

DOLFIN

Dynamic Object-oriented Library for FINite element computation

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Chalmers Finite Element Center

Introduction

- An adaptive finite element solver for PDEs
- Written by people at the Department of Computational Mathematics (Hoffman/Logg)
- Written in C++
- Highly modularized
- DOLFIN is a solver. No grid generation. No visualisation.
- Licensed under the GNU GPL
- `http://www.phy.chalmers.se/dolfin`

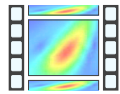
GNU and the GPL

- Makes the software free for all users
- “Free” as in “free speech”, not “free beer”
- Free to modify, change, copy, redistribute
- Derived work must also use the GPL license
- Enables sharing of code
- Simplifies distribution of the program
- Linux is distributed under the GPL license
- See <http://www.gnu.org>

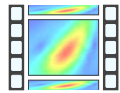
Features

- 3D or 2D
- Automatic assembling
- Tetrahedrons or triangles
- Linear elements
- Algebraic solvers: LU, GMRES, CG
- Missing: mesh refinement, adaptivity, multi-grid

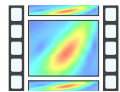
DOLFIN examples



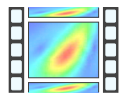
Start movie 1 (driven cavity, solution)



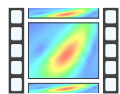
Start movie 2 (driven cavity, dual)



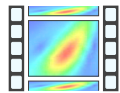
Start movie 3 (driven cavity, dual)



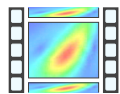
Start movie 4 (bluff body, solution)



Start movie 5 (bluff body, dual)

