

# The Flavours of OpenFOAM®

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# Outline

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

1 What is OpenFOAM?

2 Flavours of OpenFOAM

3 OpenFOAM directory structure

4 OpenFOAM and HPC

5 OpenFOAM Architecture

6 OpenFOAM examples

7 OpenFOAM learning resources

8 OpenFOAM Case structure



# What is OpenFOAM®?

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- OpenFOAM® is a general purpose Computational Mechanics (CM) suite of libraries, solvers, and utilities.
- OpenFOAM® is a registered trademark of the ESI Group.
- OpenFOAM is licensed under the GNU General Public License (GPL).
- OpenFOAM is completely free of charge and can be used in academic and commercial applications.
- Unlike commercial codes, OpenFOAM is not just a monolithic application - it is a general-purpose library that can be used to create a CM solver tuned for a particular problem.



# What is OpenFOAM®?

- FOAM is an acronym that stands for **Field Operation And Manipulation**.
- In simple terms, OpenFOAM is a library of tools that provide for the manipulation of fields.
- Examples of fields might include temperature, velocity, pressure, density, magnetic, electric fields and so on.
- Fields can be of scalar, vector or tensor types.
- The library provides a consistent set of tools to perform algebraic and differential operations on fields.
- Library is implemented in the C++ language using object orientation, inheritance, templates and programming patterns.
- The code is fully parallel and does not impose any restrictions on the number of parallel processes.

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

# What is OpenFOAM®?

- In addition to field operations, the library implements additional tools required for the solution of problems:
  - Mesh conversion and generation tools
  - Parallel communication
  - Mapping utilities for multi-physics simulations
  - Comprehensive linear algebra suite for the solution of large linear systems
- All tools are designed to work together and exchange information in order to create solvers for complex problems.
- The publicly available versions of OpenFOAM is not a polished commercial product, and some work is required if one steps out of the pre-defined examples.



Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

# Why use OpenFOAM?

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

There is one main reason.

- Freedom of open source.

The freedom that open source provides drives two motivations for OpenFOAM usage:

- Free as in "free food". You can download and install OpenFOAM without paying any money.
- Free as in "free speech". With OpenFOAM, you have the following freedoms
  - use it for any purpose,
  - modify it to suit your needs,
  - share it with anyone,
  - share your changes.



# History of OpenFOAM

- FOAM was originally developed in the 1990s at Imperial College by Henry Weller, Hrvoje Jasak and others.
- FOAM sold as commercial software 2000-2004 by Nabla Ltd. Nabla Ltd. wound up in 2006.
- FOAM was released into the public domain in 2004 as OpenFOAM.
- Henry Weller and others founded OpenCFD Ltd (2004) and later CFD Direct Ltd (2015).
- Hrvoje Jasak founded Wikki Ltd (2004).
- OpenCFD Ltd sold to SGI in 2011, and OpenFOAM Foundation was formed.
- OpenCFD Ltd sold to ESI Group in 2012.



Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

# History of OpenFOAM

## Flavours of OpenFOAM®

Background  
What?  
Why?  
History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of Models

Examples

Examples

Learning resources

Case

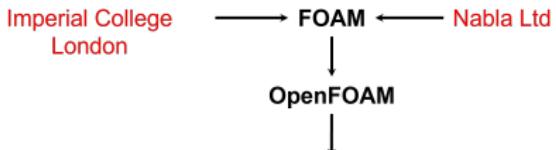
Structure

Dictionaries

system

constant

Time



## Main Flavours

OpenFOAM vX  
[www.openfoam.org](http://www.openfoam.org)

OpenFOAM Foundation  
(Copyright owners)

CFD Direct Ltd

OpenFOAM v202X  
[www.openfoam.com](http://www.openfoam.com)

OpenCFD Ltd/ESI  
(Trademark owners)

Community

foam-extend  
[foam-extend.org](http://foam-extend.org)

Wikki Ltd

Community



# Implemented capability - all flavours

- Finite Volume Method,
- Collocated polyhedral unstructured meshes,
- Second order in space and time,
- Steady and transient solvers with pressure velocity coupling via segregated methods,
- Lagrangian particle tracking,
- Automatic mesh motion with support for topological changes,
- Massive parallelism in domain decomposition mode,
- All components implemented in library form for easy re-use



Flavours of  
OpenFOAM®

Background  
What?

