

## New development in freefem++

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**Abstract** — This is a short presentation of the freefem++ software. In Section 1, we recall most of the characteristics of the software, In Section 2, we recall how to build the weak form of a partial differential equation (PDE) from the strong form. In the 3 last sections, we present different examples and tools to illustrate the power of the software. First we deal with mesh adaptation for problems in two and three dimension, second, we solve numerically a problem with phase change and natural convection, and finally to show the possibilities for HPC we solve a Laplace equation by a Schwarz domain decomposition problem on parallel computer.

**Keywords:** finite element, mesh adaptation, Schwarz domain decomposition, parallel computing, freefem++

### 1. Introduction

This paper intends to give a small presentation of the software freefem++.

A partial differential equation is a relation between a function of several variables and its (partial) derivatives. Many problems in physics, engineering, mathematics and even banking are modeled by one or several partial differential equations.

Freefem++ is a software to solve these equations numerically in dimensions two or three. As its name implies, it is a free software based on the Finite Element Method; it is not a package, it is an integrated product with its own high level programming language; it runs on most UNIX, WINDOWS and MacOs computers. Moreover freefem++ is highly adaptive. Many phenomena involve several coupled systems, for example: fluid-structure interactions, Lorentz forces for aluminum casting and ocean-atmosphere coupling. Multiphysics problems require different finite element approximations or different polynomial degrees, possibly on different meshes. Some algorithms like

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Schwarz' domain decomposition method also require data interpolation on multiple meshes within one program. Freefem++ can handle these difficulties, i.e. *arbitrary finite element spaces on arbitrary unstructured and adapted meshes*.

The characteristics of freefem++ are:

- Problem description (real or complex valued) by their variational formulations, with access to the internal vectors and matrices if needed.
- Multi-variables, multi-equations, bi-dimensional and three-dimensional static or time dependent, linear or nonlinear coupled systems; however the user is required to describe the iterative procedures which reduce the problem to a set of linear problems.
- Easy geometric input by analytic description of boundaries by pieces; however this part is not a CAD system; for instance when two boundaries intersect, the user must specify the intersection points.
- Automatic mesh generator, based on the Delaunay–Voronoi algorithm; the inner point density is proportional to the density of points on the boundaries [17].
- Metric-based anisotropic mesh adaptation. The metric can be computed automatically from the Hessian of any freefem++ function [20].
- High level user friendly typed input language with an algebra of analytic or finite element functions.
- Multiple finite element meshes within one application with automatic interpolation of data on different meshes and possible storage of the interpolation matrices.
- A large variety of triangular finite elements: linear, quadratic Lagrangian elements and more, discontinuous P1 and Raviart–Thomas elements, elements of a non-scalar type, the mini-element, etc., but no quadrangles.
- Tools to define discontinuous Galerkin finite element formulations P0, P1dc, P2dc and keywords: `jump`, `mean`, `intalledges`.
- A large variety of linear direct and iterative solvers (LU, Cholesky, Crout, CG, GMRES, UMFPACK [13], MUMPS [2], SuperLU, etc.) and eigenvalue and eigenvector solvers (ARPACK) [24], and optimization tools like Ipopt [32].

- Near optimal execution speed (compared with compiled C++ implementations programmed directly).
- Online graphics, generation of .txt, .eps, .gnu, .mesh files for further manipulations of input and output data.
- Many examples and tutorials: elliptic, parabolic and hyperbolic problems, Navier–Stokes flows, elasticity, fluid structure interactions, as well as Schwarz’s domain decomposition method, eigenvalue problem, residual error indicator, etc.
- A parallel version using MPI.

