# NEWTON'S METHOD APPLIED TO NONLINEAR BOUNDARY VALUE PROBLEMS: A NUMERICAL APPROACH

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## ABSTRACT

This work investigates the application of the Newton's method for the numerical solution of a nonlinear boundary value problem formulated through an ordinary differential equation (ODE). Nonlinear ODEs arise in various mathematical modeling contexts, where an exact solution is often unfeasible due to the intrinsic complexity of these equations. Thus, a numerical approach is employed, using Newton's method to solve the system resulting from the discretization of the original problem. The procedure involves the iterative formulation of the method, which enables the approximation of solutions and the evaluation of convergence with respect to the problem parameters. The results demonstrate that Newton's method provides a robust and efficient solution, highlighting its applicability to complex boundary value problems and reinforcing its relevance for the numerical analysis of nonlinear systems. It is concluded that the methodology discussed is suitable for solving a wide range of boundary value problems, ensuring precision and stability in the results.

Keywords Newton's method  $\cdot$  Boundary value problems  $\cdot$  Nonlinear ordinary differential equations  $\cdot$  Numerical solution.

# 1 Introduction

Ordinary differential equations (ODEs) play a central role in mathematical modeling, being used in fields ranging from physics and engineering to biology and economics [1, 2]. In particular, boundary value problems for ODEs appear in various applications, such as the description of heat transfer phenomena, fluid mechanics, and population growth [3]. However, the exact solution of these equations is not always possible, especially when the equation is nonlinear, which requires numerical methods to obtain approximate solutions [4].

Among numerical methods, Newton's method has been widely used due to its efficiency and rapid convergence, especially in nonlinear systems [5]. In the context of nonlinear ODEs, it becomes a powerful tool for solving boundary value problems by transforming the initial problem into a system of nonlinear equations, whose solution is iteratively approximated [6]. This approach is particularly relevant in situations where the behavior of the solution is complex and requires refinement of successive approximations.

The present work aims to apply Newton's method to solve a nonlinear boundary value problem, based on the formulation proposed by [4]. The differential equation addressed represents a typical example of a nonlinear problem, and the boundary conditions impose additional challenges that require robust numerical strategies. Using an interval discretization and iterative formulations, this study presents an approximate solution to the proposed problem, along with a detailed analysis of the convergence and accuracy of the results.

Thus, the objective of this work is to demonstrate the effectiveness of Newton's method in solving nonlinear boundary value problems, providing a practical approach and a theoretical basis that can be applied to a wide range of similar problems. The analysis of the results obtained illustrates the potential of this method in providing stable and reliable numerical solutions for complex ODEs, reinforcing its applicability and importance in the field of numerical analysis.

## 2 Newton's Method

Newton's method is a highly efficient iterative technique for solving nonlinear equations and finding the roots of such functions. It is applicable to both single-variable functions and systems with multiple variables. The main idea is to replace the nonlinear function with a local linear approximation to obtain a better estimate of the root.

## 2.1 Newton's Method in One Dimension

For a single-variable function f(x), Newton's method uses the tangent line to the curve of the function f at an initial point  $x_0$ , which should be relatively close to the true root  $\alpha$ . The iterative formula to obtain the next approximation  $x_{n+1}$  is given by:

$$x_{n+1} = x_n - \frac{f(x_n)}{f'(x_n)}.$$

In this case,  $f'(x_n)$  represents the derivative of f(x) evaluated at  $x_n$ . This process approximates each subsequent point through the tangent, allowing for rapid convergence towards the root  $\alpha$ , provided that  $x_0$  is sufficiently close to it and that  $f'(x_n) \neq 0$  in all iterations.

If  $x_0$  is far from  $\alpha$  or if  $f'(x_0)$  is close to zero, the method may diverge, as illustrated in Figure 1. The method has a quadratic convergence rate when  $x_0$  is chosen appropriately [7].



Figure 1: Failure of the Newton Method due to a poor initial point [7].

