Mathematical Modelling of Aluminum Reduction Cell Potshell Deformation

Marc Dupuis





Plan of the Presentation

- Introduction
 - Dewing stress-strain relationship of the sodium swelling
 - Dewing sodium diffusion coefficient
 - Historical models development background
- "Empty Shell" Potshell Model
 - Elastic mode
 - Plastic mode
- "Almost Empty Shell" Potshell Model
 - Elastic mode
 - Plastic mode
- "Half Empty Shell" Potshell Model
 - Elastic mode
 - Plastic mode
- Conclusions



Introduction



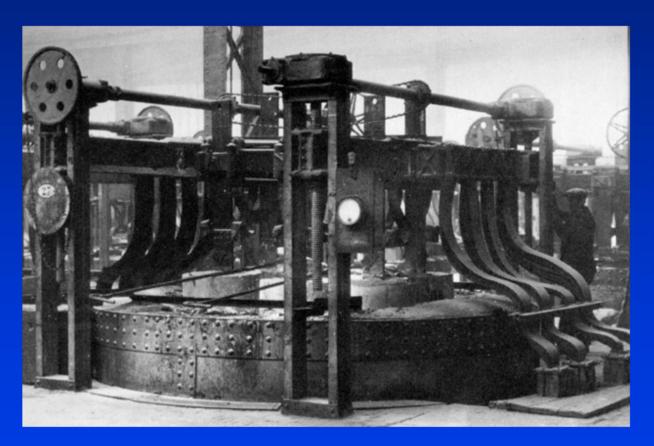
A well-designed shell is supposed to withstand the internal forces induced by thermal and chemical changes in the reduction cell.

The thermal changes are quite straightforward to assess if a complementary thermo-electric model is available.

The chemical changes is the consequence of sodium penetration into the carbon cathode blocks. This sodium penetration makes the carbon swell and hence, the cathode blocks expand chemically.



Introduction



One quite acceptable technical solution:

If the potshell want to be round, why not build it round in the first place?



Dewing Stress-Strain Relationship of the Sodium Swelling

The physic of that chemical expansion is not well understood. The key references on that subjects are Rapoport and Dewing.

According to Dewing, the stress-strain relationship of the sodium swelling of the carbon is:

$$\varepsilon = \varepsilon_0 10^{-k\sigma}$$

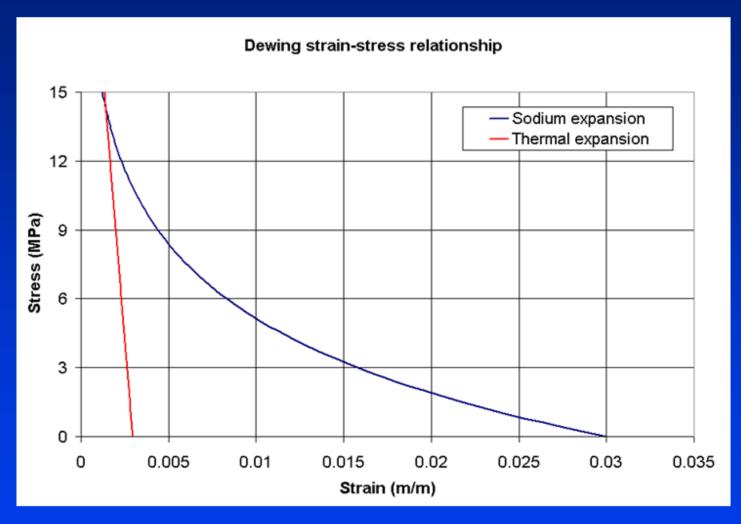
where: \mathcal{E} is the carbon strain at equilibrium

 $\boldsymbol{\mathcal{E}}_0$ is the carbon free expansion strain

σ is the compressive stress in the carbon k is a constant



Dewing Stress-Strain Relationship of the Sodium Swelling

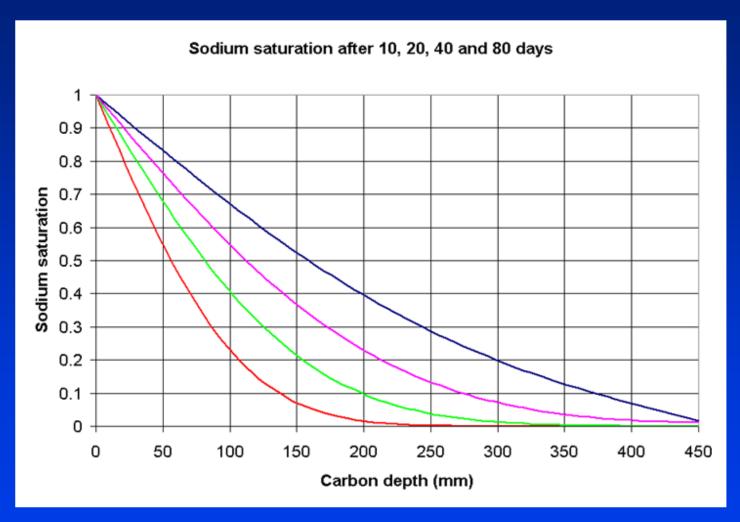


Dewing has estimated k to be equal to 6.4E-4 when expressed in PSI, which corresponds to 0.092825 when expressed in MPa.

On the other hand, the carbon free expansion strain varies a lot depending on the carbon grade and quality. 3% is a typical value for 20% semigraphitic cathode grade, which is about ten times more that the thermal expansion.



Dewing Sodium Diffusion Coefficient



Dewing has also established the sodium diffusion coefficient to be equal to 4E-5 cm2/s, which corresponds to 3.456E-4 m2/day.

