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Numerical Analysis

# Accelerating the method of finite element patches using approximately harmonic functions

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### Abstract

We present a new variant of a domain decomposition method with complete overlap and not necessarily nested grids to solve numerically elliptic problems with multi-scale data. The novelty of the method consists in restricting finite element functions on the coarse grid to be approximately harmonic inside the subdomain where a finer triangulation is applied. Numerical experiments confirm an increase in the convergence rate over a previously proposed method. *To cite this article: J. He et al., C. R. Acad. Sci. Paris, Ser. I 345 (2007).* 

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#### Résumé

Acceleration de la méthode des patches d'éléments finis en utilisant des fonctions approximativement harmoniques. Nous présentons une nouvelle variante de la méthode de décomposition de domaines avec un recouvrement complet et des maillages non nécessairement emboîtés pour la résolution numérique des problèmes elliptiques avec des données multi-échelle. La nouveauté de la méthode consiste dans la restriction de l'espace des fonctions éléments finis grossières à être des approximations des fonctions harmoniques dans le sous-domaine sur lequel une triangulation fine est appliquée. Des expériences numériques confirment une augmentation du taux de convergence par rapport à une méthode proposée précédemment. *Pour citer cet article : J. He et al., C. R. Acad. Sci. Paris, Ser. I 345 (2007).* 

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## Version française abrégée

Soit  $\Omega \subset \mathbb{R}^2$  un domaine polygonal borné et  $\Lambda \subseteq \Omega$  un autre domaine polygonal « beacoup plus petit » que  $\Omega$ . On considère le problème de trouver  $u \in H_0^1(\Omega)$  tel que

$$a(u, v) = \langle f | v \rangle, \quad \forall v \in H_0^1(\Omega),$$

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où  $a(u, v) = \int_{\Omega} \nabla u \cdot \nabla v \, d\mathbf{x}$  et  $f \in H^{-1}(\Omega)$  est une fonction donnée. On suppose que la solution u varie brutalement dans  $\Lambda$  de sorte que l'on ne puisse pas discrétiser le problème par une méthode d'éléments finis sur une grille qui est de même résolution sur  $\overline{\Omega} \setminus \Lambda$  que sur  $\overline{\Lambda}$ . On introduit donc deux espaces d'éléments finis :  $V_H$  – l'espace des fonctions continues, s'annulant sur  $\partial \Omega$  et polynômiales par morceaux sur une triangulation régulière  $\mathcal{T}_H$  de  $\overline{\Omega}$  et  $V_h$ – l'espace des fonctions continues, s'annulant sur  $\partial A$ , polynômiales par morceaux sur une triangulation régulière  $T_h$ de  $\overline{\Lambda}$  ( $h \ll H$ ) et qui sont prolongées par zéro sur  $\overline{\Omega} \setminus \Lambda$ . En posant  $V_{Hh} = V_H + V_h$  on cherche une approximation  $u_{Hh} \in V_{Hh}$  de u qui satisfait (1) pour toutes  $v \in V_{Hh}$ . En pratique, il est impossible de construire une base standard d'éléments finis pour l'espace  $V_{Hh}$  et de résoudre le problème discrétisé directement. C'est pourquoi un algorithme itératif pour calculer  $u_{Hh}$  a été proposé dans [3] (voire Algorithm 1 ci-dessous) dans lequel on ne résout que des problèmes standards sur  $V_H$  et  $V_h$ . Cet algorithme s'avère être très efficace dans les situations où la grille  $\mathcal{T}_h$  est emboîtée dans  $T_H$  (c'est-à-dire chaque triangle de  $T_h$  est contenu dans un seul triangle de  $T_H$ ), mais son taux de convergence peut être extrêmement faible pour certains placements de  $\mathcal{T}_h$  par rapport à  $\mathcal{T}_H$ . Cette observation peut être expliquée par le fait que le taux de convergence est géré par l'angle abstrait (par rapport au produit scalaire  $a(\cdot, \cdot)$ ) entre  $V_h$  et  $\tilde{V}_H$  où  $\tilde{V}_H$  est le complément *a*-orthogonal de  $V_H \cap V_h$  dans  $V_H$ . Dans le cas emboîté,  $\tilde{V}_H$  ne contient que les fonctions qui approchent des fonctions harmoniques dans  $\Lambda$ . L'espace  $\tilde{V}_H$  est donc approximativement aorthogonal à  $V_h$ , ce qui explique la rapide convergence de la méthode dans ce cas. Si, au contraire,  $T_h$  est placé sur  $\mathcal{T}_H$  de sorte à ce que  $V_H \cap V_h = 0$ , les espaces  $V_h$  et  $\tilde{V}_H = V_H$  peuvent être presque « parallèles », ce qui implique une mauvaise convergence. Le but de cette Note est de transférer partiellement les bonnes propriétés de convergence du cas emboîté au cas général. L'idée est de remplacer l'espace  $V_H$  dans l'algorithme précédent par son sous-espace  $\bar{V}_H$ qui est le complément a-orthogonal du  $V_H^0$  où  $V_H^0$  est le sous-espace de  $V_H$  composé des fonctions qui s'annulent dans  $\bar{\Omega} \setminus \Lambda$ . On obtient alors des problèmes non standards posés dans  $\bar{V}_H$  qui peuvent être résolus efficacement moyennant l'introduction de problèmes auxiliaires et peu coûteux posés dans  $V_H^0$ . Cette nouvelle variante de l'algorithme (voir Algorithm 2 ci-dessous) converge vers une approximation  $\bar{u}_{Hh} \in \bar{V}_{Hh}$  de *u* (avec  $\bar{V}_{Hh} = \bar{V}_H + V_h$ ) qui satisfait (1) pour tout  $v \in \overline{V}_{Hh}$ . On établit l'estimation a priori pour  $\overline{u}_{Hh}$  du type

