



High-level programming of OpenFOAM applications, and a first glance at C++

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Prerequisites

- You are familiar with basic usage of OpenFOAM.
- You are familiar with the directory and file structure of OpenFOAM.
- You have some programming experience.
- You are aware of the wmake command, and how it picks up the instructions for the compilation.

Learning outcomes

- You will get a suggested way of working with your own developments of applications.
- You will understand the very basic parts of a C++/OpenFOAM code.
- You will understand the wmake compilation procedure for applications, and how it is related to compilation with the g++ compiler and make.
- You will step-by-step from scratch implement and understand the purpose of the most general high-level parts of OpenFOAM solvers.



High-level programming of OpenFOAM applications, and a first glance at C++

Contents

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- We will start by setting up a code from the very scratch, and make it compile using the OpenFOAM procedures.
- We will discuss the fundamental concepts of CFD, and piece-by-piece implement a code at the very highest level - meaning that we will at this point not know exactly what is happening at the lower levels, where the actual code is found. We do not have enough knowledge at the moment to look at the lower levels. That we will get later.

The instructions are inspired by the codes found in <code>\$FOAM_APP/test</code> and of course in <code>\$FOAM_APP</code>. See also the Programmers guide, linked to from:

https://www.openfoam.com/documentation/.



A C++ code must at least have one function, named main():

```
int main()
{
return 0;
}
```

The above means that:

- The function takes no arguments (nothing between the brackets)
- The function returns an integer (int)
- The function *definition* is between the curly brackets
- The returned integer is zero (return 0).
- The code will not give any output to the terminal window.



Let's implement and compile this using the OpenFOAM procedures.

• Create a directory for the code, and go there (inspired by the original test directory):

mkdir -p \$WM_PROJECT_USER_DIR/applications/myTests/onlyMainFunction
cd \$WM_PROJECT_USER_DIR/applications/myTests/onlyMainFunction

• Add the code in the previous slide to onlyMainFunction.C (named as directory)

At this point we can actually compile the code by

g++ onlyMainFunction.C -o Test-onlyMainFunction

Run the code by

./Test-onlyMainFunction

Alternatively, put it in \$FOAM_USER_APPBIN, as when compiling OpenFOAM applications g++ onlyMainFunction.C -o \$FOAM_USER_APPBIN/Test-onlyMainFunction Run the code by

Test-onlyMainFunction

Remember that the SPATH is used to find the executable, but . is not in the SPATH



Now we will compile it the OpenFOAM way:

• Clean up

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rm Test-onlyMainFunction \$FOAM_USER_APPBIN/Test-onlyMainFunction

• Create Make/files with

onlyMainFunction.C
EXE = \$(FOAM_USER_APPBIN)/Test-onlyMainFunction

• Create an empty Make/options file:

touch Make/options

• Compile:

wmake

• Run the code by:

Test-onlyMainFunction



Have a look at the output in the terminal window from the wmake command:

g++ -std=c++11 -m64 -DOPENFOAM_PLUS=1706 -Dlinux64 -DWM_ARCH_OPTION=64 -DWM_DP -DWM_LABEL_SIZE=32 -Wall -Wextra -Wold-style-cast -Wnon-virtual-dtor -Wno-unused-parameter -Wno-invalid-offsetof -O3 -DNoRepository -ftemplate-depth-100 -IlnInclude -I. -I/home/oscfd/OpenFOAM/OpenFOAM-plus/src/OpenFOAM/InInclude -I/home/oscfd/OpenFOAM/OpenFOAM-plus/src/OSspecific/POSIX/InInclude -fPIC -Xlinker --add-needed -Xlinker --no-as-needed Make/linux64GccDPInt32Opt/onlyMainFunction.o -L/home/oscfd/OpenFOAM/OpenFOAM-plus/platforms/linux64GccDPInt32Opt/lib \ -lOpenFOAM -ldl \

-lm -o \$FOAM_USER_APPBIN/Test-onlyMainFunction

The difference from when we compiled with g++ directly is that the wmake command added a lot of flags. They are in fact not needed for this code, but they also don't hurt.

From where did the flags come?



Solving PDEs with OpenFOAM

Let's start building our code by looking at the smallest elements of CFD.

