Hydrothermal FOAM Manual

Release 1.0

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Welcome to the HydrothermalFoam Manual! Here you will find resources for using HydrothermalFoam and examples of what it can do.
INTRODUCTION

HydrothermalFoam — a combination of hydrothermal and OpenFOAM — a three-dimensional hydrothermo-transport model designed to resolve fluid flow within submarine hydrothermal circulation systems. HydrothermalFoam has been developed on the OpenFOAM platform, which is a Finite Volume based C++ toolbox for fluid-dynamic simulations and for developing customized numerical solvers that provide access to state-of-the-art parallelized solvers and to a wide range of pre- and post-processing tools. We have implemented a porous media Darcy-flow model with associated boundary conditions designed to facilitate numerical simulations of submarine hydrothermal systems. The current implementation is valid for single-phase fluid states and uses a pure water equation-of-state (IAPWS-97). We present the model formulation, OpenFOAM implementation details, and a sequence of 1-D, 2-D and 3-D benchmark tests. The source code repository further includes a number of tutorials that can be used as starting points for building specialized hydrothermal flow models.

1.1 Features

Features of the HydrothermalFoam are summarized as following:

- **Original characteristics of OpenFOAM**: HydrothermalFoam keeps all the original characteristics of OpenFoam, for example, file structure of case, syntax of all the input files and output files, mesh, utilities, even part of variable names in the source code are kept the same. This principle has two advantages, one is that it is easy to understand and to use HydrothermalFoam if you are a OpenFoam user. The other is that it is easy to compare and understand HydrothermalFoam solver and the other standard solvers in OpenFoam.

- **Mesh**: HydrothermalFoam supports both structured regular mesh and unstructured mesh. The internal structured regular mesh tool, blockMesh, is recommended for new users. While Gmsh is also an excellent open source unstructured mesh generator, and there is a utility named gmshToFoam can transfoam gmsh to OpenFoam mesh.

- **Boundary conditions**: even though OpenFoam has a lot of build-in boundary conditions, we also developed some specific boundary conditions, e.g. noFlux, HydrothermalHeatFlux for the specific problem — Hydrothermal system.
1.2 Information

- License: GNU General Public License v3.0
- GitLab repository: https://gitlab.com/gmdpapers/hydrothermalfoam
- Docker Hub repository: zguo/hydrothermalfoam

1.3 Dependent package

- OpenFOAM
- freesteam
- Gmsh
Tip: We provide a 5-minutes quick start video (on Mac OS) that can be accessed here https://youtu.be/6czcxC90gp0.

There are two ways to install HydrothermalFoam. The easiest way is installing via Docker image, see Section 2.1. Another way is building from source code if user has experience of OpenFOAM installation, see section of Build from source.

2.1 Docker image

In order to quick start to use HydrothermalFoam for non-Ubuntu users, we provided a pre-compiled docker image which can be found on Docker Hub repository, named zguo/hydrothermalfoam. It’s pretty simple to install HydrothermalFoam via Docker, all the steps are summarized below,

1. Install Docker desktop and keep it running.

2. Pull the docker image of HydrothermalFoam by running command of docker pull zguo/hydrothermalfoam.

3. Install a container from the docker image by running shell script which can be found in source code directory of docker (see also Listing 2.1 and Listing 2.2). The directory named HydrothermalFoam_runs directory is a shared folder between the container and host machine.

```
homeInHost=${1:-$HOME}
dirInContainer="/home/openfoam/HydrothermalFoam_runs"
dirInHost="${homeInHost}/HydrothermalFoam_runs"
imageName="zguo/hydrothermalfoam"
containerName="hydrothermalfoam"
# List container in docker environment
echo "*********************************************************

Listing 2.1: Script for Mac OS (installMacHydrothermalFoam.sh)"
```

(continues on next page)
4. Start the container by running command of `docker start hydrothermalfoam`.

5. Attach the container by running command of `docker attach hydrothermalfoam`.

The user now in a Ubuntu Linux environment with precompiled HydrothermalFoam tools which located at directory of `~/HydrothermalFoam`. We recommend user run HydrothermalFoam cases in the directory of `HydrothermalFoam_runs` in the container, and then the results are synchronized in the shared directory in the host, and thus can be visualized by ParaView, Tecplot or other CFD post-processing software.
2.2 Build from source

2.2.1 Install OpenFOAM

The HydrothermalFoam v1.0 is developed based on OpenFOAM-7, which can be installed according to the installation instructions (https://openfoam.org/download/) given by the development team for Ubuntu Linux, Other Linux, macOS and Windows platform, respectively.

2.2.2 Build HydrothermalFoam

Once OpenFOAM is built successfully, the source code of HydrothermalFoam be downloaded from Zenodo or from GitLab repository. The directory structure and components of HydrothermalFoam are shown in Fig. 2.1 and the components can be built follow three steps below.

**Note:** The following steps are only proper for Mac OS and Linux systems, we do not yet build HydrothermalFoam on Windows system directly. If users using ubuntu sub-system on Windows 10, the following steps could work in the sub-system.

1. **Build freesteam-2.1 library.** The freesteam project is constructed by scons, which is an open source software construction tool dependent on python 2, and based on GSL (GNU Scientific Library). Therefore python 2, scons and GSL have to be installed firstly, then change directory to freesteam-2.1 in HydrothermalFoam source code and type command of scons INSTALL_PREFIX=$FOAM_USER_LIBBIN install to compile freesteam library named libfreesteam.so. See home page of freesteam-2.1 project for more details.

2. **Build libraries of customized boundary conditions and thermo-physical model.** Change directory to libraries and type command of ./Allmake to compile the libraries named libHydroThermoPhysicalModels.so, libHydrothermalBoundaryConditions.so.

3. **Build solver of HydrothermalSinglePhaseDarcyFoam.** Change directory to HydrothermalSinglePhaseDarcyFoam and type command of wmake to compile the solver named HydrothermalSinglePhaseDarcyFoam.

**Note:** All the library files and executable application (solver) file will be generated in directories defined by OpenFOAM’ s path variables of FOAM_USER_LIBBIN and FOAM_USER_APPBIN, respectively. If build HydrothermalFoam in Mac OS, the extension of the library files is .dylib, please make a symbolic links. For example, see following command for libHydroThermoPhysicalModels.dylib

```bash
ln -s $FOAM_USER_LIBBIN/libHydroThermoPhysicalModels.dylib $FOAM_USER_LIBBIN/
```

2.2. Build from source 7
HydrothermalFoam

solvers
- HydrothermalSinglePhaseDarcyFoam

libraries
- BoundaryConditions
  - HydrothermalHeatFlux
  - SubmarinePressure \( p_{\text{seafloor}} = \rho_0 g z \)
  - noFlux \( \nabla p = \rho \vec{g} \)
  - HydrothermalMassFluxPressure

ThermoModels
- freesteam-2.1 \( \leftarrow \text{source code of freesteam2.1} \)

benchmarks
- HydrothermalFoam
  - 1d
  - 2d
  - 3d
  - HYDROTHERMAL

cookbooks
- manual

Fig. 2.1: Structure and components of the HydrothermalFoam toolbox.
CHAPTER
THREE

MODEL DESCRIPTION

3.1 Solver and equations

The solver is named HydrothermalSinglePhaseDarcyFoam, which is designed to resolve fluid flow within submarine hydrothermal circulation systems. The hydrothermal fluid flow is governed by Darcy’s law (Eqn. (3.1)), mass continuity (Eqn. (3.2)) and energy conservation (Eqn. (3.4)) equations shown below,

\[ \vec{U} = -\frac{k}{\mu_f} (\nabla p - \rho \vec{g}) \]  
\[ \varepsilon \frac{\partial \rho_f}{\partial t} + \nabla \cdot (\vec{U} \rho_f) \]  
\[ \varepsilon \rho_f \left( \beta_f \frac{\partial p}{\partial t} - \alpha_f \frac{\partial T}{\partial t} \right) = \nabla \cdot \left( \rho_f \frac{k}{\mu_f} (\nabla p - \rho_f \vec{g}) \right) \]  
\[ (\varepsilon \rho_f C_{pf} + (1 - \varepsilon) \rho_r C_{pr}) \frac{\partial T}{\partial t} = \nabla \cdot (\lambda_r \nabla T) - \rho_f C_{pf} \vec{U} \cdot \nabla T + \frac{\mu_f}{k} \| \vec{U} \|^2 - \left( \frac{\partial p}{\partial n} \rho_f \right) \frac{Dp}{Dt} \]  

where the pressure equation (3.3) is derived from continuity equation (3.2) and Darcy’s law (3.1) (see [Hasenclever et al., 2014]).

