# **Error Analysis of the Adaptive Variational Multiscale Method**

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#### **The Model Problem**

#### The Dirichlet Problem.

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$$-\nabla \cdot a \nabla u = f$$
 in  $\Omega$ ,  $u = 0$  on  $\partial \Omega$ .

where a > 0 is bounded.

Weak Form. Find  $u \in \mathcal{V} = H_0^1(\Omega)$  such that

$$a(u,v) = (a\nabla u, \nabla v) = (f,v)$$
 for all  $v \in \mathcal{V}$ .

#### **Example of a Basic Error Estimate**

If we for the moment assume a to be periodic  $a = a(x/\epsilon)$  we have (Hou),

$$||e||_{a}^{2} = a(e,e) \le C\left(\frac{h}{\epsilon}\right)^{2} ||f||^{2}$$

- h > \epsilon will give unreliable results even with exact quadrature.
- h < \epsilon will often be to computationally expensive.

#### **Conclusion of the Simple Estimate**

We need to solve PDE:s on a scale that captures the oscillations but we can not afford to do it on the entire domain.



Figure 1: Coarse  $H > \epsilon$  and fine  $h < \epsilon$  mesh. Since  $||e||_a \sim \frac{h}{\epsilon}$ ,  $||e||_a \sim H$  is reasonable.

## **First Step of Three: Spaces**

Leaving the periodic setting. We let  $\mathcal{V} = \mathcal{V}_c \oplus \mathcal{V}_f$ .

- The goal of our calculation affects the choice of  $\mathcal{V}_f$ .
- Hierarchical basis gives an approximation of the nodal interpolant of u onto  $\mathcal{V}_c$ .
- The Scott-Zhang interpolant.
- A modified hierarchical basis that increases  $L^2$  orthogonality between the scales.

Fine scale base functions needs to decay rapidly.

#### **First Step: Spaces**



Figure 2: Different choices of fine scale base functions.

#### Second Step: VMM and Decoupling

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We let  $u = u_c + u_f$  and  $v = v_c + v_f$  in the weak form (Hughes): Find  $u_c \in \mathcal{V}_c$  and  $u_f \in \mathcal{V}_f$  such that

$$\begin{aligned} a(u_c, v_c) + a(u_f, v_c) &= (f, v_c) \quad \text{for all } v_c \in \mathcal{V}_c, \\ a(u_f, v_f) &= (f, v_f) - a(u_c, v_f) \\ &:= (R(u_c), v_f) \quad \text{for all } v_f \in \mathcal{V}_f. \end{aligned}$$

Let  $\{\varphi_i\}_{i\in\mathcal{N}}$  be a partition of unity and define  $u_{f,i}\in\mathcal{V}_f$ ,

$$a(u_{f,i}, v_f) = (R(u_c), \varphi_i v_f)$$
 for all  $v_f \in \mathcal{V}_f$ .

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#### Second Step: VMM and Decoupling



#### Figure 3: The fine scale solution $u_{f,i}$ .

#### **Third Step: Approximation on Patches**

We finally solve the fine scale problems approximately on patches  $\omega_i$ : Find  $U_c \in \mathcal{V}_c$  and  $U_{f,i} \in \mathcal{V}_f^h(\omega_i)$  such that

$$a(U_c, v_c) + a(U_f, v_c) = (f, v_c) \text{ for all } v_c \in V_c,$$
  
$$a(U_{f,i}, v_f) = (R(U_c), \varphi_i v_f) \text{ for all } v_f \in V_f^h(\omega_i).$$

Fine scale problems can be solved for all coarse nodes  $\mathcal{N}$  or some  $\mathcal{F}$  where  $\mathcal{N} = \mathcal{C} \cup \mathcal{F}$ .

#### **Third Step: Approximation on Patches**



Figure 4: The fine scale solution  $U_{f,i}$  for different patches  $\omega_i$ .

One, two, and three layer stars. We get good agreement with the truth mesh solution.