$$|u - \bar{u}_{Hh}|_1 \leq C(H^r ||u||_{H^q(\Omega \setminus \bar{A})} + h^s ||u||_{H^q(A)})$$

où r et s sont les degrés des polynômes utilisés dans la construction de  $V_H$  et  $V_h$  respectivement et  $q = \max(r, s) + 1$ . On conclut la Note par les expériences numériques qui montrent une convergence très rapide de notre nouvelle méthode.

Let  $\Omega \subset \mathbb{R}^2$  be a bounded polygonal domain and  $\Lambda \Subset \Omega$  be another polygonal domain which is 'much smaller' than  $\Omega$ . Given an  $f \in H^{-1}(\Omega)$ , we consider the Poisson–Dirichlet problem of finding u such that  $-\Delta u = f$  in  $\Omega$  and u = 0 on  $\partial \Omega$ , whose weak formulation is: find  $u \in H_0^1(\Omega)$  such that

$$a(u, v) = \langle f | v \rangle, \quad \forall v \in H_0^1(\Omega), \tag{2}$$

where  $a: H_0^1(\Omega) \times H_0^1(\Omega) \to \mathbb{R}$  is symmetric, continuous and coercive bilinear form

$$a(u, v) = \int_{\Omega} \nabla u \cdot \nabla v \, \mathrm{d}\mathbf{x}.$$
(3)

We are interested in the 'multi-scale' situation where the solution u of (2) has a very strong variation inside  $\Lambda$  so that the discretization of (2) by a finite element method cannot be done using grids with the same resolution on  $\overline{\Omega} \setminus \Lambda$  and  $\overline{\Lambda}$ . Therefore we introduce two finite element spaces:  $V_H = \{g \in C_0(\overline{\Omega}) \text{ such that } g|_K \in \mathbb{P}_r(K), \forall K \in \mathcal{T}_H \text{ and } g =$ 0 on  $\partial \Omega\}$  where  $\mathcal{T}_H$  is a regular triangulation of  $\overline{\Omega}$  and  $\mathbb{P}_r(K)$  is the space of polynomials of degree  $\leq r$  on a triangle  $K \in \mathcal{T}_H$ , and  $V_h = \{g \in C_0(\overline{\Omega}) \text{ such that } g|_K \in \mathbb{P}_s(K), \forall K \in \mathcal{T}_h \text{ and } g = 0 \text{ on } \overline{\Omega} \setminus \Lambda\}$  where  $\mathcal{T}_h$  is a regular triangulation of  $\overline{\Lambda}$ . Setting  $V_{Hh} = V_H + V_h$  we search as approximation for u the function  $u_{Hh} \in V_{Hh}$  satisfying

$$a(u_{Hh}, v) = \langle f | v \rangle, \quad \forall v \in V_{Hh}.$$
<sup>(4)</sup>

