

OpenFOAM Workshop

Programming session:

*from the C++ basics to the compilation of
user libraries*

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Presentation overview:

- C++ basics
 - main, declarations, types
 - memory management: references and pointers
 - Object orientation:
 - ▶ classes
 - ▶ inheritance
 - header and source files
- Compiling applications
 - make, wmake, wmake libso. Linking libraries.
- Writing simple applications in OpenFOAM
 - general structure (includes, skim between time-dirs...)
 - examples: divergence, INDT

30 m

15 m

Presentation overview:

- **Modifying existing libraries**
 - overview of the force class
 - modifying the class
 - ▶ Change the source and modify the names
 - ▶ compile the library
 - ▶ call the library during the execution
- **Adding classes**
 - FSInterface class
 - Linking classes through pointers

20 m

25 m

PART I

C++ basics

C++ basics

Hello world code... it only prints a message on screen

```
#include <iostream>
using namespace std;

int main()
{
    cout << "Hello world..! " << endl;
    return 0;
}
```

C++ basics

Declare variables, assign values and output the results

```
#include <iostream>
using namespace std;

int main()
{
    int iA;
    float rB, rC;

    iA=10;
    rB=0.4;
    rC=0.7;

    cout << "iA= " << iA << endl;
    cout << "rB= " << rB << endl;
    cout << "rC= " << rC << endl;

    return 0;
}
```

C++ basics

Define and use of the functions

```
#include <iostream>
using namespace std;

//Declare functions BEFORE!!
float Add(float a, float b)
{
    float c;
    c=a+b;
    return c;
}

//Main code
int main()
{
    int iA;
    float rB, rC, ResAdd;

    iA=10;
    rB=0.4;
    rC=0.7;

    ResAdd = Add(rB,rC);

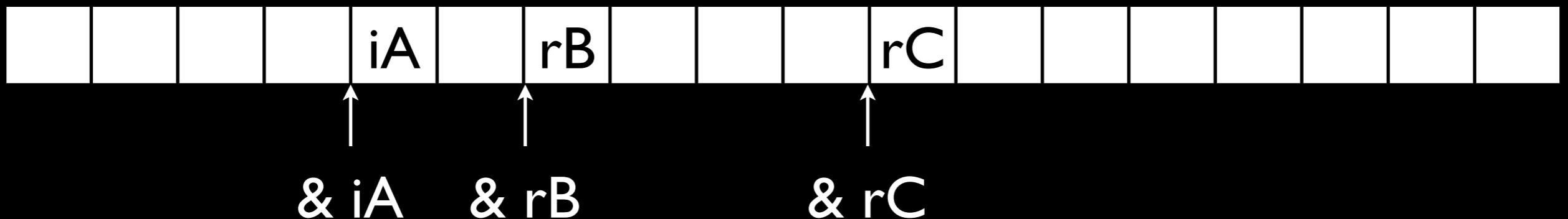
    cout<<"ResAdd= "<<ResAdd<<endl;

    return 0;
}
```

C++ basics

References and pointers

Computer memory: sequential representation



- References (`&`) are the addresses of the variables in the computer memory. References are constant.
- The value of a reference (ex: `0x7fff5fbfb604`) can be stored in a particular type of variable, called pointer.
- The pointer is a variable of the same type of the variable it points to. Integer pointer points to integer variable!

C++ basics

Declaration

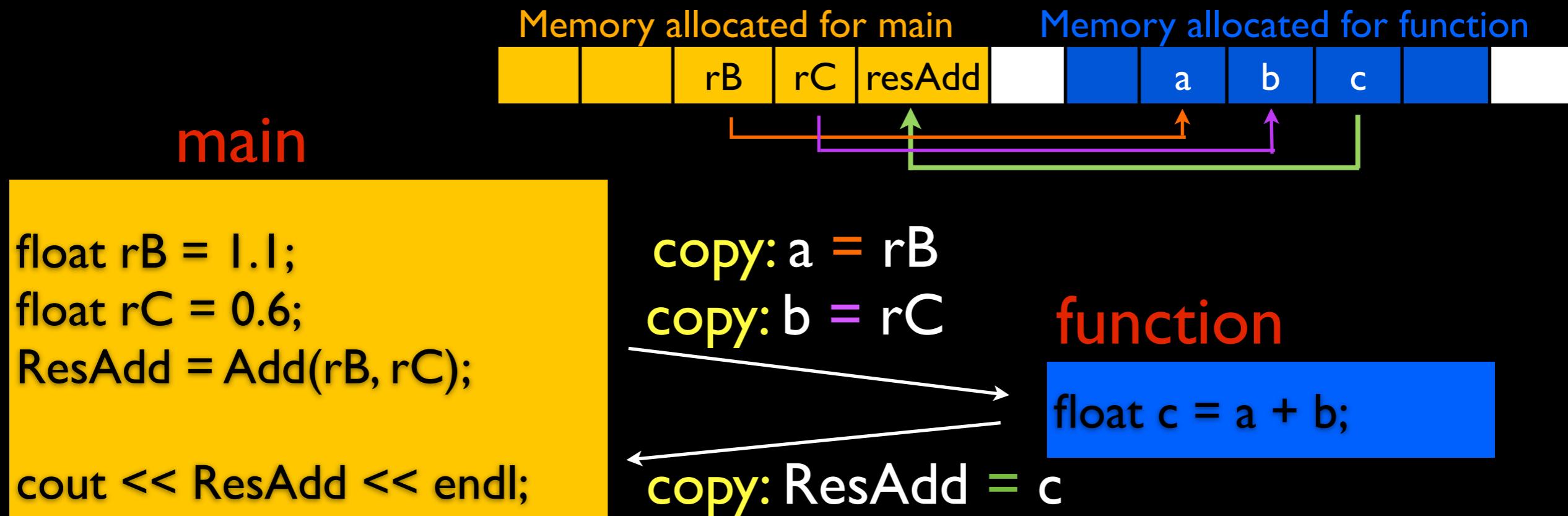
```
int * APtr = & iA
```

```
float * rPtr = & rB
```

...

Re-Assiguation

- Using references and pointers is convenient: it avoids the need of making copies when passing arguments to functions (for large arrays this is very convenient!)



C++ basics

- The **dereferencing operator** `*` restitutes the value of the variable pointed by the pointer. So:

Variable declaration and assignation:

Pointer declaration and assignation:

Pointer dereferencing:

`float rB = 1.1;`

`float * rPtr = & rB`

`cout << rB << endl;` → 1.1

`cout << * rPtr << endl;` → 1.1

`cout << & rB << endl;` → 0x7fff5fbfb604

Similar syntax, but very
different meaning!!

