

Chapter 1

Example problem: Finite-Reynolds-number flow inside an oscillating ellipse

In this example we consider our first moving-boundary Navier-Stokes problem: The flow of a viscous fluid contained in an elliptical ring whose walls perform periodic oscillations.

oomph-lib's Navier-Stokes elements are based on the Arbitrary Lagrangian Eulerian (ALE) form of the Navier-Stokes equations and can therefore be used in moving domain problems. In this example we illustrate their use in Domain - based meshes (first discussed in the example demonstrating the solution of the [unsteady heat equation in a moving domain](#)) in which MacroElements are used to update the nodal positions in response to changes in the domain boundary. In subsequent examples, we will discuss alternative, sparse mesh update techniques that are useful in problems with free boundaries and in fluid-structure interaction problems.

Finite-Reynolds-number-flow driven by an oscillating ellipse

We consider the unsteady 2D flow of a Newtonian fluid that is contained in an oscillating elliptical ring whose wall shape is parametrised by the Lagrangian coordinate ξ as

$$\mathbf{R}_w(\xi) = \begin{pmatrix} a(t) \cos(\xi) \\ a^{-1}(t) \sin(\xi) \end{pmatrix}$$

where

$$a(t) = A + \hat{A} \cos\left(\frac{2\pi t}{T}\right).$$

A represents the average half-axis of the elliptical ring in the x_1 -direction, and \hat{A} is the amplitude of its periodic variation. The ring has constant cross-sectional area – consistent with the incompressibility of the fluid whose motion is governed by the ALE form of the Navier-Stokes equations,

$$Re \left(St \frac{\partial u_i}{\partial t} + (u_j - u_j^M) \frac{\partial u_i}{\partial x_j} \right) = - \frac{\partial p}{\partial x_i} + \frac{\partial}{\partial x_j} \left(\frac{\partial u_i}{\partial x_j} + \frac{\partial u_j}{\partial x_i} \right), \quad (1)$$

and the continuity equation

$$\frac{\partial u_i}{\partial x_i} = 0,$$

where u_j^M is the mesh velocity. We exploit the symmetry of the problem and solve the equations in the quarter domain

$$D = \left\{ (x_1, x_2) \mid x_1 \geq 0, x_2 \geq 0, \left(\frac{x_1}{a(t)} \right)^2 + (x_2 a(t))^2 \leq 1 \right\},$$

shown in this sketch (for $A = 1$, $\hat{A} = 1$ and $T = 1$),

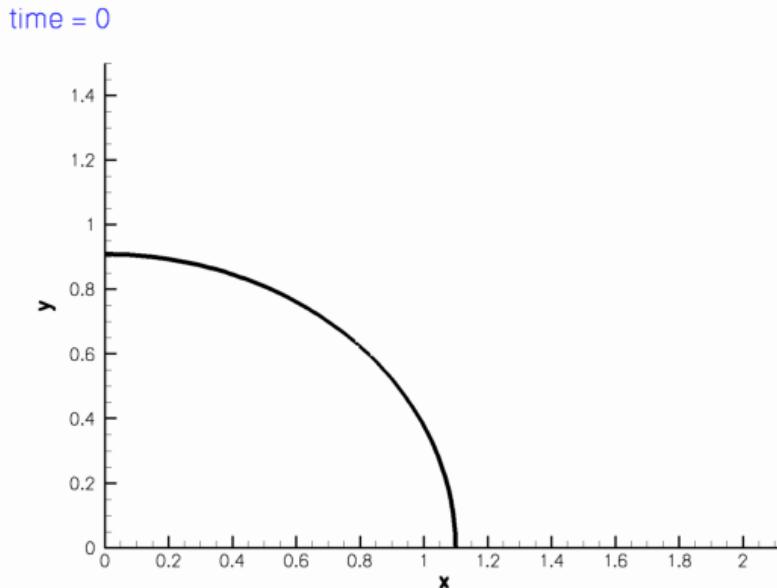


Figure 1.1 Sketch of the computational domain.

