

Isogeometric finite element analysis of time-harmonic exterior acoustic scattering problems

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Abstract

We present an isogeometric analysis of time-harmonic exterior acoustic problems. The infinite space is truncated by a fictitious boundary and (simple) absorbing boundary conditions are applied. The truncation error is included in the exact solution so that the reported error is an indicator of the performance of the isogeometric analysis, in particular of the related pollution error. Numerical results performed with high-order basis functions (third or fourth orders) showed no visible pollution error even for very high frequencies. This property combined with exact geometrical representation makes isogeometric analysis a very promising platform to solve high-frequency acoustic problems.

Keywords: Isogeometric Analysis, time-harmonic acoustics, Scattering, Helmholtz problem, high frequency, pollution error

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1. Introduction

The purpose of this paper is to investigate the performance of Non-Uniform Rational B-Splines (NURBS)-based *Isogeometric Analysis (IGA)* finite element methods (IGAFEM) for solving time-harmonic acoustic wave scattering and acoustic wave propagation problems, most particularly for small wavelengths. To fix the notations, let us consider Ω^- as a d -dimensional ($d = 1, 2, 3$) scattering bounded domain of \mathbb{R}^d , with shape $\Gamma := \partial\Omega^-$. We introduce the associated exterior (i.e. unbounded) domain of propagation $\Omega^+ := \mathbb{R}^d / \overline{\Omega^-}$. Then, solving the scattering problem leads to computing the wave field u , solution to the Boundary-Value Problem (BVP): given an incident (plane wave) field u^{inc} , find u such that

$$\begin{aligned} \Delta u + k^2 u &= 0, \quad \text{in } \Omega^+, \\ \partial_{\mathbf{n}_\Gamma} u &= \mathbf{g} := -\partial_{\mathbf{n}_\Gamma} u^{\text{inc}}, \quad \text{on } \Gamma, \\ \lim_{|\mathbf{x}| \rightarrow +\infty} |\mathbf{x}|^{(d-1)} \left(\nabla u \cdot \frac{\mathbf{x}}{|\mathbf{x}|} - iku \right) &= 0, \end{aligned} \tag{1}$$

where Δ is the Laplacian operator, ∇ the gradient operator and \mathbf{n}_Γ is the outward-directed unit normal vector to Ω^- . The spatial variable is $\mathbf{x} = (x, y)$ in 2D and $\mathbf{x} = (x, y, z)$ in 3D. The wavenumber k is related to the wavelength λ by the relation: $\lambda := 2\pi/k$. The boundary condition is a Neumann (sound-hard) boundary condition (but other boundary conditions could also be considered). Denoting by $\mathbf{a} \cdot \mathbf{b}$ the hermitian inner-product of two complex-valued vector fields \mathbf{a} and \mathbf{b} , then the last equation of system (1) is known as the Sommerfeld's radiation condition [23, 61] at infinity, which presents the outgoing wave to the domain. Various numerical methods have been developed to accurately solve the exterior scattering problem, e.g. finite elements [17, 18, 36, 40, 48, 50, 63, 69, Thompson and Pinsky], boundary elements [3, 4, 19, 21, 23, 28, 59, 61], enriched and various modified (wave-based, hybrid, asymptotic) finite elements [5, 6, 20, 33, 35, 38, 39, 43, 47, 53, 55]. A challenging and still outstanding question for numerics and applications is related to the so-called *high-frequency* regime, where the wavelength λ is very small compared with the characteristic length of the scatterer Ω^- [8, 69].

A first "exact" method consists in writing an integral equation on the surface Γ to represent the exterior field, and then solving this by means of a boundary element method, combined with fast evaluation algorithms (e.g. Fast Multilevel Multipole Methods (MFMM)[28, 29, 59], Adaptive Cross Approximation (ACA) [12, 76]) and preconditioned (matrix-free) Krylov subspace iterative

solvers (GMRES) [3, 4, 65]. This approach has the very interesting property to reduce the initial problem to a finite $(d - 1)$ -dimensional problem (set on Γ) and to be relatively stable thanks to the frequency regime (typically between $n_\lambda = 5$ to 10 points per wavelength are used, with $n_\lambda = \lambda/h$, and h the mesh-size). Nevertheless, the method is nontrivial to adapt to complex geometries, when the boundary conditions are modified or when a high-order of accuracy is required for the solution. In particular, developing efficient algorithms when the geometry is described with high accuracy remains an open question, most particularly for large wave numbers k since the number of degrees of freedom is potentially very large.

