



PyBaMM

Python Battery Mathematical Modelling



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PyBaMM's mission is to *accelerate battery modelling research* by providing an *open-source framework* for multi-institutional, interdisciplinary *collaboration*.



- Provide fast, reliable battery simulations
- Grow battery modelling community
- Facilitate development of new battery models
- Improve reproducibility of research
- Increase impact and industry engagement



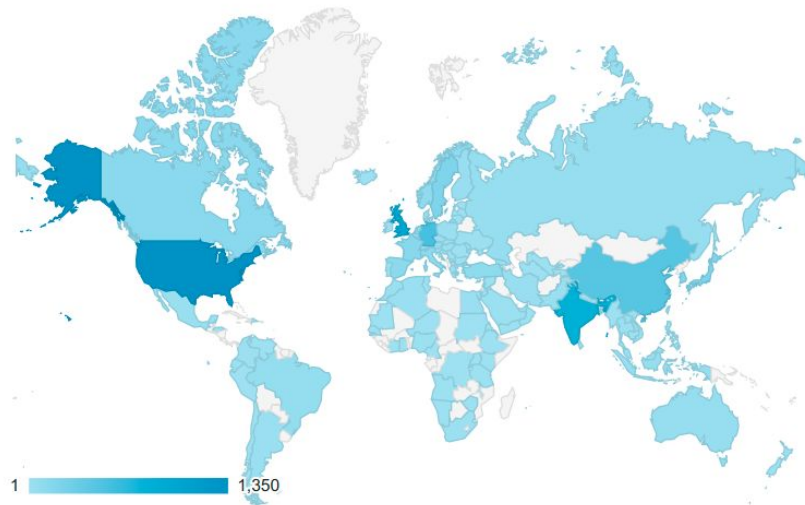
Try online



Install locally



Develop



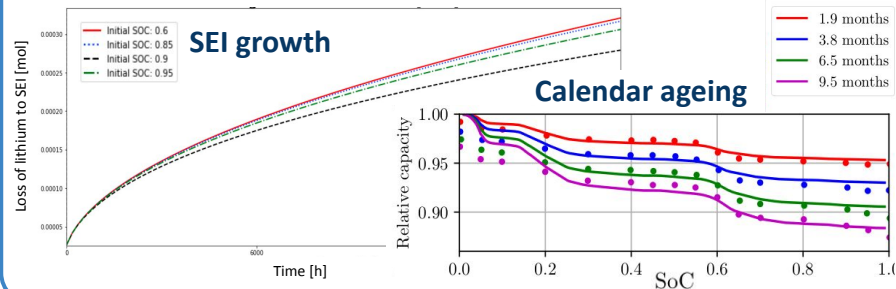
*PyBaMM has fostered an active community with almost **40 contributors** and **hundreds of users**. Our most recent training workshop attracted almost **400 participants**, over **100** of which were from **industry**.*

Python Battery Mathematical Modelling

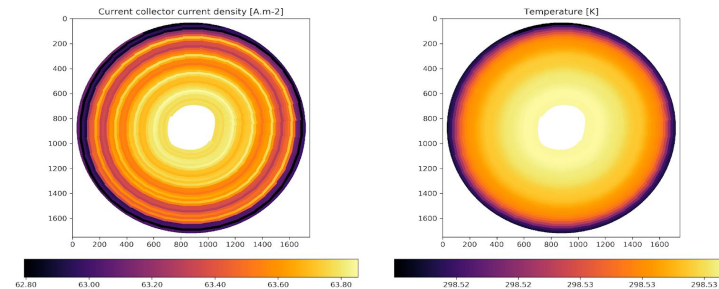
A design tool for predicting behaviour, a virtual physics lab to explore new mechanisms, a quality control environment to compare models, and a testing ground to validate model predictions.