The fluid is subject to no-slip boundary conditions on the curved wall,

$$\mathbf{u}|_{\partial D_{ellipse}} = \frac{\partial \mathbf{R}_w(\xi, t)}{\partial t}$$

and symmetry conditions on the symmetry boundaries,

$$u_1|_{x_2=0} = 0, \quad u_2|_{x_1=0} = 0.$$

The initial conditions for the velocity are given by

$$\mathbf{u}(x_1, x_2, t = 0) = \mathbf{u}_{IC}(x_1, x_2),$$

1.1 An exact solution

It is easy to show (by inspection) that the unsteady stagnation point flow

$$u_1(x_1, x_2, t) = \frac{1}{a} \frac{da}{dt} x_1 = -\frac{2\pi \hat{A} \sin(\frac{2\pi t}{T})}{T(A + \hat{A} \cos(\frac{2\pi t}{T}))} x_1 \quad \text{and} \quad u_2(x_1, x_2, t) = -\frac{1}{a} \frac{da}{dt} x_2 = -\frac{2\pi \hat{A} \sin(\frac{2\pi t}{T})}{T(A + \hat{A} \cos(\frac{2\pi t}{T}))} x_2,$$

is an exact solution of the above problem as it satisfies the Navier-Stokes equations and the velocity boundary conditions. The pressure is given by

$$p = \frac{2\pi \hat{A} Re \left(x_1^2 St \cos(\frac{2\pi t}{T}) A + x_1^2 St \hat{A} - x_1^2 \hat{A} + x_1^2 \hat{A} \cos^2(\frac{2\pi t}{T}) - x_2^2 St \cos(\frac{2\pi t}{T}) - x_2^2 St \hat{A} - x_2^2 \hat{A} + x_2^2 \hat{A} \cos^2(\frac{2\pi t}{T}) \right)}{T^2 \left(A^2 + 2A\hat{A} \cos(\frac{2\pi t}{T}) + \hat{A}^2 \cos^2(\frac{2\pi t}{T}) \right)}.$$

1.2 Results

The two figures below show two snapshots of the solution for $Re = Re St = 100$, extracted from an animation of the results computed with [Taylor-Hood](#) and [Crouzeix-Raviart elements](#). In both cases, the exact solution was used as the initial condition for the velocities. The figures show "carpet plots" of the two velocity components and the pressure, respectively, and a contour plot of the pressure, superimposed on the moving mesh. The carpet plot of the velocities clearly shows that the flow is of stagnation-point type as the horizontal velocity, u_1 , is a linear function of x_1 while the vertical velocity, u_2 , is a linear function of $-x_2$.

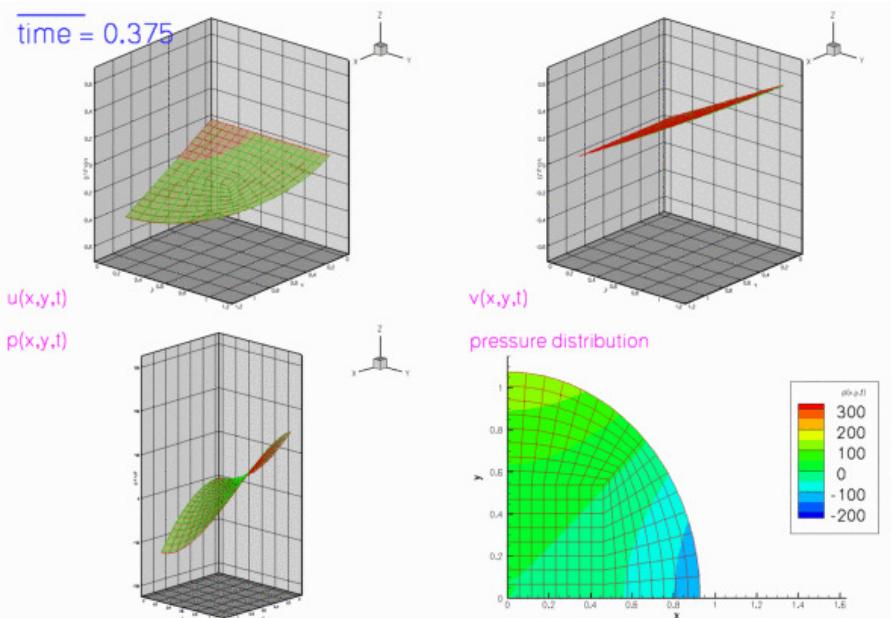


Figure 1.2 Plot of the velocity and pressure fields computed with 2D Crouzeix-Raviart elements, with $Re=100$ and $St=1$.