Note: see the manuscript (Guo, Ruepke & Tao, 2020) for details of symbols and mathematical model description.

3.2 Boundary conditions

The available boundary conditions (BCs) of temperature and pressure for HydrothermalSinglePhaseDarcyFoam are presented below.

Tip: Syntax of all the input files of OpenFOAM, thus of HydrothermalFoam is C++ style. The basic form of a boundary condition can be written as following dictionary structure,
patchName
{
    type boundaryConditionType; //compulsive
    value floatNumber; //compulsive
    option1 valueOfOption1; //optional
    option2 valueOfOption2; //optional
}

where type and value are always compulsive keys, and sometime the value key is just a placeholder but have to be there. option* represents some optional parameters for a special boundary condition, e.g. qmax for HydrothermalHeatFlux shown blow.

### 3.2.1 Basic boundary conditions

**Empty BC: empty**

The empty BC is used and only applied on non-computed patches for one-dimensional and two-dimensional models. See Listing 3.1 for example.

Listing 3.1: Example of empty BC for a 2D model

```plaintext
frontAndBack
{
    type empty;
}
```

**Fixed value BC: fixedValue**

This is the basic and commonly used Dirichlet BC in OpenFOAM, \( f = f_0 \) on \( \partial \Omega \), where \( f \) denotes some field and \( \Omega \) represents a boundary patch (*same meaning as belows*). See Listing 3.2 for example.

Listing 3.2: Example of fixedValue BC

```plaintext
bottom
{
    type fixedValue;
    value uniform 473.15; //473.15 K = 200 C
}
```

**Fixed gradient: fixedGradient**

Just like its name, fixedGradient specify a Neumann boundary condition on a patch. \( \nabla f = g_0 \) on \( \partial \Omega \). See Listing 3.3 for example.
Listing 3.3: Example of fixedGradient BC.

```cpp
bottom
{
    type fixedGradient;
    gradient 0.005; //required
}
```

**Zero gradient: zeroGradient**

It is a special case of fixedGradient. \( \nabla f = 0 \) on \( \partial \Omega \). See Listing 3.4 for example.

Listing 3.4: Example of zeroGradient BC for temperature.

```cpp
right
{
    type zeroGradient;
}
```

**Tip:** zeroGradient is always applied for permeability. Because permeability is not a primary variable and thus don’t need to solve it. But we have to regard it as a field variable just like temperature to initialize the field value, because permeability in our model is not always uniform distribution. Therefore we have to specify a boundary condition for permeability.

**Fixed flux pressure: fixedFluxPressure**

This boundary condition sets the pressure gradient to the provided value such that the flux on the boundary is that specified by the velocity boundary condition. This fixedFluxPressure BC for pressure \( p \) is commonly combined with fixedValue BC for velocity \( U \), and of course the velocity field \( U \) have to be set in \( 0 \) folder even though \( U \) is not a primary variable. see Listing 3.5 for example.

**Tip:** We highly recommend the new defined noFlux or hydrothermalMassFluxPressure (see Section 3.2.2, Section 3.2.2) for Neumann boundary condition of pressure.

Listing 3.5: Example of fixedFluxPressure BC

```cpp
right
{
    type fixedFluxPressure;
}
```
**Inlet and outlet BC: inletOutlet**

This is a generic outflow boundary condition and is commonly applied to temperature $T$ at seafloor boundary. It is a mixed boundary condition that using `zeroGradient` BC when fluid flow out of the boundary and using a `fixedValue` BC when fluid flow into the boundary. See Listing 3.6 for example and options.

**Listing 3.6: Example of inletOutlet BC**

```plaintext
{  type               inletOutlet;
    phi;             //optional
    inletValue       uniform 278.15; //required, fixed value for inflow
    value            uniform 278.15; //required, recommend the same value as...
}_inletValue
```

### 3.2.2 Customized boundary conditions for pressure

The following customized boundary conditions of pressure $p$ are designed for seafloor hydrothermal models.

**Zero mass flux BC: noFlux**

This boundary condition is always applied on impermeable insulating boundary, e.g. side wall. See Listing 3.7 for example and options.

**Listing 3.7: Example of noFlux BC**

```plaintext
left  {
    type    noFlux;
}
right {
    type    noFlux;
}
```

**Seafloor BC: submarinePressure**

This is a spatial coordinate dependent Dirichlet BC and is derived from `fixedValue` BC. The pressure boundary value is the hydrostatic pressure on the boundary patch, which is calculated from coordinate $y$ of the patch. This is designed for seafloor boundary. See Listing 3.8 for example.
Listing 3.8: Example of submarinePressure BC

top
{
    type submarinePressure;
}

Mass flux BC: hydrothermalMassFluxPressure

According Darcy’s law (3.1), gradient of pressure can be expressed by velocity, thus expressed by mass flux. See Listing 3.9 for example.

Listing 3.9: Example of hydrothermalMassFluxPressure BC

bottom
{
    type hydrothermalMassFluxPressure;
    q uniform -0.015; // inflow
}

where \( q \) in the dictionary denotes mass flux value with unit of \( \text{kg/m}^2/\text{s} \). If mass flux represents inflow through the boundary, the value is negative, otherwis is positive.

3.2.3 Customized boundary conditions for temperature

Heat flux BC: hydrothermalHeatFlux

A commonly used Neumann boundary condition of temperature \( T \) in hydrothermal modeling is heat flux (\( W/\text{m}^2 \)) BC. For example, it is applied on bottom boundary representing heat source. A basic example is shown in Listing 3.10, heat flux on the boundary patch (named heatsource) is a constant equal to 5 \( W/\text{m}^2 \).

Listing 3.10: Example of hydrothermalHeatFlux BC

heatsource
{
    type HydrothermalHeatFlux;
    shape fixed; //Optional, default is fixed.
    q uniform 5; //W/m^2
    value uniform 0; //Placeholder
}

The hydrothermalHeatFlux also support Gaussian shape heat flux, see equation (3.5) and Fig. 3.1.

\[
q_h(x, z) = q_{min} + (q_{max} - q_{min})e^{-\frac{(x-x_0)^2 + (y-y_0)^2}{2\sigma^2}}
\] (3.5)

See Listing 3.11 and Listing 3.12 for 2D and 3D example, respectively.
Listing 3.11: Example of Gaussian shape heat flux BC for 2D model (see also Fig. 3.1)

```plaintext
bottom {
    type hydrothermalHeatFlux;
    q uniform 0.05;  //placeholder
    value uniform 0;  //placeholder
    shape gaussian2d;
    x0 0;
    qmax 5;
    qmin 0.05;
    c 500;
}
```

Fig. 3.1: Gaussian shape heat flux curve corresponding to Listing 3.11.