Corresponding sodium saturation for a typical 45 cm thick cathode block for the full block thickness after 10, 20, 40 and 80 days of operation.



Dewing Sodium Diffusion Coefficient

Curves presented in previous slide can be obtained by numerically solving the diffusion partial differential equation:

$$\frac{\delta^2 C}{\delta x^2} = \frac{1}{D} \frac{\delta C}{\delta t}$$

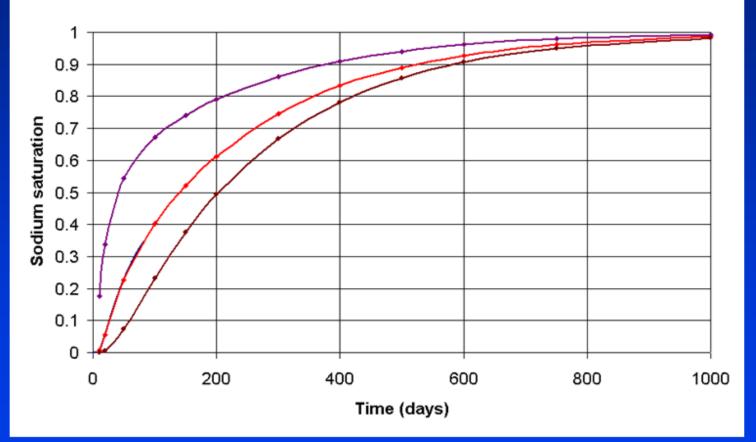
Alternatively, it is possible to use the analytical solution of that diffusion equation also presented by Dewing:

$$C = C_0 \left\{ 1 - \frac{4}{\pi} \sum_{n=0}^{\infty} \frac{1}{(2n+1)} \sin \left[\frac{(2n+1)\pi x}{2h} \right] \exp \left[-\left(\frac{(2n+1)\pi}{2n} \right)^2 Dt \right] \right\}$$



Dewing Sodium Diffusion Coefficient

Sodium saturation in cathode block as function of time and depth at 1/4 1/2 and 3/4 of the cathode block thickness

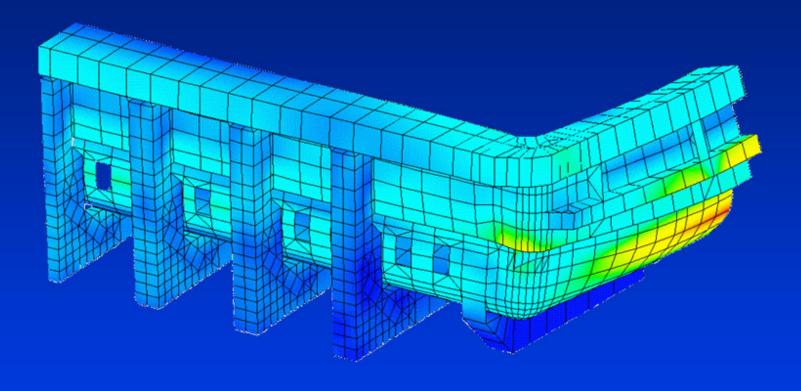


Corresponding sodium saturation for the same typical 45 cm thick cathode block, this time at $x = \frac{1}{4}, \frac{1}{2}$ and $\frac{3}{4}$ of the total block thickness as function of time as computed by that analytical equation.

It takes about 80 days for the sodium diffusion front to reach the bottom of the cathode block and about 1000 days for the cathode block to be fully saturated in sodium from top to bottom.



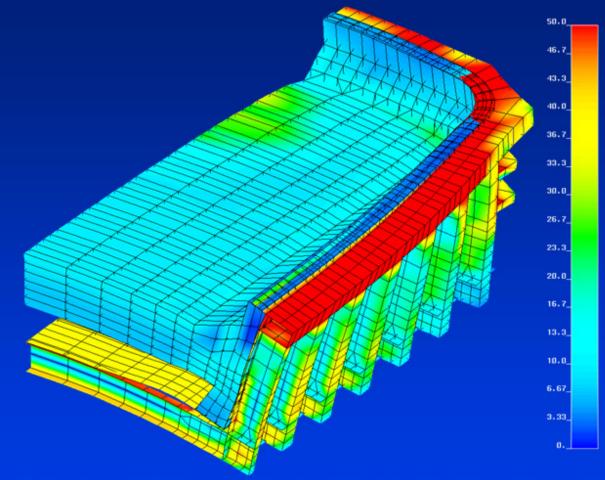
Historical Models Development Background



First "Empty shell" potshell model using plastic mode developed in 1988 and presented at the TMS in 1991



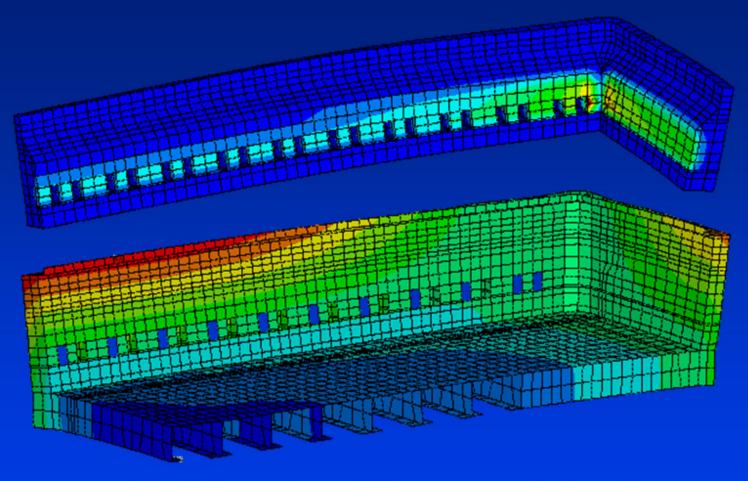
Historical Models Development Background



