

# Model Predictive Control of Axial Dispersion Tubular Reactors with Recycle: Addressing State-delay through Transport PDEs\*

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**Abstract**—This paper presents the model predictive control of an axial tubular reactor with a recycle stream, where the intrinsic time delay imposed by the recycle stream—often overlooked in prior studies—is modeled as a transport PDE. This leads to a boundary-controlled system of coupled parabolic and hyperbolic PDEs under Danckwerts boundary conditions, ideal for this reactor type. A discrete-time linear model predictive controller is designed to stabilize the system. Utilizing Caley-Tustin time discretization along with the late lumping approach, the system’s infinite-dimensional characteristics are preserved with no need for model reduction or spatial approximation. Numerical simulations demonstrate the controller’s effectiveness in stabilizing an unstable system while satisfying input constraints.

## I. INTRODUCTION

Many chemical and petrochemical processes involve states that are distributed in space and time. These systems, a.k.a. distributed parameter systems (DPS), are often modeled using partial differential equations (PDEs) to account for distributed states dynamics. Due to the infinite-dimensional nature of DPSs, the control and estimation of these systems becomes inevitably more challenging compared to the well-developed control theories for finite-dimensional systems [1], making this field an active and fertile direction of research. Two primary methods have been proposed to approach the control of DPSs in the literature. The first one is *Early Lumping*, i.e. reducing the infinite-dimensional system to a finite-dimensional one by spatial discretization in the early stages of system modeling [2]. Approximation of the system dynamics at this stage leads to standard control strategies; however, the accuracy of the model is compromised due to potential mismatch between the original system and the reduced-order model [3]. On the contrary, the second method, *Late Lumping*, aims to preserve the infinite-dimensional nature of the system. Approximation may be made only in the final numerical implementation of the controller, resulting in a more accurate yet more complex control strategy.

Numerous studies have utilized a late lumping approach for the purpose of controlling infinite dimensional systems within the field of chemical engineering, focusing on the control of either convection-reaction systems governed by first order hyperbolic PDEs, or diffusion-convection-reaction systems governed by second order parabolic PDEs. In [4],

robust control of first order hyperbolic PDEs was addressed, where a plug flow reactor system is stabilized under a distributed input. Boundary feedback stabilization using back-stepping method is proposed in [5] for a similar system of first order hyperbolic PDEs. In [6], state feedback regulator design is proposed for a countercurrent heat exchanger system, which is another example of a chemical engineering DPS governed by first order hyperbolic PDEs other than tubular reaction systems. Introducing the effect of dispersion as a prominent aspect of axial tubular reactors, the robust control of a diffusion-convection-reaction systems governed by second order parabolic PDEs is studied in [7]. A late-lumping based approach is considered in [8] to design a low-dimensional predictive controller for a diffusion-convection-reaction system, where the dominant modes of the system are captured by modal decomposition. Similar approach has been utilized to design an observer-based model predictive controller (MPC) in [9] for an axial tubular reactor, considering the effects of recycle stream as a well-known feature of industrial chemical reactors.

Delay systems are another class of infinite-dimensional systems that have been studied in the literature [10]. Commonly represented in the form of delay differential equations (DDEs), delay can also be modeled as a transport PDE, showing to be advantageous in more complex scenarios [11]. In the field of control theory for chemical engineering DPSs, input/output delay has been addressed in several works as both output measurement delays and input actuation delays are common in industrial processes. In general, such delays can be addressed by considering a transportation lag block at either the input or output of the system, resulting in a cascade PDE system [12], [13], [14]. In contrast to input/output delays, state delay is less addressed in the relevant literature, probably since not many applications in this field can be described by state delays. In one of the few attempts, a delayed-state distributed parameter system is addressed in [15], where a full-state and output feedback regulator is designed for a system of heat exchangers. The state delay in this works comes from the time it takes for a stream to leave one pass of the heat exchanger and enter the next pass. Similarly in [16], an axial tubular reactor system is considered, where the state delay is introduced as a result of the recycle delay in the system, without considering the diffusion term along the reactor. Even in [9] where the recycle stream is considered for a distributed diffusion-convection-reaction system, the recycle is assumed to be instantaneous; a simplifying assumption that leaves a gap in the literature regarding diffusion-convection-reaction systems

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with a recycle stream imposing state delay.