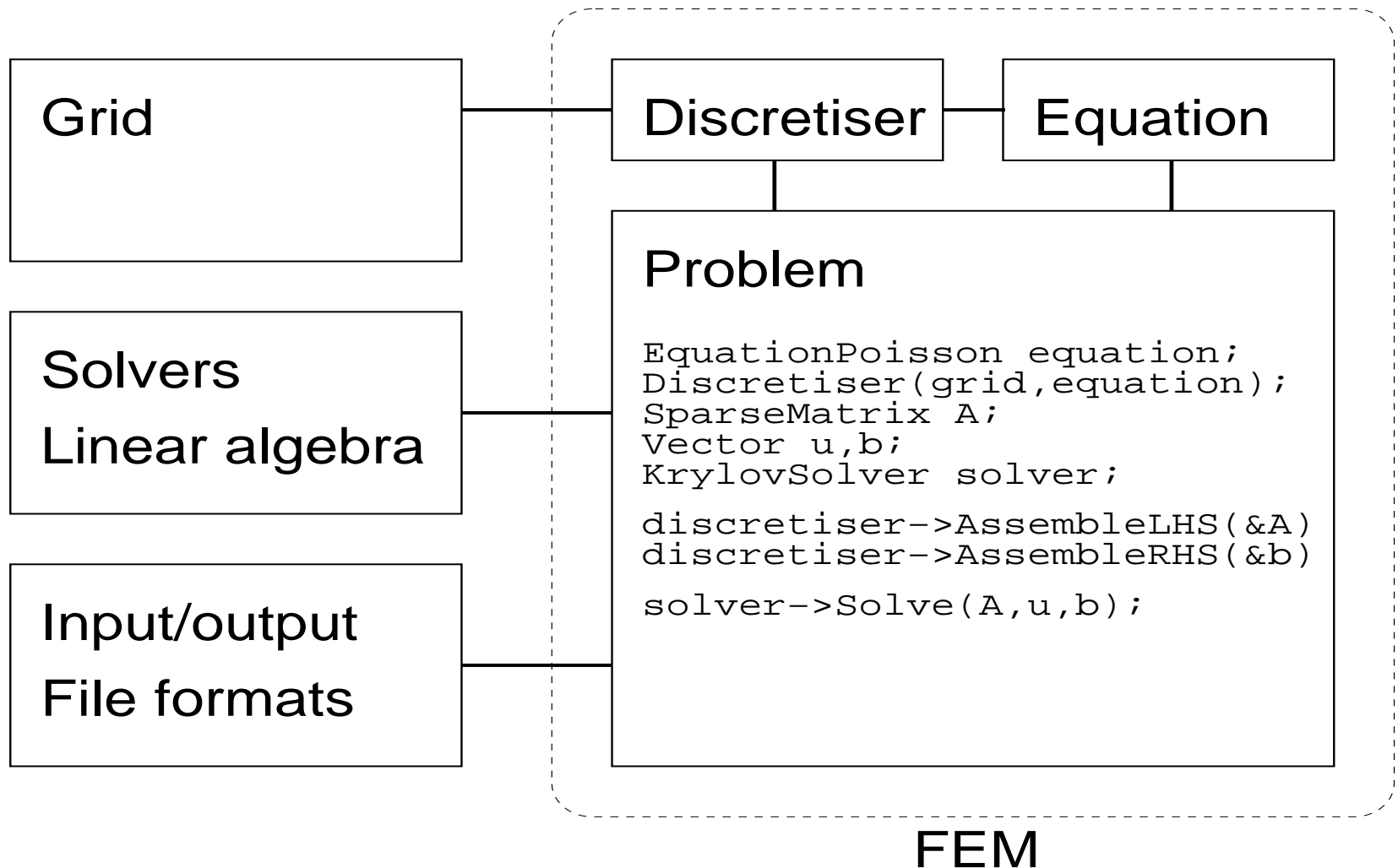


Start movie 6 (jet, solution)



Start movie 7 (transition to turbulence)

Internal structure



The finite element

Following Ciarlet and Brenner-Scott a *finite element* $(K, \mathcal{P}, \mathcal{N})$ is defined by

1. A domain $K \subseteq \mathbb{R}^N$ with piecewise smooth boundary, typically a triangle or a tetrahedron;
2. A finite-dimensional space \mathcal{P} of functions on K together with a set of basis functions (the *shape functions*);
3. A basis $\mathcal{N} = \{N_1, N_2, \dots, N_k\}$ for the dual space \mathcal{P}^* (the *degrees of freedom*).

The finite element: implementation

<code>FiniteElement</code>	The <i>finite element</i> , containing geometry and the local function space.
<code>FunctionSpace</code>	A finite-dimensional space of functions on the domain of a finite element.
<code>TriLinSpace</code> , <code>TetLinSpace</code>	Sub-classes derived from <code>FunctionSpace</code> .
<code>ShapeFunction</code>	Member of the basis for a local function space.
<code>TriLinFunction</code> , <code>TetLinFunction</code>	Sub-classes derived from <code>ShapeFunction</code> .
<code>LocalField</code>	A member of the local function space on the domain of a finite element, i.e. a linear combination of shape functions. Can also be viewed as the restriction of a <code>GlobalField</code> to the domain of a finite element.
<code>GlobalField</code>	A function (possibly vector-valued) defined on the whole of the computational domain.

Automatic assembling

- Automatic assembling is handled by *operator overloading*.
- The symmetric binary operator '*' is defined for the following classes:

* : ShapeFunction × ShapeFunction → real

* : ShapeFunction × real → real

* : ShapeFunction × LocalField → real

* : LocalField × LocalField → real

* : LocalField × real → real

Automatic assembling

For two ShapeFunctions `v` and `w`, representing two shape functions v and w on K , the operator `'*'` is defined by

$$v * w = \frac{1}{|K|} \int_K v \cdot w \, dx, \quad (1)$$

$$a * v = \frac{1}{|K|} \int_K \alpha \cdot v \, dx, \quad (2)$$

where $|K|$ is the volume (area) of the domain K and `a` represents the real number α .

Automatic assembling

$$u - \nabla \cdot (c \nabla u) = f$$

$$\int_{\Omega} \left(\frac{U^n - U^{n-1}}{k_n} \right) v \, dx + \int_{\Omega} c(\cdot, t_n) \nabla U^n \cdot \nabla v \, dx = \int_{\Omega} f(\cdot, t_n) v \, dx$$

$$(u^*v - u_p^*v) / k + c^* (u \cdot dx^*v \cdot dx + u \cdot dy^*v \cdot dy + u \cdot dz^*v \cdot dz)$$

$$f^*v$$

Automatic assembling

Exact evaluation of integrals for linear elements on triangles and tetrahedrons:

$$\frac{1}{|K|} \int_K \lambda_1^{m_1} \lambda_2^{m_2} \lambda_3^{m_3} dx = \frac{2 \cdot m_1! m_2! m_3!}{(m_1 + m_2 + m_3 + 2)!}$$