First "Half Empty shell" potshell model developed in 1989 and presented at a CRAY Supercomputing Symposium in 1990



Historical Models Development Background

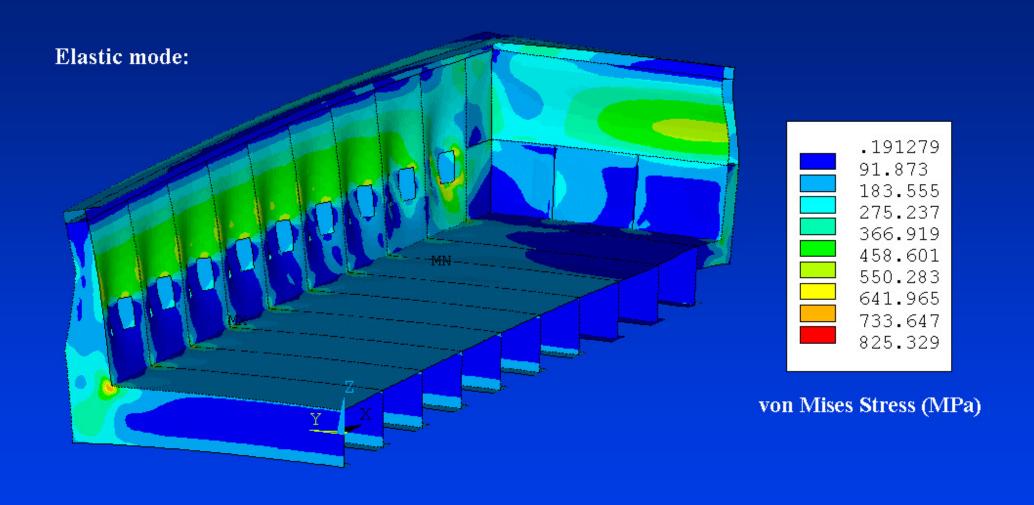


First "Almost Empty shell" quarter potshell model developed in 1992 and presented at CIM in 1993

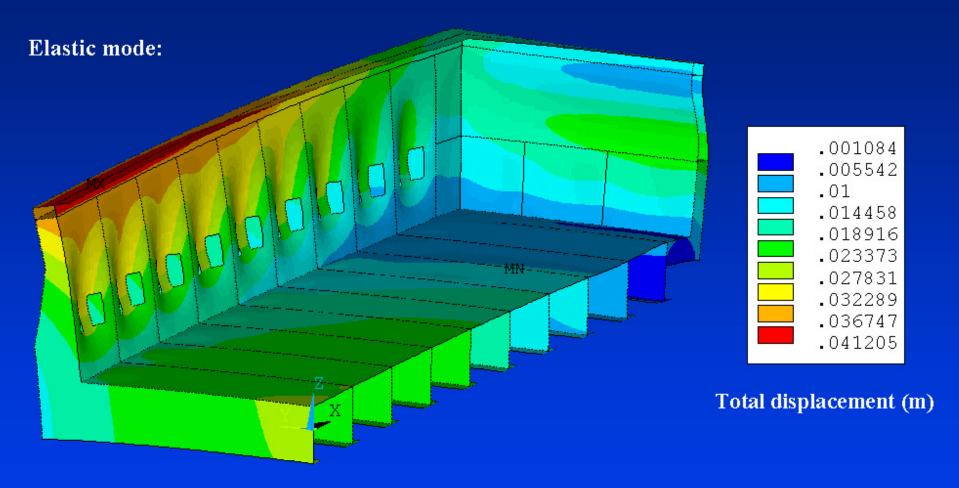


- Based on the usage of the quadrilateral Finite Strain shell element (SHELL181) in the commercial code ANSYS®.
- Temperature distribution obtained from the full cell quarter thermoelectric model applied as a body load to the entire potshell structure.
- Internal pressure (or forces) loading scheme based on the cathode block size and position relative to the potshell. That loading scheme is at best semi-empirical and is typically considered as a trade secret.
- Elastic mode: solving the mechanical problem only by considering only the elastic properties of the potshell steel.
- Plastic mode: solving the temperature dependent isotropic hardening von Mises plasticity behavior of the potshell steel structure using the MISO non-linear hardening option in ANSYS®.