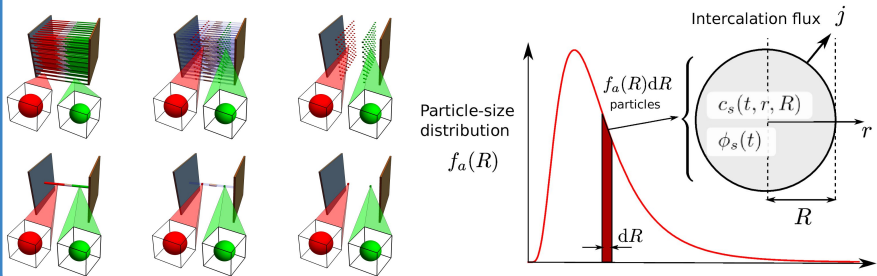
INCORPORATE "PLUG AND PLAY" PHYSICS



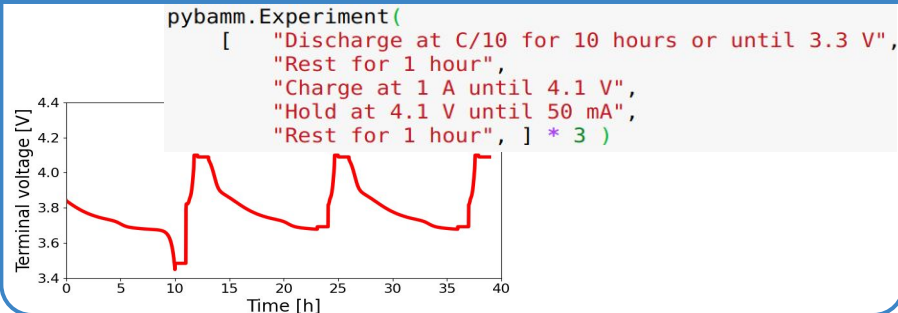
INVESTIGATE CELL DESIGN



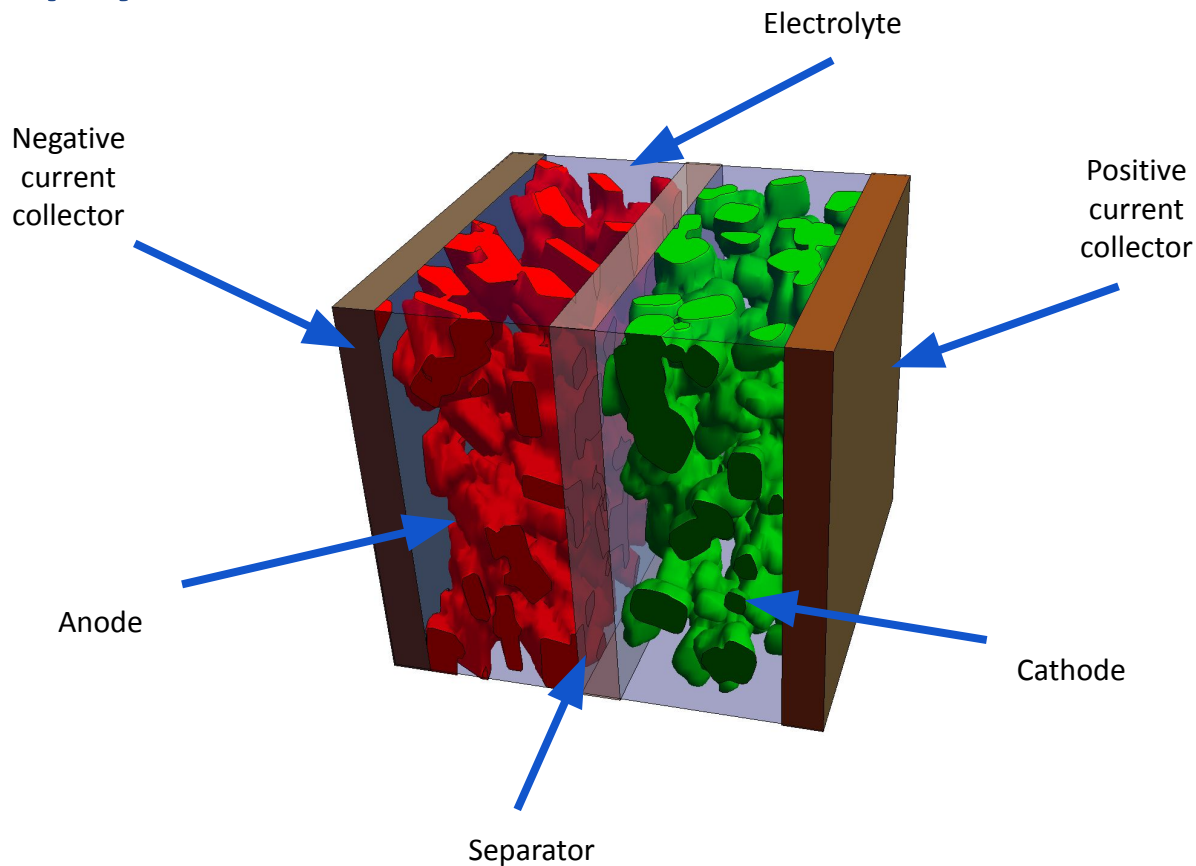