C++ basics

Optimizing the function

```
#include <iostream>
using namespace std;

//Declare functions BEFORE!!
float Add(float * a, float * b)
{
    return *a + *b;
}

//Main code
int main()
{
    int iA;
    float rB, rC, ResAdd;

    iA=10;
    rB=0.4;
    rC=0.7;

    ResAdd = Add(&rB,&rC);

    cout<<"ResAdd= "<<ResAdd<<endl;

    return 0;
}
```

2 Assign the pointer:
 $\text{float } * \text{a} = \& \text{rB}$

3 Deference and
return therefore
a float value

1 Pass references to
the function
(no copies!)

4 Assign the
returned value to
the float variable

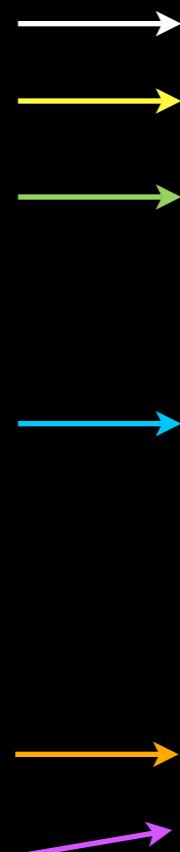
C++ basics

- CRectangle class

Class declaration
Private members
Public members
and functions

Public function
definition

Object declaration
Call the function
member of the class



```
#include <iostream>
using namespace std;

// classes example (from cplusplus.com)
#include <iostream>
using namespace std;

class CRectangle {
    int x, y;
public:
    void set_values (int,int);
    int area () {return (x*y);}
};

void CRectangle::set_values (int a, int b) {
    x = a;
    y = b;
}

int main () {
    CRectangle rect;
    rect.set_values (3,4);
    cout << "area: " << rect.area()<<endl;
    return 0;
}
```

Declarations

Definition

C++ basics

The function is called with the operator “.” : rect.setValues(3,4)

The function can also be called using a pointer:

```
CRectangle tria  
CRectangle * triaPtr
```

```
triaPtr = & tria;
```

```
cout << tria.area() << endl;  
cout << * tria.area() << endl;  
cout << triaPtr -> area() << endl
```



declare the variable
declare the pointer

assign the pointer

call fcn using object

call fcn using object, by
de-referencing pointer

call fcn using pointer

```
// classes example (from cplusplus.com)
#include <iostream>
using namespace std;
|
class CRectangle {
    int x, y;
public:
//constructor
CRectangle(); //default constructor
CRectangle(int,int);
//destructor
~CRectangle();
//member functions
int area (void);
};

CRectangle::CRectangle ()
{
    x = 1;
    y = 1;
}

CRectangle::CRectangle (int a, int b)
{
    x = a;
    y = b;
}

CRectangle::~CRectangle ()
{
//do nothing
}

int CRectangle::area()
{
    return (x*y);
}
```

C++ basics

- The constructor is a class-member function called when the object is initially build
- Dual of the Constructor is the Destructor
- As for every function, different arguments define different constructors; the constructor can also call other functions

Standard syntax for the constructor

```
int main ()
{
    CRectangle rect;
    CRectangle rect2(3,4);

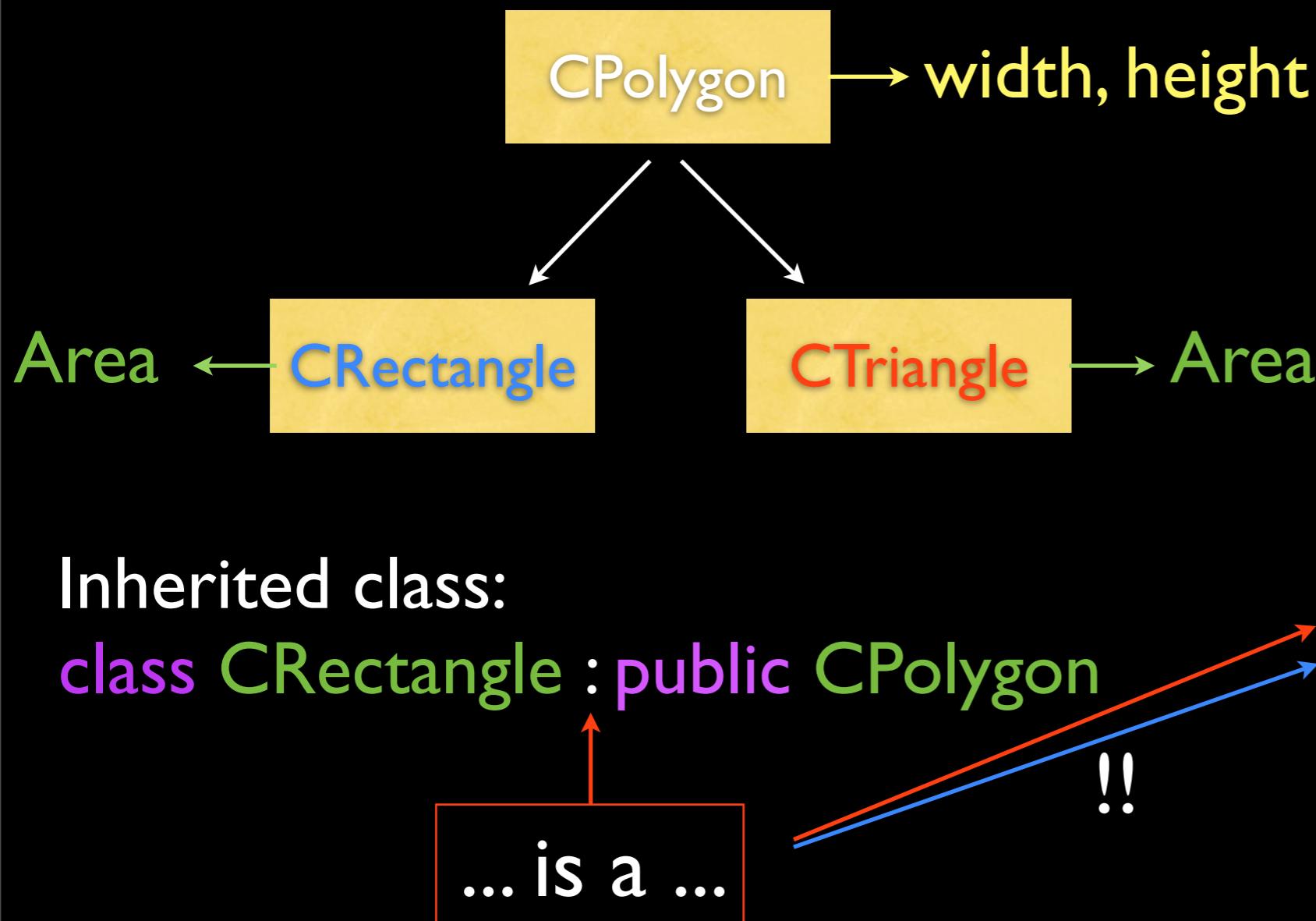
    cout << "area1: " << rect.area()<<endl;
    cout << "area2: " << rect2.area()<<endl;

    return 0;
}
```

```
CRectangle::CRectangle (int a, int b)
:
x(a),
y(b)
{
//do nothing
}
```

