

FROM RESEARCH TO INDUSTRY



Isec

THREE-DIMENSIONAL MODEL OF A NEW THIN-PLATE LEAD-ACID BATTERY

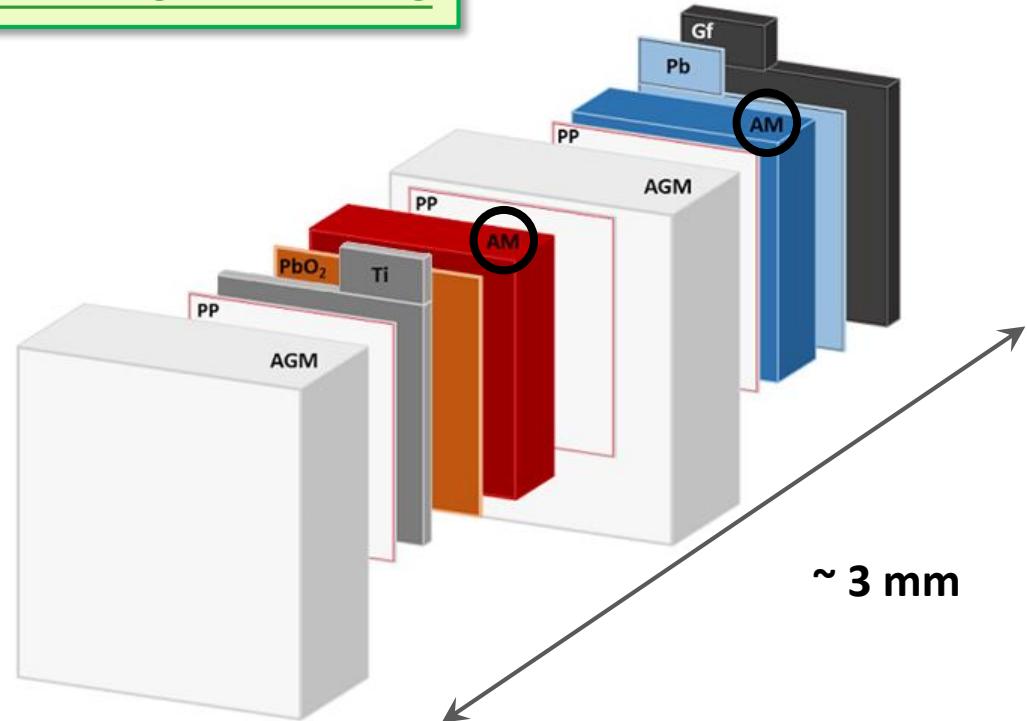
Jérémie Lannelongue

PhD Student
(with M. Cugnet & A. Kirchev)