Listing 3.12: Example of Gaussian shape heat flux BC for 3D model

```plaintext
bottom {
    type hydrothermalHeatFlux;
    q uniform 0.05;  //placeholder
    value uniform 0;  //placeholder
    shape gaussian2d;
    x0 0;
    z0 0;
    qmax 5;
    qmin 0.5;
}
```

(continues on next page)
3.2.4 Coded boundary conditions

The **fixedValue** BC mentioned above basically specifies a constant boundary condition value. OpenFOAM provided a so-called dynamic compiling mechanism which allows user define a customized fixed value boundary condition, e.g. a coordinate dependent **fixedValue** BC.

- **Coded fixed value BC**: `codedFixedValue`

**Warning**: Some programming experience of OpenFOAM, at least C++ is required to use this boundary condition.

An example of a Gaussian shape fixed temperature boundary condition is shown in Listing 3.13.

Listing 3.13: Example of Gaussian shape fixed temperature ($T$) BC of a 2D model.

```plaintext
bottom
{
    type codedFixedValue;
    value uniform 873.15; //placeholder
    name gaussShapeT;
    code #{
        scalarField x(this->patch(). Cf(). component(0));
        double wGauss=200;
        double x0=1000;
        double Tmin=573;
        double Tmax=873.15;
        scalarField T(Tmin+(Tmax-Tmin)*exp(-(x-x0)^2/(2*wGauss*wGauss)));
    # operator===(T);
    
    #};
}
```

The code shown in Listing 3.13 implements a Dirichlet boundary condition for temperature $T$ which varies along **bottom** boundary with coordinate $x$, the temperature distribution curve is shown in Fig. 3.2. In addition, there is also a **codedMixed** BC available for dynamic compiled mixed boundary condition, see OpenFOAM documentation for more details.

3.3 Properties

The parameters in the governing equations can be classified as transport and thermophysical properties.
3.3.1 Transport properties

The transport properties in HydrothermalFoam are porosity, thermal conductivity $k_r$ of rock, specific heat capacity $c_{p\_rock}$ of rock, density $\rho_{\_rock}$ of rock. All the transport properties are stored in constant/transportProperties file, see Listing 3.14 for example.

Listing 3.14: Example of transport properties.

```foam
FoamFile
{
    version 2.0;
    format ascii;
    class dictionary;
    location "constant";
    object transportProperties;
}

porosity porosity [0 0 0 0 0 0 0] 0.1;
k_r kr [1 1 -3 -1 0 0 0] 2;
c_{p\_rock} cp_rock [0 2 -2 -1 0 0 0] 880;
\rho_{\_rock} rho_rock [1 -3 0 0 0 0 0] 2700;
```

Note: The head (line 1-8 in Listing 3.14) of the transportProperties file always keep the same, users just need to modify the transport properties (line 10-12 in Listing 3.14) for a specific hydrothermal system.
3.3.2 Thermophysical properties

The thermophysical properties describe the equation of state (EOS) and thermodynamic properties of a specific fluid, e.g. water. For single phase hydrothermal circulation modeling, we developed a OpenFOAM thermophysical model (named htHydroThermo) of water based on IAPWS-IF97 and freesteam-2.1 project. Similar as other OpenFOAM thermophysical model for other specific solvers, the usage of water htHydroThermo is shown in Listing 3.15.

Listing 3.15: Usage of thermophysical model of htHydroThermo.

```FoamFile
{
  version 2.0;
  format ascii;
  class dictionary;
  location "constant";
  object thermophysicalProperties;
}

thermoType
{
  type htHydroThermo;
  mixture pureMixture;
  transport IAPWS;
  thermo IAPWS;
  equationOfState IAPWS; //Boussinesq
  specie specie;
  energy temperature; //sensibleEnthalpy
}

mixture
{
  specie
  {
    molWeight 18;
  }
}
```

Note: htHydroThermo is the unique support thermophysical model for solver of HydrothermalSinglePhaseDarcyFoam so far. Therefore, please always copy Listing 3.15 and save it in file of constant/thermophysicalProperties.

Warning: The temperature and pressure limitation of htHydroThermo are [273.15, 1073.15] K and [10⁵, 10⁸] Pa, respectively. Please make sure the temperature and pressure field value are in the valid range.

3.3. Properties
Usage of HydrothermalFoam tool is similar to any other solver of OpenFOAM, see OpenFOAM user guide for more details. The basic usage of HydrothermalFoam are shown as below.

### 4.1 Mesh generation

The are several ways to generat mesh, including OpenFOAM’s built-in blockMesh application for generating meshes of simple geometries, snappyHexMesh application for meshing complex geometries and applications that convert meshes from well known formats into the OpenFOAM format, e.g. fluentMeshToFoam, gmshToFoam. see OpenFOAM user guide-mesh generation and conversion <https://cfd.direct/openfoam/user-guide/v7-mesh/> for details.

This section we simply present how to use blockMesh and gmshToFoam for generating mesh and defining boundary patches.

#### 4.1.1 Simple mesh generation: blockMesh

The mesh is generated from a dictionary file named blockMeshDict located in the system directory of a case. blockMesh reads this dictionary, generates the mesh and writes out the mesh data to points and faces, cells and boundary files in the constant/polyMesh directory (see Fig. 4.5).

blockMesh decomposes the domain geometry into a set of 1 or more three dimensional, hexahedral blocks. Each block of the geometry is defined by 8 vertices, one at each corner of a hexahedron. An simple example block and vertices numbering is shown in Fig. 4.1. The vertices are written in a list so that each vertex can be accessed using its label, remembering that OpenFOAM always uses the C++ convention that the first element of the list has label ‘0’.

An example of blockMeshDict for describing a 2D box is shown in Listing 4.1. The key entries are vertices, blocks and boundary.

- **vertices** contains all three-dimension coordinate of each vertex (line 15-25), e.g. vertex 0 is \((0 \ ymin \ 0)\) and the vertex 5 is \((Lx \ ymin \ Lz)\), the coordinate system is shown in Fig. 4.1.

- **blocks** contains vertex connection of a hexahedron, numbers of cells in each direction and cell expansion ratios. The order of vertex connection is \((0 \ 1 \ 2 \ 3 \ 4 \ 5 \ 6 \ 7)\). \((100 \ 50 \ 1)\) means the number of cells in \(x\), \(y\) and \(z\) direction will be 100, 50 and 1 (for 2D case), respectively.
simpleGrading is typically set to (1 1 1), see OpenFOAM user guide blockMesh for more details.

• boundary contains sub-dictionaries for defining boundary patches. The name (e.g. right or whatever the user like) of the sub-dictionary is the boundary patch name which will be used in filed data (see Section 3.2). There are two key entries, type and faces in each sub-dictionary. The type is typically set to patch or empty for a non-computing boundary patch of a non-three-dimensional case. faces is a list of block faces that make up the patch with a user defined name (see lines 43-46 for right patch or lines 67-71 for frontAndBack patch shown as transparent surface in Fig. 4.1)

Listing 4.1: Example of blockMeshDict for a 2D box.

```plaintext
FoamFile
{
    version 2.0;
    format ascii;
    class dictionary;
    object blockMeshDict;
}

// * * * * * * * * * * * * * * * * * * * * * * * * * * * * * * * * * * * * * //

convertToMeters 1;
Lx 2000;   //variable definition
ymin -3000;
ymax -2000;
Lz 1;
vertices   //vertices definition
{
    (0  $ymin  0) //coordinate of vertex 0
    ($Lx  $ymin  0) //coordinate of vertex 1
    ($Lx  $ymax  0) //coordinate of vertex 2
    (0  $ymax  0) //coordinate of vertex 3
    (0  $ymin  $Lz)//coordinate of vertex 4
}
```

Fig. 4.1: The coordinate system and vertices numbering of single block for blockMesh.
blocks
{
    hex (0 1 2 3 4 5 6 7) (100 50 1) simpleGrading (1 1 1)
};
boundary
{
    left //patch name
    {
        type patch;
        faces //face list
        {
            (0 4 7 3)
        };
    }
    right
    {
        type patch;
        faces
        {
            (2 6 5 1)
        };
    }
    top
    {
        type patch;
        faces
        {
            (3 7 6 2)
        };
    }
    bottom
    {
        type patch;
        faces
        {
            (1 5 4 0)
        };
    }
    frontAndBack //patch name
    {
        type empty;
        faces //face list
        {
        }
    ]
(continues on next page)
Tip: We can also use variable in `blockMeshDict`. For example `ymin = -3000;`, it should be noted that there is no = between variable name and value, and ; is required. In addition, the `convertToMeters` keyword specifies a scaling factor by which all vertex coordinates in the mesh description are multiplied, e.g. `convertToMeters 0.01;` scales to mm.

