

## Abstract

Refinement technique is used to adapt meshes to the singularities of the numerical solution of any problem. Local refinement is applied in mesh areas where the solution must be accurate and it is necessary to dispose of reliable error indicators or estimators which define the elements that must be refined. Usually, these indicators or estimators are difficult to obtain. On the other hand, derefinement is used to remove mesh elements on which the numerical solution can be easily interpolated. This process can be guided by a parameter that indicates the required precision for the numerical solution. This paper is about using refinement/derefinement techniques to adapt meshes to the numerical solution but using global refinement instead of local ones. No error estimators or indicators are needed, and two numerical parameters are used to automatize the process.

**Keywords:** 3-D triangulations, finite element, adaptive meshes, object oriented method, adaptive refinement/derefinement, data structures, wind field models.

## 1 Introduction

Refinement/derefinement techniques are used when it is necessary to adapt meshes to the singularities of the numerical solution. Usually, local refinement is applied to elements where the solution must be adjusted. Previously, these elements must be determined using any kind of error indicator or estimator. For each mesh element, an error indicator must be computed, and thus it is possible to establish mesh zones where the refinement must be carried out. Depending on the problem, the error indicator or estimator can be more or less difficult to obtain. Moreover, it could be impossible to get a reliable one.

On the other hand, derefinement algorithms remove mesh elements where the nu-

merical solution can be obtained with the desired precision from neighbour elements. Comparing the numerical solution of an element with the interpolated solution, and if it is enough accurate, the element could be removed. Note that in this case we are dealing with numerical parameters that gives us the desired precision, and they are always possible to establish.

In this paper we propose a simple method to carry out meshes adaptation without error indicators or estimators using global refinement. With global refinement all mesh elements divided, so indicators are not needed. After that, a derefinement process will remove elements which are not necessary. Each iteration of this method implies higher computational cost than local refinement, but the total number of iterations of the proposed method could be much less if we not chose a optimal refinement strategy.

In section 2 the algorithm for refinement/derefinement will be briefly presented. Details can be found in [2, 3]. In section 3, a wind field model is commented. It has been used for implementing the algorithm and compare the proposed method with the traditional refinement. This model uses the gradient of the solution as error indicator, and it is detailed in [1, 6]. In section 4 the implementation algorithm is explained, and we present different situations that has been useful to achieve it. In section 5 can be found a problem on which the algorithm has been tested. Besides, a real problem has been implemented with a mesh of the south part of La Palma island. This mesh has been generated with [4]. Finally, in the section 6 a brief comparison between methods is commented.

## 2 Refinement/Derefinement

The refinement algorithm is based on 8-subtetrahedral subdivision, and it has been presented in [2]. From an initial triangulation  $\tau_0$ , the goal is to build a sequence of nested meshes  $T = \{\tau_0 < \tau_1 < \tau_2 < \dots < \tau_m\}$ , where  $\tau_{j+1}$  is obtained from refinement of  $\tau_j$ . Each element  $t_i \in \tau_j$  will have associated an error indicator  $\eta_i^j$ , and it will be refined if:

$$\eta_i^j \geq \gamma \eta_{max}^j \quad (1)$$

$\eta_{max}^j$  is the maximal value of the error indicator of elements in  $\tau_j$ .  $\gamma$  is the refinement parameter and  $\gamma \in [0, 1]$ .

The derefinement algorithm is the inverse of the refinement. It has been presented in [3], and it takes into account the numerical solution in the mesh nodes. Any node  $n_i \in \tau_j$  will have computed a numerical solution  $v_i^j$ . If we consider  $n_p$  and  $n_q$  nodes of the surrounding edge of  $n_i$ , then  $n_i$  can be removed if:

$$\left| v_i^j - \frac{v_p^j + v_q^j}{2} \right| < \epsilon \quad (2)$$

The derefinement parameter  $\epsilon$  is used to establish the desired precision of the solu-

tion in nodes.