Why?  
History

Flavours

Directory

HPC

Architecture

OO Concepts  
Main Concepts

Equation mimicking  
Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionarys

system

constant

Time

# Implemented capability - all flavours

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of

Models

Examples

Examples

Learning

resources

Case

Structure

Dictionaries

system

constant

Time

## Physical and multi-physics modelling:

- Turbulence modelling, RANS, DEs, LES,
- Multiphase flows and mass transfer,
- Combustion and chemical reactions,
- Rheology models,
- Thermophysical models for liquids and gases.
- Conjugate heat transfer
- Stress analysis and fluid-structure interaction
- 6 DoF
- Aero-acoustics



# Flavour differences

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

Some of the main feature differences:

- Foundation (OpenFOAM vX)
  - Modular framework
- ESI (OpenFOAM v202X)
  - Overset
  - Community add-ons, cfMesh, OpenQBMM, Petsc4Foam
- foam-extend
  - Coupled framework and solvers
  - Immersed boundary method
  - Finite area method (now ported to ESI version)
  - Overset
  - Proper Orthogonal Decomposition



# Directory Structure

The environment variable \$WM\_PROJECT\_DIR holds the OpenFOAM installation path.

```
$WM_PROJECT_DIR
├── Allwmake
├── applications ...solvers/utilities sources.
├── bin
├── build
├── CONTRIBUTORS.md
├── COPYING
├── doc
├── etc
├── META-INFO
├── modules
├── platforms ...compiled code.
├── README.md
├── site
├── src ...libraries sources.
└── tutorials ...Tutorials cases.
    └── wmake
```

# Solvers/Utilities

- OpenFOAM is not a single executable.
- Depending on what you want to do, you will need to use a specific application.
- Solvers can be found in the  
    \$WM\_PROJECT\_DIR/applications/solvers directory  
    (\$FOAM\_SOLVERS)
- Utilities can be found in the  
    \$WM\_PROJECT\_DIR/applications/utilities directory  
    (\$FOAM\_UTILITIES)
- If you want to get help on how to run an application,  
    type in terminal

```
$> application_name -help
```



# Solvers

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- Solvers are subdivided according to the physics they address.
- The solvers available in the OpenFOAM installation (version 2212):

- acoustics
- basic
- combustion
- compressible
- discreteMethods
- DNS
- electromagnetics
- financial
- finiteArea
- heatTransfer
- incompressible
- lagrangian
- multiphase
- stressAnalysis

- Each sub-directory may contain several sub-directories, one for each solver.
- Each solver directory contains a \*.C file with the same name as the directory. This file contains a short description of the solver.

# Solvers

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of

Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionarys

system

constant

Time

- For example, the *incompressible* sub-directory contains the following solvers:
  - adjointOptimisationFoam
  - adjointShapeOptimizationFoam
  - boundaryFoam
  - icoFoam
  - nonNewtonianIcoFoam
  - pimpleFoam
  - pisoFoam
  - shallowWaterFoam
  - simpleFoam

# Solvers

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

Commonly used OpenFOAM solvers and their supported physics.

Physics	bPF	cMRF	iF	pF	rPF	rPF	sF
transient	✓	✓	✓	✓	✓		
compressible	✓	✓			✓	✓	
turbulence	✓	✓	✓	✓	✓	✓	✓
heat-transfer	✓	✓			✓	✓	
buoyancy	✓	✓			✓		
combustion					✓		
multiphase				✓			
particles						✓	
dynamic mesh				✓	✓		✓
multi-region		✓					

- bPF - buoyantPimpleFoam
- cMRF - chtMultiRegionFoam
- iF - interFoam
- pF - pimpleFoam
- rPF - reactingParcelFoam
- rPF - rhoPimpleFoam
- sF - simpleFoam

