

MATHEMATICAL MODELLING FOR ALL-SOLID-STATE BATTERY: (SE|SSE)-INTERFACE

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Mathematical modelling for the next-generation All-solid-state batteries: Nucleation (SE|SSE)^(*)-interface

Rechargeable Lithium-ion battery (LIB) is at the heart of every electric vehicle (EV), portable electronic device, and energy storage system [5]. Nowadays, LIBs enable human life more efficient and help to solve global environment issues thanks to EVs' zero emission. However, conventional LIB (c-LIB) is sensible to temperature and pressure, hence, flammable and explosive, which is undesirable. This bottleneck is mainly due to **liquid-based electrolyte** found in c-LIBs.

All-solid-state battery (ASSB) is one of promising candidates to overcome bottlenecks of c-LIBs. Thanks to **solid-state electrolyte** (SSE), ASSB is highly stable towards temperature and pressure. Nevertheless, Li-metal dendrite triggered at (SE|SSE)-Interface is the main drawback of ASSB since these dendritic threads extrapolate into SSE grain boundary network, causing crevice, degradation of ionic conductivity, and the probability of short-circuit, which is unfavorable.

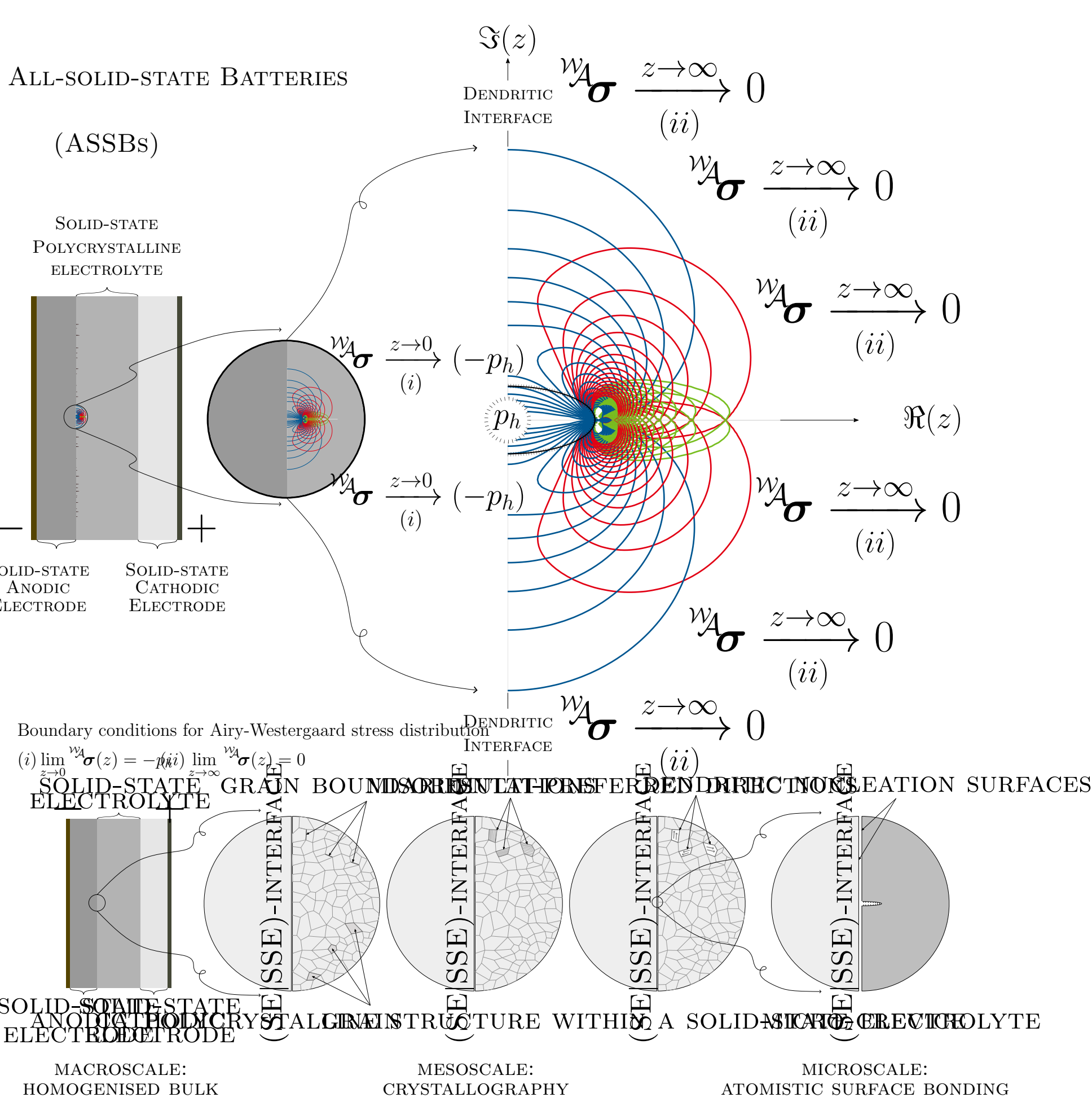
Next-generation All-solid-state battery (ng-ASSB) with a consideration of **nucleation criterion** defined by

