

An Efficient High-Order Numerical Method for Second-Order Steklov Problem in a Sectorial Domain

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Abstract— In this paper, we present a spectral Galerkin method for solving the second-order Steklov eigenvalue problem in a sectorial domain. First, by using affine coordinate transformations, we reformulate the original problem into equivalent system on a rectangular domain. Then, we establish its variational and discrete forms. Finally, numerical experiments are presented to demonstrate efficiency and spectral accuracy of the proposed algorithm.

Keywords— Steklov eigenvalue problem, Legendre-Galerkin method, sectorial domain.

I. INTRODUCTION

Steklov eigenvalue problems have important physical background and possesses broad applications across multiple fields [1-3]. Extensive researches have been conducted in both theoretical analysis and numerical computation. Tan et al investigated positivity preserving property across various domains [4-6], and the properties of the first eigenvalue have been explored in the literatures [7-10]. On the numerical side, numerous methods have been developed [11-14], existing works mainly focuses on solving the second-order Steklov eigenvalue problems on the rule regions. Notably, investigations of such problems in sector-shaped domains remain absent from the existing literatures.

Therefore, this paper presents a numerical method for solving second-order Steklov eigenvalue problems in sectorshaped domains. First, by employing affine transformation, we reformulate the original problem into equivalent system. Then through the construction of some weighted Sobolev spaces, we derive the weak formulation and its corresponding discrete scheme. Finally, extensive computational experiments demonstrate the algorithm's astringency and efficiency.

The rest of this paper is organized as follows: In Section 2, a weak form and its discrete scheme of the second-order Steklov eigenvalue problem are derived. In Section 3, we derive the corresponding matrix format for the discrete scheme. In Section 4, we present a series of numerical examples to validate the proposed method. Finally, we conclude the paper in Section 5.

EQUIVALENT COUPLED SYSTEM II.

We consider the following second-order Steklov problem:

$$\begin{cases} -\Delta w + w = 0, (x, y) \in \Omega, \\ \frac{\partial w}{\partial n} = \lambda w, (x, y) \in \partial \Omega, \end{cases}$$
(2.1)

where $\frac{\partial w}{\partial n}$ is the outward normal derivative on $\partial \Omega$, and

 $\Omega = \{(x, y) \in \mathbb{R}^2 : x = r \cos \theta, y = r \sin \theta, 0 \le r \le R, \theta_1$ $\leq \theta \leq \theta_2$ }. It is clear that the weak form of (2.1) is as follows: Find

 $(\lambda, w) \in \mathbb{R} \times H^1(\Omega)$, such that

$$A(w,v) = \lambda B(w,v), v \in H^{1}(\Omega), \qquad (2.2)$$

where

$$A(w,v) = \int_{\Omega} \nabla w \nabla v + wv dx dy, B(w,v) = \int_{\partial \Omega} wv ds.$$

Define a Laplace operator:
$$\hat{\mathbf{L}} = \frac{1}{r} \left[\partial_r (r \partial_r) + \frac{1}{r} \partial_{\theta}^2 \right].$$
 (2.3)

Let $\hat{\psi}(r,\theta) = w(x,y)$. Using the chain rule for derivatives of composite functions, we can derive the equivalent form of the gradient operator as follows:





As shown in the Figure 1, we denote the unit outward normal vector of the arc boundary AB as $n_1 = (\cos \theta, \sin \theta)$, the unit outward normal vector of the initial boundary OA as $n_2 = (\sin \theta_1, -\cos \theta_1)$, and the unit outward normal vector of the final boundary *OB* as $n_3 = (-\sin\theta_2, \cos\theta_2)$. From $\frac{\partial w}{\partial n} = \nabla w \cdot n$, we deduce that (2.1) has the following equivalent form:

$$\begin{cases} -\hat{L}\,\hat{\psi}+\hat{\psi}=0, & (r,\theta)\in D, \\ \partial_{\theta}\hat{\psi}(0,\theta)=0, \partial_{r}\hat{\psi}(R,\theta)=\lambda\hat{\psi}(R,\theta), \theta\in(\theta_{1},\theta_{2}), \\ -\frac{1}{r}\partial_{\theta}\hat{\psi}(r,\theta_{1})=\lambda\hat{\psi}(r,\theta_{1}), & r\in(0,R), \end{cases}$$
(2.4)
$$\frac{1}{r}\partial_{\theta}\hat{\psi}(r,\theta_{2})=\lambda\hat{\psi}(r,\theta_{2}), & r\in(0,R). \end{cases}$$