A priori  $V_H \cap V_h$  does not necessarily reduce to the element zero and it is impossible, practically speaking, to exhibit a 'finite element'-type basis of the space  $V_{Hh}$ . That is why the following algorithm to compute  $u_{Hh}$  was suggested in [3]:

### Algorithm 1.

- (1) Set  $u_H^0 = 0$ ,  $u_h^0 = 0$ .
- (2) For n = 1, 2, 3, ... find

  - (i)  $u_H^n \in V_H$  such that  $a(u_H^n, v) = \langle f | v \rangle a(u_h^{n-1}, v), \forall v \in V_H$ , (ii)  $u_h^n \in V_h$  such that  $a(u_h^n, v) = \langle f | v \rangle a(u_H^n, v), \forall v \in V_h$ , set  $u_{Hh}^n = u_H^n + u_h^n$ .

**Remark 1.** Algorithm 1 can be viewed as a block Gauss–Seidel method to solve (4). It was presented in [3] in a different and more general form with possible use of over-relaxation making it similar to the SOR method. The present form makes evident the connection of this algorithm to the Chimera method [1]. Note also that the coarse update (i) and fine update (ii) are swapped here in comparison with the original version of [3]. This is motivated by the numerical results presented below.

Letting  $P_h: V_{Hh} \to V_h$  and  $P_H: V_{Hh} \to V_H$  be the orthogonal (with respect to  $a(\cdot, \cdot)$ ) projectors from  $V_{Hh}$  to  $V_h$ and  $V_H$  respectively, it is easy to see that

$$u_{Hh} - u_{Hh}^n = (I - P_h)(I - P_H)(u_{Hh} - u_{Hh}^{n-1}).$$

We recall the following result from [2]:

**Lemma 1.** Let V be an inner product space of finite dimension,  $V_1$  and  $V_2$  be two subspaces of V such that V = $V_1 + V_2$  and  $P_j: V \to V_j$ , j = 1, 2, be orthogonal projectors from V to  $V_j$ . Let  $V_0 = V_1 \cap V_2$  and suppose that  $V_1 \neq V_0, V_2 \neq V_0$ . Define  $\tilde{\gamma} = \tilde{\gamma}(V_1, V_2), 0 \leq \tilde{\gamma} < 1$  to be the cosine of the abstract angle between  $V_1$  and  $V_2$  so that

$$\tilde{\gamma}(V_1, V_2) = \sup_{\substack{v_1 \in V_1 \cap V_0^{\perp}, \ v_1 \neq 0 \\ v_2 \in V_2 \cap V_0^{\perp}, \ v_2 \neq 0}} \frac{(v_1, v_2)}{\|v_1\| \|v_2\|}$$
(5)

where  $(\cdot, \cdot)$  denotes the inner product in V and  $\|\cdot\|$  is the norm induced by it. Then the spectral radius of operator  $B = (I - P_2)(I - P_1)$  is  $\rho(B) = \tilde{\gamma}^2$ .

Applying this result in the case  $V = V_{Hh}$  equipped with the inner product  $a(\cdot, \cdot)$ ,  $V_1 = V_H$ ,  $V_2 = V_h$  we readily see that Algorithm 1 converges and its asymptotic convergence rate is given by  $\tilde{\gamma}(V_H, V_h)^2$ , hence the convergence is slow if the parameter  $\tilde{\gamma}(V_H, V_h)$  is close to 1. Note, that even a more general version of Algorithm 1 using an over-relaxation can allow one to improve the convergence rate to only  $1 - \sqrt{8(1 - \tilde{\gamma})} + O(1 - \tilde{\gamma})$  as  $\tilde{\gamma} \to 1$ , see [2]. Very poor convergence was indeed observed in practice in some cases [2,6,8].