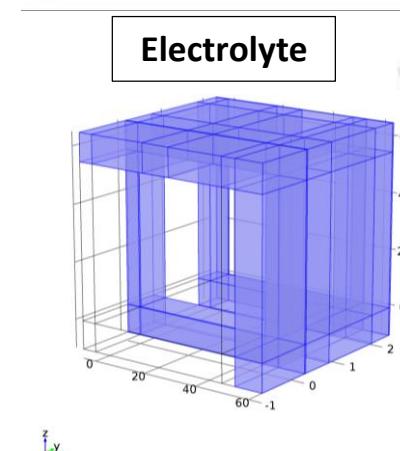
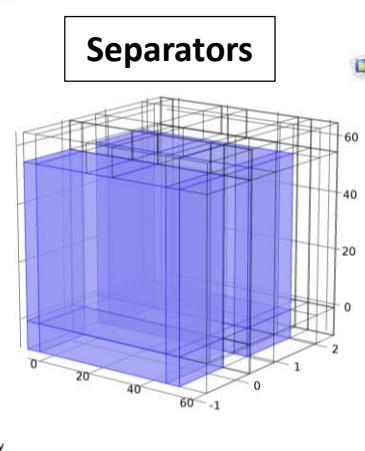
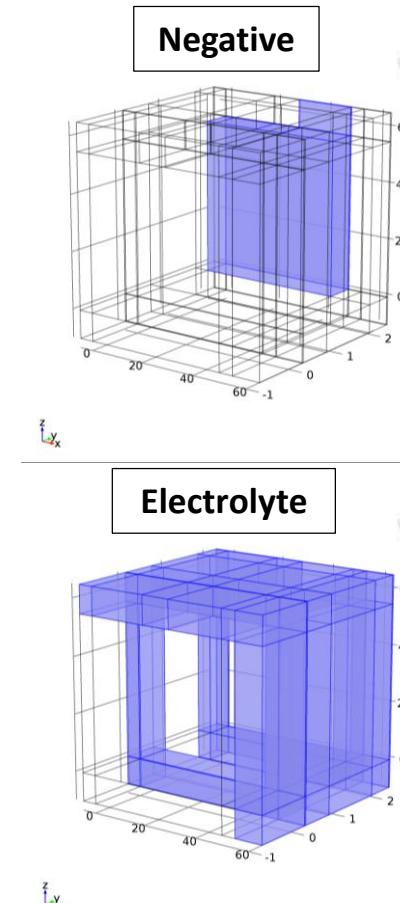
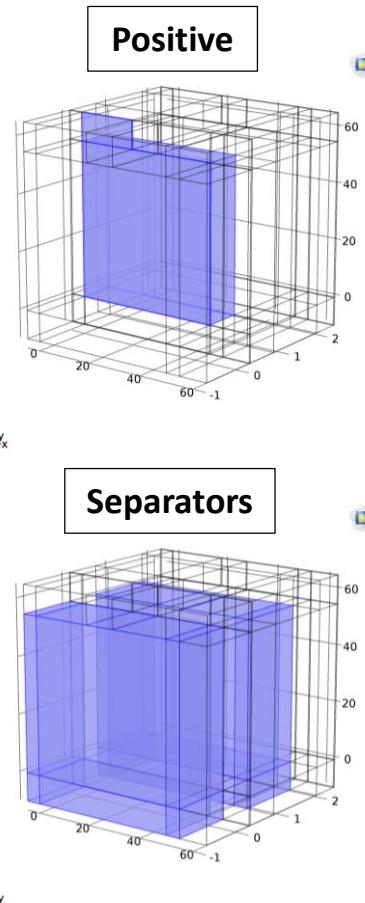
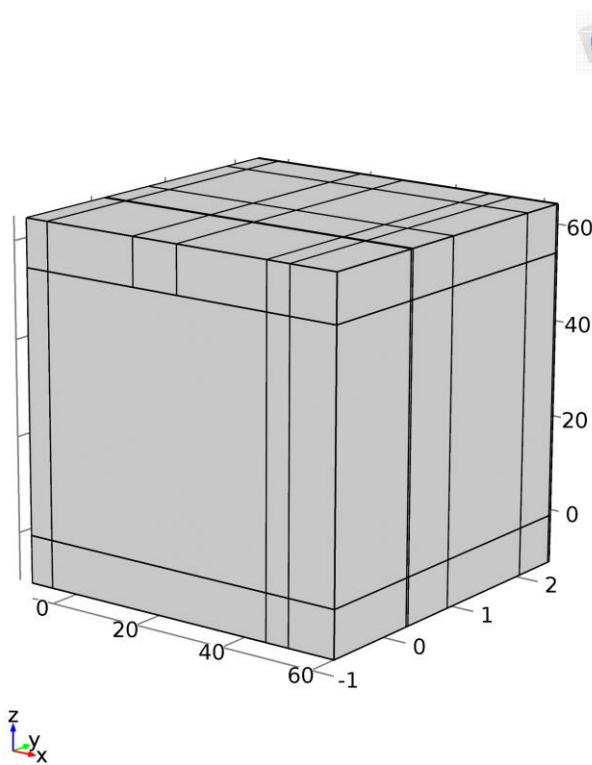
COMSOL
CONFERENCE
2015 GRENOBLE