$$\mathbf{a} * \mathbf{v} = \frac{1}{|K|} \int_K \alpha \lambda_i dx = \alpha/3$$

$$\frac{1}{|K|} \int_K \lambda_1^{m_1} \lambda_2^{m_2} \lambda_3^{m_3} \lambda_4^{m_4} dx = \frac{3! \cdot m_1! m_2! m_3! m_4!}{(m_1 + m_2 + m_3 + m_4 + 3)!}$$

$$\mathbf{a} * \mathbf{v} = \frac{1}{|K|} \int_K \alpha \lambda_i dx = \alpha/4$$

Three levels

- Simple C/C++ interface for the *user* who just wants to solve an equation with specified geometry and boundary conditions.
- New algorithms are added at *module level* by the developer or advanced user.
- Core features are added at *kernel level*.

Poisson's equation: user level

```
#include <dolfin.h>
int main(int argc, char **argv)
{
    dolfin_set_problem("poisson");

    dolfin_set_parameter("output file", "poisson.dx");
    dolfin_set_parameter("grid file", "tetgrid.inp");
    dolfin_set_parameter("space dimension", 2)

    dolfin_set_boundary_conditions(my_bc);
    dolfin_set_function("source", f);

    dolfin_init(argc, argv);
    dolfin_solve();
    dolfin_end();

    return 0;
}
```

Poisson's equation: module level

```
class EquationPoisson: public Equation{
public:
    EquationPoisson():Equation(3){
        AllocateFields(1);
        field[0] = &f;
    }

    real IntegrateLHS(ShapeFunction &u, ShapeFunction &v){
        return ( u.dx*v.dx + u.dy*v.dy + u.dz*v.dz );
    }

    real IntegrateRHS(ShapeFunction &v){
        return ( f * v );
    }

private:
    LocalField f;
};
```

Poisson's equation: module level

```
void ProblemPoisson::Solve()
{
    EquationPoisson equation;
    SparseMatrix A;
    Vector x,b;
    KrylovSolver solver;
    GlobalField u(grid,&x);
    GlobalField f(grid,"source");
    Discretiser discretiser(grid,equation);

    equation.AttachField(0,&f);
    discretiser.Assemble(&A,&b);
    solver.Solve(&A,&x,&b);

    u.SetLabel("u","temperature");
    u.Save();
}
```


