Parallel Algorithms for Elastic Systems using Multipliers-Free Domain Decomposition Methods

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Summary. This is the third article of the Minisymposium "The Multipliers-free Domain Decomposition Methods for Symmetric and Non-symmetric Matrices". In it, we apply the Multipliers-Free Domain Decomposition Method (MF-DDM) to isotropic problems of Static Elasticity. The purpose of the paper is three-fold: to exhibit the applicability of such methods to systems of equations, confirm some of the advantages that are concomitant to MF-DDMs and to establish simple procedures for developing effective codes for parallel-processing problems of elasticity. Firstly, by means a FEM formulation for Static Elasticity we derive the system-matrix associated with this kind of problems. Once this matrix is available, code development for parallel processing only requires a straightforward application of the MF-DDM matrix formulas that unify non-overlapping DDMs and were presented in paper 1 of this Minisymposium: "An Overview". The general procedures that were explained in Paper 2, "Implementation Issues", are then applied for developing the parallelprocessing codes. Such procedures can be applied straightforwardly whenever the matrix-system is available. In conclusion, this paper supplies a simple manner of developing efficient parallel codes for static elasticity, thereby demonstrating the applicability of MF-DDMs to systems of partial differential equations, and also corroborates in this particular case some of the many advantages that are concomitant to MF-DDMs.

1 Introduction

This is the third article of the Minisymposium that in the framework of DD19 was devoted to present and discuss a new methodology known as the "Multipliers-free domain decomposition methods: MF-DDMs", which is equally applicable to symmetric and non-symmetric matrices. Such methodology is based on a direct approach, without recourse to Lagrange multipliers,

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boundary conditions. Then, a weak formulation is obtained weighting Eq.2 with a vector-valued function $\underline{\psi}$ and integrating by parts:

$$A(\vec{\underline{u}},\underline{\psi}) = \int_{\vec{\Omega}} \{(\lambda+\mu)(\nabla \bullet \underline{\psi})(\nabla \cdot \underline{\vec{u}}) + \mu \nabla \underline{\psi} : \nabla \underline{\vec{u}}\} d\underline{x} = \int_{\vec{\Omega}} \underline{\psi} \bullet \underline{f}_{\vec{\Omega}} d\underline{x} \quad (3)$$

Next, a partition of the problem-domain is introduced, whose internal nodes are \underline{x}_p , p = 1, ..., N. Sometimes this partition will be referred to as the 'fine partition' because later one more partition, the 'coarse partition', will be introduced. With each node \underline{x}_p we associate a 3D-vector-valued test function to be denoted by $\underline{\psi}^{pi}$, i = 1, 2, 3. Furthermore, let $(\underline{\psi}^{pi})_j$ be the j - th component of ψ^{pi} . Then, we choose

$$\left(\underline{\psi}^{pi}\left(\underline{x}\right)\right)_{j} = \ell_{p}\left(\underline{x}\right)\delta_{ij} \tag{4}$$

Here, $\ell_p(\underline{x})$ is the Lagrange linear interpolate that is characterized by being a piecewise-linear scalar function with the property that

$$\ell_p\left(\underline{x}_q\right) = \delta_{pq}, \ p, q = 1, \dots, N \tag{5}$$

We observe that Eq.5 implies that all the test functions vanish at the nodes located on the domain-external-boundary, $\partial \dot{\Omega}$.

The set of base functions is taken to be the same as the set of test functions. Thus, we define the approximate solution of our boundary-value problem to be:

$$\frac{\overrightarrow{u}(\underline{x})}{\underline{u}(\underline{x})} \equiv \sum_{p=1}^{N} \sum_{i=1}^{3} \widehat{u}_{pi} \, \underline{\psi}^{pi}(\underline{x}) \tag{6}$$

We also define:

$$\widehat{f}_{pi} \equiv \int_{\overline{\Omega}} \underline{\psi}^{pi} \bullet \underline{f}_{\Omega} d\underline{x}, \ p = 1, ..., N \text{ and } i = 1, 2, 3$$
(7)

Eqs.6 and 3 together imply that

$$\sum_{q=1}^{N} \sum_{j=1}^{3} \widehat{A}_{qj,pi} \widehat{u}_{qj} = \widehat{f}_{pi}$$
(8)

Here:

$$\widehat{A}_{qj,pi} \equiv \int_{\widehat{\Omega}} \{ (\lambda + \mu) (\nabla \cdot \underline{\psi}^{qj}) (\nabla \cdot \underline{\psi}^{pi}) + \mu \nabla \underline{\psi}^{qj} : \nabla \underline{\psi}^{pi} \} d\underline{x}$$
(9)

Using Eq.4, it is seen that

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$$\underline{u} \bullet \underline{w} \equiv \sum_{\alpha \in \mathbb{Z}(p)} \sum_{p=1}^{N} \sum_{i=1}^{3} \frac{u\left(p, i, \alpha\right) w\left(p, i, \alpha\right)}{m\left(p\right)}, \ \forall \underline{u}, \underline{w} \in \tilde{D}\left(\bar{\Omega}\right)$$
(14)

Here $Z(p) \subset \{1, ..., E\}$ is defined by the condition that $(p, \alpha) \in \overline{\Omega}$ when $\alpha \in Z(p)$. When $\underline{\widehat{u}}, \underline{\widehat{w}} \in \overline{D}(\Omega)$ it can be seen that

$$\underline{\widehat{u}} \bullet \underline{\widehat{w}} = \tau \left(\underline{\widehat{u}}\right) \bullet \tau \left(\underline{\widehat{w}}\right)$$
(15)

The average matrix, \underline{a} , is the orthogonal projection, with respect to the Euclidean Inner product, on the subspace $\tilde{D}_{12}(\bar{\Omega})$ of continuous vectors. Its explicit expression is

$$a_{(p,i,\alpha)(q,j,\beta)} = \frac{1}{m(p)} \delta_{pq} \delta_{ij}$$
(16)

while the *jump matrix*, is $\underline{j} \equiv \underline{I} - \underline{a}$. Here, \underline{I} is the identity matrix.