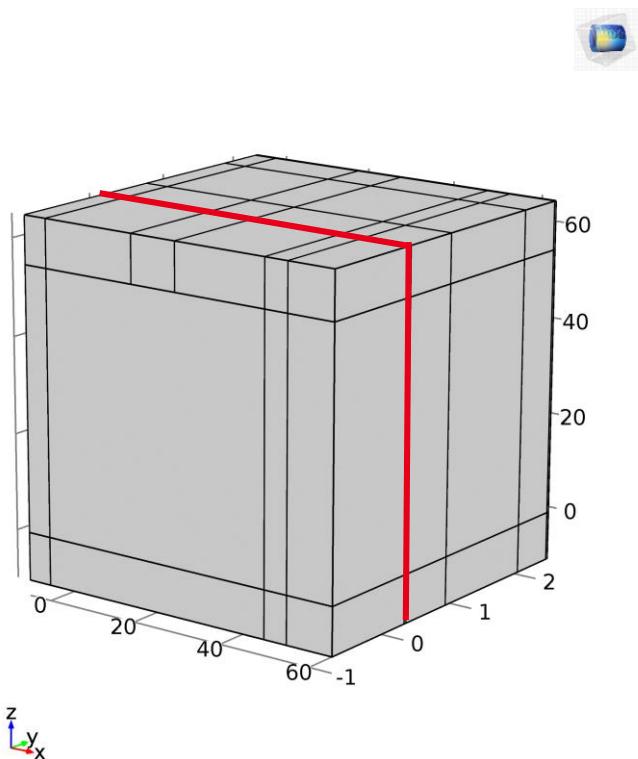


THIN-PLATE TECHNOLOGY FOR LEAD-ACID BATTERIES



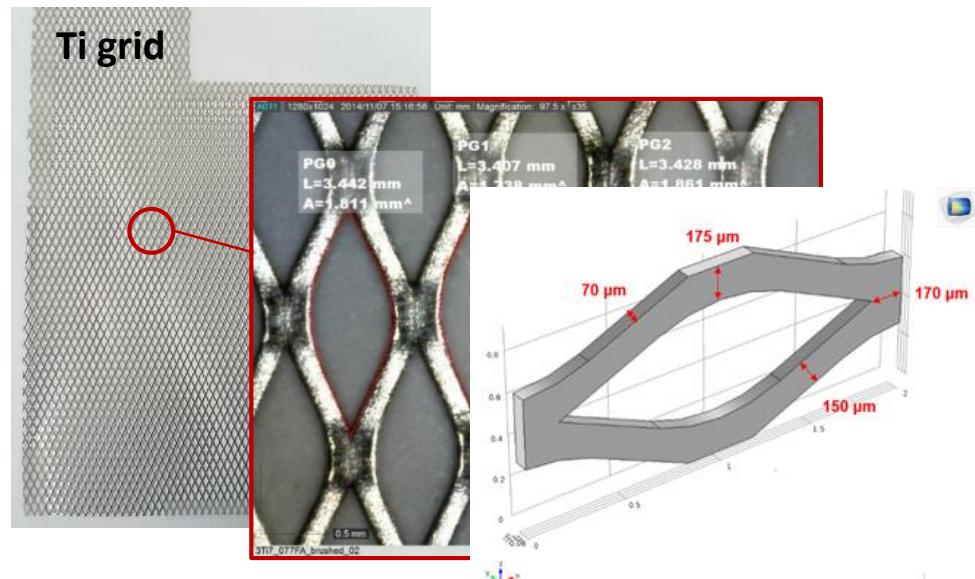
MODELISATION : GEOMETRY

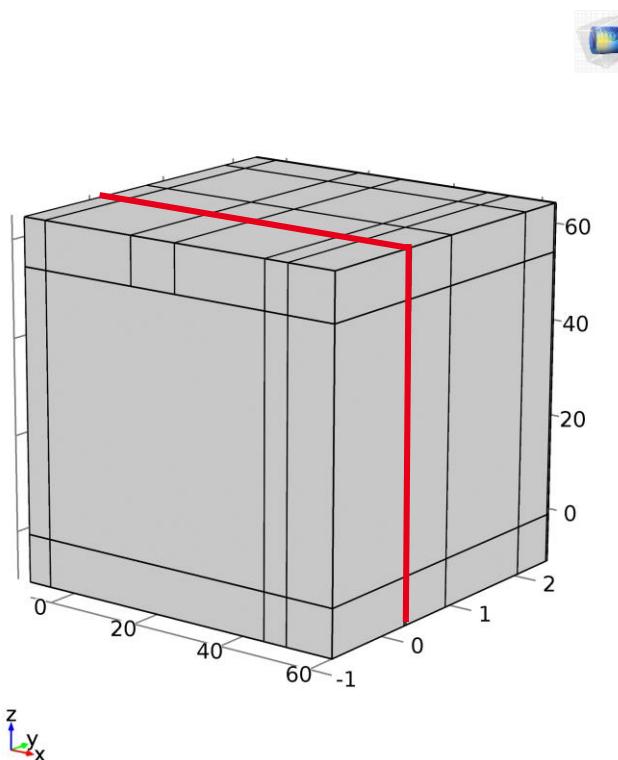




84 domains

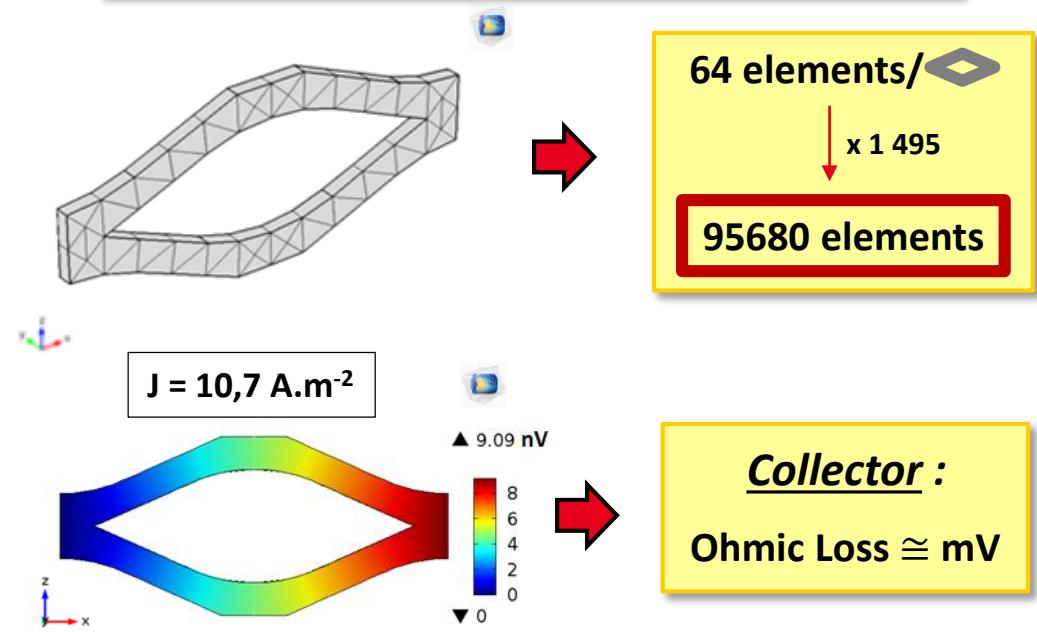
- Hypothesis : remove collectors (ex : positive)
 - High mesh cost
 - Low influence (at low rate)

