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<span id="page-0-6"></span>





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# A variable stepsize hybrid block optimized technique for integrating a class of singularly perturbed parabolic problems

<span id="page-0-5"></span><span id="page-0-4"></span>

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# A B S T R A C T

This paper presents and successfully applies an optimized hybrid block technique using a variable stepsize implementation to integrate a type of singularly perturbed parabolic convection–diffusion problems. The problem under consideration is semi-discretized by utilizing the method of lines. A few numerical experiments have been presented to ascertain the proposed error estimation and adaptive stepsize strategy. Furthermore, the comparison of the proposed method with other techniques in the literature is conducted via numerical experiments, and the results show that our method outperforms other existing methods.

# **1. Introduction**

Consider a singularly perturbed parabolic convection–diffusion problem (SPPCDP) that is expressed as:

$$
y_t - L_{x,\varepsilon}(t)y = f(x,t), \quad (x,t) \in \mathcal{D} = [0,1] \times [0,T],\tag{1}
$$

with

 $L_{x, \epsilon}(t)y \equiv \epsilon y_{xx} + b(x, t)y_{y} + d(x, t)y$ .

The above problem is subject to the initial and boundary conditions:



 $y(1,t) = q_1(t)$  on  $Q_1 = \{1\} \times [0, T],$  (4)

where the boundary layer is located near the boundary  $Q_x$ , and  $0 < \epsilon \ll 1$  is a perturbation parameter,  $Q_0, Q_1$ , and  $Q_x$  correspond to the left, right, and bottom boundaries of the domain  $D$ .

We note that the SPPCDP in [\(1\)](#page-0-4) is a problem with multi-scale solutions, in which some parts of the solution vary smoothly while others vary rapidly. It arises in problems involving convection and diffusion phenomena, where the governing equation involves a

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combination of diffusion, convection, and reaction terms. The parameter  $(\epsilon)$  in Eq. ([1](#page-0-4)) represents the perturbation parameter that determines the relative importance of the diffusion term compared to the convection term.

The singularly perturbed partial differential problem represented by Eqs.  $(1)$  $(1)$ – $(4)$  $(4)$  finds wide-ranging applications across diverse fields of engineering and applied sciences. This problem plays a crucial role in simulating real-life phenomena in areas such as aerodynamics, lubrication theory, heat transfer, biochemical reactions, physics, fluid dynamics, and engineering. Moreover, the formulation presented in Eq. ([1](#page-0-4)) captures various practical modeling applications, including semiconductor device modeling, turbulence models, laminar flow thermal boundary layer models, and heat flow modeling. Further insights, including the mentioned applications of SPPCDP in various science and engineering fields, can be found in [[1–](#page-9-0)[7\]](#page-9-1) and references therein. These sources provide valuable insights into the practical significance of the SPPCDP and its relevance for modeling numerous real-life problems in science-related areas and engineering domains.

As consequence of the presence of the small parameter  $(e)$  in [\(1\)](#page-0-4), the solution of (1)–([4](#page-0-5)) exhibits a boundary layer behavior, which is characterized by rapid changes in a narrow region near the domain boundary. The thickness of this layer is proportional to  $\sqrt{\epsilon}$ , which is small when  $\epsilon$  is small. The convection term  $b(x, t)y_x$  in this problem can cause an advection of the solution towards or away from the boundaries, while the diffusion term  $\epsilon y_{xx}$  can originate sharp gradients in the solution. The coefficient  $d(x, t)$ denotes a source or sink term that can affect the overall behavior of the solution. The initial and boundary conditions specify the values of  $y(x, t)$  at different parts of the boundary domain. The initial condition  $y(x, 0) = V(x)$  describes how the solution is initially distributed throughout the domain, while the boundary conditions  $y(0, t) = q_0(t)$  and  $y(1, t) = q_1(t)$  describe how the solution enters or exits through the boundaries. These features can lead to a complex behavior in the solution, such as oscillations and sharp gradients near the boundaries. Therefore, analytical or numerical approaches for integrating the SPPCDP in ([1\)](#page-0-4) and related problems must be carefully developed to capture these features and maintain and preserve stability and efficiency. For more detailed information on the theory and features of singularly perturbed differential problems, we refer the reader to [[5–](#page-9-2)[7\]](#page-9-1).

Numerous occasions necessitate applying numerical techniques for solving SPPCDP, either because an analytical solution is unknown or has no reasonable or practical meaning. Numerical strategies give a powerful alternative for solving SPPCDP.