4 The matrices $\underline{\underline{A}}^t$ and $\underline{\underline{A}}$

For $\gamma = 1, ..., E$, we define the linear transformations $\underline{\underline{A}}^{\gamma} : \tilde{D}(\bar{\Omega}) \to \tilde{D}(\bar{\Omega})$ by

$$A^{\gamma}_{(q,j,\beta)(p,i,\alpha)} \equiv \delta_{\alpha\gamma}\delta_{\beta\gamma} \int_{\overrightarrow{\Omega}_{\gamma}} \left\{ (\lambda+\mu)\frac{\partial\ell_p}{\partial x_i}\frac{\partial\ell_q}{\partial x_j} + \mu \sum_{r=1}^3 \frac{\partial\ell_p}{\partial x_r}\frac{\partial\ell_q}{\partial x_r}\delta_{ij} \right\} d\underline{x} \quad (17)$$

Furthermore, $\underline{\underline{A}}^{t} : \tilde{D}(\bar{\Omega}) \to \tilde{D}(\bar{\Omega})$ is defined by $\underline{\underline{A}}^{t} \equiv \sum_{\alpha=1}^{E} \underline{\underline{A}}^{\alpha}$. Next, we define the matrix $\underline{\underline{A}} : \tilde{D}(\bar{\Omega}) \to \tilde{D}(\bar{\Omega})$. To this end we choose a set $\Omega^{\pi} \subset \Omega^{\Gamma} \subset \Omega$ and define the set of 'primal nodes', $\pi \subset \Gamma \subset \bar{\Omega}$, to be defined by the condition that a *derived node*, (p, α) , belongs to π when $p \in \Omega^{\pi}$. Then, the 'dual-primal vector subspace', $\tilde{D}^{DP}(\bar{\Omega}) \subset \bar{D}(\bar{\Omega})$, is defined by requiring that its vector-members be continuous in the set π of primal nodes. Then, the transformation $\underline{a}^{\pi} : \tilde{D}(\bar{\Omega}) \to \tilde{D}(\bar{\Omega})$ is defined to be the projection of $\tilde{D}(\bar{\Omega})$ on $\tilde{D}^{DP}(\bar{\Omega})$. It can be seen that []:

$$a^{\pi}_{(p,i,\alpha)(q,j,\beta)} = \left\{ \frac{1}{m(p)} \delta_{pq} \delta^{\pi}_{pq} + \delta_{\alpha\beta} \delta_{pq} \left(1 - \delta^{\pi}_{pq} \right) \right\} \delta_{ij}$$
(18)

Here, the symbol δ_{pq}^{π} defined by

$$\delta_{pq}^{\pi} \equiv \begin{cases} 1, & \text{if } p, q \in \Omega^{\pi} \\ 0, & \text{if } p \text{ or } q \notin \Omega^{\pi} \end{cases}$$
(19)

Then, the matrix $\underline{\underline{A}}$ is defined by $\underline{\underline{A}} \equiv \underline{\underline{a}}^{\pi} \underline{\underline{Aa}}^{\pi}$. It will be assumed in what follows that π , the set of *primal nodes*, is such that $\underline{\underline{A}} : \tilde{D}^{DP}(\bar{\Omega}) \to \tilde{D}^{DP}(\bar{\Omega})$

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The Neumann – Neumann method :
$$\underline{aS}^{-1}\underline{aSu}_{\Delta} = \underline{aS}^{-1}\underline{f}_{\Delta 2}$$
 and $\underline{\underline{ju}}_{\Delta} = 0$
(28)

The non – preconditioned FETI: $\underline{\underline{S}}^{-1}\underline{\underline{j}}\underline{\underline{u}}_{\Delta} = -\underline{\underline{S}}^{-1}\underline{\underline{j}}\underline{\underline{S}}^{-1}\underline{\underline{f}}_{\Delta 2}$ and $\underline{\underline{a}}\underline{\underline{S}}\underline{\underline{u}}_{\Delta} = 0$ (29)

The preconditioned
$$FETI: \underline{\underline{S}}^{-1}\underline{\underline{j}}\underline{\underline{S}}\underline{\underline{j}}\underline{\underline{u}}_{\Delta} = -\underline{\underline{S}}^{-1}\underline{\underline{j}}\underline{\underline{S}}\underline{\underline{j}}\underline{\underline{S}}^{-1}\underline{\underline{f}}_{\Delta 2} \text{ and } \underline{\underline{a}}\underline{\underline{S}}\underline{\underline{u}}_{\Delta} = 0$$
(30)

To implement these algorithms in parallel, we simply apply the procedures of paper 2, on *implementation issues*.

5 Conclusions

By means a FEM formulation for Static Elasticity we derived the matrix system corresponding to it and, by a straightforward application of the MF-DDM, a simple procedure for developing fully parallelizable computational codes for such problems was obtained. The computational codes so derived are very robust; in particular, using object-oriented programming techniques codes that are applicable to anisotropic materials are easily constructed. Thereby, this paper confirms the applicability of the general MF-DDM matrix formulas and procedures to systems of partial differential equations, as well as other attractive features of the MF-DDMs.

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