Note that $\partial_{\theta}\hat{\psi}(0,\theta) = 0$ is an essential polar condition and $D = (0, R) \times (\theta_1, \theta_2)$. Let



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$$r = \frac{R(1+t)}{2}, \ \theta = \frac{a+b\theta}{2}, \ (t,\theta) \in D = (-1,1)^2,$$
$$L = \frac{1}{1+t}\partial_t [(1+t)\partial_t] + \frac{4}{b^2}\frac{1}{(1+t)^2}\partial_{\theta}^2, \ a = \theta_1 + \theta_2, \ b = \theta_2 - \theta_1.$$

Denoting $\psi(t, \theta) = \hat{\psi}(r, \theta)$, we further reformulate (2.4) as:

$$\begin{cases} -L \psi + \frac{R^2}{4} \psi = 0, \qquad (t, \vartheta) \in D, \\ \partial_{\vartheta} \psi(-1, \vartheta) = 0, \partial_{\imath} \psi(1, \vartheta) = \lambda \frac{R}{2} \psi(R, \vartheta), \vartheta \in (-1, 1), \\ -\frac{1}{1+t} \partial_{\vartheta} \psi(t, -1) = \lambda \frac{bR}{4} \psi(t, -1), \qquad t \in (-1, 1), \\ \frac{1}{1+t} \partial_{\vartheta} \psi(t, 1) = \lambda \frac{bR}{4} \psi(t, 1), \qquad t \in (-1, 1). \end{cases}$$

$$(2.5)$$

We define the following non-uniformly weighted Sobolev spaces:

$$L^{2}_{\omega}(D) := \left\{ \psi : \int_{D} \omega |\psi|^{2} dt d\theta < \infty \right\}, \omega = \frac{b}{2}(1+t),$$

$$H^{1}_{*}(D) := \left\{ \psi : \int_{D} \omega \partial_{t} \psi \partial_{t} h + \omega^{-1} \partial_{\vartheta} \psi \partial_{\vartheta} h + \omega \psi h dt d\vartheta < \infty, \partial_{\vartheta} \psi (-1, \vartheta) = 0 \right\}$$

the corresponding inner product and norm are given by

$$\begin{aligned} (\psi,h)_{*,1} &= \int_{D} \omega \partial_{t} \psi \partial_{t} h + \omega^{-1} \partial_{g} \psi \partial_{g} h + \omega \psi h dt d\mathcal{G}, \\ \|\psi\|_{*,1} &= \sqrt{(\psi,\psi)_{*,1}}. \end{aligned}$$

According to (2.2), a weak form of (2.5) is: Find $(\lambda, \psi) \in \mathbb{R} \times H^1_*(D)$, such that

$$\mathbf{A}(\psi,h) = \lambda \mathbf{B}(\psi,h), h \in H^1_*(D), \qquad (2.6)$$

where

$$\begin{split} \mathbf{A}(\psi,h) &= \int_{D} \omega \partial_{t} \psi \partial_{t} h + \omega^{-1} \partial_{\vartheta} \psi \partial_{\vartheta} h dt d\vartheta + \frac{bR^{2}}{8} \int_{D} (1+t) \psi h dt d\vartheta, \\ \mathbf{B}(\psi,h) &= \frac{bR}{2} \int_{-1}^{1} \psi(1,\vartheta) h(1,\vartheta) d\vartheta \\ &+ \frac{R}{2} \int_{-1}^{1} \psi(t,-1) h(t,-1) + \psi(t,1) h(t,1) dt. \end{split}$$

Define an approximation space $X_N = (P_N \times P_N) \cap H^1_*(D)$. Then, the discrete scheme corresponding to (2.6) is: Find $(\lambda_N, \psi_N) \in \mathbb{R} \times X_N$, such that

$$\mathbf{A}(\boldsymbol{\psi}_{N},\boldsymbol{h}_{N}) = \lambda_{N} \mathbf{B}(\boldsymbol{\psi}_{N},\boldsymbol{h}_{N}), \,\forall \boldsymbol{h}_{N} \in \boldsymbol{X}_{N}.$$
(2.7)

