
$$\bar{\Pi} = \frac{1}{2} \sum_e \{u\}^T \cdot [K] \cdot \{u\} - \{u\}^T \cdot \{F\}$$

FEM SOFTWARE AND SERVICES



Die große Elementbibliothek in LS-DYNA

-

Wann nimmt man was?

Dr.-Ing. Ulrich Stelzmann
CADFEM Service

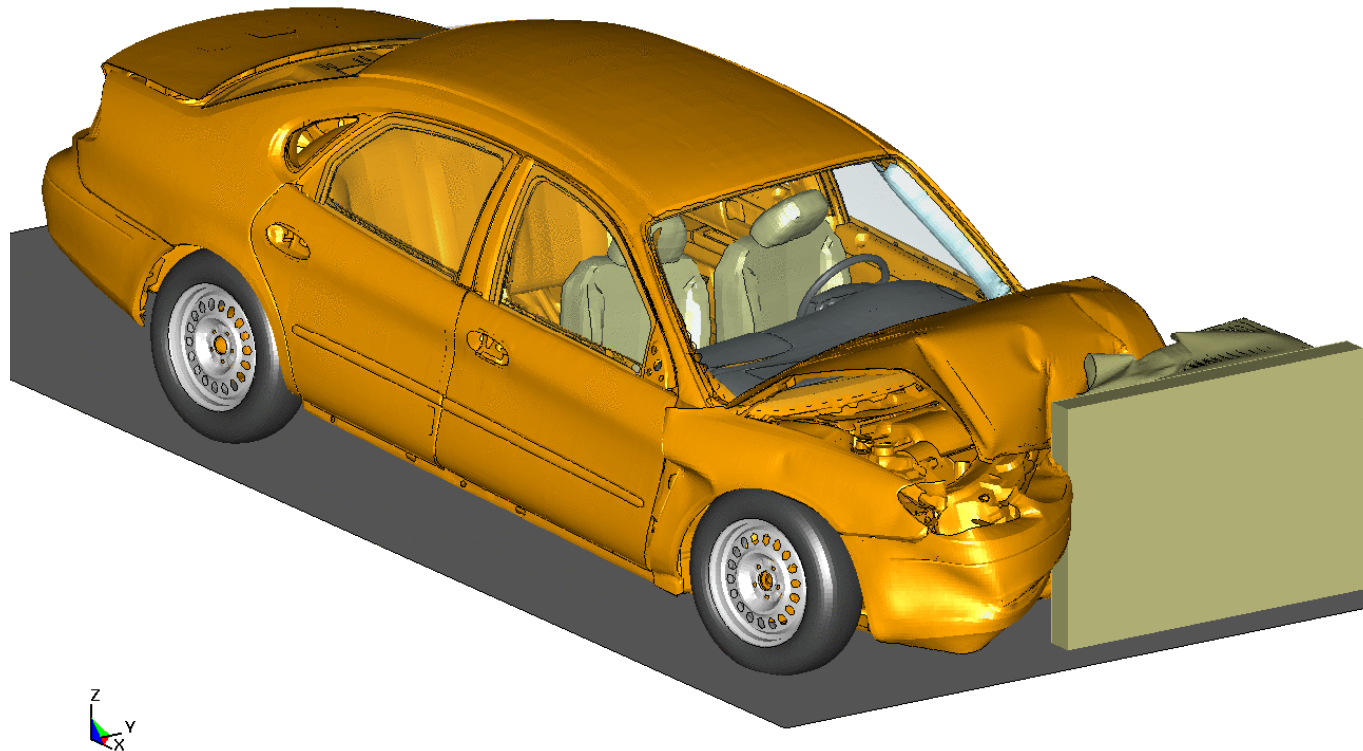


ANSYS Conference & 28th CADFEM Users' Meeting 2010
November 3-5, 2010 – Eurogress Aachen, Germany



A typical model for LS-DYNA may have Millions of elements.
But which element typ should be used for what?

Time = 0.148



*ELEMENT

Elementdefinition in LS_DYNA:

- *ELEMENT_BEAM
- *ELEMENT_DIRECT_MATRIX_INPUT
- *ELEMENT_DISCRETE
- *ELEMENT_INERTIA
- *ELEMENT_MASS
- *ELEMENT_PLOTEL
- *ELEMENT_SEATBELT
- *ELEMENT_SHELL
- *ELEMENT_SOLID
- *ELEMENT_SPH
- *ELEMENT_TSHELL

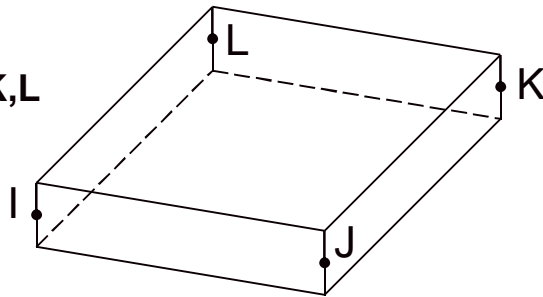
Shell Elements (thin shells)

***ELEMENT_SHELL**
***SECTION_SHELL**

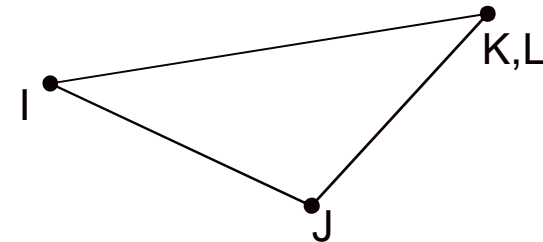
*ELEMENT_SHELL

*SECTION_SHELL:

Shell element:
with the nodes I,J,K,L



triangular
Shell element:



à Shell thickness

- the shell thickness is defined in *SECTION_SHELL, t1 until t4 (commonly four times the same value)
- additional input is possible in the element card, with *ELEMENT_SHELL_THICKNESS; this overwrites the thickness from section definition
- in order to consider thickness change of the shell due to membrane straining one has to set *CONTROL_SHELL, istupd (important for large strains)

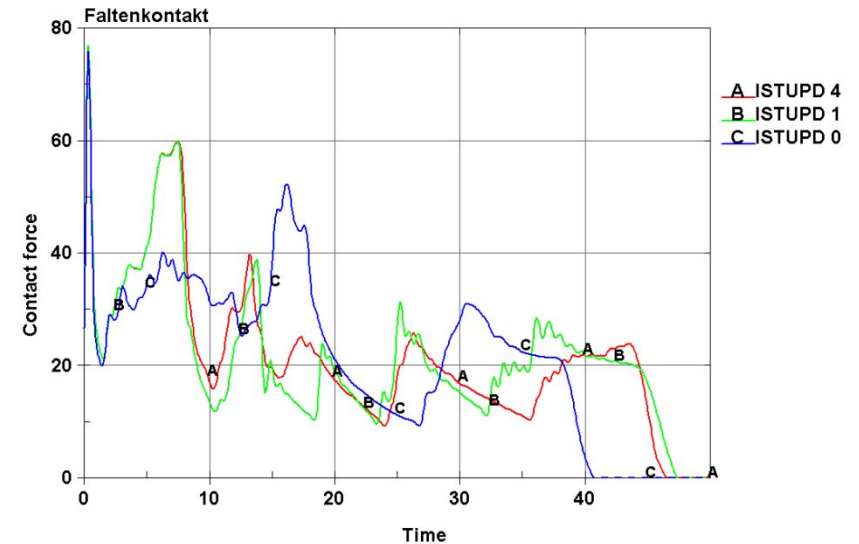
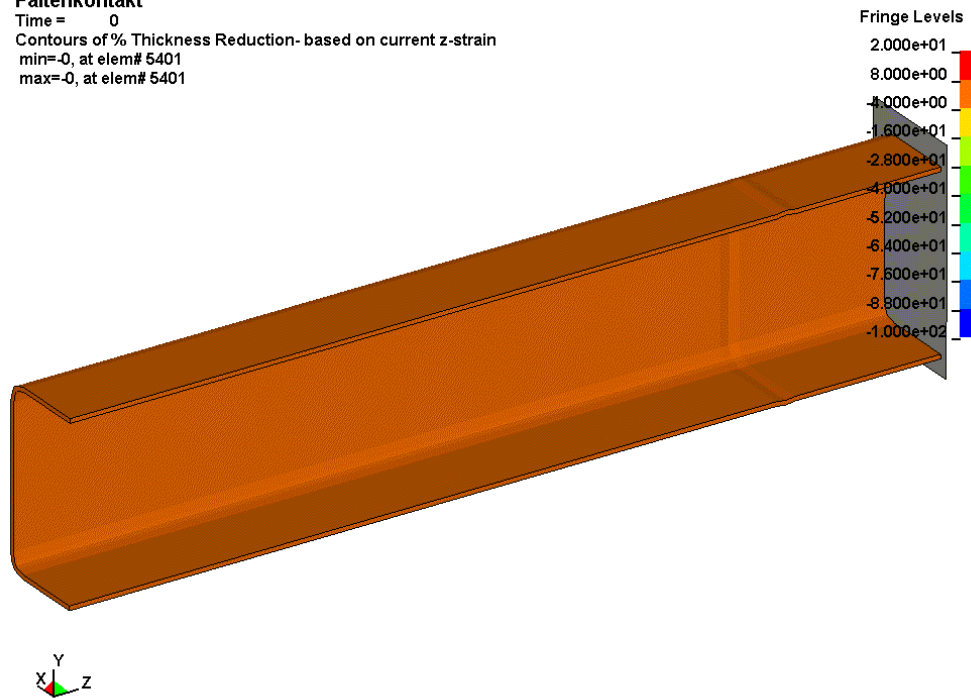
recommended: **istupd=4** (for elastic-plastic material models, only the plastic part of membrane strain will change the thickness)

*ELEMENT_SHELL

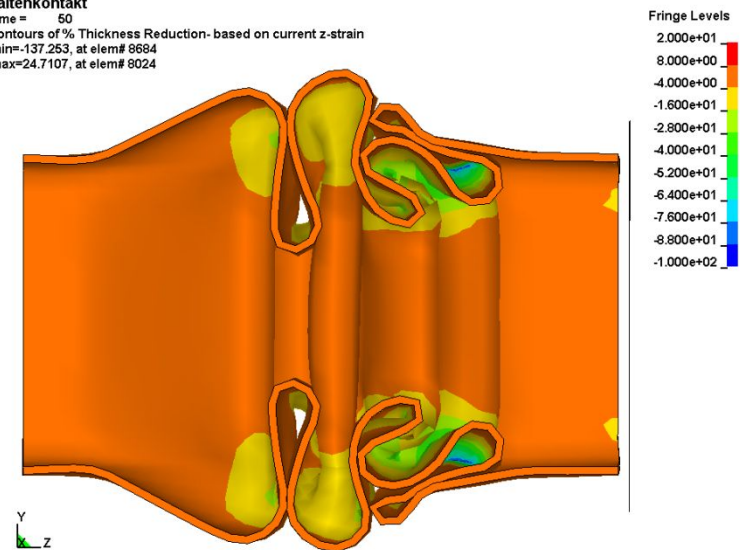
The change of shell thickness may have big influence on stiffness if large membran strain exist.

The simplest example is a tensile test.

Faltenkontakt
Time = 0
Contours of % Thickness Reduction-based on current z-strain
min=-0, at elem# 5401
max=-0, at elem# 5401



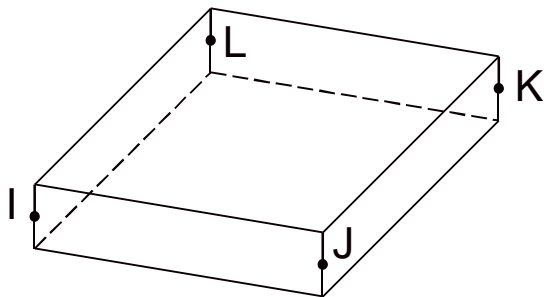
Faltenkontakt
Time = 50
Contours of % Thickness Reduction-based on current z-strain
min=-137.253, at elem# 8684
max=24.7107, at elem# 8024



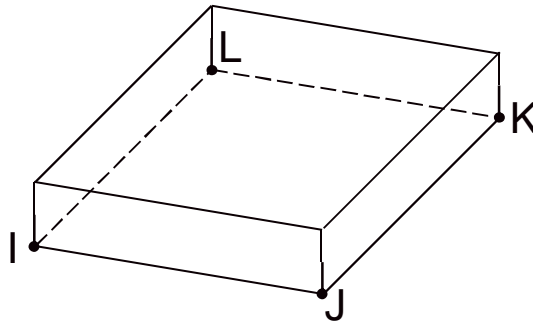
Shell Offset

Offset in shell elements: ***SECTION_SHELL, NLOC**

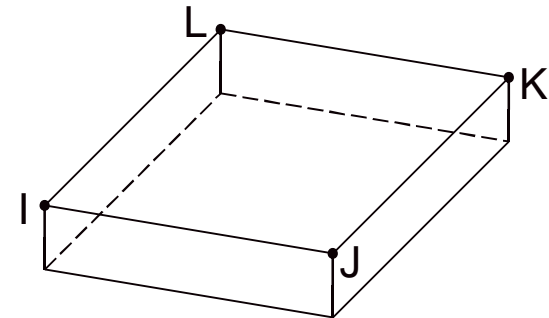
NLOC=0



NLOC=-1



NLOC=1



Since LS971 the reference surface for all shell elements can be set to an arbitrary location, not only between NLOC=-1 and 1 (lower and upper shell surface) but also outside the shell with $|NLOC| > 1$.

Since R3.2 this is considered in contact (normal and AUTOMATIC) by setting:
***CONTROL_SHELL**, third card, third field=1

LS-PREPOST can not display this offset.

*ELEMENT_SHELL

Number of integration points

- most shell elements (other than type 6, 7 and 16) have 1 integration point in plane, shell element types 6, 7 and 16 have 4 integration points in plane
- The number of the integration points across the thickness is variable and must be defined in `*SECTION_SHELL, NIP`
Default is `nip=2`, which is **NOT** sufficient for most applications.
- use the following rules to define the number of integration points throughout thickness:
 - for membranes: 1 integration point
 - for linear material: 2 integration points sufficient
 - à Attention: Stress output not accurate on shell top- and bottom surface
If stress output at top and bottom is important we recommend:
`NIP=5 (*section_shell)`
`INTGRD=1 (*control_shell, Lobatto integration)`
 - in case of non-linear material 3 to 5 (or more) integration points are needed
 - for high accuracy in springback simulations, use 7 to 9 through thickness integ. Points
 - à Attention: If NIP is larger than 10, trapezoidal integration rule is used, which is much less accurate than Gauss integration. Use this only for composite shells.

*ELEMENT_SHELL

integration rule across the thickness

- usually the Gauss integration rule is used for thickness integration
- although the outer integration points are not located on the surface, this method gives accurate results and is commonly used.
- thickness integration can be switched from Gauss to Lobatto integration by setting `*CONTROL_SHELL, intgrd=1`

In this case the inner and outer integration points are located on the shell surface. This feature is only available for 3-10 integration points throughout the thickness

Especially with `nip=3` Lobatto integration is not as good as Gauss.

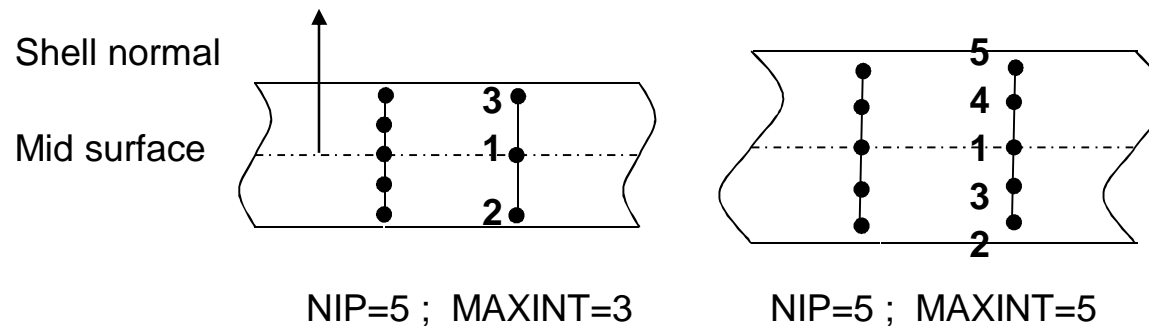
Recommend:
`intgrd=0` and
`nip=5`

GAUSS INTEGRATION RULE					
NUMBER OF GAUSSPOINT	1 POINT	2 POINT	3 POINT	4 POINT	5 POINT
# 1	.0	-.5773503	.0	-.8611363	.0
# 2		+.5773503	-.7745967	-.3399810	-.9061798
# 3			+.7745967	+.3399810	-.5384693
# 4				+.8622363	+.5384693
# 5					+.9061798
LOBATTO INTEGRATION RULE					
NUMBER OF INTEG. POINT	-	-	3 POINT	4 POINT	5 POINT
# 1			.0	-1.0	.0
# 2			-1.0	-.4472136	-1.0
# 3			+1.0	+.4472136	-.6546537
# 4				+1.0	+.6546537
# 5					+1.0

*ELEMENT_SHELL

Output of integration point results

- with `*DATABASE_EXTENT_BINARY; maxint`, declare the number of integration points, for which LS-DYNA writes results to the binary database d3plot
- à for `maxint=3` (default) the results are written for the middle and the two outermost integration points, available as middle, lower and upper surface

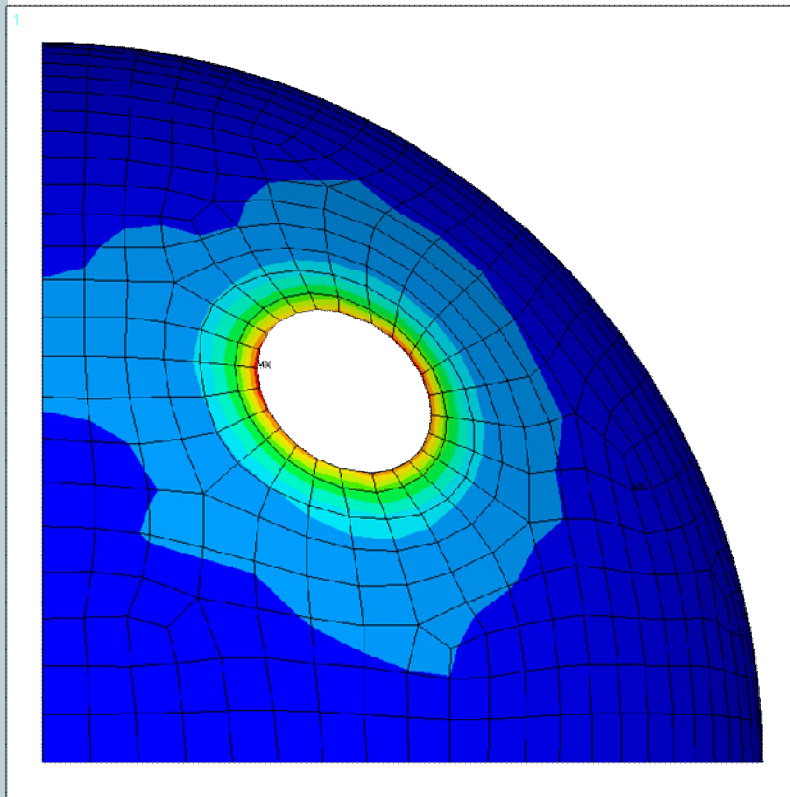


Note: LS-DYNA does not extrapolate stresses from integration points to nodes and write them to the binary database d3plot. Extrapolated stresses can be written to the ASCII file ELOUTDET for selected elements, controlled by `*DATABASE_EXTENT_BINARY; INTOUT` and `NODOUT`

Stress Output

As default for shell elements LS-DYNA writes out only one inplane integration point to the d3plot database. For fully integrated shells an averaged value is written. Because of this stress gradients can not exactly be displayed in the LS-PREPOST.