There is, however, at least one situation in which Algorithm 1 converges reasonably fast, namely when the triangulations  $\mathcal{T}_H$  and  $\mathcal{T}_h$  are nested, i.e. each triangle of  $\mathcal{T}_h$  is contained in only one triangle of  $\mathcal{T}_H$ . This experimental observation is easy to interpret in view of Lemma 1. The intersection space  $V_0 = V_H \cap V_h$  can be described in this case as  $V_0 = V_H^0$  where  $V_H^0$  is the subspace of  $V_H$  defined by

$$V_H^0 = \left\{ v_H \in V_H : v_H(x) = 0 \; \forall x \in \Omega \setminus \bar{\Lambda} \right\},\tag{6}$$

so the *a*-orthogonal compliment to  $V_0$  in  $V_H$  is

$$\bar{V}_{H} = \left\{ v_{H} \in V_{H} : a(v_{H}, \phi_{H}) = 0 \; \forall \phi_{H} \in V_{H}^{0} \right\}.$$
(7)

It consists of finite element functions of  $V_H$  that are approximately harmonic inside  $\Lambda$  on triangulation  $\mathcal{T}_H$ . The space  $V_H$  is therefore approximately *a*-orthogonal to the space  $V_h$ , since  $a(\phi_H, \phi_h)$  should be small for any  $\phi_H \in \overline{V}_H$  and any  $\phi_h \in V_h$  vanishing on the boundary  $\partial \Lambda$  by definition of  $V_h$ . Hence we can indeed expect that the parameter  $\tilde{\gamma}(V_H, V_h) = \tilde{\gamma}(V_H, V_h)$  is small in the case of nested triangulations. Note that  $V_h$  is a subspace of  $V_H^0$  so that the approximation  $u_{Hh}$  given by (4) would not change if we replace  $V_{Hh} = V_H + V_h$  by  $\bar{V}_{Hh} = \bar{V}_H + V_h$  there. Moreover, hypothetically we could replace the space  $V_H$  by  $\bar{V}_H$  on all the iterations of Algorithm 1 and this would produce the same iterates  $u_{Hh}^n$ .

We inspire ourselves from the last remark to modify Algorithm 1 by replacing the space  $V_H$  by its subspace consisting of functions which are approximately harmonic inside  $\Lambda$  in the hope to transfer partially the good convergence

properties of the nested case to the general setting of arbitrary triangulations. More specifically, we keep the notations (6) and (7) to define  $\bar{V}_{Hh} = \bar{V}_H + V_h$  also in a nonnested case. Note that  $V_H^0$  consists now of all the functions from  $V_H$  with supports inside  $\bar{\Lambda}_H$  where  $\bar{\Lambda}_H$  is the union of (closed) triangles from  $\mathcal{T}_H$  that lie completely inside  $\bar{\Lambda}$ . We are now looking for the approximation  $\bar{u}_{Hh} \in \bar{V}_{Hh}$  to u in (2) such that

$$a(\bar{u}_{Hh}, v) = \langle f | v \rangle, \quad \forall v \in V_{Hh}.$$
(8)

In order to run Algorithm 1 with  $V_H$  replaced by  $\bar{V}_H$ , we should be able to solve the problems like the following one: given  $g \in V'_H$  find  $u_H \in \overline{V}_H$  such that

$$a(u_H, v) = \langle g | v \rangle, \quad \forall v \in \bar{V}_H.$$
(9)

In practice, we do not have an explicit finite element basis for the space  $\bar{V}_H$  so that a direct implementation of (9) is impossible. To circumvent this difficulty we consider two auxiliary finite element problems:

(a) Find λ<sub>H</sub> ∈ V<sup>0</sup><sub>H</sub> such that a(λ<sub>H</sub>, μ) = ⟨g|μ⟩, ∀μ ∈ V<sup>0</sup><sub>H</sub>.
(b) Find u<sub>H</sub> ∈ V<sub>H</sub> such that a(u<sub>H</sub>, v) = ⟨g|v⟩ − a(λ<sub>H</sub>, v), ∀v ∈ V<sub>H</sub>.