# OpenFOAM and HPC

Flavours of OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of Models

Examples

Examples

Learning resources

Case

Structure

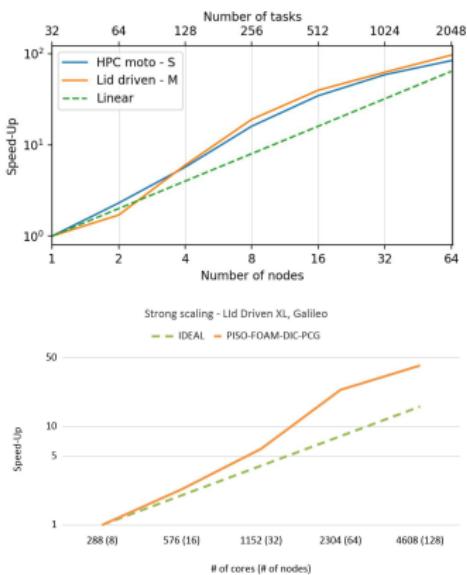
Dictionaries

system

constant

Time

OpenFOAM has been shown to scale to in excess of 4000 cores<sup>1</sup>

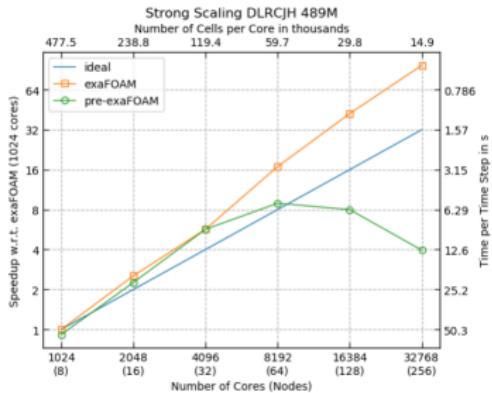


<sup>1</sup>[https://wiki.openfoam.com/images/c/c1/OpenFOAM\\_2020\\_CINECA\\_Spisso.pdf](https://wiki.openfoam.com/images/c/c1/OpenFOAM_2020_CINECA_Spisso.pdf)

# OpenFOAM and HPC

## exaFOAM

- Consortium (12 Partners) led by ESI-OpenCFD consisting of experts to work on the co-design of OpenFOAM targeting (pre)-exascale HPC architectures.
- Grant Funded by EuroHPC-03-2019 totalling €5,425,618
- Outcomes implemented in OpenFOAM V202X, example shown below<sup>2</sup>



<sup>2</sup>[https://exafoam.eu/wp-content/uploads/2023/07/exaFOAM\\_Workshop\\_GrandChallenges.pdf](https://exafoam.eu/wp-content/uploads/2023/07/exaFOAM_Workshop_GrandChallenges.pdf)

# Object Orientation in OpenFOAM

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

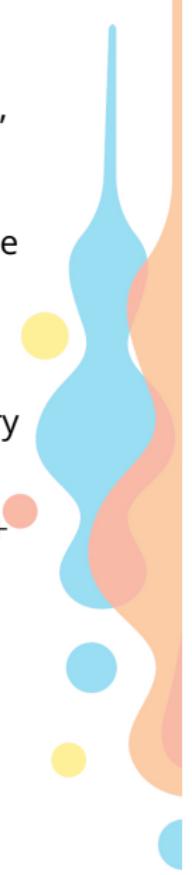
Dictionaries

system

constant

Time

- OpenFOAM is assembled from components:
  - Core libraries containing discretization, mesh handling, etc.
  - Physical modelling libraries such as thermo-physical models, turbulence models, chemical reactions, particle collision models, etc.
  - Utilities for mesh import and manipulation, parallel processing, etc.
  - Solvers implemented as a top-level code utilising library components.
- Top-level code consists of a few hundred lines of C++ code.
- In essence, top-level code defines a computational programming language.



# Object Orientation in OpenFOAM

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

## Object-Oriented analysis

- Main objects are operators, e.g. time derivative, divergence, curl, Laplacian, etc.
- Operators act on fields, and depending on the desired result of the operation in OpenFOAM, we distinguish two implementation concepts:
  - Field (explicit evaluation of the operator acting on a field)
  - Matrix (implicit evaluation of operator acting on a field)
- We also need a few more components to represent equations in OpenFOAM:
  - Computational domain (space-time representation)
  - Basic containers (scalars, vectors, and tensors)
  - Fields and their associated algebra
  - Matrices for sparse linear systems
  - Discretization methods for operators



# Computational Domain / Field Algebra

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- The computational domain is represented through space-time domain implementation:
  - Mesh primitives: points, faces, cells → point, face, cell classes
  - Space: computational mesh → polyMesh class
  - Time: time steps (database) → time class
- Field algebra and containers:
  - Tensor: list of numbers + algebra → scalar, vector, tensor classes
  - Boundary condition: list of values + condition → patchField class
  - Physical dimensions: dimension set → dimensionSet class
  - Geometric field: field + mesh + boundary conditions → geometricField class
  - Field algebra: +, -, \*, /, sin(), exp() → field operators (C++ operator overloading)

# Linear Equation Solvers / Discretization methods

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- Linear equation systems and linear solvers:
  - Linear equation matrix: table of matrix coefficients → `lduMatrix` class
  - Solvers: iterative solvers → `lduMatrix::solver`
- Discretization methods:
  - Interpolation: differencing schemes → `interpolation` class.
  - Differentiation: `ddt`, `div`, `grad`, etc. → `fvc`, `fec` classes.
  - Discretization: `ddt`, `div`, Laplacian, etc. → `fvm`, `fam` classes.
- Top-level code:
  - Model library: library → `turbulenceModel` class for example.
  - Application: `main()` → `pisoFoam` solver for example