Poisson's equation: kernel level

```
class Equation{
public:

    Equation(int nsd);
    ~Equation();

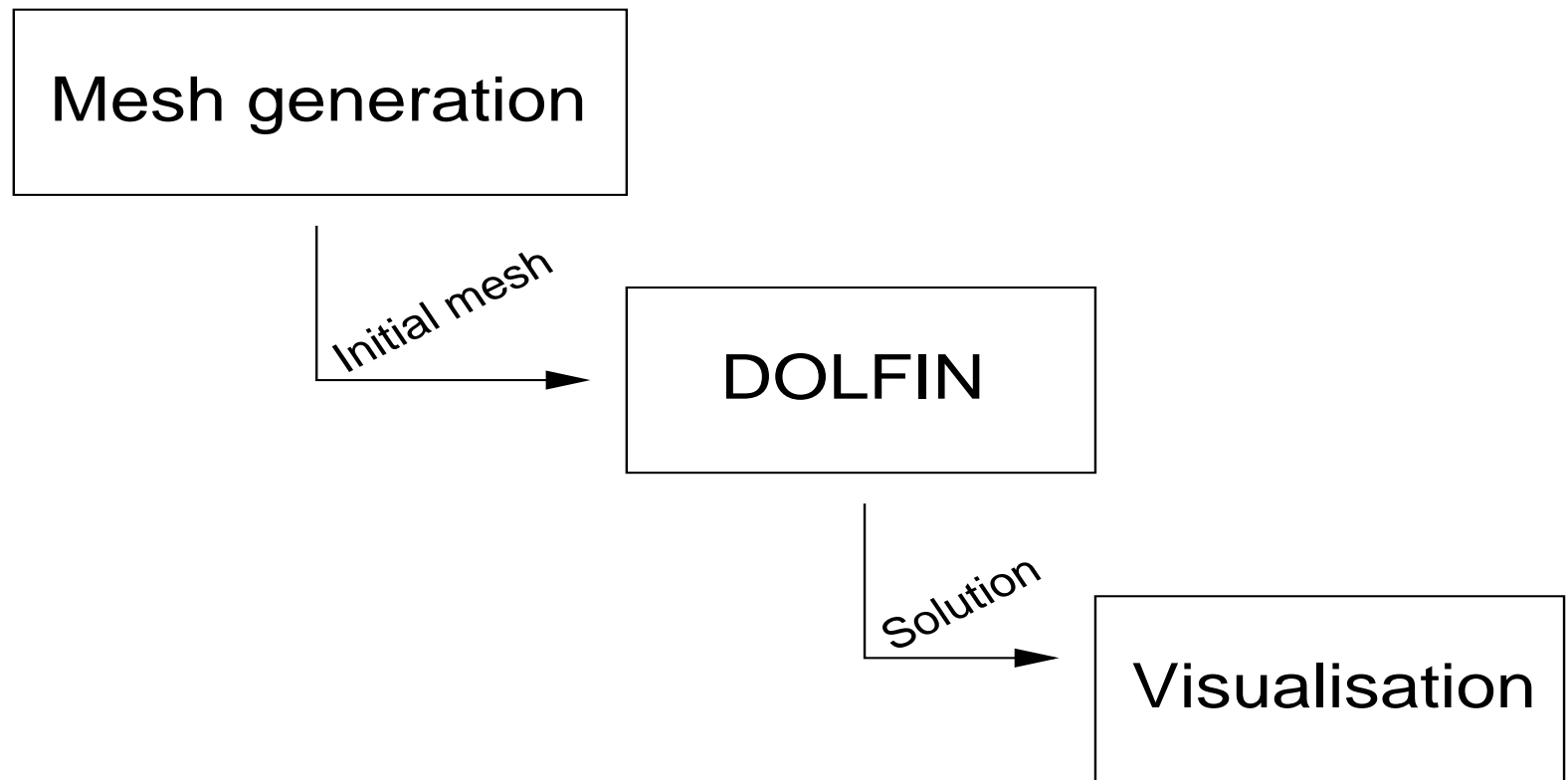
    virtual real IntegrateLHS(ShapeFunction &u, ShapeFunction &v)
    virtual real IntegrateRHS(ShapeFunction &v) = 0;

    void UpdateLHS(FiniteElement *element);
    void UpdateRHS(FiniteElement *element);
    void AttachField(int i, GlobalField *globalfield);

    ...
}
```

Input / output

The solver (DOLFIN) is the key part in the larger system containing also pre- and post-processing:



Input / output

- OpenDX: free open-source visualisation program based on IBM:s *Visualization Data Explorer*.
- MATLAB: commercial software (2000 Euros)
- GiD: commercial software (570 Euros)