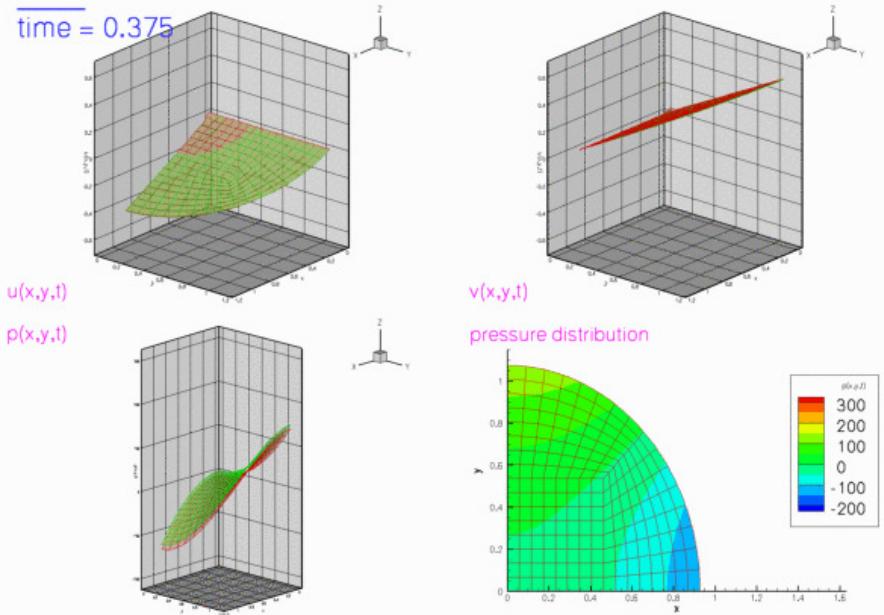


Figure 1.3 Plot of the velocity and pressure fields computed with 2D Taylor-Hood elements, with $\text{Re}=100$ and $\text{St}=1$.

1.3 The moving wall

As usual, we represent the moving wall as a `GeomObject` and define its shape by implementing the pure virtual function `GeomObject::position(...)`. The arguments to the constructor specify the mean half-axis of the ellipse, A , the amplitude of its variations, \hat{A} , and the period of the oscillation, T . We also pass the pointer to a `Time` object to the constructor and store it in a private data member, to allow the `position(...)` functions to access the current value of the continuous time.

```
=====start_of_MyEllipse=====
/// Oscillating ellipse
/// \f[ x = (A + \widehat{A} \cos(2\pi t/T)) \cos(\xi_i) \f]
/// \f[ y = \frac{\sin(\xi_i)}{A + \widehat{A} \cos(2\pi t/T)} \f]
/// Note that cross-sectional area is conserved.
=====
class MyEllipse : public GeomObject
{
public:
    /// Constructor: Pass initial x-half axis, amplitude of x-variation,
    /// period of oscillation and pointer to time object.
    MyEllipse(const double& a, const double& a_hat,
              const double& period, Time* time_pt) :
        GeomObject(1,2), A(a), A_hat(a_hat), T(period), Time_pt(time_pt) {}

    /// Destructor: Empty
    virtual ~MyEllipse() {}

    /// Current position vector to material point at
    /// Lagrangian coordinate xi
    void position(const Vector<double>& xi, Vector<double>& r) const
    {
        // Get current time:
        double time=Time_pt->time();

        // Position vector
        double axis=A+A_hat*cos(2.0*MathematicalConstants::Pi*time/T);
        r[0] = axis*cos(xi[0]);
        r[1] = (1.0/axis)*sin(xi[0]);
    }

    /// Parametrised position on object: r(xi). Evaluated at
    /// previous time level. t=0: current time; t>0: previous
    /// time level.
    void position(const unsigned& t, const Vector<double>& xi,
                  Vector<double>& r) const
    {
        // Get current time:
        double time=Time_pt->time(t);
```

```

// Position vector
double axis=A+A_hat*cos(2.0*MathematicalConstants::Pi*time/T);
r[0] = axis*cos(xi[0]);
r[1] = (1.0/axis)*sin(xi[0]);
}

private:

/// x-half axis
double A;

/// Amplitude of variation in x-half axis
double A_hat;

/// Period of oscillation
double T;

/// Pointer to time object
Time* Time_pt;

}; // end of MyEllipse