$$\partial_t \mathbf{u} + \nabla \cdot \left(\mathbb{C}^{f_{\text{allocation}}}(\lambda, \mu, \mathbf{d}_{G,i,i=1,\dots,N}^B, \mathbf{d}^E; \mathbf{x}) : \nabla \mathbf{u}^{(s)} \right) + \rho \mathbf{b} = -\rho \nabla V_e, \quad (1)$$

$$\text{s.t. } a_{\text{Griffith}} := a^* = \arg \min_{a \in \mathbb{R}} \iint_{\Omega} f(a, \mathbf{u}, \theta; \lambda, \mu, \mathbf{d}^{(*)} \otimes \mathbf{d}^{(*)}) d\Omega - \iint_{\Gamma} f(a; \gamma) d\Gamma \Big|_{\bar{\mathbf{u}}} \quad (2)$$

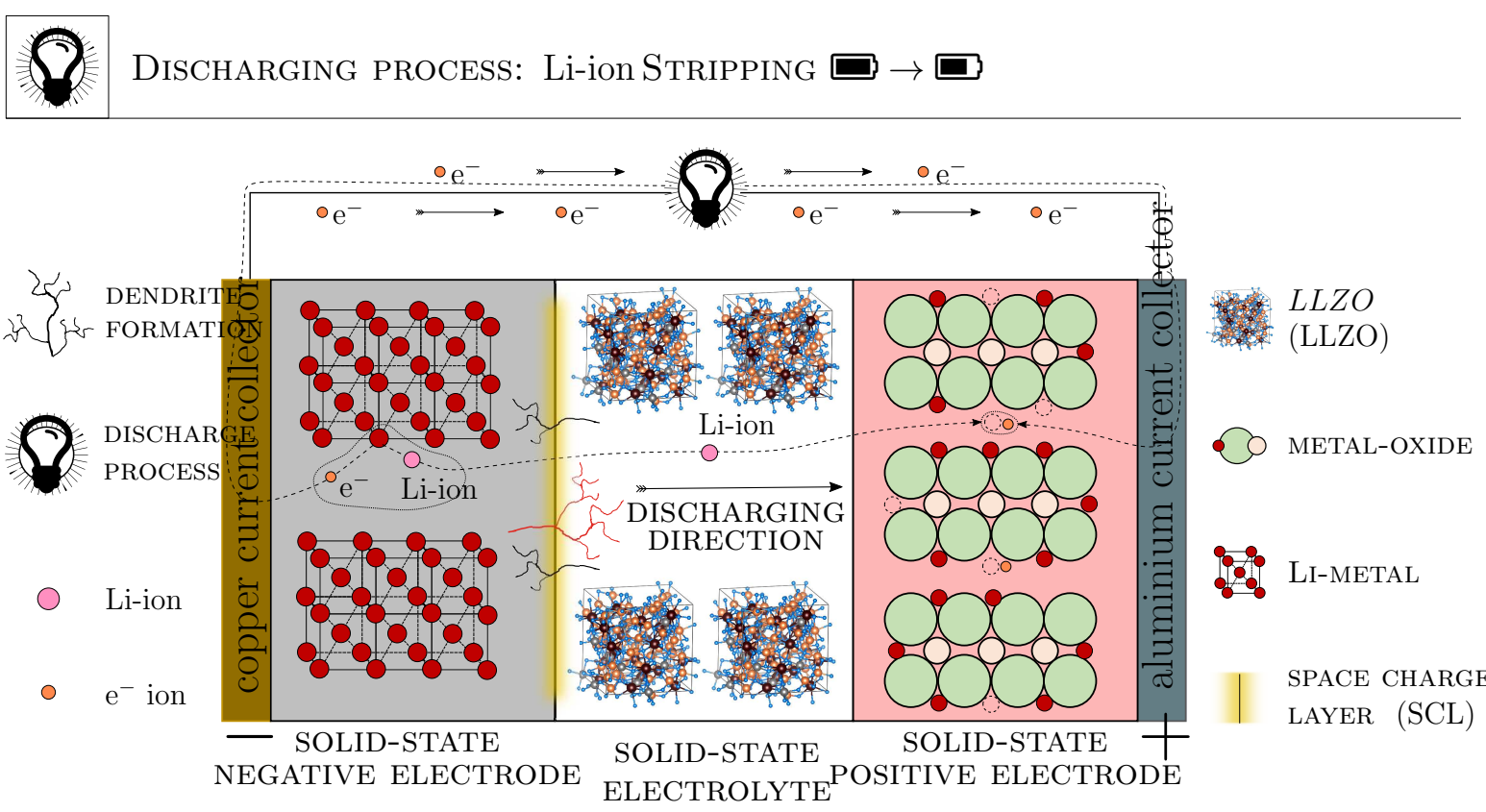
where $V_e : \mathbb{R}^3 \rightarrow \mathbb{R}$ is the electric potential applied globally on ASSB. Due to nature setting of ASSB taking the form (SE|SSE|SE) the electric potential becomes uniform. Additionally, \mathbf{u} is the displacement field, θ temperature field, a crevice length, λ, μ Lamé constants, $\mathbf{d}^{(*)} \otimes \mathbf{d}^{(*)}$ embedded misorientation structural tensor, and γ cracking-surface energy density, can help to improve ASSB performance.