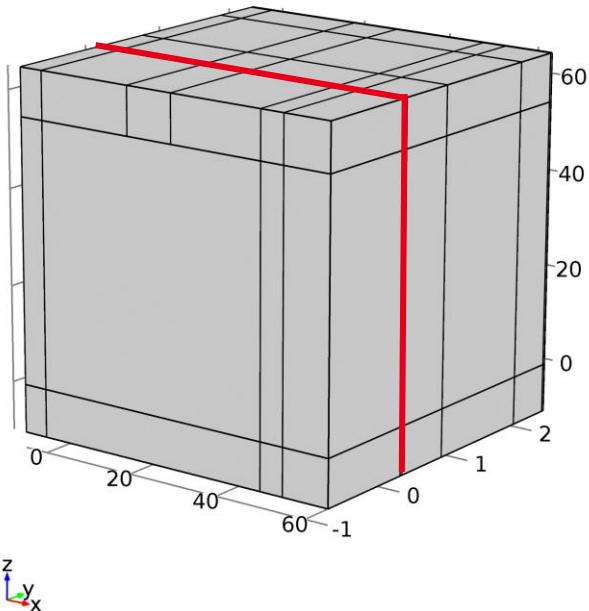




84 domains

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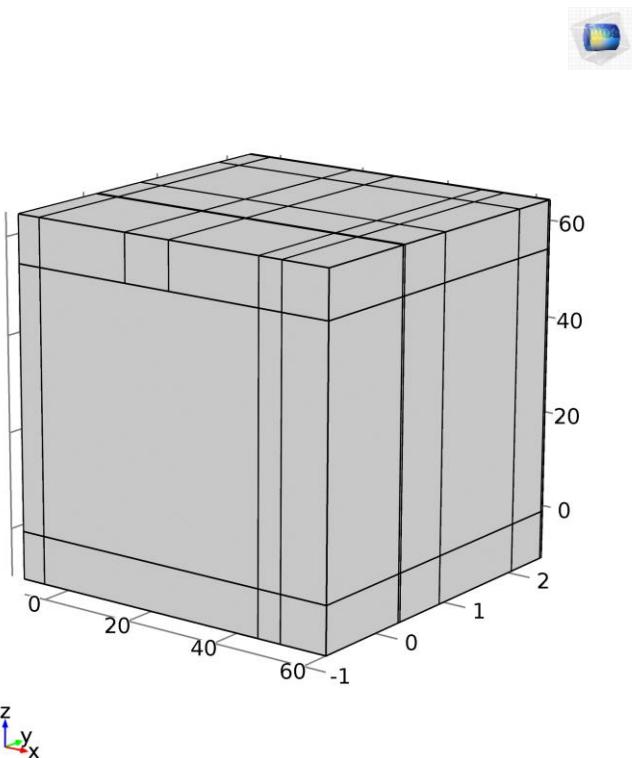




84 domains

- Hypothesis : remove collectors (ex : positive)
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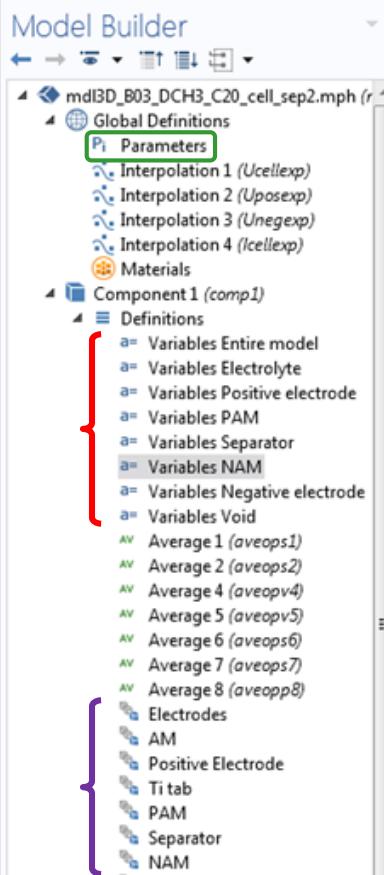