```

1.4 The global parameters

As in most previous examples, we use a namespace to define and initialise global problem parameters such as the Reynolds and Strouhal numbers:

```

=====start_of_namespace=====
// Namespace for global parameters
=====end_of_namespace=====
namespace Global_Physical_Variables
{

/// Reynolds number
double Re=100.0;

/// Womersley = Reynolds times Strouhal
double ReSt=100.0;

```

We also define and initialise the parameters that specify the motion of the domain boundary and specify the exact solution.

```

/// x-Half axis length
double A=1.0;

/// x-Half axis amplitude
double A_hat=0.1;

/// Period of oscillations
double T=1.0;

/// Exact solution of the problem as a vector containing u,v,p
void get_exact_u(const double& t, const Vector<double>& x, Vector<double>& u)
{
    using namespace MathematicalConstants;

    // Strouhal number
    double St = ReSt/Re;

    // Half axis
    double a=A+A_hat*cos(2.0*Pi*t/T);
    double adot=-2.0*A_hat*Pi*sin(2.0*Pi*t/T)/T;
    u.resize(3);

    // Velocity solution
    u[0]=adot*x[0]/a;
    u[1]=-adot*x[1]/a;

    // Pressure solution
    u[2]=(2.0*A_hat*Pi*Pi*Re*(x[0]*xx[0]*St*cos(2.0*Pi*t/T)*A +
        x[0]*xx[0]*St*A_hat - x[0]*x[0]*A_hat +
        x[0]*xx[0]*A_hat*cos(2.0*Pi*t/T)*cos(2.0*Pi*t/T) -
        x[1]*x[1]*St*cos(2.0*Pi*t/T)*A -
        x[1]*x[1]*St*A_hat - x[1]*x[1]*A_hat +
        x[1]*x[1]*A_hat*cos(2.0*Pi*t/T)*cos(2.0*Pi*t/T) ))
    / (T*T*(A*A + 2.0*A*A_hat*cos(2.0*Pi*t/T) +
        A_hat*A_hat*cos(2.0*Pi*t/T)*cos(2.0*Pi*t/T) ));
}

} // end of namespace

```

1.5 The driver code

As in most previous unsteady demo codes, we allow the code to be run in a validation mode (in which we use a coarser mesh and execute fewer timesteps). This mode is selected by specifying an (arbitrary) command line argument that we store in the namespace `CommandLineArgs`.

```
//=====start_of_main=====

/// Driver code for unsteady Navier-Stokes flow, driven by
/// oscillating ellipse. If the code is executed with command line
/// arguments, a validation run is performed.
//=====

int main(int argc, char* argv[])
{
```

```
    // Store command line arguments
    CommandLineArgs::setup(argc,argv);

    // Solve with Crouzeix-Raviart elements
    {
        // Create DocInfo object with suitable directory name for output
        DocInfo doc_info;
        doc_info.set_directory("RESLT_CR");

        //Set up problem
        OscEllipseProblem<RefineableQCrouzeixRaviartElement<2>,BDF<2> > problem;

        // Run the unsteady simulation
        problem.unsteady_run(doc_info);
    }
```

Then we repeat this process for adaptive Taylor-Hood elements.

```
    // Solve with Taylor-Hood elements
    {
        // Create DocInfo object with suitable directory name for output
        DocInfo doc_info;
        doc_info.set_directory("RESLT_TH");

        //Set up problem
        OscEllipseProblem<RefineableQTaylorHoodElement<2>,BDF<2> > problem;

        // Run the unsteady simulation
        problem.unsteady_run(doc_info);
    }

}; // end of main
```

1.6 The problem class

Most of the problem class is a straightforward combination of the problem classes employed in the simulation of the [adaptive driven cavity](#) and [Rayleigh channel](#) problems, in that the problem combines unsteadiness with spatial adaptivity (though in the present problem the adaptivity is only used to uniformly refine the very coarse base mesh; we refer to [another example](#) for the use of full spatial adaptivity in a moving-domain Navier-Stokes problem).

```
//=====start_of_problem_class=====

/// Navier-Stokes problem in an oscillating ellipse domain.
//=====

template<class ELEMENT, class TIMESTEPPER>
class OscEllipseProblem : public Problem
{

public:

    /// Constructor
    OscEllipseProblem();

    /// Destructor (empty)
    ~OscEllipseProblem() {}

    /// Update the problem specs after solve (empty)
    void actions_after_newton_solve() {}

    /// Update problem specs before solve (empty)
    void actions_before_newton_solve() {}

    /// Actions before adapt (empty)
    void actions_before_adapt() {}

    /// Actions after adaptation, pin relevant pressures
```

```

void actions_after_adapt()
{
    // Unpin all pressure dofs
    RefineableNavierStokesEquations<2>::unpin_all_pressure_dofs(mesh_pt() -> element_pt());

    // Pin redundant pressure dofs
    RefineableNavierStokesEquations<2>::pin_redundant_nodal_pressures(mesh_pt() -> element_pt());

    // Now set the first pressure dof in the first element to 0.0
    fix_pressure(0, 0, 0.0);

} // end of actions_after_adapt