## 2. Weak formulation

For the first example consider a Poisson problem on a domain  $\Omega$  with a partition of the boundary  $\partial\Omega$  in  $\Gamma_2, \Gamma_e$ : Find  $u$  such that

$$-\Delta u = 1 \quad \text{in } \Omega, \quad u = 2 \quad \text{on } \Gamma_2, \quad \frac{\partial u}{\partial \vec{n}} = 0 \quad \text{on } \Gamma_e. \quad (2.1)$$

Denote  $V_g = \{v \in H^1(\Omega) / v|_{\Gamma_2} = g\}$ , and the weak form of the equation is obtained by multiplying by  $v$  and integrating by parts (2.1). The variational formulation is: Find  $u \in V_2(\Omega)$ , such that

$$\forall v \in V_0(\Omega), \quad \int_{\Omega} \nabla u \cdot \nabla v = \int_{\Omega} 1v + \int_{\Gamma} \frac{\partial u}{\partial n} v. \quad (2.2)$$

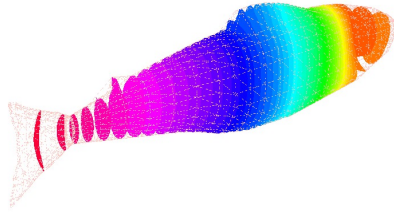
Note that due to  $\partial u / \partial \vec{n} = 0$  on  $\Gamma_2$  and  $v = 0$  the term  $\int_{\Gamma} (\partial u / \partial n) v$  disappears on  $\Gamma_2$ . Finally the finite method simply replaces  $V_g$  with a finite element space, such as

$$V_{gh} = \{v_h \in C^0(\bar{\Omega}) : \forall K \in \mathcal{T}_h, v_h|_K \in P_1 \text{ and } v_h|_{\Gamma_2} = 0\}$$

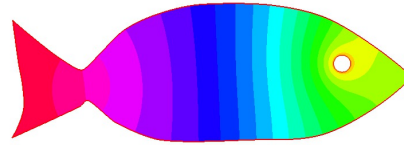
in the case of linear elements; for full detail see [10, 21].

Below, we give an example of an elaborate mesh for a 3D fish and the Poisson problem; the freefem++ code is below and the results are on Figs. 1 and 2.

```
load "msh3" load "medit" load "tetgen"
mesh3 Ths("Y5177_Fish_cut.mesh"); // read skin fish mesh...
real hh = 10; // the final mesh size
real[int] domaine = [0,0,0,1,hh^3/6.];
```



**Figure 1.** 3D Poisson solution iso-surface.



**Figure 2.** 2D Poisson solution iso-value.

```

mesh3 Th=tetg(Ths,switch="pqaAYY",regionlist=domaine);
fespace Xh(Th,P2);
real x0=-200,y0=0,z0=0;
func ball = sqrt((x-x0)^2+(y-y0)^2+(z-z0)^2) < 30;
macro Grad(u) [dx(u),dy(u),dz(u)] // EOM
Xh p,q;
solve laplace(p,q,solver=CG) =
  int3d(Th) ( ball*p*q+Grad(p)'*Grad(q) ) -
  int3d(Th) ( 1*q );
plot(p,fill=1,wait=1,value=0,nbiso=50); // see Fig.1

```

The skin mesh of the 3D fish come from the URL [http://www-roc.inria.fr/gamma/download/counter.php?dir=FISH&get\\_obj=Y5177\\_Fish\\_cut.mesh&aces=Y5177\\_Fish](http://www-roc.inria.fr/gamma/download/counter.php?dir=FISH&get_obj=Y5177_Fish_cut.mesh&aces=Y5177_Fish). The graphics solution of the problem in 2D and 3D are displayed in Figs. 2 and 1, respectively.

### 3. Mesh adaptation

In `freem++` a lot of adaptation tools are implemented; this corresponding to the work of many people [6, 9, 17, 19, 20, 25, 26, 29, 31].

All these tools are based a Delaunay–Voronoi algorithm with a distance (or Metric) between two points  $q^1, q^2$  given by  $\sqrt{{}^t(q^1 - q^2) \mathcal{M} (q^1 - q^2)}$  where the matrix  $\mathcal{M}$  is symmetric positive definite matrix field defined on the mesh  $\mathcal{T}_h$ . Consequently the length  $\ell_{\mathcal{M}}$  of a curve  $\gamma$  for  $]0, 1[ \in \mathbb{R}^d$  with respect to  $\mathcal{M}$  is

$$\ell_{\mathcal{M}} = \int_0^1 \sqrt{{}^t \gamma'(t) \mathcal{M}(\gamma(t)) \gamma'(t)} dt. \quad (3.1)$$

The computation of  $\mathcal{M}$  can be implicit of explicit; the tools to compute  $\mathcal{M}$ , in isotropic cases are scalar fields  $h$  which represent the local mesh size such that  $\mathcal{M} = I_d/h^2$ , where  $I_d$  is the identity matrix.