Aim: The study is with the purpose of getting a better insight into dendrite formation in ASSB



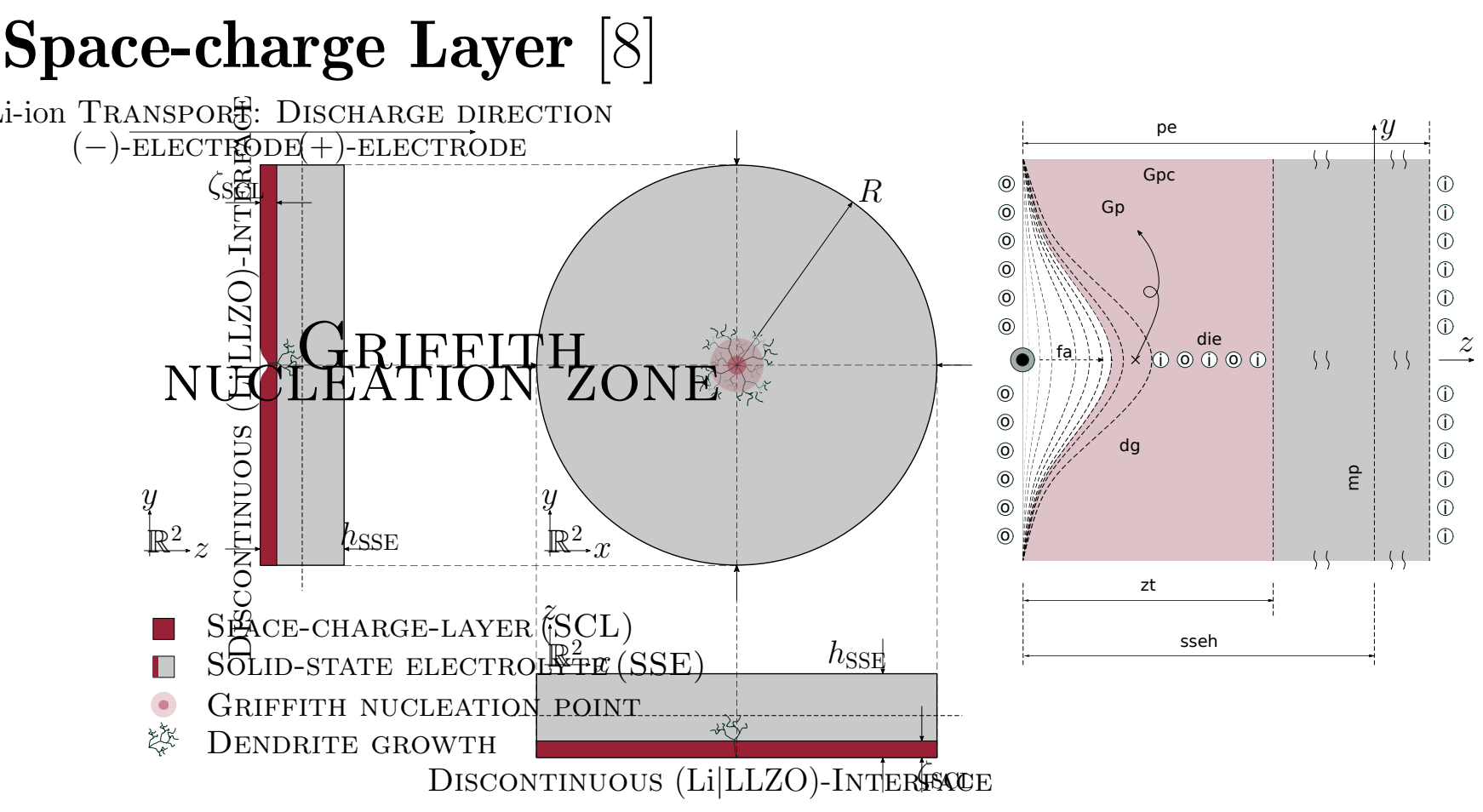
Next-generation All-solid-state battery

Griffith nucleation governs the (SE|SSE)-Interface [10]