### 3 Wind Field Model: Mass Consistent Model in 3-D

This model [6] is based on the continuity equation for an incompressible flow where the air density is constant in the domain  $\Omega$  and *no-flow-through* conditions on  $\Gamma_b$  (terrain and top) are considered

$$\vec{\nabla} \cdot \vec{u} = 0 \quad \text{in } \Omega \quad (3)$$

$$\vec{n} \cdot \vec{u} = 0 \quad \text{on } \Gamma_b \quad (4)$$

We formulate a least-square problem in  $\Omega$  with  $\vec{u}(\tilde{u}, \tilde{v}, \tilde{w})$  to be adjusted

$$E(\vec{u}) = \int_{\Omega} [\alpha_1^2 ((\tilde{u} - u_0)^2 + (\tilde{v} - v_0)^2) + \alpha_2^2 (\tilde{w} - w_0)^2] d\Omega \quad (5)$$

where the interpolated wind  $\vec{v}_0 = (u_0, v_0, w_0)$  is obtained from experimental measurements, and  $\alpha_1, \alpha_2$  are the Gauss precision moduli.

We consider Dirichlet condition for open or *flow-through* boundaries and Neumann condition for terrain and top

$$\phi = 0 \quad \text{on } \Gamma_a \quad (6)$$

$$\vec{n} \cdot T \vec{\nabla} \mu = -\vec{n} \cdot \vec{v}_0 \quad \text{on } \Gamma_b \quad (7)$$

This problem can be solved using tetrahedral finite elements (see [5]) which leads to a set of  $4 \times 4$  elemental matrices and  $4 \times 1$  elemental vectors. These are assembled to form a symmetric linear system of equations which is solved by a preconditioned conjugate gradient method.

The construction of the interpolated wind is a two step process: horizontal and vertical interpolation. Both processes are detailed in [1, 6].

In the generation of adaptive meshes, the local refinement of the domain is necessary due, on one hand, to the geometry and, on the other hand, to the numerical solution. The computation of error estimators or at least suitable error indicators of the numerical solution is carried out to determine the elements to be refined or derefined in a mesh. In this wind model we use one error indicator, proposed in [1], which takes into account the gradient of the solution in each element.

### 4 Implementation

To obtain the desired mesh, it will be necessary to compare any mesh parameter. We have chosen the number of nodes, and we define  $w_n$  as the number of nodes or  $\tau_n$ . This parameter is used because it indicates the number of points that will be used for computing the numerical solution.

The first implementation can be seen in the Algorithm 1. It was carried out attending to the  $\epsilon$  parameter. The stop criteria consist on obtaining a mesh  $\tau_n$  such  $w_n = w_{n-1}$ . This implies that all new elements added by refinement of  $\tau_{n-1}$  have been removed by derefinement, so the numerical solution is adjusted in  $\tau_n$  according to the  $\epsilon$  parameter. The Figure 1 represents the expected evolution of  $w_i$  in the resulting meshes. This number would increase slower in next steps due to it could have mesh zones with the desired precision in their numerical solution.

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**Algorithm 1** Initial approach

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Be  $\tau_0$  the initial mesh  
 $n=0$   
**repeat**  
     $n=n+1$   
     $\tau'_n = \tau_{n-1}$  globally refined  
    Compute numerical solution of  $\tau'_n$   
     $\tau_n = \tau'_n$  derefined according to  $\epsilon$  parameter  
**until**  $w_n = w_{n-1}$

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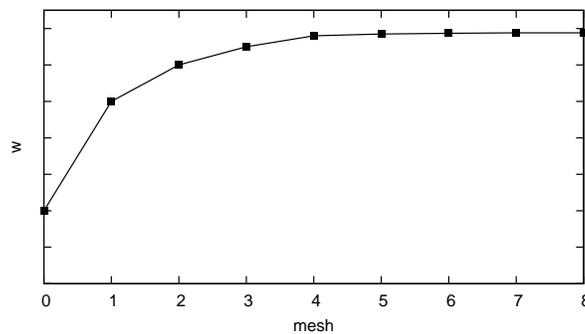


Figure 1: Expected evolution or  $w$

In the first tests the obtained results were good with values of  $\epsilon$  relatively high (6 m/s). But with lower values of  $\epsilon$  (2 m/s) the algorithm does not work properly. In Figure 2 we can see its behavior.