In this work, an axial tubular reactor equipped with recycle is addressed as a diffusion-convection-reaction DPS. First, the reactor is modeled by a second order parabolic PDE, where the recycle stream poses a state delay, resulting in a first order hyperbolic transport PDE coupled with the original PDE. Late lumping approach is utilized by obtaining the resolvent of the infinite-dimensional system in a closed operator form, with no need to perform spatial discretization. Then, to enable the implementation of MPC as a digital controller, discrete-time representation of the system is obtained using Caley-Tustin time discretization technique; i.e. a Crank-Nicolson type of discretization that preserves the energy of the system, mitigating the need for model reduction. Finally via numerical simulations, the proposed controller is shown to stabilize an unstable system within an optimal framework, given input constraints.

## II. METHODOLOGY

### A. Model representation

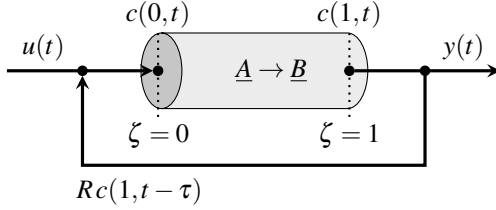


Fig. 1. Axial tubular reactor with recycle stream.

The chemical process illustrated in Figure 1 represents an axial dispersion tubular reactor, which incorporates diffusion, convection, and a first-order irreversible chemical reaction [17]. The reactor is equipped with a recycle mechanism, allowing a fraction of the product stream to re-enter the reactor to ensure the consumption of any unreacted substrate. By applying first-principle modeling through relevant mass balance relations on an infinitesimally small section of the reactor, the reactor's dynamics can be described by a second-order parabolic PDE, a common class of equations used to characterize diffusion-convection-reaction systems [18]. The resulting PDE that describes the reactor model is given by:

$$\dot{c}(\zeta, t) = D\partial_{\zeta\zeta}c(\zeta, t) - v\partial_{\zeta}c(\zeta, t) + k_r c(\zeta, t) \quad (1)$$

subject to Dankwerts boundary conditions:

$$\begin{cases} D\partial_{\zeta\zeta}c(0, t) - vc(0, t) = -v[Rc(1, t - \tau) + (1 - R)u(t)] \\ \partial_{\zeta}c(1, t) = 0 \\ y(t) = c(1, t) \end{cases} \quad (2)$$

Here,  $c(\zeta, t)$  denotes the properly scaled notion of concentration along the reactor, representing the state of the system. The physical parameters  $D$ ,  $v$ ,  $k_r$ ,  $R$ , and  $\tau$  correspond to the diffusion coefficient, flow velocity along the reactor, reaction

constant, recycle ratio, and residence time of the recycle stream, respectively. The spatial and temporal coordinates of the system are represented by  $\zeta$  and  $t$ , where  $\zeta \in [0, 1]$  and  $t \in [0, \infty)$ .

Dankwerts boundary conditions are particularly suitable for modeling axial tubular reactors, as they account for deviations from perfect mixing and piston flow, assuming negligible transport lags in connecting lines [19]. These conditions make the model more realistic for chemical reactors of this type. The input and the output of the system are also present in the boundary conditions. The system output is measured at the reactor outlet, while the input is applied at the inlet. Additionally, the delayed state resulting from the recycled portion of the flow, occurring  $\tau$  time units ago, is incorporated into the inlet; all as shown in Equation (2).

One effective method for addressing delay in systems is to represent the delay using an alternative transport partial differential equation (PDE). This approach is particularly advantageous when the problem already involves similar forms of PDEs, as is the case in the current study. To specifically address the delay in the system under consideration, the state variable  $c(\zeta, t)$  is expanded into a vector of functions  $\underline{x}(\zeta, t) \equiv [x_1(\zeta, t), x_2(\zeta, t)]^T$ , where  $x_1(\zeta, t)$  represents the concentration within the reactor, and  $x_2(\zeta, t)$  is introduced as a new state variable to account for the concentration along the recycle stream. The delay is thus modeled as a pure transport process, wherein the first state  $x_1(\zeta, t)$  is transported from the reactor outlet to the inlet, experiencing a delay of  $\tau$  time units while in the recycle stream. As a result, Equations 1 and 2 may be re-formulated as follows:

$$\partial_t \begin{bmatrix} x_1(\zeta, t) \\ x_2(\zeta, t) \end{bmatrix} = \begin{bmatrix} D\partial_{\zeta\zeta} - v\partial_{\zeta} + k_r & 0 \\ 0 & \frac{1}{\tau}\partial_{\zeta} \end{bmatrix} \begin{bmatrix} x_1(\zeta, t) \\ x_2(\zeta, t) \end{bmatrix} \quad (3)$$

$$\begin{cases} D\partial_{\zeta\zeta}x_1(0, t) - vx_1(0, t) = -v[Rx_2(0, t) + (1 - R)u(t)] \\ \partial_{\zeta}x_1(1, t) = 0 \\ x_1(1, t) = x_2(1, t) \\ y(t) = x_1(1, t) \end{cases} \quad (4)$$