To cite some of the available numerical methods for giving solutions to the SPPCDP in ([1\)](#page-0-4) and similar problems, we can cite a quintic B-spline method in [[8\]](#page-9-3), the parameter-robust numerical method in [[9](#page-9-4)], the piecewise-uniform mesh approach by Roos et al. [\[6\]](#page-9-5), the finite element method presented by Constantinou and Xenophontos [[10\]](#page-9-6), the layer-adapted strategy reported by Kumar [[11\]](#page-9-7), the hybrid block method proposed by Rufai [[12\]](#page-9-8), the parameter uniform difference scheme by Selvi and Ramanujam [[13\]](#page-9-9), the difference approximations in [\[14](#page-9-10)], a high order convergent numerical method in [[15\]](#page-9-11), the collocation method in [[16\]](#page-9-12), the continuous block method reported by Duromola et al. [[17\]](#page-9-13), a numerical simulation technique by Arora and Joshi [[18\]](#page-9-14), a parametric method in [\[19](#page-9-15)], the uniform and accelerated parameter numerical methods in [\[20](#page-9-16),[21\]](#page-9-17), optimized Nyström methods in [[22](#page-9-18)[,23](#page-9-19)], the finite difference method in [[24\]](#page-9-20), or the numerical methods reported in [\[25–](#page-9-21)[28\]](#page-9-22).

We should point out that most of the approaches for solving ([1\)](#page-0-4) cited above used constant step-size implementations (CSSI). Since the solution to the problem under consideration exhibits a multiscale behavior in the integration interval, most numerical methods with CSSI are expected to perform poorly, particularly when considering a large integration interval. In this manuscript, we describe a variable step size implementation utilizing an optimized hybrid block technique (OHBT) to overcome this issue and efficiently solve problem ([1](#page-0-4)).

This manuscript is organized as follows: In Section [2,](#page-1-0) the SPPCDP is converted into a system of ODEs using the method of lines. The proposed method is presented in Section [3](#page-2-0), the theoretical analysis is provided in Section [4,](#page-3-0) an error estimation of the proposed method is explained in Section [5](#page-4-0). In Section [6](#page-5-0), an implementation of the proposed method is described, in Section [7](#page-6-0), numerical experiments are discussed, and a brief conclusion is given in the last section.

#### **2. Problem transformation**

<span id="page-1-0"></span>In this section, the singularly perturbed parabolic convection-diffusion problem, as presented in Eq. ([1\)](#page-0-4), will be transformed into a set of ordinary differential equations (ODEs) using the method of lines for approximating the spatial derivatives by suitable finite difference schemes. We begin by dividing the spatial domain into a grid comprising  $N$  evenly spaced subintervals. This grid is given by  $x_i = i \Delta x$ , where  $i = 0, 1, \dots, N$ , and the spacing between adjacent points is represented by  $\Delta x = 1/N$ .

Next, we approximate the second-order derivative in space using a finite difference scheme. A common choice is the centered difference approximation:

$$
y_{xx}(x_i, t) \approx \frac{y_{i+1} - 2y_i + y_{i-1}}{4x^2}
$$

where  $y_i = y(x_i, t)$ .

 $\mathbf{v}$ 

Similarly, we approximate the first-order derivative in space using the centered finite difference scheme:

$$
y_x(x_i, t) \approx \frac{y_{i+1} - y_{i-1}}{2\Delta x}
$$

Substituting these approximations into the original PDE yields:

*.*

$$
\frac{dy_i}{dt} = \epsilon \frac{y_{i+1} - 2y_i + y_{i-1}}{\Delta x^2} + b(x_i, t) \frac{y_{i+1} - y_{i-1}}{2\Delta x} + d(x_i, t)y_i + f(x_i, t).
$$

We see that through the method of lines, for each interior point at the spatial grid, we obtain the following system of ODEs for  $i = 1, 2, \ldots, N - 1$ :

$$
\frac{dy_1}{dt} = \frac{\epsilon(y_2 - 2y_1 + y_0)}{\Delta x^2} + \frac{b(x_1, t)(y_2 - y_0)}{2\Delta x} + d(x_1, t)y_1 + f(x_1, t),
$$
\n
$$
\frac{dy_{N-1}}{dt} = \frac{\epsilon(y_N - 2y_{N-1} + y_{N-2})}{\Delta x^2} + \frac{b(x_{N-1}, t)(y_N - y_{N-2})}{2\Delta x} + d(x_{N-1}, t)y_{N-1} + f(x_{N-1}, t),
$$

where the values  $y_0$  and  $y_N$  are specified by the boundary conditions.

The vector form of the above system can be expressed as follows:

<span id="page-2-3"></span>
$$
\frac{dy}{dt} = Ay + f(t),
$$
\n(5)

where  $y = [y_1, \dots, y_{N-1}]^T$  is a vector of the numerical approximations to  $y(x_i, t)$  at time t, A is an  $(N - 1) \times (N - 1)$  matrix given by

$$
\mathbf{A} = \frac{\epsilon}{\Delta x^2} \mathbf{C} + \frac{1}{2\Delta x} \mathbf{B} + \mathbf{D},
$$

where **C**, **B** are  $(N - 1) \times (N - 1)$  tridiagonal matrices given by

$$
C = \begin{bmatrix} -2 & 1 & 0 & \cdots & 0 & 0 \\ 1 & -2 & 1 & \cdots & 0 & 0 \\ 0 & 1 & -2 & \cdots & 0 & 0 \\ \vdots & \vdots & \vdots & \ddots & \vdots & \vdots \\ 0 & 0 & 0 & \cdots & -2 & 1 \\ 0 & 0 & 0 & \cdots & 1 & -2 \end{bmatrix},
$$
  