C++ basics

- Inheritance: CPolygon class



Inherited class:

class CRectangle : public CPolygon

Every object has its own function definition. The function behaves differently accordingly to the type of the object

C++ basics

- Main (.cpp; .C), Header (.H) and Source (.C) files

Main code
.CPP, .C file

```
#include <iostream>
using namespace std;

// classes example (from cplusplus.com)
#include <iostream>
using namespace std;

class CRectangle {
    int x, y;
public:
    void set_values (int,int);
    int area () {return (x*y);}
};

void CRectangle::set_values (int a, int b) {
    x = a;
    y = b;
}

int main () {
    CRectangle rect;
    rect.set_values (3,4);
    cout << "area: " << rect.area()<<endl;
    return 0;
}
```

Declarations
.H file

Definition
.C file

C++ basics

- Main (.cpp; .C), Header (.H) and Source (.C) files

Main code
.CPP, .C file

```
// classes example (from cplusplus.com)
#include <iostream>
using namespace std;

#include "CRectangle.H"

int main () {
    CRectangle rect;
    rect.set_values (3,4);
    cout << "area: " << rect.area()<<endl;
    return 0;
}
```

Declarations
.H file

```
#ifndef CRectangle_H
#define CRectangle_H

class CRectangle {
    int x, y;
public:
    void set_values (int,int);
    int area () {return (x*y);}
};

#include "CRectangle.C"
#endif
```

Definition
.C file

```
void CRectangle::set_values (int a, int b) {
    x = a;
    y = b;
}
```

C++ basics

- Compiling applications: passing from human readable instructions to binaries
- Unix environment: call makefile through “make”

Set the compiler

Compiler options

Compiler Flags;
external libraries

Source files

Executable

```
CC=g++
CFLAGS=-c -Wall
LDFLAGS=

#-----
SOURCES=main.cpp
#-----

OBJECTS=$(SOURCES:.cpp=.o)

#-----
EXECUTABLE=program
#-----

all: $(SOURCES) $(EXECUTABLE)
    $(EXECUTABLE): $(OBJECTS)
        $(CC) $(LDFLAGS) $(OBJECTS) -o $@
    .cpp.o:
        $(CC) $(CFLAGS) $< -o $@
```

- The “make” command is overwritten in OpenFOAM by “wmake” and “wmake libso”

References

Several books are available on C++. For example:

Deitel, Deitel
C++, How to program
ed. Prentice Hall

A book is generally better (especially for C++)...
but on-line useful guidance can be also found:

<http://www.cplusplus.com/doc/tutorial/>

PART 2

Writing simple applications in OpenFOAM

General structure of an OpenFOAM application

```
# include "fvCFD.H"
int main(int argc, char *argv[])
{
    # include "setRootCase.H"
    # include "createTime.H"
    # include "createMesh.H"
    # include "createFields.H"
    ...A lot of nice code...
    return(0);
}
```

```
//
// createTime.H
// ~~~~~
Foam::Info<< "Create time\n" << Foam::endl;
Foam::Time runTime
(
    Foam::Time::controlDictName,
    args.rootPath(),
    args.caseName()
);
```

src/finiteVolume/cfdTools/general/include/fvCFD.H
several other include: classes for time, mesh, geometry, math constants...

src/OpenFOAM/include createTime.H
Declares runTime, object of the class
Foam::Time.

Constructor defined in
src/OpenFOAM/db/Time/Time.H, line 193

//- Construct given dictionary, rootPath and casePath
Time
(
 const dictionary& dict,
 const fileName& rootPath,
 const fileName& caseName,
 const word& systemName = "system",
 const word& constantName = "constant"
)

General structure of an OpenFOAM application

```
# include "fvCFD.H"

int main(int argc, char *argv[])
{
    # include "setRootCase.H"
    # include "createTime.H"
    # include "createMesh.H"
    # include "createFields.H"
    ...A lot of nice code...
    return(0);
}

volScalarField divergence
(
    IOobject
    (
        "divergence",
        runTime.timeName(),
        mesh,
        IOobject::READ_IF_PRESENT,
        IOobject::AUTO_WRITE
    ),
    mesh
);
```

File in the source directory
Declares a **VolScalarField** called *divergence* to be written in every time-step folder

Calculating the divergence $\nabla \cdot \vec{v}$

$$\nabla \cdot \vec{v} = \frac{\partial v_1}{x_1} + \frac{\partial v_2}{x_2} + \frac{\partial v_3}{x_3} = \delta_{ik} v_{i,k}$$

```
instantList timeDirs = timeSelector::select0(runTime, args);  
forAll(timeDirs, timeI)  
{  
    runTime.setTime(timeDirs[timeI], timeI);  
    Info<< "Time = " << runTime.timeName() << endl;  
    Info<< "Reading field U\n" << endl;
```

```
//Reading field  
volVectorField U  
{  
    IObject  
{  
        "U",  
        runTime.timeName(),  
        mesh,  
        IObject::MUST_READ,  
        IObject::AUTO_WRITE  
    },  
    mesh  
};
```

```
divergence=fvc::div(U);  
divergence.write();
```

Check for existing time dirs
OpenFOAM version of the
for loop. Equivalent to:

```
for(int timel=0; i<timeDirs.size(); timel++)  
{...}
```

Declares and reads the
field U from the selected
time directory

calculates the field

Compiling the application

Working directory:

SolutionDivergence.C

createFields.H

Make

files

options

solutionDivergence.C

EXE = \$(FOAM_USER_APPBIN)/DivU

EXE_INC = \

-I\$(LIB_SRC)/finiteVolume/lnInclude

EXE_LIBS = -lfiniteVolume

Source file

Compile application

include headers for FV

include FV library

The application is compiled typing at terminal the command: wmake

Calculating the Normalised invariant of the deformation tensor

$$D = \frac{S_{ij}S_{ij} - W_{ij}W_{ij}}{S_{ij}S_{ij} + W_{ij}W_{ij}}; S_{ij} = \frac{1}{2} \frac{\partial u_i}{\partial x_j} + \frac{\partial u_j}{\partial x_i}; W_{ij} = \frac{1}{2} \frac{\partial u_i}{\partial x_j} - \frac{\partial u_j}{\partial x_i}$$

