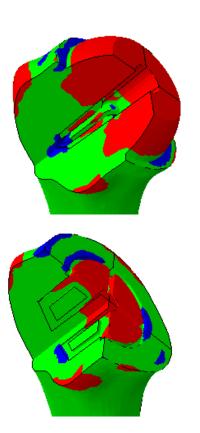


Tissue Biomechanics

2º semester, 2017-2018



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Introduction

Material properties

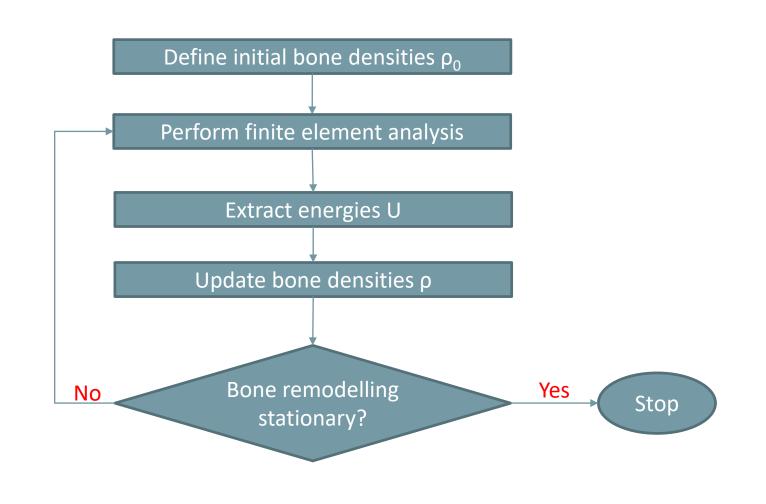
History output

Loads

Bone densities

Job

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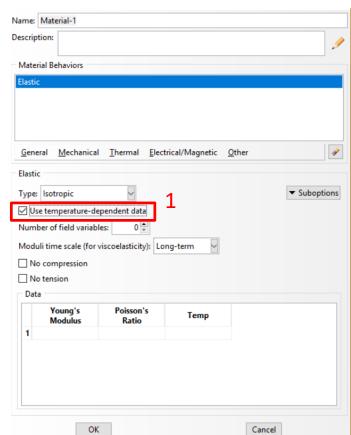
Bone densities

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- **Bone densities** are defined at the **nodes** of the finite element mesh as if they were temperatures.
- Defining bone densities as temperatures simplifies the definition of the bone material properties as Abaqus allows the definition of temperature-dependent properties (1).
- A table describing how the material properties of bone change with bone density (temperature) needs to be defined.

$$E = E_0 \rho^p$$







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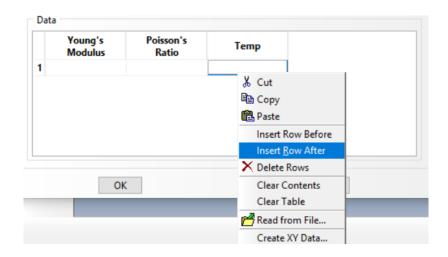
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 To create a table, add rows to the data section by right-clicking the mouse over this section and selecting Insert Row After.



- As the material law for bone ($E = E_0 \rho^p$) cannot be defined analytically in Abaqus, several data points of the law need to be defined to better approximate the function.
- Between data points, Abaqus interpolates the data using linear interpolation.





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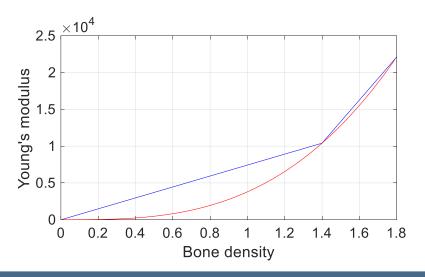
Python script

Bone remodeling implementation

• For example, if three points were used,

Modulus Ratio	Temp	
0.1 0.3	0.001	
2 10000 0.3	1.4	
20000 0.3	1.8	
20000 0.3	1.8	

- Abaqus would define the Young's modulus of bone using the blue curve instead of the true curve in red.
- Notice that the bone material properties are defined at this point but the bone densities throughout the bone still have to be defined.