84 domains

- Hypothesis : remove collectors (ex : positive)
- Physics implementing
 - Ohm's law (solid (1)/liquid)
 - Porosity (2)
 - Material balance

$$(1) i_{soln} - \epsilon_a^{exm} \sigma \frac{\delta \varphi_{solid}}{\delta x} - I = 0$$

$$(2) \frac{\partial \epsilon}{\partial t} - \frac{\delta i_{soln}}{\delta x} \frac{1}{nF} (\bar{V}_a s_{a,e} + \bar{V}_b s_{b,e}) + \frac{R_{b,e}}{s_{b,s}} (\bar{V}_a s_{a,s} + \bar{V}_b s_{b,s}) = 0$$



Variables

Label: **Variables NAM**

Geometric Entity Selection

Geometric entity level: Domain

Selection: **NAM**

Active

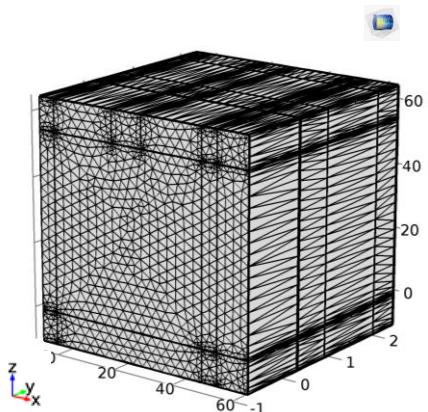
Name	Expression	Unit	Description
$S_{electron}$	$-A_a_{neg} * i0_{neg} * (\exp((1-a_{neg}) * n * F_{const} * \eta) - \exp(-a_{neg} * n * F_{const} * \eta))$	A/m^3	Electron current source
S_{ion}	$-S_{electron}$	A/m^3	Ion current source
η	$\phi1 - \phi2 - U_{neg}$	V	Overpotential
U_{neg}	$-0.294606[V] - 0.073595[V] * \log10(m[kg/mol]) - 0.030432[V] * \log10(m[kg/mol])^{2-0.5}$	V	Negative equilibrium voltage
S_{por}	$-(PMV_{PbSO4} - PMV_{Pb}) / (2 * F_{const}) * S_{ion}$	1/s	Source term of the porosity
e_0	e_{neg_0}		Initial porosity
e_{neg}	e		Negative electrode porosity
s_{NAM}	$s_{Pb} * e^{0.5}$	S/m	Effective NAM electronic surface
A_a_{neg}	$A_a_{neg} * 0 * (1 - \exp(-0.5 * (e - 0.448) / (e_0 - 0.448))) / (1 - \exp(-0.5))$	$1/m$	Effective active area
S_{tds}	$(1 - 2 * tp0) / (2 * F_{const}) * S_{ion} - c * S_{por}$	$mol/(m^3)$	Source term of material b...
$i0_{neg}$	$i0_{neg} * (c / c_{ref})^{1/2}$	A/m^2	Exchange current density
n_{Pb}	$\rho_{Pb} * V_{nam} * e_{Pb} / MW_{Pb}$	mol	Moles of Pb

84 domains

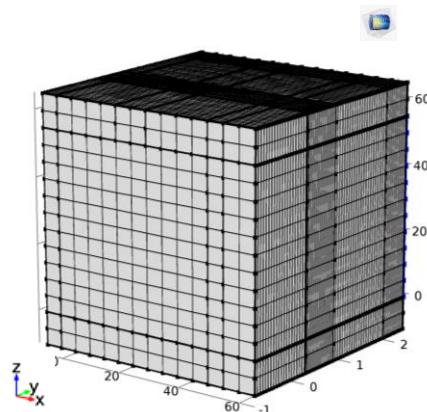
- Hypothesis : remove collectors
- Physics implementing

