◆ **Model** from `AFFE_MODELE`
 - ◆ Pay attention to the initial state if you change the model !
- ◆ **Loads** from `AFFE_CHAR_*` and `EXCIT`
- ◆ **Elementary characteristics** from `AFFE_CARA_ELEM`
- ◆ **Mechanical behaviours** from `COMPORTEMENT` keyword:
 - ◆ From `ELASTIC` to non-linear behavior laws => OK !
 - ◆ From a non-linear behavior laws to another => Not OK (because of internal state variables)

You cannot change the mesh from one computation to another :

- ➔ you have to project results (use `PROJ_CHAMP` with `ECLA_PG` for Gauss fields)
- ➔ you have to re-define everything (models, loads, ...)

Note : for excavation or layer-based structures, see U2.04.06

Multi-step analysis

- ▶ Initial state is defined by:
 - ▶ Displacements
 - ▶ Stresses
 - ▶ Internal state variables
 - ▶ +more specific fields for certain applications such as XFEM or multifiber beam elements
- ▶ By default: initial state is empty (everything is zero)
- ▶ The initial index to access data fields (`NUME_ORDRE`) is always 0 except when the result datastructure was extended (`reuse`) => **CANNOT BE CHANGED BY THE USER**
- ▶ The initial time may be defined by `ETAT_INIT/INST_ETAT_INIT` keyword if necessary
- ▶ Careful ! In `reuse` mode, `INST_ETAT_INIT` must be superior to the previous time step

Multi-step analysis

▶ Defining a specific initial state by `ETAT_INIT` keyword

- ▶ From previous `STAT_NON_LINE:EVOL_NOLI + NUME_ORDRE` or `INST` or `{nothing}`
 - ▶ If `{nothing}` => last computed step in datastructure
- ▶ From individual fields or manually assembled datastructures
 - ▶ Field by field: displacement (`DEPL`), stress (`SIGM`), internal state variables (`VARI`)
 - ▶ Complete datastructure: from `LIRE_RESU` or `CREA_RESU` command

Multi-step analysis

▶ For non-linear behavior laws:

- ▶ The initial stress field is used to update stresses during behavior law integration
- ▶ Depend on laws: **some of them (hyperelastic for large strains) DON'T take into account initial stresses** => you should use internal state variables (`SIMO_MIEHE` for instance)
- ▶ For `GDEF_LOG` strains measure => stresses are not Cauchy but logarithmic ones (see `ssnp159b` for instance)
- ▶ To have initial strains: use `AFFE_CHAR_MECA/PRE_EPSI`
- ▶ To have initial stresses that are not used to update stresses (*e.g. during behavior law integration*): use `AFFE_CHAR_MECA/PRE_SIGM`

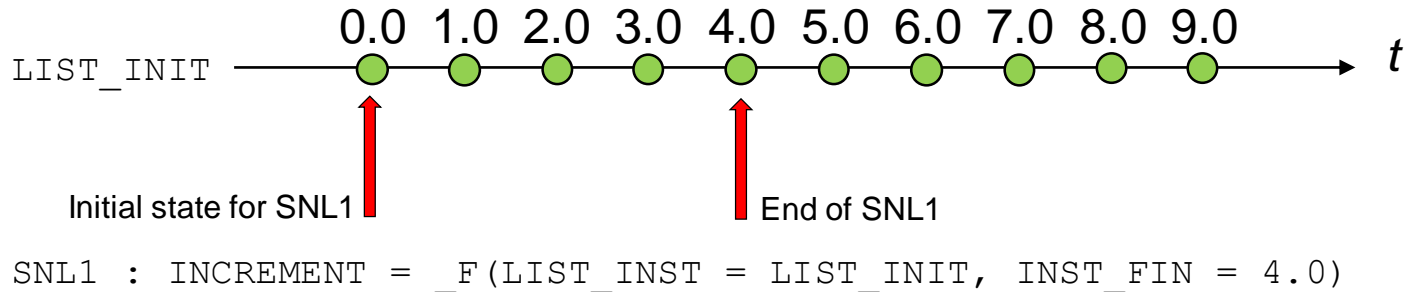
▶ For contact (CONTINUE formulation):

- ▶ Initial contact behavior is activated by default in `DEFI_CONTACT` command (`CONTACT_INIT='INTERPENETRE'`)
- ▶ Friction history is constructed from displacement field (`DEPL`) => careful !