Note that:

1. There are meshes with  $w_k < w_{k-1}$ . Refinements improve the numerical solution in certain zones and derefinement removes elements introduced many steps before.
2. There are two groups of meshes:
  - $w_{11} = w_{13} = w_{15} = \dots$
  - $w_{12} = w_{14} = w_{16} = \dots$

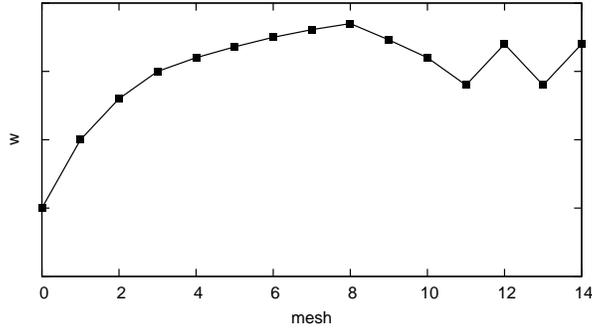


Figure 2:  $w$  for test with  $\epsilon = 2$  m/s

The second aspect implies that the condition of the Algorithm 1,  $w_n = w_{n-1}$ , will be never reached. To solve this problem, we have redefined the comparison between meshes. The stop criteria can not be the mentioned before. Instead of that, two meshes will be considered similar if the difference between  $w_n$  and  $w_k$  ( $k < n$ ) is lesser than a user defined percentage, that is, it will be compared  $w_n$  with all the previous  $w$  to find two similar meshes.

In final test, with  $\epsilon = 1.5$  m/s, we have obtained graphics like Figure 3. This means that the numerical solution will be never adjusted. This problem is a particular case of the wind field model, in elements close the terrain. The difference between the numerical solution of the terrain elements and their adjacent presents low variation with refinements. If new elements are introduced, the difference with this elements is similar than previous, so the numerical solution is not being improved. The derefinement process would not remove that elements, and they will be refined time and again.

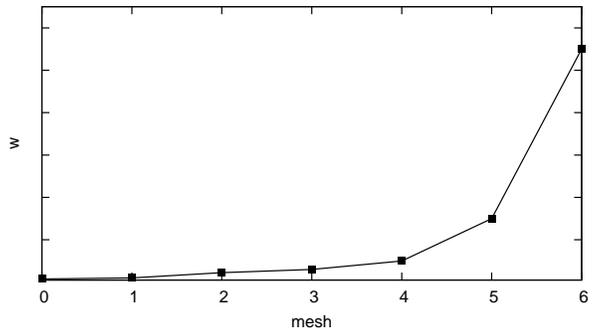


Figure 3:  $w$  for test with  $\epsilon = 1.5$  m/s

To prevent this, we introduce a new parameter  $\delta$ . It indicates the minimal size for the edges of any mesh. The derefinement process will be carried out attending to both,  $\epsilon$  and  $\delta$  parameters. If an element has any edge lesser than  $\delta$ , it will be removed. The final implementation can be seen in Algorithm 2.

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**Algorithm 2** Implemented

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Be  $\tau_0$  the initial mesh
n=0
loop
  n=n+1
   $\tau'_n = \tau_{n-1}$  globally refined
  Compute numerical solution of  $\tau'_n$ 
   $\tau_n = \tau'_n$  derefined according to  $\delta$  and  $\epsilon$  parameters
  for i=0 to n-1 do
    Exit when  $w_n \approx w_i$ 
  end for
end loop
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## 5 Applications

All the executions were run in a XEON dual processor, with 2 Gb of RAM, under linux and programs compiled with GNU C++. To stop the process we have defined that two meshes are similar if their  $w_i$  are different in less than 1%:

$$\frac{w_k}{w_n} \in [0.99, 1.01], k < n \quad (8)$$

The first mesh used is a 3D gauss curve as shown in Figure 4. It consists on 1680 nodes and 7645 elements for a simulated domain of  $10000 \times 10000 \times 10000 m^3$ .