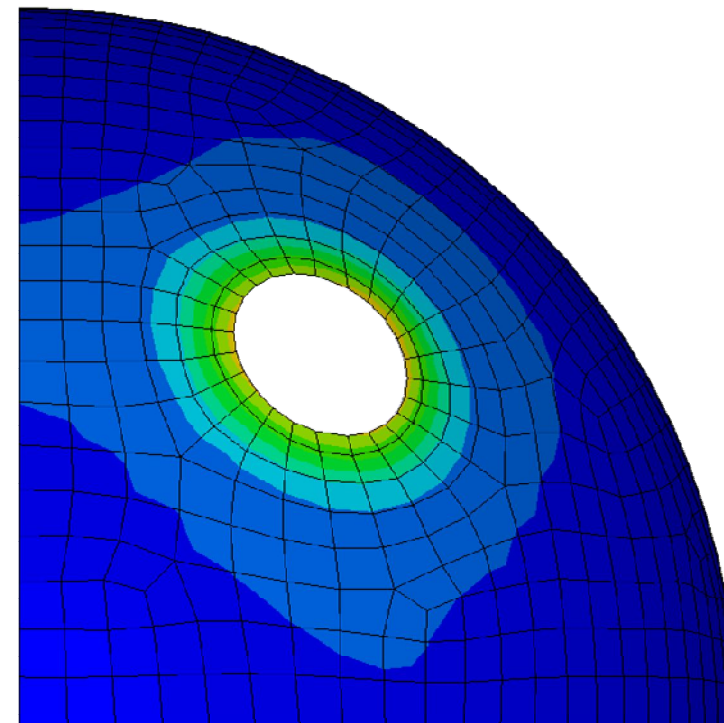
ANSYS/POST1



ANSYS 11.0
NODAL SOLUTION
STEP=1
SUB =1
TIME=1
SEQV (AVG)
MIDDLE
SMN =147.531
SMX =312.876

135
153
171
189
207
225
243
261
279
297
315

LS-PREPOST



Fringe Levels

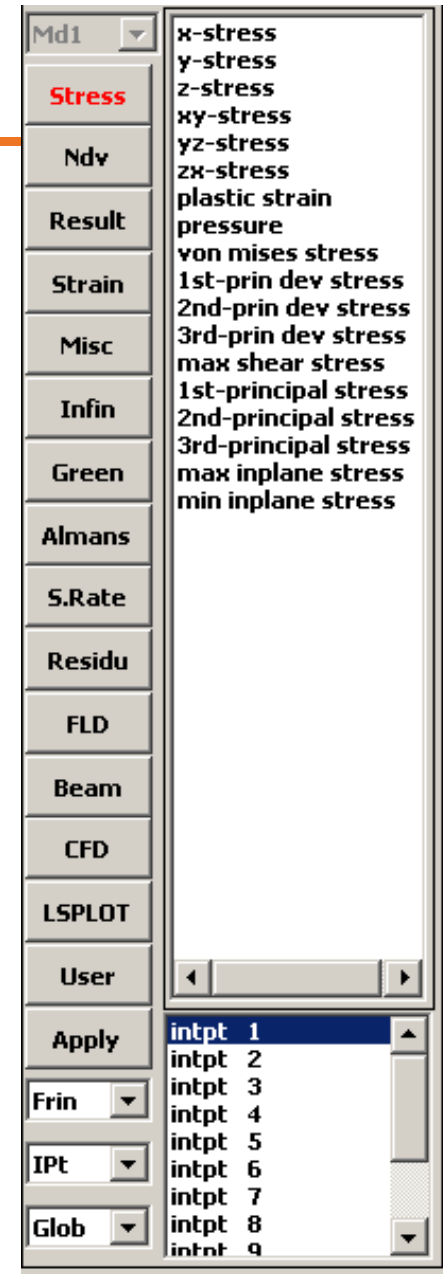
3.150e+02
2.970e+02
2.790e+02
2.610e+02
2.430e+02
2.250e+02
2.070e+02
1.890e+02
1.710e+02
1.530e+02
1.350e+02

Stress Output

Since LS971_R3.2.1 four inplane integration points for shells can be written to the d3plot database by setting *DATABASE_EXTENT_BINARY, MAXINT as a negative number. The absolute value of MAXINT describes the number of through thickness integration points.

For a fully integrated shell element with five through thickness integration points and MAXINT=-3, 12 integration points are written. Current version 2.4 and 3.0 of LS-PREPOST is now able to use all four inplane integration points for averaged nodal stresses. But without averaging again only a constant value is displayed for each element.

Third party Postprocessors may show 12 thickness integration points in this case if they are not updated – like LSPP 2.3 ...

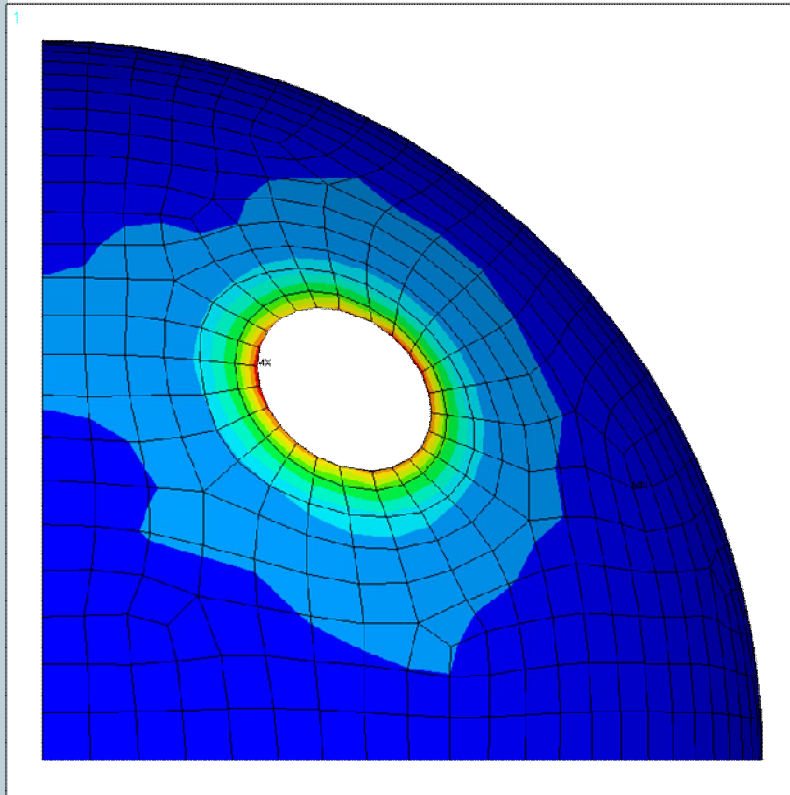


Stress Output

Now stress gradients and stress spots can be visualized more accurately in the LS-PREPOST.

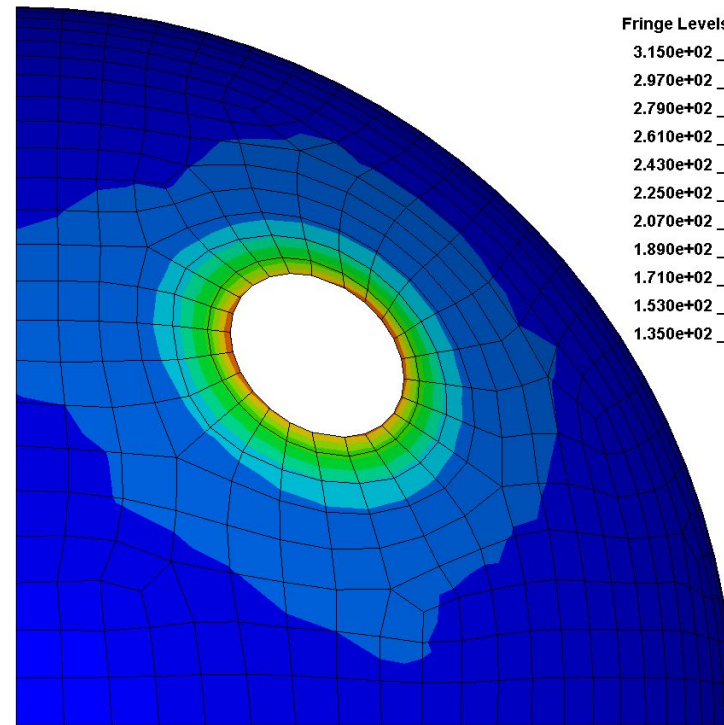
ANSYS/POST1

LS-PREPOST



ANSYS 11.0
NODAL SOLUTION
STEP=1
SUB =1
TIME=1
SEQV (AVG)
MIDDLE
SMN =147.531
SMX =312.876

135
153
171
189
207
225
243
261
279
297
315



Fringe Levels

3.150e+02
2.970e+02
2.790e+02
2.610e+02
2.430e+02
2.250e+02
2.070e+02
1.890e+02
1.710e+02
1.530e+02
1.350e+02

*ELEMENT_SHELL

*SECTION_SHELL: ELFORM

EQ.1: Hughes-Liu	à expensive
EQ.2: Belytschko-Tsay	à default, recommended
EQ.3: BCIZ triangular shell	
EQ.4: co-rotational C0, triangular shell	à indirect use
EQ.5: Belytschko-Tsay membrane	à FABRIC only
EQ.6: S/R Hughes-Liu	} Very expensive
EQ.7: S/R co-rotational Hughes-Liu	
EQ.8: Belytschko-Leviathan shell	
EQ.9: fully integrated Belytschko-Tsay membrane	à FABRIC only
EQ.10: Belytschko-Wong-Chiang	
EQ.11: fast (co-rotational) Hughes-Liu	à expensive
EQ.12: plane stress (x-y plane)	} Only 2D
EQ.13: plane strain (x-y plane)	
EQ.14: axisymmetric solid (y-axis of symmetry) – area weighted	
EQ.15: axisymmetric solid (y-axis of symmetry) – volume weighted	
EQ.16: fully integrated shell element with EAS-formulation	à recommended
EQ.17: fully integrated DKT, triangular shell element	à indirect use
EQ.18: fully integrated linear DK quadrilateral/triangular shell	} Only for linear implicit!
EQ.20: fully integrated linear assumed strain C0 shell	
EQ.21: fully integrated linear assumed strain C0 shell	
EQ.25: Belytschko-Tsay shell with thickness stretch	} Under development
EQ.26: fully interated shell (EAS formulation) with thickness stretch	
EQ.27: C0 triangular shell with thickness stretch	

*ELEMENT_SHELL

Belytschko-Tsay-Shell (Type 2)

- standard element with one point integration
- very fast
- problems in case of warping and large shear deformation
- very efficient: moderate accuracy (often sufficient) in combination with high speed
- Quality can be improved by:

Belytschko-Wong-Chiang warping stiffness: *CONTROL_SHELL, bwc=1, proj=1
(ca. 20% more CPU time)

à Recommended shell element if speed is desired

Fully integrated shell (Type 16)

- fully integrated element with EAS-formulation and without Hourglass modes
- very fast for a fully integrated element (2.5 times more expensive than type 2)
- new standard element of Belytschko-Tsay group for increased accuracy
- Bathe/Dvorkin method for improvement of transversal shear
- behaviour of warped elements can be improved by

*HOURGLASS, IHQ=8 (15% speed penalty)

à Recommended shell element if accuracy is desired

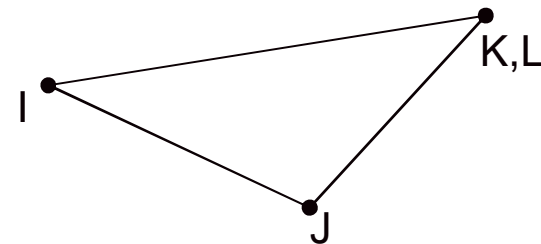
*ELEMENT_SHELL

C0 Triangular shell (Type 4)

- one point integrated special triangular element, no hourglass forms
- by setting *CONTROL_SHELL, esort=1, all triads use this formulation automatically
- only a small number of triads recommended in a quad dominated mesh (too stiff)
- 0.85 times less expensive than type 2, but two times more elements

DKT Triangular Shell (Typ 17)

- fully integrated Discrete Kirchhoff Triangular-Element
- sometimes better than triads type 4, particularly in bending, but needs twice calculation time
- by setting *CONTROL_SHELL, esort=2, all triads use this formulation automatically
- may become unstable for thick elements → not generally recommended



*ELEMENT_SHELL

Belytschko-Wong-Chiang (Type 10)

- slightly slower than type 2 (1.2 times more expensive than type 2)
- nearly identical to type 2 if *CONTROL_SHELL, bwc=1, proj=1 is set

Belytschko-Leviathan (Type 8)

- calculation time and accuracy comparable to type 10 (1.4 times more expensive than type2)
- physical Hourglass control, i.e. no input of Hourglass parameters needed
- for linear material it should be as accurate as an fully integrated element

Hughes-Liu-Shell (Type 1)

- first shell element in LSDYNA (DYNA3D)
- developed from continuum model, one point integration
- high accuracy (also in case of twisted elements)
- highly expensive (2.5 times more expensive than type 2)

selective reduced Hughes-Liu-Shell (Type 6,7)

- most costly shell element (10–20 times more expensive than type 2) → never recommended
- only shear part with reduced integration, otherwise 4 integration points in plane
→ thus only one Hourglass mode
- use of *CONTROL_SHELL; CSTYP=2 (unique normal orientation) is approved

*ELEMENT_SHELL

Belytschko-Tsay- Membrane (Type 5)

- membrane element without bending stiffness
- only 1 integration point throughout the thickness
- one integration point in the element plane (Hourglass modes possible)
- Use only with *MAT_FABRIC à not generally recommended

Fully integrated Belytschko-Tsay- Membrane (Type 9)

- same as Type 5, but 4 integration points in the element plane (no Hourglass modes)
- Use only with *MAT_FABRIC à not generally recommended

Recommendation for general membrane element: Typ 2 or Typ 16 with NIP=1 !

*ELEMENT_SHELL

Newest development: thin shell elements with thickness stress

Classical shell elements have the assumption that the stress in thickness direction is zero. The new element formulations 25-27 consider a fully three dimensional stress state, like in solid elements. Therefore 3d material models are necessary for these elements. To apply appropriate forces – currently by contact only - these elements have 8 additional nodes: four nodes at the top surface and four nodes at the bottom surface. These nodes are generated automatically and are invisible.

Belytschko-Tsay Shell with thickness stretch (Typ 25)

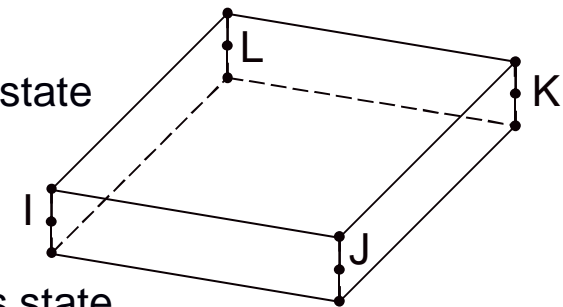
- extended formulation of type 2 shell with 3 dimensional stress state
- one inplane integration point

Fully integrated Shell with thickness stretch (Typ 26)

- extended formulation of type 16 shell with 3 dimensional stress state
- four inplane integration points, no hourglass modes

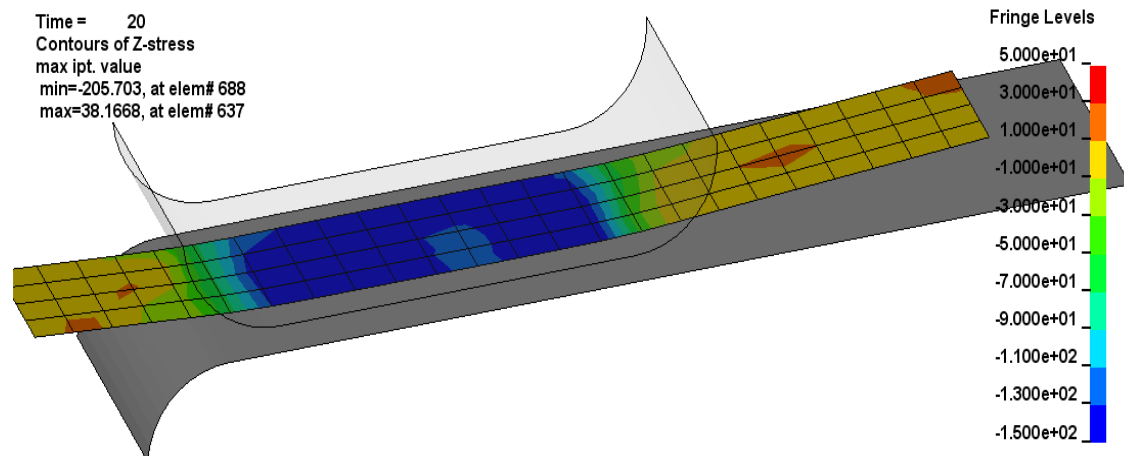
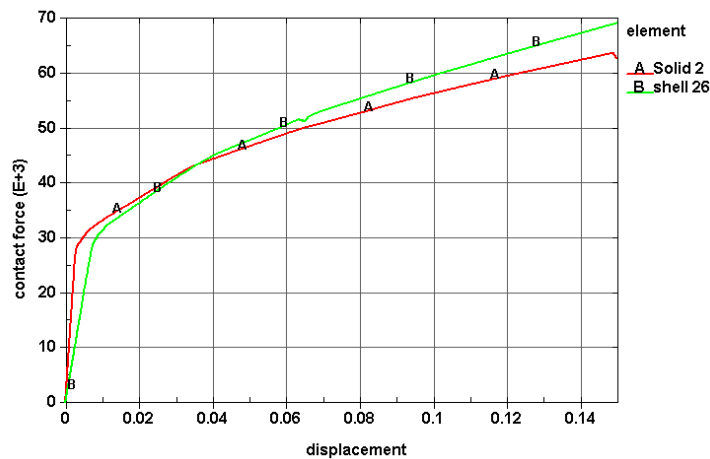
C0 Triangular Shell with thickness stretch (Typ 27)

- extended formulation of type 4 triad shell with 3 dimensional stress state
- one inplane integration point



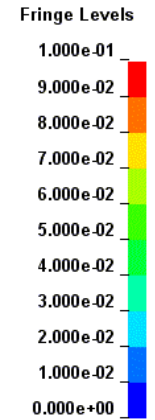
Shell elements with thickness stretch

- LS-DYNA shell element types 25-27 are available starting R4 version. These elements are locking like 4 noded shells. Internally additional nodes (scalar nodes) are generated to represent the upper and lower surface. With IDOF in *SECTION_SHELL it can be decided if each element has its unique scalar nodes (IDOF=2: discontinuous thickness field) or connected by the neighbor scalar nodes (IDOF=1: continuous thickness field).
- In R5 the standard shell type 16 gets an similar option by setting IDOF=3 in *SECTION_SHELL. This implementation is much easier than elform 25-27. Only contact forces between shell lower and upper surface can cause thickness stress here.
- In opposite to solid elements: time step size is not based on thickness!