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Huiskes' bone remodelling law is given by:

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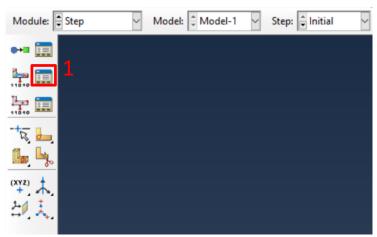
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$$\frac{d\rho}{dt} = \begin{cases} B\left(\frac{U}{\rho} - (1-s).k\right), & if \frac{U}{\rho} < (1-s).k\\ B\left(\frac{U}{\rho} - (1+s).k\right), & if \frac{U}{\rho} > (1+s).k\\ 0, & otherwise \end{cases}$$

• By default, Abaqus does not report the strain energy density U. To report U, the Field Output needs to be edited using the Field Output Manager (1) in the step module.







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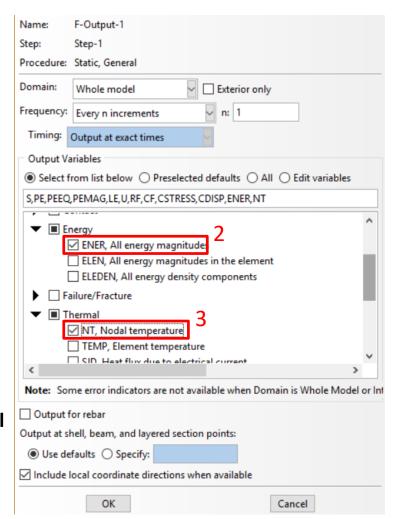
Python script

Bone remodeling implementation

• Click edit (1):



- Select the ENER, All energy magnitudes
 (2) option.
- To be able to see the bone density distribution (temperature), the NT, Nodal Temperature (3) option also needs to be selected.







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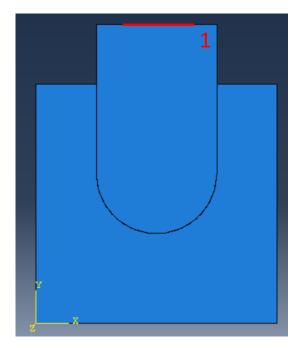
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• If a concentrated force is to be distributed over a certain region (for example, region (1)), the force can be applied on an attachment point and be distributed to the region using a coupling constraint.



 An attachment point is a point that does not belong to any part but may transmit force to the model.





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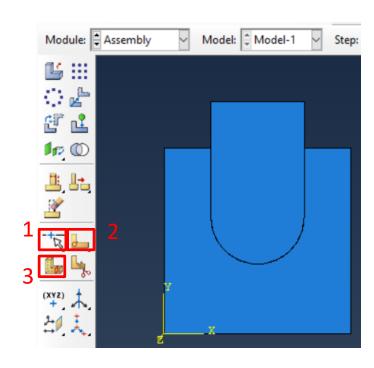
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• To define the **region for force distribution**, **partitions** may have to be created using the **Partition edge: Specify Parameter by Location (1)** at the assembly module.



Partitions on surfaces and solids can be created using the features Partition Face
 (2) and Partition Cell (3), respectively.





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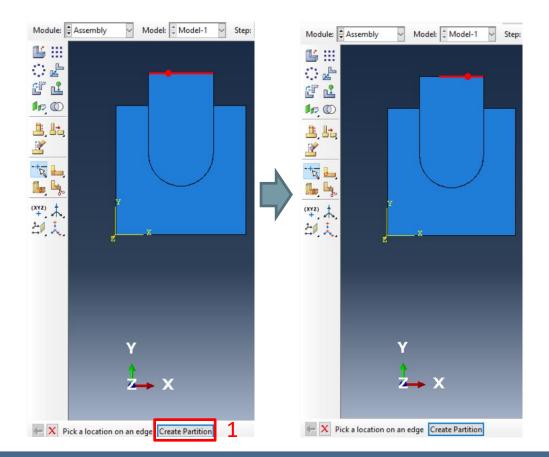
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- Using the **Partition edge: Specify Parameter by Location** tool, select the first point of the region to split the edge into two edges and click on **Create Partition (1)**.
- Repeat the procedure for the opposite extremity of the region.