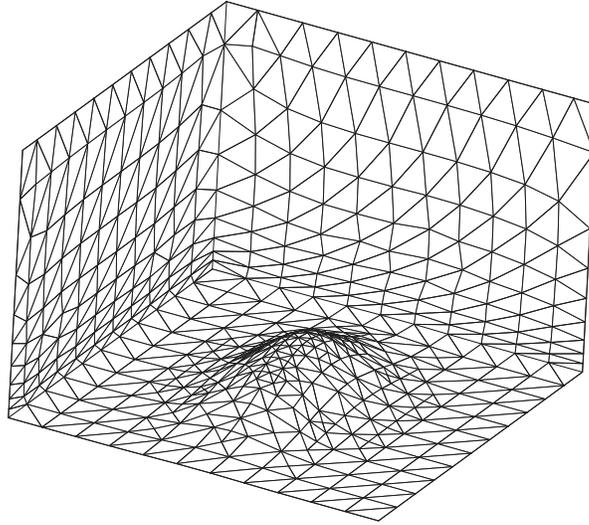


Figure 4:  $\tau_0$

The first test was made with parameters  $\epsilon = 2$  m/s and  $\delta = 40$  m. Five steps were necessary to reach the adjusted mesh (only few meshes are shown in Figure 6(a) and

6(b)). In Figure 1 it can be seen the graphic for  $w$ , and in Table 1(b) are printed CPU time for each process.

Figure 5: Meshes obtained from  $\tau_0$  with  $\epsilon = 2$  m/s and  $\delta = 40$  m

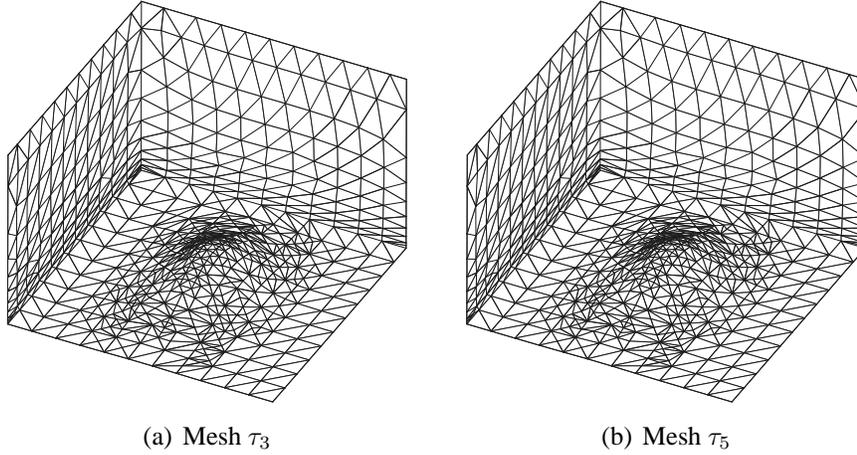
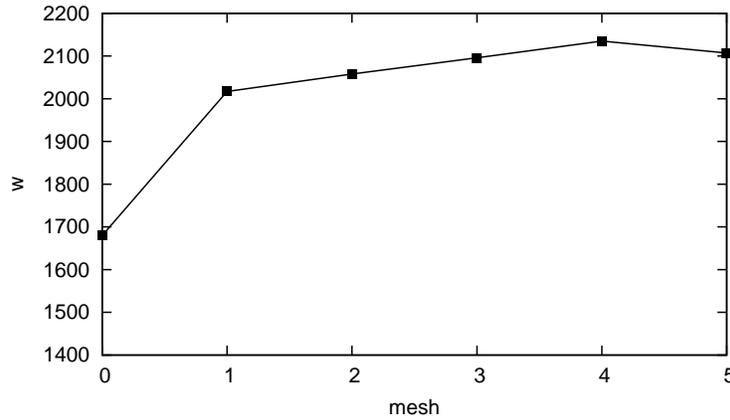


Table 1: Data for  $\tau_0$  with  $\epsilon = 2$  m/s and  $\delta = 40$  m (Figure 5)