- In CFD we solve partial differential equations (PDEs)
- The PDEs we wish to solve involve derivatives of tensor fields with respect to time and space
- The PDEs must be discretized in time and space before we solve them
- We will start by having a look at algebra of tensors in OpenFOAM at a single point
- We will then have a look at how to generate tensor fields from tensors, and how to relate those fields to a mesh
- Finally we will see how to discretize the PDEs and how to set boundary conditions using high-level coding in OpenFOAM
- For further details, see the ProgrammersGuide (found at https://www.openfoam.com/documentation/)



• OpenFOAM provides pre-defined classes for tensors of rank 0-3:

Rank	Common name	Basic name	Access function
0	Scalar	scalar	
1	Vector	vector	x(), y(), z()
2	Tensor	tensor	xx(), xy(), xz(),

A tensor
$$T = \begin{bmatrix} 11 & 12 & 13 \\ 21 & 22 & 23 \\ 31 & 32 & 33 \end{bmatrix}$$
 is defined line-by-line in OpenFOAM tensor T(11, 12, 13, 21, 22, 23, 31, 32, 33);

However, we can't just put this line in the main() function definition, since C++ does not know what a tensor is. You can try if you like!

We need to *declare* the tensor class and *link* to its definition, which is in the shared object file of the library to which the tensor class belongs.



We can find the implementation of the tensor class by

find \$FOAM_SRC -name tensor

which gives

\$FOAM_SRC/OpenFOAM/primitives/Tensor/tensor

We see that the tensor class belongs to the OpenFOAM library, since it is in the OpenFOAM directory we find the Make directory - showing that it is the top level of that library. In the OpenFOAM directory we also find the InInclude directory, in which there is a link to

\$FOAM_SRC/OpenFOAM/primitives/Tensor/tensor.H

For all declaration files in OpenFOAM you find both a link in an InInclude directory and the file itself: find \$FOAM_SRC -name tensor.H yields

```
$FOAM_SRC/OpenFOAM/lnInclude/tensor.H
$FOAM_SRC/OpenFOAM/primitives/Tensor/tensor/tensor.H
```

That file contains the *declaration* of the class, and we need to include that file at the header of our code. We also have to link to the <code>OpenFOAM</code> library, where the compiled definition of the class can be found.

Let's implement this...



mkdir -p \$WM_PROJECT_USER_DIR/applications/myTests/myTensor cd \$WM_PROJECT_USER_DIR/applications/myTests/myTensor

Add the following code to myTensor.C

```
#include "tensor.H"
using namespace Foam;
int main()
{
  tensor T( 11, 12, 13, 21, 22, 23, 31, 32, 33);
Info << "T: " << T << endl;
Info << "Txz: " << T.xz() << endl;
return 0;
}</pre>
```

- The first line includes the declaration of the tensor class, so that it can be used below.
- The second line says that we should use the namespace Foam, in which everything in Open-FOAM is implemented. We will get back to this later.
- The OpenFOAM Info output stream allows us to write to the terminal window, which we will get back to later. The second Info line uses a call to the access function xz(), which belongs to the tensor class. This is how you typically call functions of a class.



Now we need to set up the Make directory so that wmake finds the tensor. H file and links to the OpenFOAM library.

 $Create \; \texttt{Make/files with}$

myTensor.C
EXE = \$(FOAM_USER_APPBIN)/Test-myTensor

Create an empty Make/options file:

touch Make/options

Compile:

wmake

Run the code by:

Test-myTensor

MAGIC? How does it find tensor.H and the OpenFOAM library? Have a look again at the default wmake flags, showing that the compilation is now done in two steps. The first step is for compilation of the intermediate object file, and the second step is for the linking. In the first step the flag -I/home/oscfd/OpenFOAM/OpenFOAM-plus/src/OpenFOAM/lnInclude tells wmake where to find tensor.H. In the second step the flag

-L\$WM_PROJECT_DIR/platforms/linux64GccDPInt32Opt/lib tells wmake where the libraries can be found, and the flag -lOpenFOAM tells wmake to link to that library. NO MAGIC!

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Algebraic tensor operations in OpenFOAM

- Tensor operations operate on the entire tensor entity instead of a series of operations on its components
- The OpenFOAM syntax closely mimics the syntax used in written mathematics, using descriptive functions or symbolic operators

	Operation	Comment	Mathematical	Description		
			description	in OpenFOAM		
	Addition		a + b	a + b		
	Outer product	Rank $\mathbf{a}, \mathbf{b} \ge 1$	ab	a * b		
	Inner product	Rank $\mathbf{a}, \mathbf{b} \ge 1$	a · b	a & b		
Framplace	Cross product	Rank $\mathbf{a}, \mathbf{b} = 1$	$\mathbf{a} \times \mathbf{b}$	a^b		
Examples:	Operations exclusive to tensors of rank 2					
	Transpose		\mathbf{T}^{T}	T.T()		
	Determinant		$\det \mathbf{T}$	det(T)		
	Operations exclusive to scalars					
	Positive (boolean)		$\mathbf{s} \ge 0$	pos(s)		
	Hyperbolic arc sine		asinh s	asinh(s)		