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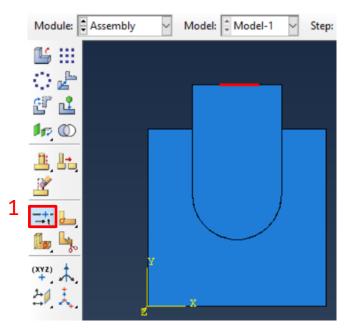
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- Considering that the attachment point should be placed at the centroid of the region, a node at the middle of the region may be created to help define the attachment point.
 - For that, select the tool **Partition Edge: Enter Parameter (1)**, which appears after clicking a few seconds on the Partition Edge tool.
- After selecting the edge, click **Done**.
- Then input 0.5 and click Create
 Partition. Notice that t = 0 is the first point of the edge and t = 1 is the last.











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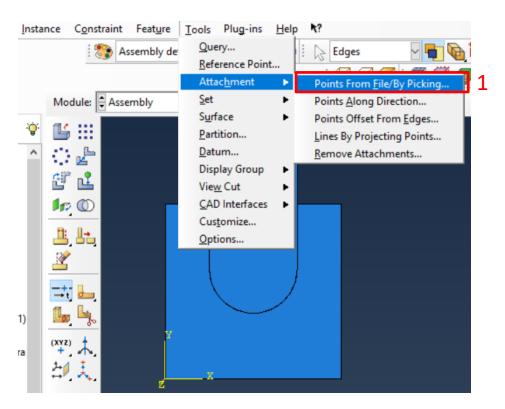
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To create an attachment point, click Tools -> Attachment -> Points From File / By Picking... (1).



Other, intuitive, options are also available to define attachment points.





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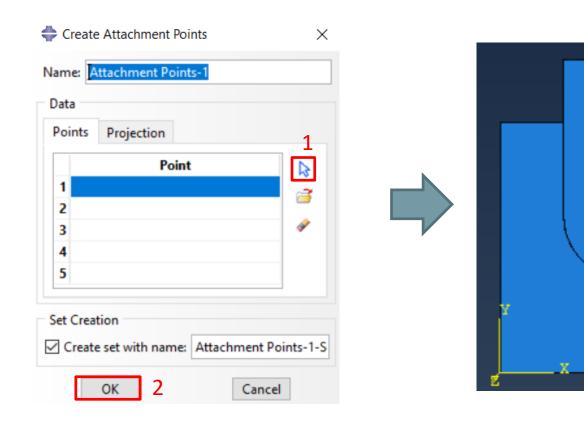
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• Then click on the arrow (1), select the middle node of the edge, and click OK (2).



The attachment point appears as a green plus sign.





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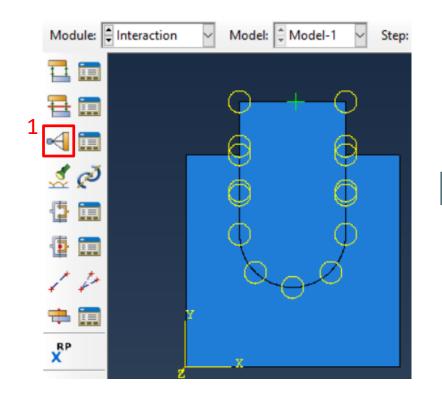
Bone densities

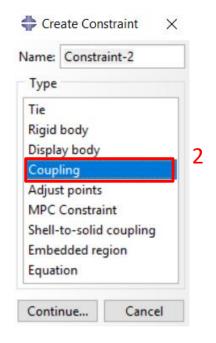
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• To define a **coupling interaction** between the attachment point and the region, go to the **Interaction module**, click on **Create Constraint (1)**, and select **Coupling (2)**.