#### 2.2 Newton's Method for Nonlinear Systems

The concept of Newton's method can be extended to solve systems of nonlinear equations, where we seek simultaneous solutions for multiple functions of multiple variables. Consider a system with two nonlinear equations:

$$f_1(x_1, x_2) = 0,$$
  
 $f_2(x_1, x_2) = 0.$ 

Newton's method for this system begins with an initial estimate  $(x_1^{(0)}, x_2^{(0)})$ . In each iteration, each function is approximated by its tangent plane, and the solution of the resulting linear system provides the next approximation of

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the root. The vector formulation to obtain this new estimate is:

$$J(\mathbf{X}^{(i)})\Delta\mathbf{X}^{(i)} = -\mathbf{F}(\mathbf{X}^{(i)})$$

where  $J(\mathbf{X}^{(i)})$  is the Jacobian matrix of the system, composed of the partial derivatives of the functions with respect to the variables. In the case of two variables, the Jacobian matrix is:

$$J(\mathbf{X}^{(i)}) = \begin{bmatrix} \frac{\partial f_1}{\partial x_1} & \frac{\partial f_1}{\partial x_2} \\ \frac{\partial f_2}{\partial x_1} & \frac{\partial f_2}{\partial x_2} \end{bmatrix}$$

After solving this system for  $\Delta \mathbf{X}^{(i)} = (\Delta x_1^{(i)}, \Delta x_2^{(i)})$ , we update the solution:

$$x_1^{(i+1)} = x_1^{(i)} + \Delta x_1^{(i)}, \quad x_2^{(i+1)} = x_2^{(i)} + \Delta x_2^{(i)}.$$

This process is repeated until the vector  $\mathbf{F}(\mathbf{X}^{(i)})$  is close to zero or the variations in the variables are sufficiently small, indicating convergence. For a system with *n* equations and *n* variables, the generalized notation is:

$$J(\mathbf{X})\Delta\mathbf{X} = -\mathbf{F}(\mathbf{X}),$$

where  $J(\mathbf{X})$  is the  $n \times n$  Jacobian matrix. The solution update is:

$$\mathbf{X}^{(i+1)} = \mathbf{X}^{(i)} + \Delta \mathbf{X}^{(i)}.$$

This extension of Newton's method to nonlinear systems is powerful, but it requires that the Jacobian matrix be nonsingular and that the initial solution choice be close to the true value to ensure quadratic convergence [7].

As discussed, Newton's method involves iteratively solving a system of nonlinear equations using linear approximations. The process can be described by the following pseudocode:

Algorithm 1: Pseudocode for Newton's Method for Nonlinear Systems

**Input:** *F*: The nonlinear function we want to solve **Input:** JF: The function that calculates the Jacobian of F (can be empty) **Input:**  $\mathbf{X}_0$ : The initial guess vector for the solution **Input:** *tol*: The tolerance for the stopping criterion Input: maxit: The maximum number of allowed iterations Function Newton's Method:  $\mathbf{X} \leftarrow \mathbf{X}_0$ ; 2 iter  $\leftarrow 1$ ; 3  $h \leftarrow 10^{-6}$  // Small step for finite differences 4 while iter < maxit do 5 if JF is empty then 6  $J \leftarrow$  calculate Jacobian numerically using F, **X**, and h; 7 else 8  $J \leftarrow \text{call function } JF \text{ with } \mathbf{X};$ 9  $\Delta \mathbf{X} \leftarrow -J^{-1}F(\mathbf{X}) / /$  Newton step 10  $\mathbf{X}_n \leftarrow \mathbf{X} + \Delta \mathbf{X}$  // Update the solution 11  $err \leftarrow maximum$  absolute value of  $(\mathbf{X}_n - \mathbf{X}) / /$  Calculate the error 12 13 if err < tol then | return  $\mathbf{X}_n$  // Convergence achieved 14 else 15  $| \mathbf{X} \leftarrow \mathbf{X}_n;$ 16  $iter \leftarrow iter + 1;$ 17 Error: Newton's method did not converge; 18 **return** Last calculated solution  $\mathbf{X}_n$ 19

# **3** Problem Definition

This study is based on the work of [4], which explores the solution of boundary value problems for nonlinear ordinary differential equations. These problems are relevant in various applications, from modeling physical phenomena to

engineering and material science. The complexity of nonlinear problems, particularly in the context of fixed boundary conditions, requires robust numerical approaches to find reliable approximate solutions.