Then just type `blockMesh` command in the terminal in the root directory of the case after finishing `blockMeshDict`.

4.1.2 Convert gmsh to OpenFOAM: `gmshToFoam`

The OpenFOAM built-in meshing utility, `blockMesh`, can generate some simple mesh, but it is difficulty to describe complex geometry in `blockMeshDict` dictionary file. In this section, we introduce how to use `gmsh` to generate mesh and then use `gmshToFoam` convert the mesh to OpenFOAM format.

Create geometry and define boundary patches

Here we present a example for creating 2D and 3D box geometry, and defining boundary patches. See `gmsh manual` (pdf) for more details about geometry definition and mesh generation.

- Geometry description

The geometry descriptions are stored in a gmsh geometry script file with extension `.geo`. Gmsh script files support both C and C++ style comments and the syntax is similar to C/C++ as well. The user can use gmsh GUI to create a geometry script file, and also can write the geometry script file directly in your favorite text editor, e.g. Visual Studio Code with gmsh extension. An gmsh geometry script file of a 3D box is shown in `Listing 4.2 (box.geo)`.

```plaintext
// 0. define some variables
xmin=0;
xmax=600;
ymin=-200;
ymax=0;
zmin=0;
zmax=100;
lc=10;
```

(continues on next page)
// 1. define points
Point(1) = {xmin, ymax, zmin, lc};
Point(2) = {xmax, ymax, zmin, lc};
Point(3) = {xmax, ymin, zmin, lc};
Point(4) = {xmin, ymin, zmin, lc};

// 2. define lines
Line(1) = {1, 2};
Line(2) = {2, 3};
Line(3) = {3, 4};
Line(4) = {4, 1};

// 3. define line loop and surface
Line Loop(6) = {4, 1, 2, 3};
Plane Surface(6) = {6};

// 4. extrude 2D surface to a 3D volume
Extrude {0, 0, zmax} {
Surface(6);
Layers(1); //set layer number to 1 for 2D model
Recombine;
}

The geometry script file can be opened and visualized by gmsh GUI, see Fig. 4.2. Gmsh supports three-dimensional interactive operation and there are mouse tooltips of number and other informations for each point, line, surface and volume. Therefor the user can easily get the number of each boundary patch and volume (cell region).

Fig. 4.2: Gmsh GUI display a 2D box geometry.

• Boundary patches and cell regions definition

The user can specify a specific name, e.g. `bottom`, for each Surface (gmsh keyword) or a name for boundary patches group, e.g. `frontAndBack`. These boundary patches name will be used to specify
boundary conditions in field data file (see Section 3.2). In addition, the user can also specify specific name for each Volume (gmsh keyword). This volume name will be used to set field distribution, e.g. permeability (see Section 4.3). The specification of boundary patch name and cell region name can be done by Physical keyword, see

```
// 5. define boundary patches via Physical keyword
Physical Surface("frontAndBack") = {28, 6};
Physical Surface("bottom") = {27};
Physical Surface("left") = {15};
Physical Surface("top") = {19};
Physical Surface("right") = {23};
// 6. specify a name for cell region which is used for 'setFields'
Physical Volume("internal") = {1};
```

Tip: Gmsh will specify some default colors for every surfaces and volumes. The user can also set specific color for each patch and volume (see below) to check boundary mesh.

```
// 7. specify different color for different boundary patches
Color Gray{Surface{28, 6};}
Color Red{Surface{27};}
Color Purple{Surface{15};}
Color Pink{Surface{23};}
Color Blue{Surface{19};}
Color Green{Volume{1};}
```

Generate mesh

Mesh generation process is pretty easy, the user can do it using gmsh GUI Mesh -> 3D. The mesh result is shown in Fig. 4.3. Because gmshToFoam in OpenFOAM-7 can only read gmsh .msh file in format of version 2, the user have to export (File -> Export -> Gmsh MESH) the mesh file to Version 2 ASCII format (see Fig. 4.4).

Alternately, the mesh generation and save file can be done by a single line of command (see Listing 4.4).

```
Listing 4.4: Meshing and saving command of gmsh.
gmsh gmsh/mesh.geo -3 -o gmsh/mesh.msh -format msh22
```

Convert mesh

If the mesh file is generated successfully, the user can convert the mesh to OpenFOAM format by running command of gmshToFoam mymesh.msh in the root directory of a case. Then a directory named
Fig. 4.3: Gmsh GUI display a 2D box mesh.

Fig. 4.4: Export msh to Version 2.
polyMesh will be generated in constant directory. All the defined boundary patches are converted to polyMesh/boundary file (see Listing 4.5). If the case is three-dimensional, now the mesh generation process is done. But for 2D case, the user have to set empty boundary patches by modifying the boundary file directly.

Listing 4.5: polyMesh/boundary converted by gmshToFoam.

```plaintext
5
{
    frontAndBack
    {
        type patch;
        physicalType patch;
        nFaces 6240;
        startFace 4600;
    }
    left
    {
        type patch;
        physicalType patch;
        nFaces 20;
        startFace 10840;
    }
    top
    {
        type patch;
        physicalType patch;
        nFaces 60;
        startFace 10860;
    }
    right
    {
        type patch;
        physicalType patch;
        nFaces 20;
        startFace 10920;
    }
    bottom
    {
        type patch;
        physicalType patch;
        nFaces 60;
        startFace 10940;
    }
}
```

Tip: We provide a python script (setEmptyPatch.py) to modify a boundary patch to empty by patch name. For example, set frontAndBack patches to empty by running the following command at case
root directory.

```python
python setEmptyPatch.py frontAndBack
```

### 4.2 Case setup

A model case of HydrothermalFoam basically consists of time directory (e.g. 0 for initial state), constant and system directory (see Fig. 4.5). The sub-directory of polyMesh in constant folder consists mesh files, which are generated by meshing utility, e.g. blockMesh, gmshToFoam, snappyHexMesh, ..., see Section 4.1.

![Diagram of directory structure](image)

Fig. 4.5: Directory structure of a case of HydrothermalFoam.
4.2.1 Initial state

The field data files of initial state are commonly stored in 0 time directory. Of course it can be another time directory, e.g. 1000, which is specified by key of startFrom in controlDict in system directory (see Section 4.2.3). In the initial state directory, the field file of primary variable $T$, $p$ and permeability are compulsive, and $U$ is also required if fixedFluxPressure is applied on a boundary patch for pressure (see Section 3.2.1). An example field data file of $T$ is shown in Listing 4.6, which is a basic structure of dictionary file of a field data. A field data file basically contains its variable type (line 5), object name (line 6), dimension (line 9), internal field value (line 10) and boundary field value (boundary conditions, line 11-42).