```

The key new feature in the current problem is the presence of the moving domain which requires updates of

1. all nodal positions
2. the prescribed velocities on the moving wall via the no-slip condition.

before every timestep. Since the nodal positions of the QuarterCircleSectorMesh are determined via its MacroElement / Domain representation (which updates the nodal position in response to changes in the geometry of the GeomObjects that define its boundaries), the former task may be accomplished by executing the Mesh::node_update() function; the update of the no-slip condition may be performed by calling the function FSI_functions::apply_no_slip_on_moving_wall(Node* node_pt), a helper function, defined in the namespace FSI_functions, which updates the velocity components $u_1, u_2, [u_3]$ according to the no-slip boundary condition

$$\mathbf{u}_{Node} = \frac{\partial \mathbf{x}_{Node}}{\partial t}$$

where the time-derivative of the nodal positions is evaluated by the Node's positional timestepper. [Note: The function FSI_functions::apply_no_slip_on_moving_wall(...) assumes that the velocity components are stored in the Node's first 2 [3] values. This is consistent with the storage of the velocity component in all existing Navier-Stokes elements. If you develop your own Navier-Stokes elements and use a different storage scheme you use this function at your own risk.]

Here is the implementation of these tasks:

```

/// Update the problem specs before next timestep
void actions_before_implicit_timestep()
{
    // Update the domain shape
    mesh_pt() -> node_update();

    // Ring boundary: No slip; this implies that the velocity needs
    // to be updated in response to wall motion
    unsigned ibound=1;
    unsigned num_nod=mesh_pt() -> nboundary_node(ibound);
    for (unsigned inod=0; inod<num_nod; inod++)
    {
        // Which node are we dealing with?
        Node* node_pt=mesh_pt() -> boundary_node_pt(ibound, inod);

        // Apply no slip
        FSI_functions::apply_no_slip_on_moving_wall(node_pt);
    }
}

/// Update the problem specs after timestep (empty)
void actions_after_implicit_timestep(){}

```

The remaining functions are similar to those used in our previous Navier-Stokes examples and require no further explanation.

```

/// Doc the solution
void doc_solution(DocInfo& doc_info);

/// Timestepping loop
void unsteady_run(DocInfo& doc_info);

/// Set initial condition
void set_initial_condition();

private:

    /// Fix pressure in element e at pressure dof pdof and set to pvalue
    void fix_pressure(const unsigned &e, const unsigned &pdof,
                     const double &pvalue)
{

```

```

//Cast to proper element and fix pressure
dynamic_cast<ELEMENT*>(mesh_pt() ->element_pt(e)) ->
    fix_pressure(pdof,pvalue);
} // end_of_fix_pressure

/// Pointer to GeomObject that specifies the domain boundary
GeomObject* Wall_pt;

}; // end of problem class

```

1.7 The problem constructor

We start by creating a timestepper of the type specified by the Problem's template parameter and add (a pointer to) it to the Problem's collection of Timesteppers. Recall that this function also creates the Problem's Time object.

```

//=====start_of_constructor=====
// Constructor for Navier-Stokes problem on an oscillating ellipse domain.
//=====
template<class ELEMENT, class TIMESTEPPER>
OscEllipseProblem<ELEMENT,TIMESTEPPER>::OscEllipseProblem()
{
    //Create the timestepper and add it to the problem
    add_time stepper_pt(new TIMESTEPPER);
}

// Build geometric object that forms the curvilinear domain boundary:
// an oscillating ellipse

// Half axes
double a=Global_Physical_Variables::A;

// Variations of half axes
double a_hat=Global_Physical_Variables::A_hat;

// Period of the oscillation
double period=Global_Physical_Variables::T;

// Create GeomObject that specifies the domain boundary
Wall_pt=new MyEllipse(a,a_hat,period,Problem::time_pt());

// Start and end coordinates of curvilinear domain boundary on ellipse
double xi_lo=0.0;
double xi_hi=MathematicalConstants::Pi/2.0;

// Now create the mesh. Separating line between the two
// elements next to the curvilinear boundary is located half-way
// along the boundary.
double fract_mid=0.5;
Problem::mesh_pt() = new RefineableQuarterCircleSectorMesh<ELEMENT>(
    Wall_pt,xi_lo,fract_mid,xi_hi,time stepper_pt());

