

# Projection-based Model Order Reduction techniques

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## 1 Description of the model

Consider the problem of finding the function  $u : \Omega \rightarrow \mathbb{R}$  which satisfies the following Partial Differential Equation (PDE)

$$\begin{aligned} -\nabla \cdot \kappa \nabla u + \alpha \cdot \nabla u &= f & \text{in } \Omega, \\ u &= 0 & \text{on } \partial\Omega. \end{aligned} \quad (1)$$

The domain  $\Omega = [0, 1]^2$  is a square which is partitioned into four subdomains

$$\Omega = \begin{array}{c} 1 \\ \begin{array}{|c|c|} \hline \Omega_1 & \Omega_4 \\ \hline \Omega_2 & \Omega_3 \\ \hline \end{array} \\ 0 \qquad 0.5 \qquad 1 \end{array}$$

The diffusion field  $\kappa = \kappa(x)$  and the advection field  $\alpha = \alpha(x)$  are parametrized by means of a 5-dimensional vector  $x = (x_1, \dots, x_5)$  as follow

$$\kappa(x) = \kappa(s, x) = \sum_{i=1}^4 x_i 1_{\Omega_i}(s) \quad \text{and} \quad \alpha(x) = \alpha(s, x) = x_5 \begin{pmatrix} 1/2 - s_2 \\ s_1 - 1/2 \end{pmatrix}, \quad (2)$$

where  $s = (s_1, s_2) \in \Omega$  denotes the spatial variable and  $1_{\Omega_i}$  the indicator function<sup>1</sup> associated with  $\Omega_i$ . The parameter set is defined by

$$\mathcal{P} = \left\{ x \in \mathbb{R}^5 \text{ s.t. } \begin{array}{l} 0.05 \leq x_1, \dots, x_4 \leq 1 \\ -100 \leq x_5 \leq 100 \end{array} \right\}.$$

When it is considered at random, the parameter  $X$  is a random vector uniformly distributed over  $\mathcal{P}$ .

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<sup>1</sup>The indicator function  $1_{\Omega_i}$  is such that  $1_{\Omega_i}(s) = 1$  if  $s \in \Omega_i$  and  $1_{\Omega_i}(s) = 0$  if  $s \notin \Omega_i$ .

The source term  $f = f(s)$  is equal to  $f(s) = +1$  if  $s \in [0.05, 0.2] \times [0.4, 0.6]$ ,  $f(s) = -1$  if  $s \in [0.05, 0.2] \times [0.4, 0.6]$  and  $f(s) = 0$  elsewhere. The source term  $f$  is independent  $x$ . Finally, the scalar quantity of interest  $Y(x)$  is defined as the mean value of the solution on the subdomain  $\Omega_3$ , meaning

$$Y(x) = \frac{1}{|\Omega_3|} \int_{\Omega_3} u(x) d\Omega. \quad (3)$$

## 1.1 Weak solution and finite element approximation

The weak formulation of (1) consists in finding  $u = u(x) \in V = H_0^1(\Omega)$  such that

$$\underbrace{\int_{\Omega} \kappa(x) \nabla u(x) \cdot \nabla v \, d\Omega + \int_{\Omega} \alpha(x) \cdot \nabla u(x) v \, d\Omega}_{= a(u(x), v; x)} = \underbrace{\int_{\Omega} f v \, d\Omega}_{= \ell(v)}, \quad (4)$$

for all  $v \in V$ . We recall that the weak formulation is obtained by multiplying (1) by a test function  $v \in V$ , by integrating over  $\Omega$  and by doing an integration by part. The Finite Element Method (FEM) consists in replacing  $V$  by a finite dimension space  $V^h = \text{span}\{\phi_1, \dots, \phi_N\} \subset V$ ,  $N \gg 1$ , which consists of piecewise polynomial functions  $\phi_i$  (sometimes called “hat” functions, or “shape” functions). With this discretization, the variational formulation (4) becomes an algebraic equation

$$A(x)u(x) = b,$$

where  $A_{i,j}(x) = a(\phi_j, \phi_i; x)$ ,  $b_i = \ell(\phi_i)$  and where  $u(x)$  is now a vector containing the coefficients of the approximation on the finite element basis  $\{\phi_1, \dots, \phi_N\}$ . For more details on the FEM, see the book [1]. The quantity of interest  $Y(x)$  defined by (3) can be expressed as a linear form of the solution vector  $u(x)$  as

$$Y(x) = q^T u(x),$$

where  $q \in \mathbb{R}^N$  is the vector defined by  $q_i = \frac{1}{|\Omega_3|} \int_{\Omega_3} \phi_i d\Omega$ .