Another standard and very attractive engineering-type approach consists in truncating the exterior domain Ω^+ through the introduction of a fictitious boundary/layer Σ that surrounds the scatterer Ω^- , resulting in a bounded computational domain Ω_b with inner boundary Γ and outer boundary/layer Σ . When Σ is a boundary, an *Absorbing Boundary Condition (ABC)* [2, 11, 41, 42, 52, 60, 71, 72] is set on the fictitious boundary which introduces a truncation error even at the continuous level. This can be reduced by considering high-order (local) ABCs. A very popular alternative is, rather, to consider a surrounding absorbing layer and modifying the nonphysical media to obtain a Perfectly Matched Layer [13–15, 22].

In this paper, since our goal is to understand the quality of the NURBS-approximations used in Isogeometric Analysis (IGA) to approximate the wave field, we will consider simple boundary conditions where we can separate the continuous truncation error from the numerical approximation. High-order ABCs as well as PML will be treated in future work. When the problem is truncated, any numerical method can be used such as e.g. finite-element methods, the resulting linear system being then solved through well-adapted domain decomposition techniques [16, 31, 32, 37]. In the high frequency regime, it is well-known that the numerical solution suffers from a phase shift due to numerical dispersion which is called the *pollution error* [1, 8, 49, 51] of the finite element method. To maintain a prescribed pollution error, the mesh density n_λ should be increased faster than the wave number leading to high computational cost for high frequency problems. Therefore, the FEM is limited to a upper frequency bound for which the computational cost becomes prohibitive and increasing the order of the polynomial basis functions used in conventional FEM is required to reduce the pollution error, but it does not fully eliminate it.

Original techniques were recently proposed to reduce the pollution error. One approach is to replace the basis functions, usually a polynomial, with plane wave functions [39, 55]. Since pollution error can be viewed as a phase shift, another hybrid approach is to asymptotically approximate the phase of the solution and to reformulate the problem into an equivalent problem with a slowly varying unknown [5–7, 9, 10, 19, 38, 73]. Nevertheless, until now, the only viable solution for complex problems is to consider a sufficiently high-order polynomial basis into the FEM together with high-order meshes to represent accurately the geometry Γ .

The aim of the present paper is to analyze the quality of alternative approx-

imation methods based on B-splines and NURBS-based *Isogeometric Analysis (IGA)* and to compare the properties of these approximation schemes to more traditional, Lagrange-based high-order FEM. The goal of the paper is, specifically, to investigate the properties of the methods for large wave numbers k . Isogeometric analysis (IGA) was introduced over 10 years ago to streamline the transition from Computer-Aided Design to Analysis [27, 44]. The central idea of the approach is to use the same shape functions to approximate the field variables as those used to describe the geometry of the domain. As the geometry of computational domains is typically provided by Computer-Aided Design (CAD) software, it is sensible to use NURBS shape functions, which are the most commonly used functions in CAD.

Since 2005, IGA has been developed within the finite element, boundary element [57, 67, 68] and was enriched through partition of unity [62] IGA benefits from exact (and smooth) geometry representation and no loss of boundary details. This is fundamental in problems involving wave propagation or turbulence, as spurious wave scattering and vortices typically appear close to rough boundaries. IGA therefore alleviates the meshing burden as if the CAD model is modified, any modifications are immediately inherited by the shape functions used for analysis. When coupled to boundary element approaches, this suppresses completely the generation and regeneration of the mesh without any regeneration of the mesh [56, 57]. When used within a finite element approach, volume parameterization is still required, but the shape optimization process is simplified [34, 74]. Since CAD models are generated from the boundary data, *IGA Boundary Element Method (IGABEM)* is promising in bridging the gap between CAD and analysis offering a unique shape optimization package [24, 54, 58, 66].

IGA additionally offers higher inter-patch continuity, large support sizes and superior refinement capabilities compared to Lagrange-based approximation schemes. A recent review of IGA and its computer implementation aspects is available in [62]. GEOPDE is another open source IGA research tool [30] which provides a flexible coding structure. The application of IGA and its performance in solving one- and two-dimensional eigenvalue and interior Helmholtz problems have been studied [25, 26, 45, 46] in which IGA showed superiority to conventional FEM. However, the performance of IGA in solving exterior scattering problems and related pollution error is still not thoroughly investigated.

In this paper, we study the performance of IGA for solving time-harmonic exterior scattering problems in one, two and three dimensions. First, in Section 2, we briefly present the Isogeometric Analysis and the related NURBS and B-Spline basis functions. Then in Section 3, we consider a simple but meaningful one-dimensional scattering problem and provide a pseudo-code to solve it in IGA context. In Section 4, we consider two-dimensional problems and study the performance of IGA in estimating the scattered field of a duct and a circular cylinder. The in-house IGA and Lagrange-based FEM codes developed in this study were written in Matlab [®] following the GEOPDE platform for two- and three-dimensional examples. Finally, section 5 conclude the paper.