```

Both velocity components on the curvilinear mesh boundary are determined by the no-slip condition and must therefore be pinned,

```

// Fluid boundary conditions
//-----
// Ring boundary: No slip; this also implies that the velocity needs
// to be updated in response to wall motion
unsigned ibound=1;
{
    unsigned num_nod= mesh_pt() ->nboundary_node(ibound);
    for (unsigned inod=0;inod<num_nod;inod++)
    {
        // Pin both velocities
        for (unsigned i=0;i<2;i++)
        {
            mesh_pt() ->boundary_node_pt(ibound,inod) ->pin(i);

```

```

        }
    } // end boundary 1
whereas on the symmetry boundaries only one of the two velocity components is set to zero:
// Bottom boundary:
ibound=0;
{
    unsigned num_nod= mesh_pt()->nboundary_node(ibound);
    for (unsigned inod=0;inod<num_nod;inod++)
    {
        // Pin vertical velocity
        {
            mesh_pt()->boundary_node_pt(ibound,inod)->pin(1);
        }
    }
} // end boundary 0

// Left boundary:
ibound=2;
{
    unsigned num_nod= mesh_pt()->nboundary_node(ibound);
    for (unsigned inod=0;inod<num_nod;inod++)
    {
        // Pin horizontal velocity
        {
            mesh_pt()->boundary_node_pt(ibound,inod)->pin(0);
        }
    }
} // end boundary 2

```

Finally, we pass the pointers to *Re*, *Re St* and the global Time object (automatically created by the Problem when the timestepper was passed to it at the beginning of the constructor) to the elements, pin the redundant nodal pressure degrees of freedom (see the discussion of the [adaptive driven-cavity problem](#) for more details), pin one pressure degree of freedom, and set up the equation numbering scheme.

```

// Complete the build of all elements so they are fully functional
//-----
// Find number of elements in mesh
unsigned n_element = mesh_pt()->nelement();

// Loop over the elements to set up element-specific
// things that cannot be handled by constructor
for(unsigned i=0;i<n_element;i++)
{
    // Upcast from FiniteElement to the present element
    ELEMENT *el_pt = dynamic_cast<ELEMENT*>(mesh_pt()->element_pt(i));

    //Set the Reynolds number, etc
    el_pt->re_pt() = &Global_Physical_Variables::Re;
    el_pt->re_st_pt() = &Global_Physical_Variables::ReSt;
}

// Pin redundant pressure dofs
RefineableNavierStokesEquations<2>::
pin_redundant_nodal_pressures(mesh_pt()->element_pt());
// Now set the first pressure dof in the first element to 0.0
fix_pressure(0,0,0.0);

// Do equation numbering
cout << "Number of equations: " << assign_eqn_numbers() << std::endl;

} // end of constructor

```

1.8 Assigning the initial conditions

This function assigns "history values" for the velocities and the nodal positions from the exact solution. It is implemented in exactly the same way as in the solution of the [unsteady heat equation in a moving domain](#). Note that because the domain is moving, the nodal positions must be updated (according to the position of the domain boundary at the relevant previous timestep), before evaluating the exact solution at the nodal position.

```

=====start_of_set_initial_condition=====
/// Set initial condition: Assign previous and current values
/// from exact solution.
=====template<class ELEMENT, class TIMESTEPPER>
void OscEllipseProblem<ELEMENT,TIMESTEPPER>::set_initial_condition()
{
    // Backup time in global timestepper
    double backed_up_time=time_pt()->time();

    // Past history for velocities must be established for t=time0-deltat, ...
    // Then provide current values (at t=time0) which will also form

