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A Modified Spectral Relaxation Method for Some Emden-Fowler Equations

Gerald Tendayi Marewo

Abstract

In this chapter, we present a modified version of the spectral relaxation method for solving singular initial value problems for some Emden-Fowler equations. This study was motivated by the several applications that these equations have in Science. The first step of the method of solution makes use of linearisation to solve the model problem on a small subinterval of the problem domain. This subinterval contains a singularity at the initial instant. The first step is combined with using the spectral relaxation method to recursively solve the model problem on the rest of the problem domain. We make use of examples to demonstrate that the method is reliable, accurate and computationally efficient. The numerical solutions that are obtained in this chapter are in good agreement with other solutions in the literature.

Keywords: Emden-Fowler equations, Lane-Emden equations, singular initial value problem, spectral relaxation method, numerical method

1. Introduction

The singular initial value problem

$$\begin{aligned} \frac{d^2y}{dx^2} + \frac{\gamma}{x} \frac{dy}{dx} + r(x, y) &= s(x), 0 < x, \gamma < \infty \\ y(0) &= \alpha, \frac{dy}{dx}(0) = 0, \alpha \in \mathbb{R} \end{aligned} \quad (1)$$

for the Lane-Emden Eq. (1) models several phenomena such as the thermal behaviour of a spherical cloud of gas acting under the mutual attraction of its molecules [1], the temperature variation of a self gravitating star, the kinetics of combustion [2], thermal explosion in a rectangular slab [3] and the density distribution in isothermal gas spheres [4]. Moreover, Eq. (1) has been used many a time as a benchmark for new methods.

A particular case of Eq. (1) is the Emden-Fowler equation of the first kind:

$$\frac{d^2y}{dx^2} + \frac{2}{x} \frac{dy}{dx} + y^m = 0, y(0) = 1, \frac{dy}{dx}(0) = 0, m \in \mathbb{N} \quad (2)$$

As mentioned in [5], Eq. (2) represents the dimensionless form of the governing equation for the gravitational potential of a Newtonian self-gravitating, spherically

symmetric, polytropic fluid. The equation provides a useful approximation for stars.

A more general form for Eq. (2) is the Emden-Fowler equation

$$\frac{d^2y}{dx^2} + \frac{\gamma}{x} \frac{dy}{dx} + f(x)g(y) = 0 \quad (3)$$

which can be written as

$$(py')' + qy = h(x, y, y') \text{ where } ()' = \frac{d}{dx} (), \quad (4)$$

an equation which was discussed in [6]. An existence result for the solution is given therein under certain conditions on $p(x)$, $q(x)$ and $h(x, y, y')$.

Exact solutions are available for particular cases of Eq. (3) [7], but not for the general case according to the best of our knowledge. This is motivation enough for seeking approximate solutions. To this end several approximate analytical methods were used by other researchers to solve Eq. (3). Van Gorder [8] made use of the Homotopy analysis method (HAM) and its variant, the Optimal homotopy analysis method, to solve a boundary value problem for the Lane-Emden equation of the second kind. The two respective analytical solutions that they obtained were in strong agreement. The Homotopy perturbation method (HPM) is another variant of the HAM that was used by Chowdhury and Hashim [9] to solve an initial value problem for Eq. (3). Their analytical solutions were the same as those that were obtained by Wazwaz [10] using the Adomian decomposition method (ADM). Chowdhury and Hashim observed that the HPM was less computationally expensive than the ADM. Wazwaz [11] made use of the variational iteration method (VIM) to solve both initial value problems and boundary value problems for Eq. (3) and for some inhomogeneous Emden-Fowler equations. The results that they obtained demonstrated the reliability and effectiveness of the VIM.