\n
$$
B = \begin{bmatrix} 0 & b(x_1, t) & 0 & \cdots & 0 & 0 \\ -b(x_2, t) & 0 & b(x_2, t) & \cdots & 0 & 0 \\ 0 & -b(x_3, t) & 0 & \cdots & 0 & 0 \\ \vdots & \vdots & \vdots & \ddots & \vdots & \vdots \\ 0 & 0 & 0 & \cdots & 0 & b(x_{N-2}, t) \\ 0 & 0 & 0 & \cdots & -b(x_{N-1}, t) & 0 \end{bmatrix},
$$

and **D** is a diagonal matrix with entries given by the coefficients  $d(x_i, t)$ ,  $i = 1, 2, ..., N - 1$ . Finally,

$$
\mathbf{f}(t) = \left[ \left( \frac{\epsilon}{\Delta x^2} - \frac{b(x_1, t)}{2\Delta x} \right) y_0 + f(x_1, t), f(x_2, t), \dots, f(x_{N-2}, t), \left( \frac{\epsilon}{\Delta x^2} + \frac{b(x_{N-1}, t)}{2\Delta x} \right) y_N + f(x_{N-1}, t) \right]^T
$$

is a vector related to the boundary conditions and the evaluation of  $f$  at selected spatial grid points. Furthermore, the initial values at  $t = 0$  can be incorporated into the system of ODEs by specifying appropriate values for the entries of y at time  $t = 0$  at the grid points of the spatial domain. The resulting system of ODEs can then be solved numerically using an ODE solver to obtain the numerical solution  $y(t)$  at each time step. The solution  $y(t)$  can then be used to approximate the solution  $y(x_i, t)$  of the original PDE at time t. For solving the resulting ODE system we will consider the block method derived in the following section.

#### **3. Derivation of the proposed block method**

<span id="page-2-0"></span>Following the optimization procedure in [[29,](#page-9-23)[30](#page-9-24)], we begin the derivation of the block method by assuming that the solution of a first order differential equation of the form  $y'(t) = f(t, y(t))$  can be approximated by the following polynomial  $\rho(t)$ , with  $t_{n+1} = t_n + h$ and *ℎ* is the stepsize

$$
y(t) \simeq \rho(t) = \sum_{k=0}^{4} a_k t^k,
$$
 (6)

whose first derivative is

<span id="page-2-2"></span><span id="page-2-1"></span>
$$
y'(t) \simeq \rho'(t) = \sum_{k=1}^{4} a_k k t^{k-1},\tag{7}
$$

where the coefficients  $a_k$  represent unknown values that must be determined when collocation conditions are imposed at suitable points.

The OHBT method is developed based on the optimization of two points, namely  $t_{n+c_1} = t_n + c_1 h$  and  $t_{n+c_2} = t_n + c_2 h$ , within the interval  $[t_n, t_{n+1}]$ . We proceed to evaluate the equation in ([6\)](#page-2-1) at  $t_n$  and the equation in [\(7\)](#page-2-2) at  $t_n$ ,  $t_{n+c_1}$ ,  $t_{n+c_2}$ , and  $t_{n+1}$ . This leads to a system of five equations comprising the five unknowns,  $a_k$  for  $k = 0, 1, \ldots, 4$ . The equations are as follows: The first equation sets the value of the function  $\rho$  at the point  $t_n$  equal to the approximation of the solution  $y_n$ . The second equation involves the first derivative of the function  $\rho$  and evaluates it at four different points. Specifically, we approximate the derivative of the solution  $y$  at  $t_n$  by  $f_n$ , and we approximate it at three other points, namely  $t_{n+c_1}$ ,  $t_{n+c_2}$ , and  $t_{n+1}$ , by  $f_{n+c_1}$ ,  $f_{n+c_2}$ , and  $f_{n+1}$ , respectively. In this context,  $y_{n+j}$  and  $f_{n+j}$  are approximations of the solution  $y(t_{n+j})$  and its derivative  $f(t_{n+j}, y(t_{n+j}))$ , respectively, where j is either 0,  $c_1$ ,  $c_2$ , or 1. The system of equations is solved for the unknowns  $a_k$ , and the matrix form of the obtained system is

$$
\begin{pmatrix}\n1 & t_n & t_n^2 & t_n^3 & t_n^4 \\
0 & 1 & 2t_n & 3t_n^2 & 4t_n^3 \\
0 & 1 & 2t_{n+c_1} & 3t_{n+c_1}^2 & 4t_{n+c_1}^3 \\
0 & 1 & 2t_{n+c_2} & 3t_{n+c_2}^2 & 4t_{n+c_2}^{3} \\
0 & 1 & 2t_{n+1} & 3t_{n+1}^2 & 4t_{n+1}^3 \\
0 & 1 & 2t_{n+1} & 3t_{n+1}^2 & 4t_{n+1}^3 \\
0 & 1 & 2t_{n+1} & 3t_{n+1}^2 & 4t_{n+1}^3\n\end{pmatrix}\n\begin{pmatrix}\na_0 \\
a_1 \\
a_2 \\
a_3 \\
a_4\n\end{pmatrix}\n=\n\begin{pmatrix}\ny_n \\
f_n \\
f_{n+c_1} \\
f_{n+c_2} \\
f_{n+1}\n\end{pmatrix}.
$$