III. ALGORITHMIC IMPLEMENTATION

In this section, we shall derive the equivalent matrix form of the discrete format (2.7). Let

$$\begin{split} \phi_{i}(t) &= L_{i}(t) - L_{i+2}(t), 0 \le i \le N - 2, \phi_{N-1}(t) = \frac{t+1}{2}, \phi_{N}(t) = 1, \\ \hat{\phi}_{i}(\mathcal{G}) &= L_{i}(\mathcal{G}) - L_{i+2}(\mathcal{G}), 0 \le i \le N - 2, \hat{\phi}_{N-1}(\mathcal{G}) = L_{0}(\mathcal{G}), \hat{\phi}_{N}(\mathcal{G}) = L_{1}(\mathcal{G}), \\ S_{N} &= \operatorname{span}\{\phi_{0}(t), \phi_{1}(t), \cdots, \phi_{N}(t)\}, \hat{S}_{N} = \operatorname{span}\{\hat{\phi}_{0}(\mathcal{G}), \hat{\phi}_{1}(\mathcal{G}), \cdots, \hat{\phi}_{N}(\mathcal{G})\}. \\ \text{Then the approximate space is:} \end{split}$$

$$X_{N} = S_{N} \times \hat{S}_{N}.$$

Let us expand ψ_N as follows:

$$\psi_N = \sum_{i,j=0}^N u_{ij} \phi_i(t) \hat{\phi}_j(\mathcal{G}).$$
(3.1)

We denote by

$$\mathbf{U} = \begin{pmatrix} u_{00} & u_{01} & \cdots & u_{0N} \\ u_{10} & u_{11} & \cdots & u_{1N} \\ \vdots & \vdots & \ddots & \vdots \\ u_{N0} & u_{N1} & \cdots & u_{NN} \end{pmatrix}.$$

Let \overline{U} be a column vector of a length $(N+1)^2$ obtained by expanding U. Denote

$$\begin{split} e_{kj} &= \int_{-1}^{1} \hat{\phi}_{j}(\mathcal{G}) \hat{\phi}_{k}(\mathcal{G}) d\mathcal{G}, g_{kj} = \int_{-1}^{1} \hat{\phi}_{j}'(\mathcal{G}) \hat{\phi}_{k}'(\mathcal{G}) d\mathcal{G}, m_{kj} = \hat{\phi}_{j}(-1) \hat{\phi}_{k}(-1), \\ p_{kj} &= \hat{\phi}_{j}(1) \hat{\phi}_{k}(1), \ q_{kj} = \phi_{j}(1) \phi_{k}(1), \\ \hat{e}_{kj} &= \int_{-1}^{1} \phi_{j}(t) \phi_{k}(t) dt, \ \hat{g}_{kj} = \int_{-1}^{1} \frac{1}{1+t} \phi_{j}(t) \phi_{k}(t) dt, \\ \hat{m}_{kj} &= \int_{-1}^{1} (1+t) \phi_{j}(t) \phi_{k}(t) dt, \ \hat{p}_{kj} = \int_{-1}^{1} (1+t) \phi_{j}'(t) \phi_{k}'(t) dt. \end{split}$$

By substituting (3.1) into (2.7), and then allowing h_N across all basis functions in X_N , we obtain the following linear eigenvalue system:

$$W_1\overline{U}=\lambda_NW_2\overline{U}$$
 , ,

where

$$\begin{split} W_{1} &= \frac{b}{2} E(k,:) \otimes \hat{P}(l,:) + \frac{2}{b} G(k,:) \otimes \hat{G}(l,:) + \frac{bR^{2}}{8} E(k,:) \otimes \hat{M}(l,:), \\ W_{2} &= \frac{bR}{2} E(k,:) \otimes Q(l,:) + \frac{R}{2} (M+P)(k,:) \otimes \hat{E}(l,:), \\ E &= (e_{ij})_{i,j=0}^{N}, G = (g_{ij})_{i,j=0}^{N}, M = (m_{ij})_{i,j=0}^{N}, P = (p_{ij})_{i,j=0}^{N}, Q = (q_{ij})_{i,j=0}^{N}, \\ \hat{E} &= (\hat{e}_{ij})_{i,j=0}^{N}, \hat{G} = (\hat{g}_{ij})_{i,j=0}^{N}, \hat{M} = (\hat{m}_{ij})_{i,j=0}^{N}, \hat{P} = (\hat{p}_{ij})_{i,j=0}^{N}. \end{split}$$

IV. NUMERICAL EXPERIMENT

In this section, a series of numerical experiments will be presented to demonstrate the efficiency and spectral accuracy of our algorithm. Our program is compiled and executed in MATLAB R2016b.