Some numerical methods have been used by other researchers to approximate solutions to Eq. (3). Many of these numerical methods fall in the class of *collocation methods*. Examples of these collocation methods include but are not limited to the Chebyshev wavelet finite difference method (CWFD) [12], the Haar wavelet collocation method (HWCM) [13], the Taylor wavelet method (TWM) [14] and the Radial basis function - differential quadrature method (RDF-DQM) [15]. One distinct feature of these collocation methods is the choice of collocation points for discretizing the problem domain. Another distinct feature is the choice of basis functions that are used either for constructing numerical solutions or for numerical differentiation. The CWFD makes use of Chebyshev-Gauss-Lobatto collocation points and Chebyshev wavelet finite difference basis functions. The HWCM makes use of the collocation points

$$x_j = (j - 0.5)/2M, j = 1, \dots, 2M$$

on the problem domain $[0, L]$, and the method uses integrals of Haar wavelets as basis functions. The TWM uses roots of shifted Legendre polynomials as collocation points, and as basis functions the method uses Taylor wavelets which are special functions that defined in terms of Taylor polynomials. Convergence results for the Taylor wavelet solution were presented in [14]. The RDF-DQM uses collocation points

$$x_j = \frac{2}{L} \left(1 - \cos \left(\frac{j-1}{N-1} \right) \right), j = 1, \dots, N$$

on the problem domain $[0, L]$ and the method makes use of Radial basis functions. Unlike making use of collocation methods for solving Eq. (3) Van Gorder and Vajravelu [1] used the Runge–Kutta–Fehlberg 4-5 (RKF45) method to validate the analytical solutions that they obtained from using the HAM and from using the traditional power series method. The RKF45 method is an embedded Runge–Kutta–pair which makes use of an adaptive stepsize to control the method and to ensure stability properties such as A -stability. See [16] for more details on the RKF45 method.

In this chapter we make use of a modified version of the spectral relation method (SRM) to solve an initial value problem for Eq. (3). We denote our method by MSRM. The SRM was successfully used to solve fluid flow problems by for example Motsa [17], Motsa *et al.* [18], Shateyi *et al.* [19] and Gangadhar *et al.* [20] to mention a few. In [17–20] the SRM was shown to be accurate, computationally efficient and easy to implement. Moreover, the SRM was applied only after transforming the governing partial differential equations to ordinary differential equations. However, not so long ago the SRM was modified in such a way that it was directly applicable to partial differential equations. See for example [21]. The SRM was used to solve other types of problems such as hyperchaotic systems [22]. It is to the best of our knowledge that the MSRM has not been used in existing literature. We chose the MSRM because it is not computationally intensive and it is easy to implement.

In Section 2 we describe the MSRM for the model problem. In Section 3 we make use of examples to demonstrate the accuracy and computational efficiency of the MSRM. Section 4 concludes this chapter.

2. The MSRM for the model problem

We seek an approximate solution to

$$y'' + \frac{\gamma}{x}y' + f(x)g(y) = 0, 0 < x \leq L, y(0) = \alpha, y'(0) = 0 \quad (5)$$

where $\gamma > 0, L, \alpha$ are given constants and f and g are given functions.

We follow the idea behind the solution method by Ramos [23], where the singularity at $x = 0$ is isolated in a sufficiently small subinterval $I_\varepsilon = [0, \varepsilon]$ of $I = [0, L]$ where $\varepsilon > 0$. The point $x = \varepsilon$ splits interval I into two subintervals: I_ε and $I - I_\varepsilon = [\varepsilon, L]$. A linearisation method is used to solve Eq. (5) restricted to I_ε , i.e., *near the singularity* at $x = 0$. In order to improve the accuracy of the method on the subinterval $I - I_\varepsilon$ we form a partition

$$I - I_\varepsilon = \bigcup_{m=1}^M I_m \text{ where } I_m = [x_{m-1}, x_m], x_0 = \varepsilon \text{ and } x_M = L \quad (6)$$

The method by Ramos proceeds by using the same linearisation method on the subintervals $I_m, (m = 1, \dots, M)$ of $I - I_\varepsilon$, i.e. *away from the singularity*. To this end, at the interface x_{m-1} of I_{m-1} and I_m we make use the solution of Eq. (3) restricted to I_{m-1} to generate the initial conditions for Eq. (3) restricted to I_m . This ensures continuity of the solution. In this chapter we avoid linearisation on $I - I_\varepsilon$ by making use of the SRM on the subintervals of $I - I_\varepsilon$. However, as was done in [23] we make use of linearisation on I_ε . This approach results in the MSRM. A detailed description of the MSRM is given in Sections 2.1 and 2.2.