#### **Error Equation**

We introduce the coarse and fine scale error  $e_c = u_c - U_c$ ,  $e_{f,i} = u_{f,i} - U_{f,i}$ , and  $e = e_c + \sum_{i \in \mathcal{N}} e_{f,i}$ . We have the following orthogonality properties:

$$a(e_c, v_c) + a(e_f, v_c) = 0, \text{ for all } v_c \in \mathcal{V}_c$$

#### and

$$a(e_{f,i}, v_f) + a(e_c, \varphi_i v_f) = 0, \quad \text{for all } v_f \in \mathcal{V}_f^h(\omega_i).$$

$$||e||_{a}^{2} = a(e, e) = a(e, e - \pi_{c}e)$$
  
=  $(f, e - \pi_{c}e) - a(U_{c}, e - \pi_{c}e) - a(U_{f}, e - \pi_{c}e)$   
=  $\sum_{i \in \mathcal{N}} (f, \varphi_{i}(e - \pi_{c}e)) - a(U_{c}, \varphi_{i}(e - \pi_{c}e))$   
 $- a(U_{f,i}, e - \pi_{c}e)$ 

#### Remember

$$a(U_{f,i}, v_f) + a(U_c, \varphi_i v_f) - (f, \varphi_i v_f) = 0$$
, for all  $v_f \in \mathcal{V}_f^h(\omega_i)$ 



Figure 5: We study  $z = e - \pi_c e$ . We have  $z - \pi_f^0 z$  to the left,  $z - \pi_f z$  in the middle and  $\pi_f z - \pi_f^0 z$  to the right.

We work with these two parts separately. We start with  $z - \pi_f z$  where  $z = e - \pi_c e$ ,

$$\sum_{i \in \mathcal{N}} (f, \varphi_i(z - \pi_f z)) - a(U_c, \varphi_i(z - \pi_f z))$$
$$- a(U_{f,i}, z - \pi_f z)$$
$$\leq \left(\sum_{i \in \mathcal{N}} C_a \|h\mathcal{R}_i(U_{f,i})\|_{\omega_i}^2\right)^{1/2} \|e\|_a.$$

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Next we study the boundary part  $\pi_f z - \pi_f^0 z$ ,

$$\sum_{i \in \mathcal{N}} (f, \varphi_i(\pi_f z - \pi_f^0 z)) - a(U_c, \varphi_i(\pi_f z - \pi_f^0 z))$$
$$- a(U_{f,i}, \pi_f z - \pi_f^0 z)$$
$$\leq \left(\sum_{i \in \mathcal{N}} C_a \|\sqrt{H} \Sigma(U_{f,i})\|_{\partial \omega_i}^2\right)^{1/2} \|e\|_a.$$

 $\Sigma(U_{f,i})$  is a variational approximation of  $\partial_n U_{f,i}$ .

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$$\|e\|_a^2 \leq \sum_{i \in \mathcal{N}} C_a \left( \|h\mathcal{R}_i(U_{f,i})\|_{\omega_i}^2 + \|\sqrt{H}\Sigma(U_{f,i})\|_{\partial\omega_i}^2 \right)$$

## If we just solve local problems on part of the domain ${\mathcal F}$ we get,

$$\begin{aligned} \|e\|_a^2 &\leq \sum_{i \in \mathcal{C}} C_a \|H\mathcal{R}(U_c)\|_{\omega_i}^2 \\ &+ \sum_{i \in \mathcal{F}} C_a \left( \|h\mathcal{R}_i(U_{f,i})\|_{\omega_i}^2 + \|\sqrt{H\Sigma}(U_{f,i})\|_{\partial\omega_i}^2 \right) \end{aligned}$$

The boundary part  $\|\sqrt{H}\Sigma(U_{f,i})\|_{\partial\omega_i}^2$ , where  $\Sigma(U(f,i))$  is an approximation of  $\partial_n U_{f,i}$  decays rapidly on  $\partial\omega_i$ .



Figure 6: One two and three layer stars.