Example 1: We take R = 1, $\theta_1 = 30^\circ$, $\theta_2 = 90^\circ$. In Table 1, we list the first four approximate eigenvalues λ_N^i (i = 1,2,3,4) for varying values of N. To further validate the computational efficiency of our algorithm, we used the numerical solution obtained at N = 60 as a reference value. In Figure 2, we plotted the error curves of the eigenvalues for different values of N. Additionally, in Figure 3, we presented the image of the eigenfunction corresponding to the minimum eigenvalue at N = 60, along with the error image between the eigenfunctions corresponding to the minimum eigenvalues at N = 50 and N = 60.

TABLE 1. The n	umerical results of the f	irst four eigenvalues for	different N.
			-

N	λ_N^1	λ_N^2	λ_N^3	λ_N^4
10	0.167847123435502	1.358638641239197	1.792266897280079	4.040399080553113
20	0.167847123322805	1.358638633801590	1.792266818716613	4.040398943579751





Figure 2: Error curves (left) between the numerical eigenvalues and the reference eigenvalues and their error curves (right) under semi-log scale.



Figure 3: Image (left) of the eigenfunction $w_{60}(x, y)$ and the error image (right) between eigenfunctions $w_{60}(x, y)$ and $w_{50}(x, y)$ corresponding to the minimum eigenvalue.

As observed in Table 1, the computed eigenvalues achieve a precision of at least 10-digit when $N \ge 30$. Additionally, as illustrated in Figures 2-3, our algorithm demonstrates both convergence and spectral accuracy. *Example 2.* We take $R = 1, \theta_1 = 60^\circ, \theta_2 = 180^\circ$. In Table 1, we list the first four approximate eigenvalues λ_N^i (i = 1, 2, 3, 4) for varying values of N. We used the numerical solution obtained at N = 60 as a reference value to further validate the computational efficiency of our algorithm. In Figure 4, we plotted the error curves of the eigenvalues for different values of N. Additionally, in Figure 5, we presented the image of the eigenfunction corresponding to the minimum eigenvalue at N = 60, along with the error image between the eigenfunctions corresponding to the minimum eigenvalues at N = 50 and N = 60.

As observed in Table 2, the computed eigenvalues achieve a precision of at least 10-digit when $N \ge 40$. Additionally, as illustrated in Figures 4-5, our algorithm demonstrates both convergence and spectral accuracy.

TABLE 2. The numerical results of the first four eigenvalues for different N.

N	λ_N^1	λ_N^2	λ_N^3	λ_N^4
10	0.245318192635769	1.054242994892192	2.094981780998436	2.497636674150089
20	0.245318191456520	1.054242728944644	2.094981574949448	2.497636503216148
30	0.245318191449657	1.054242726021611	2.094981573880636	2.497636502299467
40	0.245318191449364	1.054242725798506	2.094981573835764	2.497636502261011
50	0.245318191449334	1.054242725762994	2.094981573831236	2.497636502257186





Figure 4: Error curves (left) between the numerical eigenvalues and the reference eigenvalues and their error curves (right) under semi-log scale.



Figure 5: Image (left) of the eigenfunction $W_{60}(x, y)$ and the error image (right) between eigenfunctions

 $W_{60}(x, y)$ and $W_{50}(x, y)$ corresponding to the minimum eigenvalue.

V. CONCLUSION

In this paper, an efficient Legendre spectral method is proposed and studied for second-order Steklov eigenvalue problems in sectorial domain. By utilizing affine technology, a fresh weak formulation and its corresponding discrete variational form are formulated. In addition, numerical results validate the effectiveness of the algorithm. The algorithm proposed in this paper can also be generalized to eigenvalue problems on more complex geometric domains, which is the goal of our future research.

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