Find more examples in the Programmer's guide

(linked to at https://www.openfoam.com/documentation/)



Algebraic tensor operations in OpenFOAM

Add the following to our myTensor.C file (before return), compile and run:

```
// Line to remove copy-paste problem
tensor t1(1, 2, 3, 4, 5, 6, 7, 8, 9);
tensor t2(1, 2, 3, 1, 2, 3, 1, 2, 3);
tensor t3 = t1 + t2;
Info<< t3 << endl;
tensor t4(3,-2,1,-2,2,0,1, 0, 4);
Info<< inv(t4) << endl;
Info<< (inv(t4) & t4) << endl;
Info<< t1.x() << t1.y() << t1.z() << endl;
Info<< t1.T() << endl;
Info<< det(t1) << endl;
Info<< det(t1) << endl;
Info<< pos(s1) << endl;
Info<< Foam::asinh(s1) << endl;</pre>
```

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We do not have to add an include line for scalar, since that is done through the tensor.H file (through Tensor.H -> Vector.H -> VectorSpace.H).

The Foam:: at the final line is because the compiler complained that the function is ambiguous, i.e. it exists in more than one namespace. Therefore we explicitly say that we want to use the function that belongs to namespace Foam. It has to be done without using namespace Foam.

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A quick note on the tensor class

This slide is actually a bit early since it requires a bit more knowledge, but you can see it as a teaser... It shows how you can learn what a tensor is and how you can learn more on its usage.

- In tensor.H, Tensor.H is included (located in \$FOAM_SRC/OpenFOAM/primitives/Tensor), which declares the access functions and operators and includes TensorI.H, where they are defined. The capital T means that it is a template class. The capital I means that the file contains inline functions.
- See also vector, symmTensorField, sphericalTensorField and many other examples.



Dimensional units in OpenFOAM

• OpenFOAM checks the dimensional consistency

Declaration of a tensor with dimensions:

```
dimensionedTensor sigma
  (
    "sigma",
    dimensionSet( 1, -1, -2, 0, 0, 0, 0),
    tensor( 1e6, 0, 0, 0, 1e6, 0, 0, 0, 1e6)
  );
```

The values of dimensionSet correspond to the powers of each SI unit:

No.	Property	Unit	Symbol		
1	Mass	kilogram	kg		
2	Length	metre	m		
3	Time	second	S		
4	Temperature	Kelvin	Κ		
5	Quantity	moles	mol		
6	Current	ampere	А		
7	Luminous intensity	candela	cd		
views than has the dimension $[ka/ma^2]$					

sigma then has the dimension $\left[kg/ms^2\right]$



Dimensional units in OpenFOAM

 Add the following to myTensor.C: Before main(): #include "dimensionedTensor.H" Before return(0):

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```
dimensionedTensor sigma
(
    "sigma",
    dimensionSet( 1, -1, -2, 0, 0, 0, 0),
    tensor( 1e6, 0, 0, 0, 1e6, 0, 0, 0, 1e6)
);
Info<< "Sigma: " << sigma << endl;</pre>
```

• Compile, run again, and you will get:

```
Sigma: sigma [1 -1 -2 0 0 0] (1e+06 0 0 0 1e+06 0 0 0 1e+06)
```

You see that the object sigma belongs to the dimensionedTensor class that contains both the name, the dimensions and values.

• See \$FOAM_SRC/OpenFOAM/dimensionedTypes/dimensionedTensor



Dimensional units in OpenFOAM

• Try some member functions of the dimensionedTensor class:

```
Info<< "Sigma name: " << sigma.name() << endl;
Info<< "Sigma dimensions: " << sigma.dimensions() << endl;
Info<< "Sigma value: " << sigma.value() << endl;</pre>
```

• You now also get:

Sigma name: sigma Sigma dimensions: [1 -1 -2 0 0 0 0] Sigma value: (1e+06 0 0 0 1e+06 0 0 0 1e+06)

• Extract one of the values:

Info<< "Sigma yy value: " << sigma.value().yy() << endl; Note here that the value() member function first converts the expression to a tensor, which has a yy() member function. The dimensionedTensor class does not have a yy() member function, so it is not possible to do sigma.yy().

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Construction of a tensor field in OpenFOAM

- A tensor field is a list of tensors
- The use of typedef in OpenFOAM yields readable type definitions: scalarField, vectorField, tensorField, symmTensorField, ...
- Algebraic operations can be performed between different fields, and between a field and a single tensor, e.g. Field U, scalar 2.0: U = 2.0 * U;
- Add the following to myTensor.C: Before main(): #include "tensorField.H" Before return(0):