One can easily see that  $u_H$  in (b) gives the solution of (9) and both problems (a) and (b) are straightforward to implement. We can now announce a modified version of Algorithm 1 in which  $V_H$  is implicitly replaced by  $\bar{V}_H$  as:

## Algorithm 2.

(1) Set  $u_H^0 = 0$ ,  $u_h^0 = 0$ .

(2) For n = 1, 2, 3, ... find

- (i)  $\lambda_H^n \in V_H^0$  such that  $a(\lambda_H^n, \mu) = (f|\mu) a(u_h^{n-1}, \mu), \ \forall \mu \in V_H^0$ , (ii)  $u_H^n \in V_H$  such that  $a(u_H^n, v) = \langle f|v \rangle a(u_h^{n-1}, v) a(\lambda_H^n, v), \ \forall v \in V_H$ , (iii)  $u_h^n \in V_h$  such that  $a(u_h^n, v) = \langle f|v \rangle a(u_H^n, v), \ \forall v \in V_h$ , set  $u_{Hh}^n = u_H^n + u_h^n$ .

Note that auxiliary problem (i) in the algorithm above is normally very cheap and its matrix is just a submatrix of that in (ii) corresponding to the basis functions of  $V_H$  whose supports are inside  $\bar{A}_H$ . Applying Lemma 1 with  $V_1 = V_H$ ,  $V_2 = V_h$  we see immediately that Algorithm 2 converges to the solution  $\bar{u}_{Hh}$  of (8). Although the space  $\bar{V}_{Hh}$  in (8) is poorer than  $V_{Hh}$  in (4), its solution  $\bar{u}_{Hh}$  approximates u with the optimal order as shows the following proposition the proof of which will be found in [4]:

**Proposition 2.** Let  $q = \max(r, s) + 1$  and suppose that the solution u of (2) is in  $H^q(\Omega)$ . If the patch  $\Lambda$  is convex then the solution  $\bar{u}_{Hh}$  of (8) satisfies

$$|u - \bar{u}_{Hh}|_{1} \leqslant C \left( H^{r} \|u\|_{H^{q}(\Omega \setminus \bar{\Lambda})} + h^{s} \|u\|_{H^{q}(\Lambda)} \right), \tag{10}$$

where H and h are the maximum diameters of the triangles in  $T_H$  and  $T_h$  respectively and C is a constant independent of H and h.

Algorithm 2 is particularly efficient in the case  $\bar{\Lambda} = \bar{\Lambda}_H$ . This situation will be referred to as 'boundary conforming' triangulations. We prove in [4] that the convergence rate of Algorithm 2 can be estimated independently of H and hin this case.

Numerical example. We illustrate the two algorithms presented above on the example of the Poisson–Dirichlet problem (2) in  $\Omega = (-1, 1)^2$  with the exact solution

$$u = \cos(k\pi x_1)\cos(k\pi x_2) + \eta\chi(r)\exp\left(\frac{1}{\epsilon^2} - \frac{1}{|\epsilon^2 - r^2|}\right)$$

where  $r = \sqrt{x_1^2 + x_2^2}$  and  $\chi(r) = 1$  if  $r \leq \epsilon$ ,  $\chi(r) = 0$  if  $r > \epsilon$ ;  $k = \frac{1}{2}$ ,  $\eta = 20$  and  $\epsilon = 0.3$ . In a region close to (0, 0)where the solution is peaking, we apply a patch with a finer grid. We choose  $P_1$  finite elements for both  $V_H$  and  $V_h$ . The software Freefem++ [5] is used to generate the grids and to implement the algorithm. The 'mixed terms' of the



Fig. 1. (a) Grids  $\mathcal{T}_H$  and  $\mathcal{T}_h$  in the boundary conforming situation; (b) zoom of (a) to the rectangle  $(-0.4, 0.4)^2$ ; (c) Grids  $\mathcal{T}_H$  and  $\mathcal{T}_h$  in the boundary nonconforming situation; (d) zoom of (c) to the rectangle  $(-0.4, 0.4)^2$ .

type  $a(u_H, v_h)$  with  $u_H \in V_H$  and  $u_h \in V_h$  are calculated exactly using the numerical integration on the auxiliary grid representing the intersections of  $\mathcal{T}_H$  and  $\mathcal{T}_h$  and constructed with the aid of software Triangle [7].