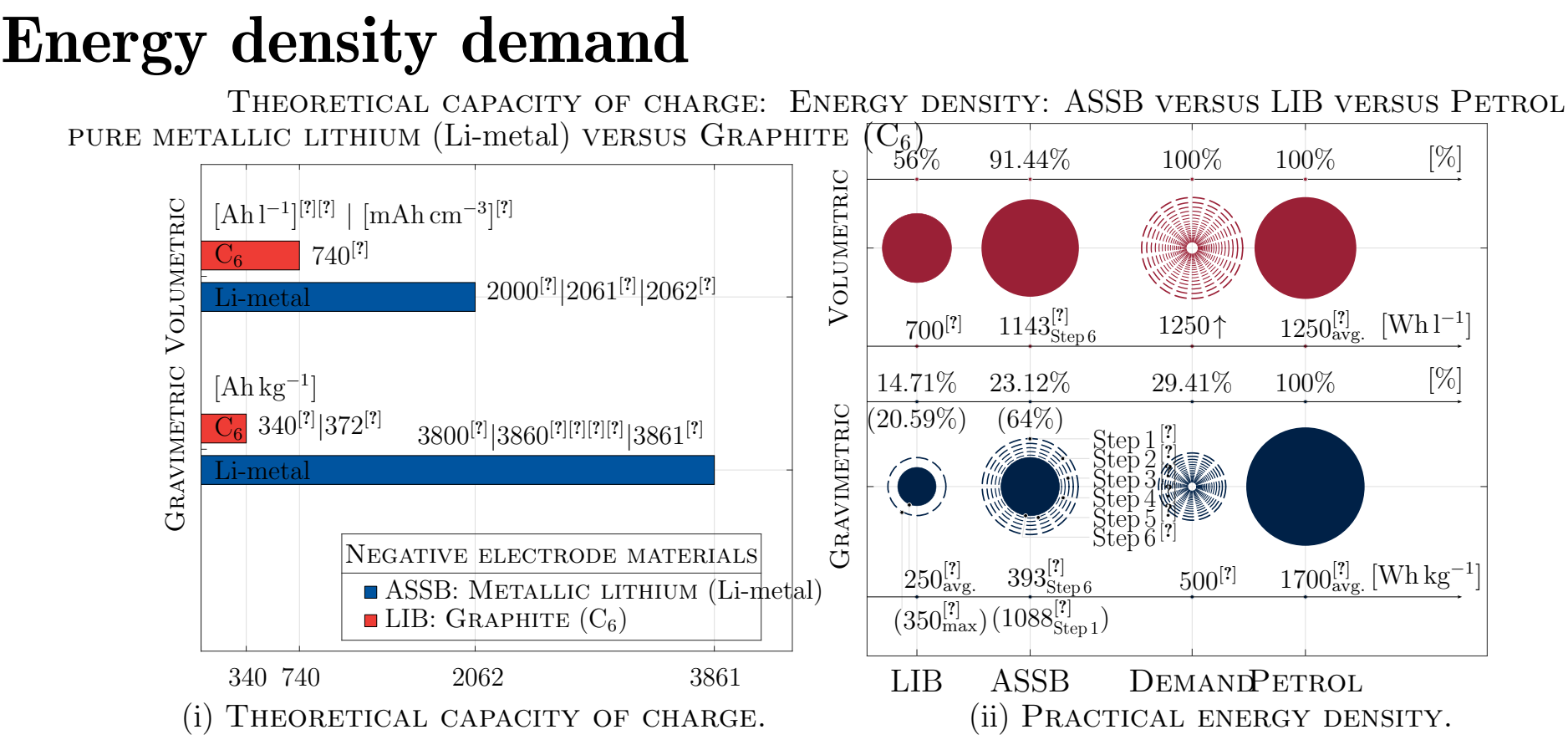


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Space-charge Layer



Energy density demand landscape



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Artificial Neural Networks

Application: Steel property prediction [9]