With all state variables now expressed explicitly at a specific time instance  $t$ —in contrast to the previous representation where states at  $t$  were directly involved with states at  $(t - \tau)$ —the open-loop system can be described in the standard state-space form of an infinite-dimensional linear time-invariant (LTI) system as  $\dot{\underline{x}} = \mathfrak{A}\underline{x}$ . Here,  $\mathfrak{A}$  is a linear operator  $\mathcal{L}(X)$  acting on a Hilbert space  $X : L^2[0, 1] \times L^2[0, 1]$  and  $\underline{x}(\zeta, t)$ , as defined previously, is the vector of functions describing the states of the system. The operator  $\mathfrak{A}$  and its domain are defined in detail as shown in Equation (5):

$$\mathfrak{A} \equiv \begin{bmatrix} D\partial_\zeta \zeta - v\partial_\zeta + k_r & 0 \\ 0 & \frac{1}{\tau}\partial_\zeta \end{bmatrix}$$

$$D(\mathfrak{A}) = \left\{ \underline{x} = [x_1, x_2]^T \in X : \underline{x}(\zeta), \partial_\zeta \underline{x}(\zeta), \partial_{\zeta\zeta} \underline{x}(\zeta) \text{ a.c.,} \right.$$

$$D\partial_\zeta x_1(0) - vx_1(0) = -v[Rx_2(0) + (1-R)u],$$

$$\left. \partial_\zeta x_1(1) = 0, x_1(1) = x_2(1) \right\} \quad (5)$$

The eigenvalue problem for  $\mathfrak{A}$  is formulated as:

$$\mathfrak{A}\underline{\phi}_i(\zeta) = \lambda_i \underline{\phi}_i(\zeta) \quad (6)$$

where  $\lambda_i \in \mathbb{C}$  is the  $i^{\text{th}}$  eigenvalue. The characteristics equation of the system is obtained by solving the equation  $\det(\mathfrak{A} - \lambda_i I) = 0$  for  $\lambda_i$ . Attempts to analytically solve this equation has failed; therefore, it is solved numerically using the parameters in Table I. The resulting eigenvalue distribution is depicted in Figure 2 in the complex plane.

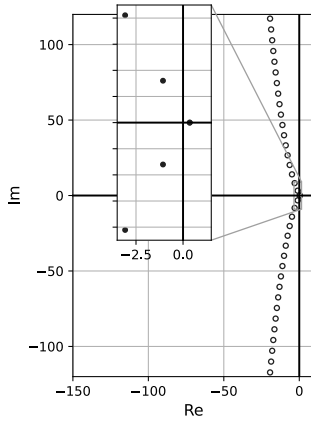


Fig. 2. Eigenvalues of operator  $\mathfrak{A}$ .

TABLE I  
PHYSICAL PARAMETERS FOR THE SYSTEM

Parameter	Symbol	Value	Unit
Diffusivity	$D$	$2 \times 10^{-5}$	$\text{m}^2/\text{s}$
Velocity	$v$	0.01	$\text{m}/\text{s}$
Reaction Constant	$k_r$	1.5	$\text{s}^{-1}$
Recycle Residence Time	$\tau$	80	$\text{s}$
Recycle Ratio	$R$	0.3	—

Following the calculation of eigenvalues, the eigenfunctions  $\{\phi_i(\zeta), \psi_i(\zeta)\}$  (for  $\mathfrak{A}$  and  $\mathfrak{A}^*$ , respectively) may be obtained. Once the adjoint system properties are determined, the eigenfunctions can be properly scaled to form a bi-orthonormal basis for the Hilbert space  $X$ ; which will be later used in the controller design.

B. Resolvent operator

C. Adjoint system

D. Caley-Tustin time discretization

E. Model predictive control design

### III. RESULTS AND DISCUSSION

#### IV. CONCLUSION

#### ACKNOWLEDGMENT

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