Input / output: examples

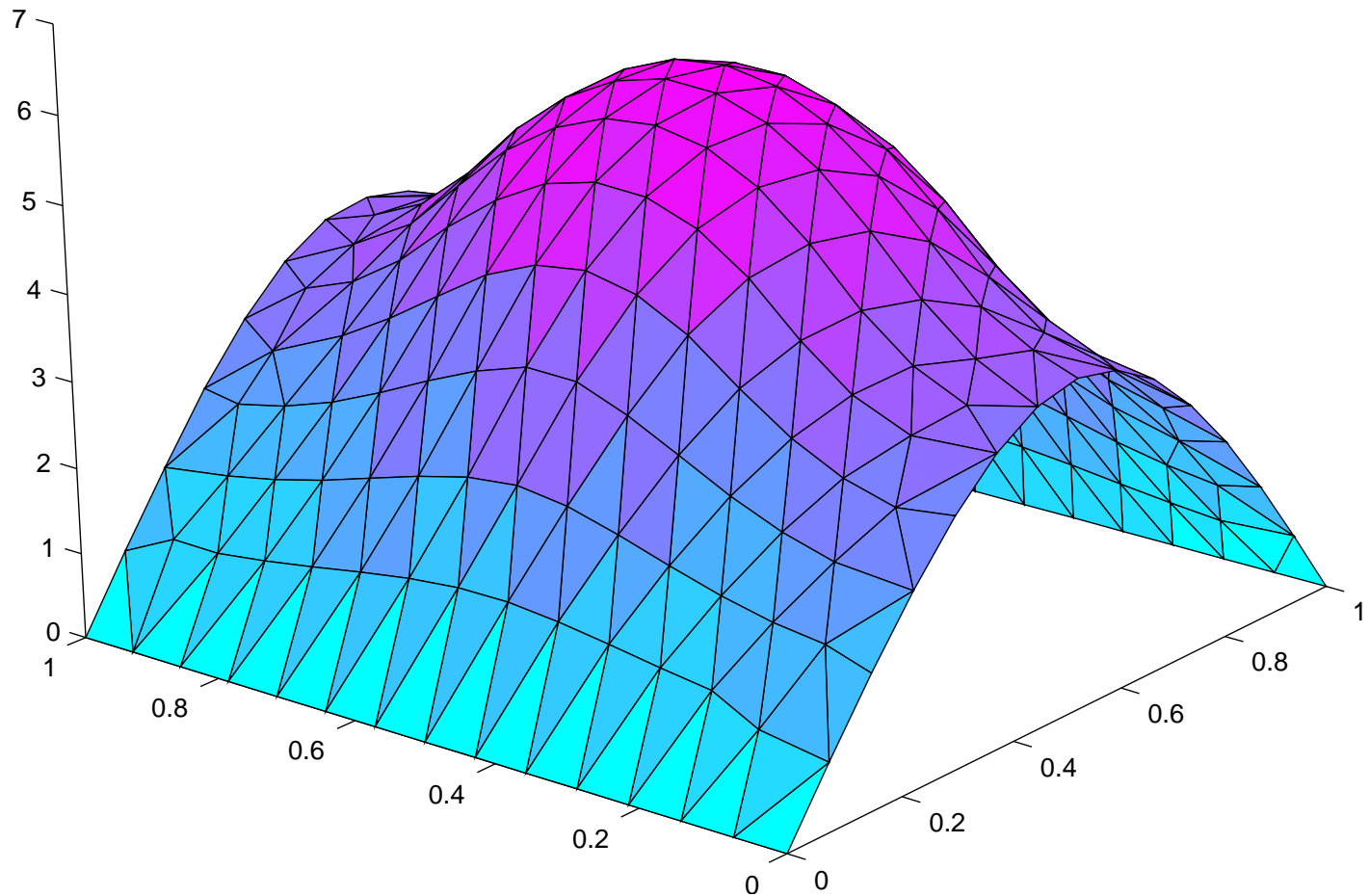
Poisson's equation:

$$-\Delta u(x) = f(x), \quad x \in \Omega, \quad (3)$$

on the unit square $\Omega = (0, 1) \times (0, 1)$ with the source term f localised to the middle of the domain.

Grid generation with **GiD** and visualisation using the `pdesurf` command in **MATLAB**.