(a) Evolution of  $w$



(b) CPU Time in seconds

Initial $\tau_{n-1}$	$w_{n-1}$	Refine	$w'_n$	Compute	Derefine	Final $\tau_n$	$w_n$
$\tau_0$	1680	3.72	11787	5.93	2.68	$\tau_1$	2017
$\tau_1$	2017	5.18	12591	5.31	3.16	$\tau_2$	2058
$\tau_2$	2058	5.39	12865	5.52	3.45	$\tau_3$	2096
$\tau_3$	2096	5.76	13084	5.51	3.83	$\tau_4$	2135
$\tau_4$	2135	6.09	13326	5.98	4.00	$\tau_5$	2107

Another test was made with parameters  $\epsilon = 1.5$  m/s and  $\delta = 80$  m. For this problem, twelve steps were necessary to adjust the mesh (Figure 6 and Table 2).

In both test the algorithm worked properly, and they were useful to validate and adjust the method.

Figure 6: Meshes obtained from  $\tau_0$  with  $\epsilon = 1.5$  m/s and  $\delta = 80$  m

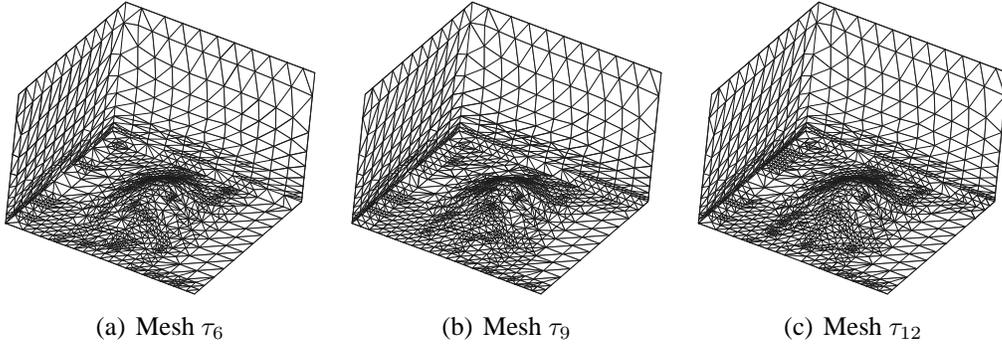
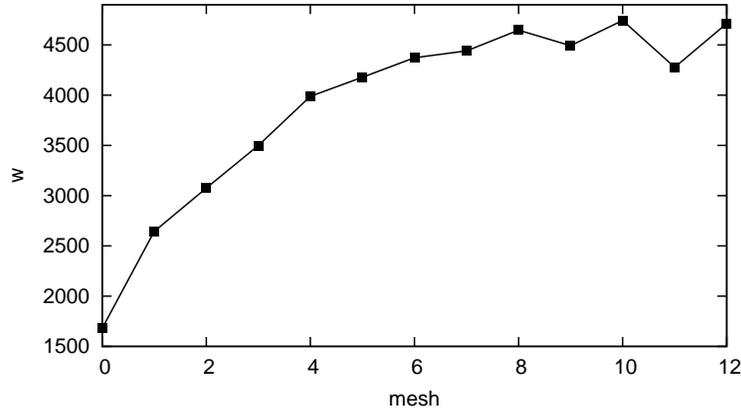


Table 2: Data for  $\tau_0$  with  $\epsilon = 1.5$  m/s and  $\delta = 80$  m (Figure 6)

(a) Evolution of  $w$



(b) CPU Time in seconds

Initial $\tau_{n-1}$	$w_{n-1}$	Refine	$w'_n$	Compute	Derefine	Final $\tau_n$	$w_n$
$\tau_0$	1680	3.78	11787	6.01	2.94	$\tau_1$	2644
$\tau_1$	2644	6.84	15857	7.43	4.57	$\tau_2$	3075
$\tau_2$	3075	8.32	18775	9.69	7.05	$\tau_3$	3498
$\tau_3$	3498	9.82	21564	11.67	8.02	$\tau_4$	3987
$\tau_4$	3987	11.73	24532	14.40	8.15	$\tau_5$	4178
$\tau_5$	4178	12.26	25982	16.70	10.46	$\tau_6$	4372
$\tau_6$	4372	13.36	27250	16.85	10.77	$\tau_7$	4440
$\tau_7$	4440	13.96	27597	18.23	11.43	$\tau_8$	4647
$\tau_8$	4647	14.64	29158	20.26	12.35	$\tau_9$	4492
$\tau_9$	4492	14.76	28123	18.52	11.49	$\tau_{10}$	4743
$\tau_{10}$	4743	16.12	29617	18.55	12.43	$\tau_{11}$	4277
$\tau_{11}$	4277	14.60	26783	16.73	11.61	$\tau_{12}$	4713