Listing 4.6: Example field data of temperature.

```
FoamFile
{
    version 2.0;
    format ascii;
    class volScalarField;
    object T;
}

dimensions [0 0 0 1 0 0];
internalField uniform 278.15; //5 C
boundaryField
{
    left
    {
        type zeroGradient;
    }
    right
    {
        type zeroGradient;
    }
    top
    {
        type fixedValue;
        value uniform 278.15;
    }
    bottom
    {
        type hydrothermalHeatFlux;
        q uniform 0.05;
        value uniform 0;
    }
    heatsource
    {
        type hydrothermalHeatFlux;
        q uniform 5;
    }
```
• **Variable type** (class) and **object name** (object). The See the manuscript(Guo, Ruepke & Tao, 2020) for variable type and object name index.

• **Internal field** (internalField) can be set as uniform (just like line 10 in Listing 4.6) or non-uniform (see Section 4.3).

• **Boundary conditions** (boundaryField). The boundary patch name, e.g. bottom, is defined in mesh file of constant/polyMesh/boundary (see Fig. 4.5 and Section 4.1). The available boundary conditions and their usage can be found in Section 3.2.

---

**Tip:** To avoid making mistakes, user should copy the field data files form any case in cookbooks or benchmarks and then make some changes.

### 4.2.2 Constant

For HydrothermalFoam, the constant directory always contains three files named $g$, thermophysicalProperties, transportProperties respectively, and one folder named polyMesh (see Fig. 4.5).

• **$g** contains gravitational acceleration constant. It is same for basically all cases of OpenFOAM-based solvers. Therefor, please just copy this file from any existed case to a new case.

• **thermophysicalProperties** contains thermophysical model. It is the same for all cases of HydrothermalSinglePhaseDarcyFoam solver, see Section 3.3.2.

• **transportProperties** contains constant parameters about transport, see Section 3.3.1.

• **polyMesh** directory contains necessary mesh files which are automatically generated by meshing utility, e.g. blockMesh, gmshToFoam (see Section 4.1).

### 4.2.3 System

As shown in Fig. 4.5, the **system** directory contains three compulsive directory files of controlDict, fvSchemes and fvSolution, and optional dictionary file, e.g. blockMeshDict for OpenFOAM built-in meshing utility blockMesh.
Time and data input/output control

The controlDict dictionary sets time and data input/output control, the commonly used entries for HydrothermalSinglePhaseDarcyFoam solver are shown in Listing 4.7. See OpenFOAM user guide-controlDict for more details of each entry. A few entries that need to be emphasized are highlighted in the Listing 4.7.

- **application** specifies solver name, which should be HydrothermalSinglePhaseDarcyFoam.

- **libs** contains some shared libraries using in the case. These three libraries listed in lines 46-48 must be included in libs sub-dictionary for customized thermophysical model and boundary conditions.

Listing 4.7: Example entries from a controlDict dictionary.

```plaintext
FoamFile
{
    version 2.0;
    format ascii;
    class dictionary;
    object controlDict;
}

application HydrothermalSinglePhaseDarcyFoam;

startFrom latestTime;
startTime 0;
stopAt endTime;
endTime 6912000000; //86400000000
deltaT 864000;
adjustTimeStep yes;
maxCo 0.8;
maxDeltaT 86400000;
writeControl adjustableRunTime;
writeInterval 86400000;
purgeWrite 0;
writeFormat ascii;
writePrecision 6;
writeCompression off;
timeFormat general;
timePrecision 14;
runTimeModifiable true;

libs
{
    "libfreesteam.so"
    "libHydrothermalBoundaryConditions.so"
    "libHydroThermoPhysicalModels.so"
};
```

Note: There are three options (firstTime, startTime, latestTime) available for the
startFrom keyword entry. startTime specifies start time for the simulation when startFrom startTime;

**Numerical schemes**

The fvSchemes dictionary in the system directory sets the numerical schemes for terms, such as derivatives in equations, that are calculated during a simulation. The commonly used entries of fvSchemes for HydrothermalSinglePhaseDarcyFoam solver are shown in Listing 4.8. See OpenFOAM user guide-fvSchemes for more available scheme options of each terms.

As shown in Listing 4.8, user have to specify numerical scheme for transient (lines 11-14), gradient of pressure p (line 19) and temperature T (line 20), laplacian terms (line 33-34), surface interpolation (lines 37-40), surface gradient (lines 42-45). In addition, the fluxRequired sub-dictionary have to be specified for reconstructing Darcy velocity from flux after solving pressure. See Model development section in the manuscript(Guo, Ruepke & Tao, 2020) or source code the solver for more detail implementation of each terms.

Listing 4.8: Example entries from a fvSchemes dictionary.

```plaintext
FoamFile
{
    version 2.0;
    format ascii;
    class dictionary;
    location "system";
    object fvSchemes;
}

ddtSchemes
{
    default Euler;
}
gradSchemes
{
    default none;
    grad(p) Gauss linear;
    grad(T) Gauss linear;
}
divSchemes
{
    default none;
    div(phi,T) Gauss upwind;
    div((phi*interpolate(Cp)),T) Gauss vanLeer;
}
laplacianSchemes
{
    default none;
}
```
Tip: The basic numerical schemes of HydrothermalSinglePhaseDarcyFoam solver are shown in Listing 4.8, user can other available schemes of each term (see OpenFOAM-fvSchemes), e.g. Gauss vanLeer for div(\phi, T).

Solution and algorithm control

The equation solvers, tolerances and algorithms are controlled from the fvSolution dictionary in the system directory. An example set of entries from the fvSolution dictionary required for the HydrothermalSinglePhaseDarcyFoam solver is shown in Listing 4.9. See OpenFOAM user guide-fvSolution for more details of each entry.

As shown in Listing 4.9, user have to specify linear equations solver, preconditioner, tolerance and relTol for pressure and temperature in sub-dictionary of p and T, respectively. In addition, a sub-dictionary named in format of *Final is required as well (see lines 20-24 and 32-36). The prefix of * denotes primary variable name, e.g. p for pressure. The key entry relTol is typically set to 0. For HydrothermalSinglePhaseDarcyFoam solver, key entry PIMPLE is also required besides solvers, because we adopt pimple.correctNonOrthogonal() for non-orthogonal correction. The key entry nNonOrthogonalCorrectors in PIMPLE dictionary specifies repeated solutions of the pressure equation, used to update the explicit non-orthogonal correction of Laplacian term (see OpenFOAM user guide-Surface normal gradient schemes for more details). nNonOrthogonalCorrectors is typically set to 0 or 1.

Listing 4.9: Example entries from a fvSolution dictionary.
format ascii;
class dictionary;
location "system";
object fvSolution;

solvers
{
    p
    {
        solver PCG;
        preconditioner DIC;
        tolerance 1e-12;
        relTol 0;
    }
    pFinal
    {
        $p;
        relTol 0;
    }
    T
    {
        solver PBiCG;
        preconditioner DILU;
        tolerance 1e-06;
        relTol 0;
    }
    TFinal
    {
        $T;
        relTol 0;
    }
}
PIMPLE
{
    nNonOrthogonalCorrectors 0;
}

Warning: The entries and values shown in Listing 4.9 are the recommended options, it is unnecessary to modify them unless the user understands the source code of the solver and OpenFOAM and then wants to test other available parameters of fvSolution.
4.3 Set fields

The user can set a specific value for a field (e.g. permeability) in a specific cell region using utility setFields, which reads dictionary file setFieldsDict in system directory. An example of setFieldsDict is shown in Listing 4.10 (see Section 5.2).