Input / output: examples



Input / output: examples

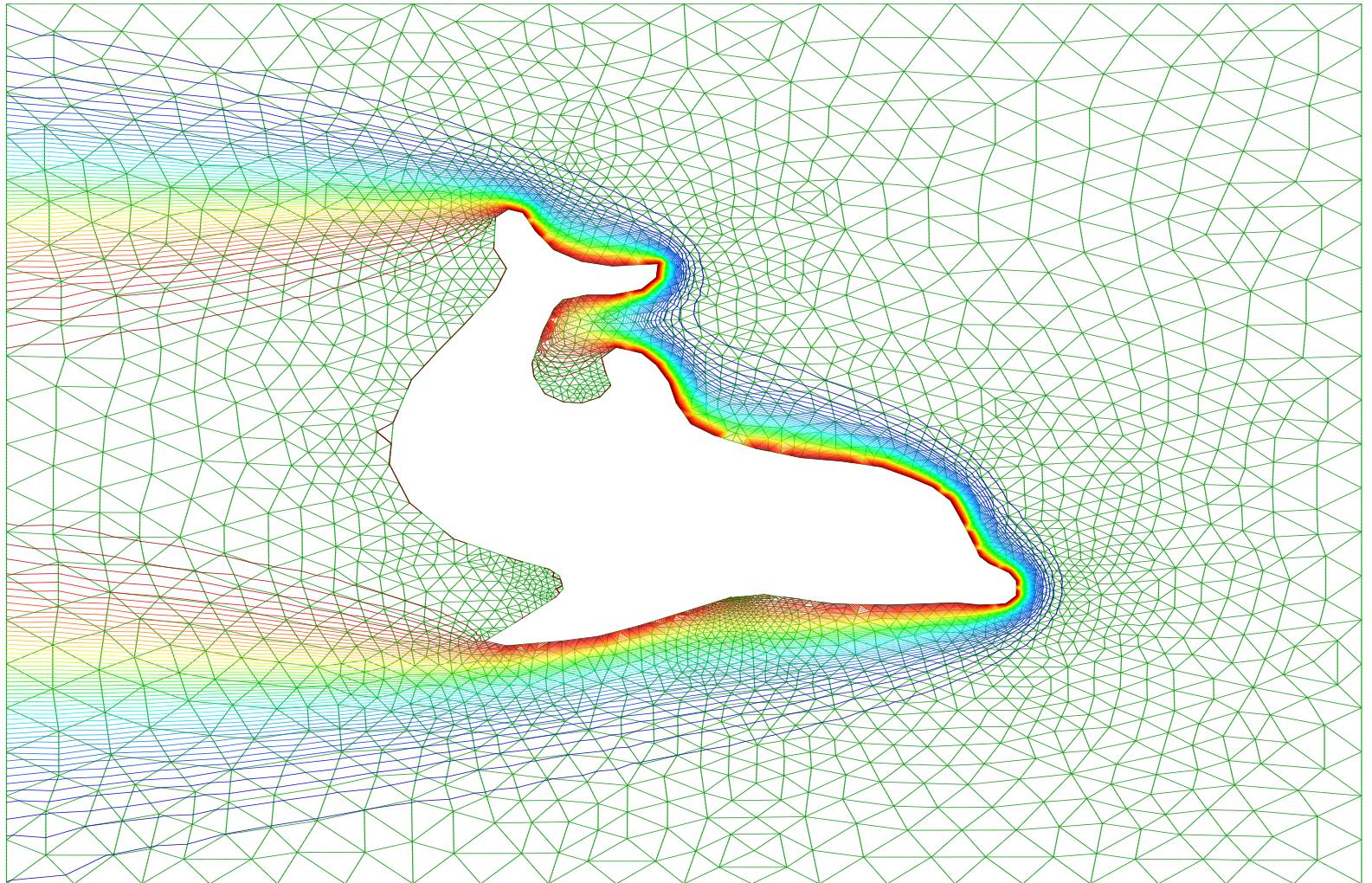
Convection–diffusion:

$$\dot{u} + b \cdot \nabla u - \nabla \cdot (\epsilon \nabla u) = f, \quad (4)$$

with $b = (-10, 0)$, $f = 0$ and $\epsilon = 0.1$ around a warm dolphin.

Grid generation with **MATLAB** and visualisation using *contour lines* in **GiD**.

