# MODEL REDUCTION FOR COUPLED AND NONLINEAR PDE SETS UNDER DIRICHLÉT BOUNDARY EXCITATIONS

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**ABSTRACT** Solutions of most Partial Differential Equations (PDEs) are dominated by coherent modes that can be extracted through the use of decomposition techniques. This paper focuses on a benchmark example for low dimensional modeling, named 2D Burgers flow. The boundary excitations are Dirichlét type and the technique results in a reduced order model of the infinite dimensional process. The issues influencing the performance are discussed and it is seen that the results obtained from the low order model are compliant with the desired results.

### INTRODUCTION

Finding a representative dynamic model for spatially continuous systems is a major problem in heat and fluid flows. The reason that lies behind is tightly associated with control system synthesis, i.e. the design of best external excitation (boundary condition) such that the the manner in which the system behaves is the prescribed behavior. From a boundary controller design point of view, it is seen that PDE systems do not enjoy the classical tools of control theory directly, instead, the PDE system is preprocessed in such a way that the essential behavior is reproduced by a set of Ordinary Differential Equations (ODE) associated with some spatial eigenfunctions. Naturally, the obtained model is an approximation to the original system dynamics. The purpose of this paper is to show how complicated PDEs could be approximated by finite dimensional ODEs that can be used for boundary control system synthesis.

The two-dimensional (2D) Burgers equation is a good example to study the difficulties encountered in low dimensional modeling of infinite dimensional systems. The reason is that the involved dynamics is governed by two coupled nonlinear PDEs. This fact indicates that the efforts towards a suitable model are to deal with a significant amount of computational complexity. Referring to [5, 1], the 2D Burgers equation is described by

$$w_t + \epsilon (w \cdot \nabla) w = \mu \nabla^2 w \tag{1}$$

where  $w(x, y, t) := (u(x, y, t) \ v(x, y, t))^{\mathsf{T}}$ , and  $(x, y, t) \in [0, 1] \times [0, 1] \times [0, T]$  with T being some final time. The similarity of the vector PDE set in (1) with the Navier-Stokes equations makes the low dimensional modeling efforts worthwhile particularly for the boundary control ap-

plications in fluid dynamics. A striking example is the problem of reducing the skin friction of air vehicles through active flow control, (See [2]). However, the 2D Burgers equation is a turbulence free cartoon for Navier-Stokes equations and has been studied in the past for modeling traffic flows, shock waves and acoustic transmission. In [14, 16, 18], some variants of 2D Burgers equation have been considered with the goal of finding exact solutions under certain circumstances. Blender, on the other hand, postulates a method to obtain the solution of the PDE set in (1) iteratively, [20]. Güngör demonstrates that if suitable subalgebras can be defined, the PDE could be converted into an ODE, but it is a major problem to find such subalgebras particularly for boundary control purposes, [9]. Nishinari et al. focus on cellular automaton, which is extensively studied for developing models of traffic flow, fluids and immune systems, [10], and therefore a good model to work on is a variant of Burgers equation. In [17], the dynamics that arises upon discretization of 2D Burgers equation is analyzed. The effects of chosen time step  $(\Delta t)$  for getting physically reasonable numerical solutions are elaborated. Wescott et al. present a computational technique to obtain the numerical solutions of PDEs having nonlinear convection terms like 2D Burgers equation and Navier-Stokes equations, [11]. The goal of the paper is to reduce the computation time without giving concessions from the accuracy. Boules et al., [6], obtain the solution for a specific boundary regime and initial conditions. Using a truncated Fourier series expansion yields an autonomous ODE set, the solution of which approximates the numerical solution, and the derived model rebuilds the situation implied by the chosen initial and boundary conditions. Except [6], the works on 2D Burgers equation emphasize the similar difficulties as the motivating factors and focus on the solutions and solvability issues. The current paper, on the other hand, derives a non-autonomous ODE model that has external inputs explicitly and that is valid for some set of boundary conditions with zero initial conditions.

When the 1D version is taken into consideration, it is seen that the 1D Burgers equation has previously been considered for modeling and control system design purposes, and it has been shown in [3] and the references therein that the task is achievable, yet there are very few results reporting the modeling issues for vector PDE sets and higher dimensional cases as emphasized above. This paper fills the gap between very simple models such as 1D heat flow or Burgers system and very complicated systems like those reported in [2, 12, 13, 15]. Seemingly, the presented work is a step to-

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wards the goal of modeling and control of more complicated PDE systems.