The idea is to build a meshes with edges of equal length with respect of  $\mathcal{M}$ , i.e. the length with respect to  $\mathcal{M}$  of all the mesh edges should be closed to one so as to get an equi-repartition of the error close to  $\text{err}$  in norm  $L^\infty$ . So for  $P_1$  continuous finites elements. The metric can be defined by

$$\ell_{\mathcal{M}} = \frac{1}{\text{err}} |\partial_h^2 u_h| \quad (3.2)$$

where  $|\partial_h^2 u_h| = \sqrt{(\partial_h^2 u_h)^2}$  and  $\partial_h^2 u$  is an approximation of the Hessian matrix of  $u_h$ . Freefem++ does it automatically by default with `adaptmesh` function (2D), `MetricPk` (2D), or with `mshmet` (2D).

An example of  $L^\infty$  error mesh adaptation with metric for the Poisson problem in an L-shape domain  $\Omega = ]0, 1[^2 \setminus [1/2, 1[^2$ . Find  $u \in H^1(\Omega)$  such that

$$-\Delta u = (x-y) \quad \text{in } \Omega, \quad \frac{\partial u}{\partial \mathbf{n}} = 0 \quad \text{on } \partial\Omega, \quad \int_{\Omega} u = 0 \quad (3.3)$$

we add do a small stabilization term  $\varepsilon u$  to remove the constraint  $\int_{\Omega} u$ , so the weak form is: Find  $u_\varepsilon \in H^1(\Omega)$  such that

$$\forall v \in H^1(\Omega), \quad \int_{\Omega} u_\varepsilon v + \varepsilon u_\varepsilon v - \int_{\Omega} (x-y)v = 0. \quad (3.4)$$

You can easily prove that  $\|u_\varepsilon - u\|_{H^1} < \varepsilon C$ , the discretization with  $P_1$  Lagrange finite element and  $\varepsilon = 10^{-10}$ , is:

```
int[int] lab=[1,1,1,1];
mesh Th = square(6,6,label=lab);
Th=trunc(Th,x<0.5 | y<0.5, label=1);
fespace Vh(Th,P1);
Vh u,v;
real error=0.01;
problem Poisson(u,v,solver=CG,eps=1.0e-6) =
  int2d(Th,qforder=2)(u*v*1.0e-10+ dx(u)*dx(v) + dy(u)*dy(v)) +
  int2d(Th,qforder=2)((x-y)*v);
for (int i=0; i<4; i++)
{
  Poisson;
  plot(u,wait=1);
  Th=adaptmesh(Th,u,err=error);
  plot(Th,wait=1);
  u=u; // reinterpolation of u on new mesh Th.
  error = error/2;
};
```

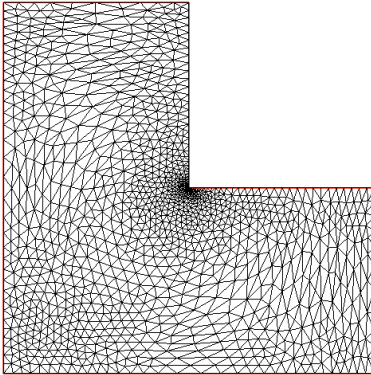


Figure 3. Mesh.