## 1.2 Affine decomposability

Using the parametrization (2) of  $\kappa(x)$  and  $\alpha(x)$ , the bilinear form  $a(\cdot, \cdot; x)$  can be written as

$$a(u(x), v; x) = \sum_{k=1}^4 x_k \underbrace{\int_{\Omega_i} \nabla u(x) \cdot \nabla v \, d\Omega}_{= a_k(u(x), v)} + x_5 \underbrace{\int_{\Omega} \left( \frac{1/2 - s_2}{s_1 - 1/2} \right) \cdot \nabla u(x) v \, d\Omega}_{= a_5(u(x), v)},$$

which, after FEM discretization, yields the decomposition

$$A(x) = \sum_{k=1}^5 x_k A_k, \quad (5)$$

where  $(A_k)_{i,j} = a_k(\phi_j, \phi_i)$ . Thus,  $A(x)$  admits an affine decomposition with respect to the parameter

## 2 Tasks

The above described benchmark is implemented in the file `parametrizedPDE.m`. It is an object oriented code which uses Matlab<sup>®</sup> classes. The file `ProjectionBasedMOR.m` provides a tutorial to learn how to use this class. Read and run the `ProjectionBasedMOR.m` one section after the other to learn how to:

- instantiate an object: `model = parametrizedPDE();`
- draw a random parameter: `X = model.randX();`
- compute the a solution: `u = model.u(X);`
- compute the quantity of interest: `Y = model.q'*u;`
- plot the solution: `model.plotSol(u)`
- get the FEM (full) system: `model.A(X)` and `model.b`
- get the matrices of the affine decomposition (5):

$$A_1 = \text{model.K1}, \dots, \quad A_4 = \text{model.K4}, \quad A_5 = \text{model.Ad},$$

### Task#1: Offline-online efficient galerkin projection

Using the affine decomposition of the operator  $A(x)$ , implement the offline-online efficient algorithm presented in the Lecture Note, see Algorithm 2. Tasks:

- Observe the speedup of the online phase and the slowdown of the offline phase.
- Estimate the computation time of the offline phases and of the online phases (per sample evaluation) of Algorithm 1 and of Algorithm 2. After how many sample evaluations does Algorithm 2 becomes more interesting to use compared to Algorithm 1?
- How this evolves with respect to  $r$ ? (repeat the previous analysis for, say,  $r \in \{5, 10, 20, 50, 100\}$ )

## Task#2: POD

In the `script.m` file, the reduced space  $V_r$  (see the matrix  $\mathbf{Vr}$ ) is built in a naive way using random snapshots of the solution. A better way to build  $V_r$  is to use the Proper Orthogonal Decomposition (POD) or the Reduced Basis (RB) method, see Sections 2 and 3 of the Lecture notes. Implement the POD (Algorithm 3). Hint: according to Remark 2.2, the function `[U,D,V]=svd(A)` can be useful...

- Compare the quality of the POD basis and the one of the naive algorithm. For instance, one can plot the  $L^2$ -error as a function of  $r \in \{5, 10, 20, 50, 100\}$ .
- Compare the  $L^2$ -error of the POD with the sum of the eigenvalues  $\sqrt{\sum_{i \geq r+1} \hat{\sigma}_i^2}$ , see Equation (7) in the Lecture notes.

## Task#3: Reduced basis

Implement and test the RB method (Algorithm 4). How does this compare to the POD in terms of number of solution evaluations? of reduced space dimension? of  $L^2$  versus  $L^\infty$  error?

## References

- [1] T. J. HUGHES, *The finite element method: linear static and dynamic finite element analysis*, Courier Corporation, 2012.