The current paper approaches the modeling problem from a control specialist's point of view, i.e. a suitable model reduction associated with a set of well-defined system inputs and a well-defined range of operating region. This process contains three major issues that need to be addressed appropriately. First issue is to collect the representative data and to exploit decomposition techniques for obtaining a set of ODEs. The next issue is to separate the effect of external stimuli from the other terms by utilizing the boundary conditions. The last issue is to validate the model. The process is continuous over a physical domain  $(\Omega := [0, 1] \times [0, 1])$ , the corners of which are the possible entries of external stimuli for both u(x, y, t) and v(x, y, t). Choosing an adequately dense grid, say  $\Omega_d$ , makes it possible to obtain a finite element representation of the process w(x, y, t) over  $\Omega_d$ . When the content of the observed data, say w(x, y, t), is decomposed into spatial and temporal constituents  $(u(x, y, t) \approx$  $\langle \Phi(x,y), \alpha(t) \rangle_{\Omega_d}$  and  $v(x,y,t) \approx \langle \Psi(x,y), \alpha(t) \rangle_{\Omega_d}$ , the essence of spatial behavior appears as a set of spatially varying gains  $(\Phi(x,y) = {\Phi_1(x,y), \Phi_2(x,y), \dots, \Phi_{R_L}(x,y)}$ and  $\Psi(x,y) = \{\Psi_1(x,y), \Psi_2(x,y), \dots, \Psi_{R_L}(x,y)\}$  with  $R_L$ being a positive integer), and the essence of temporal evolution,  $\alpha(t)$ , appears as the solution of a set of ODEs obtained after utilizing the orthogonality properties of the spatial basis functions, i.e. the eigenfunctions.

The dimension reduction in systems having high orders can be done by utilizing Proper Orthogonal Decomposition (POD) or Singular Value Decomposition (SVD) in cooperation with Galerkin projection [12, 13, 15] or balancing methods (for linear systems) as discussed in the survey of Gügercin, [4]. The decomposition based methods exploit the exemplar solutions obtained from the process and yield a set of temporal variables associated with a set of spatial basis functions. In order to obtain a useful approximation, the data, which is the raw information entering the modeling process, should contain coherent modes. The procedure, if it succeeds, yields a set of autonomous ODEs synthesizing the aforementioned temporal values.

The motivation of this paper is to draw a clear path between a given PDE system and the representative finite dimensional non-autonomous ODE model. With this in mind, the paper is organized as follows: The second section presents briefly the POD technique and its relevance to the modeling strategy. In the third section, obtaining the ODE model for the 2D Burgers equation is demonstrated. The justification of the model, results and the contributions are discussed in Section 4. Concluding remarks constitute the last part of the paper.

#### PROPER ORTHOGONAL DECOMPOSITION

Consider the ensemble  $P_i$ ,  $i = 1, 2, ..., N_s$ , where  $N_s$  is the number of elements. Every element of this set corresponds to a snapshot observed from a process, say for example, 2D Burgers equation given in (1) or

$$u_t = \mu u_{xx} + \mu u_{yy} - \epsilon u u_x - \epsilon v u_y$$
  
$$v_t = \mu v_{xx} + \mu v_{yy} - \epsilon u v_x - \epsilon v v_y, \qquad (2)$$

with  $\epsilon$  and  $\mu$  being known constants, and the subscripts x, y and t refer to the partial differentiation with respect to x, y and time, respectively. The continuous time process takes place over the physical domain  $\Omega := \{(x, y) | (x, y) \in$ 

 $[0, 1] \times [0, 1]$  and the solution is obtained on a grid denoted by  $\Omega_d$ , which describes the coordinates of the pixels of every snapshot in the ensemble.

The goal is to find an orthonormal basis set letting us to write the solution as

$$\begin{pmatrix} u(x, y, t) \\ v(x, y, t) \end{pmatrix} \approx \begin{pmatrix} \hat{u}(x, y, t) \\ \hat{v}(x, y, t) \end{pmatrix}$$
$$= \sum_{i=1}^{R_L} \alpha_i(t) \begin{pmatrix} \Phi_i(x, y) \\ \Psi_i(x, y) \end{pmatrix}$$
(3)

where  $\alpha_i(t)$  is the temporal part,  $\begin{pmatrix} \Phi_i(x,y) \\ \Psi_i(x,y) \end{pmatrix}$  is the spatial part,  $\begin{pmatrix} \hat{u}(x,y,t) \\ \hat{v}(x,y,t) \end{pmatrix}$  is the finite element approximate of the infinite dimensional PDE and  $R_L$  is the number of independent basis functions that can be synthesized from the given ensemble, or equivalently the set of eigenfunctions that spans the space described by the ensemble. It will later be clear that if the elements of the basis set,  $\begin{pmatrix} \Phi_i(x,y) \\ \Psi_i(x,y) \end{pmatrix}$ , are orthonormal for  $i = 1, 2, \ldots, R_L$ , then the modeling task can exploit Galerkin projection technique. More explicitly, the inner product operator defined in over the basis functions should function as,

$$\begin{pmatrix} \Phi_i \\ \Psi_i \end{pmatrix}, \begin{pmatrix} \Phi_j \\ \Psi_j \end{pmatrix} \rangle_{\Omega} :=$$
$$\iint_{\Omega} \left( \Phi_i \Phi_j + \Psi_i \Psi_j \right) \, d\Omega = \delta_{ij}$$
(4)

where  $\delta_{ij} = 1$  when i = j and zero otherwise, i.e. the Kronecker delta. With these definitions, the POD procedure can be summarized as follows:

**Step 1.** Define the concatenated process snapshot captured at time t as

 $\begin{array}{l} P_t := \left( \begin{array}{c} U_t \\ V_t \end{array} \right) \text{, where } U_t \text{ and } V_t \text{ are } R \times R, P_t \text{ is } 2R \times R \text{ and } R \text{ determines the spatial resolution. Without loss of generality, } t \text{ could be an integer that is used to index the snapshots. Start calculating the } N_s \times N_s \text{ dimensional correlation matrix } L, \text{ the } (ij)\text{-th entry of which is } L_{ij} := \langle P_i, P_j \rangle_{\Omega_d}, \text{ where } \langle \ldots, \rangle_{\Omega_d} \text{ is the inner product operator defined over the chosen spatial grid } \Omega_d. \text{ Notice that the basis vectors } \left( \begin{array}{c} \Phi_i(x,y) \\ \Psi_i(x,y) \end{array} \right) \text{ are defined over } \Omega, \text{ whereas the bases that } \text{ are obtained numerically (the sampled forms) } \left( \begin{array}{c} \phi_i \\ \psi_i \end{array} \right) \text{ are } \end{array}$ 

defined over  $\Omega_d$  and,  $\phi_i$  and  $\psi_i$  are  $R \times R$  matrices. Therefore, we need the equivalent form of the used inner product, which is given as

$$\begin{pmatrix} \begin{pmatrix} \phi_i \\ \psi_i \end{pmatrix}, \begin{pmatrix} \phi_j \\ \psi_j \end{pmatrix} \rangle_{\Omega_d} := \frac{1}{N_s} (\phi_i \star \phi_j + \psi_i \star \psi_j) = \frac{1}{N_s} \sum_{p=1}^R \sum_{q=1}^R \phi_i(p,q) \phi_j(p,q) + \psi_i(p,q) \psi_j(p,q) = \delta_{ij},$$
 (5)

where  $\star$  stands for the sum of all elements of a matrix that is obtained through elementwise multiplication of two matrices. **Step 2.** Find the eigenvectors denoted by  $\xi_i$  and the associated eigenvalues  $(\lambda_i)$  of the symmetric matrix L. Sort them in a descending order in terms of the magnitudes of  $\lambda_i$ . Note that every  $\xi_i$  is an  $N_s \times 1$  dimensional vector satisfying  $\xi_i^{\mathrm{T}} \xi_i = \frac{1}{\lambda_i}$ , here, for simplicity of the exposition, we assume that the eigenvalues are distinct.

Step 3. Construct the basis set by utilizing the snapshots

$$\begin{pmatrix} \phi_i \\ \psi_i \end{pmatrix} = \sum_{k=1}^{N_s} \xi_{ik} P_k = \sum_{k=1}^{N_s} \begin{pmatrix} \xi_{ik} U_k \\ \xi_{ik} V_k \end{pmatrix}, \quad (6)$$

where  $\xi_{ik}$  is the k-th entry of the eigenvector  $\xi_i$ , and  $i = 1, 2, ..., R_L$ , where  $R_L = \operatorname{rank}(L)$ . It can be shown that  $\langle \begin{pmatrix} \phi_i \\ \psi_i \end{pmatrix}, \begin{pmatrix} \phi_j \\ \psi_j \end{pmatrix} \rangle_{\Omega_d} = \delta_{ij}$  with  $\delta_{ij}$  being the Kronecker delta function. Notice that the basis functions are admixtures of the snapshots, [3, 13].