Input / output: examples



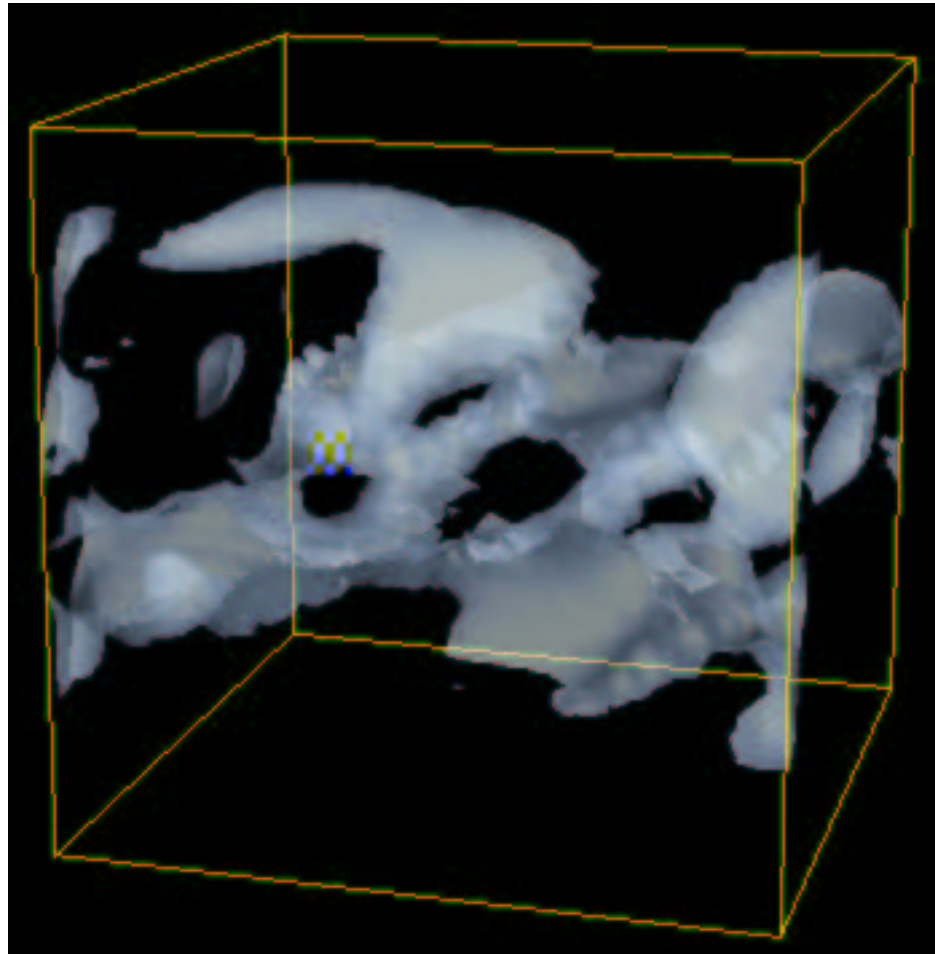
Input / output: examples

Incompressible Navier–Stokes:

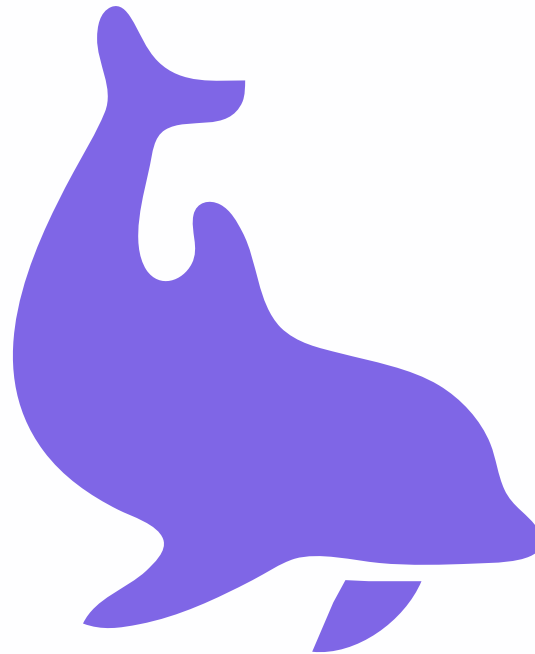
$$\begin{aligned} \dot{u} + u \cdot \nabla u - \nu \Delta u + \nabla p &= f, \\ \nabla \cdot u &= 0, \end{aligned} \tag{5}$$

Visualisation in **OpenDX** of the isosurface for the velocity in a computation of transition to turbulence in shear flow on a mesh consisting of 1,600,000 tetrahedral elements.

Input / output: examples



Web page



- `www.phi.chalmers.se/dolphin`
- `www.freshmeat.net/projects/dolphin`

Organisation of the code

```
doc
src/io
src/la
src/fem
src/grid
src/init
src/test
src/utils
src/common
src/config
src/modules/navier-stokes
src/modules/poisson
src/problems/navier-stokes/benchmark
src/problems/navier-stokes/jet
src/problems/navier-stokes/...
src/problems/poisson
data/grids
```

src/io

Display

Terminal

Curses

Value

Input

Output

inp.h

opendx.h

matlab.h

gid.h

«

Vector

DenseMatrix

SparseMatrix

DirectSolver

SISolver

KrylovSolver

«

src/fem

Discretiser

Equation

EquationSystem

FiniteElement

FunctionSpace

GlobalField

LocalField

Problem

ShapeFunction

TetLinFunction

TetLinSpace

TriLinFunction

TriLinSpace

«

src/grid

Grid

Cell

CellType

Node

Point

Tetrahedron

Triangle

«

src/init

dolphin.h
dolphin.C

«

src/common

Parameter

ParameterList

Settings

Globals

«

src/modules/poisson

EquationPoisson

ProblemPoisson

«

src/problems/poisson

main.C

«

data/grids

```
tetgrid_1_1_1.inp  
tetgrid_4_4_4.inp  
tetgrid_8_8_8.inp  
tetgrid_24_8_8.inp
```

...