We have also used a real geometry generated with [4, 5]. It represents part of the south of La Palma island, which can be seen in Figure 7. It consists on 4535 nodes and 21137 elements for a real domain of  $45600 \times 31200 \times 6000$  m<sup>3</sup>.

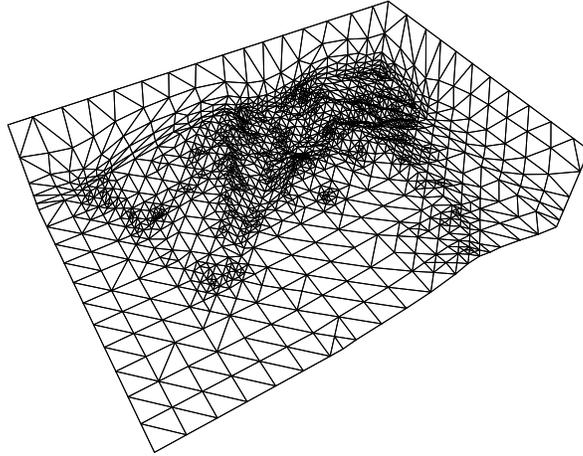


Figure 7:  $lp_0$

The goal was to obtain an adjusted mesh according to  $\epsilon = 4$  m/s and  $\delta = 40$  m. Only five steps were necessary. In Figure 8 we can see meshes generated and in Table 3 information about the process.

On the other hand, we have used the same mesh  $lp_0$  with the traditional method, that is, local refinement - error estimation - derefinement. The  $\gamma$  parameter for refinement was adjusted to 0.6 and  $\epsilon$  parameter was equal than the above run (4 m/s).  $\delta$  parameter was not necessary. Meshes obtained are similar that shown in Figure 8, but in Tables 4 and 5 we can see that number of steps is higher.

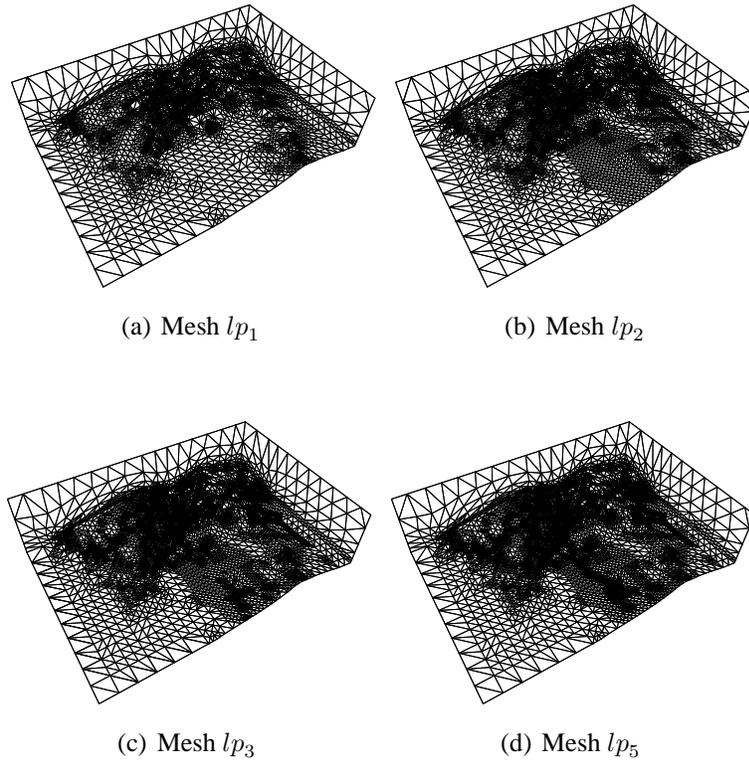
## 6 Conclusion

To adjust meshes to a numerical solution with local refinement is necessary to establish any error indicator or estimator, and not always is easy. The number of steps depends on both, position and number of singularities of numerical solution. The computational cost in each step is lower than the proposed method due to the number of elements treated.