CreateFields.h

```
volScalarField D
(
    I0object
    (
        "D",
        runTime.timeName(),
        mesh,
        I0object::READ_IF_PRESENT,
        I0object::AUTO_WRITE
    ),
    mesh
);
```

Core of the main code

```
volTensorField gradU = fvc::grad(U);

volSymmTensorField S = 0.5*symm(gradU); // symmetric part of tensor
volTensorField W = 0.5*skew(gradU); // anti-symmetric part

volScalarField SS = S&&S;
volScalarField WW = W&&W;

volScalarField D
(
    I0object
    (
        "D",
        runTime.timeName(),
        mesh,
        I0object::NO_READ,
        I0object::NO_WRITE
    ),
    (SS - WW) / (WW + SS)
);

D.write();
```

———— Double inner product operator,
see Programmer guide P23

PART 3

Modifying existing libraries

Constructor of the class forces

```
Foam::forces::forces
(
    const word& name,
    const objectRegistry& obr,
    const dictionary& dict,
    const bool loadFromFiles
):
    name_(name),
    obr_(obr),
    active_(true),
    log_(false),
    patchSet_0,
    pName_(word::null),
    UName_(word::null),
    rhoName_(word::null),
    directForceDensity_(false),
    fDName_(""),
    rhoRef_(VGREAT),
    pRef_(0),
    CofR_(vector::zero),
    forcesFilePtr_(NULL)
{
    ...
    read(dict);
}
```

Path:

<OpenFOAM/src/postProcessing/functionObjects/forces/forces/Force.C>

Reference to the Object Registry. This is a list of the entities pertaining to an object

Reference to the controlDict

Call the member function forces::read
Read the entries in the controlDict

Reading entries from the controlDict

```
void Foam::forces::read(const dictionary& dict)
{
    log_ = dict.lookupOrDefault<Switch>("log", false);
    const fvMesh& mesh = refCast<const fvMesh>(obr_);
    patchSet_ =
        mesh.boundaryMesh().patchSet(wordList(dict.lookup("patches")));
    ...
    // Optional entries U and p
    pName_ = dict.lookupOrDefault<word>("pName", "p");
    UName_ = dict.lookupOrDefault<word>("UName", "U");
    rhoName_ = dict.lookupOrDefault<word>("rhoName", "rho");
    ...
    // Reference density needed for incompressible calculations
    rhoRef_ = readScalar(dict.lookup("rhoInf"));
    // Reference pressure, 0 by default
    pRef_ = dict.lookupOrDefault<scalar>("pRef", 0.0);
    // Centre of rotation for moment calculations
    CofR_ = dict.lookup("CofR");
}
```

patches on which forces
will be integrated

system/controlDict:
functions

```
(  
    forces  
    {  
        type forces;  
        functionObjectLibs ("libforces.dylib");  
        outputControl outputTime;  
        patches (wing);  
        pName p;  
        Uname U;  
        rhoName rhoInf;  
        rhoInf 1.2; //Reference density  
        pRef 0;  
        CofR (0 0 0); //Origin for moments  
    }  
)
```

Calculating the forces

The **virtual** function `write()` is called during the execution. This calls `forces::calcForcesMoment()`, where the calculation is performed

```
forAllConstIter(labelHashSet, patchSet_, iter) {  
    label patchi = iter.key();  
  
    vectorField Md = mesh.C().boundaryField()[patchi] - CofR_;  
  
    vectorField pf = Sfb[patchi]*(p.boundaryField()[patchi] - pRef);  
  
    fm.first().first() += rho(p)*sum(pf);  
    fm.second().first() += rho(p)*sum(Md ^ pf);  
  
    vectorField vf = Sfb[patchi] & devRhoReffb[patchi];  
  
    fm.first().second() += sum(vf);  
    fm.second().second() += sum(Md ^ vf);  
}
```

OpenFOAM iterator.
it corresponds to a for
cycle

mesh is object of the class fvMesh
The expression returns a vector
with the cell centres of the chosen
patch

const DimensionedField< scalar, volMesh > & V00 () const	<i>Return old-old-time cell volumes.</i>
const surfaceVectorField & Sf () const	<i>Return cell face area vectors.</i>
const surfaceScalarField & magSf () const	<i>Return cell face area magnitudes.</i>
const surfaceScalarField & phi () const	<i>Return cell face motion fluxes.</i>
const volVectorField & C () const	<i>Return cell centres as volVectorField.</i>
const surfaceVectorField & Cf () const	

Calculating the forces

The **virtual** function `write()` is called during the execution. This calls `forces::calcForcesMoment()`, where the calculation is performed

```
forAllConstIter(labelHashSet, patchSet_, iter)
{
    label patchi = iter.key();

    vectorField Md = mesh.C().boundaryField()[patchi] - CofR_;

    vectorField pf = Sfb[patchi]*(p.boundaryField()[patchi] - pRef);

    fm.first().first() += rho(p)*sum(pf);
    fm.second().first() += rho(p)*sum(Md ^ pf);

    vectorField vf = Sfb[patchi] & devRhoReffb[patchi];
    fm.first().second() += sum(vf);
    fm.second().second() += sum(Md ^ vf);
}
```

Sfb is the (reference to) the face area vector

It is here multiplied for the pressure boundaryField => pf returns the vector of forces on the chosen patch

$$F = \rho \int pdA = \rho \sum p_i A_i$$

$$M = F \times r = \rho \sum f_i \times r_i$$

Re-compiling the forces library

The basic idea of the openFOAM environment is:
find something similar and modify it as you like, but
DO NOT TOUCH THE ORIGINAL SOURCES!

STEP 1: copy the forces directory from the original location into another directory

STEP 2: copy also the Make folder

STEP 3: substitute strings and modify (all) the file names

(sed ‘s/forces/Myforces/g’ forces.C > Myforces.C)

STEP 4: modify the local functionObject.H file (add the new class to the list of loadable functions)

```
#include "Myforces.H"  
...  
namespace Foam  
{  
    typedef OutputFilterFunctionObject<Myforces>forcesFunctionObject;  
}  
...
```

Re-compiling the forces library

STEP 4: Modify the Make/files:

```
Myforces.C  
forcesFunctionObject.C
```

```
LIB = $(FOAM_USER_LIBBIN)/LibMyforces
```

STEP 5: modify the Make/options file:

```
EXE_INC = \  
.... all what was already there ...  
-I$(LIB_SRC)/postProcessing/functionObjects/forces/lnInclude
```

include all what was needed by the original library!

