Inverse iteration method for higher eigenvalues of the p-Laplacian

Timur Galimov

23.06

12:45-13:00

Institute of Mathematics with Computing Centre, Ufa Federal Research Centre, RAS

galim.timerxan@yandex.ru

Let $1 and let <math>\Omega \subset \mathbb{R}^D$ be a domain of finite measure, $D \ge 1$. Consider the problem of finding an eigenvalue $\lambda \in \mathbb{R}$ and an eigenfunction $u \in W_0^{1,p}(\Omega) \setminus \{0\}$ such that

$$\int_{\Omega} |\nabla u|^{p-2} \langle \nabla u, \nabla v \rangle \, dx = \lambda \int_{\Omega} |u|^{p-2} uv \, dx \quad \text{for any } v \in W_0^{1,p}(\Omega).$$

This is a weak form of the Dirichlet eigenvalue problem for the p-Laplace operator

$$\Delta_p(\cdot) := \operatorname{div}(|\nabla(\cdot)|^{p-2}\nabla(\cdot)),$$

turning into the classical Dirichlet Laplace eigenvalue problem for p=2. Eigenfunctions are precisely critical points of the Rayleigh quotient functional $u\mapsto R[u]$ defined on $W_0^{1,p}(\Omega)\setminus\{0\}$ as

$$R[u] := \frac{\int_{\Omega} |\nabla u|^p \, dx}{\int_{\Omega} |u|^p \, dx},$$

and eigenvalues are critical levels of R.

It is known (see, for example, [Lin]) that there exists the smallest eigenvalue $\lambda_1(\Omega,p)$, called the *first* eigenvalue, and that $\lambda_1(\Omega,p)$ is also the only eigenvalue admitting sign-constant eigenfunctions (all of which being constant multipliers of one-another). The properties of $\lambda_1(\Omega,p)$ allow to develope different numerical methods for its numerical approximation, among which algorithms based on inverse iteration schemes play an important role. All the other eigenvalues are called *higher*, and they are significantly less studied from both the theoretic and the numeric viewpoints.

We propose a novel algorithm using inverse iterations to approximate some higher eigenvalue alongside the corresponding eigenfunctions. Given an arbitrary initial guess $u_0 \in L^{\infty}(\Omega)$ such that

$$\min\left\{\|u_0^+\|_p,\|u_0^-\|_p\right\}>0\quad\text{and}\quad\left|\{x\in\Omega:\,u_0(x)=0\}\right|=0,$$

this algorithm generates a sequence $\{u_{k+1}\}\subset W^{1,p}_0(\Omega)\setminus\{0\}$ by solving consecutive p-Poisson equations with the property that their solutions u_{k+1} satisfy

 $R[u_{k+1}] = R[u_{k+1}^+] = R[u_{k+1}^-],$

where for any $u: \Omega \to \mathbb{R}$ we let $u^{\pm} := \max\{\pm u, 0\}$. We prove the following convergence results for $\{u_{k+1}\}$.

Theorem 1. Let \mathcal{U} be the set of all strong limit points of the sequence $\{u_k\}$ in $W_0^{1,p}(\Omega)$. Then the following assertions hold:

1. The sequence $\{R[u_k]\}$ monotonically decreases towards a higher eigenvalue R^* :

$$R^* := \inf_{k \geqslant 1} R[u_k] = \lim_{k \to \infty} R[u_k].$$

- 2. $\mathcal{U} \neq \emptyset$ and it is a subset of the (sign-changing) eigenfunctions corresponding to the eigenvalue R^* .
- 3. Any subsequence $\{u_{k_n}\}$ contains a sub-subsequence having a strong limit in $W_0^{1,p}(\Omega)$. If $\{u_{k_n}\}$ converges strongly to some $u \in \mathcal{U}$, then, for any $i \in \mathbb{N}$, the shifted subsequence $\{u_{k_n+i}\}$ also converges strongly to u.
- 4. $\lim_{k\to\infty} \rho(u_k,\mathcal{U})=0$, where ρ stands for the distance function in $W_0^{1,p}(\Omega)$.
- 5. U either consists of a single eigenfunction or has no isolated elements. In particular, if R^* is a simple eigenvalue, then the whole sequence $\{u_k\}$ converges to the corresponding eigenfunction.

The talk is based on the paper [BG25].

- [BG25] V. Bobkov and T. Galimov, Inverse iteration method for higher eigenvalues of the p-Laplacian, 2025, arXiv: 2504. 12836.
- [Lin] P. Lindqvist, *A nonlinear eigenvalue problem*, Topics in mathematical analysis, pp. 175–203.