# Equation mimicking

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- Uses equation mimicking
- Consider PDE for the transport of turbulent kinetic energy:

$$\frac{\partial k}{\partial t} + \nabla \cdot (\mathbf{u} k) - \nabla \cdot [(\nu + \nu_t) \nabla k] = \nu_t \left[ \frac{1}{2} (\nabla \mathbf{u} + \nabla \mathbf{u}^T) \right]^2 - \frac{\varepsilon_0}{k_0} k$$

- Implementation:

solve

(

```
fvm::ddt(k)
+ fvm::div(phi,k)
- fvm::laplacian(nu() + nut, k)
== nut*magSqr(symm(fvc::grad(U)))
- fvm::Sp(epsilon/k, k)
```

) ;

# Model implementation

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- Model-to-model interaction is handled through common interfaces.
- New components do not disturb existing code, resulting in fewer bugs.
- Run-time selection tables for dynamic binding of new functionality.
- This approach is used for every implementation:
  - Differencing schemes
  - Gradient calculations
  - Boundary conditions
  - Linear equation solvers
  - Physical models
  - Mesh motion algorithms

The key aspect of model implementation is class inheritance and factory mechanism.



# Model implementation

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- Consider the turbulence model:

```
class turbulenceModel
{
    virtual volTensorField R() const = 0;
    virtual fvVectorMatrix divR
    (
        volVectorField& U
    ) const = 0;
    virtual void correct() = 0;
};
```

- New turbulence model implementation  
`funkyTurbulence`:

```
class funkyTurbulence :
    public turbulenceModel {};
```

# Model implementation

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- Main code sees only virtual interface:

```
autoPtr<turbulenceModel> turbulence
(
    turbulenceModel::New(U, phi, laminarTransport)
);
```

```
fvVectorMatrix UEqn
(
    fvm::ddt(rho, U)
    + fvm::div(phi, U)
    + turbulence->divR(U)
    ==
    - fvc::grad(p)
);
```

- Turbulence model selected by user at run-time.

# Laplacian Scalar Equation

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- PDE for Laplacian scalar equation:

$$\frac{\partial T}{\partial t} - \nabla^2(\mathcal{D}_T T) = 0$$

- Code:

```
for(int nonOrth=0;  
     nonOrth<=nNonOrthCorr;  
     nonOrth++)  
{  
    solve  
    (  
        fvm::ddt (T) - fvm::laplacian(DT, T)  
    );  
}
```

# Scalar Transport Equation

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of

Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- PDE for scalar transport:

$$\frac{\partial T}{\partial t} + \nabla \cdot (\mathbf{U} T) - \nabla^2 (\mathcal{D}_T T) = 0$$

- Code:

```
solve
(
    fvm::ddt (T)
    + fvm::div(phi, T)
    - fvm::laplacian(DT, T)
);
```

# Examples

- The following slides are only a part of what is possible with OpenFOAM
- Chosen for (personal) interest and illustration of the range of capabilities rather than an exhaustive illustration of the range of capabilities
- All results shown are from simulations conducted either by Laminar2 Turbulent or Applied CCM.



Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

## Examples - Wind engineering

Flavours of  
OpenFOAM®

## Background

What?

Why?