With global refinement all singularities are treated simultaneously. We avoid the error estimation and the method only depends on numerical parameters:  $\epsilon$  and, in our case,  $\delta$ . Of course, due to all mesh elements are involved in each step the computational cost is higher, but the number of steps is lower.

If the number of step is very high with local refinement, this method could be advantageous not only for using numeric parameters but CPU time.

Figure 8: Meshes obtained from  $lp_0$  with  $\epsilon = 4$  m/s and  $\delta = 40$  m



## Acknowledgements

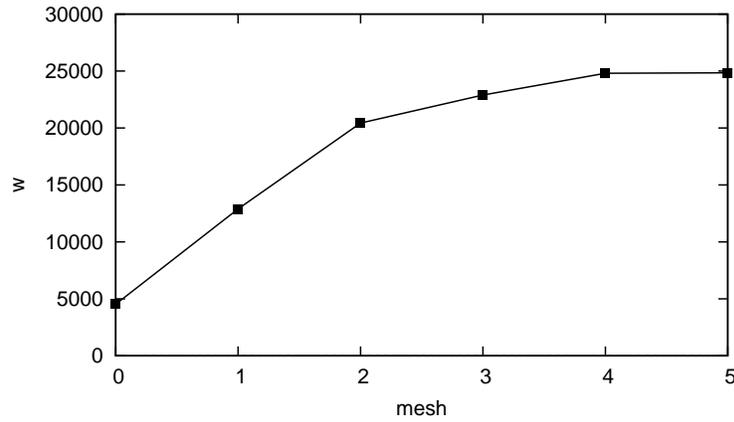
This work has been supported by the Spanish Government, “Ministerio de Educacin y Ciencia” and FEDER, grant contacts: CGL2004-06171-C03-02/CLI.

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Table 3: Data for  $lp_0$  with  $\epsilon = 4$  m/s and  $\delta = 40$  m (Figure 8)

(a) Evolution of  $w$



(b) CPU Time in seconds

Initial $lp_{n-1}$	$w_{n-1}$	Refine	$w_n$	Compute	Derefine	Final $lp_n$	$w_n$
$lp_0$	4535	9.04	32139	23.16	8.37	$lp_1$	12898
$lp_1$	12898	29.83	83648	99.84	27.71	$lp_2$	20426
$lp_2$	20426	46.58	125989	177.27	42.19	$lp_3$	22887
$lp_3$	22887	52.47	139159	211.89	53.05	$lp_4$	24806
$lp_4$	24806	57.92	150243	203.95	59.67	$lp_5$	24845

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Table 4: Data for  $lp_0$  with  $\gamma = 0.6$  and  $\epsilon = 4$  m/s

Initial $lp_{n-1}$	$w_{n-1}$	Refine	$w'_n$	Compute	Derefine	Final $lp_n$	$w_n$
$lp_0$	4535	5.90	18200	10.59	5.03	$lp_1$	12768
$lp_1$	12768	19.75	44340	35.01	19.00	$lp_2$	17593
$lp_2$	17593	28.38	45780	39.78	19.62	$lp_3$	18773
$lp_3$	18773	37.23	42330	33.26	22.35	$lp_4$	19108
$lp_4$	19108	38.88	41944	31.12	23.72	$lp_5$	19304
$lp_5$	19304	45.69	49265	44.49	27.29	$lp_6$	19519
$lp_6$	19519	48.27	52984	50.30	29.66	$lp_7$	20058
$lp_7$	20058	56.93	55388	57.96	31.93	$lp_8$	20294
$lp_8$	20294	59.61	54401	45.93	32.72	$lp_9$	20536
$lp_9$	20536	52.58	51416	49.49	32.04	$lp_{10}$	20505

Table 5: Evolution of  $w$  in adjusting of  $lp_0$  using global (Table 3) and local (Table 4) refinement