```

```

// the initial guess for first solve at t=time0+delta_t
// Vector of exact solution value
Vector<double> soln(3);
Vector<double> x(2);
//Find number of nodes in mesh
unsigned num_nod = mesh_pt()->nnode();
// Get continuous times at previous timesteps
int nprev_steps=time stepper_pt()->nprev_values();
Vector<double> prev_time(nprev_steps+1);
for (int itime=nprev_steps;itime>=0;itime--)
{
    prev_time[itime]=time_pt()->time(unsigned(itime));
}
// Loop over current & previous timesteps (in outer loop because
// the mesh also moves!)
for (int itime=nprev_steps;itime>=0;itime--)
{
    double time=prev_time[itime];

    // Set global time (because this is how the geometric object refers
    // to continous time )
    time_pt()->time()=time;

    cout << "setting IC at time =" << time << std::endl;

    // Update the mesh for this value of the continuous time
    // (The wall object reads the continous time from the
    // global time object)
    mesh_pt()->node_update();

    // Loop over the nodes to set initial guess everywhere
    for (unsigned jnod=0;jnod<num_nod;jnod++)
    {
        // Get nodal coordinates
        x[0]=mesh_pt()->node_pt(jnod)->x(0);
        x[1]=mesh_pt()->node_pt(jnod)->x(1);

        // Get exact solution (unsteady stagnation point flow)
        Global_Physical_Variables::get_exact_u(time,x,soln);

        // Assign solution
        mesh_pt()->node_pt(jnod)->set_value(itime,0,soln[0]);
        mesh_pt()->node_pt(jnod)->set_value(itime,1,soln[1]);

        // Loop over coordinate directions
        for (unsigned i=0;i<2;i++)
        {
            mesh_pt()->node_pt(jnod)->x(itime,i)=x[i];
        }
    }
    // end of loop over previous timesteps
    // Reset backed up time for global timestepper
    time_pt()->time()=backed_up_time;
} // end of set initial condition

```

1.9 Post processing

The function `doc_solution(...)` is similar to that in the `unsteady heat examples` and the previous Navier-Stokes examples. We add dummy zones and tecplot geometries to facilitate the post-processing of the results with tecplot.

```

//=====start_of_doc_solution=====
// Doc the solution
//=====
template<class ELEMENT, class TIMESTEPPER>
void OscEllipseProblem<ELEMENT,TIMESTEPPER>::doc_solution(DocInfo& doc_info)
{
    ofstream some_file;
    char filename[100];

    // Number of plot points
    unsigned npts;
    npts=5;

    // Output solution
    //-----
    snprintf(filename, sizeof(filename), "%s/soln%i.dat", doc_info.directory().c_str(),
             doc_info.number());
    some_file.open(filename);
    mesh_pt()->output(some_file,npts);
    some_file << "TEXT X=2.5,Y=93.6,F=HELV,HU=POINT,C=BLUE,H=26,T=\"time = "
               << time_pt()->time() << "\"";
    some_file << "GEOMETRY X=2.5,Y=98,T=LINE,C=BLUE,LT=0.4" << std::endl;
    some_file << "1" << std::endl;
    some_file << "2" << std::endl;

```

```

some_file << " 0 0" << std::endl;
some_file << time_pt()->time()*20.0 << " 0" << std::endl;

// Write dummy zones that force tecplot to keep the axis limits constant
// while the domain is moving.
some_file << "ZONE I=2,J=2" << std::endl;
some_file << "0.0 0.0 -0.65 -0.65 -200.0" << std::endl;
some_file << "1.15 0.0 -0.65 -0.65 -200.0" << std::endl;
some_file << "0.0 1.15 -0.65 -0.65 -200.0" << std::endl;
some_file << "1.15 1.15 -0.65 -0.65 -200.0" << std::endl;
some_file << "ZONE I=2,J=2" << std::endl;
some_file << "0.0 0.0 0.65 300.0" << std::endl;
some_file << "1.15 0.0 0.65 300.0" << std::endl;
some_file << "0.0 1.15 0.65 300.0" << std::endl;
some_file << "1.15 1.15 0.65 300.0" << std::endl;

some_file.close();

// Output exact solution
//-----
snprintf(filename, sizeof(filename), "%s/exact_soln%i.dat", doc_info.directory().c_str(),
         doc_info.number());
some_file.open(filename);
mesh_pt()->output_fct(some_file,npts,time_pt()->time(),
                        Global_Physical_Variables::get_exact_u);
some_file.close();

// Doc error
//-----
double error,norm;
snprintf(filename, sizeof(filename), "%s/error%i.dat", doc_info.directory().c_str(),
         doc_info.number());
some_file.open(filename);
mesh_pt()->compute_error(some_file,
                           Global_Physical_Variables::get_exact_u,
                           time_pt()->time(),
                           error,norm);
some_file.close();

// Doc solution and error
//-----
cout << "error: " << error << std::endl;
cout << "norm : " << norm << std::endl << std::endl;

// Plot wall posn
//-----
snprintf(filename, sizeof(filename), "%s/Wall%i.dat", doc_info.directory().c_str(),
         doc_info.number());
some_file.open(filename);
unsigned nplot=100;
for (unsigned iplot=0;iplot<nplot;iplot++)
{
    Vector<double> xi_wall(1), r_wall(2);
    xi_wall[0]=0.5*MathematicalConstants::Pi*double(iplot)/double(nplot-1);
    Wall_pt->position(xi_wall,r_wall);
    some_file << r_wall[0] << " " << r_wall[1] << std::endl;
}
some_file.close();
// Increment number of doc
doc_info.number()++;

} // end of doc_solution