2.1 Near the singularity

Let $\varepsilon \in (0, L)$ be a sufficiently small number. Restrict problem (5) to $[0, \varepsilon]$ and re-arrange to get

$$y'' = - \underbrace{\left[\frac{\gamma}{x} y' + f(x)g(y) \right]}_{u(x,y,y')}, 0 < x \leq \varepsilon, y(0) = \alpha, y'(0) = 0 \quad (7)$$

If we Taylor expand u about $\varepsilon_0 = \left(\varepsilon, \underbrace{y(0)}_{y_0}, \underbrace{y'(0)}_{y'_0} \right)$ and neglect higher order

terms we get

$$u(x,y,y') = \underbrace{u(\varepsilon_0)}_{u_0} + \underbrace{u_y(\varepsilon_0)}_{H_0}(y - y_0) + \underbrace{u_{y'}(\varepsilon_0)}_{G_0}(y' - y'_0) + \underbrace{u_x(\varepsilon_0)}_{L_0}(x - \varepsilon)$$

where u_x denotes $\frac{\partial u}{\partial x}$. Consequently, Eq. (7) can be replaced by

$$y'' = u_0 + H_0(y - y_0) + G_0(y' - y'_0) + L_0(x - \varepsilon), 0 \leq x \leq \varepsilon, y(0) = \alpha, y'(0) = 0, \quad (8)$$

where the differential equation now holds at $x = 0$ because the singularity there is no longer present. If $H_0 \neq 0$ and $R_0 := (G_0/2)^2 + H_0 > 0$, then problem (8) has analytical solution [23]

$$y(x) = A_0 \exp(\lambda_0^+(x - \varepsilon)) + B_0 \exp(\lambda_0^-(x - \varepsilon)) + C_0(x - \varepsilon) + D_0 \quad (9)$$

for $0 \leq x \leq \varepsilon$, where $\lambda_0^\pm = G_0/2 \pm \sqrt{R_0}$, $C_0 = -L_0/H_0$,

$$D_0 = -(G_0 C_0 + P_0)/H_0, P_0 = u_0 - H_0 y_0 - G_0 y'_0,$$

$$A_0 = \frac{\exp(\lambda_0^+ \varepsilon)}{\lambda_0^+ - \lambda_0^-} (y'_0 - C_0 - \lambda_0^- (y_0 + C_0 \varepsilon - D_0)),$$

$$B_0 = \frac{\exp(\lambda_0^- \varepsilon)}{\lambda_0^+ - \lambda_0^-} (\lambda_0^+ (y_0 + C_0 \varepsilon - D_0) - (y'_0 - C_0))$$

Thus

$$y(\varepsilon) = \underbrace{A_0 + B_0 + D_0}_{\alpha_0} \text{ and } y'(\varepsilon) = \underbrace{\lambda_0^+ A_0 + \lambda_0^- B_0 + C_0}_{\alpha'_0} \quad (10)$$

We take α_0 and α'_0 as initial values for problem (7) restricted to $[\varepsilon, L]$ in the next section.

2.2 Away from the singularity

We seek $y(x)$ satisfying

$$y'' + \frac{\gamma}{x} y' + f(x)g(y) = 0, \varepsilon \leq x \leq L, y(\varepsilon) = \alpha_0, y'(\varepsilon) = \alpha'_0 \quad (11)$$

In this section we begin by describing the SRM for problem (11). In practical applications it is usually important to obtain a solution to (11) which possesses a prescribed degree of accuracy. To this end we make use of the SRM on a partition of the problem domain $[\varepsilon, L]$. This is our last task in this section.