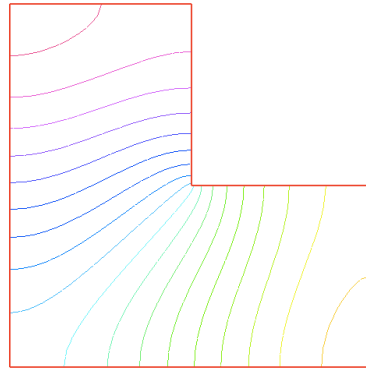


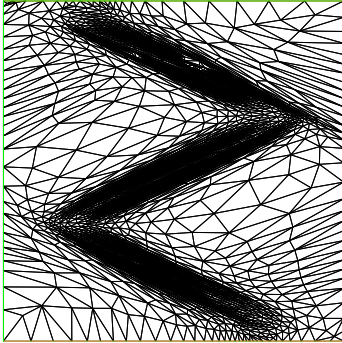
Figure 4. Solution.

The final mesh adaption and associated solution will be show in Figs. 3 and 4

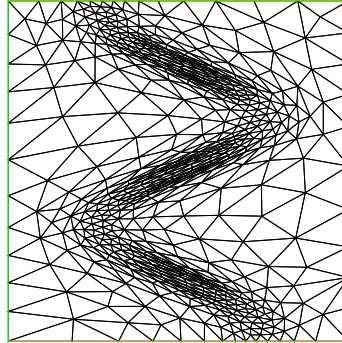
With the `MetricPk` plugin of J.-M. Mirebeau we can build a metric in  $\mathbb{R}^2$  for the Lagrangian triangular continuous approximation of degree  $k$ ,  $P_k$ , with respect to the norm  $W^{r,p}$  with  $r = 0$  or  $1$ ; the output is an adapted mesh with  $Nt$  element with a metric which can be optimal or sub optimal; the lost of optimality is due to constraint on acute angle in triangulation in case of  $r = 1$  (see [26, 27] or J.-M. Mirebeau these details):

```
load "MetricPk"
real Nt = 1000, r=1, k=3, p=1; // def of the parameter...
mesh Th=square(20,20,[2*x-1,2*y-1]);
func f = x^2*y + y^3 + tanh(5*(-2*x + sin(5*y)));
fespace Metric(Th,[P1,P1,P1]);
Metric [m11,m12,m22];
for(int i=0; i<5; i++) {
  plot(Th,wait=true); // see Figs. 6 and 5
  [m11,m12,m22]=[0,0,0]; // resize metric array
  m11 []=MetricPk(Th,f, kDeg=k,rDeg=r,pExp=p, mass=Nt/Z,
    TriangulationType=0); // 0 sub optimal
    // 1 Optimal (=>no acute angle)
  Th = adaptmesh(Th,m11,m12,m22,IsMetric=true); }
```

In 3D we can use `meshmet` a plugin by J. Morice for the `mshmet` library of P. Frey [16]. To compute an isotropic adapted metric on can use `tetgen`, another plugin documented in [30]. The following listing is a full test of this idea.



**Figure 5.** Optimal metric in norm  $W^{1,1}$  for  $10^3$  triangle but it is wrong mesh because the acute constraint is miss.



**Figure 6.** Suboptimal but not acute constraint for  $10^3$  triangle.

```

load "msh3" load "tetgen" load "mshmet" load "medit"
int n = 6;
int[int] l1=[1,1,1,1],l01=[0,1],l11=[1,1];
// label numbering for boundary condition
mesh3 Th3=buildlayers(square(n,n,region=0,label=l1),
    n, zbound=[0,1], labelmid=l11, labelup = l01,
    labeldown = l01);
Th3 = trunc(Th3,(x<0.5) | (y < 0.5) | (z < 0.5),
    label=1); // remove the ]0.5,1[^3 cube
// end of build initial mesh
fespace Vh(Th3,P1);
Vh u,v,usol;
macro Grad(u) [dx(u),dy(u),dz(u)] // EOM
solve Poisson(u,v,solver=CG) = int3d(Th3)( 1e-6*u*v +
    Grad(u)'*Grad(v) ) - int3d(Th3)( (z-y/2-x/2)*v );
real errm=1e-2; // level of error
for(int ii=0; ii<5; ii++) {
    Vh h; h[]=mshmet(Th3,u,normalization=1,aniso=0,nbregul=1,
        hmin=1e-3,hmax=0.3,err=errm);
    errm *= 0.6; // change the level of error
    Th3=tetgreconstruction(Th3,switch="raAQ",
        sizeofvolume=h*h*h/6.);
    Poisson;
    plot(u,wait=1,nbiso=15); } // see Fig.8
medit("U3",Th3,u); // see Fig.7