Shell elements with thickness stretch

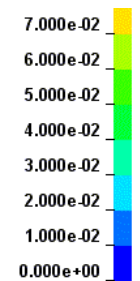
Time = 0
Contours of Effective Plastic Strain
max ipt. value
min=0, at elem# 706
max=0, at elem# 706



2 Solid elements
over thickness
elform=1,
IHQ=6



Shell element
elform=26



*ELEMENT_SHELL

Conclusion:

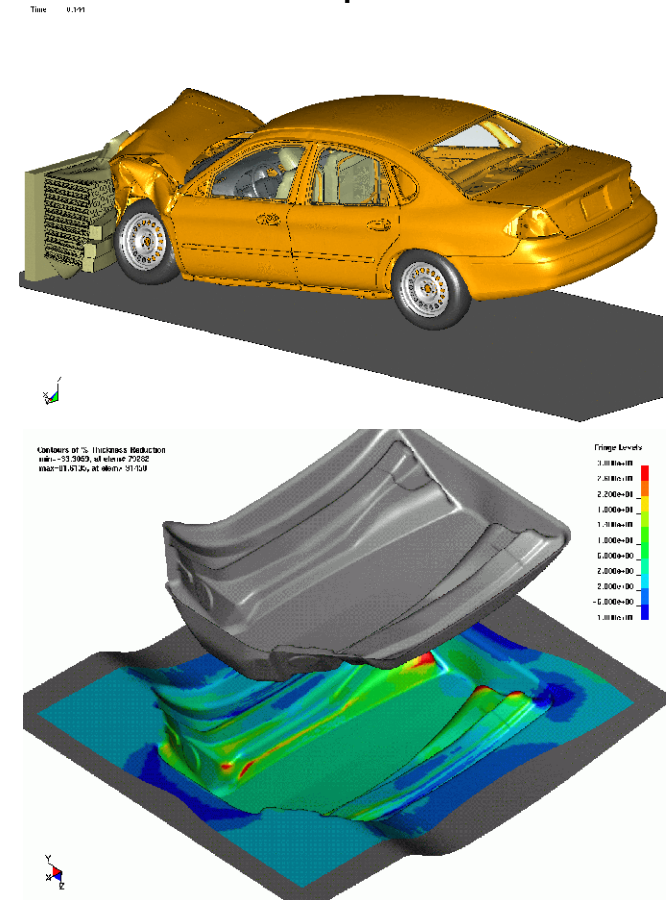
For most structural applications only two element formulations are important:

Prefer speed:

- *SECTION_SHELL; ELFORM=2, NIP=3
- *HOURGLASS; IHQ=4
- *CONTROL_SHELL; ISTUPD=4, BCW=1, PROJ=1

Prefer accuracy:

- *SECTION_SHELL; ELFORM=16, NIP=5
- *HOURGLASS; IHQ=8
- *CONTROL_SHELL; ISTUPD=4



Thick Shell Elements

***ELEMENT_TSHELL**

***SECTION_TSHELL**



*ELEMENT_TSHELL

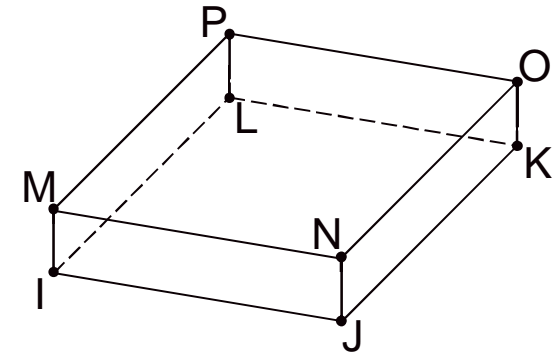
*SECTION_TSHELL: ELFORM

EQ.1: one point reduced integration

EQ.2: selective reduced 2 x 2 in plane integration

EQ.3: assumed strain 2 x 2 in plane integration

EQ.5: assumed strain reduced integration



8 Node Shell elements are rarely used in LS-DYNA

*ELEMENT_TSHHELL

One point integrated thick shell (Type 1)

- 8 node shell with 2d stress state (like thin shell)
- The thickness of the element is constrained by a penalty function between top and bottom nodes, thickness can only change by membran strain (like thin shell)
- Hourglass modes possible, mostly much to soft à not recommended

Fully integrated thick shell (Type 2)

- 8 node shell with 2d stress state (like thin shell)
- The thickness of the element is constrained by a penalty function between top and bottom nodes, thickness can only change by membran strain (like thin shell)
- No Hourglass modes
- Accuracy comparable to thin shell but much slower: 7-8 times more than type 2 shell
- Time step size is based on all three dimensions
- à Typically no advantages over thin shells

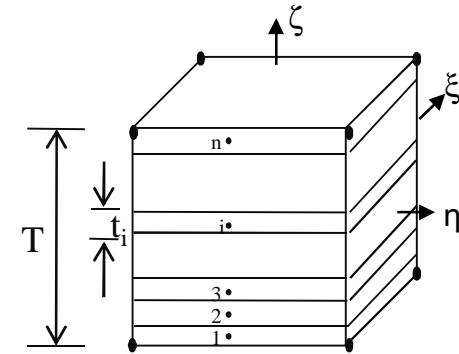
Assumed strain thick shell (Type 3)

- 8 node shell/solid with 3d stress state
- thickness can change by thickness stress
- extremely slow: 65 times more than type 2 shell
- not really a shell: at least two elements over thickness are necessary so solve for bending
- à Not meaningful in explicit

*ELEMENT_TSHELL

Layered solid (Type 5)

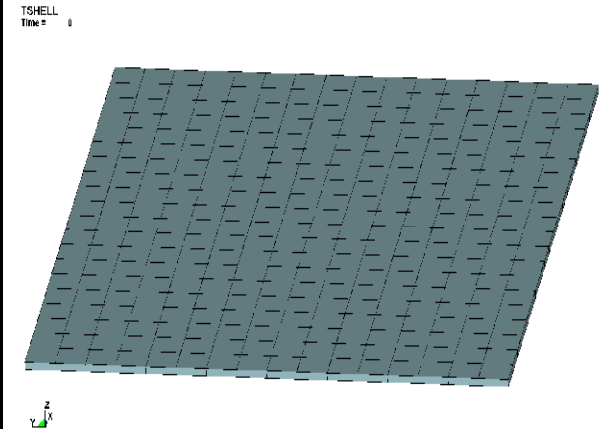
- 8 node layered solid with 3d stress state
- new Feature in LS971 R5
- access to *MAT_ELASTIC and composite material models *MAT_22, *MAT_59
- shear locking and hourglass stabilization by build in assumed strain method, *HOURGLASS setting has no meaning
- efficient in computation: only 1.5 times more than type 2 shell
- solves bending problem with only one element over thickness, also for bad aspect ratio, of course more elements are possible
- laminated shell theory available (*CONTROL_SHELL, lamsht=4), but makes only sense if only one element over the thickness
- developed to model thick composite structures



*ELEMENT_TSHELL

Element type	Time step	Number of cycles	Time (sec)	Element cost factor	Reaction force (N)
Shell type 2	1.66e-3	18135	5	1	0.20
Shell type 16	1.66e-3	18133	12	2.4	0.20
Tshell type 1	3.3e-4	90760	43	1.7	2.8
Tshell type 2	3.3e-4	90705	193	7.7	0.20
Tshell type 3	2.82e-4	106540	1950	66.4	0.14
Solid 1, IHQ=6	2.82e4	106570	37	1.25	0.203
Tshell type 5 (MLS)	2.82e-4	106570	41	1.5	0.206

Rectangular plate under concentrated load



Dimension: 200*160*2
 Element length: 10mm
 Thickness: 2mm

Solid Elements

***ELEMENT_SOLID**

***SECTION_SOLID**

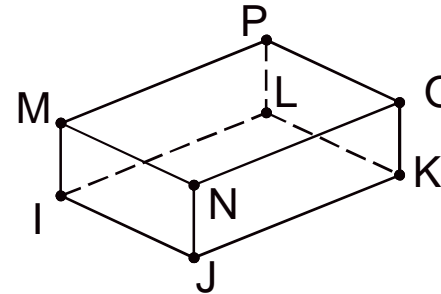


*ELEMENT_SOLID

Element shapes:

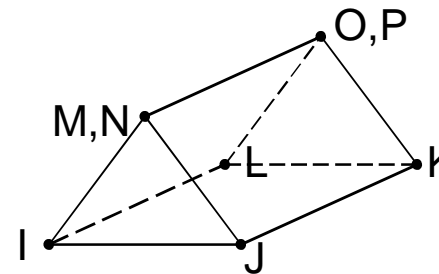
Hexahedron:

(favoured solid element)



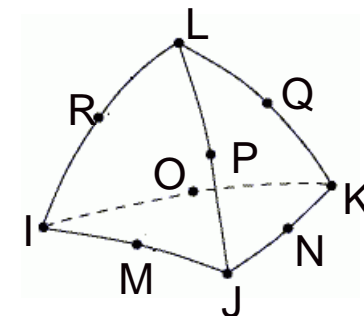
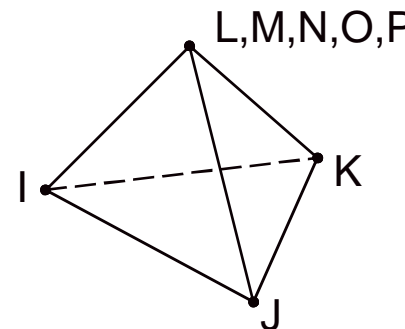
Pentahedron:

(extrusion in depth of triangles)



Tetrahedron:

- 4-noded without rotation: very stiff, only used for foams
- 4-noded with rotation: compromise between effort and accuracy
- 10-noded very accurate but also very costly in terms of computation time



*ELEMENT_SOLID

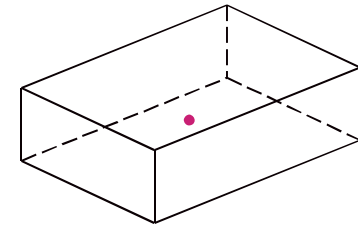
*SECTION_SOLID: ELFORM

- EQ.1: constant stress hexahedron element à default
- EQ.2: fully integrated S/R hexahedron
- EQ.-1: fully integrated S/R hexahedron without shear locking (simplified)
- EQ.-2: fully integrated S/R hexahedron without shear locking
- EQ.3: fully integrated quadratic 8 node hexahedron with nodal rotations
- EQ.4: **S/R quadratic tetrahedron element with nodal rotations**
- EQ.5: 1 point ALE hexahedron
- EQ.6: 1 point Eulerian hexahedron
- EQ.7: 1 point Eulerian ambient hexahedron
- EQ.8: acoustic hexahedron
- EQ.9: 1 point corotational hexahedron for *MAT_MODIFIED_HONEYCOMB
- EQ.10: **1 point tetrahedron**
- EQ.11: 1 point ALE multi-material element, hexahedron
- EQ.12: 1 point integration with single material and void, hexahedron
- EQ.13: **1 point tetrahedron with nodal pressure**
- EQ.14: 8 point acoustic hexahedron
- EQ.15: 2 point pentahedron element
- EQ.16: **5 point 10 noded quadratic tetrahedron with mid side nodes**
- EQ.17: 10 noded composite tetrahedron with mid side nodes (new in ls971)
- EQ.18: 8 point enhanced strain hexahedron element for linear statics only
- EQ.19: 4 point cohesive element
- EQ.20: 4 point cohesive element with offsets for use with shells
- EQ.41: Mesh-free hexahedron - EFG
- EQ.42: Mesh-free tetrahedron - EFG

*ELEMENT_SOLID

standard element (Type 1)

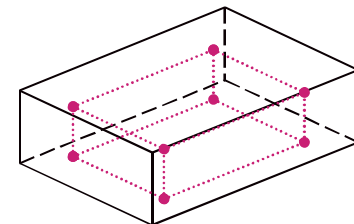
- 8-node hexahedron solid element with tri-linear shape functions
- reduced integration, i.e. stresses are calculated only in one integration point in the middle of the element
- Hourglass modes possible



1 integration point

fully integrated element (Type 2)

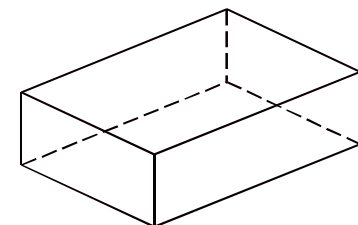
- 8-node hexahedron solid element with tri-linear shape functions
- fully integrated with 8 integration points
 - à no Hourglass modes
- 2-3 times more expensive than type 1
- helpful, if Hourglass modes are a problem
- handicap: lower deformations obtained as with type 1
- handicap: shear locking if bad aspect ratio



8 integration points

fully integrated quadratic 8 node element with nodal rotations (Type 3)

- 8-node hexahedron solid element with quadratic shape function
- 6 degrees-of-freedom per node: translations and rotations
- 14 integration points
- not useful for plasticity or material with Poisson ratio close to 0.5
- very expensive in cpu time (3 times more expensive than type 2)
- à not generally recommended

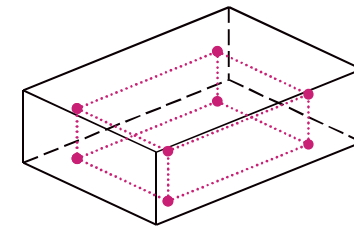


14 integration points

*ELEMENT_SOLID

fully integrated S/R hexahedron without shear locking (Typ -2)

- extension of type 2 hexahedron: accurate formulation
- selectively reduced integration, 8 integration points
 - à no Hourglass forms
- 5 times more expensive than type 2
- Advantage over type 2: no shear locking, also in bad aspect ratios



8 integration points

fully integrated S/R hexahedron without shear locking (Typ -1)

- extension of type 2 hexahedron: simplified, fast formulation
- selectively reduced integration, 8 integration points
 - à may have some remaining hourglass forms
- 1.2 times more expensive than type 2
- Advantage over type 2: no shear locking, also in bad aspect ratios
- Advantage over type -2: much less expensive and in most cases sufficient

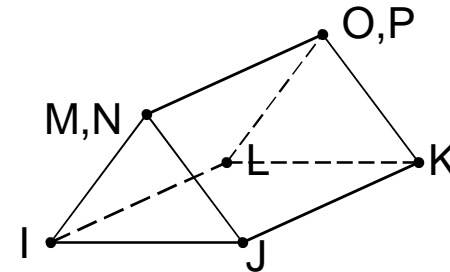
*ELEMENT_SOLID

Pentahedron element (Type 15)

- 6-noded element with trilinear displacement ansatz and 2 integration points
- element typically is generate when triangular surface element is extruded into the depth
- with the input:

*CONTROL_SOLID, esort=1

all 6-noded solid elements get automatically this element formulation



Solid element - (type 8 and type 14)

- for acoustic simulation (sound distribution within fluids)
- nodes only have a pressure degree of freedom
- One point (type 8) and fully integrated (type 14)

1 point corotational for *MAT_MODIFIED_HONEYCOMB (Type 9):

- special hexahedron element for extra large deformations in combination with material law 126 only (*MAT_MODIFIED_HONEYCOMB)
- keeps stable in large element distortion, prevent “negative volume”, but limited accuracy

*ELEMENT_SOLID

Tetrahedron element (Type 10)

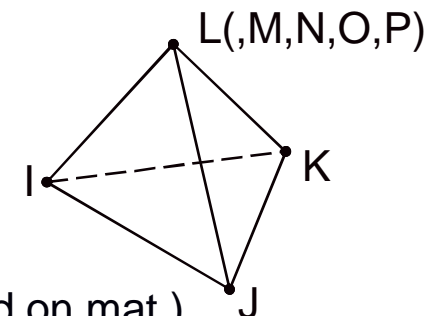
- 4-nodes tetrahedron element with tri-linear shape functions and 1 integration point
- in general much too stiff
 - à often used in combination with foam material, where compression dominates

Tetrahedron element (Type 13)

- like tetrahedron element type 10, but with additional pressure degree of freedom at nodes
- 10-40% more expensive than type 10 – based on material law
- nearly no volumetric locking
 - à recommended for incompressible material like rubber or plastic flow in bulk metal forming
- Shear locking is not much better than tet type 10
- only available for a few material models, for all other materials it switch back type 10

S/R quadratic tetrahedron element with nodal rotations (Type 4)

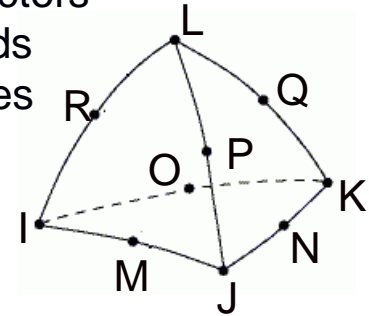
- 4-node tetrahedron solid element with quadratic shape functions
- 6 degrees-of-freedom per node: translations and rotations
- 5 integration points
- expensive in cpu time (1.5-5 times more expensive than type 10 – based on mat.)
- accuracy better than tetrahedron type 10, but less than hexahedron type 2
- sometimes double precision is needed if element rotates



*ELEMENT_SOLID

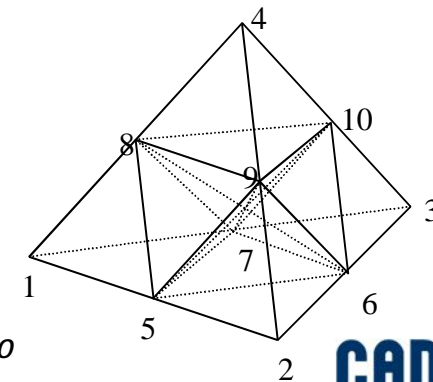
10-noded tetrahedron element (Type 16)

- tetrahedron element with midside nodes and quadratic shape function
- 4+1 integration points
- needs approx. the same computation time as type 4 but time step size is halved
- because of shape function the midside nodes have different weighting factors than the corner nodes → midside nodes should not be used to apply loads
- midside nodes must stay “near” the straight line between the corner nodes
→ not applicable for large deformation



10-noded composite tetrahedron element (Type 17)

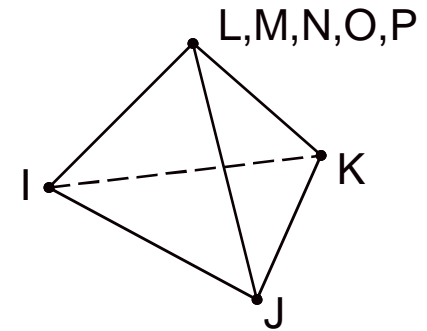
- tetrahedron element with midside nodes and quadratic shape function
- the word “composite” means, that one element internally consists of a lot of smaller 4 noded elements
- because of this, also midside nodes have the same weighting factors than the corner nodes
- 4+1 integration points
- same time step size as type 16 but much more cpu time
- still under development, not always stable



TET13

*SECTION_SOLID, elform=13

4 noded, 1 point tetrahedron with linear shape function and additional pressure degree of freedom.