The first step of the SRM for (11) is to let $v = y$ and $w = v'$ so that we obtain the equivalent problem

$$v' = w, v(\varepsilon) = \alpha_0, \quad (12)$$

$$w' + \frac{\gamma}{x}w + f(x)g(v) = 0, w(\varepsilon) = \alpha'_0 \quad (13)$$

for $\varepsilon \leq x \leq L$ which upon making use of the change of variable

$$x(\eta) = \frac{L + \varepsilon}{2} + \frac{L - \varepsilon}{2}\eta$$

becomes

$$\beta \frac{dv}{d\eta} = w, v(-1) = \alpha_0 \quad (14)$$

$$\beta \frac{dw}{d\eta} + \frac{\gamma}{x(\eta)}w + f(x(\eta))g(v) = 0, w(-1) = \alpha'_0 \quad (15)$$

for $-1 \leq \eta \leq 1$ where $\beta = 2/(L - \varepsilon)$. As described in [19] the next step of the SRM mimicks the Gauss-Seidel method for linear systems and it yields the iteration

$$\beta \frac{dv_{r+1}}{d\eta} = \underbrace{w_r}_{R^{(1)}(\eta)}, v_{r+1}(\eta_N) = \alpha_0 \quad (16)$$

$$\beta \frac{dw_{r+1}}{d\eta} + \frac{\gamma}{x}w_{r+1} = \underbrace{-f(x(\eta))g(v_{r+1})}_{R^{(2)}(\eta)}, w_{r+1}(\eta_N) = \alpha'_0 \quad (17)$$

for $-1 = \eta_N \leq \eta \leq \eta_0 = 1$ where $r = 0, 1, \dots$ and on $[-1, 1]$ we formed a grid consisting of the Chebyshev-Gauss-Lobatto collocation points

$$\eta_i = \cos\left(\frac{\pi i}{N}\right), \quad i = 0, 1, \dots, N. \quad (18)$$

If the initial approximations v_0 and w_0 to v and w , respectively, are prescribed then Eqs. (16) and (17) generate sequences $\{v_r\}_{r=1}^{\infty}$ and $\{w_r\}_{r=1}^{\infty}$ of consecutive approximations. To this end we assume that v_r and w_r are known at the end of the r th iteration. Once Eq. (16) is solved for v_{r+1} the right hand side $R^{(2)}$ of Eq. (17) becomes known and we solve this equation for w_{r+1} . Since v_{r+1} and w_{r+1} are now known, we proceed in a similar manner to compute v_{r+2} and w_{r+2} , and so on. As done in [19] we solve Eqs. (16) and (17) by making use of Chebyshev differentiation [24] to obtain

$$\underbrace{\hat{D}}_{A_1} \mathbf{v}_{r+1} = \mathbf{R}^{(1)}, v_{r+1}(\eta_N) = \alpha_0, \quad (19)$$

$$\underbrace{(\hat{D} + \gamma \text{diag}[1/x(\eta_0), \dots, 1/x(\eta_N)])}_{A_2} \mathbf{w}_{r+1} = \mathbf{R}^{(2)}, w_{r+1}(\eta_N) = \alpha'_0 \quad (20)$$

where $\hat{D} = \beta D$, D is the $N \times N$ Chebyshev differentiation matrix,

$$\text{diag}[b_0, \dots, b_N] = \begin{pmatrix} b_0 & & \\ & \ddots & \\ & & b_N \end{pmatrix} \quad (21)$$

is a diagonal matrix, and $\mathbf{R}^{(i)} = [R^{(i)}(\eta_0), \dots, R^{(i)}(\eta_N)]^T$ with $i = 1, 2$. Moreover,

$$\mathbf{v}_r = [v_r(\eta_0), \dots, v_r(\eta_N)]^T, \mathbf{w}_r = [w_r(\eta_0), \dots, w_r(\eta_N)]^T, \quad r = 0, 1, \dots \quad (22)$$

and $[\]^T$ denotes matrix transpose.

We prescribe v_0 and w_0 by requiring that

$$v_0(\varepsilon) \equiv y_0(\varepsilon) = \alpha_0 \text{ and } w_0(\varepsilon) \equiv y'_0(\varepsilon) = \alpha'_0$$

Moreover, we assume that

$$\lim_{r \rightarrow \infty} v_r = v \text{ and } \lim_{r \rightarrow \infty} w_r = w$$

The initial conditions

$$v_{r+1}(\eta_N) = \alpha_0 \text{ and } w_{r+1}(\eta_N) = \alpha'_0$$

are included in the iterative scheme consisting of Eqs. (19) and (20) as shown below.