```

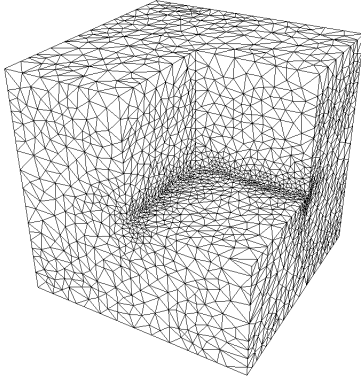


Figure 7. 3D adapted mesh.

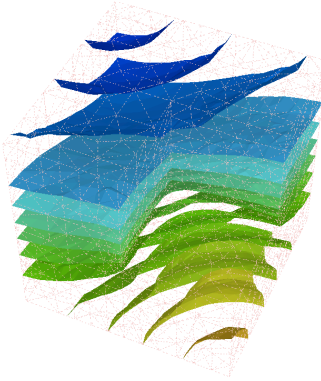


Figure 8. Iso surface.

Finally, tools like `splitmesh` (only in 2D) can also be used; it splits triangle uniformly into sub triangles. In 3D the plugin `mng3d-v4` moves 3D meshes or build 3D anisotropic meshes inside each element (see [15]).

#### 4. Phase change with natural convection

This example illustrates the coupling of natural convection modeled by the Boussinesq approximation and liquid to solid phase change in  $\Omega = ]0, 1[^2$ . No slip condition for the fluid are applied at the boundary and adiabatic condition on upper and lower boundary and given temperature  $\vartheta_r$  (resp  $\vartheta_l$ ) at the right and left boundaries. The starting point of the problem is Brainstorming session (part I) of the third `freefem++` days in December, 2012, this is almost the Orange Problem is described in URL <http://www.ljll.math.upmc.fr/hecht/ftp/ff++days/2011/Orange-problem.pdf>, we present the validation part with paper [34].

The main interest of this example is to show the capability of `freefem++` to solve complex problem with different kind of non-linearity. You can find other complex example in [1, 5, 7, 8, 11, 12, 23] for example, (thanks to I. Danalia to the help in the modelization).

So the full model is: Find  $\mathbf{u} = (u_1, u_2)$ , the velocity field,  $p$  the pressure field and  $\vartheta$  the temperature field in domain  $\Omega$  such that

$$\left\{ \begin{array}{ll} \mathbf{u} \text{ given} & \text{in } \Omega_s \\ \partial_t \mathbf{u} + (\mathbf{u} \nabla) \mathbf{u} + \nabla \cdot \mu \nabla \mathbf{u} + \nabla p = -c_T \mathbf{e}_2 & \text{in } \Omega_f \\ \nabla \cdot \mathbf{u} = 0 & \text{in } \Omega_f \\ \partial_t \vartheta + (\mathbf{u} \nabla) \vartheta + \nabla \cdot k_T \nabla \vartheta = \partial_t S(T) & \text{in } \Omega. \end{array} \right. \quad (4.1)$$

Where  $\Omega_f$  is the fluid domain and the solid domain is  $\Omega_s = \Omega \setminus \Omega_f$ . The enthalpy of the change of phase is given by the function  $S$ ,  $\mu$  is the relative vis-



cosity,  $k_T$  the thermal diffusivity.

In  $\Omega_f = \{x \in \Omega; \vartheta > \vartheta_f\}$ , with  $\vartheta_m$  the melting temperature the solid has melt.

We modeled, the solid phase as a fluid with huge viscosity, so:

$$\mu = \begin{cases} \vartheta < \vartheta_f \sim 10^6 \\ \vartheta \geq \vartheta_m \sim \frac{1}{\text{Re}}. \end{cases}$$

This removes movement in the solid phase, and here  $\text{Re}$  is the Reynolds number of the liquid phase.