2. Isogeometric Analysis (IGA)

In IGA, both the physical model and the solution space are constructed by *B-spline/NURBS* functions. The model is constructed using a set of control points and *knot vectors* in each spatial direction. A knot vector is defined as a set of non-decreasing coordinates and represented by $\boldsymbol{\xi} = \{\xi_1, \xi_2, \dots, \xi_{(n+p+1)}\}$, where ξ_i is the i^{th} knot and i is the knot index, $i = 1, 2, \dots, n + p + 1$, where p is the polynomial order and n is the number of basis functions. Usually an *open knot vector* is used in IGA where the first and the last knot values appear $p + 1$ times. The elements in IGA are defined as non-zero knot intervals. If the knots are spaced equally within the knot vector, the knot vector is called *uniform* otherwise it is *non-uniform*.

2.1. B-Splines

A B-Spline basis function is defined recursively by the Cox-de Boor recursion formula starting with the zeroth order ($p = 0$) basis function:

$$\text{for } p = 0, N_{i,p}(\xi) = \begin{cases} 1 & \xi_i \leq \xi \leq \xi_{i+1}, \\ 0 & \text{otherwise,} \end{cases} \quad (2)$$

$$\text{for } p = 1, 2, 3, \dots \quad N_{i,p}(\xi) = \frac{\xi - \xi_i}{\xi_{i+p} - \xi_i} N_{i,p-1}(\xi) + \frac{\xi_{i+p+1} - \xi}{\xi_{i+p+1} - \xi_{i+1}} N_{i+1,p-1}(\xi), \quad (3)$$

where $0/0$ is defined to be zero. To compute the tangent and the normal functions, it is required to calculate the derivative of B-Splines. Efficient algorithms to calculate B-Splines and their derivatives can be found in the NURBS Book by Piegl and Tiller [64]. The first-order B-Spline functions are identical to their Lagrangian (FEM) counterparts. Also, the B-Splines number of required shape functions for a specific order is similar to that of the FEM. For example, three shape functions are required to construct second-order Lagrangian or B-Spline shape functions. In contrast to FEM, the shape functions in IGA are non-negative. The first derivative of the B-spline basis function is calculated using the following recursive formula:

$$\frac{d}{d\xi} N_{i,p}(\xi) = \frac{p}{\xi_{i+p} - \xi_i} N_{i,p-1}(\xi) - \frac{p}{\xi_{i+p+1} - \xi_{i+1}} N_{i+1,p-1}(\xi). \quad (4)$$

These functions exhibit C^{p-m_i} continuity across knot ξ_i where m_i is the multiplicity of the knot ξ_i . The B-Spline basis function corresponding to $p = 1, \dots, p = 5$ are shown in Fig. 1 where the knot vectors are defined to represent one element (knot span)

$$\boldsymbol{\xi} = \{\underbrace{0, \dots, 0}_{p+1 \text{ times}}, \dots, \underbrace{1, \dots, 1}_{p+1 \text{ times}}\}. \quad (5)$$

The quadratic B-Spline basis function for a 1D domain divided into four ele-

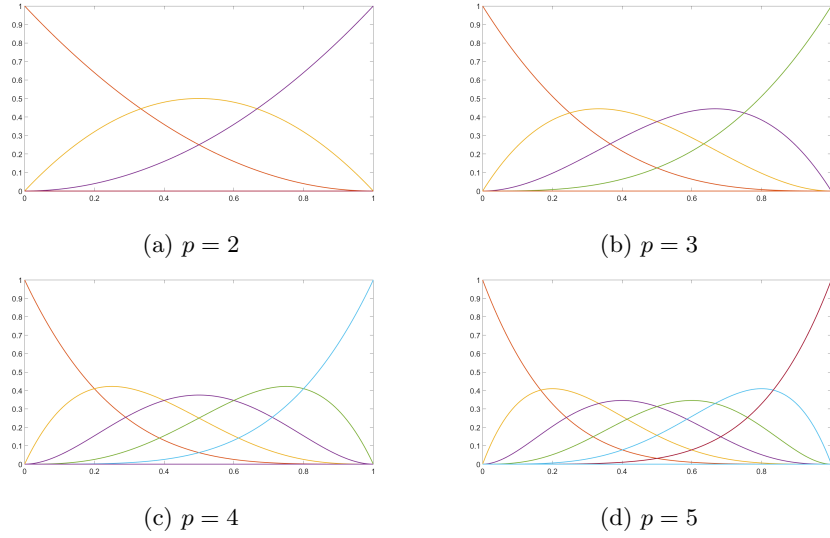


Figure 1: The B-Spline basis functions, $p = 1 \dots, p = 5$.

ments is shown in Fig. 2 where the knot vector is defined as

$$\xi = \{0, 0, 0, 1/4, 1/2, 3/4, 1, 1, 1\}.$$

The four non-zero knot spans correspond to the four elements of this one-dimensional domain.