Multi-step analysis

Management of time list:

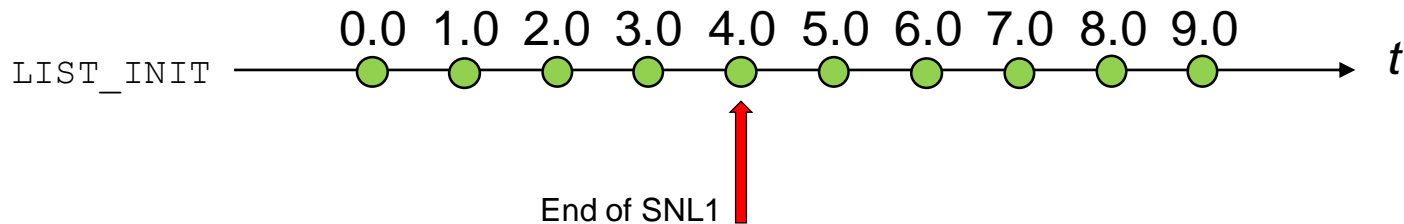
- Defined in `INCREMENT/INST_INCREMENT`
- You can manage the start/end of the time list with `NUME_INST_INIT / INST_INIT / NUME_INST_FIN / INST_FIN`
- To end at `INST=4.0`



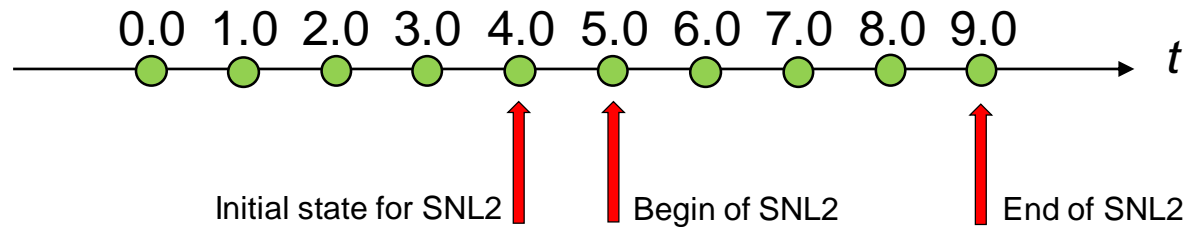
Multi-step analysis

Management of time list:

- « Simple » continuation of a non-linear computation



```
SNL1 : INCREMENT = _F(LIST_INST = LIST_INIT, INST_FIN = 4.0)
```



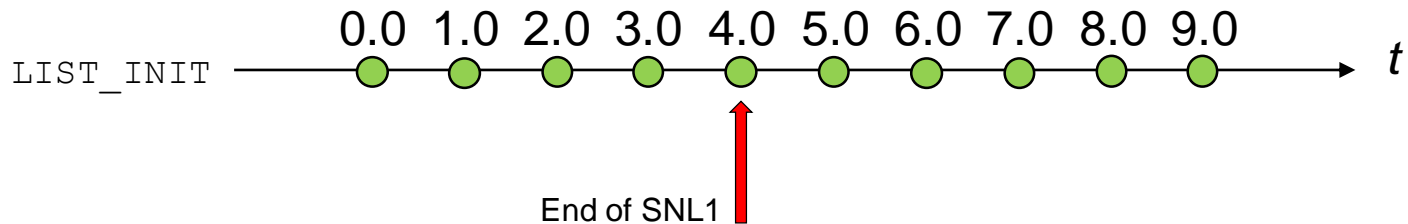
Standard reuse:

```
SNL2 : reuse = SNL1  
      ETAT_INIT = _F(EVOL_NOLI=SNL1),  
      INCREMENT = _F(LIST_INST = LIST_INIT)
```

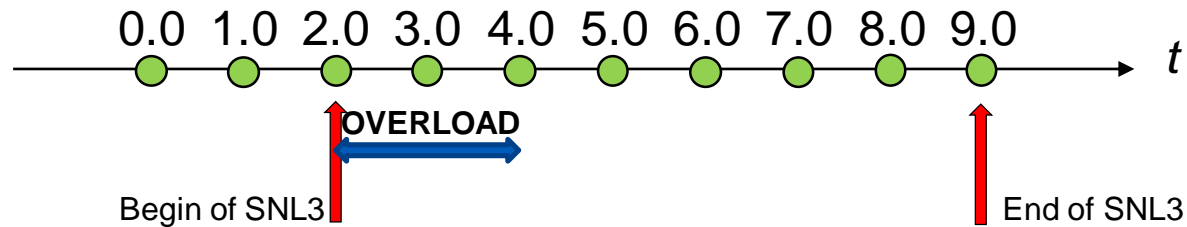
Multi-step analysis

Management of time list:

- Overload a non-linear computation



```
SNL1 : INCREMENT = _F(LIST_INST = LIST_INIT, INST_FIN = 4.0)
```



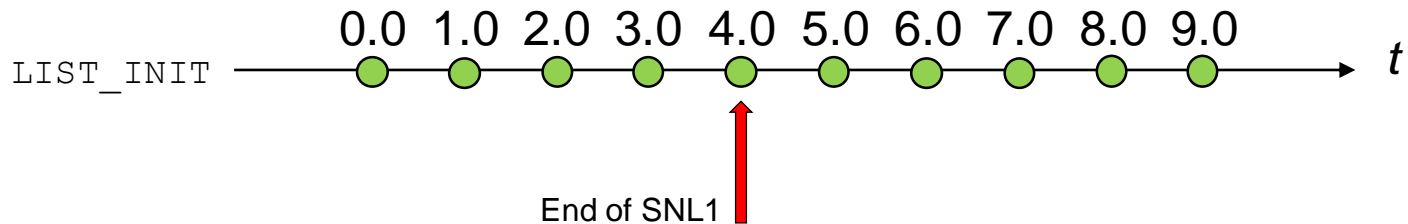
Overload :

```
SNL3 : reuse = SNL1  
      ETAT_INIT = _F(EVOL_NOLI=SNL1),  
      INCREMENT = _F(LIST_INST = LIST_INIT, INST_INIT = 2.0)
```

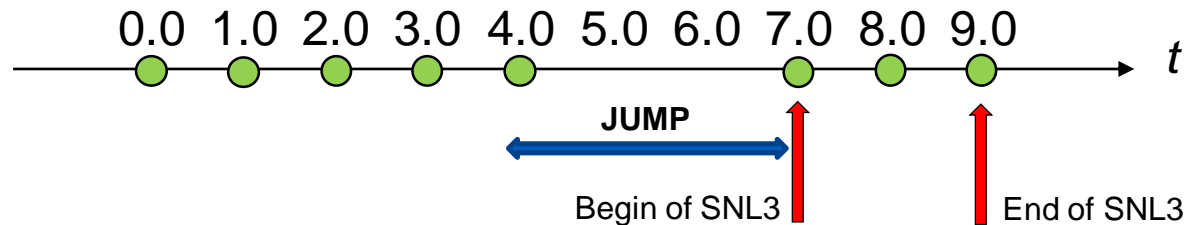
Multi-step analysis

Management of time list:

- Jump over a non-linear computation



```
SNL1 : INCREMENT = _F(LIST_INST = LIST_INIT, INST_FIN = 4.0)
```



Jump :

```
SNL4 : reuse = SNL1  
ETAT_INIT = _F(EVOL_NOLI=SNL1),  
INCREMENT = _F(LIST_INST = LIST_INIT, INST_INIT = 7.0)
```

Advanced algorithm controls

Advanced algorithm controls

◆ Global non-linear solver – Keyword `METHODE`

- ◆ Newton and quasi-Newton method – `METHODE = 'NEWTON'`
- ◆ Implicit/explicit method (Oliver and *al.*) – `METHODE = 'IMPLEX'`
 - ◆ For damaged concrete analysis (to use with special time step adaptation)
- ◆ Newton-Krylov method – `METHODE = 'NEWTON_KRYLOV'`
 - ◆ To use with an iterative linear solver (`PETSC` or `GCPC`), adaptative convergence criterion in the iterative solver → very efficient if a good preconditionner is available

Advanced algorithm controls

◆ Newton solver — Keyword `METHODE= 'NEWTON'`

- ◆ Quasi-Newton options: see `REAC_INCR` and `REAC_ITER` to update matrix
- ◆ Choice of matrix for prediction phase (Euler):
 - ◆ `PREDICTION= 'ELASTIQUE' / 'TANGENTE'` => prefer `PREDICTION= 'ELASTIQUE'` for severe non-linear analysis (especially contact) and unloading
 - ◆ `PREDICTION= 'DEPL_CALCULE'` => useful to run post-critical analysis or damage
- ◆ Choice of matrix for correction phase (Newton):
 - ◆ `MATRICE= 'TANGENTE'` => exact tangent matrix (if available from the constitutive law), should generally be preferred (default)
 - ◆ `MATRICE= 'ELASTIQUE'` => will converge for generalized material (very slowly though)
 - ◆ To swap from `TANGENTE` to `ELASTIQUE` => use `PAS_MINI_ELAS`
 - ◆ For damaged structure with elastic matrix => use `REAC_ITER_ELAS`
 - ◆ Force symmetric matrix with `MATR_RIGI_SYME= 'OUI'`