$$\begin{pmatrix} A_1 \\ 0 \dots 0 \ 1 \end{pmatrix} \begin{pmatrix} v_{r+1}(\eta_0) \\ \vdots \\ v_{r+1}(\eta_{N-1}) \\ v_{r+1}(\eta_N) \end{pmatrix} = \begin{pmatrix} R^{(1)}(\eta_0) \\ \vdots \\ R^{(1)}(\eta_{N-1}) \\ \alpha_0 \end{pmatrix} \quad (23)$$

$$\begin{pmatrix} A_2 \\ 0 \dots 0 \ 1 \end{pmatrix} \begin{pmatrix} w_{r+1}(\eta_0) \\ \vdots \\ w_{r+1}(\eta_{N-1}) \\ w_{r+1}(\eta_N) \end{pmatrix} = \begin{pmatrix} R^{(2)}(\eta_0) \\ \vdots \\ R^{(2)}(\eta_{N-1}) \\ \alpha'_0 \end{pmatrix} \quad (24)$$

The larger L is the less reliable is the SRM. As a workaround to this problem is we subdivide interval $[\varepsilon, L]$ into a disjoint union of non-overlapping subintervals as detailed in Eq. (6). Given the the model problem on I_1 , we use the SRM to compute estimate \tilde{y} to y on I_1 . We make use of \tilde{y} to compute initial values for the problem on I_2 . We repeat this procedure for the problem on I_2 and continue in a similar manner until we exhaust $[\varepsilon, L]$. Shown in Algorithm 2.2 is an outline of the MSRM for problem (5).

Algorithm 1: Putting the MSRM together.

1. Let $[0, L] = I_\varepsilon \cup [\varepsilon, L]$ where $I_\varepsilon := [0, \varepsilon]$ for some sufficiently small number $\varepsilon > 0$.
2. Replace nonlinear Eq. (5) with linear Eq. (8) on I_ε and compute solution $y_\varepsilon(x)$ to Eq. (7) on I_ε .
3. Set $\alpha_0 = y_\varepsilon(\varepsilon)$, set $\alpha'_0 = y'_\varepsilon(\varepsilon)$ and let $\varepsilon = x_0 < x_1 < \dots < x_M = L$.

4. **for** $m = 1, 2, \dots, M$ **do**

5. Define $I_m := [x_{m-1}, x_m]$ and make use of the SRM to solve

$$y'' = u(x, y, y'), x \in I_m, y(x_{m-1}) = \alpha_0, y'(x_{m-1}) = \alpha'_0$$

for the estimate \tilde{y} to y on I_m .

6. Set $\alpha_0 = \tilde{y}(x_m)$ and set $\alpha'_0 = \tilde{y}'(x_m)$.

7. **end**

2.3 Examples

In this section we make use of some examples to investigate the accuracy and computational efficiency of the MSRM.

Example 1 We look for a numerical solution to the problem

$$y'' + \frac{1}{x}y' + 3y^5 - y^3 = 0, 0 < x \leq 10, y(0) = 1, y'(0) = 0 \quad (25)$$

which Wazwaz [11] solved using the VIM and obtained the approximate analytical solution

$$y(x) = \frac{1}{\sqrt{1+x^2}} \quad (26)$$

When we apply the MSRM on (25) we get the following components for constructing the numerical solution.

1. Coefficients

$$u_0 = -2, L_0 = 0, H_0 = -12, G_0 = -10^4 \quad (27)$$

for problem (8).

2. Initial values

$$y(\varepsilon) := \alpha_0 = 9.999999926 \times 10^{-1} \text{ and } y'(\varepsilon) := \alpha'_0 = -1.264241093 \times 10^{-4} \quad (28)$$

for problem (11).

3. Entries

$$R_i^{(1)} = w(\eta_i) \text{ and } R_i^{(2)} = -(3v_{r+1}^5(\eta_i) - v_{r+1}^3(\eta_i)), i = 0, \dots, N$$

in the vectors on the right hand sides of Eqs. (19) and (20).

The MSRM generates a numerical solution to problem (25) which is plotted together with the analytical solution (26) in **Figure 1(a)**. **Figure 1(a)** shows a good agreement between the numerical and analytical solutions. For a more detailed comparison of the two solutions see **Table 1**. **Table 1** and **Figure 1(b)** show that the absolute error of the numerical solution is no more than 0.5×10^{-6} . Thus the numerical solution agrees with the analytical solution in the first 6 decimal places.