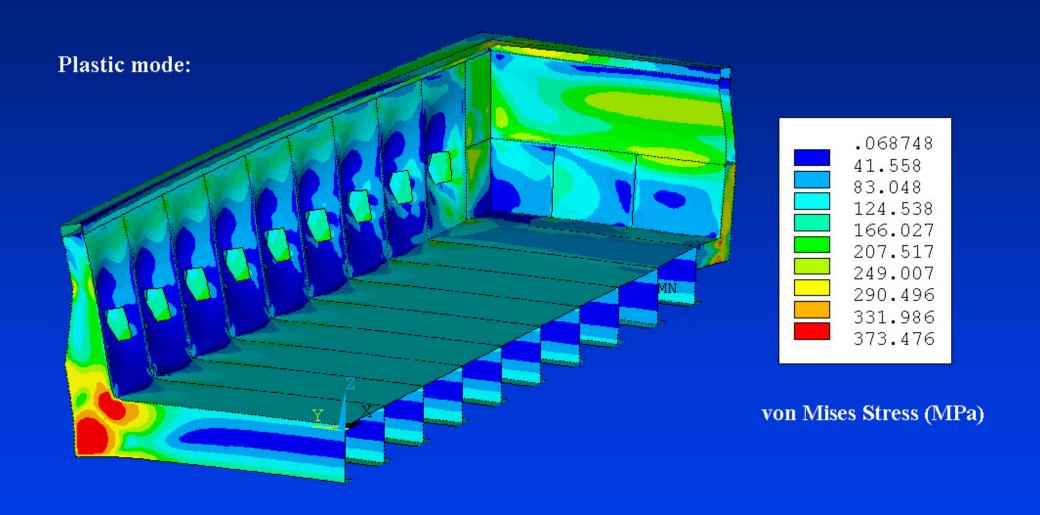




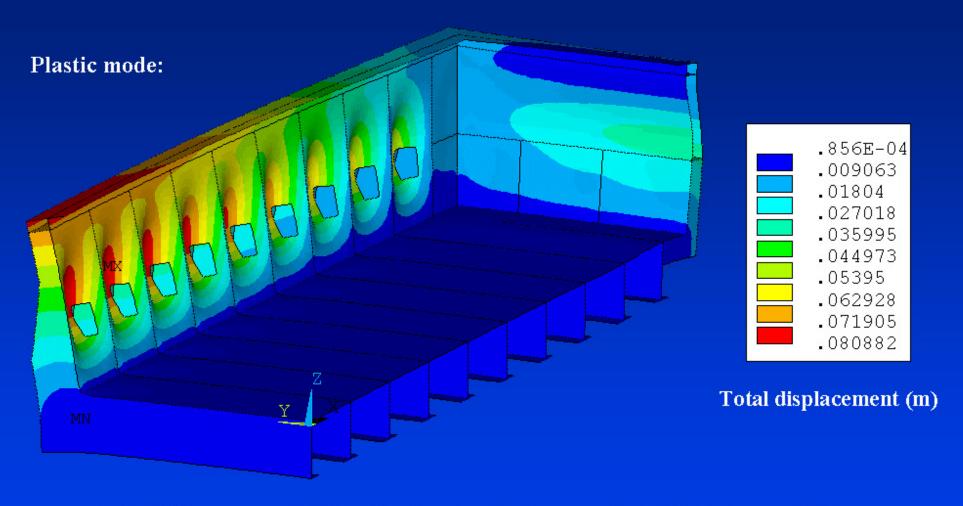


Took only 150 CPU seconds to solve on a 64 bits dual core Intel Centrino T9300 Dell Precision M6300 portable computer running ANSYS® 11.0 version.







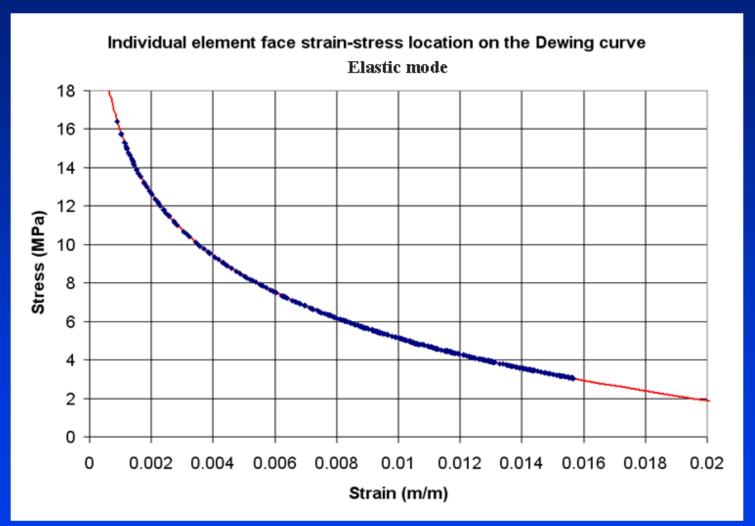


Took 11 times more CPU to solve than in elastic mode, but still this is only 1706 CPU seconds or 28 CPU minutes.



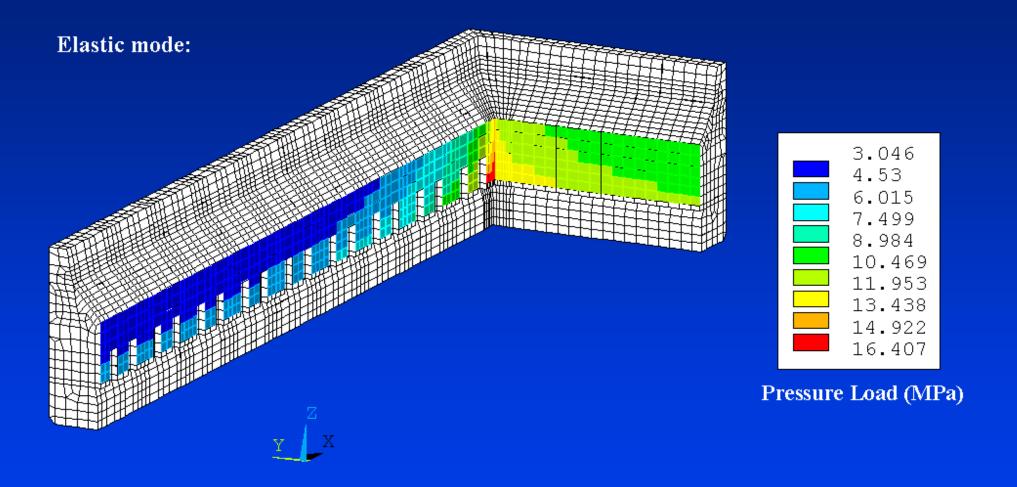
- Based on the usage of the quadrilateral Finite Strain shell element (SHELL181) in the commercial code ANSYS®.
- Temperature distribution obtained from the full cell quarter thermoelectric model applied as a body load to the entire potshell structure.
- Adding the lining geometry between the potshell walls and the cathode blocks all around the potshell and applying a pressure loading as boundary condition at the carbon block/side lining interface that is lying on the Dewing strain-stress relationship.
- Elastic mode: solving the mechanical problem only by considering only the elastic properties of the potshell steel.
- Plastic mode: solving the temperature dependent isotropic hardening von Mises plasticity behavior of the potshell steel structure using the MISO non-linear hardening option in ANSYS®.