COMPARE MODELS FOR ACCURACY AND SPEED



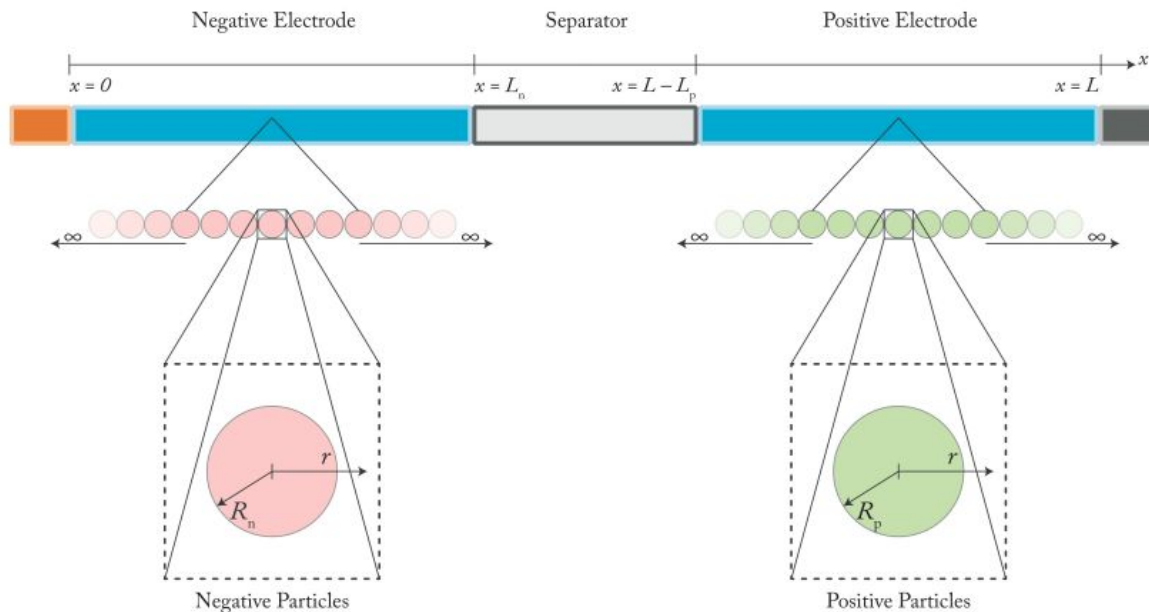
SIMULATE EXPERIMENTAL PROTOCOLS



What is a physics-based model?