Advanced algorithm controls

▶ Storage management (ARCHIVAGE keyword)

▶ Default:

- ▶ All time steps are saved with displacements, stresses and internal state variables

▶ User-defined storage:

- ▶ Use `ARCHIVAGE + LIST_INST/INST/PAS_ARCH` to store only part of the results. Initial state and last converged step are **ALWAYS** saved (to allow restart after errors, lack of CPU time for instance)
- ▶ Use `ARCHIVAGE + CHAM_EXCLU` to exclude some fields in the saving process
- ▶ Warning: saving all fields for a lot of time steps and a large model => VERY LARGE MEMORY

Advanced algorithm controls

▶ Output management (`AFFICHAGE` keyword)

- ▶ The convergence table is printed at each time step
- ▶ To save this table in a CSV file => `AFFICHAGE/UNITE`
- ▶ To print only for some steps (useful when there's a lot of time steps) => `AFFICHAGE/PAS`
- ▶ To print where residuals are maximum in convergence table => `AFFICHAGE/INFO_RESIDU='OUI'`
- ▶ To print CPU time for each Newton iteration => `AFFICHAGE/INFO_TEMPS='OUI'`

Advanced algorithm controls

◆ Statistics management (MEASURE keyword)

- ◆ Some statistics (CPU time, number) are printed at each time step and at the end of computation
- ◆ To save this table in a CSV file => MEASURE/UNITE
- ◆ To generate a TABLE in Aster format => MEASURE/TABLE='OUI'

```
RESU = STAT_NON_LINE (  
    MEASURE = _F (TABLE='OUI', UNITE=50), )  
STAT = RECU_TABLE (CO=RESU, NOM_TABLE='STAT', )  
IMPR_TABLE (TABLE=STAT)
```

Real-time monitoring

Real-time monitoring

▶ Extract values for monitoring analysis (OBSERVATION/SUIVI_DDL keywords)

▶ Select field:

- ▶ DEPL, VITE, ACCE, TEMP
- ▶ SIEF_ELGA, VARI_ELGA, EPSI_ELGA
- ▶ CONT_NOEU
- ▶ FORC_NODA, REAC_NODA
- ▶ *_ABSOLU

▶ Select where:

- ▶ TOUT/GROUP_MA/MAILLE for ELEM fields + Gauss point / sub-point
- ▶ TOUT/GROUP_MA/MAILLE/GROUP_NO/NOEUD for NODE fields

▶ Select what:

- ▶ Value of component (ex. DX, DY, SIXX, V1, etc.) or formula (ex. $\sqrt{DX^2+DY^2+DZ^2}$)
- ▶ On selected or global on several (ABS, MAX, MIN, ...)

Real-time monitoring

◆ Save values in TABLE (OBSERVATION keyword)

- ◆ Extract (field, where, what)
- ◆ Which step time: OBSERVATION+LIST_INST/INST/PAS_OBSE
- ◆ Save in TABLE
- ◆ Get from results

```
RESU = STAT_NON_LINE (  
  OBSERVATION = _F (NOM_CHAM='SIEF_ELGA',  
                    EVAL_CMP='VALE',  
                    NOM_CMP='M11',  
                    EVAL_ELGA='MAX',  
                    TOUT='OUI'), )  
  
OBSV = RECU_TABLE (CO=RESU, NOM_TABLE='OBSERVATION', )  
  
IMPR_TABLE (TABLE=OBSV)
```


Real-time monitoring

◆ Print values in message file (SUIVI_DDL keyword)

- ◆ Extract (field, where, what)
- ◆ Print in message file (convergence table)
- ◆ Warning: number of SUIVI_DDL is limited (only 15 columns in convergence table)
- ◆ As is convergence table => save in message file (AFFICHAGE keyword) in real-time

```
RESU = STAT_NON_LINE (  
    SUIVI_DDL= ( _F (NOM_CMP='PRE1',  
                    NOM_CHAM='DEPL',  
                    EVAL_CHAM='MAXI_ABS',  
                    NOEUD= ('NO1'), ), ), )
```

End of presentation

Is something missing or unclear in this document?
Or feeling happy to have read such a clear tutorial?

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