## **Adaptive Algorithm**

$$\|e\|_a^2 \leq \sum_{i \in \mathcal{N}} C_a \left( \|h\mathcal{R}_i(U_{f,i})\|_{\omega_i}^2 + \|\sqrt{H\Sigma}(U_{f,i})\|_{\partial\omega_i}^2 \right)$$

- 1. Start with given r and L where  $h = H/2^r$ .
- 2. Calculate U using AVMM.
- **3.**  $E_r^i = \|h\mathcal{R}_i(U_{f,i})\|_{\omega_i}^2 E_L^i = \|\sqrt{H}\Sigma(U_{f,i})\|_{\partial\omega_i}^2$ .
- 4. Stop if  $E_h^i$  and  $E_L^i$  are small enough else if  $E_h^i > E_L^i$  let  $r_{new} := 2r$  and if  $E_L^i > E_h^i$  let  $L_{new} = 2L$  end return to 2.

#### **Adaptive Algorithm**

#### Example using the Adaptive Algorithm,





Figure 7: Periodic *a* (0.05 blue 1 white) with  $\epsilon = H$  and  $129 \times 129$  coarse nodes.

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We introduce the dual problem: find  $\phi$  such that

$$a(v,\phi_c) + a(v,\phi_f) = (v,\psi), \text{ for all } v \in \mathcal{V}.$$

From the orthogonality showed earlier on we have  $a(e, \phi_c) = 0$ . We get

$$(e, \psi) = a(e, \phi) = a(e, \phi_f) = (f, \phi_f) - a(U, \phi_f)$$
$$= \sum_{i \in \mathcal{C}} (f, \varphi_i \phi_f) - a(U_c, \varphi_i \phi_f)$$
$$+ \sum_{i \in \mathcal{F}} (f, \varphi_i \phi_f) - a(U_c, \varphi_i \phi_f) - a(U_{f,i}, \phi_f).$$

We need to calculate an approximation of  $\phi_f$  numerically.

- If possible  $\phi_f$  can be calculated by a global calculation on a mesh where  $h < h_{\phi} < H$ .
- Or  $\phi_f$  can be calculated by AVMM. We need to keep track on neighboring patches to form  $\phi_f$  locally.



Figure 8: The dual solution on the red patches affects the error calculated on the black patch.

We consider an example where we seek a very accurate solution in part of the domain.



Figure 9: Geometry to the left with interesting region marked  $\psi = I_{\{0 \le x, y \le 0.5\}}$ . Solution to the right, a = f = 1.

We solve the dual problem with  $\psi = I_{\{0 \le x, y \le 0.5\}}$ .



Figure 10: Dual solution  $\phi$  to the left and  $\phi_f$  to the right.

We remember the error estimate,

$$(e, \psi) = \sum_{i \in \mathcal{C}} (f, \varphi_i \phi_f) - a(U_c, \varphi_i \phi_f) + \sum_{i \in \mathcal{F}} (f, \varphi_i \phi_f) - a(U_c, \varphi_i \phi_f) - a(U_{f,i}, \phi_f).$$

- 1. Start with all nodes in C. Calculate U and  $\phi_f$ .
- 2. Calculate error estimators, solve local problems or increase number of layers.
- 3. Stop if the error is small enough else go to 2.



Figure 11: Local problems are solved using two and three layer stars in marked nodes.



Figure 12: Error compared to reference mesh for standard Galerkin (left) and two iterations of AVMM as described above (right).

#### **Standard A Posteriori Error Estimates**

For duality based algorithms we know that:

- The dual solution  $\phi$  need to be approximated but not in  $\mathcal{V}_c$ .
- Regular refinement or higher order methods allocate lots of memory.

Instead we solve the dual problem by AVMM in each coarse node,

$$(e,\psi) = \sum_{i=1}^{n} (R(U), \Phi_{f,i}) + a(U, \phi_f - \Phi_f).$$

#### **Standard A Posteriori Error Estimates**

The second term can be estimated in the following way,

$$a(e, (\phi_f - \Phi_f)) \le ||e||_a ||\phi_f - \Phi_f||_a$$
  
$$\le ||e||_a ||\phi - (\Phi_c + \Phi_f)||_a.$$

Both these terms can be estimated.

We can use standard Galerkin on the primal and AVMM on the dual and have control on the error committed by not solving the dual exactly.

#### **Ongoing Projects on This Subject**

- Adaptivity for the local problems.
- More tests on how the split between  $V_c$  and  $V_f$  affects the algorithm.
- More layers than two.
- 3D implementation.