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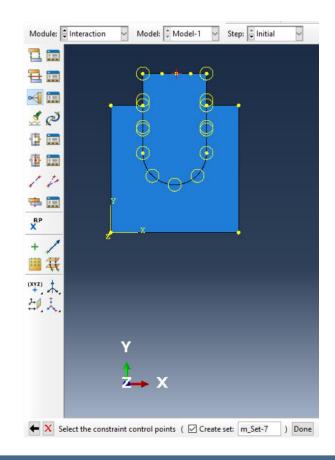
Bone densities

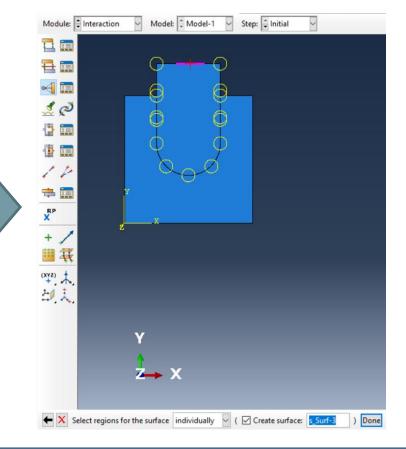
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• Select the attachment point and click Done. Then, select Surface as the region type, identify the region, and click Done.









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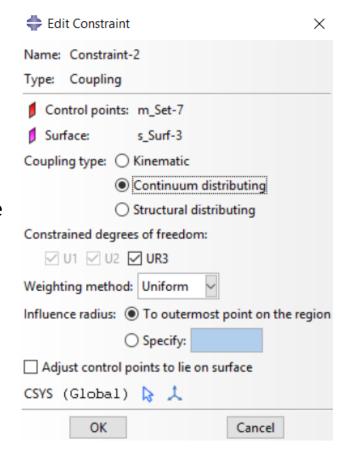
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- The **coupling type** must be **Continuum distributing.** The remaining options may be the default.
- The Weighting method defines how the force is distributed.
 - Uniform means the force is distributed uniformly over the region.
 - The remaining options decrease the magnitude of the forces applied on the nodes according to their distance to the attachment point.







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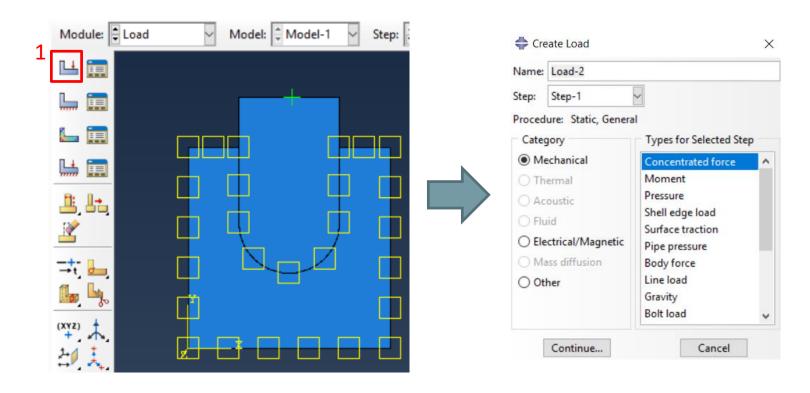
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 Once the coupling interaction between the attachment point and the region is defined, the concentrated force should be defined at the attachment point using the Create Load (1) tool at the Load module.