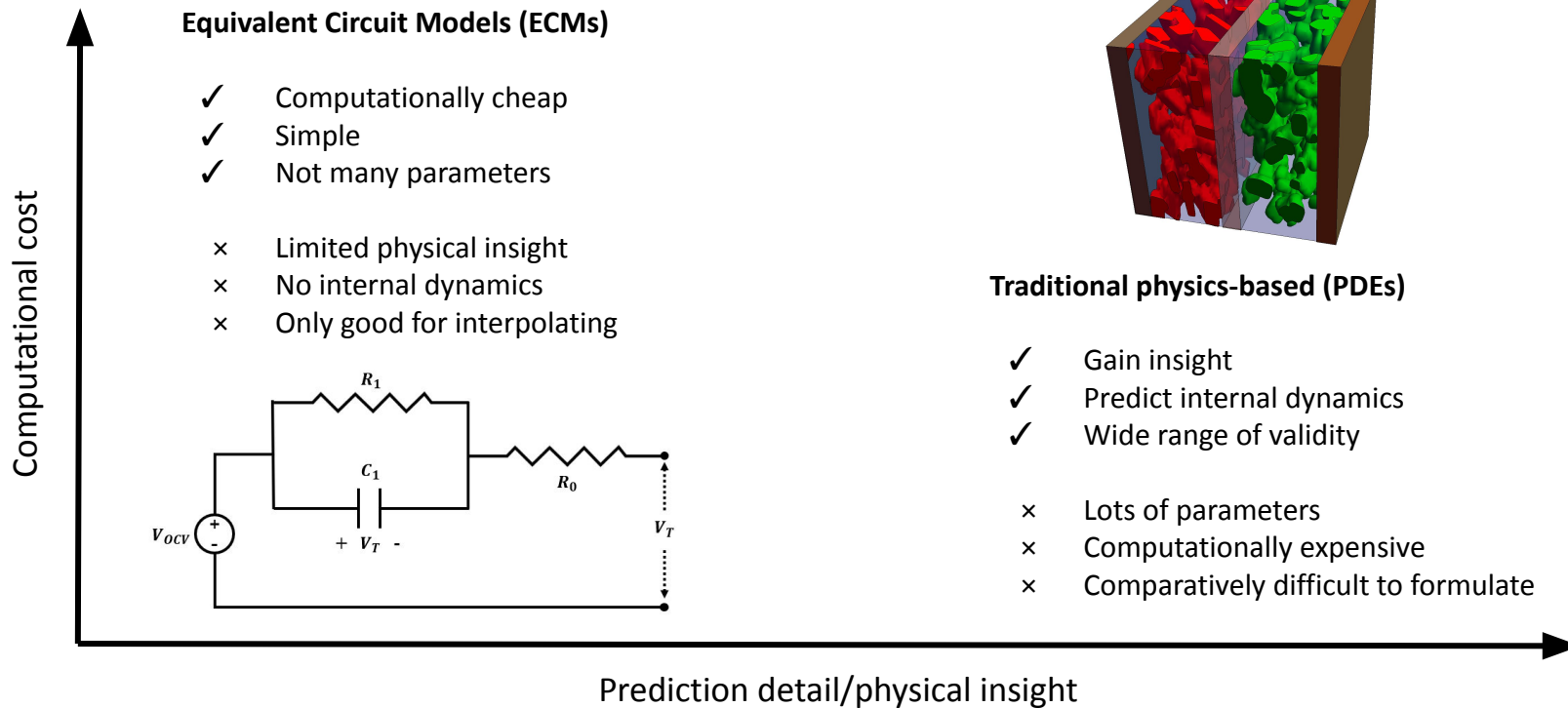


What is a physics-based model?