Once the values of  $a_k$ ,  $k = 0(1)4$  have been obtained, we can modify the variable *t* as  $t = t_n + zh$ , and represent the polynomial in ([6](#page-2-1)) using  $\alpha_0(z)$  and  $\beta_i(z)$ , where  $i = 0, c_1, c_2, 1$ . Then, the polynomial in (6) may be rewritten as

<span id="page-3-1"></span>
$$
\rho(t_n + zh) = \alpha_0(z)y_n + h\left(\beta_0(z)f_n + \beta_{c_1}(z)f_{n+c_1} + \beta_{c_2}(z)f_{n+c_2} + \beta_1(z)f_{n+1}\right),\tag{8}
$$

where  $\alpha_0(z) = 1, \{\beta_i(z)\}\$  $_{i=0,c_1,c_2,1}$  and the values of  $\beta_i(z)$  are dependent on  $c_1$  and  $c_2$ . We optimize the local truncation error (LTE) of [\(8](#page-3-1)) in order to obtain optimal values of  $c_1$  and  $c_2$  for approximating  $y(t_{n+1})$ . This LTE is given by

<span id="page-3-2"></span>
$$
L[y(t_{n+1}), h] = \frac{(5c_2 + c_1 (5 - 10c_2) - 3) h^5 y^{(5)}(t_n)}{1440} + \mathcal{O}(h^6).
$$
\n(9)

By setting the leading term in the LTE  $(9)$  to zero, we arrive at the following equation

<span id="page-3-3"></span>
$$
5c_2 + c_1 (5 - 10c_2) - 3 = 0,\tag{10}
$$

which we can solve to obtain a relationship between  $c_1$  and  $c_2$ . Specifically, we can express  $c_2$  in terms of  $c_1$  as

$$
c_2 = \frac{5c_1 - 3}{10c_1 - 5}.\tag{11}
$$

It is worth noting that there are infinitely many pairs of values for  $c_1$  and  $c_2$  that satisfy Eq. ([10\)](#page-3-3) and the condition  $0 < c_1 < c_2 < 1$ . However, by optimizing the LTE in [\(9\)](#page-3-2) for approximating  $y(t_{n+1})$ , we can choose a specific pair of values for  $c_1$  and  $c_2$  that yields an accurate approximation. In this case, one such optimized pair is given by  $c_1 = \frac{1}{3}$  and  $c_2 = \frac{4}{5}$ .

Substituting  $z = 1$  in Eq. ([8](#page-3-1)) yields the formula for approximating the solution  $y(t_{n+1})$ . This is because  $t_{n+1} = t_n + h$ , and setting  $z = 1$  corresponds to evaluating the polynomial at  $t_{n+1}$ . The resulting formula is given as follows:

<span id="page-3-4"></span>
$$
y_{n+1} = y_n + \frac{h}{336} \left( 35f_n + 162f_{n+c_1} + 125f_{n+c_2} + 14f_{n+1} \right).
$$
 (12)

We also proceed to evaluate the polynomial  $\rho(t)$  at the points  $t_{n+c_1}, t_{n+c_2}$  (that is, taking  $z = c_1, c_2$  in ([8](#page-3-1))) to get the three formulas that form the new OHBT. Here are the remaining formulas:

$$
y_{n+c_1} = y_n + \frac{h}{9072} \left( 1169f_n + 2214f_{n+c_1} - 625f_{n+c_2} + 266f_{n+1} \right),
$$
  
\n
$$
y_{n+c_2} = y_n + \frac{2h}{2625} \left( 133f_n + 648f_{n+c_1} + 325f_{n+c_2} - 56f_{n+1} \right).
$$
\n(13)

Formulas [\(12](#page-3-4))–[\(13](#page-0-6)) are the block method that will be used to solve the ordinary differential system in ([5\)](#page-2-3).

#### **4. Theoretical analysis**

<span id="page-3-0"></span>Here, the mathematical properties and theoretical aspects of the OHBT are discussed in detail. The analysis in this section provides a rigorous understanding of the OHBT method and its performance, which is essential for its practical application. The analysis includes the calculation of the local truncation errors (LTEs), the determination of the order of accuracy, and the discussion of stability and convergence properties.

#### *4.1. Local truncation errors, consistency, and convergence of the OHBT*

The formulas in Eqs.  $(12)$  $(12)$ – $(13)$  $(13)$  can be written as a recurrence relation as follows:

$$
\bar{R}Y_n = h\,\bar{S}\,Y_n',\tag{14}
$$

in which the constant matrices  $\bar{R}$ ,  $\bar{S}$  represent the coefficients of the formulas ([12\)](#page-3-4)–([13\)](#page-0-6), and the vectors  $Y_n$  and  $Y'n$  represent the values of the solution and its derivative at the nodes  $t_n$ ,  $t_{n+c_1}$ ,  $t_{n+c_2}$ , and  $t_{n+1}$ .