History

## Flavours

## Directory

HPC

Architecture

## Main Concepts

Equation mimicking

## Models

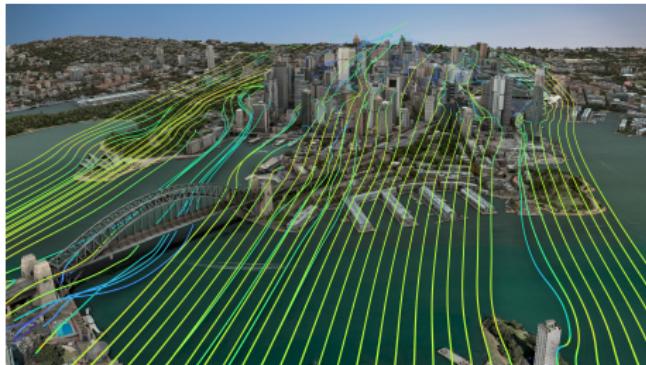
## Examples

## Learning resources

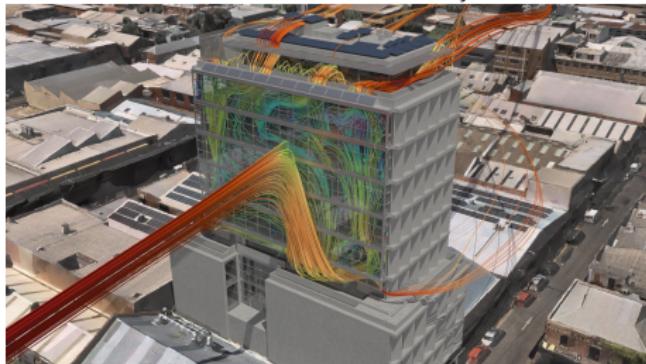
Case

Structure

constant



## Pedestrian comfort and safety



## Solar-heated vented facade simulation

# Examples - External aerodynamics

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of

Models

Examples

Examples

Learning  
resources

Case

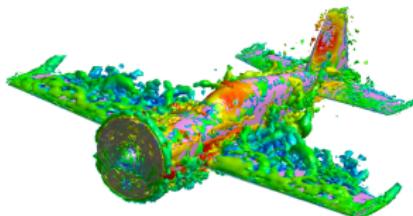
Structure

Dictionaries

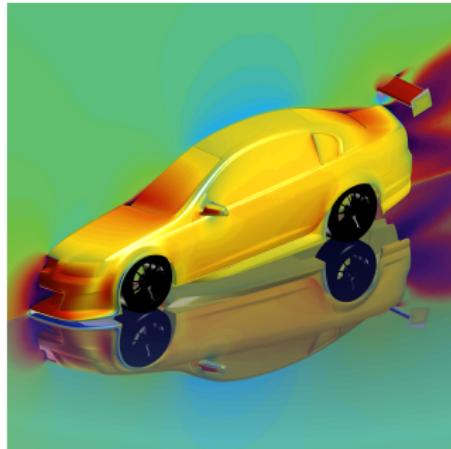
system

constant

Time



LES simulation



Steady-state RANS

# Examples - Marine hydrodynamics

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

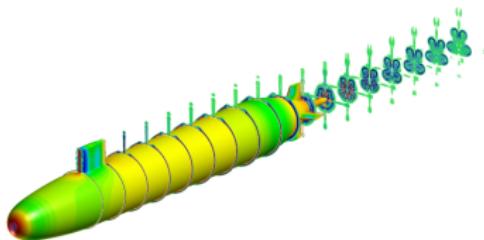
Structure

Dictionaries

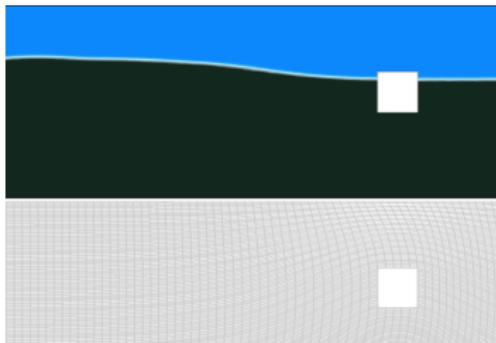
system

constant

Time



Steady ship resistance and propulsion



Unsteady 6-DoF motions with wave excitation

## Examples - Turbomachinery

Flavours of  
OpenFOAM®

## Background

What?

Why?

History

## Flavours

## Directory

HPC

Architecture

Main Concepts

### Equation mimicking

Implementation Models

### Examples

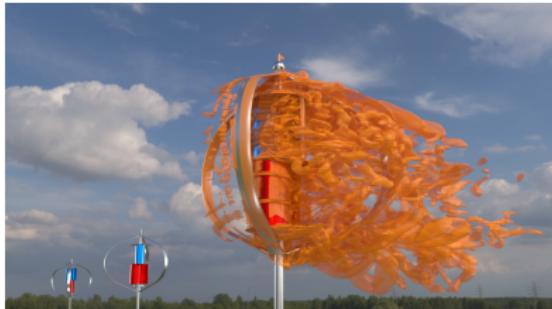
## Examples

## Learning resources

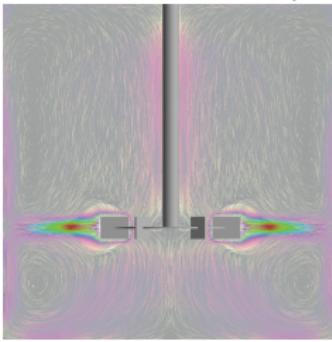
Case

## Dictionaries

system



Vertical axis wind turbine development



## Ruston turbine in a mixed tank



# Examples - Multiphase

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

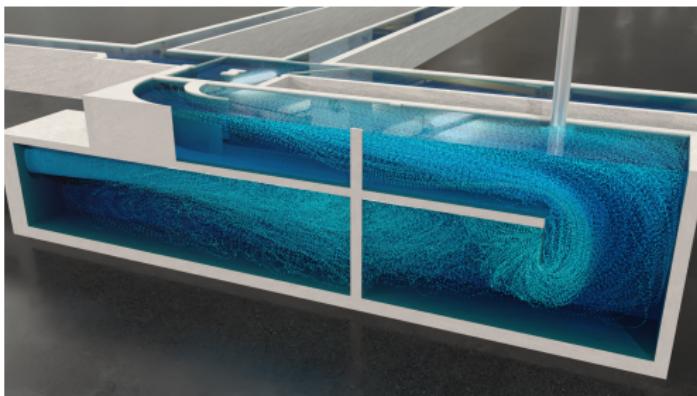
Structure

Dictionaries

system

constant

Time



Ozone contactor tank, multiphase flow mixing



Cyclone separator

# OpenFOAM learning resources

## Flavours of OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of Models

Examples

Examples

Learning resources

Case

Structure

Dictionaries

system

constant

Time

## Official sources:

- User guide:

<https://www.openfoam.com/documentation/user-guide>

- Tutorial guide:

<https://www.openfoam.com/documentation/tutorial-guide>

- Extended Code guide (new in 2023):

<https://doc.openfoam.com/2306/>

- Tutorial Wiki: <https://wiki.openfoam.com/Tutorials>



# OpenFOAM learning resources

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

## Unofficial sources:

- OpenFoam Wiki:  
[https://openfoamwiki.net/index.php/Main\\_Page](https://openfoamwiki.net/index.php/Main_Page)
- Chalmers University CFD with OpenSource Software course:  
[https://www.tfd.chalmers.se/~hani/kurser/OS\\_CFD/](https://www.tfd.chalmers.se/~hani/kurser/OS_CFD/)
- CFD Online forums:  
<https://www.cfd-online.com/Forums/openfoam/>
- The OpenFOAM Technology Primer book by Mooney, Maric and Höpken:  
[https://www.researchgate.net/publication/267569764\\_The\\_OpenFOAM\\_Technology\\_Primer](https://www.researchgate.net/publication/267569764_The_OpenFOAM_Technology_Primer)
- The Finite Volume Method in Computational Fluid Dynamics book by Moukalled, Magani and Darwish:  
<https://link.springer.com/book/10.1007/978-3-319-16874-6>



# Case Structure

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

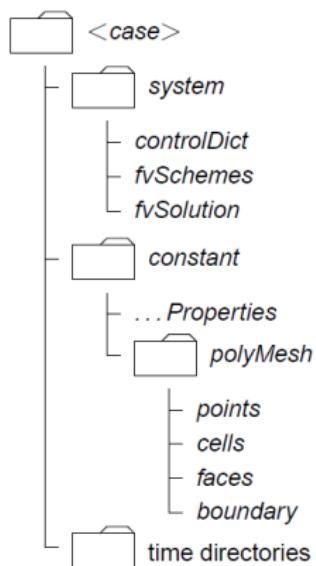
Dictionaries

system

constant

Time

- A simulation (case) in OpenFOAM is a hierarchy of directories and text files:



## Case Structure



- Any solver run will expect to have the following three directories in the case directory:
    - system
    - constant
    - time directories (e.g. 0, 0.1, etc)
  - Each directory contains a series of files and sub-directories that help define the computational problem.

# Case Structure

- Computational data is classified either as heavy- or light-weight data.
- Different formats are used for storing different types of data.
- The `system` directory contains light-weight data required to control execution of the code.
- The `constant` directory contains a mixture of light- and heavy-weight data.
- `Time` directories almost exclusively store heavy-weight data.



Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionarys

system

constant

Time

# Dictionaries

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- A dictionary is an entity that contains data entries that can be retrieved by the I/O system by means of keywords.
- The general format for dictionary entry:

```
<keyword> <dataEntry1> ... <dataEntryN>
```

- A sub-dictionary is identified by the name followed by data entries within curly braces:

```
<dictionaryName>
{
    ... keyword entries ...
}
```

# File dictionary

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- Each data file must have a description entry:

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

# Dictionary examples

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- An example of a file dictionary,  
*transportProperties*:

```
transportModel Newtonian;  
nu           1.5e-05;
```

- An example of a sub-dictionary:

```
kEpsilonCoeffs  
{  
    Cmu          0.09;  
    C1           1.44;  
    C2           1.92;  
    alphaEps     0.76923;  
}
```

# Dimensions

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- OpenFOAM is fully dimensional with dimensions required for each field and physical property
- Dimensions are checked for consistency in all field operations
- Some entries assume dimensions and are not required, such as viscosity.

No	1	2	3	4	5	6	7
Property	Mass	Length	Time	Temperature	Quantity	Current	Luminous Intensity
Unit	Kilogram	meters	second	Kelvin	moles	ampere	candela

- Kinematic viscosity  $m^2/s$  would be represented as

[ 0 2 -1 0 0 0 0 ] ;

# system directory

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- The `system` directory is the repository for files responsible for the control of execution of the solver.
- Files in `system` directory are run-time changeable, the solver reads their contents if they were changed during the run.
- `system` files:
  - `controlDict` is responsible for the solver execution.
  - `fvSchemes` is responsible for the specification of the discretization schemes.
  - `fvSolution` is responsible for linear and non-linear solver controls.