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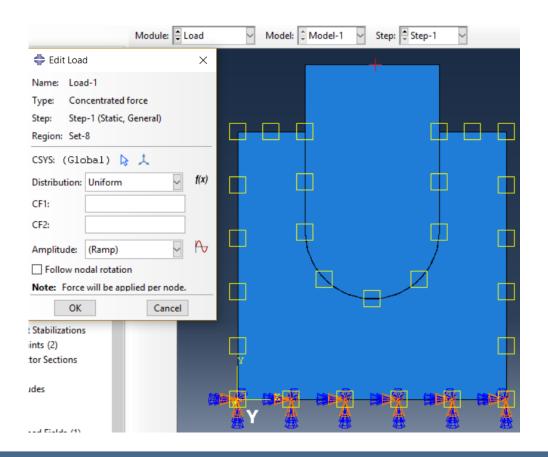
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After selecting the attachment point, input the components of the force (CF1 and CF2).







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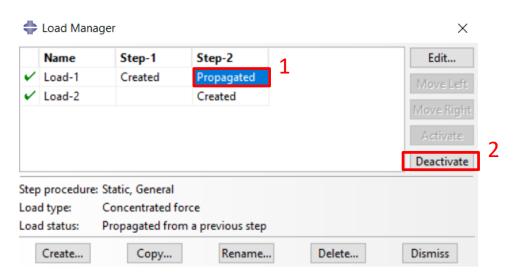
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Bone remodeling implementation

• If more than one load case exists, you can define different steps, one for each load case.



- By default, Abaqus **propagates the loads** from one step to the **following steps** (notice the propagated instruction in (1)).
- To prevent Abaqus from propagating the loads along the steps, click on the first **Propagated instruction** for the load case and click on **Deactivate (2)**.





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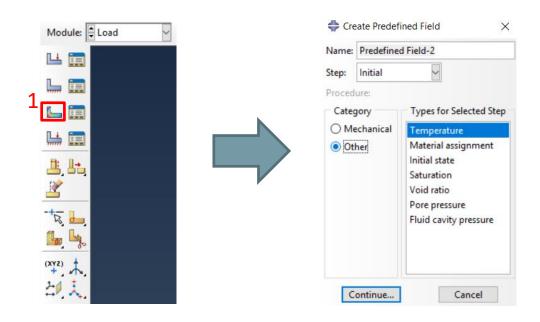
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• To define how the **bone density changes** within the bone, a **predefined field** needs to be defined (1).



 As bone density will be simulated as temperature, select Category Other and Type Temperature. For the step, select Initial.





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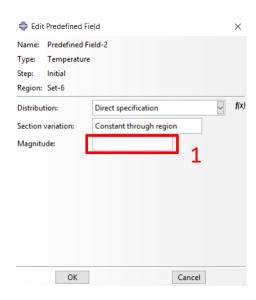
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• After selecting the regions for the field, i.e., the bone, define the bone density (temperature) at (1).



- At the **Abaqus interface**, only a **constant bone density** can be defined within the bone.
- A variable bone density distribution needs to be defined later by changing the
 instructions of the input file regarding the definition of the constant predefined
 field.





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• To change the input file, search for the keyword **Initial**:

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** Name: Predefined Field-1 Type: Temperature

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*Initial Conditions, type=TEMPERATURE

Set-5, 1.

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 To define a variable bone density distribution, to be read from a text file called dens.txt, change the line following the instruction *Initial, i.e.:

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** Name: Predefined Field-1 Type: Temperature

*Initial Conditions, type=TEMPERATURE

*Include, input=dens.txt





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 The density file should contain in each line the bone density for each node of the finite element mesh using the following structure:

NameOfInstance.NumberOfNode, NodalDensity

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where NameOfInstance is the name of the instance, NumberOfNode is the number of the node, and NodalDensity is the density at the node NumberOfNode.

• For example:

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Part-3-1.1, 0.631353

Part-3-1.2, 0.256001

Part-3-1.3, 0.050299

Part-3-1.4, 0.175944

Part-3-1.5, 0.141576

Part-3-1.6, 0.104506

Part-3-1.7, 1.800000

Part-3-1.8, 0.631713





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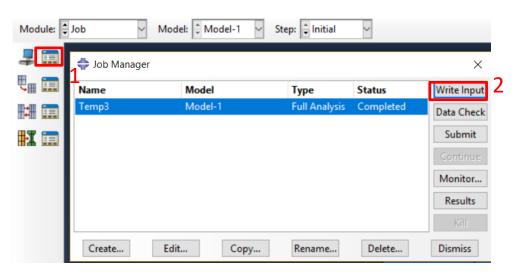
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• Once the finite element model is defined, the *.inp file needs to be written. For that, click on the Job Manager (1) -> Write Input (2) in the Job module.