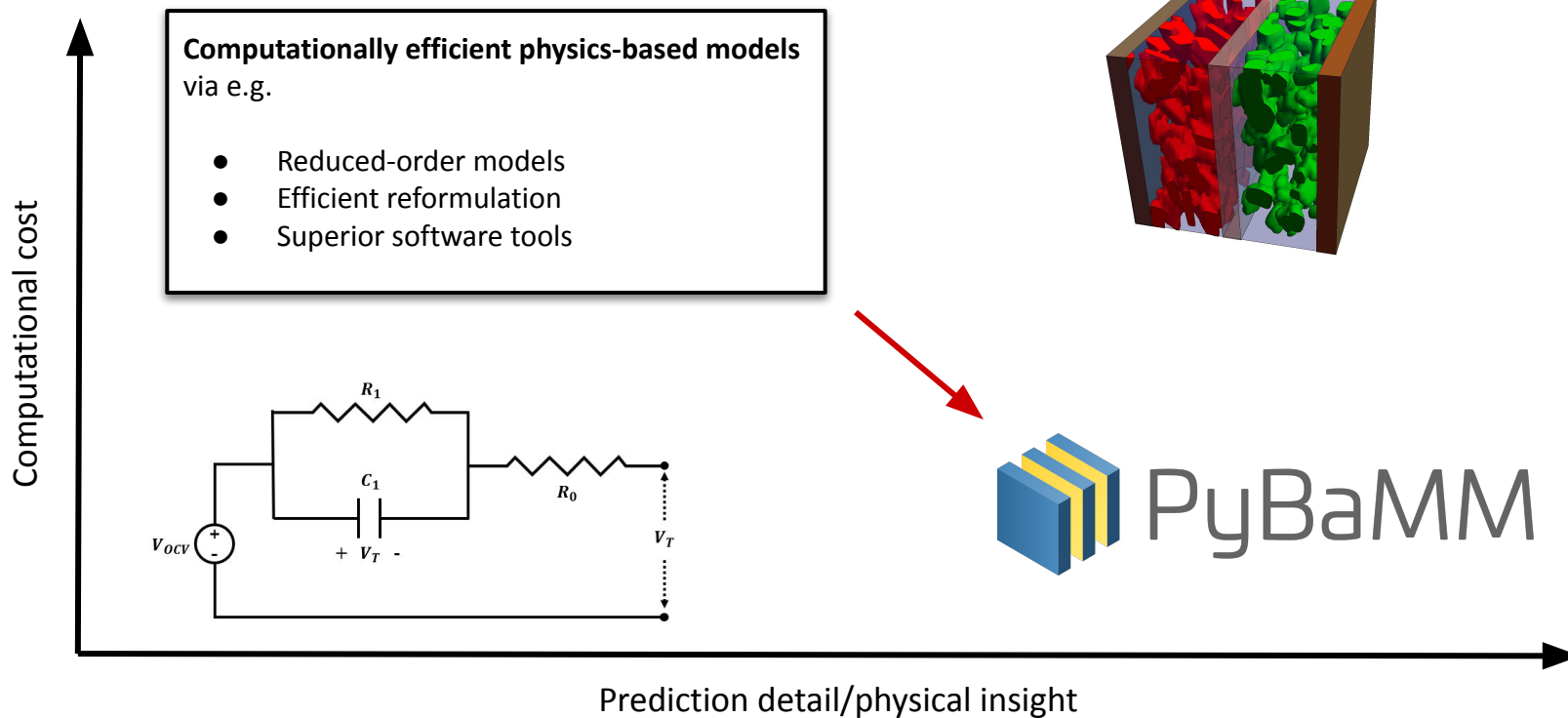


“Doyle-Fuller-Newman model” or “Newman model” or “Pseudo two-dimensional model”

Why physics-based models?



Why physics-based models?

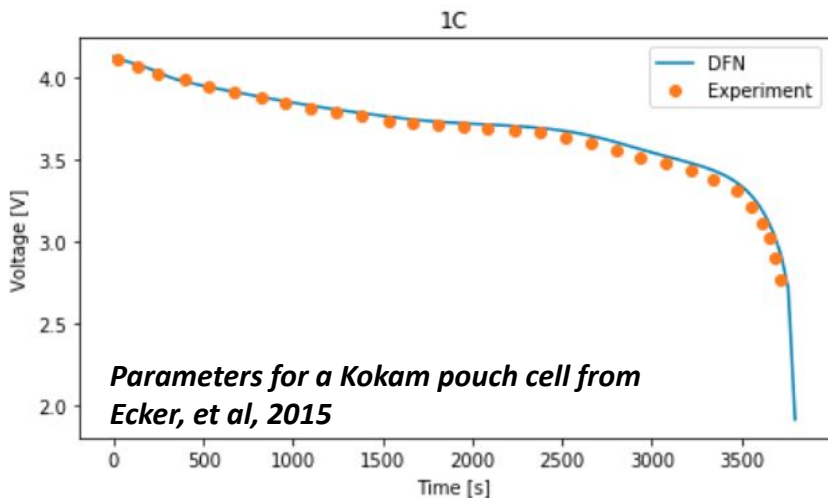


Fast, reliable simulations via a simple interface

PyBaMM's user friendly interface makes simulating battery behaviour easy

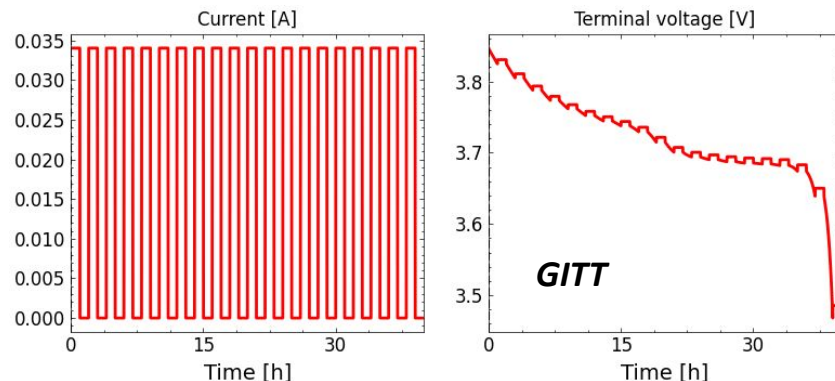


TEST & VALIDATE MODELS & PARAMETERISATIONS



*1 hour discharge of the DFN model with **over 20,000 states** solves in less than **3s** on a standard laptop. With **300 states** the solve time is just **35ms**.*