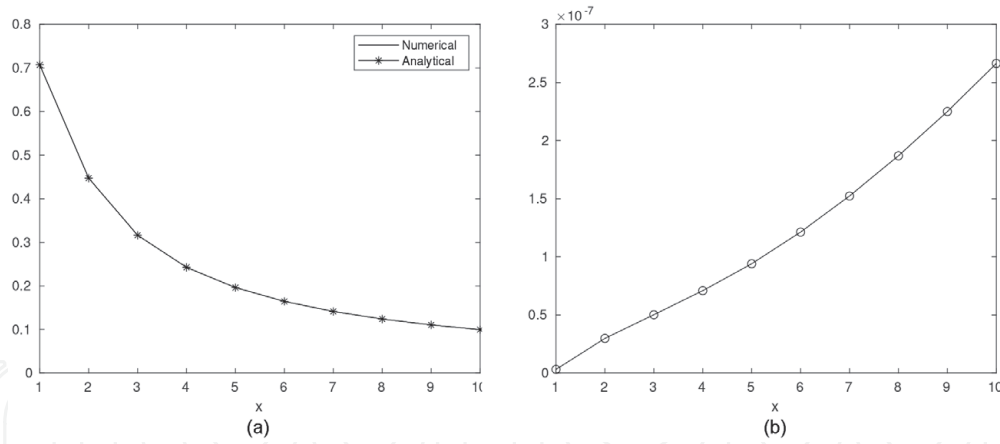


Figure 1. (a) Solutions to problem (25). (b) Absolute error of the MSRM solution to (25).

x	MSRM	Solution (26)	Absolute error
1	0.70710678	0.70710678	3.1×10^{-9}
2	0.44721363	0.44721360	3×10^{-8}
3	0.31622782	0.31622777	5×10^{-8}
4	0.24253570	0.24253563	7.1×10^{-8}
5	0.19611623	0.19611614	9.4×10^{-8}
6	0.16439911	0.16439899	1.2×10^{-7}
7	0.14142151	0.14142136	1.5×10^{-7}
8	0.12403492	0.12403473	1.9×10^{-7}
9	0.11043175	0.11043153	2.3×10^{-7}
10	0.09950399	0.09950372	2.7×10^{-7}

Table 1. Comparison of numerical values of y with solution (26).

This degree of accuracy was achieved by the MSRM upon choosing $N = 60, M = 10$, setting the maximum number of iterations as $r_{max} = 20$ and imposing the stopping criterion

$$\max \{ \|\mathbf{v}_{r+1} - \mathbf{v}_r\|_\infty, \|\mathbf{w}_{r+1} - \mathbf{w}_r\|_\infty \} \leq \varepsilon, r = 0, 1, \dots$$

for the iterative method, where $\varepsilon = 10^{-6}$ is the tolerance and $\|\mathbf{w}\|_\infty = \max_{0 \leq i \leq N} |w_i|$ is the infinity norm. It took only $r = 5$ iterations for the MSRM to achieve the given degree of accuracy.

Example 2 We consider the initial value problem

$$y'' + \frac{5}{x}y' + 8(e^y + 2e^{y/2}) = 0, 0 < x \leq 10, y(0) = y'(0) = 0 \quad (29)$$

that was solved by Wazwaz [10] using the ADM to obtain the approximate analytical solution

$$y(x) = -2 \ln(1 + x^2) \quad (30)$$

Applying the MSRM on (29) produces the following building blocks for constructing the numerical solution.

1. Coefficients

$$u_0 = -24, L_0 = 0, H_0 = -16, G_0 = -5 \times 10^4 \quad (31)$$

for problem (8).

2. Initial values

$$y(\varepsilon) := \alpha_0 = -3.846468396 \times 10^{-8} \text{ and } y'(\varepsilon) := \alpha'_0 = -4.767657761 \times 10^{-4} \quad (32)$$

for problem (11).

