

1. We have developed the initial version of our in-house code based on deal.II tutorial step-46. This code is capable of addressing fluid-structure interaction (FSI) problems within the hp-nothing formulation framework.

First, familiarize yourself with the fundamental concepts of deal.II.

<https://www.dealii.org/current/doxygen/deal.II/index.html>

<http://dealii.org/current/doxygen/deal.II/Tutorial.html>

Then, familiarize yourself with the member functions used in step-46

[https://www.dealii.org/current/doxygen/deal.II/step\\_1.html](https://www.dealii.org/current/doxygen/deal.II/step_1.html)

[https://www.dealii.org/current/doxygen/deal.II/step\\_2.html](https://www.dealii.org/current/doxygen/deal.II/step_2.html)

[https://www.dealii.org/current/doxygen/deal.II/step\\_3.html](https://www.dealii.org/current/doxygen/deal.II/step_3.html)

[https://www.dealii.org/current/doxygen/deal.II/step\\_4.html](https://www.dealii.org/current/doxygen/deal.II/step_4.html)

[https://www.dealii.org/current/doxygen/deal.II/step\\_6.html](https://www.dealii.org/current/doxygen/deal.II/step_6.html)

[https://www.dealii.org/current/doxygen/deal.II/step\\_8.html](https://www.dealii.org/current/doxygen/deal.II/step_8.html)

[https://www.dealii.org/current/doxygen/deal.II/step\\_22.html](https://www.dealii.org/current/doxygen/deal.II/step_22.html)

[https://www.dealii.org/current/doxygen/deal.II/step\\_27.html](https://www.dealii.org/current/doxygen/deal.II/step_27.html)

[https://www.dealii.org/current/doxygen/deal.II/step\\_46.html](https://www.dealii.org/current/doxygen/deal.II/step_46.html)

2. If you decide to use the external mesh file generated by GMSH (complex geometry case), you may need to refer to

<https://www.dealii.org/current/doxygen/deal.II/classGridIn.html>

[https://www.dealii.org/current/doxygen/deal.II/step\\_49.html](https://www.dealii.org/current/doxygen/deal.II/step_49.html)

<https://gmsh.info/doc/texinfo/gmsh.html>

3. If you've understood all the material mentioned above, you should be able to follow the in-house code. I assumed that you have read our JCP paper: <https://www.sciencedirect.com/science/article/abs/pii/S0021999121007464?via%3Dihub>

In our code, we have numerous template functions that can be roughly divided into three complexity levels based on either their conceptual difficulty or their code complexity.

Level 1:

**Identify fluid and solid regions.**

This includes template functions such as `cell_is_in_fluid_domain()` and `cell_is_in_hydrogel_domain()`

**Create mesh and finite element discretizations in each corresponding region.**

It involves template functions such as, `make_grid()`, `create_fe_multiplicities()`, and `set_active_fe_indices()`

**Distribute degrees of freedom.**

Template function: `setup_dofs()`.

**Set up boundary conditions.**

Template function: `make_boundary_constraints_hp.`

**Output the results.** `output_results()`

Level 2:

**Assemble and solve the system**

Template function:

`local_assemble_volume()`

`copy_local_to_global_volume()`

`assemble_volume_system_workstream()`

`solve_volume()`

`local_assemble_hp()`

`copy_local_to_global_hp()`

`assemble_system_workstream()`

`newton_iteration()`

Level 3:

**Setup constraints for the system.**

`get_physical_face_points()`

`add_interface_constraints()`

`make_flux_constraints()`

`update_constraints()`

`set_interface_dofs_flag()`