We consider two situations:

(i) Boundary conforming case with an unstructured grid  $\mathcal{T}_H$  having the grid size  $H_b$  on the boundary and structured grid  $\mathcal{T}_h$  that covers the patch  $\Lambda = (-0.2, 0.2)^2$ . The grid  $\mathcal{T}_H$  is constructed so that  $\partial \Lambda$  is a union of edges of some triangles of  $\mathcal{T}_H$ , see Fig. 1(a), (b) with  $H_b = 1/10$  and  $\mathcal{T}_h$  consisting of  $24 \times 24$  nodes. The integrals on the right-hand side  $\langle f | v_h \rangle$  for  $v_h \in V_h$  are approximated by 7-point quadrature formulas on  $\mathcal{T}_h$ . Their coarse triangulation

Table 1

Relative  $H^1$ - and  $L^2$ -error on iterations and for the converged solution, number of iterations until convergence and numerically computed asymptotic convergence rate in the case of boundary conforming grids

H <sub>b</sub> iter.	Algorithm 1						Algorithm 2						
	1/10		1/20		1/40		1/10		1/20		1/40		
	$H^1$	$L^2$	$H^1$	$L^2$									
0.5	3.82E-1	1.44E-1	2.77E-1	6.42E-2	1.41E-1	1.78E-2	9.98E-1	8.46E-1	9.98E-1	8.48E-1	9.98E-1	8.49E-1	
1.5	4.31E-2 7.86E-2	9.58E-3 4.33E-2	4.58E-2 5.14E-2	2.60E-3 1.49E-2	2.43E-2 2.61E-2	7.03E-4 5.01E-3	1.76E-2 9.59E-3	9.50E-3 4.39E-3	4.18E-2 2.28E-3	1.94E - 3 1.02E - 3	9.12E-4 5.33E-4	5.30E-4 2.50E-4	
2	3.83E-2	7.37E-3	4.30E-2	2.30E-3	2.38E-2	5.45E-4	8.18E-3	4.21E-3	1.99E-3	1.01E-3	5.16E-4	2.49E-4	
2.5	7.06E-2	3.92E-3	4.81E-2	1.40E-3	2.54E-2	4.88E-3	7.94E-3	4.23E-3	1.95E-3	1.00E-3	5.14E-4	2.50E-4	
conv. iter.	1.12E-2 54	3.69E-3	1.25E-2 68	1.97E-3	1.22E-2 65	5.03E-4	7.87E-3 5	4.20E-3	1.94E-3 4	1.00E-3	5.13E-4 3	2.49E-4	
$\tilde{\gamma}^2$	0.9565		0.9927		0.9967		0.2006		0.2046		0.2046		