The Stefan enthalpy  $S_c$  with defined by  $S_c(\vartheta) = H(\vartheta)/S_{th}$  where  $S_{th}$  is the Stefan number, and  $H$  is the Heaviside function (0, if  $(\vartheta < 0)$ , 1 else). with use the following smooth the enthalpy:

$$S(\vartheta) = \frac{\tanh(50(\vartheta - \vartheta_m))}{2S_{te}}.$$

We apply a fixed point algorithm for the phase change part (the domain  $\Omega_f$  is fixed at each iteration) and a full no-linear Euler implicit scheme with a fixed domain for the rest. We use a Newton method to solve the non-linearity.

Remark, if we don't make mesh adaptation the Newton method do not converge, and if we use explicit method this diverge too, and finally if we implicit the dependence in  $\Omega_s$  the method also diverge. This is a really difficult problem.

The finite element space to approximate  $u_1, u_2, p, \vartheta$  is defined by:

```
fespace Wh(Th, [P2,P2,P1,P1]);
```

We do mesh adaptation a each time step, with the following code:

```
Ph ph = S(T), pph=S(Tp);
Th = adaptmesh(Th,T,Tp,ph,pph, [u1,u2], err=errh,
               hmax=hmax,hmin=hmax/100,ratio = 1.2);
```

This mean, we adapt with all variable plus the 2 melting phase a time  $n + 1$  and  $n$  and we smooth the metric with a ratio of 1.2 to account for the movement of the melting front.

The Newton loop and the fixed point are implemented as follows:

```
real err=1e100, errp;
for (int kk=0; kk<2; ++kk) // 2 step of fixed point on  $\Omega_s$ 
```

```

{ nu = nuT; // recompute the viscosity in  $\Omega_s, \Omega_f$ 
  for (int niter=0; niter<20; ++niter) // newton loop
  { BoussinesqNL;
    err = u1w[].linfty;
    cout << niter << " err NL " << err << endl;
    u1[] -= u1w[];
    if(err < tolNewton) break; } // convergence...
}

```

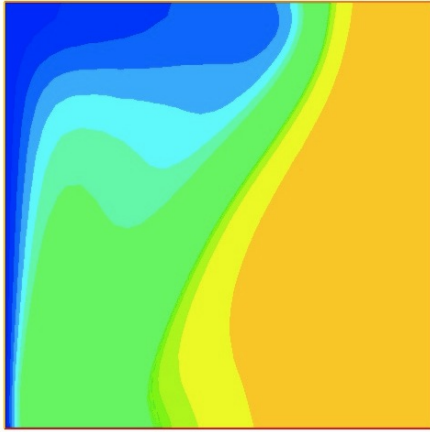
The linearized problem is:

```

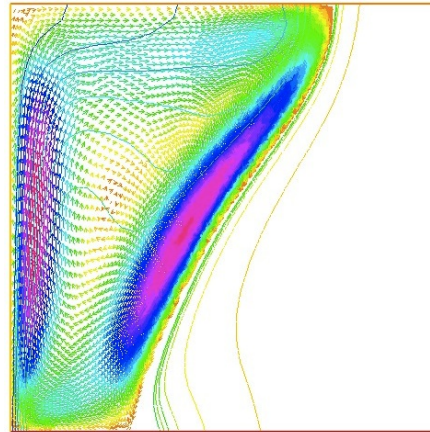
problem BoussinesqNL([u1w,u2w,pw,Tw],[v1,v2,q,TT])
= int2d(Th) (
  [u1w,u2w,Tw]'*[v1,v2,TT]*cdt
  + UgradV(u1,u2,u1w,u2w,Tw)' * [v1,v2,TT]
  + UgradV(u1w,u2w,u1,u2,T)' * [v1,v2,TT]
  + ( Grad(u1w,u2w)'*Grad(v1,v2)) * nu
  + ( Grad(u1,u2)'*Grad(v1,v2)) * dnu* Tw
  + cmT*Tw*v2
  + grad(Tw)'*grad(TT)*kT
  - div(u1w,u2w)*q -div(v1,v2)*pw - eps*pw*q
  + dS(T)*Tw*TT*cdt
)
- int2d(Th) (
  [u1,u2,T]'*[v1,v2,TT]*cdt
  + UgradV(u1,u2,u1,u2,T)' * [v1,v2,TT]
  + ( Grad(u1,u2)'*Grad(v1,v2)) * nu
  + cmT*T*v2
  + grad(T)'*grad(TT)*kT
  - div(u1,u2)*q -div(v1,v2)*p
  - eps*p*q // stabilization term
  + S(T)*TT*cdt
  - [u1p,u2p,Ip]' * [v1,v2,TT]*cdt
  - S(Tp)*cdt*TT
)
+ on(1,2,3,4, u1w=0,u2w=0) + on(2,Tw=0) + on(4,Tw=0);