- Less volumetric locking, therefore recommended for applications where incompressible material behavior is important:

hyper-elastic material (Rubber)

isochoric plastic deformation (Forging)

- Only less improvement in shear locking in comparison to standard linear tetrahedron elform=10

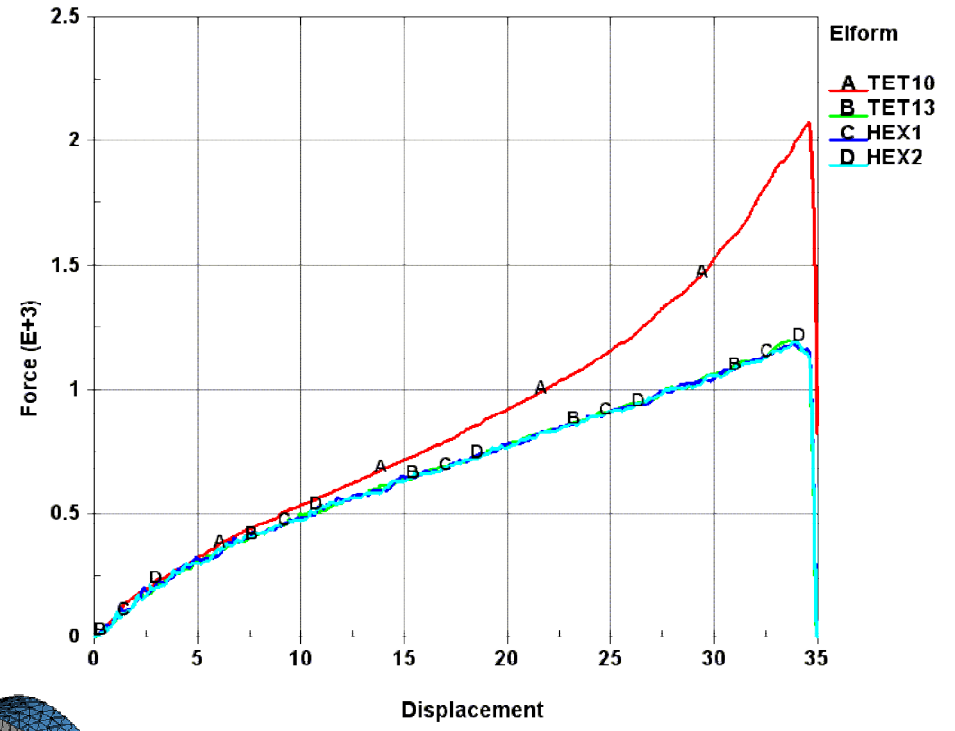
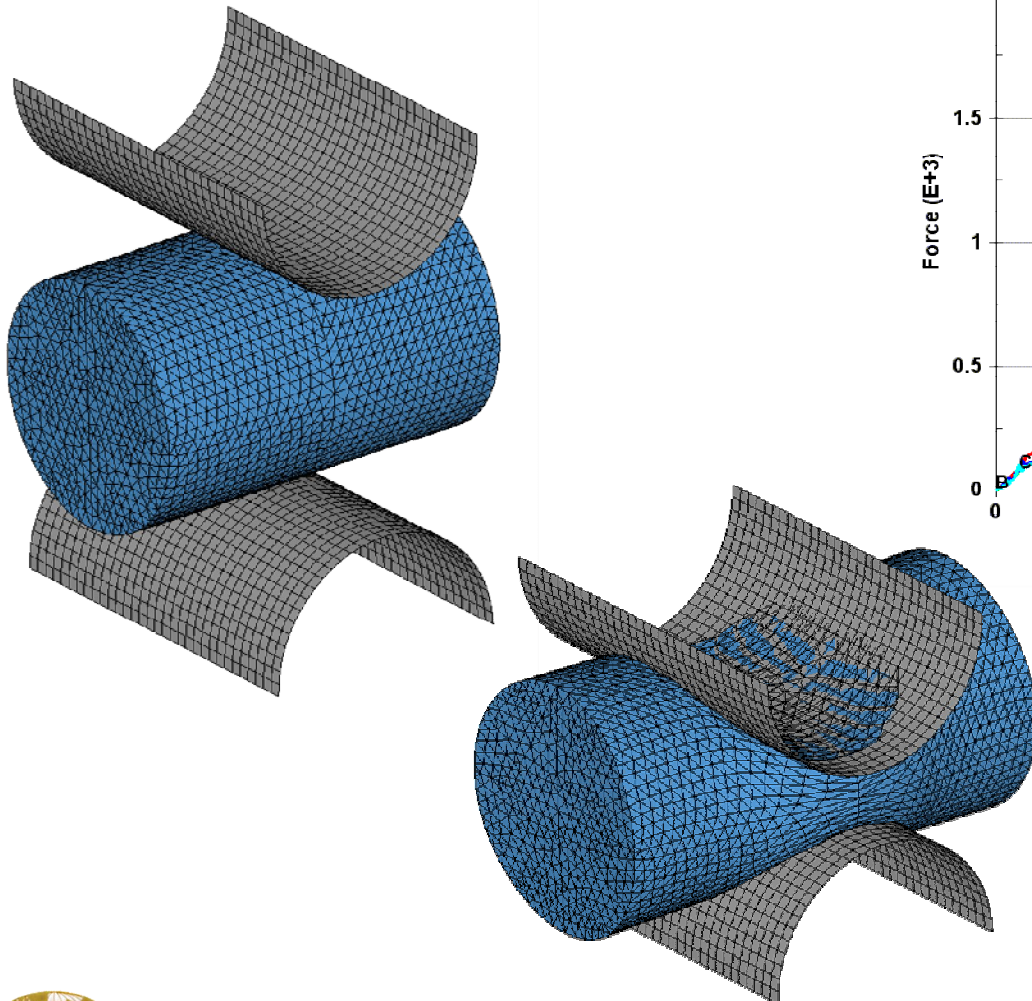
Only available for material models:

1,3,6,24,27,77,81,82,91,92,106,120,123,124,128,129,181,183,225,244

For all other material models this element is identical to elform=10

TET13

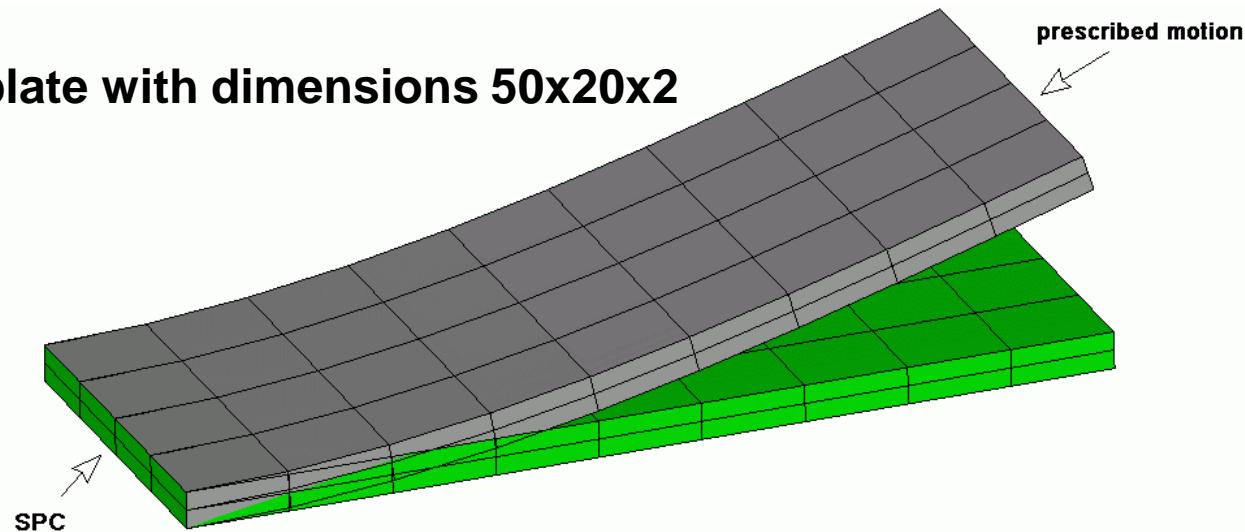
Example: Bulk metal forming



Elform	TET10	TET13	HEX1	HEX2
CPU time	46s	65s	13s	36s

Mesh for thin walled structures

Test model: plate with dimensions 50x20x2



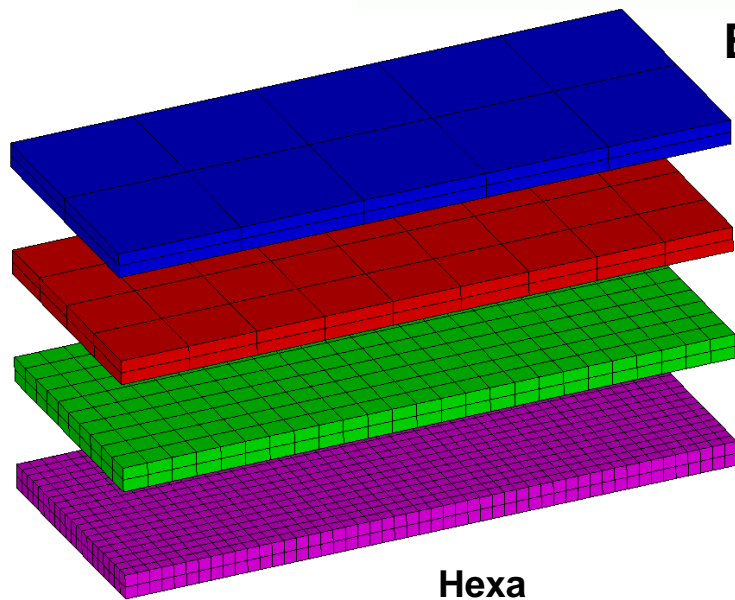
Element size in plane:

10 mm

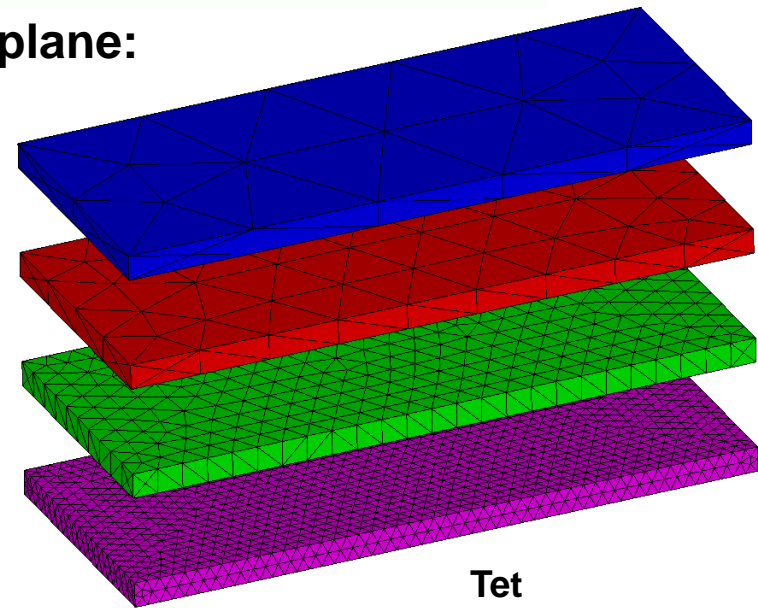
5 mm

2 mm

1 mm



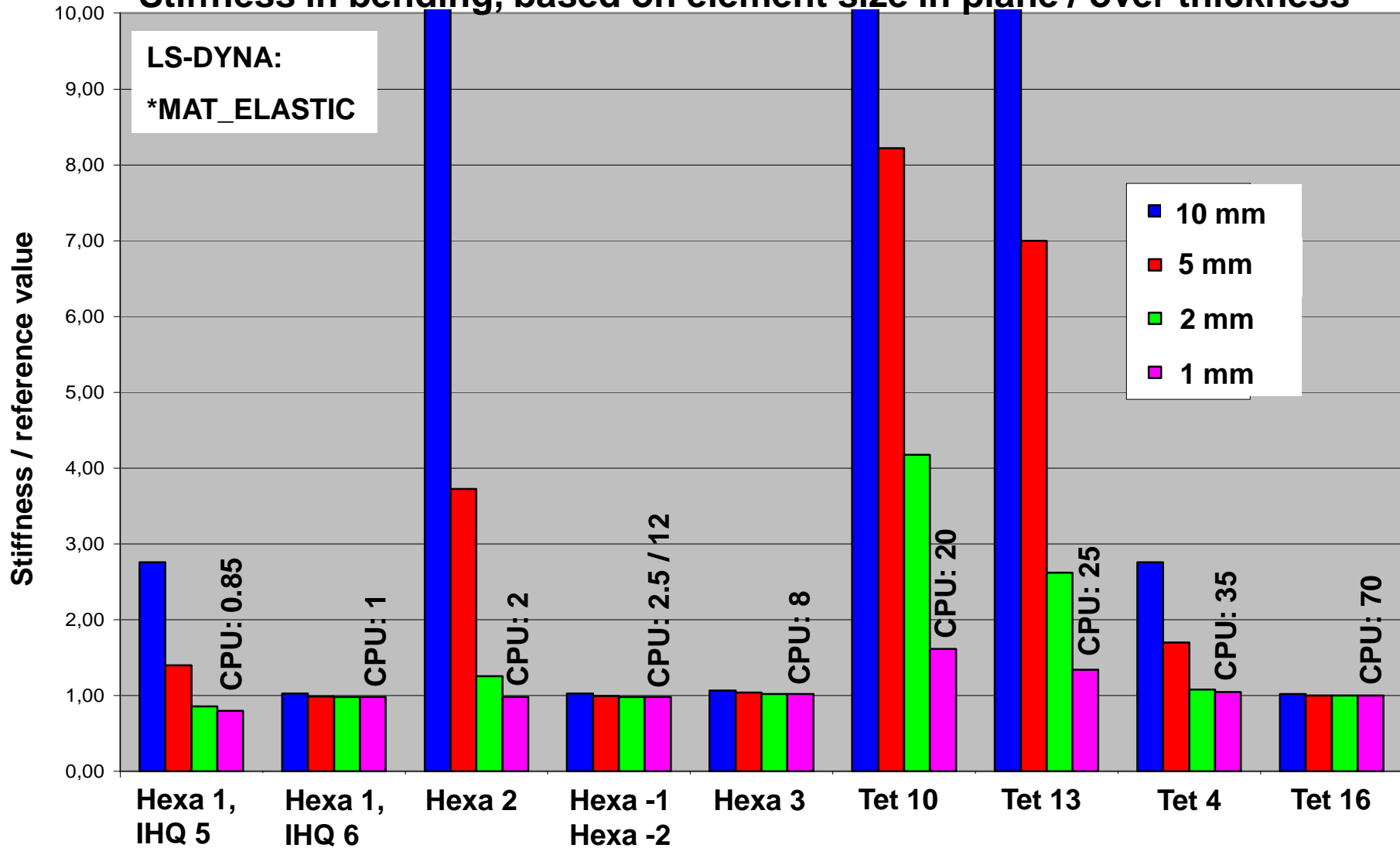
Hexa



Tet

Mesh for thin walled structures

Stiffness in bending, based on element size in plane / over thickness



Mesh for thin walled structures

Conclusions:

- Solid type 2 is very sensitive in aspect ratio for bending problems - shear locking
- Solid type 1 in combination with Hourglass control type 6 (IHQ=6, QM=1) solves linear bending problem very well but is too stiff for plasticity with this setting.
- Solid types -1 and -2 do not show shear locking effects. Type -1 is approx. 30% more expensive than type 2 and type -2 is 5 times more expensive than type 2.
- Linear tetrahedron type 10 is much too stiff in bending. Type 13 is a little bit better, but only available for a limited number of material models. Type 13 is favored for large plastic deformation and rubber material.
- The 10 noded quadratic tetrahedron type 16 gives very good bending response also if aspect ratio is poor. Less elements are necessary. But this element type is more sensitive because of its midside nodes and may have problems with large distortions.
- The older 4 noded tetrahedron type 4 (with nodal rotations) gives only acceptable results if aspect ratio is good. This element type may be recommended only if mesh density is always fine, e.g. to allow large local deformation. This element was improved in R4 regarding large strains and large rotation.

Mesh for thin walled structures

Shells or Solids ?

- 1) if thickness stress is important, then Solids
- 2) if transversal shear is important, then Solids
- 3) if only membrane stiffness is important, then Shells
- 4) for bending: if bending radius R greater than three times thickness
 $R > 3d$: then Shells
if bending radius smaller than thickness
 $R < d$: then 5 Solids over thickness

A new alternative are the new shell elements with thickness stretch:

*SECTION_SHELL, elform=25-27

A complete 3d stress state is used in these elements, so thickness stress is considered in the material law. But for bending the cross section always remain rectangular.

Mesh for thin walled structures

Should I use tetrahedrons ?

Advantage:

- Tetrahedron meshing is easy and fast.

Disadvantage:

- Run time is large.

If the same mesh density is desired, 6 tetrahedron elements are needed instead of one hexahedron element. à 6 times more effort

One tetrahedron element (type 4 or 16) needs 1.5-5 times more cpu time than one hexahedron (type 1). à 9 - 30 times more effort

The 10 noded tetrahedron type 16 needs half the time step size because of mid side nodes à 18 - 60 times more effort

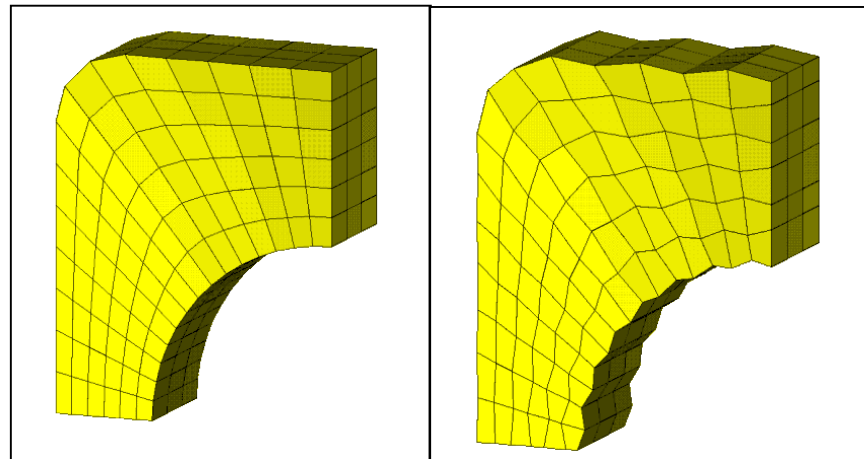
Hourglass control

For hexahedron solid elements Hourglass control is important.
Fully integrated solid elements (*SECTION_SOLID, elform=2) have some disadvantages:

- § higher computation time (approx. factor 2.5)
- § shear locking if aspect ratio is bad
- § may fail if element distortion is too large (negative jacobian)

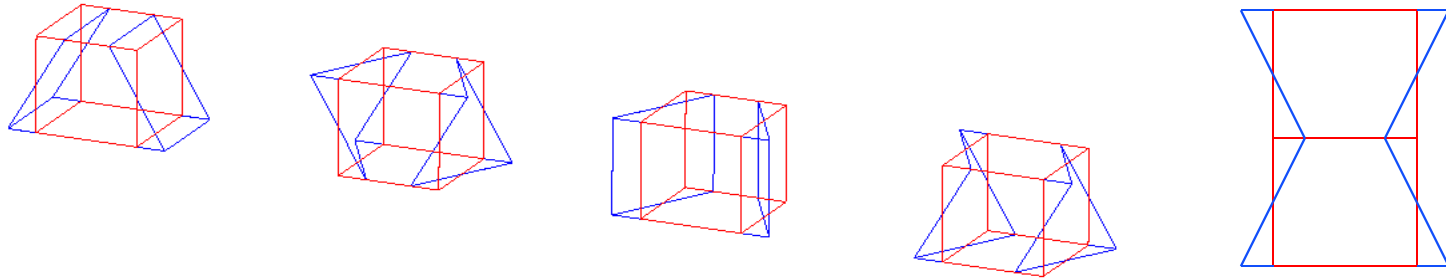
One point integrated element have only two disadvantages:

- § only constant stress over element
- § zero energy deformation possible (Hourglass modes)



Hourglass control

- Hourglassing is a state of strain, which is free of energy (ZEM: Zero Energy Mode) and can emerge in case of one-point-integrated solid- (hexahedrons) and shell elements



- Hourglass modes are mostly caused by:
 - concentrated loads
 - contact (contact force at several nodes)



- in LS-DYNA there are 2 possibilities to prevent Hourglassing:

à using the automatic stabilisation against this deformation with

- ***HOURGLASS** (input for each part) or
- ***CONTROL_HOURGLASS** (global control)

à using a fully integrated element type

disadvantages: - more computation time

- more sensitive with respect to large element deformations

Hourglass control

Recommendation for ***HOURGLASS** and. ***CONTROL_HOURGLASS**

for shell elements	ihq=4	(stiffness form, default settings)
for solid elements (in general)	ihq=5	(stiffness form, default settings)
for solid elements (foam)	ihq=3	(viscous form, default settings)
for solid elements (honeycomb)	ihq=3	(viscous form, default settings)
for solid elements (elastic)	ihq=6,qm=1.0	(stiffness form)
for solid elements (plastic)	ihq=6,qm=0.01-0.001	(stiffness form)
for solid elements (rubber)	ihq=7,qm=1	(stiffness form)
for solid elements (viscoelast.)	ihq=7,qm=1	(stiffness form)

Note: ihq=6 and 7 is a special solid element formulation according to Belytschko-Bindemann

Danger: Default **ihq=1** it not orthogonal to rigid body rotation
à do not use, except for Eulerian elements !