↳ **Parameters, Variables & Selections**

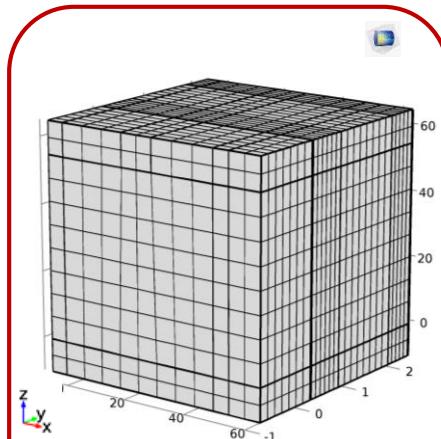
MODELISATION : MESH



Physics-controlled mesh
Coarser size
57 107 elements



User-controlled mesh
Mapped 1 & Distribution 1
13 860 elements
1h10min

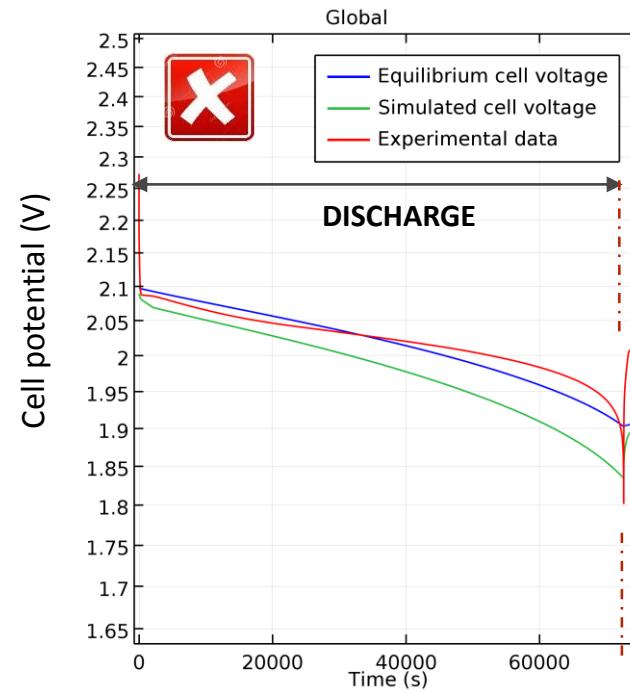
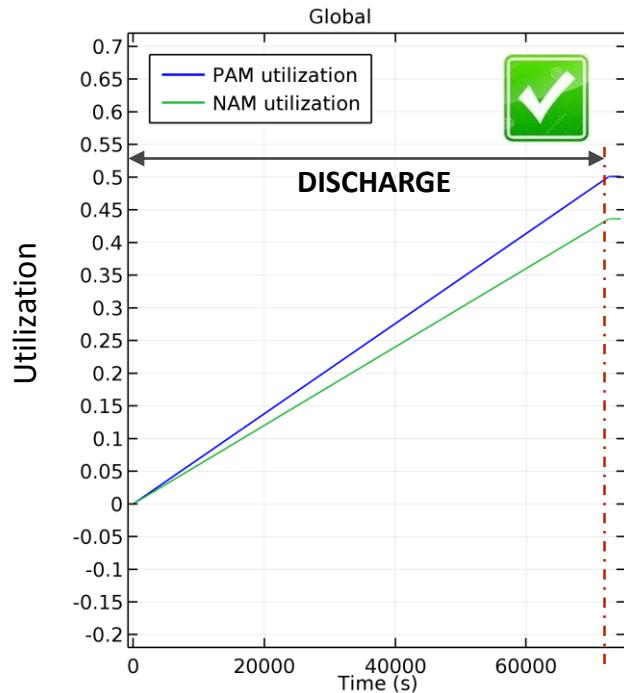


User-controlled mesh
Mapped 2 & Distribution 2
4 032 elements
15min

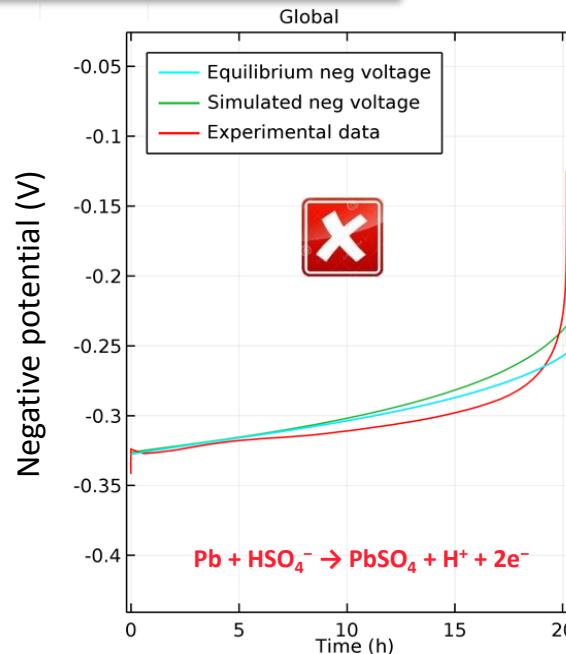
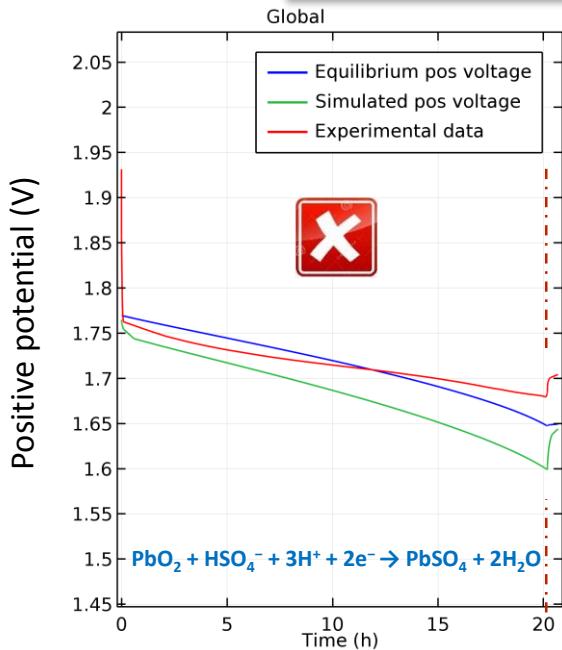
EFFICIENT MESH = TIME GAIN