Listing 4.10: Example of setFieldsDict file.

```
FoamFile
{
    version 2.0;
    format ascii;
    class dictionary;
    location "system";
    object setFieldsDict;
}
defaultFieldValues
{
    volScalarFieldValue permeability 1e-15
};
regions
{
    zoneToCell
    {
        name "layer2A";
        fieldValues
        {
            volScalarFieldValue permeability 4e-14
        }
    }
    zoneToCell
    {
        name "layer2B";
        fieldValues
        {
            volScalarFieldValue permeability 1e-15
        }
    }
};
```

The option (e.g. layer2A in line 17) of key entry name in zoneToCell sub-dictionary in Listing 4.10 is defined as Physical Volume("xxx")=...; in gmsh geometry (.geo) file (see Section 4.1.2). Another commonly used key entry is boxToCell which sets field value in a box defined by the two ends of the diagonal, see Listing 4.11.
4.4 Run Case

The user can run a case just by running command of HydrothermalSinglePhaseDarcyFoam in the root directory of a case. But several processes mentioned above, e.g. mesh generation and/or mesh conversion, empty boundary type modification, case setup and initial field setting, have to be done before running a case. All the pre-processing steps can be assembled into a bash file, e.g. run.sh shown in Listing 4.12.

Listing 4.11: Example of boxToCell in setFieldsDict file.

```plaintext
regions
{
    boxToCell
    {
        box (0 0 0) (10 10 10);
        fieldValues
        {
            volScalarFieldValue permeability 4e-14
        }
    }
}
```

Listing 4.12: Example of command set(run.sh) to run a case.

```plaintext
#!/bin/sh

cd ${@%/*} || exit 1  # Run from this directory

# 1. Source tutorial run functions
. $WM_PROJECT_DIR/bin/tools/RunFunctions

# 2. get solver name
application=`getApplication`

# 3. clean case if necessary
./clean.sh

# 4. generate mesh file using gmsh command
gmsh gmsh/mesh.geo -3 -o gmsh/mesh.msh -format msh22

# 5. convert gmsh to OpenFOAM format
gmshToFoam gmsh/mesh.msh

# 6. set empty patchcs for 2D case
python setEmptyPatch.py frontAndBack

# 7. set fields
runApplication setFields

# 8. run a case using single thread
runApplication $application
```

(continues on next page)
Listing 4.13: Example of command set(clean.sh) to clean a case.

```bash
#!/bin/sh
cd $(0%/*) || exit 1 # run from this directory

# Source tutorial run functions
. $WM_PROJECT_DIR/bin/tools/CleanFunctions
cleanCase
```

**Note:** If the user want to run a case in parallel, just need to comment line 20 in Listing 4.12 and uncomment lines 23-25 in Listing 4.12. In addition, the user have to setup the decomposeParDict dictionary file in system directory (see Listing 4.14 for example). See OpenFOAM user guide for more details about parallel computing.


```plaintext
FoamFile
{
  version 2.0;
  format ascii;
  class dictionary;
  location "system";
  object decomposeParDict;
}

numberOfSubdomains 4;
method scotch;

simpleCoeffs
{
  n (2 2 1);
  delta 0.001;
}

hierarchicalCoeffs
{
```

(continues on next page)
Note: It is of course possible to set up input files for cases completely from scratch. However, in practice, it is often simpler to go through the list of cookbooks already provided and find one that comes close to what you want to do. You would then modify this cookbook until it does what you want to do. The advantage is that you can start with something you already know works, and you can inspect how each change you make – changing the details of the geometry, changing the boundary conditions, or changing initial field distribution – affects what you get.

4.5 Post-processing and visualization

The commonly used post-processing tool is ParaView. There is a built-in utility paraFoam, which is based on ParaView, can read and render generic results of OpenFOAM case. The user can run the paraFoam command in the root directory of case direction, or run command of paraFoam -case <caseDir> in any other directory.

Tip: The new version (e.g. 5.5.0) of ParaView has OpenFOAM case reader which will read a file with extension of .foam in the root directory of a case. Therefore, if the user install OpenFOAM or HydrothermalFoam tool via Docker (see Section 2.1 and video), the results can be visualized in host through the shared folder. The results will saved in the shared folder when running a case in the container, and then create a empty file with extension of .foam, e.g. results.foam, in the shared folder. The user can open file results.foam in the shared folder in host by ParaView to display the results.
5.1 Convection in a box

In this first example, let us consider a simple situation: a 2d box that is heated from below, insulated at the side walls, and cooled from the top. We will start from a 2D box example and then make some changes to show its features.

5.1.1 Hello World: 2D box

All the input files can be found in `cookbooks/helloworld` directory.

This is the first simplest example to show the basic steps to run a case. The geometry and boundary conditions are shown in figure Fig. 5.1. The following steps present how to setup a case from scratch, see `helloworld` case.

![Fig. 5.1: The geometry and boundary conditions of the 2D box model.](image-url)
Step 1: create controlDict file

First thing we have to do is create a case directory, e.g. named helloworld, and sub-directory system (see Listing 5.1).

Listing 5.1: Create case directory.

```bash
1   mkdir helloworld
2   cd helloworld
3   mkdir system
```

The controlDict file is the first and most important file, which is located at system directory. Create controlDict file in system directory with script shown in Listing 5.2.

Listing 5.2: controlDict.

```plaintext
FoamFile
{
    version 2.0;
    format ascii;
    class dictionary;
    object controlDict;
}

application HydrothermalSinglePhaseDarcyFoam;
startFrom latestTime;
startTime 0;
stopAt endTime;
endTime 25920000000;
deltaT 86400000;
writeControl adjustableRunTime;
writeInterval 86400000;
purgeWrite 0;
writeFormat ascii;
writePrecision 6;
writeCompression off;
timeFormat general;
timePrecision 14;
runTimeModifiable true;
libs
{
    "libfreesteam.so"
    "libHydrothermalBoundaryConditions.so"
    "libHydroThermoPhysicalModels.so"
};
```

Note: The key entry application has to be set to HydrothermalSinglePhaseDarcyFoam which
is the solver name.

### Step 2: mesh generation

In this step we create dictionary `blockMeshDict` (just copy Listing 4.1) in `system` to setup geometry of the 2D box. Then we just run command `blockMesh` to generate mesh (see Fig. 5.1). The directory `constant` will be created automatically and the mesh files will be generated in `polyMesh` directory.

**Note:** After step 2, the mesh is basically generated and we can display using `paraFoam` utility default or using ParaView for Docker user. The mesh result is shown in Fig. 5.2.

![Fig. 5.2: The screenshot of 2D box mesh in ParaView.](image)

### Step 3: field data in 0 folder

In this step, we need to create I/O data files of field `p`, `T` and permeability in 0 time directory. This three files are shown in `lst:helloworld:p`, `lst:helloworld:T` and `lst:helloworld:permeability`, respectively.

**Listing 5.3: Data field p.**

```plaintext
FoamFile
{
    version 2.0;
    format ascii;
    class volScalarField;
    object p;
}
```

(continues on next page)
dimensions [1 -1 -2 0 0 0 0];
internalField uniform 300e5; //300e5 Pa = 300 bar

boundaryField
{
  left
  {
    type noFlux;
  }
  right
  {
    type noFlux;
  }
  top
  {
    type fixedValue;
    value uniform 300e5;
  }
  bottom
  {
    type noFlux;
  }
  frontAndBack
  {
    type empty;
  }
}

Listing 5.4: Data field T.

FoamFile
{
  version 2.0;
  format ascii;
  class volScalarField;
  object T;
}

dimensions [0 0 0 1 0 0 0];
internalField uniform 278.15;  //278.15 K = 5 C

boundaryField
{
  left
  {
    type zeroGradient;
  }
  right
  {
    type zeroGradient;
  }
  top
  {
    type inletOutlet;
    phi phi;
    inletValue uniform 278.15;
  }
  bottom
  {
    type fixedValue;
    value uniform 873.15;
  }
  frontAndBack
  {
    type empty;
  }
}

// ************************************************************************* //

FoamFile
{
  version 2.0;
  format ascii;
  class volScalarField;
  location "0";
  object permeability;
}

// ************************************************************************* //

dimensions [0 2 0 0 0 0 0];

internalField uniform 1e-14;

Listing 5.5: Data field permeability.