Pentahedrons and Tetrahedrons are not influenced by ***HOURGLASS**.

Hourglass control

For elastic plastic material models the Belytschko-Bindeman assumed strain formulation (*HOURGLASS, IHQ=6) is successfully used.

Unfortunately the Hourglass stiffness is controlled by the elastic material properties only. This tends to that Hourglass control is very accurate for pure elastic deformation but may be much too stiff if material becomes plastic. The user should scale down the Hourglass coefficient manually. A factor of QM=0.1 or QM=0.01 is common.

With LS971 the Puso enhanced assumed strain formulation is also possible (*HOURGLASS, IHQ=9). The behaviour is comparable to the latter one, with one exception: For IHQ=9 the Hourglass coefficient QM can be defined as a negative number. Then the absolute value scales the Hourglass stiffness based on the current material properties. If deformation is purely elastic, the young's modulus is used, in case of plastic deformation the tangent modulus is used.

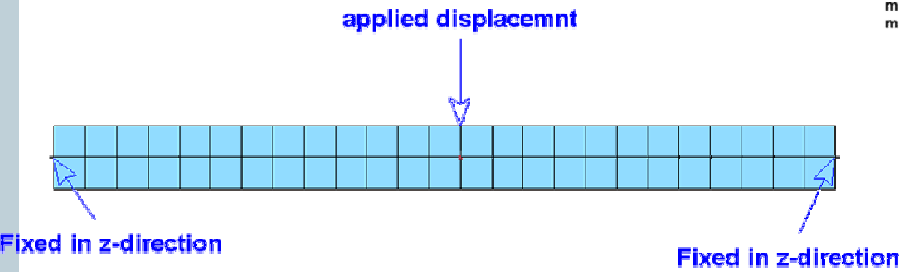
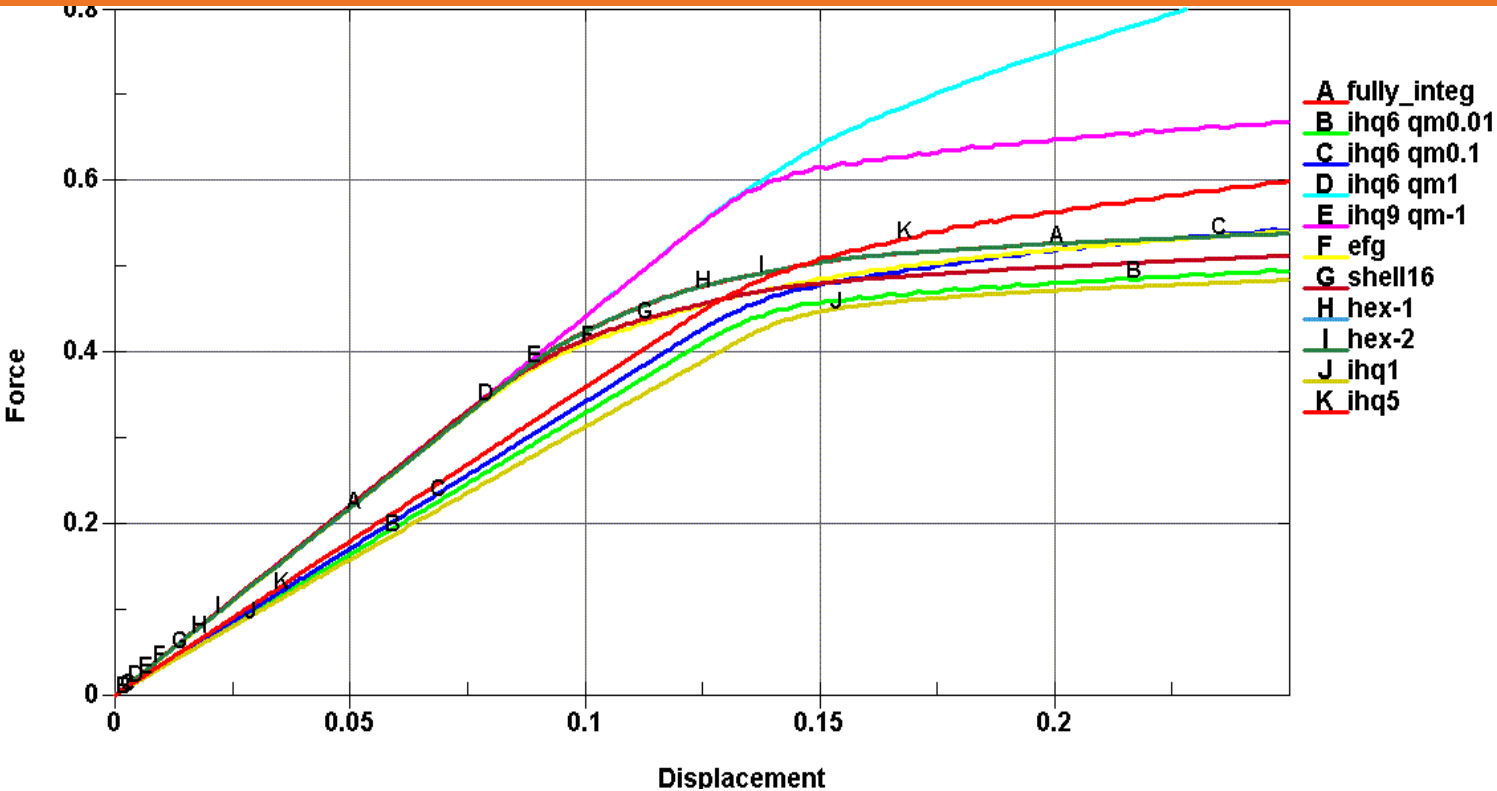
This automatic correction is only available for material models:

- *MAT_PIECEWISE_LINEAR_PLASTICITY (24)
- *MAT_PLASTIC_KINEMATIC (3)
- *MAT_POWER_LAW_PLASTICITY (18)

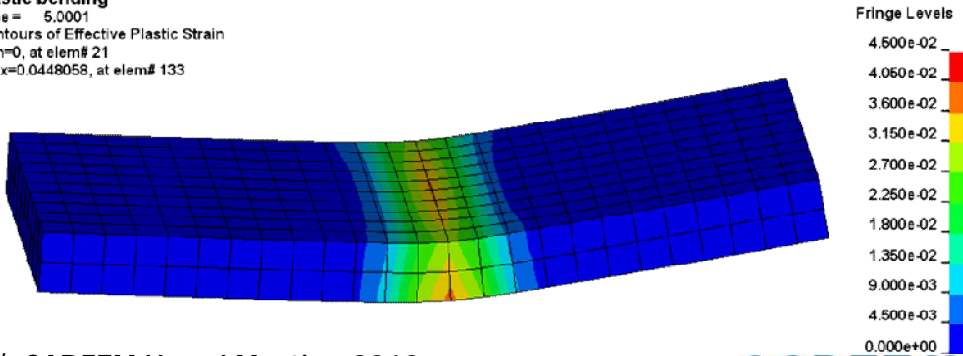


Plastic Bending – HOURGLASS control

Test:
Plastic bending with two elements over thickness

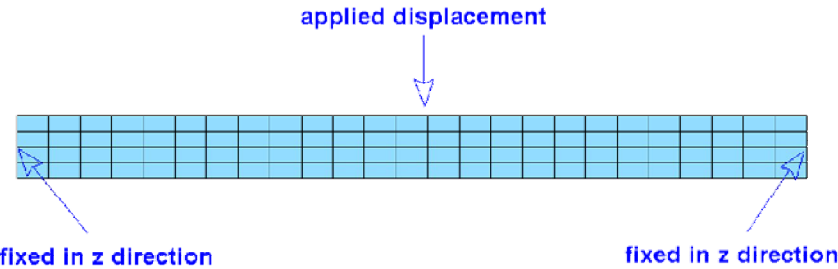
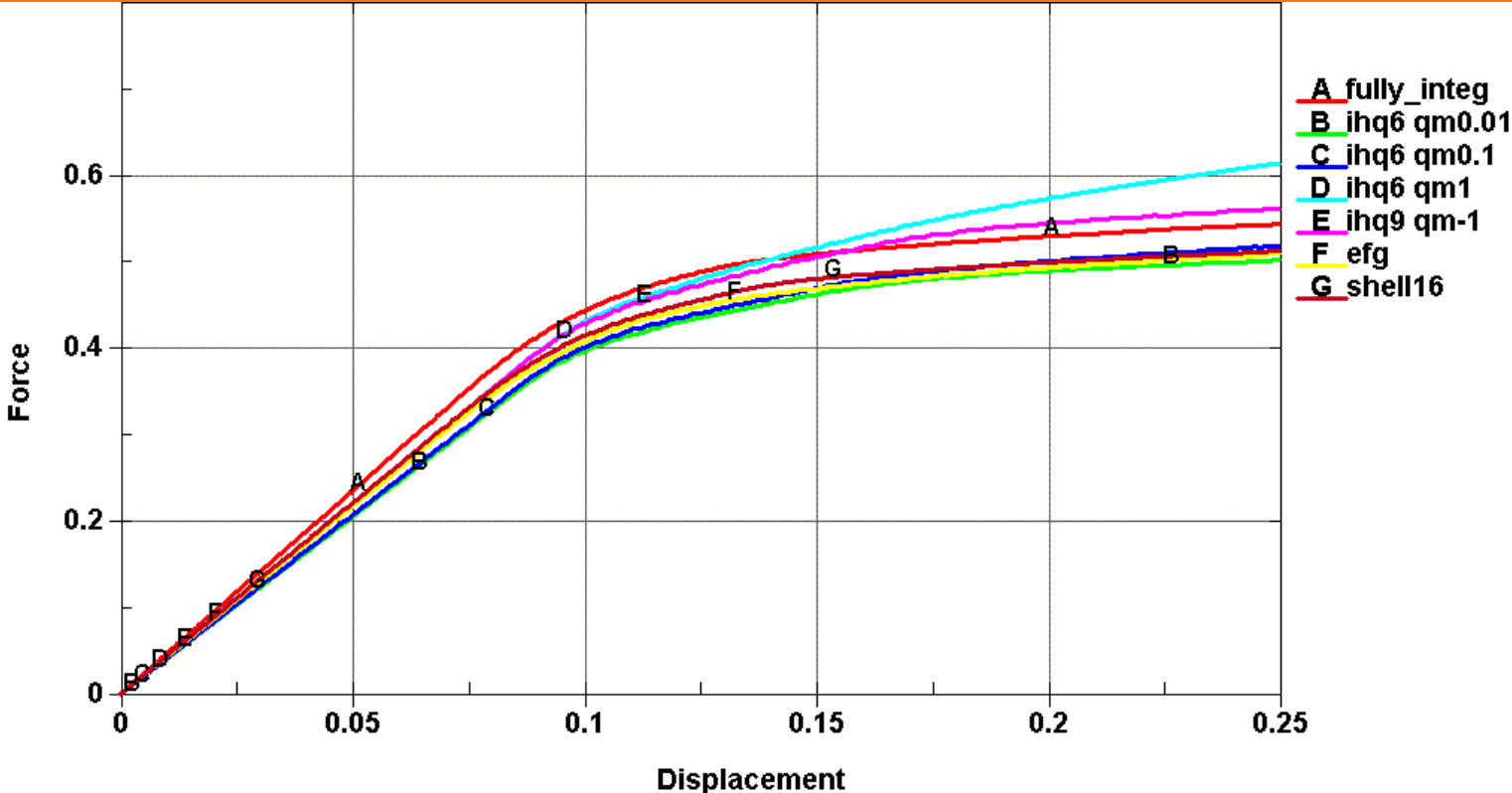


plastic bending
 Time = 5.0001
 Contours of Effective Plastic Strain
 min=0, at elem# 21
 max=0.0448058, at elem# 133

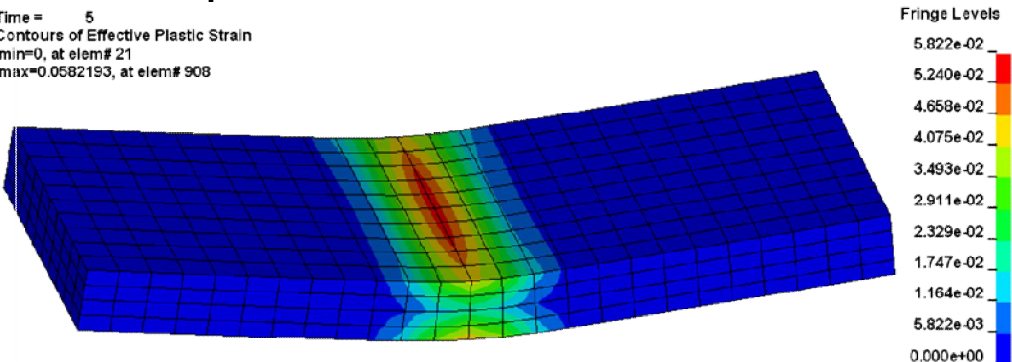


Plastic Bending – HOURGLASS control

Test:
 Plastic bending with four elements over thickness



Time = 5
 Contours of Effective Plastic Strain
 min=0, at elem# 21
 max=0.0582193, at elem# 908



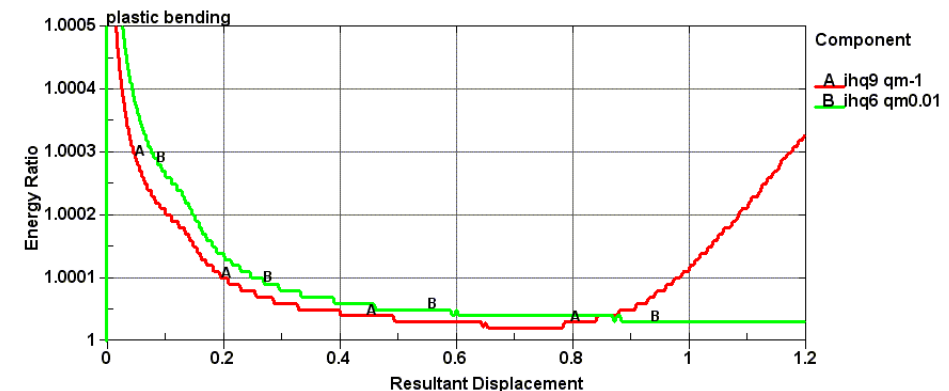
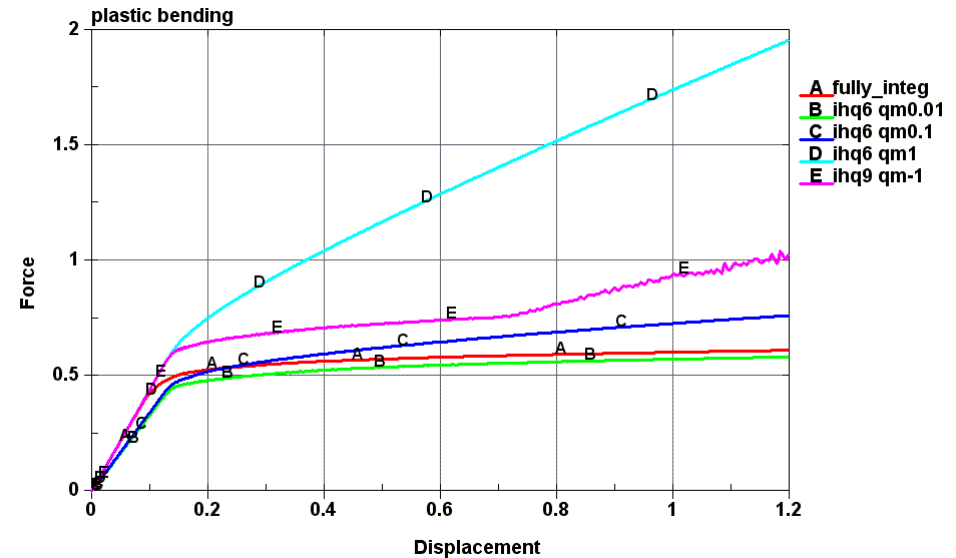
Hourglass control

In the current program version, the Puso Hourglass control sometimes tends to become unstable if element deformation becomes too large. This can be identified by an increasing energy ratio.