In this paper, we consider the following boundary value problem involving a nonlinear ordinary differential equation:

$$y'' = \frac{1}{8}(32 + 2x^3 - yy'), \quad 1 \le x \le 3$$
<sup>(1)</sup>

The imposed boundary conditions are:

$$y(1) = 17, \quad y(3) = 14.333333$$
 (2)

This problem describes the relationship between the dependent variable y, its first derivative y', and its second derivative y'', subject to specific values of y at x = 1 and x = 3. The nonlinear function of the derivative and the function itself makes the analytical solution unattainable, motivating the application of numerical methods for a satisfactory approximation.

To solve the problem, we choose to discretize the interval [1,3] using a step size h = 0.1, which generates a sequence of uniformly spaced points  $x_i$ . This approach allows for the formulation of a system of nonlinear equations from the discretization of the differential equation, which will then be solved iteratively using Newton's method. The choice of h = 0.1 is a compromise between accuracy and computational cost, ensuring that the resulting system is sufficiently detailed to capture the solution's dynamics while remaining computationally feasible.

## 4 Numerical Method

The numerical solution of the proposed boundary value problem is carried out using the interval discretization method and an iterative approximation technique. The steps of the adopted method are detailed below.

#### 4.1 Step 1: Interval Discretization

The first step consists of discretizing the interval [1, 3], where the value of x varies between 1 and 3. To do this, we divide the interval into N + 1 = 20 subintervals of equal width, with a step size h = 0.1. This means that the distance between two consecutive points  $x_i$  and  $x_{i+1}$  will be 0.1. Each value  $x_i$  is calculated by:

$$x_i = a + i \times h, \quad i = 0, 1, 2, \dots, 20$$
 (3)

where a = 1 is the lower bound of the interval, and h = 0.1 is the step size. The resulting values of  $x_i$  for each *i* are presented in Table 1, which illustrates the discretization of the interval.

i	$x_i$	
0	1.0	
1	1.1	
2	1.2	
3	1.3	
4	1.4	
5	1.5	
6	1.6	
7	1.7	
8	1.8	
9	1.9	
10	2.0	
11	2.1	
12	2.2	
13	2.3	
14	2.4	
15	2.5	
16	2.6	
17	2.7	
18	2.8	
19	2.9	
20	3.0	
for the	e inter	val $[1, 3]$ with $h = 0.1$
	$ \begin{array}{r} i\\ 0\\ 1\\ 2\\ 3\\ 4\\ 5\\ 6\\ 7\\ 8\\ 9\\ 10\\ 11\\ 12\\ 13\\ 14\\ 15\\ 16\\ 17\\ 18\\ 19\\ 20\\                                   $	$\begin{array}{c ccccccccccccccccccccccccccccccccccc$

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Tab

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## 4.2 Step 2: Definition of Boundary Conditions

The boundary conditions for the problem are provided as values of y at the interval's endpoints. In this case, we have:

 $w_0 = 17$  and  $w_{20} = 14.333333$ 

These values are used as starting points for the numerical solution and will be applied in the iteration process.

## 4.3 Step 3: Initial Approximation

To start the iterative process, an initial approximation  $w^{(0)}$  for the values of y at all discretized points  $x_i$  must be defined. The initial approximation is given by:

	(15.6666)	t	
	15.6666		
	15.6666		
	15.6666		
	15.6666		
	15.6666		
	15.6666		
	15.6666		
	15.6666		
$w^{(0)} =$	15.6666		(4)
	15.6666		
	15.6666		
	15.6666		
	15.6666		
	15.6666		
	15.6666		
	15.6666		
	15.6666		
	\15.6666/	1	

The initial approximation for the values  $w_i$ , with i = 1, 2, ..., 19, was obtained from the average of the boundary values  $w_0 = 17$  and  $w_{20} = 14.33$ , as defined by the boundary conditions.