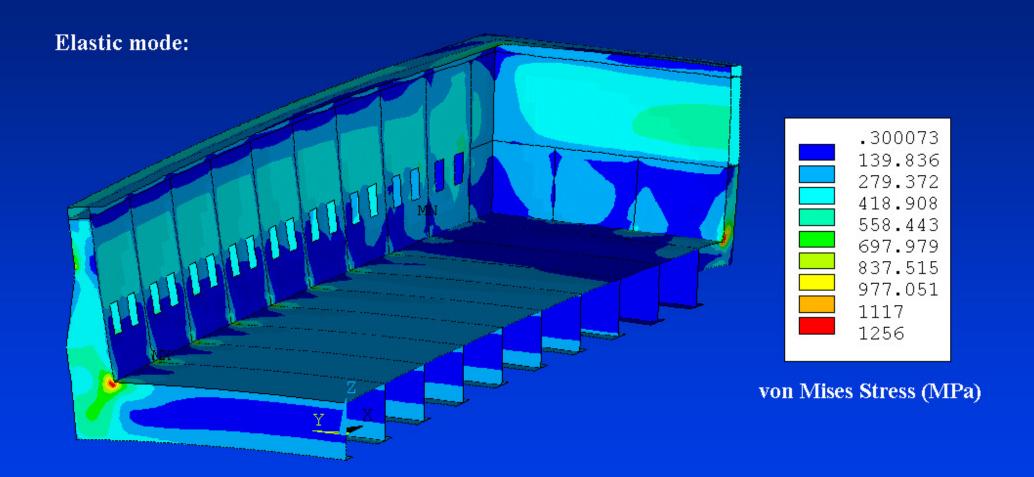


A numerical scheme external to the ANSYS solver must be setup, starting from an assumed initial internal load, the task of that external numerical scheme is to converge toward that cathode block equilibrium condition pressure loading for each element face of the carbon block/side lining interface.