STEP 6: compile with wmake libso

Re-compiling the forces library

STEP 7: Add the entries in the controlDict, in order for the library to be loaded and used:

```
libs ( "libMyforces.dylib" ); ← Load the library (.dylib on MAC; .so on Linux)  
functions  
(  
Myforces ← Search in the library for the entry called  
{  
    type Myforces;  
    functionObjectLibs ("libMyforces.dylib");  
    outputControl outputTime;  
    patches (BottWall);  
    pName p;  
    Uname U;  
    rhoName rhoInf;  
    rhoInf 1.0;  
    pRef 0;  
    CofR (0 0 0);  
}  
);
```

Use the library

References

Several examples can be found on-line:

http://www.tfd.chalmers.se/~hani/kurser/OS_CFD_2007/

http://www.tfd.chalmers.se/~hani/kurser/OS_CFD_2008/

http://www.tfd.chalmers.se/~hani/kurser/OS_CFD_2009/

http://www.tfd.chalmers.se/~hani/kurser/OS_CFD_2010/

See for example the work of A. Gonzales.

But have a deep look at the whole web-site,
there's a lot of enlightening material!!

PART 4

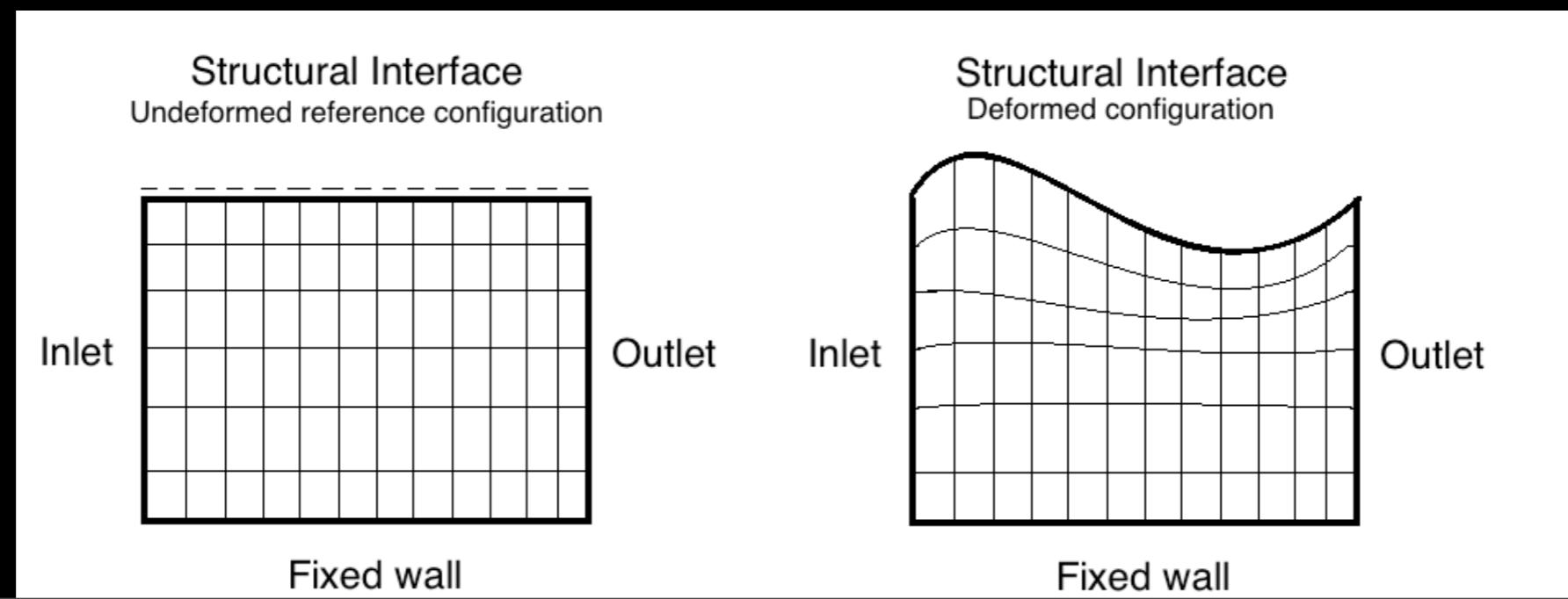
Adding new classes in OpenFOAM the FSInterface class

Scope of this class

- The class is designed to manage the mesh motions. It is used in the main of the solver `pimpleDyMFOAM`:
 - establish the communication (if needed) with the external solver trough MPI
 - send pressure data and retrieve mesh displacement data
 - communicate with the `AitkenControl` class, in charge for calculating the dynamic relaxation factor:

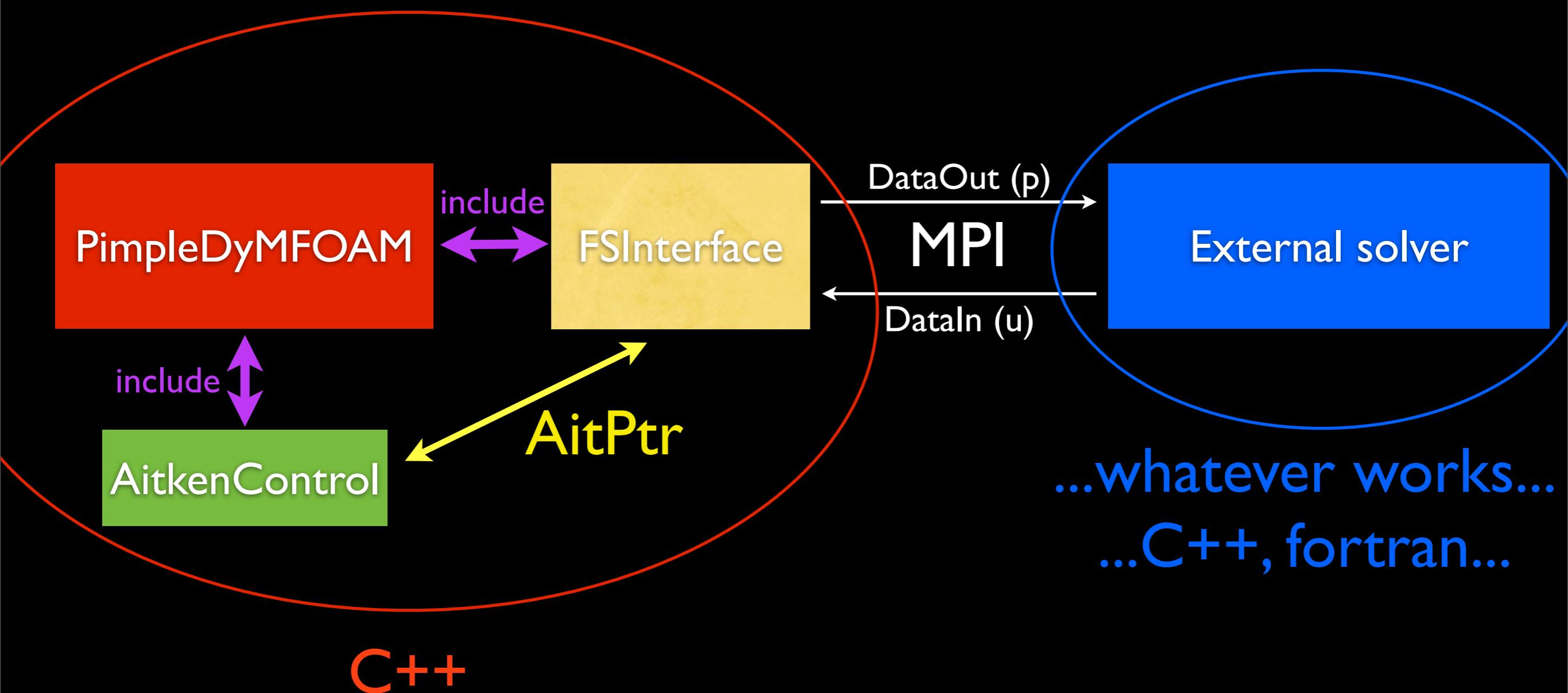
$$u_{k+1} = \omega_k \tilde{u}_{k+1} + (1 - \omega_k) u_k$$