**Step 4.** Calculate the temporal coefficients. When  $t = t_k$ , taking the inner product of both sides of (3) with  $\begin{pmatrix} \Phi_i \\ \Psi_i \end{pmatrix}$ , the orthonormality property leads to

$$\begin{aligned}
\alpha_i(t_k) &= \langle \begin{pmatrix} \Phi_i(x,y) \\ \Psi_i(x,y) \end{pmatrix}, \begin{pmatrix} \hat{u}(x,y,t_k) \\ \hat{v}(x,y,t_k) \end{pmatrix} \rangle_{\Omega} \\
&= \langle \begin{pmatrix} \phi_i \\ \psi_i \end{pmatrix}, \begin{pmatrix} U_{t_k} \\ V_{t_k} \end{pmatrix} \rangle_{\Omega_d},
\end{aligned} \tag{7}$$

Note that the temporal coefficients satisfy orthogonality properties over the discrete set  $t_k \in \{t_1, t_2, \ldots, t_{N_s}\}$  (See (8)). For a more detailed discussion on the POD method, the reader is referred to [12, 13, 15] and the references therein,

$$\sum_{i=1}^{N_s} \langle \begin{pmatrix} U_i \\ V_i \end{pmatrix}, \begin{pmatrix} \phi_k \\ \psi_k \end{pmatrix} \rangle_{\Omega_d}^2 \approx \sum_{i=1}^{N_s} \alpha_i^2(t_i) = \lambda_k.$$
(8)

Fundamental Assumption: The majority of works dealing with POD and model reduction applications presume that the flow is dominated by coherent modes and the quantities  $\begin{pmatrix} u(x, y, t) \\ v(x, y, t) \end{pmatrix}$  and  $\begin{pmatrix} \hat{u}(x, y, t) \\ \hat{v}(x, y, t) \end{pmatrix}$  are indistinguishable, [15, 12, 13]. Because of the dominance of coherent modes, the typical spread of the eigenvalues of the correlation matrix L turns out to be logarithmic and the terms decay very rapidly in magnitude. This fact further enables to assume that a reduced order representation, say with M modes ( $M \leq \min(R_L, N_s)$ ) can also be written as an equality

$$\begin{pmatrix} \hat{u}(x,y,t)\\ \hat{v}(x,y,t) \end{pmatrix} = \sum_{i=1}^{M} \alpha_i(t) \begin{pmatrix} \Phi_i(x,y)\\ \Psi_i(x,y) \end{pmatrix}, \quad (9)$$

and the reduced order model is derived under the assumption that (9) satisfies the governing PDE set. Unsurprisingly, such an assumption results in a model having uncertainties, however, one should keep in mind that the goal is to find a model, which matches the infinite dimensional system in some sense of approximation with typically  $M \ll R_L \leq N_s$ . To represent how good such an expansion is, a percent energy measure is defined as follows

$$E = 100 \frac{\sum_{i=1}^{M} \lambda_i}{\sum_{i=1}^{R} \lambda_i},\tag{10}$$

where the tendency of  $E \rightarrow 100\%$  means that the model captures the dynamical information contained in the snapshots well. Conversely, an insufficient model will be obtained if E is far below 100%. In the next section, we demonstrate how the boundary condition is transformed to an explicit control input in the ODEs.

#### **OBTAINING THE ODE MODEL**

According to the underlying assumption of POD based model reduction scheme, the approximate solution in (9) must satisfy the PDE in (2). Substituting (9) into (2) and taking the inner product of both sides with  $\begin{pmatrix} \Phi_k \\ \Psi_k \end{pmatrix}$  yields,

$$\dot{\alpha}_{k} = \mu \sum_{i=1}^{M} \alpha_{i} \langle \begin{pmatrix} \Phi_{xxi} + \Phi_{yyi} \\ \Psi_{xxi} + \Psi_{yyi} \end{pmatrix}, \begin{pmatrix} \Phi_{k} \\ \Psi_{k} \end{pmatrix} \rangle_{\Omega} -\epsilon \sum_{i=1}^{M} \sum_{j=1}^{M} \alpha_{i} \alpha_{j} \langle \begin{pmatrix} \Phi_{i} \Phi_{xj} \\ \Phi_{i} \Psi_{xj} \end{pmatrix}, \begin{pmatrix} \Phi_{k} \\ \Psi_{k} \end{pmatrix} \rangle_{\Omega}. -\epsilon \sum_{i=1}^{M} \sum_{j=1}^{M} \alpha_{i} \alpha_{j} \langle \begin{pmatrix} \Psi_{i} \Phi_{yj} \\ \Psi_{i} \Psi_{yj} \end{pmatrix}, \begin{pmatrix} \Phi_{k} \\ \Psi_{k} \end{pmatrix} \rangle_{\Omega}.$$

$$(11)$$

Notice that, the orthonormality property of the basis vectors leaves the  $\dot{\alpha}_k$  term alone on the left hand side. Equivalently, by using the numerical quantities, the expression above can be rewritten as follows:

$$\dot{\alpha}_{k} = \mu \sum_{i=1}^{M} \alpha_{i} \langle \begin{pmatrix} \phi_{xxi} + \phi_{yyi} \\ \psi_{xxi} + \psi_{yyi} \end{pmatrix}, \begin{pmatrix} \phi_{k} \\ \psi_{k} \end{pmatrix} \rangle_{\Omega_{d}} \\ -\epsilon \sum_{i=1}^{M} \sum_{j=1}^{M} \alpha_{i} \alpha_{j} \langle \begin{pmatrix} \phi_{i} \circ \phi_{xj} \\ \phi_{i} \circ \psi_{xj} \end{pmatrix}, \begin{pmatrix} \phi_{k} \\ \psi_{k} \end{pmatrix} \rangle_{\Omega_{d}} \\ -\epsilon \sum_{i=1}^{M} \sum_{j=1}^{M} \alpha_{i} \alpha_{j} \langle \begin{pmatrix} \psi_{i} \circ \phi_{yj} \\ \psi_{i} \circ \psi_{yj} \end{pmatrix}, \begin{pmatrix} \phi_{k} \\ \psi_{k} \end{pmatrix} \rangle_{\Omega_{d}}$$

$$(12)$$

where  $\circ$  stands for the elementwise multiplication of two matrices.

Although it is straightforward to conclude with the ODEs given in (11) and (12), it is apparent that these equations do not have the boundary conditions (external inputs) explicitly. Chosen initial conditions and boundary excitation regime determine the solution, and the ODEs above resynthesize the temporal variables,  $\alpha_i(t)$ , associated to that particular solution (See (9)). It is clear that such an ODE model is useless as it is specific to the chosen boundary conditions. Our goal is to obtain a model that has external inputs explicitly and that can be used for the boundary conditions other than the used ones. For this purpose, a method needs to be postulated for separating the boundary excitations appropriately. According to the definition of the inner product operator, it should be obvious that

$$\langle \begin{pmatrix} \phi_i \\ \psi_i \end{pmatrix}, \begin{pmatrix} \phi_j \\ \psi_j \end{pmatrix} \rangle_{\Omega_d} = \langle \begin{pmatrix} \phi_i \\ \psi_i \end{pmatrix}, \begin{pmatrix} \phi_j \\ \psi_j \end{pmatrix} \rangle_{\Omega_d \setminus \partial \Omega_d} + \langle \begin{pmatrix} \phi_i \\ \psi_i \end{pmatrix}, \begin{pmatrix} \phi_j \\ \psi_j \end{pmatrix} \rangle_{\partial \Omega_d}$$
(13)

where  $\partial \Omega_d$  indicates the boundaries of the considered domain. Apparently, the last term in (13) can enjoy the boundary conditions that are specified freely. Denote  $(x_c, y_c)$  as one of the points at which the solution is independently specified (i.e. the boundary), and  $(p_{x_c}, q_{y_c})$  as the row and column numbers of this location in matrices  $\phi_i$ and  $\psi_i$ . Note that the prescribed solution in (9) must be satisfied also at  $(x_c, y_c)$ , i.e. we have,

$$\hat{u}(x_{c}, y_{c}, t) := \gamma_{x_{c}y_{c}u}(t) = \sum_{i=1}^{M} \alpha_{i}(t)\phi_{i}(p_{x_{c}}, q_{y_{c}})$$
$$\hat{v}(x_{c}, y_{c}, t) := \gamma_{x_{c}y_{c}v}(t) = \sum_{i=1}^{M} \alpha_{i}(t)\psi_{i}(p_{x_{c}}, q_{y_{c}})$$
(14)

or equivalently

$$\alpha_{k}(t)\phi_{k}(p_{x_{c}}, q_{y_{c}}) = \gamma_{x_{c}y_{c}u}(t) - \sum_{i=1}^{M} (1 - \delta_{ik}) \alpha_{i}(t)\phi_{i}(p_{x_{c}}, q_{y_{c}}),$$
(15)

$$\alpha_{k}(t)\psi_{k}(p_{x_{c}},q_{y_{c}}) = \gamma_{x_{c}y_{c}v}(t) - \sum_{i=1}^{M} (1-\delta_{ik})\,\alpha_{i}(t)\psi_{i}(p_{x_{c}},q_{y_{c}}).$$
(16)