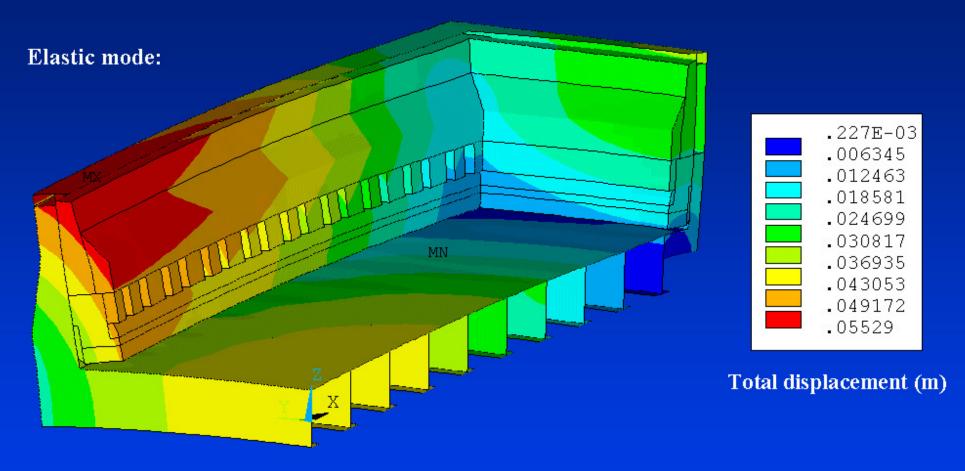






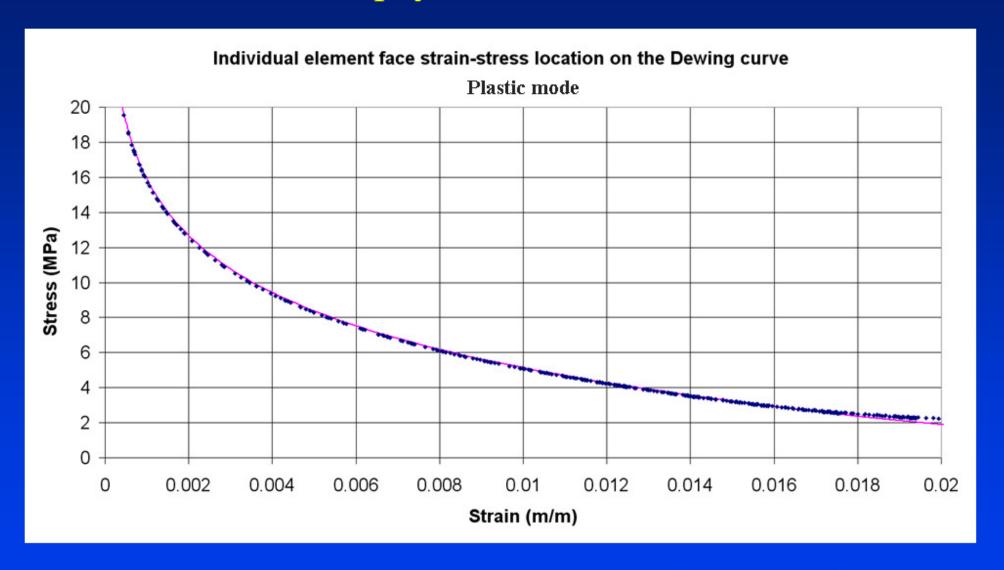




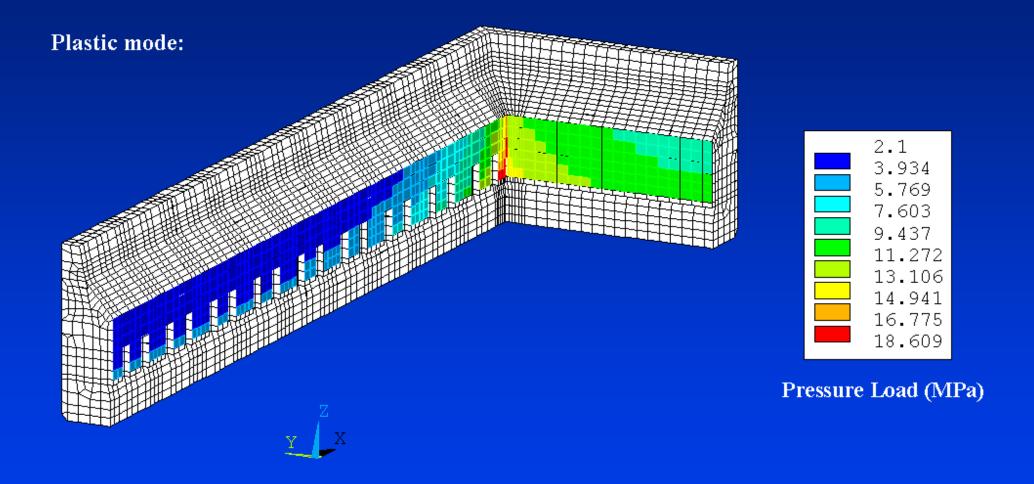


The external under-relaxation numerical scheme converges in about 20 iterations adding about the same factor to the required CPU time, in this case it took 3898 seconds CPU or 65 minutes CPU to solve.