To analyze the error of the OHBT method, we define an operator  $L$  that is related to the method presented in Eqs.  $(12)$  $(12)$ – $(13)$  $(13)$ . This operator L takes as input the function  $y(t)$  and the stepsize *h* as follows:

$$
L[y(t_n);h] = \sum_{j\in I} \left[\Theta_j y\left(t_n + jh\right) - h\Psi_j y'\left(t_n + jh\right)\right],\tag{15}
$$





<span id="page-4-1"></span>

where the terms  $\Theta_j$  and  $\Psi_j$  are vector columns within the matrices  $\bar{R}$  and  $\bar{S}$ , respectively. The set *I* is defined as  $0, c_1, c_2, 1$ , which are the indices used in the OHBT method. Assuming that  $y(t)$  is sufficiently differentiable, we expand  $y(t_n + jh)$  and  $y'(t_n + jh)$  in a Taylor series about the point  $t_n$ , where *j* represents the index in the set *I*. Doing this, we get

$$
L[y(t_n); h] = \theta_0 y(t_n) + \theta_1 h y'(t_n) + \theta_2 h^2 y''(t_n) + \dots + \theta_q h^q y^q(t_n) + \dots,
$$
\n(16)

where

$$
\theta_q = \frac{1}{q!} \left[ \sum_{j \in I}^{k} j^q \Theta_j - q \sum_{j \in I}^{k} j^{q-1} \Psi_j \right], \ q = 0, 1, 2, \dots
$$
\n(17)

By applying the definitions of order and LTEs as provided in  $[17,31]$  $[17,31]$  $[17,31]$ , we can derive the order  $(p)$  and expressions of the LTEs for the obtained formulas.

[Table](#page-4-1) [1](#page-4-1) presents the LTEs and orders of the obtained formulas. It is important to note that the OHBT is a one-step method; therefore, OHBT is a zero-stable method. Additionally, the method is consistent because the order  $p$  is greater than 1. Furthermore, the OHBT is a convergent method due to its zero-stability and consistency.

# *4.2. Linear-stability analysis*

Following the guidelines of the method proposed by [[31\]](#page-9-25), we analyze the linear stability of the proposed optimized method. Applying the proposed OHBT to  $y' = vy$ ,  $Re(v) < 0$ , we get

$$
Y_{n+1} = M(z)Y_n, \quad z = \nu h,\tag{18}
$$

where v is a complex number and  $M(z)$  is indeed the stability matrix, which can be expressed as

$$
M(z) = \left(C^0 + zE\right)^{-1} \left(C^1 + zD\right),\tag{19}
$$

with

$$
C^{0} = \begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{bmatrix}, \qquad C^{1} = \begin{bmatrix} 0 & 0 & 1 \\ 0 & 0 & 1 \\ 0 & 0 & 1 \end{bmatrix},
$$
  
\n
$$
E = \begin{pmatrix} 0 & 0 & 1 + \frac{167}{1296} \\ 0 & 0 & 1 + \frac{38}{375} \\ 0 & 0 & 1 + \frac{5}{48} \end{pmatrix},
$$
  
\n
$$
D = \begin{pmatrix} 1 - \frac{41}{168} & \frac{625}{9072} & -\frac{19}{648} \\ -\frac{432}{875} & 1 - \frac{26}{105} & \frac{16}{375} \\ -\frac{27}{56} & -\frac{125}{336} & 1 - \frac{1}{24} \end{pmatrix}.
$$

The obtained eigenvalues of the stability matrix  $M(z)$  for the OHBT are  $\left\{0,0,\frac{z^3+15z^2+84z+180}{-2z^3+21z^2-96z+180}\right\}$ .

We note that the above eigenvalues are essential because the stability of the numerical solution depends on those eigenvalues. Evaluating the stability of the OHBM requires considering the spectral radius, denoted as  $\rho(M(z))$ , which provides valuable insights into the stability behavior. The region where the absolute value of the spectral radius is less than one, that is,  $\mathcal{S}$  =  ${z \in \mathbb{C} : |\rho(M(z))| < 1}$  ${z \in \mathbb{C} : |\rho(M(z))| < 1}$  ${z \in \mathbb{C} : |\rho(M(z))| < 1}$ , represents the resulting region of absolute stability. [Fig.](#page-5-1) 1 visually represents the stability region of the OHBT method, revealing that the entire left half of the complex plane lies within this region. This observation indicates that the OHBT method is A-stable, signifying its ability to solve the differential problem under consideration.