```
tensorField tf1(2, tensor::one);
Info<< "tf1: " << tf1 << endl;
tf1[0] = tensor(1, 2, 3, 4, 5, 6, 7, 8, 9);
Info<< "tf1: " << tf1 << endl;
Info<< "2.0*tf1: " << 2.0*tf1 << endl;</pre>
```

• However, this kind of tensor field is not related to any mesh, and can therefore not be discretized.



#include "fvCFD.H"

- OpenFOAM uses the finite volume method (fvm) to discretize the PDEs, and there are many classes in OpenFOAM that are related to fvm.
- OpenFOAM provides the header file fvCFD.H that only includes other header files related to fvm, including the tensor classes we have discussed. It can therefore be used to reduce the number of header files. It in fact also ends with using namespace Foam.
- Exchange *all* the lines before the main() function with:

```
#include "fvCFD.H"
```

• That file is located in a sub-directory of \$FOAM_SRC/finiteVolume, so add the following to Make/options (Note that we do not have to link. Why?):

```
EXE_INC = \
    -I$(LIB_SRC)/finiteVolume/lnInclude \
    -I$(LIB_SRC)/meshTools/lnInclude
```

The meshTools line is needed for a cyclicAMILduInterface.H file that is included through fvCFD.H, and obviously includes something from the meshTools library.

• Compile and test!



Discretization of a tensor field in OpenFOAM

- OpenFOAM uses the finite volume method (fvm) to discretize the PDEs of the tensor fields on a mesh.
- The tensor fields must thus be related to a mesh.

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- The polyMesh class can be used to construct a polyhedral mesh using the minimum information required. I.e. that in <case>/constant/polyMesh
- The fvMesh class extends the polyMesh class to include additional data needed for the fvm discretization (see test/mesh)
- The geometricField class relates a tensor field to an fvMesh (can also be typedef volField, surfaceField, pointField)
- A geometricField inherits all the tensor algebra of its corresponding field, has dimension checking, and can be subjected to specific discretization procedures

We will now investigate a polyMesh, an fvMesh, and a geometricField, but for that we need a base code...



Base code

- For the base code we are inspired both by \$FOAM_APP/test/mesh/Test-mesh.C and \$FOAM_SOLVERS/incompressible/icoFoam/icoFoam.C
- We need a code that starts with (explanations coming later):

```
<header files>
int main(int argc, char *argv[])
{
    #include "setRootCase.H"
    #include "createTime.H"
```

where the <header files> can preferrably be fvCFD.H.

• We use a special functionality to generate a template application:

```
cd $WM_PROJECT_USER_DIR/applications/myTests
foamNewApp meshAndField
cd meshAndField
wmake
meshAndField
```

It complains that it can't find a system/controlDict. We will see why soon.

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Base code

The base code (meshAndField.C) contains a commented header, and:

#include "fvCFD.H"

Some explanations in the coming slides...



Base code

- int argc, char *argv[] The arguments to the main function are the number of flags and the flags supplied when running the code.
- setRootCase.H (find it with find \$FOAM_SRC -name setRootCase.H) Not a header file, just a piece of code:

```
Foam::argList args(argc, argv);
if (!args.checkRootCase())
{
    Foam::FatalError.exit();
}
```

This constructs the object args, and uses it to check that we are running the code in a case directory. It simply checks if there is an appropriate system/controlDict(!)

• createTime.H (find it with find \$FOAM_SRC -name createTime.H) Not a header file, just a piece of code:

```
Foam::Info<< "Create time\n" << Foam::endl;
Foam::Time runTime(Foam::Time::controlDictName, args);</pre>
```

This writes some text and constructs the runTime object of the class Time.



Base case

We need a case to run our code (here the cavity case with only four cells):

cp -r \$FOAM_TUTORIALS/incompressible/icoFoam/cavity/cavity \$FOAM_RUN/cavityFourCells
sed -i s/"20 20 1"/"2 2 1"/g \$FOAM_RUN/cavityFourCells/system/blockMeshDict
blockMesh -case \$FOAM_RUN/cavityFourCells
meshAndField -case \$FOAM_RUN/cavityFourCells

Now we can move on to examine a polyMesh and an fvMesh, following by a geometricField, in the form of a volScalarField...



Examine a polyMesh

Add after the line with #include "createTime.H" (from test/mesh):