• To perform a **finite element analysis in Matlab**, use the following instruction:

system('abaqus job=NameInpFile inter');

where **NameInpFile** is the **name of the finite element model** (Temp3 in the example above).





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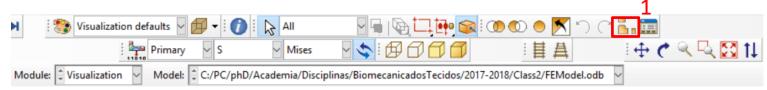
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- Once a finite element analysis is completed, the strain energy densities need to be extracted from the output file. For that purpose, a python script needs to be created to have the strain energy densities reported automatically into a text file.
- To create the python script, complete a finite element analysis of the model developed once and open the corresponding *.odb file.
- Once the *.odb file is open, Create a Display Group (1) to make bone the only part visible in the results. If only the bone is visible, the text file will only contain data for this part.



To obtain a text file with the strain energy densities at the nodes, click Report -> Field
Output (2).







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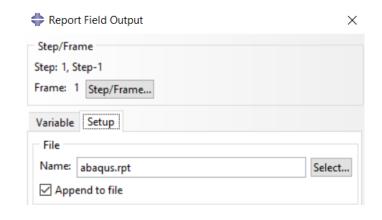
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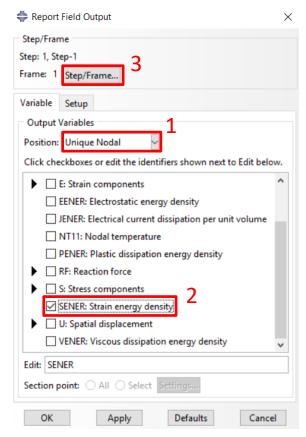
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- Then select Position Unique Nodal (1) and check the box SENER: Strain energy density (2).
- The results are reported for a specific step. You can select the step wanted by clicking on Step/Frame (3).
- At the **Setup** tab, the name of the ouput file can be defined, as well as the option to append the data (or not).









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- Once the report is created, **close Abaqus**. At the working folder, there should be an **abaqus.rpy** file that contains all the instructions performed.
- Rename the file abaqus.rpy to OutputExtraction.py, which will be used in Matlab to create automatically an output file with the strain energy densities from the finite element analyses.
- To run the Python script in Matlab, use the following instruction:

system('abaqus viewer noGUI=OutputExtraction.py');

• Once the output file is created, the strain energy densities need to be read from the file to evaluate the bone remodelling law.





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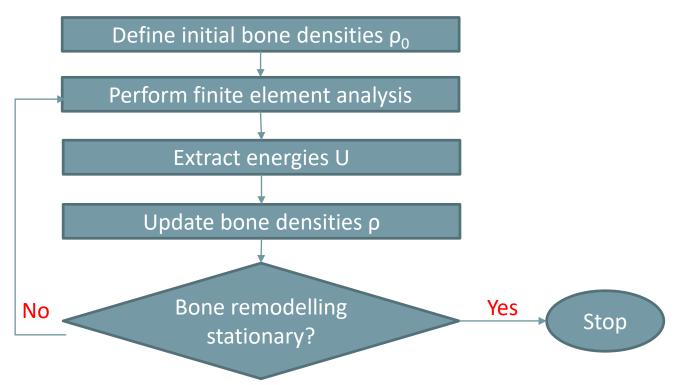
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• Using the tools and procedures described, the bone remodelling process needs to be implemented in Matlab.



 If more than one load case exists, caution is needed to extend the procedures described to such conditions.