- User-controlled mesh
- Mapped & Swept
- Distribution

Discharge + rest : first results



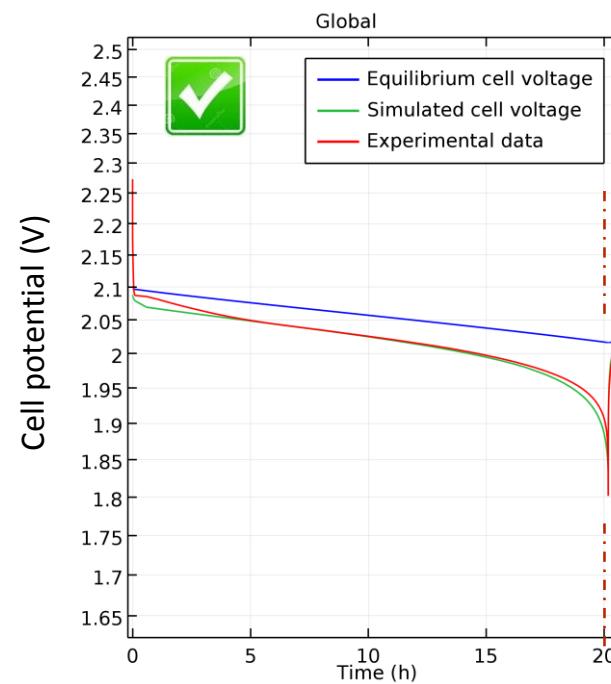
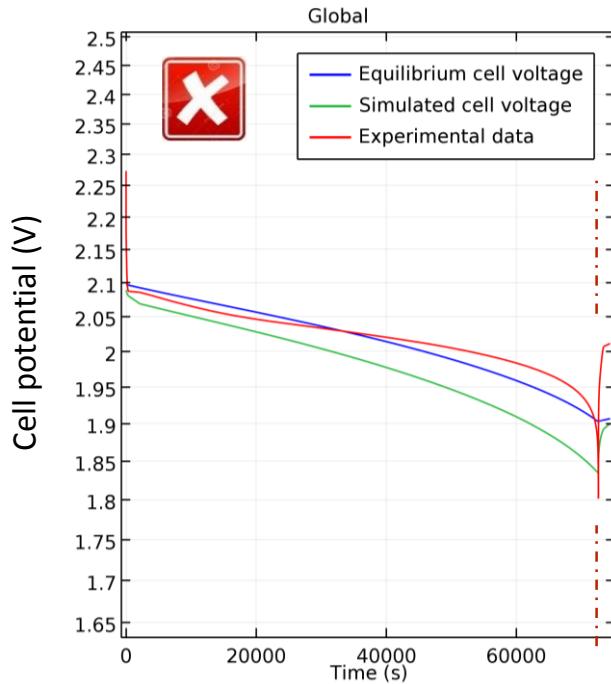
Discharge + rest : first results



- Lack of acid (+ & -)
- Inappropriate surface coverage (-)

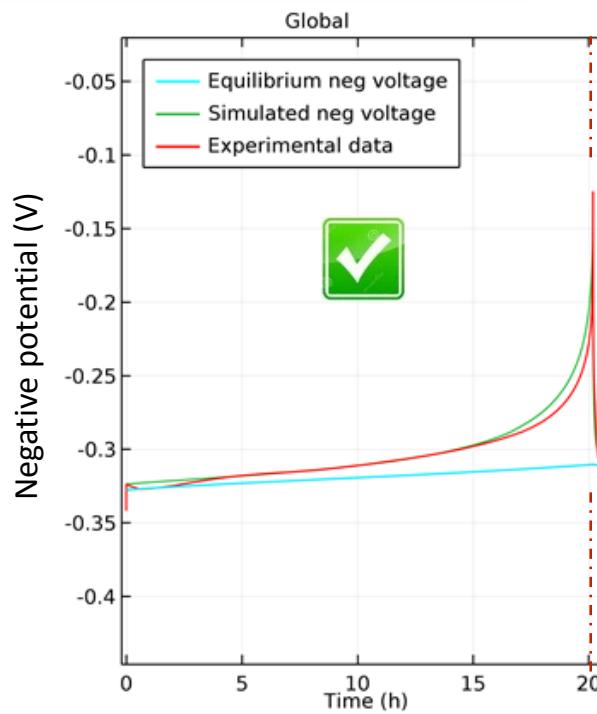
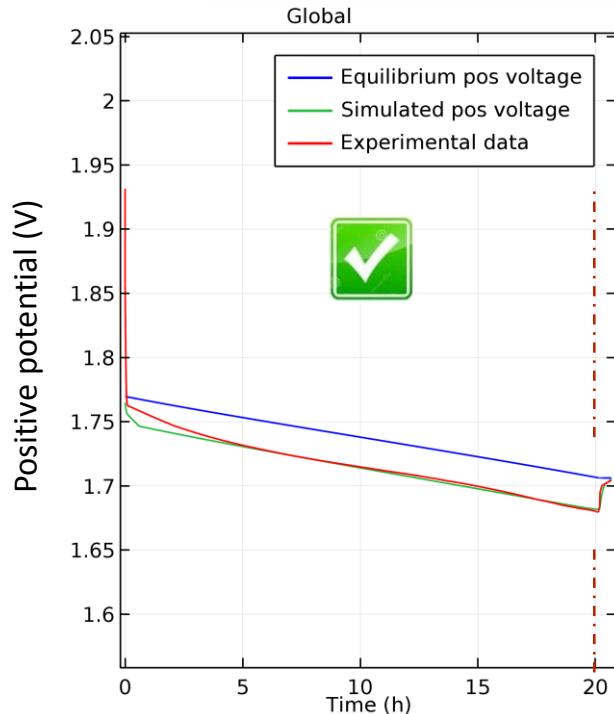
Adjustments