# **5. Error estimation and control strategy of the OHBT**

<span id="page-4-0"></span>A variable stepsize formulation (VSSF) for the OHBT will now be presented in this section. Using VSSF, we can obtain a more efficient numerical approximation of the transformed singularly perturbed problem presented in Eq. ([5\)](#page-2-3). The introduction of this manuscript emphasizes the importance of changing the stepsize to obtain a more reliable solution to a differential problem. We used



<span id="page-5-3"></span><span id="page-5-2"></span>Fig. 1. The region in the complex z-plane where the OHBT method is absolutely stable.

<span id="page-5-1"></span>an embedding-like strategy to estimate the local truncation error (LTE) robustly. The lower-order approach used for error estimation of the OHBT is as follows:

$$
y_{n+1}^* = y_n + \frac{h}{56} \left( 7f_n + 24f_{n+c_1} + 25f_{n+c_2} \right),\tag{20}
$$

with  $LTE = \frac{h^4 y^{(4)}(t_n)}{1080} + O(h^5)$ . Thus, the local error is estimated through the quotient

$$
EST = \frac{\|y_{n+1}^* - y_{n+1}\|}{\left(1 + \|y_{n+1}\|\right)}.\tag{21}
$$

Before continuing our calculations, it is necessary to verify that the absolute value of the estimated error (EST) is equal to or less than the specified relative tolerances (RET). The results are accepted if  $EST \leq RET$ , while the stepsize is increased to minimize the computational cost. If *EST* > RET, the results are rejected, and the stepsize is adjusted using the following stepsize selection strategy to improve the accuracy

$$
h_{\text{vary}} = \eta h_{\text{old}} \left( \frac{RET}{\|EST\|} \right)^{1/(p+1)}.
$$
\n(22)

In Eq. ([22\)](#page-5-2),  $p = 3$  represents the order of the lower order approach in ([20](#page-5-3)), and we prevent failed steps by using a safety factor  $\eta$ . In our numerical experiment, we set  $\eta$  to 0.90, and we initialized the stepsize that will be changed by the algorithm if needed using the method described earlier to ensure accuracy and efficiency.

# **6. Implementation of the proposed method**

<span id="page-5-0"></span>Eq. [\(1\)](#page-0-4) is discretized using approximations for  $y_x$  and  $y_{xx}$  at  $(x_i, t)$  as given in Section [2.](#page-1-0) Then, we rewrite the system in [\(12](#page-3-4))–[\(13](#page-0-6)) for the discretized problem obtained from [\(5\)](#page-2-3) as  $F(y) = 0$ , where the unknowns are represented by the following vector  $\tilde{Y}$ 

$$
\tilde{\mathbf{Y}} = \left(y_{1,n+c_1}, y_{2,n+c_1}, \dots, y_{N-1,n+c_1}, y_{1,n+c_2}, y_{2,n+c_2}, \dots, y_{N-1,n+c_2}, y_{1,n+1}, y_{2,n+1}, \dots, y_{N-1,n+1}\right).
$$

We use Newton's method (NM) to solve the obtained nonlinear equations. The NM is an iterative method that starts with an initial guess of the solution and then updates the guess by computing the solution of a linear system at each iteration until convergence. For the OHBT scheme, the NM takes the form of

$$
\tilde{Y}^{i+1} = \tilde{Y}^i - \left(J_0\right)^{-1} F^i,
$$

 $\overline{a}$ 

where  $\tilde{Y}^i$  is the current guess of the solution,  $F^i$  is the corresponding nonlinear function evaluated at  $\tilde{Y}^i$ , and  $J_0$  is the frozen Jacobian matrix of F evaluated at  $\tilde{Y}$ <sup>i</sup>. To obtain a starting point for the NM iterations, we can use an approximation of the solution based on the values of the previous time step. In the OHBT scheme, we use the values of  $y_n$  and  $f_n$  (which are known) to compute the starting values for the unknowns in each iteration, given by  $y_{n+j} = y_n + (jh)f_n$ , where *h* is the time step size and  $j = c_1, c_2, 1$  correspond to the three different sets of unknowns in the system. These starting values help speed up the convergence of the NM iterations, reducing the computational cost of solving the system at each time step. Following is an outline of the steps that are taken for the proposed method to be used in integrating problem ([5](#page-2-3)):

- 1. Initialize the algorithm by setting the initial stepsize  $(h)$ , the initial condition  $(t_0, y_0)$ , the number of spatial subintervals N, and the end-point of the integration interval  $T$ .
- 2. Define the function  $f(t, y(t))$  and initialize  $t_n$  and  $y_n$ .
- 3. Check if the current value of  $t_n$  is greater than or equal to the end-point T. If true, then exit the algorithm.
- 4. If  $t_n + h$  is greater than T, reduce the stepsize ( $h$ ) to ensure that the next step does not exceed the end-point.
- 5. While  $t_n$  is less than T, solve the system of equations in  $(12)$  $(12)$ – $(13)$  $(13)$  using the current step size  $h_{old}$  to obtain the value of  $y_{n+1}$ .
- 6. Compute an estimate of the error using  $EST = \frac{||y_{n+1}^* y_{n+1}||}{(1 + ||y_{n+1}||)}$  $\frac{n+1}{(1+\|y_{n+1}\|)}$ .
- 7. If *EST* is less than or equal to the user-defined relative tolerance RET, accept the results and increase the stepsize (h) by a factor of 2. Set  $t_n$  to  $t_n + h$ , set  $y_n$  to  $y_{n+1}$ , and go to step 3.
- 8. If *EST* is greater than *RET*, reject the results, compute a new stepsize (*h*) using Eq. [\(22](#page-5-2)), and go back to step 5.