Maybe further improvements are necessary.

$$\text{Energy ratio} = \frac{\Sigma E(t)}{\Sigma E(t=0) + W_{\text{ext}}(t)}$$

$$\Sigma E = \text{kinetic} + \text{internal} + \text{hourglass} + \text{damping} + \text{sliding}$$



EFG

EFG
Element Free Galerkin



*ANSYS Conference & 28th CADFEM Users' Meeting 2010
November 3-5, 2010 – Eurogress Aachen, Germany*



EFG

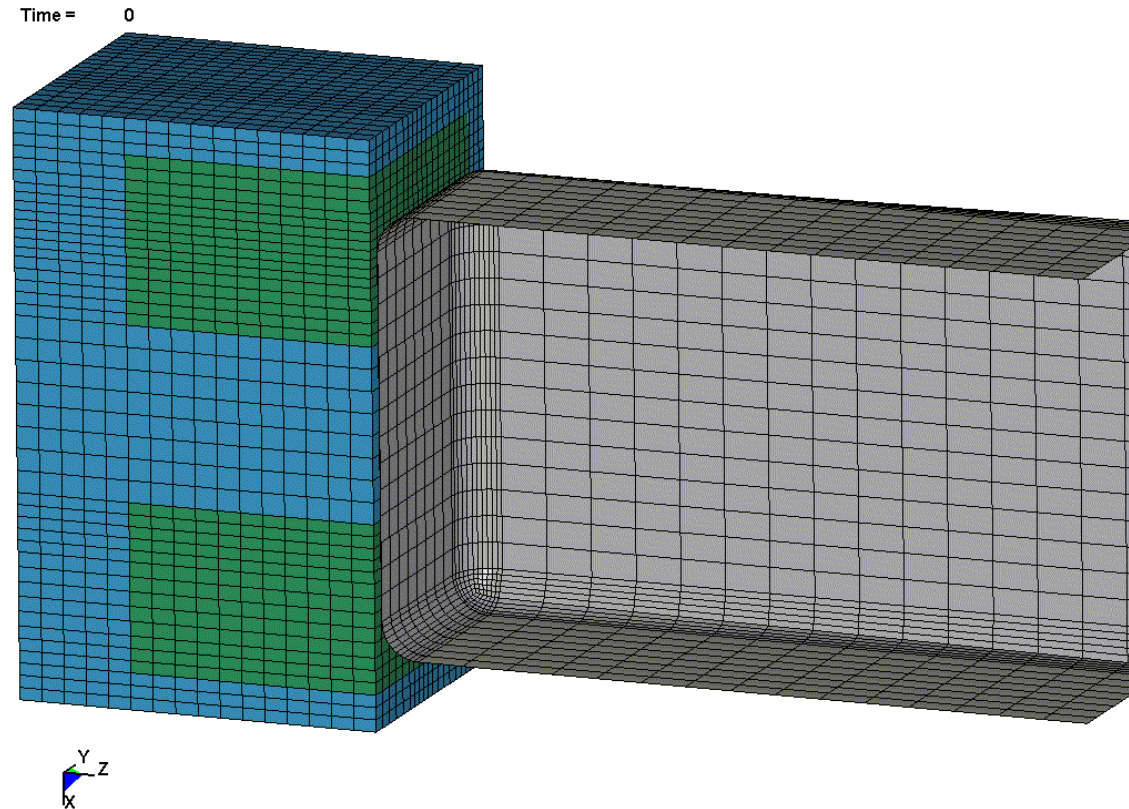
Although EFG is a particle method, it can be used in LS-DYNA like standard elements, because EFG works with a background mesh.

Simply use *SECTION_SOLID_EFG and set ELFORM=41 (hex) or 42 (tet)

EFG parts and FEM parts can share nodes.

EFG Background mesh is only needed in initialization phase, for contact and for postprocessing.

The main advantage of EFG over FEM is robustness and accuracy in large distorted elements. EFG has less danger of „negative volume“.



EFG

Disadvantage of EFG over FEM:

- Larger CPU time
- More memory needed
(automatic memory allocation may fail, define memory by hand)
- a lot of settings possible, results are dependent on these settings
(e.g. DX, DY, DZ influencing CPU time and stiffness)

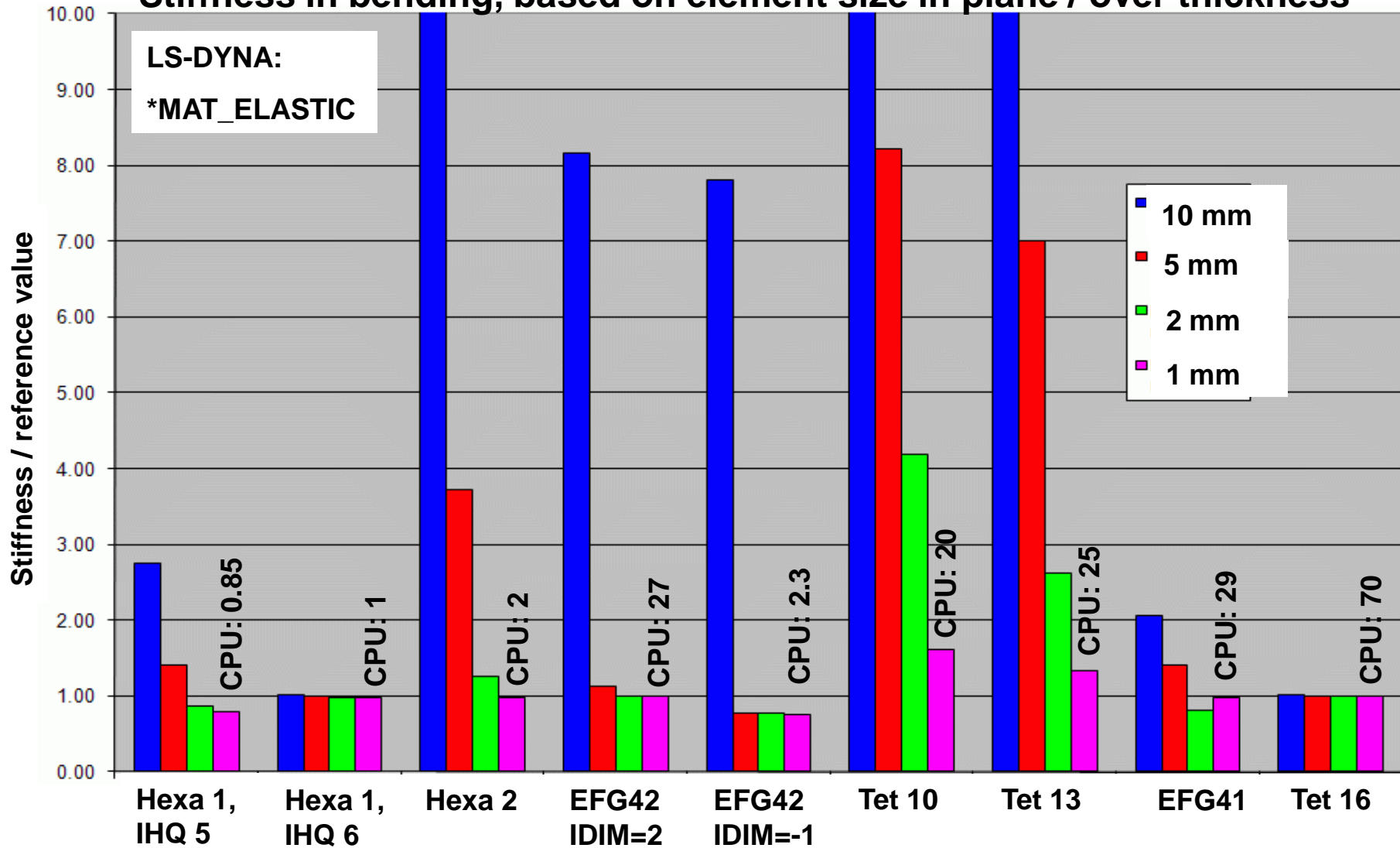
*SECTION_SOLID_EFG

Card 2 and Card 3 define only for the EFG option

Variable	DX	DY	DZ	ISPLINE	IDILA	IEBT	IDIM	TOLDEF
Type	F	F	F	I	I	I	I	F
Default	1.01	1.01	1.01	0	0	-1	2	0.01
Variable	IGL	STIME	IKEN	SF	MID	IBR	DS	ECUT
Type	I	F	I	F	I	I	F	F
Default	0	1.e+20	0	0.0		1	1.01	0.1

EFG for thin walled structures

Stiffness in bending, based on element size in plane / over thickness



EFG

Recomendations

Metal materials in Forging:

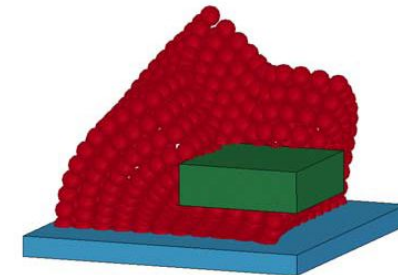
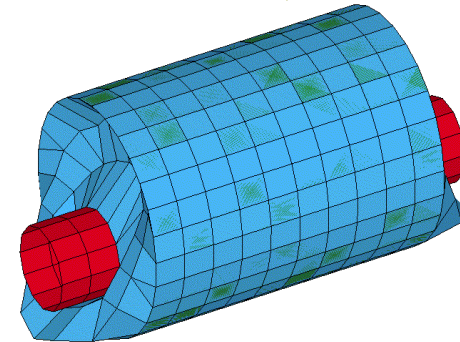
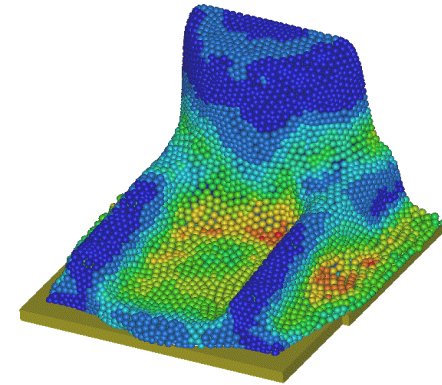
- Lagrangian kernel (TOLDEF=0)
- Maximum Entropy approximation (IEBT=7)
- Two point Gauss integration (IDIM=2)

Rubber materials:

- Lagrangian kernel (TOLDEF=0)
- Maximum Entropy approximation (IEBT=7)
- Stabilized domain integration (IDIM=-1)

Foam materials:

- Semi-Lagrangian kernel (TOLDEF>0)
- Fast transformation (IEBT=4)
- Default domain integration (IDIM=1)

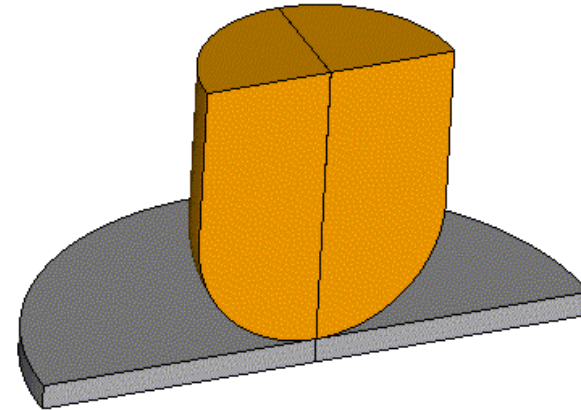


EFG

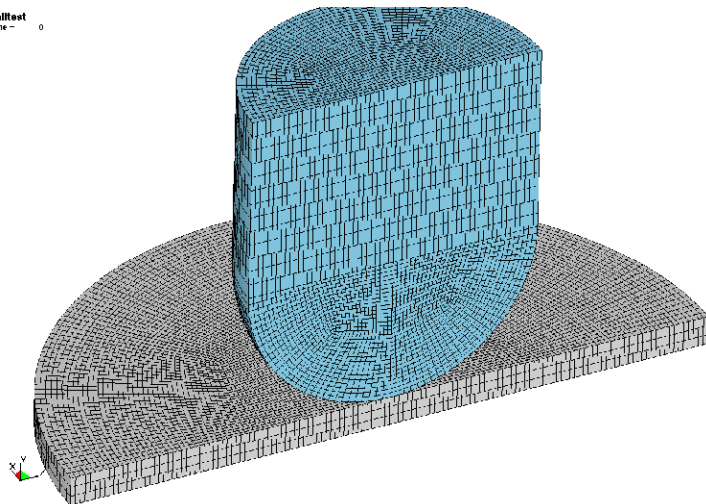
Falltest
Time = 0

Example:
Rubber Impact with rigid projectile

- EFG type 41 (hexahedron)
- Maximum Entropy approximation
IEBT=7
- Stabilized method IDIM=-1 (very fast)
- *MAT_SIMPLIFIED_RUBBER



Falltest
Time = 0

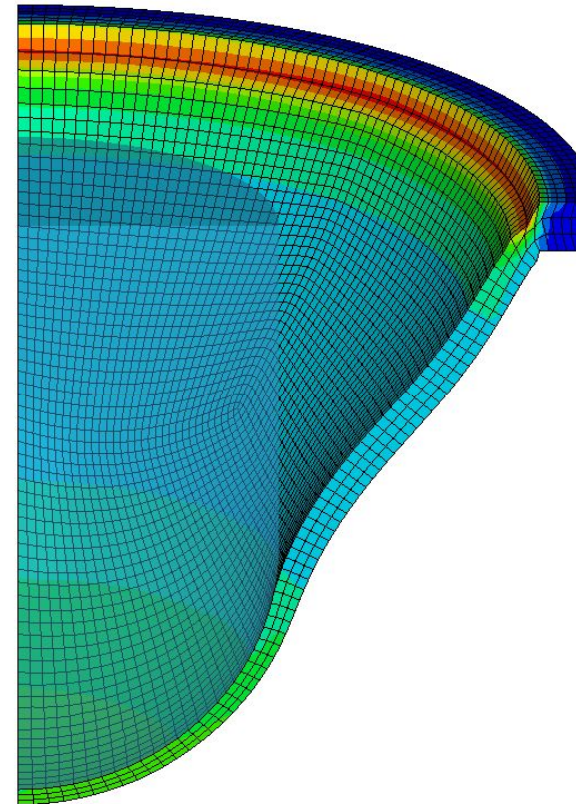
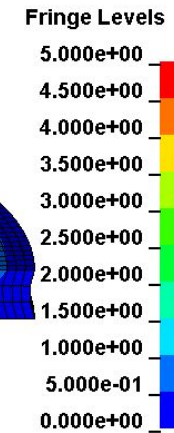
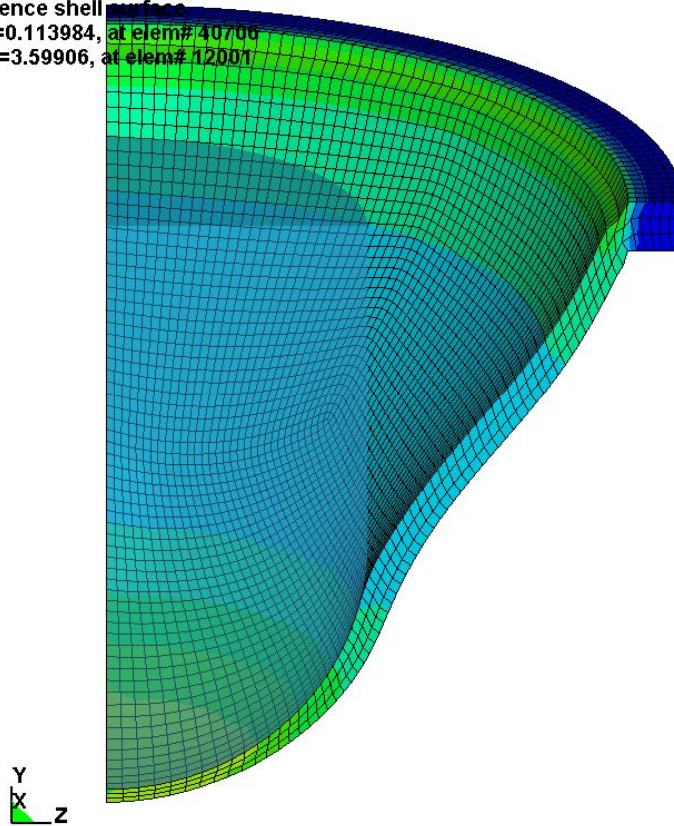


EFG

Comparisons: EFG Type 41

FEM Type 1

Falltest
Time = 0.00079988
Contours of Effective Stress (v-m)
reference shell surface
min=0.113984, at elem# 40706
max=3.59906, at elem# 12001

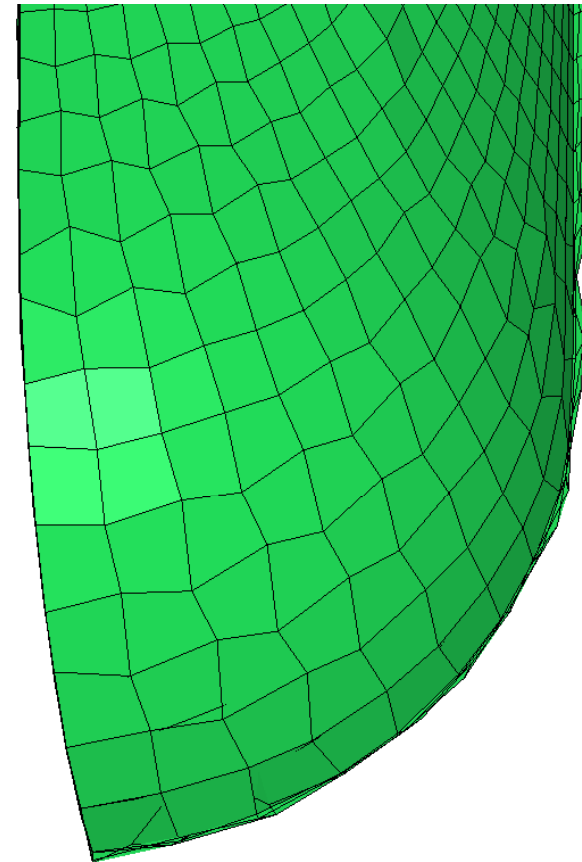
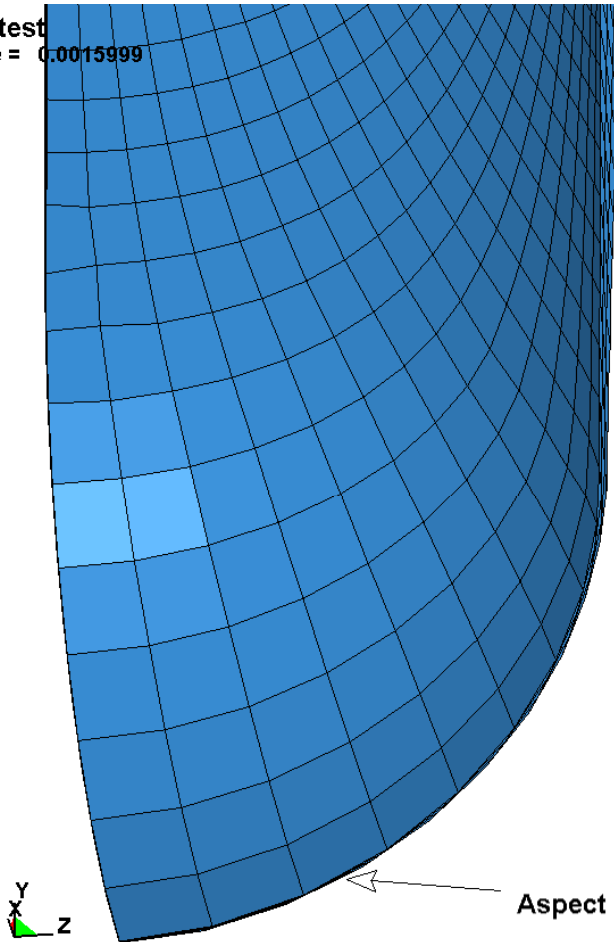


EFG

Comparisons: EFG Type 41

FEM Type 1 (last state)

Falltest
Time = 0.0015999



Aspect Ratio=1:100

ALE

ALE
Arbitrary Lagrangian Eulerian



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CADFEM

ALE

Lagrangian- formulation:

- is used in structure mechanics
- material and elements are bonded together
- large deformation induces element distortion

Eulerian- formulation:

- is used in fluid mechanics
- mesh of elements is fixed in space
- material 'flows' through the elements
- variable boundary conditions are complicated

ALE: Arbitrary- Lagrangian- Eulerian:

- both formulations in combination: the mesh can move and deform, the material can flow through the mesh
- two possible kinds of applications:

à REZONING: large deformation in structure mechanics;
mesh must be corrected

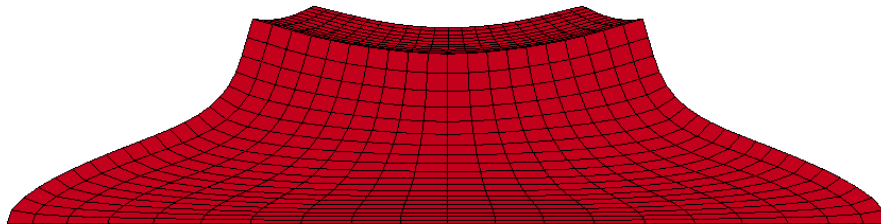
à FLUID-STRUCTURE-INTERACTION: Airbag inflation, Tank sloshing



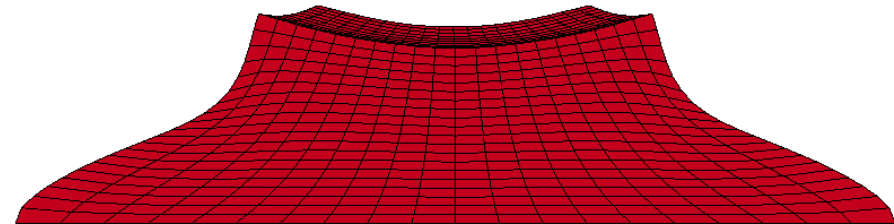
ALE

ALE :=Material(s) can flow through the mesh which itself can move

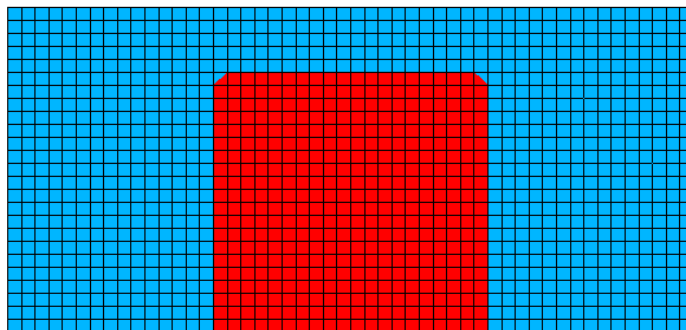
Lagrangian formulation



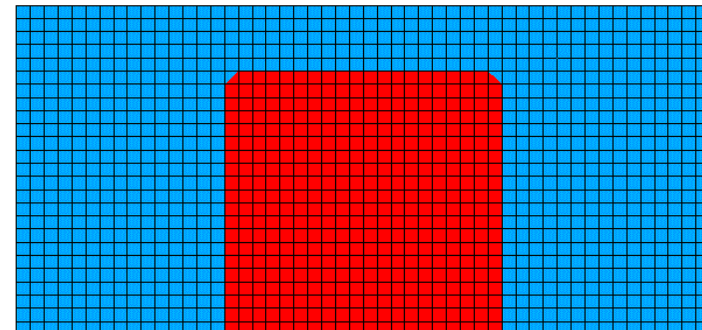
Single material ALE



Multi-material Eulerian



Multi-material ALE formulation



ALE hexahedron (type 5)

- 8 noded hexahedron element with trilinear shape function with reduced integration
- single material Arbitrary Eulerian Lagrangian element
 - à coupling of lagrangian and eulerian formulation: materials flow through elements
- useful for simulations with large element distortion but remeshing not necessary
- approx. 2 times slower than a lagrangian element elform=1

Single material and void ALE - (type 12)

- Eulerian element which can be filled by one material or void
- use in combination with void-definition (12)

Multi material ALE - (type 11)

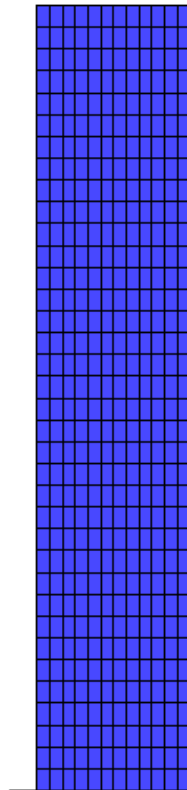
- Eulerian element which can be filled by an arbitrary number of materials
- use in combination with *ALE_MULTI-MATERIAL_GROUP (11)
- approx. 2.5 times slower than a lagrangian element elform=1

Testcase: Taylor beam

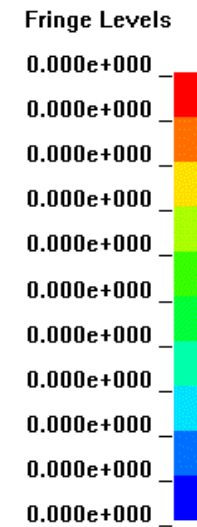
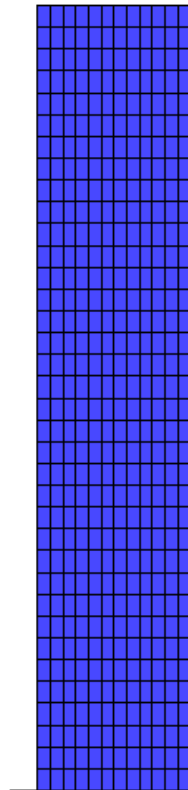
Hexahedron elform=1

Single material euler
element elform=5

TAYLOR BAR
Time = 0
Contours of Effective Plastic Strain
max ipt. value
min=0, at elem# 1
max=0, at elem# 1



TAYLOR BAR
Time = 0
Contours of Effective Plastic Strain
max ipt. value
min=0, at elem# 1
max=0, at elem# 1



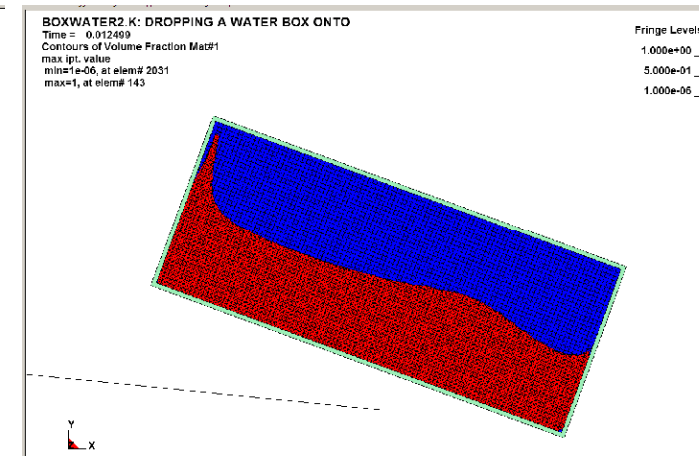
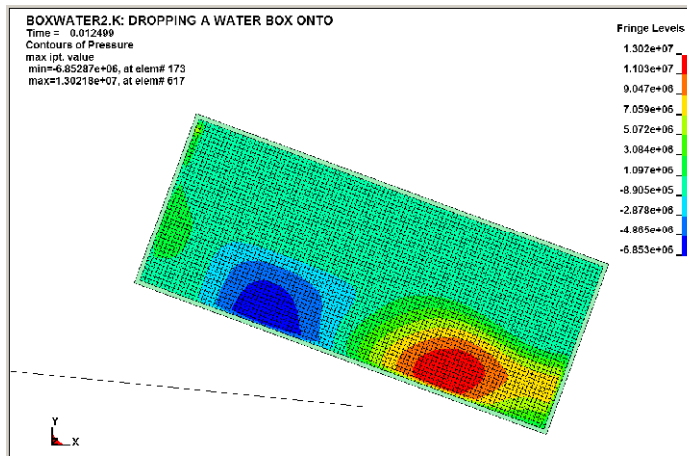
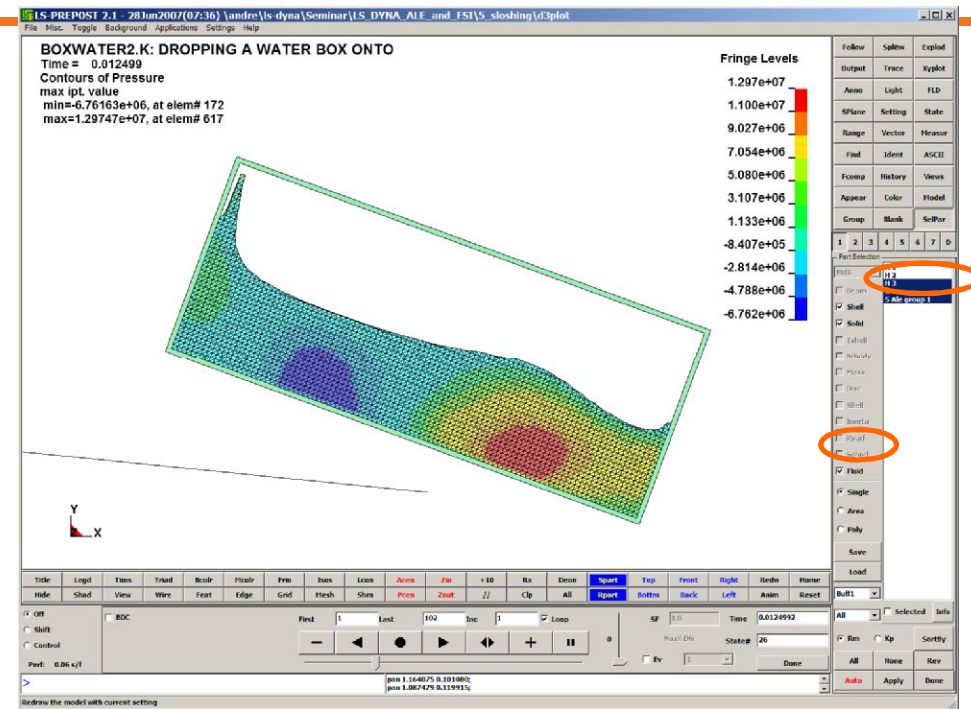
ALE

Postprocessing

Display of interface reconstruction for multi material euler elements

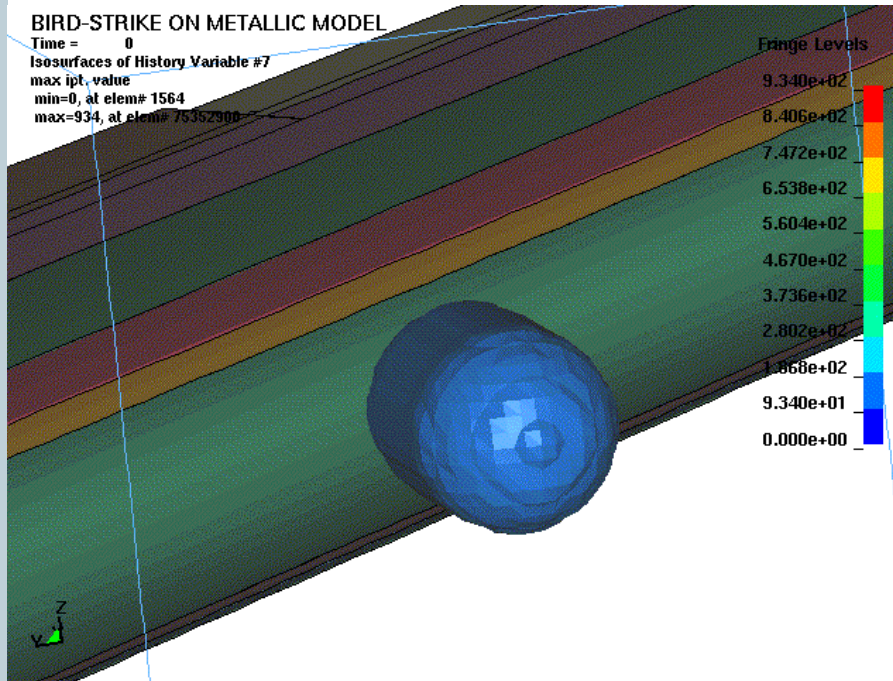
Part selection: Fluid
à additional ALE parts

- reconstruction of geometry, mesh with tetrahedrons
- results on ALE material outer surface

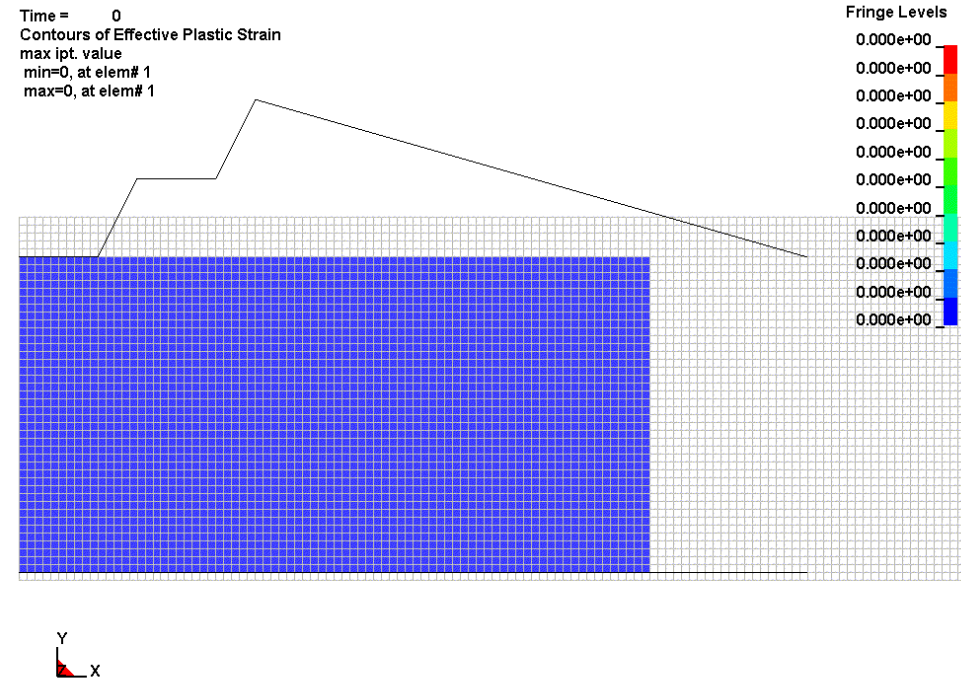


Examples for multi material euler:

Bird strike



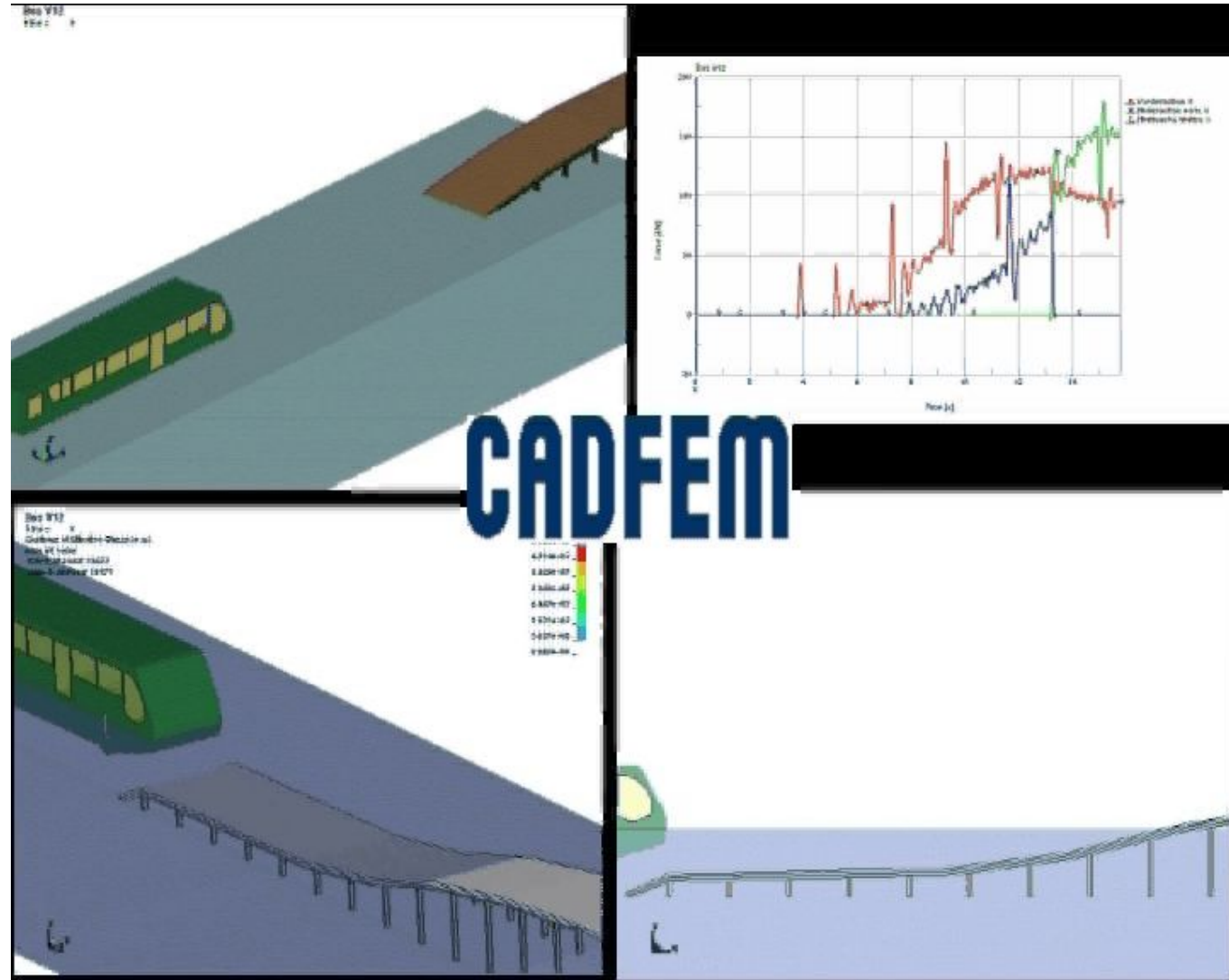
2d forging



ALE

Example

Fluid structure interaction.
Eulerian elements for water.