5.1. Convection in a box
Tip: Now we can display the initial field of \( p \), \( T \) and permeability using ParaView. Screenshot of field \( T \) is shown in Fig. 5.3. In order to view the initial field, the user have to uncheck the checkbox of Skip Zero Time (see red rectangle in the figure).

**Step 4: constant property files**

In this step we need to copy three constant files \( g \) (shown in Listing 5.6), thermophysicalProperties (Listing 3.15) and transportProperties (Listing 3.14) into the constant directory.

Listing 5.6: Constant property file \( g \).

```
FoamFile
{
    version 2.0;
    format ascii;
    class uniformDimensionedVectorField;
    location "constant";
```

(continues on next page)
Fig. 5.3: The screenshot of field $T$ in ParaView.

(continued from previous page)

```
object g;

// * * * * * * * * * * * * * * * * * * * * * * * * * * * * * * * * * * * * * //
dimensions [0 1 -2 0 0 0 0 0];
value (0 -9.81 0);
// ************************************************************************* //
```

Tip: The user can change some propertie values in `transportProperties`, e.g. porosity, according to a specific modeling problem.

### Step 4: setup numerical schemes and solution control

In this step we will create numerical schemes and solution control dictionary file `fvSchemes` and `fvSolution` in `system` directory. These two files can be found in Listing 4.8 and Listing 4.9 in Section 4.2.3.

### Step 5: run the case

Now the simple case setup or pre-processing has been completed. The we just run the following simple command (Listing 5.7) to run the case.

5.1. Convection in a box
Listing 5.7: Command of running the 2D box case.

```
1. blockMesh # mesh generation
2. HydrothermalSinglePhaseDarcyFoam # execute the solver application
```

**Step 6: display results**

We can use paraFoam or ParaView (just like step 3) to display results which saved in time directories. The time directories name is dependent on key entries `writeInterval`, `purgeWrite` and `timePrecision`. The temperature is shown in figure **Fig. 5.4**.

![Temperature Result](image)

**Fig. 5.4: Temperature result of the hello world model.**

**Tip:** If users want to run a case for a long time, there is an excellent tool named tmux, which is a terminal multiplexer. It lets you switch easily between several programs in one terminal, detach them (they keep running in the background) and reattach them to a different terminal.

### 5.1.2 Nonuniform fixed temperature BC

All the input files can be found in `cookbooks/nonUniformFixedValueBC` directory.

This example, based on **Hello World: 2D box**, presents how to set a nonuniform fixed boundary condition by using `codedFixedValue BC` type (see **Listing 5.8**).
Listing 5.8: Example of nonuniform boundary condition: codedFixedValue.

```plaintext
bottom
{
    type        codedFixedValue;
    value       uniform 873.15; //placeholder
    name        gaussShapeT;
    code #{
        scalarField x(this->patch().Cf().component(0));
        double wGauss=200;
        double x0=1000;
        double Tmin=573;
        double Tmax=873.15;
        scalarField T(Tmin+(Tmax-Tmin)*exp(-(x-x0)*(x-x0)/
        (∕(2*wGauss*wGauss))));
        operator==T; #}
}
```

The model geometry and boundary conditions are shown in Fig. 5.5.

![Model geometry and boundary conditions](image)

Fig. 5.5: The geometry and boundary conditions of the Nonuniform fixed temperature BC model.

The temperature result is shown in Fig. 5.6.

### 5.1.3 Time-dependent permeability

All the input files can be found in `cookbooks/timeDependentPerm` directory.

The change of permeability over time, e.g. due to mineral precipitation, is an important process real hydrothermal systems. This example, based on *Hello World: 2D box*, presents how to set time-dependent
Fig. 5.6: Temperature result of the Nonuniform fixed temperature BC model.

permeability at run time in HydrothermalFoam tools. This can be reached using function sub-dictionary in controlDict file (see Listing 5.9).

Listing 5.9: Change permeability at run time: controlDict.

```plaintext
functions
{
    changePermeability
    {
        libs            ("libutilityFunctionObjects.so");
        type            coded;
        enabled         true;
        writeControl    runTime;
        writeInterval   86400000;
        name            changePermeability;
        codeWrite
        #{
            //1. Get simulation time
            double time = mesh().time().value();

            //2. Get modifiable pointer of a field variable by name
            volScalarField& perm_ = const_cast<volScalarField&>
            (mesh().lookupObject<volScalarField>("permeability");

            //3. Increase permeability from 1e-14 to 1e-13 after 500 years in region of y>2400 m,
            double year2sec = 86400*365;
            double t0 = 500*year2sec; // 500 years
            double wGauss = 100*year2sec; // 300 years
            double kmax = 14, kmin = 13;
            double y0 = -2400;
            if (time > t0)
```
The model geometry and boundary conditions are shown in Fig. 5.7, the permeability in shallow region (depth < 2.4 km) will be increase after 500 years (the inset curve).

Fig. 5.7: The geometry, BCs and permeability of the Time-dependent permeability model.

The temperature and permeability results are shown in Fig. 5.8.

5.1.4 Gmsh

All the input files can be found in cookbooks/gmsh directory.

This example is based on Hello World: 2D box, the only difference is the mesh generation. We use Gmsh to generate an unstructured triangular mesh. The gmsh script box.geo is shown in Listing 5.10.
Fig. 5.8: Temperature and permeability result of the *Time-dependent permeability* model.
Listing 5.10: Gmsh geometry script of the Gmsh model.

```
// 0. define some variables
xmin=0;
xmax=2000;
ymin=-3000;
ymax=-2000;
zmin=0;
zmax=10;
lc=20;

// 1. define points
Point(1) = {xmin, ymax, zmin, lc};
Point(2) = {xmax, ymax, zmin, lc};
Point(3) = {xmax, ymin, zmin, lc};
Point(4) = {xmin, ymin, zmin, lc};

// 2. define lines
Line(1) = {1, 2};
Line(2) = {2, 3};
Line(3) = {3, 4};
Line(4) = {4, 1};

// 3. define line loop and surface
Line Loop(6) = {4, 1, 2, 3};
Plane Surface(6) = {6};

// 3.1 make regular mesh
Transfinite Surface {6};
Recombine Surface {6};

// 4. extrude 2D surface to a 3D volume
Extrude {0, 0, zmax} {Surface(6); Layers(1); //set layer number to 1 for 2D model
Recombine;

// 5. define boundary patches via Physical keyword
Physical Surface("frontAndBack") = {28, 6};
Physical Surface("bottom") = {27};
Physical Surface("left") = {15};
Physical Surface("top") = {19};
Physical Surface("right") = {23};

// 6. specify a name for cell region which is used for 'setFields'
Physical Volume("internal") = {1};
```

Tip: The Gmsh can generate regular mesh as well, see lines 23-24 in Listing 5.10.

The model geometry, boundary conditions and mesh structure are shown in Fig. 5.9.
The temperature result is shown in Fig. 5.10

5.1. Convection in a box
Fig. 5.9: The geometry, boundary conditions and mesh structure of the *Gmsh* model.

Fig. 5.10: Temperature result of the *Gmsh* model.
5.1.5 3D box

All the input files can be found in `cookbooks/3Dbox` directory.

This example is based on models in Section 5.1.2 and Section 5.1.4. We just to make the following three changes to make a 3D box model,

1. For mesh generation, we just need to change \( z_{\text{max}} = x_{\text{max}} \); and change the extrude layers to a number greater than 1, e.g. 40, see line 7 and line 28 in Listing 5.10.

2. For boundary conditions, we just need to change `type` of `frontAndBack` boundary to be consistent with the `left` and `right` boundary, rather than empty. And then change `bottom` boundary condition of \( T \) similar to example Section 5.1.2, see Listing 5.11.