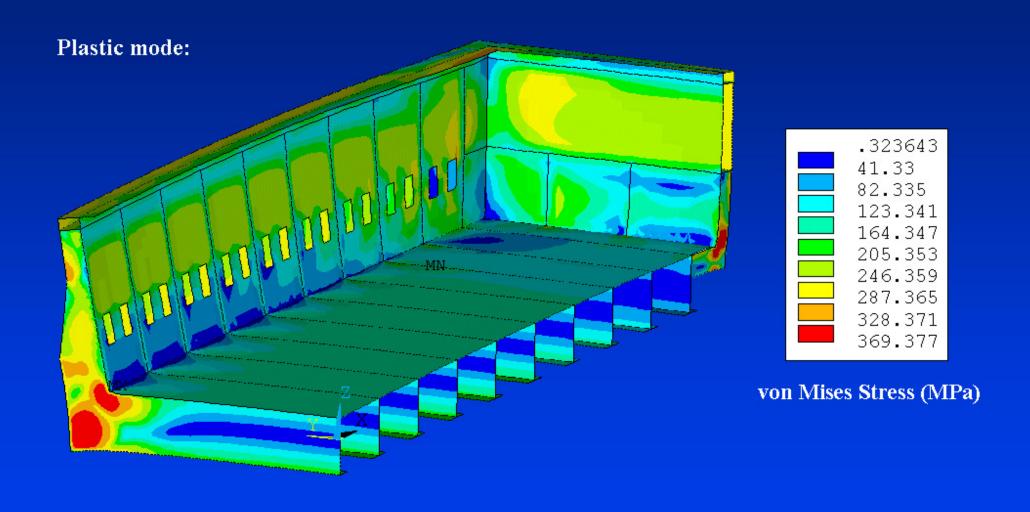




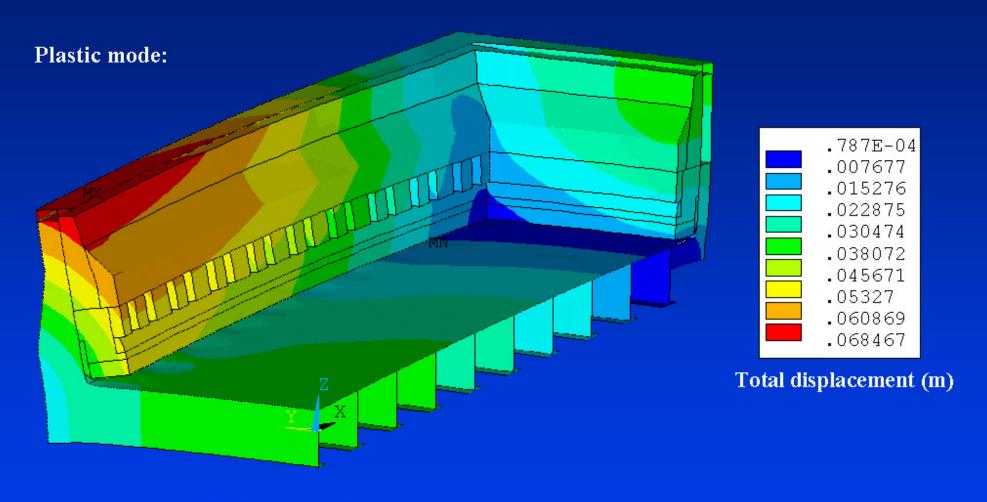








GENISIM

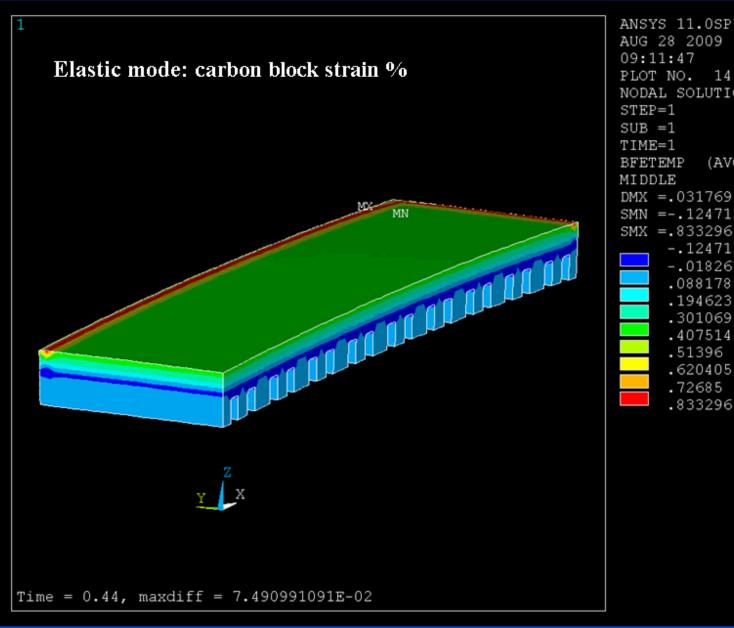


Took 27524 seconds CPU or 7.6 hours CPU to solve.



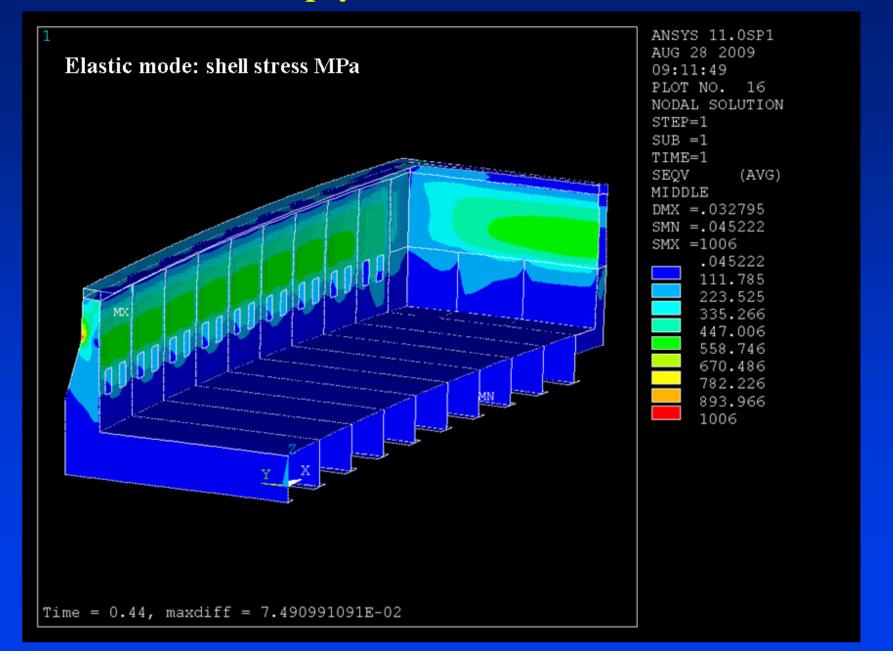
- Based on the usage of the quadrilateral Finite Strain shell element (SHELL181) in the commercial code ANSYS®.
- Temperature distribution obtained from the full cell quarter thermoelectric model applied as a body load to the entire potshell structure.
- The geometry the lining material located between the potshell walls and the cathode carbon of the cathode blocks themselves is also includes, the Dewing sodium expansion behavior of the cathode blocks is treated in ANSYS® as a "creep-like" behavior.
- Elastic mode: solving the mechanical problem only by considering only the elastic properties of the potshell steel.
- Plastic mode: solving the temperature dependent isotropic hardening von Mises plasticity behavior of the potshell steel structure using the MISO non-linear hardening option in ANSYS®.

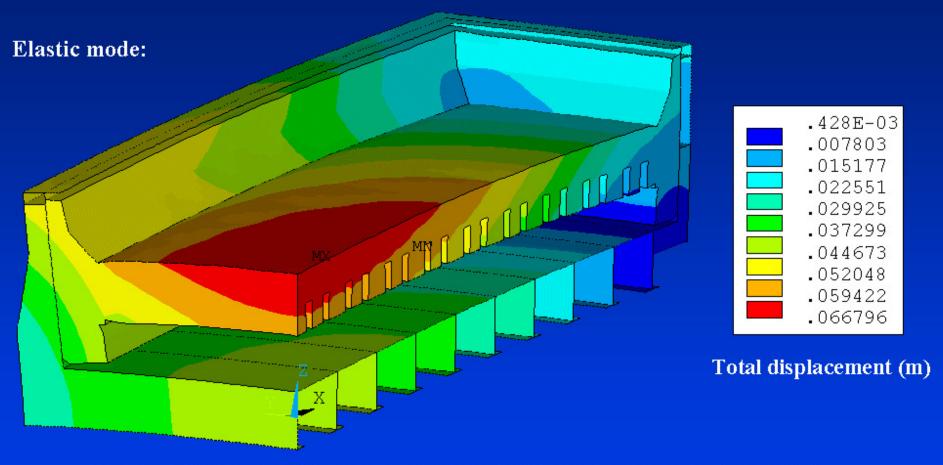




ANSYS 11.0SP1 AUG 28 2009 09:11:47 PLOT NO. 14 NODAL SOLUTION STEP=1 SUB = 1TIME=1 BFETEMP (AVG) MIDDLE DMX = .031769SMN =-.124713 SMX = .833296-.124713 -.018267 .088178 .194623 .301069 .407514 .51396 .620405 .72685