EASILY DEFINE EXPERIMENTAL PROTOCOLS OR PROVIDE DRIVE CYCLE DATA



```
import pybamm

experiment = pybamm.Experiment(
    [("Discharge at C/20 for 1 hour", "Rest for 1 hour")] * 20,
)
model = pybamm.lithium_ion.DFN()
sim = pybamm.Simulation(model, experiment=experiment)
sim.solve()
sim.plot()
```

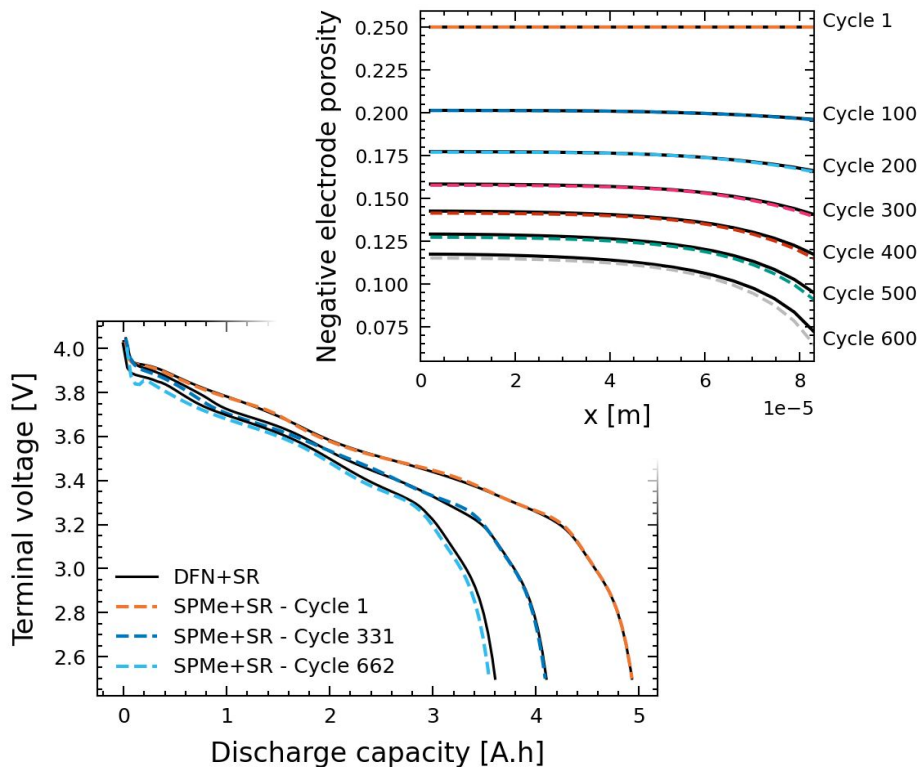

Advanced multi-physics battery models

PyBaMM's flexible submodel structure enables true plug and play physics



- **Comprehensive** model library, including full and reduced-order models
- **Plug and play** submodels describe **key physics** and are all **fully coupled**
- Connections between submodels allow **multiple physical effects and their interactions to be understood**
- Predict cell **performance and lifetime** through coupled **electrochemical-thermal-mechanical degradation** models
- Track changes in behaviour through various **health metrics**, e.g. capacity fade, LLI, LAM

Extra options include SEI growth, lithium plating, mechanics, particle size distributions, and more!



Accessing support



PyBaMM doesn't do something you think it should?

Not sure how to use existing features or implement your own model?

Interested in commercial or academic collaboration?

Please get in touch!

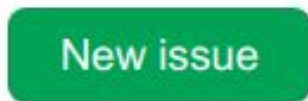
Get in touch on Slack or email

Use the #technical-questions channel, send a direct message, or send us an email



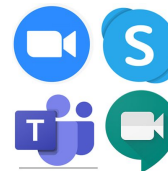
Create an issue or discussion

Head over to GitHub and post an issue or start a discussion



Implementation Sessions

One-on-one or group session to workshop your ideas, discuss implementation details and help with any issues



Accessing support



PyBaMM doesn't do something you think it should?



*New pack modelling software in PyBaMM ecosystem
currently under active development!*

Visit <https://github.com/pybamm-team/liionpack>

send a direct message, or send us an email

start a discussion

workshop your ideas, discuss implementation details and help with any issues



New issue