3. Elements

$$R_i^{(1)} = w(\eta_i) \text{ and } R_i^{(2)} = -8 \left(e^{v_{r+1}(\eta_i)} + 2e^{v_{r+1}(\eta_i)/2} \right), i = 0, \dots, N$$

of the vectors on the right hand sides of Eqs. (19) and (20).

A comparison of the MSRM solution to problem (29) with the analytical solution (30) is shown in **Figure 2(a)**. The graph suggests that the two solutions are exactly the same. However, a closer look at the two solutions is provided in **Table 2** and we observe that the analytical and numerical solutions are slightly different. A plot of the absolute error of the MSRM solution over a grid on the problem domain $[0, 10]$ is shown in **Figure 2(b)**. We observe that the absolute error does not exceed 0.5×10^{-7} . Hence the MSRM solution and the analytical solution agree to within 7 decimal places. For the MSRM to achieve this degree of accuracy we chose $N = 60, M = 10, r_{max} = 20$ and $\varepsilon = 10^{-6}$. We observed that the MSRM stopped at iteration $r = 5$ for problem (29).

Example 3 We seek a numerical solution to

$$y'' + \frac{2}{x}y' - 6y - 4y \ln y = 0, 0 < x \leq 1, y(0) = 1, y'(0) = 0, \quad (33)$$

where

$$y(x) = e^{x^2} \quad (34)$$

is the exact solution [25]. We restrict the problem to a relatively small interval $[0, 1]$ because the solution (34) grows rapidly.

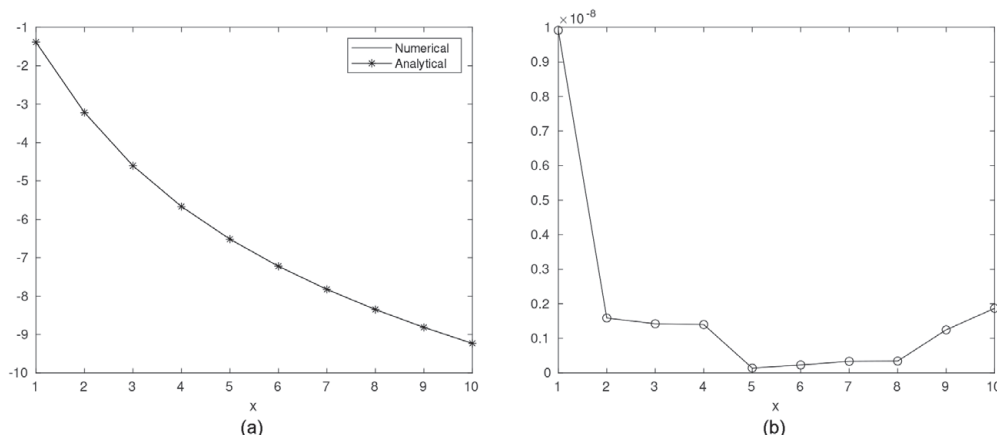


Figure 2.
 (a) Solutions to problem (29). (b) Absolute error of the MSRM solution to (29).

x	MSRM	Solution (34)	Absolute error
1	-1.38629437	-1.38629436	9.9×10^{-9}
2	-3.21887583	-3.21887582	1.6×10^{-9}
3	-4.60517018	-4.60517019	1.4×10^{-9}
4	-5.66642669	-5.66642669	1.4×10^{-9}
5	-6.51619308	-6.51619308	1.4×10^{-10}
6	-7.22183583	-7.22183583	2.3×10^{-10}
7	-7.82404601	-7.82404601	3.4×10^{-10}
8	-8.34877454	-8.34877454	3.4×10^{-10}
9	-8.81343849	-8.81343849	1.2×10^{-9}
10	-9.23024103	-9.23024103	1.9×10^{-9}

Table 2.
Comparison of numerical values of y with solution (30).

The MSRM for (34) gives the following components for constructing the numerical solution.

1. Coefficients

$$u_0 = 6, L_0 = 0, H_0 = 10 \text{ and } G_0 = -2 \times 10^4 \quad (35)$$

for problem (8).

2. Initial values

$$\alpha_0 = 1.000000017 \text{ and } \alpha'_0 = 2.593994191 \times 10^{-04} \quad (36)$$

for problem (11).