Semiconductor

Application: Start/Stop Starter [7]

Nucleation interface: Taking place at the critically dendritic (SE|SSE)-interface

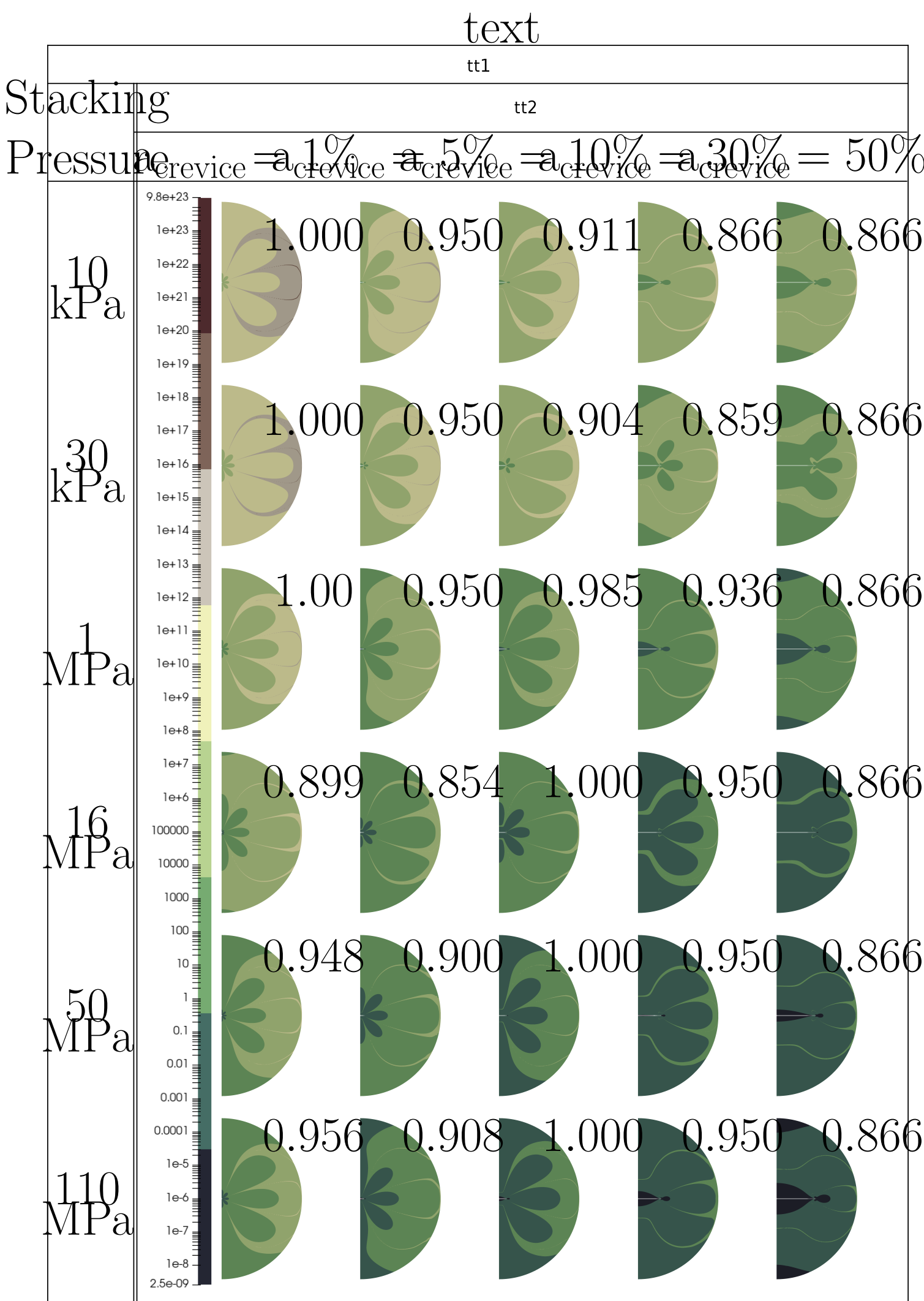
Coupled fields are Displacement field \mathbf{u} and temperature field θ ;