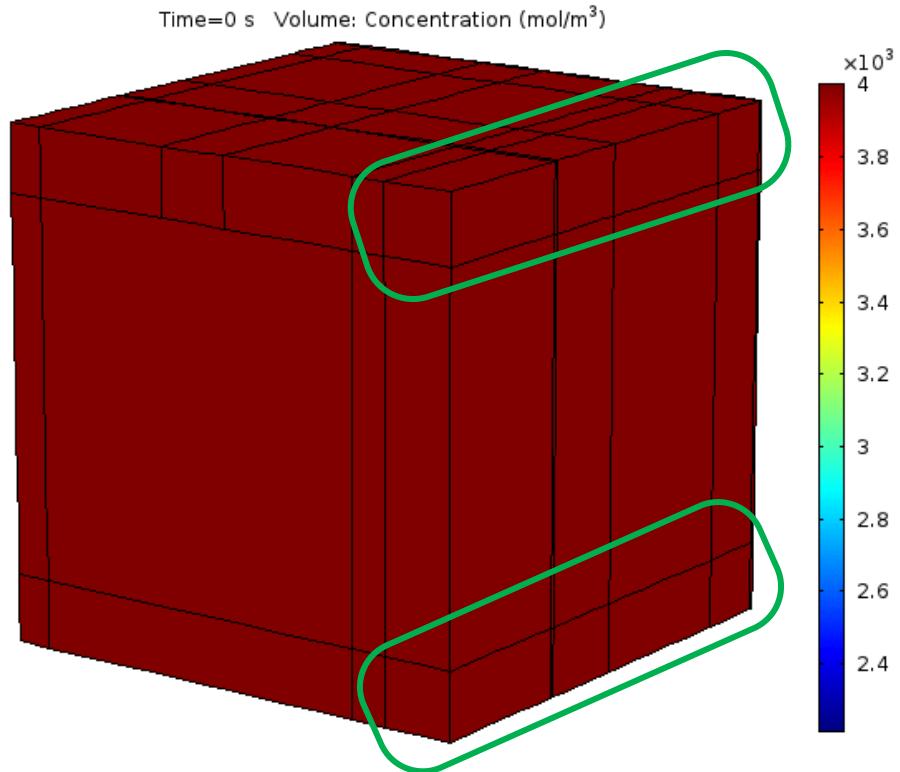
- PAM & NAM thicknesses / Porosity
- Electrolyte volume
- Coverage factor



Adjustments

- PAM & NAM thicknesses / Porosity
- Electrolyte volume
- Coverage factor





Acid concentration

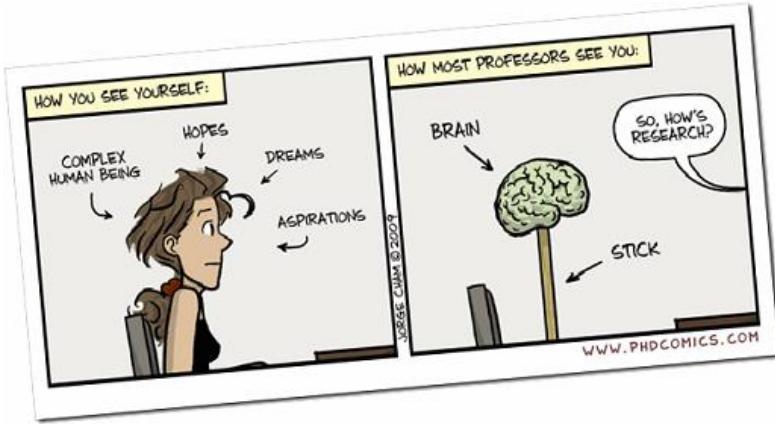
- Heterogeneous consumption
- Critical volume

Conclusions

- Importance of Hypothesis / Mesh / Parameters
- Determine precisely the key parameters :
 - AM thicknesses
 - Porosity
 - Exchange current etc
- Better understanding of the reaction (Acid consumption)

Perspectives

- Add Physics (heat transfer, nucleation and growth)
- Modelisation of Discharge at high rate
- Move to 2D ?



Thank you for your attention

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