3. Right hand sides

$$R_i^{(1)} = w_r(\eta_i) \text{ and } R_i^{(2)} = 6v_{r+1}(\eta_i) + 4v_{r+1}(\eta_i) \ln(v_{r+1}(\eta_i)), i = 0, \dots, N \quad (37)$$

for linear systems (19) and (20).

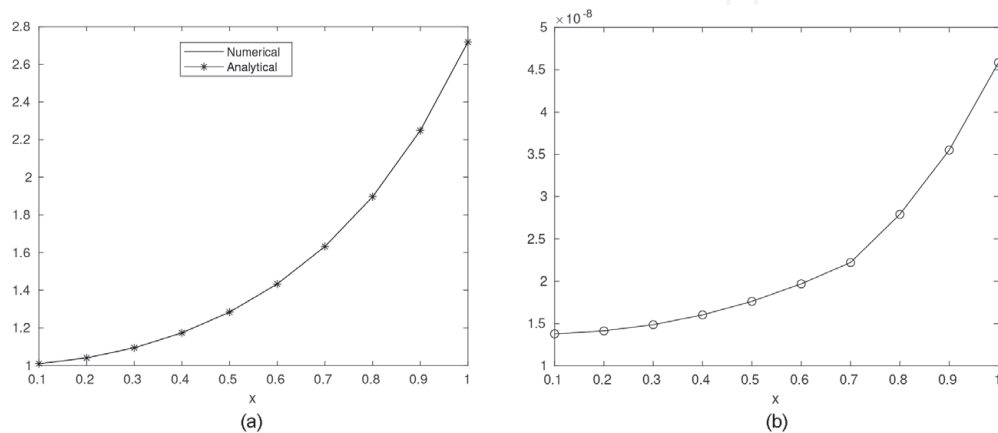


Figure 3.
(a) Solutions to problem (33). (b) Absolute error of MSRM solution to (33).

x	MSRM	Solution (34)	Absolute error
0.1	1.01005018	1.01005017	1.4×10^{-8}
0.2	1.04081079	1.04081077	1.4×10^{-8}
0.3	1.09417430	1.09417428	1.5×10^{-8}
0.4	1.17351089	1.17351087	1.6×10^{-8}
0.5	1.28402543	1.28402542	1.8×10^{-8}
0.6	1.43332943	1.43332941	2×10^{-8}
0.7	1.63231624	1.63231622	2.2×10^{-8}
0.8	1.89648091	1.89648088	2.8×10^{-8}
0.9	2.24790802	2.24790799	3.6×10^{-8}
1	2.71828187	2.71828183	4.6×10^{-8}

Table 3.
 Comparison of numerical values of y with solution (34).

Figure 3(a) shows that the numerical solution agrees well with the analytical solution (34). For a closer look at how the numerical and analytical solutions compare see **Table 3**. We observe that the MSRM solution agrees with the analytical solution (34) on the problem domain $[0, 1]$ to within at least 7 decimal places. A plot of the absolute error at these grid points on $[0, 1]$ is shown in **Figure 3(b)**. We observe that the absolute error in the MSRM increases as we move away from the singular point $x = 0$, but the absolute error never exceeds 0.5×10^{-7} . This degree of accuracy is achieved with $N = 60, M = 10, r_{max} = 20$ and $\epsilon = 10^{-6}$. Moreover, only 6 iterations were required to achieve this degree of accuracy.

3. Conclusions

In this chapter we presented a modified spectral relaxation method (denoted by MSRM) for solving singular initial value problems for some Emden-Fowler equations. We made use of some examples of the model problem to demonstrate that the MSRM is reliable, accurate and computationally efficient. The method provided a reliable treatment of the singular point. The MSRM solutions were compared with analytical solutions that were obtained using other methods, i.e., the Variational iteration method and the Adomian decomposition method. There was agreement between the solutions that were compared in the first 6 decimal places. A possible way of increasing the degree of accuracy of the MSRM would be to increase the tolerance for the method. This and other ways for optimizing the method could constitute future work. In all the examples that were considered, it took at most 6 iterations for the MSRM to converge. Hence the method exhibited rapid convergence.

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