- move the fluid mesh (ALE framework)



Scope of this class

Multiple Program Multiple Data type environment, the external solver is “spawned” during the execution time. This generates a communicator we can use for exchanging data (white arrows)



Use of the class

In the main solver: Include and declare

new classes definition

Foam classes needed for the mesh motion

Declare the new objects
using the right arguments!

Constructor of FSInterface:

header (.H) file

```
namespace Foam
{
    class FSInterface
    {
        //private data
        dynamicFvMesh& mesh;
        volScalarField& p;
        volVectorField& U;
        dimensionedScalar& rhoFluid;
    };
}
```

... a lot of other stuff...

```
// Constructor from components
FSInterface(dynamicFvMesh &mesh,
            volScalarField &p,
            volVectorField &U,
            dimensionedScalar &rhoFluid,
            Time &runTime,
            AitkenControl & alfa);
```

References are constant: they **MUST** be initialised at the creation. In this case this is done by passing the values to the constructor

source (.C) file

```
//- Constructor from components
FSInterface::FSInterface(dynamicFvMesh &mesh_,
                        volScalarField &p_,
                        volVectorField &U_,
                        dimensionedScalar &rhoFluid_,
                        Time &runTime_,
                        AitkenControl & alfa_)

    :
    mesh(mesh_),
    p(p_),
    U(U_),
    rhoFluid(rhoFluid_),
    runTime(runTime_),
    alfa2(alfa_)

{
    initialize();
}
```

Arguments passed to the class

Assign values to the class members

Execute other fcns

linking to the AitkenControl class:

An object of the type AitkenControl is created right before the object FSInterface. A reference to this object is passed to the constructor. This reference is stored in a pointer main code:

```
AitkenControl alfa2(mesh, runTime);
FSInterface interface(mesh,p,U,rhoFluid,runTime,alfa2);
```

alfa2 is instantiated in the constructor, the pointer is referenced also in the constructor:

```
FSInterface::FSInterface(dynamicFvMesh &mesh_,
                          volScalarField &p_,
                          volVectorField &U_,
                          dimensionedScalar &rhoFluid_,
                          Time &runTime_,
                          AitkenControl & alfa_)

:
mesh(mesh_),
p(p_),
U(U_),
rhoFluid(rhoFluid_),
runTime(runTime_),
alfa2(alfa_)

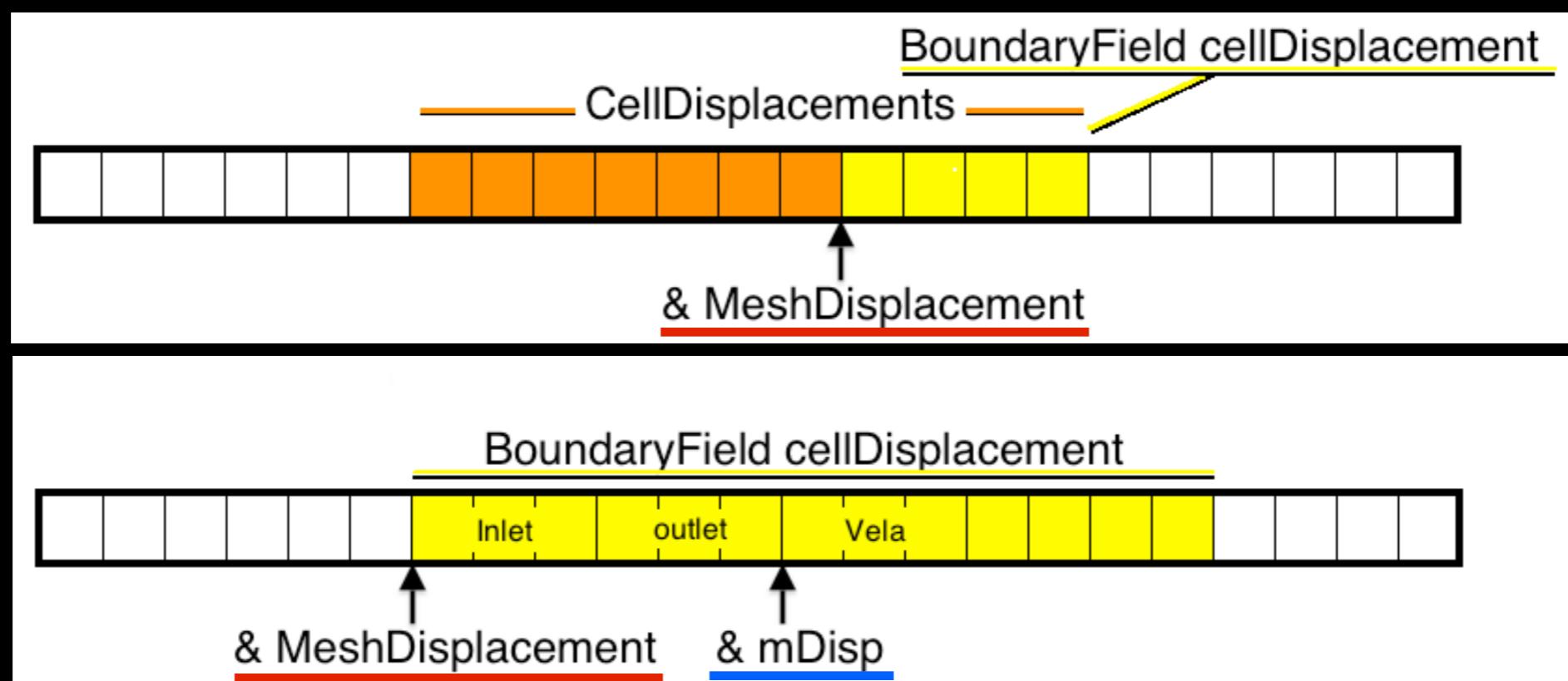
{
    AitPtr = & alfa2; //Pointer points to the object
}
```

Searching the mesh motion entries

Depending upon the motion solver, the mesh motion is stored in a field “`pointDisplacements`”, “`cellDisplacement`” or “`motionU`”.