## 4.4 Step 4: Definition of the Nonlinear System

The nonlinear system that describes the problem consists of 19 equations, which relate the values of  $w_i$  to their respective approximations. Each equation in the system F(w) = 0 has the general form:

$$F(w) = \begin{pmatrix} 2w_1 - w_2 + 0.01 \left(4 + 0.33275 + \frac{w_1(w_2 - 17)}{1.6}\right) - 17 &= 0\\ -w_1 + 2w_2 - w_3 + 0.01 \left(4 + 0.432 + \frac{w_2(w_3 - w_1)}{1.6}\right) &= 0\\ -w_2 + 2w_3 - w_4 + 0.01 \left(4 + 0.432 + \frac{w_3(w_3 - w_2)}{1.6}\right) &= 0\\ -w_3 + 2w_4 - w_5 + 0.01 \left(4 + 0.686 + \frac{w_4(w_5 - w_4)}{1.6}\right) &= 0\\ -w_4 + 2w_5 - w_6 + 0.01 \left(4 + 0.84375 + \frac{w_5(w_6 - w_4)}{1.6}\right) &= 0\\ -w_5 + 2w_6 - w_7 + 0.01 \left(4 + 1.024 + \frac{w_6(w_7 - w_5)}{1.6}\right) &= 0\\ -w_6 + 2w_7 - w_8 + 0.01 \left(4 + 1.22825 + \frac{w_7(w_8 - w_6)}{1.6}\right) &= 0\\ -w_7 + 2w_8 - w_9 + 0.01 \left(4 + 1.458 + \frac{w_8(w_0 - w_7)}{1.6}\right) &= 0\\ -w_8 + 2w_9 - w_{10} + 0.01 \left(4 + 1.71475 + \frac{w_9(w_{10} - w_8)}{1.6}\right) &= 0\\ -w_1 + 2w_{11} - w_{12} + 0.01 \left(4 + 2.31525 + \frac{w_{11}(w_{12} - w_{10})}{1.6}\right) &= 0\\ -w_{11} + 2w_{12} - w_{13} + 0.01 \left(4 + 2.662 + \frac{w_{12}(w_{13} - w_{11})}{1.6}\right) &= 0\\ -w_{13} + 2w_{14} - w_{15} + 0.01 \left(4 + 3.04175 + \frac{w_{14}(w_{15} - w_{13})}{1.6}\right) &= 0\\ -w_{15} + 2w_{16} - w_{17} + 0.01 \left(4 + 4.390625 + \frac{w_{15}(w_{16} - w_{13})}{1.6}\right) &= 0\\ -w_{16} + 2w_{17} - w_{18} + 0.01 \left(4 + 4.92075 + \frac{w_{17}(w_{18} - w_{16})}{1.6}\right) &= 0\\ -w_{16} + 2w_{17} - w_{18} + 0.01 \left(4 + 5.488 + \frac{w_{18}(w_{17} - w_{15})}{1.6}\right) &= 0\\ -w_{18} + 2w_{19} + 0.01 \left(4 + 6.09725 + \frac{w_{19}(w_{13} - w_{13})}{1.6}\right) - 14.333333 &= 0 \end{pmatrix}$$

This system is derived from the discretization of the original differential equation, considering interactions between adjacent points  $w_i$  and the boundary conditions. The objective is to find a numerical solution that satisfies this system of equations.

## **4.5** Step 5: Calculation of the Jacobian J(w)

The Jacobian J(w) of the nonlinear system is a matrix that describes the partial derivatives of F(w) with respect to each component  $w_i$ . This matrix is used in the Newton method to update the solution approximations. The Jacobian matrix J(w) is structured as follows:

	$\left[2 + 0.01\left(\frac{w_2 - 17}{1.6}\right)\right]$	$-1 + 0.05(\frac{1}{8}w_1)$	0	• • •	•••	• • •		0
J(w) =	$-1 - 0.05(\frac{1}{8}w_2)$	$2 + 0.01(\frac{w_3 - w_1}{1.6})$	$-1 + 0.05(\frac{1}{8}w_2)$	0	•••	•••		0
	0	$a_{32}$	a <sub>33</sub>	$a_{34}$	0	•••	•••	0
	÷	0	·	·	۰.	0		0
	÷	÷	0	·	·	·	0	0
	÷	÷	÷	0	·	·.	·	0
	÷	÷	÷	÷	0	۰.	·	$a_{i-1j}$
	L 0	0	0	0	0	0	$a_{i,j-1}$	$a_{ij}$

The Jacobian matrix is essential to the iterative process, as it is used to calculate the corrections needed for the  $w_i$  values in each iteration.