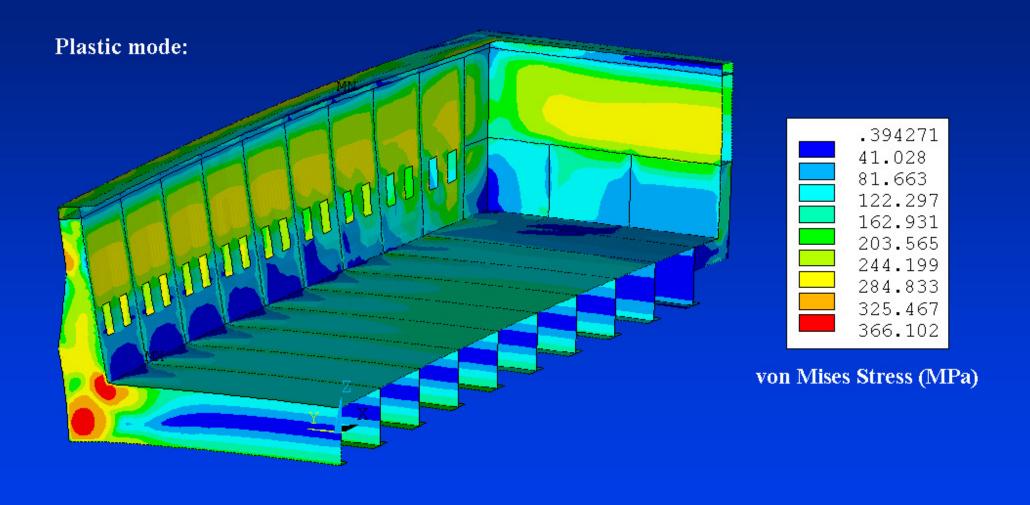
By definition, this means that it is required to solve that model in transient mode following the build-up of the sodium concentration in the cathode blocks from start-up to around 1000 days of operation where the cathode blocks get fully saturated in sodium.



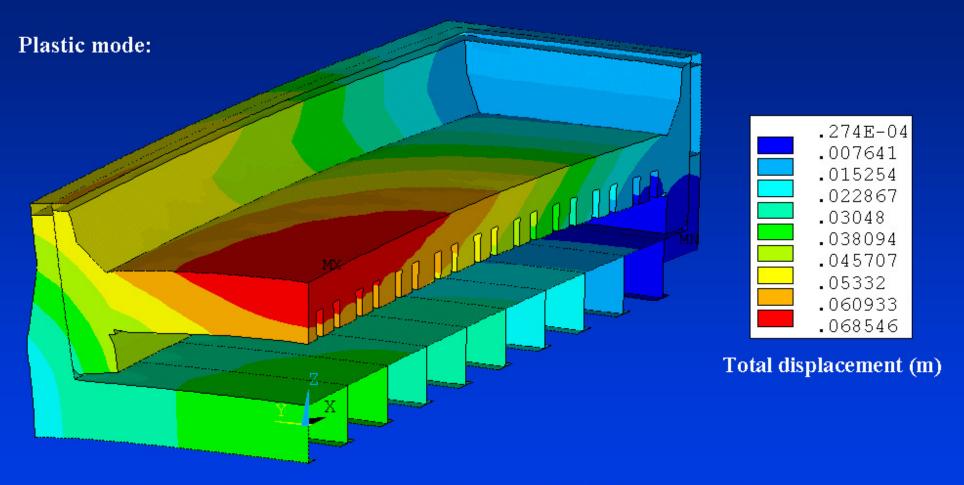


Solving the "half empty shell" potshell model in elastic properties mode took 25,335 CPU seconds or 7.0 CPU hours which is 6.5 times more than what was required to solve the "almost empty shell" potshell model.









Took 103842 CPU seconds or 1.2 CPU days which is 3.8 times more than what was required to solve the "almost empty shell" potshell model.



Conclusions

- Three types of ANSYS® based thermo-mechanical potshell models have been presented and described in details.
- The "empty shell" potshell model type has by far the fastest turn around time which is very convenient to be able to quickly investigate potshell design alternatives, but its ability to make accurate predictions outside of the narrow range of conditions where it has been validated is very questionable.
- The "almost empty shell" potshell model type is offering a much wider applicability range as it will automatically adjust the internal load as function of the grade of cathode blocks selected and of the rigidity of the potshell structure, it is offering an overnight turn-around time for a very little loss of accuracy.
- The "half empty shell" potshell model type is the most accurate model type of the three presented because it incorporates all the known physics of the cathode block sodium expansion phenomena affecting the potshell deformation in addition to the thermal expansion, unfortinately its turn-around time of 1.2 CPU days is already exceeding the overnight turn-around time that could be considered as the practical upper limit for an efficient design tool.



Conclusions

• Finally, until the day where the turn-around time of the "half empty shell" potshell model type would be reduced to mere minutes as it will become maybe 20 years from now (after all it was 80 days on a CRAY X-MP/24 supercomputer 20 years ago), it could be argued that all three types of potshell models have their places in the potshell designer toolbox as sometimes turn-around time matter more than accuracy and sometimes not.



Acknowledgements

The author wishes to thanks Dr Daniel Richard of Hatch and Mr Lalit Mishra of Dubal for their much needed assistance in using ANSYS® 11.0 thermo-mechanical features, especially in plastic properties mode.