# controlDict

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- `controlDict` file contains entries that are used to control run time behaviour of the solver.
- In general, there are several sections in `controlDict` file responsible for the control of different aspects of the solver execution:
  - time control
  - data writing
  - data reading
  - run-time loadable functionality

## Time controls:

- `startFrom` - controls the start time of the simulation
  - `firstTime` - earliest time step from the set of time directories
  - `startTime` - time specified by the `startTime` keyword entry
  - `latestTime` - most recent time step from the set of time directories
- `startTime` - start time for the simulation with  
`startFrom startTime`



## Time controls:

- **stopAt**
  - `endTime` - time specified by the `endTime` keyword entry
  - `writeNow` - stops simulation on completion of current time step and writes data
  - `noWriteNow` - stops simulation on completion of current time step and does not write out data
  - `nextWrite` - stops simulation on completion of next scheduled write time, specified by `writeControl`
- **endTime** - end time for the simulation when `stopAt endTime;` is specified
- **deltaT** - time step of the simulation

## Data writing

- `writeControl` - controls the timing of write output to file:
  - `timeStep` - writes data every `writeInterval` time steps
  - `runtime` - writes data every `writeInterval` seconds of simulated time
  - `adjustableRunTime` - Writes data every `writeInterval` seconds of simulated time, adjusting the time steps to coincide with the `writeInterval` if necessary. Most useful in cases with automatic time step adjustment.
  - `cpuTime` - writes data every `writeInterval` seconds of CPU time
  - `clockTime` - writes data out every `writeInterval` seconds of real time

- `writeInterval` - integer used in conjunction with `writeControl` described above
- `purgeWrite` - integer representing a limit on the number of time directories that are stored by overwriting time directories on a cyclic basis
- `writeFormat` - specifies the format of the data files.
  - `ascii` - ASCII format, written to `writePrecision` significant figures
  - `binary` - Binary format.
- `writePrecision` - integer used in conjunction with `writeFormat` described above (default = 6)
- `writeCompression` - specifies the compression of the data files
  - `uncompressed` - no compression
  - `compressed` - gzip compression

- **timeFormat** - Choice of format of the naming of the time directories:
  - **fixed** -  $+/-m.aaaaaaaa$  where number a is set by timePrecision
  - **scientific** -  $+/-m.aaaaaaaaa +/-xx$  where number a is set by timePrecision
  - **general** - specifies scientific format if the exponent is less than -4 or greater than or equal to that specified by timePrecision
- **timePrecision** - integer used in conjunction with timeFormat described above (default = 6)

## Data reading:

- `runTimeModifiable` - yes/no switch for whether dictionaries, e.g. `controlDict`, are re-read by solver at the beginning of each time step
- `libs` - list of additional libraries to be loaded at run-time
- `functions` - list of functions, e.g. probes to be loaded at run-time

# fvSolution

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- **fvSolution:** controls linear and non-linear solver execution
- Consists of at least two sections: **solvers** and **PISO**, **SIMPLE**, or **PIMPLE**.

- Linear solver controls:

```
solvers
{
    p
    {
        solver          PCG;
        preconditioner DIC;
        tolerance       1e-06;
        relTol          0;
    };
    U
    {
        solver          PBiCG;
        preconditioner DILU;
        tolerance       1e-05;
        relTol          0;
    };
}
```

- Non-linear controls:

PISO

{

```
nCorrectors      2;  
nNonOrthogonalCorrectors 0;  
pRefCell          0;  
pRefValue         0;
```

}

# fvSchemes

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- **fvSchemes**: controls discretization methods for various terms fields and PDE operators
- All entries are given in the form of dictionaries that are parsed at the run time.
- The following dictionaries are required for a typical flow problem:
  - `dttSchemes`
  - `gradSchemes`
  - `divSchemes`
  - `laplacianSchemes`
  - `interpolationSchemes`
  - `snGradSchemes`

- Time discretization scheme is defined as:

```
ddtSchemes
{
    default      Euler;
}
```

- Other possibilities are: CoEuler,  
CrankNicholson, Euler, SLTS, backward,  
bounded, localEuler, steadyState

# fvSchemes

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- Computations of gradients are defined as:

```
gradSchemes
{
    default          Gauss linear;
    grad(p)         Gauss linear;
}
```

- Other possibilities are: cellLimited,  
cellLimited<Venkatakrishnan>,  
cellLimited<cubic>, cellMDLimited,  
edgeCellsLeastSquares, faceLimited,  
faceMDLimited, leastSquares, fourth

# fvSchemes

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of

Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- Divergence discretization is defined as:

```
divSchemes
{
    default          none;
    div(phi,U)      Gauss linear;
}
```

- Find out other possibilities by “asking” the solver:

```
divSchemes
{
    default          none;
    div(phi,U)      Gauss ?;
}
```

- Please observe that only Gauss integration is possible.