The above algorithm iteratively computes an approximate solution on a discrete mesh by adapting the stepsize based on the error estimate. The algorithm increases the stepsize if the error is small and reduces the stepsize if the error is large. The algorithm terminates when the end-point of the integration interval is reached. The user-defined relative tolerance  $RET$  controls the accuracy of the solution. The smaller the RET, the more accurate the solution, but the longer the algorithm takes to compute. The stepsize (*h*) is critical to the accuracy and efficiency of the algorithm. A large stepsize may result in inaccurate solutions, while a small stepsize may result in excessive computational time. Therefore, the algorithm uses an adaptive stepsize for better accuracy and efficiency.

#### **7. Numerical experiments**

<span id="page-6-0"></span>Some test problems are presented here to evaluate the effectiveness of the proposed OHBT method for solving the problem under consideration. These problems are used to compare the numerical results obtained by the OHBT method with those obtained by other existing numerical methods. By comparing the OHBT with other methods, one can determine the accuracy and reliability of the proposed method. In order to compare the accuracy of different methods, the maximum absolute error (MAE) is calculated using the following formula:

$$
MAE = \max_{i=0,...,N, j=0,...,N} ||y(x_i, t_j) - Y(x_i, t_j)||_{\infty},
$$

where  $y(x_i, t)$  represents the reference solution, and  $Y(x_i, t_j)$  is the solution obtained by the OHBT method at each time point  $t_j$ .

In the Tables and Figures presented in this paper, we use the following abbreviations to represent different methods and performance measures. The optimized hybrid block technique proposed in this manuscript is denoted by OHBT. The implicit Gauss– Legendre method of order six formulated in variable stepsize is represented by G6P and is cited from [[32\]](#page-9-26). The uniformly convergent B-spline method is denoted by UCBSM and is described in [[4](#page-9-27)]. The higher-order uniformly convergent method is referred to as HOMCM and is reported in [[33\]](#page-9-28).

We also use various performance measures to evaluate the accuracy and efficiency of the methods. The number of steps taken by the solver is denoted by NOS, while the number of rejected steps is represented by RS. The number of Newton iterations used is denoted by NI, and the total number of Jacobian evaluations is denoted by TNJ. The total number of function evaluations is represented by FE. The initial step size is denoted by *h<sub>ini</sub>*. The computational time taken by the solver is denoted by CT and measured in seconds. Finally, we also compare our proposed method OHBT with the built-in package solver in Mathematica, which we refer to as NDSolve.

## *7.1. Example 1*

Consider the following SPPCDP described in [[4](#page-9-27)]

<span id="page-6-1"></span>
$$
\epsilon y_{xx} + x^p y_x - y_t - y = x^2 - 1, (x, t) \in D = [0, 1] \times [0, 1],
$$
  
\n
$$
y(0, t) = 1 + t^2, y(1, t) = 0, y(x, 0) = (1 - x)^2.
$$
\n(23)

The problem has boundary conditions specified at  $x = 0$  and  $x = 1$ , and an initial condition specified at  $t = 0$ . However, the exact solution to this problem is not known.

# *7.2. Example 2*

Consider the SPPCDP discussed in [\[4\]](#page-9-27)

<span id="page-6-2"></span>
$$
\epsilon y_{xx} + x^p y_x - y_t - (x + p)y = p(x^2 - 1) \exp(-t), (x, t) \in \mathcal{D} = [0, 1] \times [0, 1],
$$
  
\n
$$
y(0, t) = 1 + t^2, y(1, t) = 0, y(x, 0) = (1 - x)^2.
$$
\n(24)

It should be noted that no exact solution has been found for the above problem.

#### **Table 2**

<span id="page-7-0"></span>Comparison of numerical results for Problem [\(23](#page-6-1)) with  $h_{ini} = 10^{-1}, \epsilon = 2^{-8}, p = 1, N = 29$ .



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<span id="page-7-1"></span>**Table 3**<br>Comparison of numerical results for Problem [\(23](#page-6-1)) with  $h_{ini} = 10^{-2}$ ,  $\varepsilon = 2^{-16}$ ,  $p = 1$ ,  $N = 29$ .





<span id="page-7-2"></span>**Fig. 2.** Surface solution of Problem [\(23\)](#page-6-1) obtained with NDSolve and the approximate solution obtained with the OHBT method (red dots) using  $h_{ini} = 10^{-2}$ ,  $\epsilon =$  $2^{-7}$ ,  $p = 3$ .