«

Grid

```
class Grid{
public:
    ...
    void Init();
    void Clear();

    int GetNoNodes();
    int GetNoCells();

    Node* GetNode(int node);
    Cell* GetCell(int cell);

    void Read(const char *file);
    void Write(const char *file);
    ...
}
```



KrylovSolver

```
class KrylovSolver{
public:

    KrylovSolver();
    ~KrylovSolver(){}

    void SetMethod( KrylovMethod km );

    void Solve(Vector* x, Vector* b);
    void SolveCG(Vector* xvec, Vector* b);
    void SolveGMRES(Vector* xvec, Vector* b);

    ...
}
```



SISolver

```
class SISolver{
public:

    SISolver();
    ~SISolver(){}

    void Solve(SparseMatrix *A, Vector *x, Vector *b);

private:

    void IterateRichardson (SparseMatrix *A, Vector *x, Vector *b);
    void IterateJacobi      (SparseMatrix *A, Vector *x, Vector *b);
    void IterateGaussSeidel (SparseMatrix *A, Vector *x, Vector *b);
    void IterateSOR         (SparseMatrix *A, Vector *x, Vector *b);
    ...
}
```

Vector

```
class Vector{
public:

    Vector (int n);
    ~Vector ();

    void Add (real a, Vector *v);
    real Dot (Vector *v);
    real Norm ();
    ...
}
```

«

SparseMatrix

```
class SparseMatrix{
public:

    SparseMatrix (int m, int n, int *ncols);
    ~SparseMatrix ();

    void Mult(Vector* x, Vector* Ax);

    ...

```

«

Equation

```
class Equation{
public:

    Equation(int noeq, int nsd);

    virtual real IntegrateLHS(TrialFunction &u, TestFunction &v) = 0;
    virtual real IntegrateRHS(TestFunction &v) = 0;

    ...

```

«

Discretiser

```
class Discretiser{
public:

    Discretiser(Grid *grid, Equation *equation);
    ~Discretiser();

    void AssembleLHS(SparseMatrix *A);
    void AssembleRHS(Vector *b);

    ...
}
```

«

Problem

```
class Problem{
public:

    Problem(Grid *grid);
    ~Problem();

    virtual const char *Description() = 0;

    virtual void Solve() = 0;

    ...
}
```

«

ParameterList

```
class ParameterList{
public:

    ...

    void Add(const char *identifier, Type type, ...);
    void Set(const char *identifier, ...);
    void Get(const char *identifier, ...);

    void Save(const char *filename);
    void Load(const char *filename);

    ...
}
```

OpenDX

```
class OpenDX{
public:

    OpenDX (const char *filename, int n, ...);
    ~OpenDX ();

    void SetLabel (int component, const char *label);
    void AddFrame (Grid *grid, Vector *u, real t);

    ...
}
```

«

dolfin.h

```
void dolfin_init (int argc, char **argv);
```

```
void dolfin_end ();
```

```
void dolfin_solve ();
```

```
void dolfin_set_problem (const char *problem);
```

```
void dolfin_set_parameter (const char *identifier, ...);
```

```
void dolfin_get_parameter (const char *identifier, ...);
```

```
void dolfin_save_parameters (const char *filename);
```

```
void dolfin_load_parameters (const char *filename);
```

```
void dolfin_set_boundary_conditions
```

```
    (dolfin_bc (*bc)(real x, real y, real z, int node, int component));
```

```
void dolfin_set_function
```

```
    (const char *identifier, real (*f)(real x, real y, real z, real t));
```

«

main.C

```
#include <dolfin.h>

int main(int argc, char **argv)
{
    kw_set_problem("poisson");

    kw_set_parameter("problem description", "Poisson's equation on the
    kw_set_parameter("grid file",          "../.../data/grids/tetgri
    kw_set_parameter("output file prefix",  "poisson");

    kw_init(argc,argv);
    kw_solve();
    kw_end();

    return 0;
}
```