CADFEM

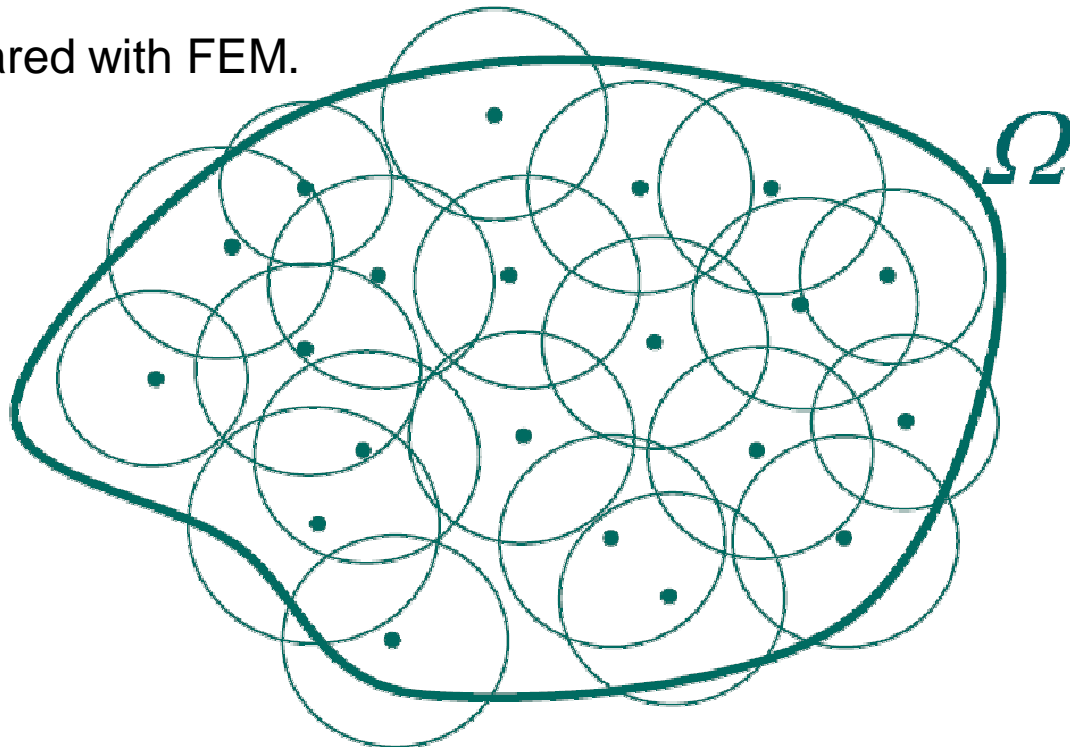
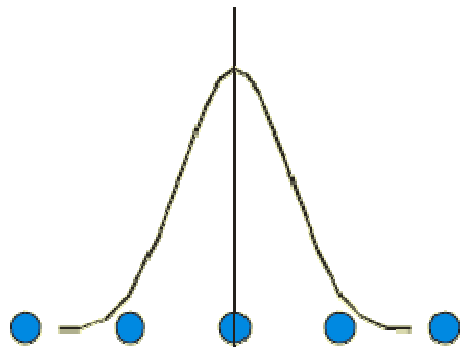
SPH

**Not really Elements, only particles
- but defined with *ELEMENT_SPH**

SPH Smooth Particle Hydrodynamics

SPH

- A collection of nodes (particles) discretize a continuum.
- These are mass points that describe the motion of the continuum.
- SPH is a lagrangian collocative method, i.e. the nodes and the integration points are the same. Shape functions are centered on the particles.
- Accuracy is not high compared with FEM.



SPH

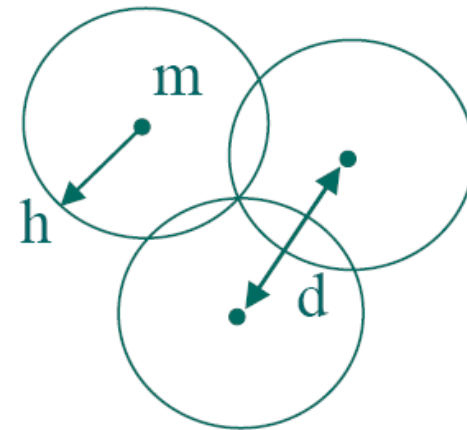
Particle has mass m – has to be defined independently from density

2 parameters of discretization:

d is distance between particles

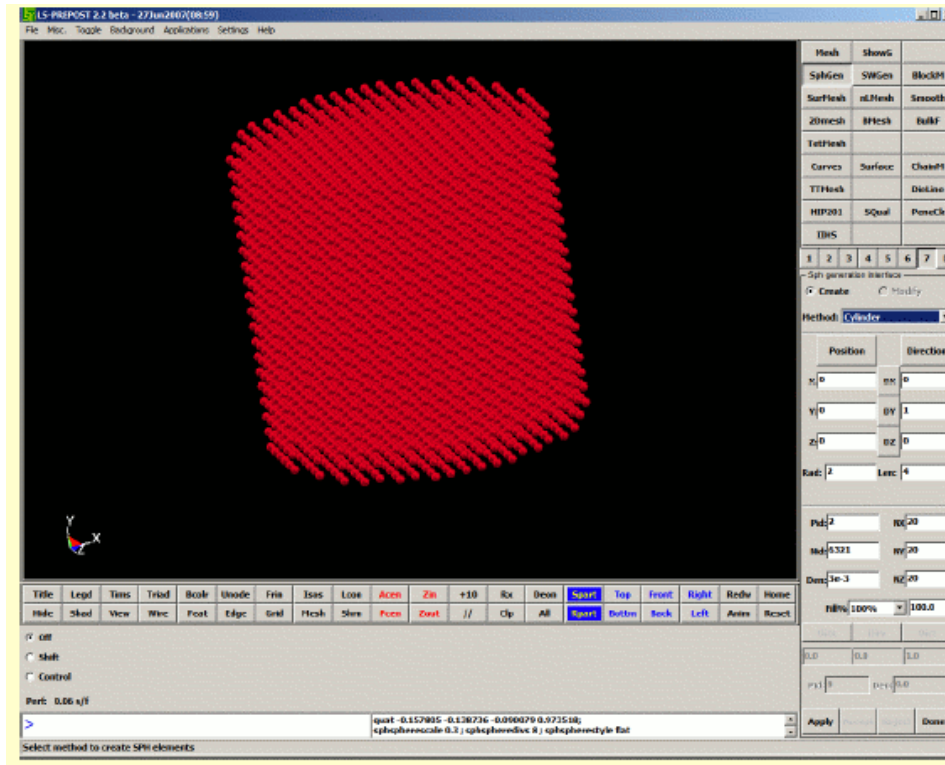
h is smoothing length

– different than classical methode

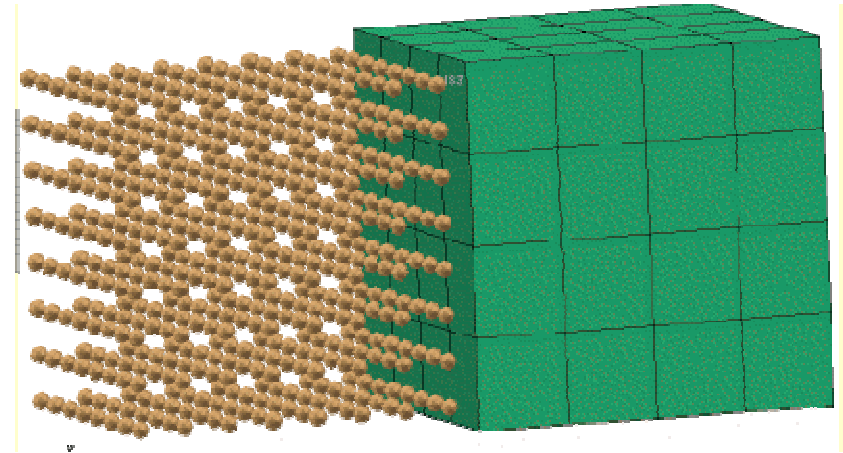


SPH

- Minimum of four particles per one solid element face
- SPH mesh should always be regular



Meshing



SPH meshing in LS-PREPOST:
Page 7 à SphGen

Computational cost:

- Time step size is comparable to solidelements
- CPU time approx. 2.5-3 times slower than Hexahedron type 1

Regular contact with:

- *CONTACT_AUTOMATIC_NODES_TO_SURFACE
- SPH nodes on the slave side

Tied contact:

- *CONTACT_TIED_NODES_TO_SURFACE
- SPH nodes on the slave side
- Note: SPH nodes do not have any thickness in contact!

Contact between SPH parts:

- see *CONTROL_SPH, parameter CONT

Other recommendations:

- reduce time step scale factor (*CONTROL_TIMESTEP, tssfac<0.8
- do not use mass scaling
- use a SPH box (*CONTROL_SPH, boxid), particles flowing outside of this box not longer computed
- use negative MEMORY value (*CONTROL_SPH) is often more stable
- look in the material section of Keyword users manual for valid material model

SPH

Postprocessing

File Misc. Toggle Background Applications Settings Help

TEST OF MATERIAL 63 - CRUSHABLE FOAM
Time = 67.996
Contours of Effective Stress (v-m)
max ipt. value
min=0, at node# 1001003
max=1.79938e-05, at node# 1000045

Fringe Levels
1.799e-05
1.619e-05
1.440e-05
1.260e-05
1.080e-05
8.997e-06
7.198e-06
5.398e-06
3.599e-06
1.799e-06
0.000e+00

Follow Splitw Explod
Output Trace Xyplot
Anno Light FLD
SPlane **Setting** State
Range Vector Measur
Find Ident ASCII
Fcomp History Views
Appear Color Model
Group Blank SelPar

1 2 3 4 5 6 7 D
Set Display Options
 Axes/Surface
 Disp. scale factor
 Reflections
 Hic/Csi const.
 SPH nodes
 Thickness Scale Factor
 Fringe Scale Factor
 Concrete Crack Width
 Local Coord System
 FLD E'Strain
 Local B'Strain

SPH Radius Scale:
0.4 0.3

SPH Sphere divs:
12 12
Style: flat
 Fixed Radius

Aply Clr Rset Done

Title Legd Tims Triad Bcolr Mcolr Frin Isos Leon Acen Zin +10 Rx Deon Spart Top Front Right Retw H...
Hide Shad View Wire Feat Edge Grid Mesh Shrn Peen Zout // Clp All Rpart Bottm Back Left Anim Reset

◆ Off BDC First 1 Last 152 Inc 1 Loop
◆ Shift
◆ Control
Perf: 0.00 s/f

SF 1.0 Time 67.996
No. of Div State# 69
 Ev 1 Done

quat 0.250055 -0.481241 -0.752187 0.374292

SPH

Examples

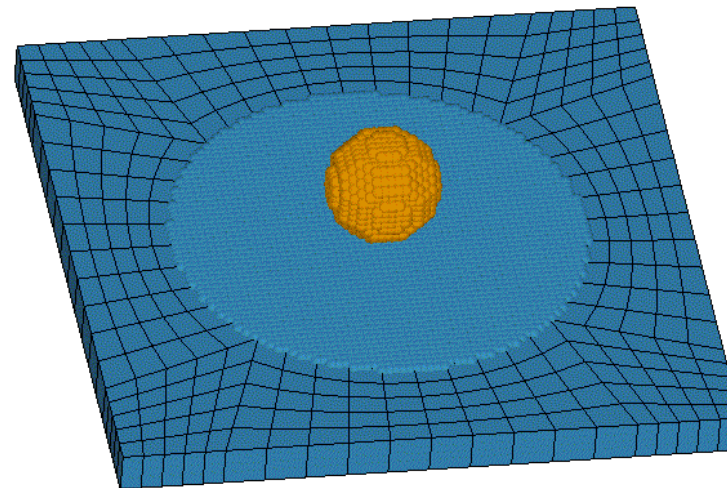
IMPACT 6.18 KM/S ALU/ALU
Time = 0

Testmodel:

High velocity impact:
 $v=6.18$ km/s

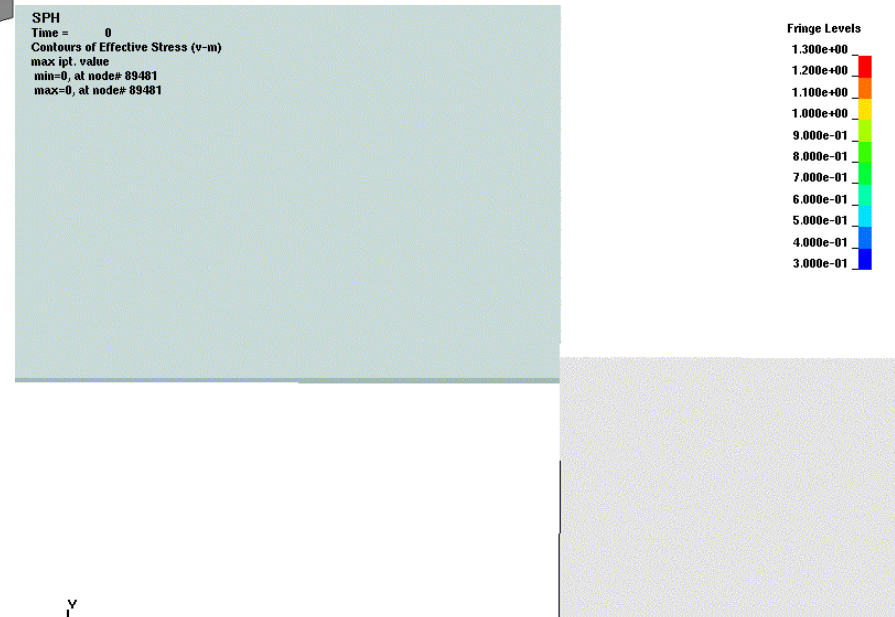
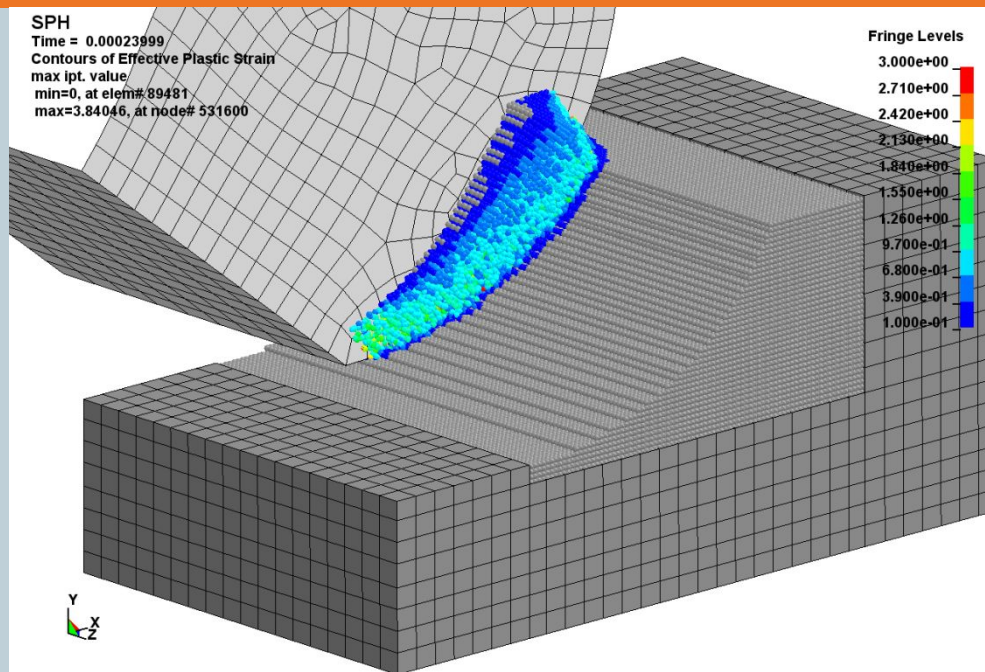
19.000 SPH elements
300 Solid Elements
600 Cycles

2min on this laptop



SPH

Examples



Testmodel:

- 130.000 SPH elements
- 5.000 Solid Elements
- 60.000 Cycles

4h 30min on a quad core CPU



CPM

**Looks like SPH but is completely
different:**

-

CPM

Corpuscular Particle Method

for gases only



CPM

Basic idea: Model the gas as a set of rigid particles in random motion

Advantages:

- no field equations – easy and numerically robust
- contact with solid parts simple with *contact_nodes_to_surface

The theory describes the interaction between gas molecules on a microscopic level.

Original application: Airbag deployment

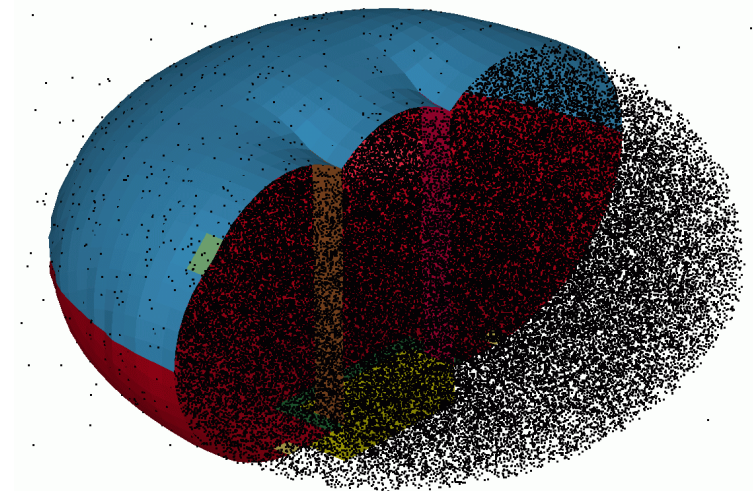
Problem:

A deployed airbag is filled up by roughly 10^{24} molecules.

Solution:

Reduce system from many molecules to a “few” particles

Now a airbag is filled by 500.000 particles



CPM

*airbag_particle

IAIR=2: Initial gas in Bag

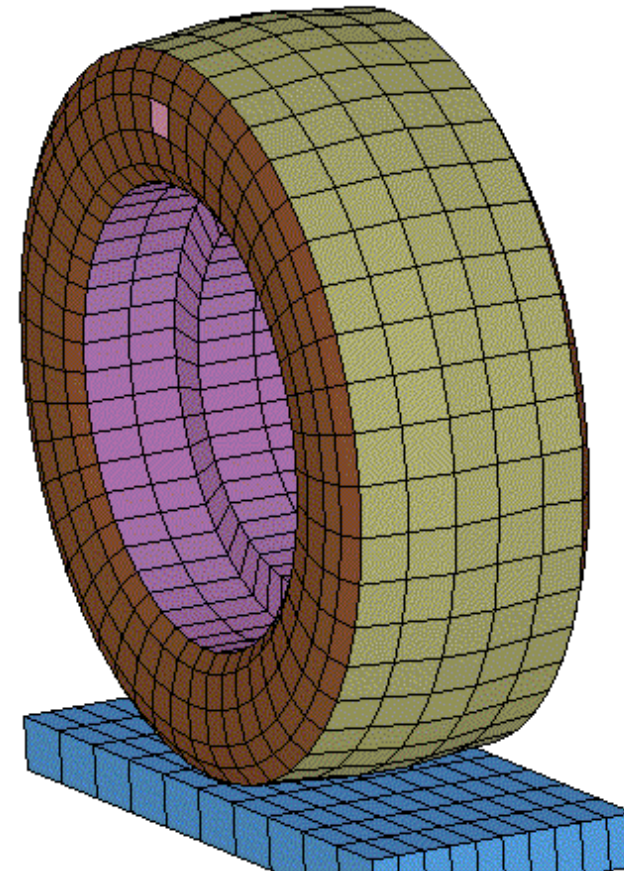
Application:

Initialize pressure in a closed volume

- Airbags
- Door cavity
- Tires

This makes the method open for much more than airbags.

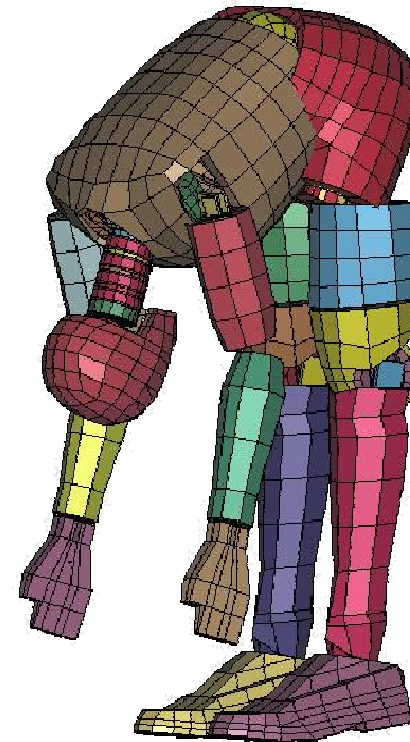
tire leak using particle method
Time = 0



Kontakt

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Vielen Dank für Ihre Aufmerksamkeit!