- Surface normal gradient is prescribed as:

```
snGradScheme
{
    default           corrected;
}
```

- Laplacian Schemes prescribed as:

```
laplacianSchemes
{
    default           none;
    laplacian(nuEff,U) Gauss linear corrected;
    laplacian((1|A(U)),p) Gauss linear corrected;
    laplacian(1,p) Gauss linear limited 1;
}
```

# Structure of constant directory

- As the directory's name indicates, the `constant` directory is a place for files and directories considered constant during the run.
- While solvers have the ability to change their settings at run time, the `constant` directory is not a subject to that look-up.
- The `constant` directory is a placeholder for the mesh files and material properties data needed for the run.
- In the case of laminar flows, the `transportProperties` file is required to specify the kinematic viscosity.



Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

# polyMesh

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

ictionaries

system

constant

Time

- Mesh and topological data is stored in the `constant/polyMesh` directory:
  - `boundary`: contains a list of boundary patches with their index offsets.
  - `faces`: contains the list of faces together with points describing the face positively oriented in counter-clock direction.
  - `owner/neighbour`: contains the list of faces consecutively numbered from 0 to  $N_{faces} - 1$  by their position in the list with index of the neighbour/owner cell.
  - `points`: contains the list of points consecutively numbered by the position in the list together with the coordinate position of the point as a triplet of values.

# Time directories

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of

Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- 0 directory is a special case of a general time directory.
- A general time directory stores the boundary, field and mesh data (case of dynamic mesh) required for post-processing and solver restarts.
- Therefore, field initialization, computed field values and corresponding boundary conditions are located in these directories.
- In order to start the solver run, at least one time directory is required.

# Time directories

- In the case of time dependent simulations starting from the time 0, the time directory is typically called “/ 0” directory containing the corresponding field files required by a given solver.
- Each solver requires different data to be specified on boundaries and within the computational domain.

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

# Types of Boundaries

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

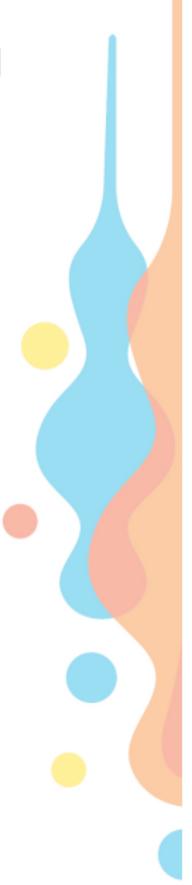
Dictionaries

system

constant

Time

- OpenFOAM has a concept of primitive and numerical boundary conditions.
  - Primitive boundaries are the actual surface patches where the numerical condition is applied.
  - Numerical boundary conditions assign a field value at the surface patch.
- Several types of primitive boundaries are available in OpenFOAM
  - **Constrained** - Applies a constraint to a field.
  - **Unconstrained** - Assigns field values through a numerical condition.



# Types of Boundaries

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- Constrained primitive boundaries

- `symmetry`
- `cyclic`
- `empty`
- `wedge`
- `processor`

- `symmetry` and `cyclic` boundaries are used to put constraints on fields.
- `empty` used for 2D simulations.
- `wedge` used for axisymmetric simulations.
- `processor` boundary is used for inter-processor communication.

# Types of Boundaries

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of

Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

- Unconstrained primitive boundaries
  - [patch](#)
  - [wall](#)
- In general, [patch](#) boundary is commonly used for inlets, outlets, and slip walls.
- [wall](#) boundary is used for no-slip walls. This boundary condition is not contained in the [patch](#) boundary condition because specialised modelling options can be used on this boundary condition, such as turbulence.
- Numerical boundary conditions can be one of the following types:
  - fixed value (Dirichlet)
  - specified gradient (Neumann)
  - mixed condition (Robin = Dirichlet + Neumann)

# Questions

Flavours of  
OpenFOAM®

Background

What?

Why?

History

Flavours

Directory

HPC

Architecture

OO Concepts

Main Concepts

Equation mimicking

Implementation of  
Models

Examples

Examples

Learning  
resources

Case

Structure

Dictionaries

system

constant

Time

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