```

The parameters of the computation are taken from [34] case 2,  $\vartheta_m = 0$ ,  $Re = 1$ ,  $S_{le} = 0.045$ ,  $P_r = 56.2$ ,  $R_a = 3.27 \cdot 10^5$ ,  $\vartheta_l = 1$ ,  $\vartheta_r = -0.1$  so in this case  $cmT = c_T = -R_a/P_r$ ,  $kT = k_T = 1/P_r$ ,  $eps = 10^{-6}$ , time step  $\delta t = 10^{-1}$ ,  $cdt = 1/\delta t$ , and with a good agreement this is Fig. 6 of [34] at time  $t = 80$  (see Figs. 9 and 10).



**Figure 9.** Iso-value of the temperature at time  $t = 80$  of problem (4.1).



**Figure 10.** Velocity at time 80 of problem (4.1).

## 5. A Schwarz domain decomposition example in parallel

The following is an explanation of the scripts `DDM-Schwarz-*.edp` of the distribution. This is Schwarz domain decomposition in parallel with a complexity almost independent of the number of process (with a coarse grid preconditioner), thanks to F. Nataf for the all the discussion to implementation of this algorithm (see [28] for the theory).

To solve the following Poisson problem on domain  $\Omega$  with boundary  $\Gamma$  in  $L^2(\Omega)$ :

$$-\Delta u = f \quad \text{in } \Omega, \quad u = g \quad \text{on } \Gamma$$

where  $f$  and  $g$  are two given functions of  $L^2(\Omega)$  and of  $H^{1/2}(\Gamma)$ .

Let  $(\pi_i)_{i=1, \dots, N_p}$  be a regular partition of unity of  $\Omega$ , i.e.:

$$\pi_i \in \mathcal{C}^0(\Omega) : \quad \pi_i \geq 0, \quad \sum_{i=1}^{N_p} \pi_i = 1.$$

Denote  $\Omega_i$  the sub domain which is the support of  $\pi_i$  function and also denote  $\Gamma_i$  the boundary of  $\Omega_i$ . The parallel Schwarz method is as follows.

Let  $\ell = 0$  be the iterator and  $u^0$  an initial guess respecting the boundary condition (i.e.  $u^0|_{\Gamma} = g$ ):

$$-\Delta u_i^\ell = f \quad \text{in } \Omega_i, \quad u_i^\ell = u^\ell \quad \text{on } \Gamma_i \setminus \Gamma, \quad u_i^\ell = g \quad \text{on } \Gamma_i \cap \Gamma \quad (5.1)$$

$$u^{\ell+1} = \sum_{i=1}^{N_p} \pi_i u_i^\ell \quad \forall i = 1, \dots, N_p. \quad (5.2)$$