A field in OpenFOAM is defined as: `internalField` + `boundaryField`

Imposing the motion of a boundary means writing the motion in the correspondent entry of the `boundaryField`. For example, “`&MeshDisplacement`” is the address of the `BoundaryField`, while “`&mDisp`” is the address of the mesh interface we want to move



Searching the mesh motion entries

```
IOdictionary couplingDict
```

```
(
```

```
IOobject
```

```
(
```

```
"CouplingDict",
```

```
runTime.constant(),
```

```
mesh,
```

```
IOobject::MUST_READ,
```

```
IOobject::NO_WRITE
```

```
)
```

```
);
```

```
word temp(couplingDict.lookup("fluidPatch"))  
word interface = temp;
```

```
label fluidPatchID = mesh.boundaryMesh().findPatchID(interface);
```

Opens the FOAM dictionary named “couplingDict”, to be read only. The dict must be placed in the folder constant

Searches and reads the entry “fluidPatch” in CouplingDict

Searches the the entry “fluidPatchName” in the mesh boundaryField. Returns a (integer) label: the id of the patch in the mesh order

Searching the mesh motion entries

We need now to find the references to the mesh motion boundaryField. Using displacementLaplacian...

```
→ pointVectorField & PointDisplacement =  
  const_cast<pointVectorField&>(mesh.objectRegistry::  
    lookupObject<pointVectorField>("pointDisplacement"));  
  
pDispPtr = & ( refCast<vectorField>(PointDisplacement.boundaryField()[fluidPatchID])); ←
```

Search in the objectRegistry of the mesh an object of the type:
pointVectorField called pointDisplacement. Whatever its type, impose
(const_cast) to be a reference of the type PointVectorField

The entry number “fluidPatchID” is the reference to the interface mesh motion. Store this reference into a pointer

The mesh motion is imposed using the surcharged operator == :

```
PointDisplacement.boundaryField()[ fluidSidel ] == U_kpl
```

Searching the mesh motion entries

We need now to find the references to the mesh motion boundaryField. Using laplaceFaceDecomposition...

```
const fvMesh& motionMesh =
    runTime.objectRegistry::lookupObject<fvMesh>(movingRegionName);

tetPointVectorField& motionU =
    const_cast<tetPointVectorField&>
    ( motionMesh.objectRegistry::lookupObject<tetPointVectorField>("motionU") );

tetPolyMesh& tetMesh = const_cast<tetPolyMesh&>(motionU.mesh());

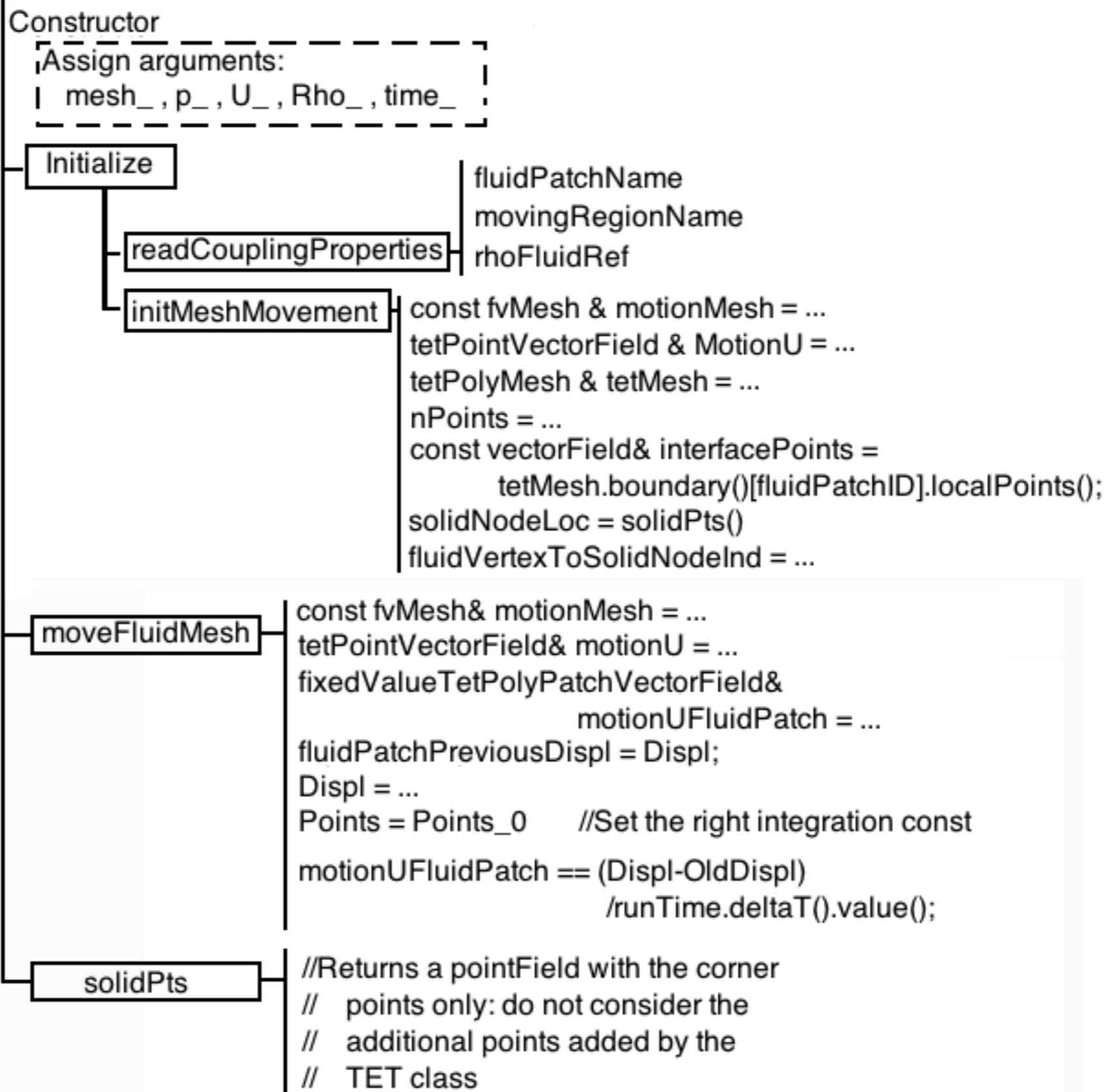
motionUFluidPatchPtr = &
    refCast<fixedValueTetPolyPatchVectorField>
    ( motionU.boundaryField()[fluidPatchID] );
```