3. Unlike 2D model (e.g.), we don’t need to set `frontAndBack` to empty patch in `polyMesh/boundary` file. See Section 4.1.2 for empty boundary of 2D model.

Listing 5.11: Bottom boundary condition of \( T \) of the 3D box model.

```plaintext
bottom
{
    type codedFixedValue;
    value uniform 873.15; // placeholder
    name gaussShapeT;
    code #{
        scalarField x(this->patch().Cf().component(0));
        scalarField z(this->patch().Cf().component(2));
        double wGauss=200;
        double x0=1000;
        double z0=1000;
        double Tmin=573;
        double Tmax=873.15;
        scalarField T(Tmin+(Tmax-Tmin)*exp(-((x-x0)*(x-x0)+(z-z0)*(z-z0))/(2*wGauss*wGauss)));
        operator=(T);
    };
}
```

The mesh is shown in Fig. 5.11 and the bottom boundary condition of temperature is shown in Fig. 5.12. The isothermal surface of 300 °C, flow arrows and stream lines are shown below.

5.1.6 Parallel computing

All the input files can be found in `cookbooks/3Dbox_par` directory.

Parallel computing is one of features of OpenFOAM, and the HydrothermalFoam fully inherits this feature. There are two steps to run a case in Parallel,

1. Add `decomposeParDict` dictionary file into `system` directory (see Listing 4.14).
2. Decomposition of mesh and initial field data.
3. Replace `runApplication $application` with `runParallel $application`

4. Reconstructing mesh and data.

The last three steps can be assembled in `run.sh` which is highlighted in Listing 5.12

Listing 5.12: Command set of running a case in parallel.

```bash
#!/bin/sh

cd $(0%/*) || exit 1  # Run from this directory

# Source tutorial run functions
.
$WM_PROJECT_DIR/bin/tools/RunFunctions

application=`getApplication`
.
clean.sh

# generate mesh using gmsh

gmsh gmsh/box.geo -3 -o gmsh/box.msh -format msh22

# convert gmsh to OpenFOAM format

gmshToFoam gmsh/box.msh

# run solver in parallel

runApplication decomposePar
runParallel $application
```

(continues on next page)
Fig. 5.12: The bottom boundary condition of temperature of the 3D box model.

5.1. Convection in a box
Fig. 5.13: The isothermal surface of 300 °C = 573.15 K, flow arrows and stream lines of the 3D box simulation at 265 year. The unit of temperature in this figure is K.

```
runApplication reconstructPar
```

The mesh will be decomposed into multiple connected regions, the decomposed region number is defined in `decomposeParDict` file. The mesh of this example, which is based on 3D box model in Section 5.1.5, is decomposed into 4 regions by using `decomposePar` command in step 2.

The decomposed mesh of the 3D box model is shown in Fig. 5.14.

### 5.2 Pipe model

#### 5.2.1 Two-dimensional pipe

All the input files can be found in `cookbooks/pipe` directory.

The pipe model, illustrated in Fig. 5.15, could mimic a simplified scenario of hydrothermal circulation in oceanic crust.

There are five highlights in the pipe model.

1. To simulate a focused upflow zone in the deeper crust (e.g. layer 2B) where permeability is sufficiently low \( k = 10^{-15} \text{ m}^2 \) to allow for high-temperature fluid flow, a `hydrothermalMassFluxPressure` \( \phi = 1 \text{ g/m}^2/\text{s} \) boundary condition for pressure and
Fig. 5.14: The decomposed mesh of the 3D box.

Layer 2A, \( k = 4 \times 10^{-14} \text{ m}^2 \)

Layer 2B, \( k = 10^{-15} \text{ m}^2 \)

Side walls
p: noFlux
T: zeroGradient

Top, \( p = 300 \) bar
T: inletOutlet

Bottom, \( T = 400 \) °C
p: flux=1 g/m²/s

Fig. 5.15: Mesh and boundary conditions of the pipe model.

5.2. Pipe model
fixedValue \( (T_{\text{bot}} = 400 \, ^{\circ}\text{C}) \) boundary condition for temperature are applied on the pipe bottom.

2. Hybrid mesh, triangular mesh and regular mesh, is generated by \texttt{gmsh}. The mesh in the central zone is regular.

3. Different resolution of mesh is set by \texttt{gmsh}.

4. Permeability field is specified with different value for layer 2A and layer 2B by \texttt{zoneToCell} keyword in \texttt{setFieldDict} file.

5. To simulate hydrothermal flow discharge out of seafloor, a \texttt{inletOutlet} boundary condition of temperature is applied on the pipe top.

The temperature result is shown in Fig. 5.16.

![Fig. 5.16: Temperature result of the Two-dimensional pipe model.](image)

5.2.2 Three-dimensional pipe

All the input files can be found in \texttt{cookbooks/pipe_3D} directory.

The three-dimensional pipe model, illustrated in Fig. 5.17, is the result of a two-dimensional model (see Section 5.2.1) rotating around the central axis. Of course one can extrude mesh of Two-dimensional pipe model to get a three-dimensional pipe model (see Fig. 5.18).

\textbf{Note:} The boundary conditions setup is similar to the Two-dimensional pipe model. One can also run this example in parallel, see Section 5.1.6. The result is shown below.

The isothermal surface of 200 \(^{\circ}\text{C}\), flow arrows and stream lines are shown below.

5.3 Single pass model

5.3.1 Two-dimensional single pass model

All the input files can be found in \texttt{cookbooks/singlepass} directory.
Fig. 5.17: Mesh of the 3D pipe model.

Fig. 5.18: Alternative mesh of the 3D pipe model.
Fig. 5.19: The isothermal surface (red) of $200 \, ^\circ C = 473.15 \, K$, flow arrows and stream lines of the 3D pipe simulation at 80 year. The unit of temperature in this figure is K.
One- and two-limb single-pass models are usually used to determine vent field characteristics such as mass flow rate $Q$, bulk permeability in the discharge zone $k_d$, thickness of the conductive boundary layer at the base of the system $d$, magma replenishment rate, and residence time in the discharge zone [Lowell et al., 2013].

**One-limb classical single pass model**

A schematic of the one- and two-limb single-pass model are shown in Fig. 5.20.

![Fig. 5.20: Schematic of the one-limb (A) and two-limb (B) single-pass model (reproduced from [Lowell et al., 2013]).](image)

Here we present the one-limb single pass mode simulation using HydrothermalFoam, the model geometry, mesh and boundary conditions are shown in Fig. 5.21.

The model setup is similar to [Lowell et al., 2007]. The temperature evolution and streamlines of the **Two-dimensional single pass model** are in Fig. 5.22.

If we consider the recharge zone on the right side in single pass model shown in Fig. 5.21, the hydrothermal circulation pattern would be different, a schematic of the full single pass model is shown in

Here we present the full single pass mode simulation using HydrothermalFoam, the model geometry, mesh and boundary conditions are shown in Fig. 5.24.

The temperature evolution and streamlines of the full **Two-dimensional single pass model** are in Fig. 5.22.

**Two-limb single pass model**

All the input files can be found in cookbooks/singlepass_twolimb directory. The model geometry, mesh and boundary conditions are shown in Fig. 5.26.
Fig. 5.21: Model geometry, mesh and boundary conditions of the 2D single pass model.

Fig. 5.22: Temperature result of the Two-dimensional single pass model.
Fig. 5.23: Schematic of the full single-pass model (reproduced from [Lowell et al., 2014]).

Fig. 5.24: Model geometry, mesh and boundary conditions of the improved 2D single pass model.

Fig. 5.25: Temperature result of the Two-dimensional single pass model.
Fig. 5.26: Model geometry, mesh and boundary conditions of the two-dimensional two-limb single pass model.

The temperature evolution and streamlines of the two-limb single pass model are in Fig. 5.27.

Fig. 5.27: Temperature result of the Two-dimensional single pass model.