Table 2 Same results as in Table 1 for the case of boundary nonconforming grides

H <sub>b</sub> iter.	Algorithm 1						Algorithm 2						
	1/10		1/20		1/40		1/10		1/20		1/40		
	$H^1$	$L^2$	$H^1$	$L^2$	$H^1$	$L^2$	$H^1$	$L^2$	$H^1$	$L^2$	$H^1$	$L^2$	
0.5	4.14E-1	1.44E-1	2.72E-1	6.41E-2	1.39E-1	1.78E-2	1.00	8.50E-1	9.98E-1	8.48E-1	9.98E-1	8.49E-1	
1	7.55E-2	7.81E-3	3.94E-2	1.97E-3	2.28E-2	5.77E-4	1.03E-2	6.18E-3	2.18E-3	1.25E-3	6.08E-4	4.05E-4	
1.5	8.87E-2	3.55E-2	4.71E-2	1.54E - 2	2.49E-2	5.17E-3	1.01E - 2	5.22E-3	2.36E-3	1.27E-3	6.42E-4	4.36E-4	
2	6.85E-2	7.94E-3	3.71E-2	1.96E-3	2.22E - 2	5.09E-4	8.93E-3	4.79E-3	2.10E-3	1.11E-3	5.56E-4	3.01E-4	
2.5	7.91E-2	3.18E-2	4.41E-2	1.45E-2	2.42E-2	5.04E-3	9.40E-3	5.09E-3	2.16E-3	1.17E-3	5.91E-4	3.72E-4	
conv.	1.53E-2	4.11E-3	1.17E-2	2.32E-3	1.16E-2	4.31E-4	8.72E-3	4.89E-3	2.09E-3	1.09E-3	5.51E-4	2.87E-4	
iter.	76		61		61		11		4		3		
$\tilde{\gamma}^2$	0.9687		0.9907		0.9969		0.8236		0.9339		0.9698		

counterparts  $\langle f | v_H \rangle$  for  $v_H \in V_H$  are approximated using both coarse triangulation  $\mathcal{T}_H$  in  $\overline{\Omega} \setminus \Lambda$  and fine triangulation  $\mathcal{T}_h$  in  $\Lambda$ .

(ii) Boundary nonconforming case with the same unstructured grid  $\mathcal{T}_H$  as above and structured grid  $\mathcal{T}_h$  that covers the patch  $\Lambda = (-0.27, 0.27)^2$ , see Fig. 1(c), (d) with  $H_b = 1/10$  and  $\mathcal{T}_h$  consisting of  $31 \times 31$  nodes so that the grid size *h* is approximately the same as in the case (i). Unlike situation (i), the integrals  $\langle f | v_H \rangle$  for  $v_H \in V_H$  are approximated using only coarse triangulation  $\mathcal{T}_H$  because of algorithmic difficulties with the evaluation of integrals on  $\Lambda \setminus \overline{\Lambda}_H$ .

Table 1 represents the results obtained in situation (i) on the grids as in Fig. 1(a), (b) and two consecutive twofold refinements thereof. We report the evolution of the error  $e^n = u_{Hh}^n - u$  and  $e_{Hh}^{n-1/2} = u_H^n + u_h^{n-1} - u$  in  $H^1$  and  $L^2$  norms on first 2 iterations, the error after convergence, the stopping criterion being  $|u_{Hh}^n - u_{Hh}^{n-1}|_1 < 10^{-4}|u_{Hh}^n|_1$  and the number of iterations needed to reach the convergence. In order to determine numerically the asymptotic convergence rate we have also run our algorithms with f = 0 starting from some non-zero  $u_H^0$ ,  $u_h^0$  until the error reduction per iteration quotient attains a stable value (with variation below  $10^{-6}$ ). This quantity is equal to  $\tilde{\gamma}^2$  according to Lemma 1 and we report it in the last line of Table 1. Numerical results clearly demonstrate the superiority of our new Algorithm 2. Note also that the error in Algorithm 1 tends to be smaller after the fine update than after the coarse update (look at the iterations with numbers n - 1/2).

Analogous numerical experiments in boundary nonconforming situation (ii) are reported in Table 2. The numerically calculated abstract angle  $\tilde{\gamma}$  governing the asymptotic convergence rate of Algorithm 2 is no longer as good as in the boundary conforming case. However, the experiments with 'multi-scale' non-zero right-hand side f show again the superiority of the new algorithm both in terms of the number of iterations needed to reach convergence and the accuracy of the converged solution.

Algorithm 2 can be easily extended to general elliptic problems in  $\mathbb{R}^d$  for any positive integer d and also to the situations with multiple patches. We note finally that another way to increase the convergence rate of Algorithm 1 is to use it as a preconditioner for a Conjugate Gradient (CG) method [6]. One can expect therefore to obtain yet faster convergence by incorporating Algorithm 2 into the CG approach.

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