Since we consider the problem on a square domain,  $\Omega_d$ , for both states, every corner can be a possible entry for the external excitations, i.e. we may have at most eight distinct inputs for this system. Once the Dirichlét type corner conditions are specified, the numerical solutions u(x, y, t) and v(x, y, t) on x = 0, y = 0, x = 1 and y = 1 segments of  $\Omega_d$ are obtained by setting the relevant partial derivatives to zero. For example, we solve

$$u_t = \mu u_{yy} - \epsilon v u_y$$
  
$$v_t = \mu v_{yy} - \epsilon v v_y$$
(17)

along x = 1 segment. For the simplicity of the exposition, assume  $x_c = 0$  and  $y_c = 0$   $((p_{x_c}, q_{y_c}) = (1, 1))$  is the chosen corner, and rewrite (12) as follows:

$$\dot{\alpha}_{k} = \frac{\mu}{N_{s}} \sum_{i=1}^{M} \alpha_{i} \left( \zeta_{i} \star \phi_{k} + \theta_{i} \star \psi_{k} \right)$$
$$-\frac{\epsilon}{N_{s}} \sum_{i=1}^{M} \sum_{j=1}^{M} \alpha_{i} \alpha_{j} \left( \left( \phi_{i} \circ \phi_{xj} + \psi_{i} \circ \phi_{yj} \right) \star \phi_{k} \right)$$
$$-\frac{\epsilon}{N_{s}} \sum_{i=1}^{M} \sum_{j=1}^{M} \alpha_{i} \alpha_{j} \left( \left( \phi_{i} \circ \psi_{xj} + \psi_{i} \circ \psi_{yj} \right) \star \psi_{k} \right)$$
(18)

where  $\zeta_i := \phi_{xxi} + \phi_{yyi}$  and  $\theta_i := \psi_{xxi} + \psi_{yyi}$ . Define  $\phi'_k = \{\phi'_k | \phi'_k(i,j) = \phi_k(i,j) \text{ when } i \neq p_{x_c}, j \neq q_{y_c}, \text{ and } \phi'_k(p_{x_c}, q_{y_c}) = 0\}$  and  $\zeta'_k = \{\zeta'_k | \zeta'_k(i,j) = \zeta_k(i,j) \text{ when } i \neq p_{x_c}, j \neq q_{y_c}, \text{ and } \zeta'_k(p_{x_c}, q_{y_c}) = 0\}$  and so on. By this means, the matrices used in the derivation have zero values corresponding to the external excitation

entries. Now we can explicitly write the first term in (18) as follows:

$$\sum_{i=1}^{M} \alpha_i \zeta_i \star \phi_k = \sum_{i=1}^{M} \alpha_i \left( \zeta'_i \star \phi'_k \right) + \alpha_k \zeta_k (1, 1) \phi_k (1, 1) + \sum_{i=1}^{M} (1 - \delta_{ik}) \alpha_i \zeta_i (1, 1) \phi_k (1, 1).$$
(19)

Utilizing (15) for the term  $\alpha_k \zeta_k(1,1) \phi_k(1,1)$ , we get

$$\alpha_k \zeta_k(1,1) \phi_k(1,1) = \gamma_{00u}(t) \zeta_k(1,1) - \sum_{i=1}^M (1-\delta_{ik}) \alpha_i \zeta_k(1,1) \phi_i(1,1).$$
(20)

Inserting (20) into (19) yields the following

$$\sum_{i=1}^{M} \alpha_i \zeta_i \star \phi_k = \gamma_{00u}(t) \zeta_k(1,1) + \sum_{i=1}^{M} \alpha_i \left( \zeta_i \star \phi_k - \zeta_k(1,1) \phi_i(1,1) \right).$$
(21)

For the second term in the first summation of (18), this result implies the equality in (22), and the concatenated form is given in (23),

$$\sum_{i=1}^{M} \alpha_i(\theta_i \star \psi_k) = \gamma_{00v}(t)\theta_k(1,1) + \sum_{i=1}^{M} \alpha_i\left(\theta_i \star \psi_k - \theta_k(1,1)\psi_i(1,1)\right).$$
(22)

$$\sum_{i=1}^{M} \alpha_i (\zeta_i \star \phi_k + \theta_i \star \psi_k)$$
(23)  
=  $\gamma_{00u}(t) \zeta_k(1, 1) + \gamma_{00v}(t) \theta_k(1, 1)$   
+  $\sum_{i=1}^{M} \alpha_i (\zeta_i \star \phi_k) +$   
-  $\sum_{i=1}^{M} \alpha_i (\zeta_k(1, 1) \phi_i(1, 1) + \theta_k(1, 1) \psi_i(1, 1)).$ 