```
Info<< "Create mesh" << endl;</pre>
polyMesh mesh
    IOobject
        fvMesh::defaultRegion,
         runTime.timeName(),
        runTime,
        IOobject::MUST READ
);
Info<< "Cell centres" << nl << mesh.cellCentres() << endl;</pre>
Info<< "Cell volumes" << nl << mesh.cellVolumes() << endl;</pre>
Info<< "Cell face centres" << nl << mesh.faceCentres() << endl;</pre>
```

Compile and run. Some descriptions follow...



Examine a polyMesh

 \bullet polyMesh mesh (...)

An object named mesh is constructed from the polyMesh class. There is only one argument to the constructor, which is an IOobject. The IOobject itself takes four arguments that are not obvious. We can make the educated guess that it reads the mesh from constant/polyMesh. At some point we have to stop figuring out exactly how things are done. Now we are fine relying on how OpenFOAM reads from files.

- cellCentres() gives the center of all cells and boundary faces.
- cellVolumes() gives the volume of all the cells.
- faceCentres() gives the center of all the faces.

A detailed note is that the above functions are actually inherited from the primitiveMesh class.

However, a polyMesh only has the very basic information of the mesh...



Examine an fvMesh

An fvMesh builds on top of a polyMesh, and has additional attributes.

- Let us examine an fvMesh: Start by changing the mesh class from polyMesh to fvMesh in meshAndField.C sed -i s/'polyMesh mesh'/'fvMesh mesh'/g meshAndField.C
- Compile and run.
- Note that we are using the same member functions as before. They still belong to the polyMesh class. However, since the fvMesh class inherits from the polyMesh class they also belong to the fvMesh class. We will discuss inheritance more later.
- Add the following after our previous insertion in meshAndField.C, compile and run again:

Info<< mesh.C() << endl; Info<< mesh.V() << endl; Info<< mesh.Cf() << endl;</pre>

This could not be done with the polyMesh, since those member functions are defined in the sub-class fvMesh. They give the corresponding information as before, but with additional information.



Examine an fvMesh

The construction of the fvMesh is done the same way for all solvers that use a static mesh. Instead of replicating the call for the fvMesh constructor in all the solvers they just have after #include "createTime.H" the line (see e.g. icoFoam.C)

#include "createMesh.H"

This is as well not a real header file, but just inserts the code in: find \$FOAM_SRC -name createMesh.H Clean up your code by doing this, and check that it still works!



Examine a volScalarField

Now we are at a point where we can start implementing a particular PDE solver. Depending on the problem at hand we want to solve equations for different fields. I.e. the fields to be constructed depends on the solver. That is why all the solvers include a local file named createFields.H. See e.g. the original icoFoam directory:

```
createFields.H icoFoam.C Make
```

Inside the main(...) function of icoFoam.C we find:

```
#include "createFields.H"
```

This refers to the local file in the same directory, which is found by the compiler thanks to the compiler flag -I.

Add such a line below #include "createMesh.H", create the file:

touch createFields.H

and we will continue constructing a field in that file.



Examine a volScalarField

We are here inspired by icoFoam/createFields.H, and construct a volScalarField (which is typedef of a geometricField, and therefore relates to a mesh).

Add in createFields.H:

```
volScalarField p
(
    IOobject
    (
        "p",
        runTime.timeName(),
        mesh,
        IOobject::MUST_READ,
        IOobject::AUTO_WRITE
    ),
        mesh
);
Info<< p << endl;
Info<< p.boundaryField()[0] << endl;</pre>
```

Compile, run and have a look at the output! Note that the above Info statement is in createFields.H, which is in the beginning of the execution.

Although we are not interested in all the details here we will still have a quick look...



Examine a volScalarField

volScalarField p(IOobject (...) , mesh);

The object p is constructed as a volScalarField, and the constructor takes two arguments.

- IOobject ("p",runTime.timeName(),mesh,IOobject::MUST_READ,IOobject::AUTO_WRITE)
 - The first argument is an internal name p, so that we at the end can ask our object for its name (i.e. the internal name should be the same as the object name).
 - The second argument uses the runTime object to determine which time directory to read from (according to settings in system/controlDict).
 - The third argument states that the volScalarField should be related to the fvMesh corresponding to our object mesh.
 - The fourth argument states that it must be read.
 - The fifth argument states that the field should be written in the coming time directories according to the settings in <code>system/controlDict</code>.
- mesh

If the volScalarField is read from a file, as in this case, you just specify mesh here. I haven't been interested in checking the details of this, but I guess it is used to specify the size and structure of the field. If the volScalarField is NOT read from a file the second argument can be an existing volScalarField that corresponds to the same mesh.