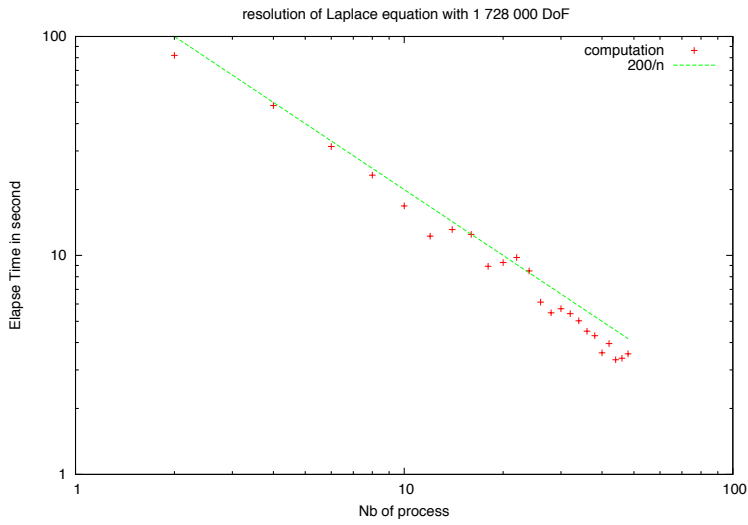


Figure 11. Resolution of Poisson equation with 1 728 000 DoF, scalability curve.

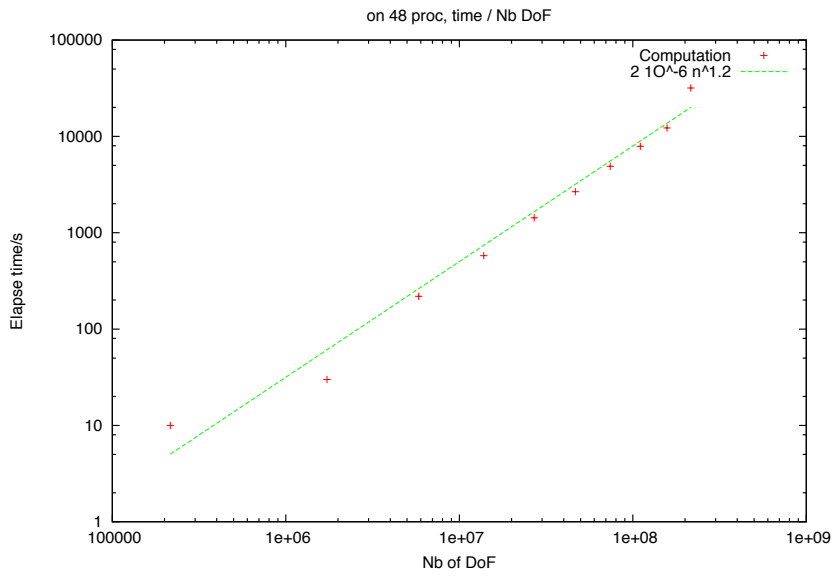


Figure 12. On 48 processors, test on number of DoF from 216 10<sup>3</sup> to 216 10<sup>6</sup>, complexity curve.

After discretization with the Lagrange finite element method, with a compatible mesh  $\mathcal{T}_{h_i}$  for  $\Omega_i$ , there exists a global mesh  $\mathcal{T}_h$  such that  $\mathcal{T}_{h_i}$  is included in  $\mathcal{T}_h$ .

**Remark 5.1.** We avoid using finite element spaces associated to the full domain  $\Omega$  because it is too expensive.

Now let us look at this academic example:

$$-\Delta u = 1 \quad \text{in } \Omega = ]0, 1[^3, \quad u = 0 \quad \text{on } \partial\Omega. \quad (5.3)$$

In the following test we use a GMRES version preconditioned by the Schwarz algorithm ( $P_s$ ) and with a coarse grid solver ( $P_c$ ) on a coarse mesh. To build a new preconditioner  $IP$  from two preconditioner  $P_c$  and  $P_s$  we use the following approximation  $0 \sim I - AP^{-1} = (I - AP_c^{-1})(I - AP_s^{-1})$  and after a simple calculus we have  $P^{-1} = P_c^{-1} + P_s^{-1} - P_c^{-1}AP_s^{-1}$ .

## 6. Conclusion

Freefem++ is a continuously evolving software because it is easy to add new tools, finite elements. We have tried to illustrate this with three examples. In the future we expect to include more tools for three dimensional anisotropic meshes adaptive, automatic differentiation by operator overloading, seamless integration of parallel tools to free the user off low level programming.

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