The same reasoning can be applied to the terms seen in the second and third lines of (18). Due to the space limit, we skip repeating the same step. Such a separation technique lets us obtain the low dimensional model for the 2D Burgers equation given by

$$\dot{\mathcal{X}}(t) = \mathcal{A}\mathcal{X}(t) - \mathcal{B}\left(\mathcal{X}(t)\right) + \mathcal{C}\Gamma(t) - \mathcal{D}(\mathcal{X}(t), \Gamma(t)), \quad (24)$$

where  $\mathcal{X}(t) = (\alpha_1(t) \ \alpha_2(t) \ \dots \ \alpha_M(t))^{\mathsf{T}}, \ \Gamma(t) = (\gamma_{00u}(t) \ \gamma_{00v}(t))^{\mathsf{T}}, \ \mathcal{A} \text{ is } M \times M, \ \mathcal{B} \text{ is } M \times 1, \ \mathcal{C} \text{ is } M \times 2$ and  $\mathcal{D} \text{ is } M \times 1$ . From (23), we can write the (ki)-th entry of matrix  $\mathcal{A}$  and k-th row of matrix  $\mathcal{C}$  as given below:

$$(\mathcal{A})_{ki} = \frac{\mu}{N_s} \left( \zeta_i \star \phi_k + \theta_i \star \psi_k \right)$$
(25)  
$$-\frac{\mu}{N_s} \left( \zeta_k(1, 1) \phi_i(1, 1) + \theta_k(1, 1) \psi_i(1, 1) \right),$$

$$\left(\mathcal{C}\right)_{k} = \frac{\mu}{N_{s}} \left(\zeta_{k}(1,1) \quad \theta_{k}(1,1)\right).$$
(26)

where  $k, i = 1, 2, \ldots, M$ . Similarly,

$$\mathcal{B}(\mathcal{X}) = \left(\begin{array}{ccc} \mathcal{X}^{\mathsf{T}} B_1 \mathcal{X} & \mathcal{X}^{\mathsf{T}} B_2 \mathcal{X} & \dots & \mathcal{X}^{\mathsf{T}} B_M \mathcal{X} \end{array}\right)^{\mathsf{T}}$$
(27)

where the *j*-th of matrix  $B_k$  is

$$(B_k)_{ij} = \frac{\epsilon}{N_s} \left( \phi'_i \circ \phi'_{xj} \right) \star \phi'_k + \frac{\epsilon}{N_s} \left( \psi'_i \circ \phi'_{yj} \right) \star \phi'_k + \frac{\epsilon}{N_s} \left( \phi'_i \circ \psi'_{xj} \right) \star \psi'_k + \frac{\epsilon}{N_s} \left( \psi'_i \circ \psi'_{yj} \right) \star \psi'_k,$$
(28)

and the k-th row entry of vector  $\mathcal{D}$  is computed as

$$(\mathcal{D})_{k} = \gamma_{00u} \frac{\epsilon}{N_{s}} \sum_{j=1}^{M} \alpha_{j} \left( \phi_{k}(1,1)\phi_{xj}(1,1) \right) + \gamma_{00u} \frac{\epsilon}{N_{s}} \sum_{j=1}^{M} \alpha_{j} \left( \psi_{k}(1,1)\psi_{xj}(1,1) \right) + \gamma_{00v} \frac{\epsilon}{N_{s}} \sum_{j=1}^{M} \alpha_{j} \left( \phi_{k}(1,1)\phi_{yj}(1,1) \right) + \gamma_{00v} \frac{\epsilon}{N_{s}} \sum_{j=1}^{M} \alpha_{j} \left( \psi_{k}(1,1)\psi_{yj}(1,1) \right)$$
(29)

or

$$\mathcal{D} = D_u \mathcal{X} \gamma_{00u} + D_v \mathcal{X} \gamma_{00v} \tag{30}$$

where  $D_u$  and  $D_v$  are  $M \times M$  matrices and the (kj)-th entry is computed as

$$(D_u)_{kj} = \frac{\epsilon}{N_s} \left( \phi_k(1,1)\phi_{xj}(1,1) + \psi_k(1,1)\psi_{xj}(1,1) \right), \quad (31)$$

and

$$(D_v)_{kj} = \frac{\epsilon}{N_s} \left( \phi_k(1,1)\phi_{yj}(1,1) + \psi_k(1,1)\psi_{yj}(1,1) \right).$$
(32)

with k, j = 1, 2, ..., M. According to the derivation discussed in this section, once the initial and boundary conditions are specified, one can get a dynamic model that captures the essential features contained by the solution. Although we have derived the model by assuming the external excitation enters at a single point for u and v dynamics, it is straightforward to apply the scheme for obtaining a model having up to eight inputs. In the next section, justification of the model is presented through some exemplar cases.