# *7.3. Discussion of results*

The numerical results obtained using the suggested OHBT approach are compared to those obtained using the G6P, UCBSM, and HOMCM methods. We used NDSolve to provide the reference solution in order to determine MAE for numerical Experiments 1 and 2. Notably, NDSolve is a numerical method for solving ordinary or partial differential equations implemented in Mathematica. [Tables](#page-7-0) [2](#page-7-0)[–5](#page-8-0) show that the numerical results obtained with our current approach is more accurate than the G6P method. The reported MAEs for Problems [\(23](#page-6-1))–[\(24](#page-6-2)) with  $NOS = 128$ ,  $p = 1$ ,  $\epsilon = 2^{-8}$  and  $NOS = 128$ ,  $p = 1$ ,  $\epsilon = 2^{-10}$  for the B-spline based method in [\[4\]](#page-9-27)

#### **Table 5**

<span id="page-8-0"></span>Comparison of numerical results for Problem [\(24](#page-6-2)) with  $h_{ini} = 10^{-2}$ ,  $\epsilon = 2^{-20}$ ,  $N = 29$ .

<b>RET</b>		Methods	<b>NOS</b>	<b>RS</b>	NI	FE	<b>TNJ</b>	$\sqrt{\text{T}}$	<b>MAE</b>
$10^{-5}$ $10^{-5}$		OHBT G6P	12 22		26 456	48 1332	13 228	1.4063 59.4844	$6.1794 \times 10^{-7}$ $3.7334 \times 10^{-4}$
$10^{-5}$ $10^{-5}$	10 10	OHBT G6P	13 271		30 554	52 1626	15 277	1.6563 77.0000	$1.5295 \times 10^{-7}$ $3.9088 \times 10^{-4}$



<span id="page-8-1"></span>**Fig. 3.** Surface solution of Problem [\(24\)](#page-6-2) obtained with NDSolve and the approximate solution obtained with the OHBT method (red dots) using  $h_{i,j} = 10^{-1}$ ,  $\epsilon =$  $2^{-8}$ ,  $p = 6$ ,  $N = 29$ .

are  $3.96300 \times 10^{-5}$  and  $5.4298 \times 10^{-5}$ , respectively, while for the OHNT with  $NOS = 10$ ,  $p = 1$ ,  $\varepsilon = 2^{-8}$  and  $NOS = 9$ ,  $p = 1$ ,  $\varepsilon = 2^{-10}$ are 1*.*2579 × 10−7 and 1*.*2934 × 10−7, respectively, confirming that OHNT method performs better than the B-spline technique in [[4\]](#page-9-27).

In addition, from [Tables](#page-7-1) [3](#page-7-1) and [5](#page-8-0), we can see the MAEs for Problems ([23\)](#page-6-1)–([24\)](#page-6-2) with different values of  $p$  (specifically  $p = 2$  and  $p = 6$ ). We emphasize that the numerical techniques presented in [[4,](#page-9-27)[33\]](#page-9-28), using the same values of  $p$ , report MAEs of 4.7157 × 10<sup>-5</sup>, 2.2990 × 10<sup>-4</sup> and 4.3795 × 10<sup>-4</sup>, 8.4190 × 10<sup>-4</sup>, respectively, for Problem ([23\)](#page-6-1), while for the OHNT method, the obtained MAEs for the same  $p$  values are  $7.3002 \times 10^{-6}$  $7.3002 \times 10^{-6}$  $7.3002 \times 10^{-6}$ ,  $4.9189 \times 10^{-6}$ , respectively, indicating the superior performance of the OHBT method. [Figs.](#page-7-2) [2–](#page-7-2)3 display surface plots of the proposed OHBT method and NDSolve on Problems  $(23)$  $(23)$ – $(24)$  $(24)$  to demonstrate the good agreement between the OHBT method and NDSolve solutions.

# **8. Conclusion**

This paper introduces a new approach to solve a singularly perturbed parabolic convection–diffusion problem. The method is called the optimized hybrid block technique (OHBT), and we present a formulation with variable stepsizes to improved accuracy. To evaluate the effectiveness of OHBT, we compare its results with those obtained using other existing numerical methods. The comparison is based on the numerical solutions of the problem in Eqs. ([1](#page-0-4))–[\(4\)](#page-0-5). The reported numerical solutions in [Tables](#page-7-0) [2](#page-7-0)[–5](#page-8-0) and [Figs.](#page-7-2) [2](#page-7-2)[–3](#page-8-1) indicate that the proposed OHBT outperforms different existing numerical techniques utilized for comparison. Future research work could investigate the possibility of solving the time-dependent partial differential of Lane–Emden–Fowler equations using the type of pairwise methods reported in [[34,](#page-9-29)[35\]](#page-9-30).

#### **Declaration of competing interest**

We declare that the work presented has not been published previously, and it is not under consideration for publication elsewhere, that authors approve its publication and that it will not be published elsewhere in the same form, in English or any other language, including electronically without the written consent of the copyright holder.

We confirm that the authors have no conflict of interest.

#### **Data availability**

No data was used for the research described in the article.

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