```

1.10 The timestepping loop

The timestepping loop is extremely straightforward: We choose a timestep and the overall length of the simulation, initialise the timestepper(s) by calling `Problem::initialise_dt(...)` and assign the initial condition.

```

//=====start_of_unsteady_run=====
// Unsteady run
//=====
template<class ELEMENT, class TIMESTEPPER>
void OscEllipseProblem<ELEMENT,TIMESTEPPER>::unsteady_run(DocInfo& doc_info)
{

    // Specify duration of the simulation
    double t_max=3.0;

    // Initial timestep
    double dt=0.025;

    // Initialise timestep

```

```

initialise_dt(dt);

// Set initial conditions.
set_initial_condition();

```

Next we set the number of timesteps for a normal run.

```

// Alternative initial conditions: impulsive start; see exercise.
//assign_initial_values_impulsive();

// find number of steps
unsigned nstep = unsigned(t_max/dt);

```

We over-write this number and perform a single uniform mesh refinement if the code is run in self-test mode (indicated by a non-zero number of command line arguments),

```

// If validation: Reduce number of timesteps performed and
// use coarse-ish mesh
if (CommandLineArgs::Argc>1)
{
    nstep=2;
    refine_uniformly();
    cout << "validation run" << std::endl;
}

```

otherwise we refine the mesh three times and output the initial conditions

```

else
{
    // Refine the mesh three times, to resolve the pressure distribution
    // (the velocities could be represented accurately on a much coarser mesh).
    refine_uniformly();
    refine_uniformly();
    refine_uniformly();
}

// Output solution initial
doc_solution(doc_info);

```

Finally we execute the proper timestepping loop and document the solution after every timestep

```

// Timestepping loop
for (unsigned istep=0;istep<nstep;istep++)
{
    cout << "Timestep " << istep << std::endl;
    cout << "Time is now " << time_pt()>time() << std::endl;

    // Take timestep
    unsteady_newton_solve(dt);

    //Output solution
    doc_solution(doc_info);
}

} // end of unsteady run

```

1.11 Comments and Exercises

1. Compare the results of the numerical simulation in which \mathbf{u}_{IC} is given by the exact solution (an unsteady stagnation point flow) to that obtained from an "impulsive start" where $\mathbf{u}_{IC} = \mathbf{0}$. (This is most easily implemented by replacing the call to `set_initial_condition()` with a call to `Problem::assign_initial_values_impulsive()`.)

Why do we obtain the same velocity with both initial conditions and why does the pressure take a few timesteps (How many exactly? Compare simulations with `BDF<4>` and `BDF<2>` timesteppers.) to "catch up" with the exact solution? [Hint: The unsteady stagnation point flow is a potential flow, therefore the viscous terms in the Navier-Stokes equations disappear. See also chapter 3.19 in Volume 2 of Gresho & Sani's wonderful book "Incompressible Flow and the Finite Element Method".]

1.12 Source files for this tutorial

- The source files for this tutorial are located in the directory:

```
demo_drivers/navier_stokes/osc_ellipse/
```

- The driver code is:

```
demo_drivers/navier_stokes/osc_ellipse/osc_quarter_ellipse.cc
```

1.13 PDF file

A [pdf version](#) of this document is available. \