# 5 Results

The iterative method was applied successively until the difference between consecutive iterations became sufficiently small, indicating the solution's convergence. The convergence criterion was set as  $||w^{(k)} - w^{(k-1)}|| \le \epsilon$ , where  $\epsilon$  is a predefined tolerance value, signifying that the solution achieved satisfactory accuracy. Table 2 displays the  $w^{(k)}$  values for each iteration k, until convergence was reached, with the condition  $||w^{(4)} - w^{(3)}|| = 0$  demonstrating that the w values in successive iterations stabilized.

$x_i$	$w_i$	$\mathbf{w}^{(0)}$	$\mathbf{w}^{(1)}$	$\mathbf{w}^{(2)}$	$\mathbf{w}^{(3)}$	$\mathbf{w}^{(4)}$
1.0	$w_0$	17.0000	17.0000	17.0000	17.0000	17.0000
1.1	$w_1$	15.6666	16.7657	16.7605	16.7605	16.7605
1.2	$w_2$	15.6666	16.5183	16.5135	16.5134	16.5134
1.3	$w_3$	15.6666	16.2664	16.2589	16.2589	16.2589
1.4	$w_4$	15.6666	16.0102	15.9974	15.9974	15.9974
1.5	$w_5$	15.6666	15.7504	15.7299	15.7298	15.7298
1.6	$w_6$	15.6666	15.4879	15.4578	15.4577	15.4577
1.7	$w_7$	15.6666	15.2240	15.1830	15.1829	15.1829
1.8	$w_8$	15.6666	14.9609	14.9084	14.9083	14.9083
1.9	$w_9$	15.6666	14.7011	14.6376	14.6375	14.6375
2.0	$w_{10}$	15.6666	14.4483	14.3751	14.3750	14.3750
2.1	$w_{11}$	15.6666	14.2070	14.1267	14.1266	14.1266
2.2	$w_{12}$	15.6666	13.9835	13.8994	13.8993	13.8993
2.3	$w_{13}$	15.6666	13.7852	13.7019	13.7018	13.7018
2.4	$w_{14}$	15.6666	13.6220	13.5444	13.5443	13.5443
2.5	$w_{15}$	15.6666	13.5060	13.4392	13.4391	13.4391
2.6	$w_{16}$	15.6666	13.4525	13.4010	13.4010	13.4010
2.7	$w_{17}$	15.6666	13.4804	13.4475	13.4475	13.4475
2.8	$w_{18}$	15.6666	13.6132	13.5999	13.5999	13.5999
2.9	$w_{19}$	15.6666	13.8801	13.8843	13.8843	13.8843
3.0	$w_{20}$	14.3333	14.3333	14.3333	14.3333	14.3333

Table 2: Iteration results for the value of  $w_i$ 

# 6 Conclusion

After applying the iterative method with the Jacobian matrix, an approximate solution to the boundary value problem was obtained through successive iterations. The approximate values for  $w_i$  across the discretized interval are given by the sequence:

$$\overline{\mathbf{w}} = \begin{pmatrix} 17.0000, \\ 16.7605, \\ 16.5134, \\ 16.2589, \\ 15.9974, \\ 15.7298, \\ 15.4577, \\ 15.1829, \\ 14.9083, \\ 14.6375, \\ 14.3750, \\ 14.1266, \\ 13.8993, \\ 13.7018, \\ 13.5443, \\ 13.4391, \\ 13.4010, \\ 13.4475, \\ 13.5999, \\ 13.8843, \\ 14.3333 \end{pmatrix}$$
(6)

Each value  $w_i$  represents an approximation of  $y(x_i)$ , with  $x_i = a + i \times h$  and i = 0, 1, 2, ..., N + 1, where h is the spacing between discretized points. The Newton method, when applied to the system of discretized differential equations, effectively converged to a solution for this boundary value problem, as demonstrated by the stabilization of  $w_i$  values in the final iterations.

Below, we present a graph illustrating the evolution of the vector  $\overline{\mathbf{w}}$  over the iterations. This graph visually demonstrates how the values  $w_i$  stabilize, reflecting the convergence of the method to the approximate solution.



Figure 2: Plot of the vector  $\overline{\mathbf{w}}$  over the discretized interval.

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