The mesh motion is imposed using the surcharged operator == :

```
* motionUFluidPatchPtr == ( U_kp1 - U_old ) / runTime.deltaT().value();
```

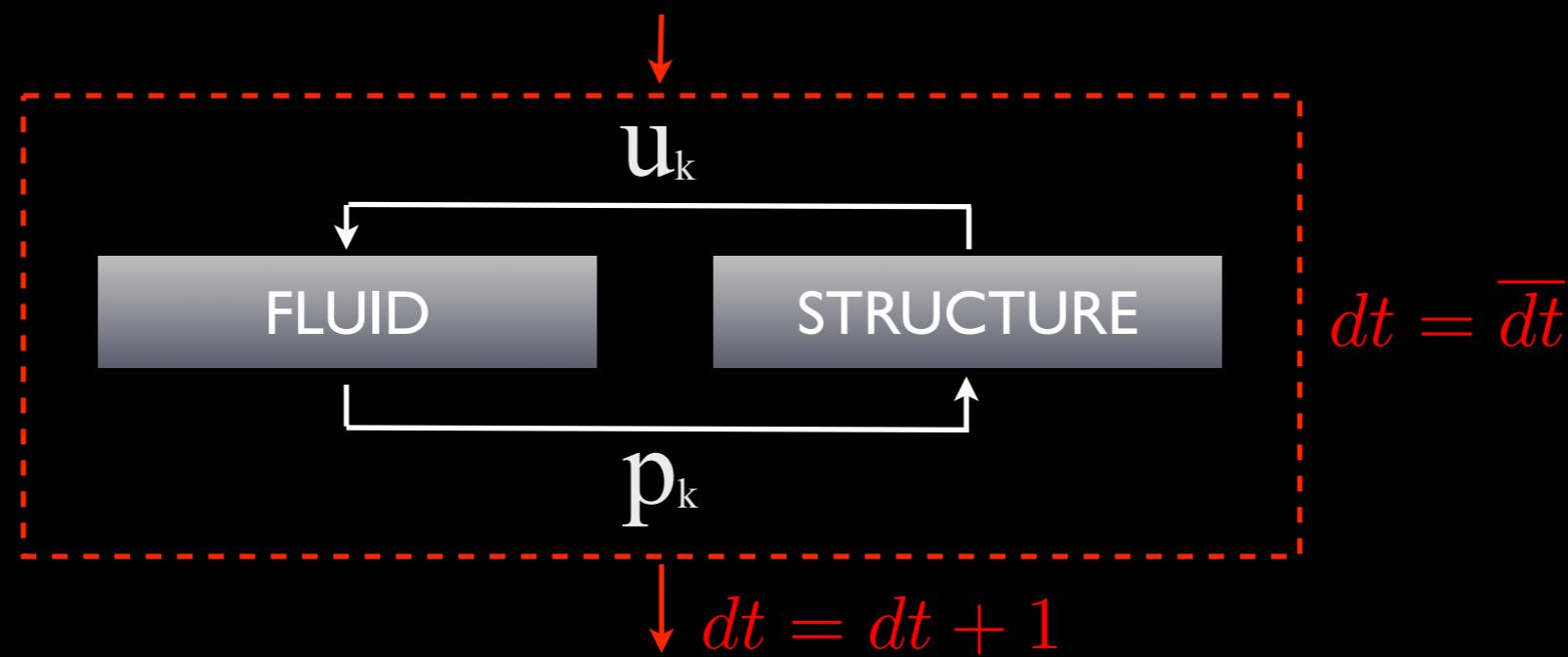
Scheme of the class

```
FSInterface (dynamicFvMesh & mesh, volScalarField & p, volVectorField & U,  
dimensionedScalar & rhoFluid, Time & runTime)
```



Algorithm

Implicit coupling: the equilibrium within the time-step is verified using a fixed-point between the fluid and the structural solver. Although this algorithm is computationally expensive, it is unconditionally stable



This is realised in OpenFOAM using a while loop within the time-step. The convergence criterion is verified by a function of the class `FSInterface`

Algorithm (main)

```
int main(int argc, char *argv[])
{
# include "setRootCase.H"
...
AitkenControl alfa2(mesh, runTime);
FSInterface interface(mesh,p,U,rhoFluid,runTime,alfa2);
while (runTime.run()) {
...
    runTime++;
    interface.MPISpawn(); //this is called the first time only
    int iterCnt=0, advance=0;
    do{
        iterCnt++;
        if (iterCnt == 1 )
        {advance=0;}
        interface.setCntr(iterCnt);
        if(runTime.value() > interface.FSI_init )
        {
            interface.sendPressures();
            interface.moveFluidMesh(); //CAREFUL: IT MUST BE BEFORE
                                         //THE FLUID CALCULATION
        }
        else
        {advance=1;}
        ... // --- PIMPLE loop
        ... // --- PISO loop
        if(runTime.value() > interface.FSI_init )
        { advance = interface.AdvContrl(); }
    }while(advance==0);
    runTime.write();
}
Info<< "End\n" << endl;
interface.makeClean();
```

time loop

Aitken iterations loop

Launch the structural solver

Send the number of iteration
to the interface (different
calculations if iterCntr == 1)

Send the pressures, retrieve
the displacements

Convergence is checked by
interface

Compiling

As for every FOAM application, we need to edit:

Make/files:

```
FSIfluidFOAMtet.C
```

→ Source file

```
EXE = $(FOAM_USER_APPBIN)/FSIfluidFOAMtetMPI
```

→ Name and path of the
compiled application

Make/options:

```
include $(RULES)/mplib$(WM_MPLIB)
```

```
EXE_INC = \
```

...all remains as in the original file...

```
EXE_LIBS = \
```

...all remains as in the original file, but add:

```
-lmpi \
```

```
$(WM_DECOMP_LIBS)
```

And finally some results!

FLUID DENSITY	FLUID VISCOSITY	STRUCTURE's THICKNESS	STRUCTURE's DENSITY	STRUCTURE's POISSON RATTIO	STRUCTURE's YOUNG'S MODULUS
ρ_f	ν_f	t_s	ρ_s	ν_s	E_s
Kg/m^3	m^2/s	m	Kg/m^3	-	N/m^2
1.0	0.01	0.002	500	0	250

$$u = 1 - \cos(2\pi t/5)$$