### JUSTIFICATION OF THE MODEL

In order to obtain the model, the 2D Burgers equation in (1) is solved for the boundary conditions given as



Figure 1: Boundary signals used for model derivation



Figure 2: First set of boundary signals that are used for model validation

$$\gamma_{00u}(t) = \sin(1000\pi t(T-t)) \gamma_{00v}(t) = \cos(1000\pi t(\frac{T}{2}-t)).$$
(33)

The time plots and the Fast Fourier Transforms (FFT) of the signals above are depicted in Figure 1. The reason that drives us to choose such signals is the spectral richness. If the spectral content of the excitations are rich enough, the resulting model is more likely to operate properly over the covered frequency range, [3]. The other important parameters of the simulation are tabulated in Table 1. The numerical solution is obtained through Crank-Nicholson method with zero initial conditions, [19], and after the application of the modeling procedure discussed in the previous section, a model is obtained in the form of (24).

Table 1: Simulation Settings

R	25
M	8
$\Delta T$	0.1msec.
Т	$0.2 \mathrm{sec.}$
$N_s$	201
$\epsilon$	1
$\mu$	5



Figure 3: The desired values (thick curves) of  $\alpha_i(t)$  and the obtained values (thin curves) for the first set of test conditions

It has been observed that the eigenvalues  $(\lambda_i)$  decay very rapidly, and the captured energy content described by (10) is 99.9021%, which is found acceptable. The justification has been done with the same settings as shown in Table 1, and the first set of boundary excitations that are used in the model validation phase are

$$\gamma_{00u}(t) = \sin(700\pi t(T-t)) + 0.2\sin(1700\pi t(T-t)),$$
  

$$\gamma_{00v}(t) = \cos(500\pi t(\frac{T}{2}-t)) + 0.1\cos(1500\pi t(\frac{T}{2}-t)).$$
(34)

Figure 2 illustrates these signals and the low frequency appearance of their FFT magnitude plots. The PDE is solved for this new case and the  $\alpha_i(t)$  values are obtained by using (7), which yield the desired values. On the other hand, the model had already been developed, and it is simulated for the test boundary conditions in (34) with zero initials. The outcome is expected to approximate the desired ones if the algorithm succeeds. The results are shown in Figure 3, where the first remark of us is the number of ODEs that let us obtain this result. With M = 8 modes (ODEs), the task can be achieved to the extent seen in the figure. For the first five modes, the match is quite good yet as the mode number increases the dissimilarity between the desired and generated values become more distinguishable. Since the dominance of the corresponding modes decrease logarithmically, as seen from the figure as well, so do their effect on the overall result. Therefore the similarity of the first few modes is more substantial than the similarity of modes having high index numbers. A rough look at the eight subplots of Figure 3 altogether gives the idea of a successful approximation from a higher dimensionality to low orders, which is the goal of this paper.

We have repeated our tests for many other test signals but some of them have driven us to conclude with the relevance of model performance and spectral content of the external excitations. The descriptive nature of the signals used in the model derivation is inherited by the developed dynamical model, and the signals that do not resemble to the model derivation conditions make the system fail depending on the level of dissimilarity between the model derivation Efe

portant conclusion if the model is to be used in a feedback system synthesis. This claim has been justified by choosing signals that are spectrally similar to and dissimilar from the test conditions.

According to the results of this paper, it is fair to claim that the dynamic model in (24) functions properly up to 100 Hz. Considering the results obtained in [6], it is seen that the presented work achieves the modeling goal with a few ODEs and associated spatial eigenfunctions.

A natural issue that needs to be highlighted is the ways of improving obtained results. Expectedly, increasing the grid fineness, decreasing  $\Delta t$ , increasing the number of snapshots entering the POD procedure  $(N_s)$ , increasing the number of modes (M) are the alternatives that result in better model performance yet the price paid for this improvement is the increased computational requirements.

#### CONCLUSIONS

The research on flow control is in its infancy due to the unavailability of widely accepted and standardized frameworks for design. One central issue if the model reduction POD is one alternative among many others (see e.g. [4, 8, 7] and the references therein). Its capability of capturing the essential dynamics dominating the entire physical phenomena makes POD preferable in flow modeling and control research.

This paper focuses on the low dimensional modeling of 2D Burgers equation. The driving fact for focusing on this system is its similarity to Navier-Stokes equations, and its nonlinear, coupled and vector valued PDE nature. Once the POD algorithm is implemented, it is seen that the resulting ODE model is an autonomous one and a method to overcome this problem needs to be developed. One major contribution of this paper is on this issue, i.e. the separation of boundary terms to obtain a non-autonomous ODE model is demonstrated step by step on such a complicated system. The second contribution of the paper is its emphasis on the locality of the developed low order models. It is seen that the conditions used in the model derivation are critically important and the POD procedure has a natural propensity to build models that are valid on particular conditions. With these results, the present work advances the subject area to the establishment of a clear connection between state-space methods of control theory and complex infinite dimensional systems of PDEs. The fact that this connection is built through the numerical observations from the infinite dimensional process is worthwhile to stress the applicability of the bottom-up modeling effort for other systems of PDEs.

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