$$\mathbf{u} : \begin{cases} \Omega \times \mathbb{R}_+ \rightarrow \mathbb{R}^3, \\ (\mathbf{x}, t) \mapsto \mathbf{u}(\mathbf{x}, t), \end{cases} \quad \theta : \begin{cases} \Omega \times \mathbb{R}_+ \rightarrow \mathbb{R}, \\ (\mathbf{x}, t) \mapsto \theta(\mathbf{x}, t), \end{cases}$$

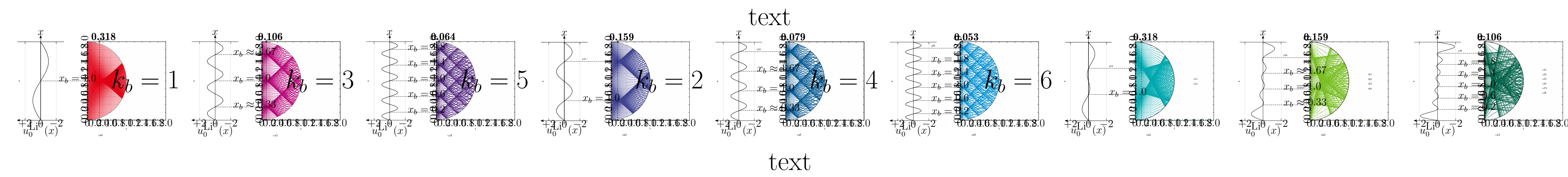
Governing conservation equations used to describe balance of mass, conservation of linear momentum, conservation of angular momentum, and conservation of energy with $\rho(\mathbf{x}, t)$ is mass density per unit volume (puv); $\mathbf{b}(\mathbf{x}, t)$ body force puv; $\mathbf{v}(\mathbf{x}, t)$ velocity; $e(\mathbf{x}, t)$ internal energy puv; $\mathbf{q}(\mathbf{x}, t)$ heat flux; $r(\mathbf{x}, t)$ heat source puv; $\boldsymbol{\sigma}$ Cauchy stress and $\boldsymbol{\varepsilon}$ infinitesimal strain.

Strain energy is based on the deformation of SSE due to dendrite formation at (SE|SSE)-interface

Surface energy is analysed based on the open crevice cracking at (SE|SSE)-interface affected by prescribed pressure



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Acknowledgments

T.Vo would like to extend heartfelt gratitude to the JARA-CSD research project, spearheaded by Dr.S.Braun and Dr.C.Hüter, for their invaluable support. Additionally, T.Vo would like to say thank to Prof.Dr.R.Spatschek for the host at IEK-2 in FZ Jülich. Furthermore, T.Vo's teaching assistant experience in the courses of Mathematische Grundlagen I, II, III, IV, and NumPDE at Computational Engineering Science (CES) at RWTH Aachen University was made possible thanks to the guidance and mentorship of Prof. Dr. M. Torrilhon, Prof. Dr. M. Schlottke-Lakemper, Dr. R. Speck, Prof. Dr. B. Berkels, Dr. S. Braun, Prof. Dr. J. Kowalski, Dr. A. Jha, Prof. Dr. B. Stamm, Dr. A. Litvinenko, and Dr. M. Kirchhart, spanning from the Summer Semester of 2019 to the Summer Semester of 2023.