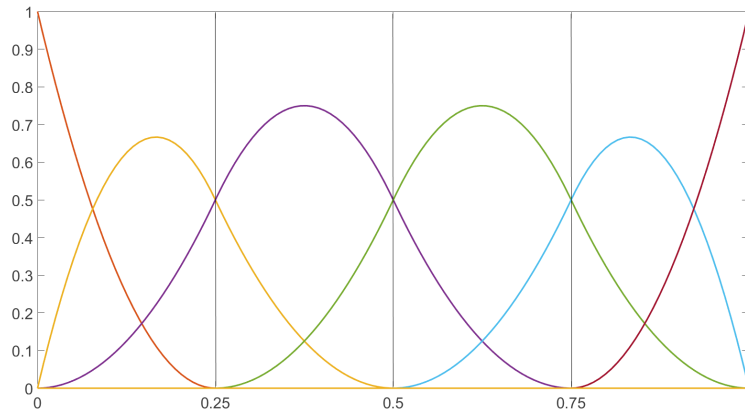


Figure 2: The quadratic ($p = 2$) basis function of a 1D domain defined over $\xi = \{0, 0, 0, 1/4, 1/2, 3/4, 1, 1, 1\}$ and representing four elements.

2.1.1. B-Spline curves

The piecewise-polynomial B-Spline curves in \mathbb{R}^d are constructed as a linear combination of B-Spline basis functions:

$$\mathbf{C}(\xi) = \sum_{i=1}^n N_{i,p}(\xi) \mathbf{B}_i. \quad (6)$$

The vector-valued coefficients of the basis functions in Eq.6 are control points ($\mathbf{B}_i \in \mathbb{R}^d, i = 1, 2, \dots, n$). Since $\mathbf{C}(\xi)$ is representing a curve, the corresponding control points, \mathbf{B}_i , are analogous to nodal coordinates in FEM. The resulting curve is C^{p-1} -continuous everywhere except at the location of the repeated knots with m_i multiplicity, where it is C^{p-m_i} -continuous. Therefore, when utilizing open knot vectors, the resulting curve is C^0 -continuous, tangent to the control polygon, and interpolatory at the start and the end of the curve and where the multiplicity is equal to the polynomial order. The support of a B-Spline basis function is local. In other words, moving a single control point can affect the geometry of no more than $p + 1$ elements of the curve.

2.1.2. B-Spline surfaces and volumes

A tensor product B-Spline surface is defined as:

$$\mathbf{S}(\xi, \eta) = \sum_{i=1}^n \sum_{j=1}^m N_{i,p}(\xi) M_{j,q}(\eta) \mathbf{B}_{i,j}, \quad (7)$$

where $\mathbf{B}_{i,j}, i = 1, 2, \dots, n, j = 1, 2, \dots, m$ are control points, and $N_{i,p}(\xi)$ and $M_{j,q}(\eta)$ are univariate B-Spline functions of orders p and q corresponding to the knot vectors $\Xi^1 = \{\xi_1, \xi_2, \dots, \xi_{n+p+1}\}$ and $\Xi^2 = \{\eta_1, \eta_2, \dots, \eta_{m+q+1}\}$, respectively. Similarly, a B-Spline volume is constructed as a tensor product of three B-Spline basis functions:

$$\mathbf{V}(\xi, \eta, \zeta) = \sum_{i=1}^n \sum_{j=1}^m \sum_{k=1}^l N_{i,p}(\xi) M_{j,q}(\eta) L_{k,r}(\zeta) \mathbf{B}_{i,j,k}. \quad (8)$$

2.2. NURBS

There are certain geometries, such as circles and ellipsoids, which B-Spline functions and Lagrange polynomials cannot represent exactly. In the CAD community, B-Spline functions are replaced with a more general form to overcome this shortcoming. This new form is called *Non-Uniform Rational B-Splines* (*NURBS*). NURBS shape functions are defined as:

$$R_i^p(\xi) = \frac{N_{i,p}(\xi)w_i}{W(\xi)} = \frac{N_{i,p}(\xi)w_i}{\sum_{i=1}^n N_{i,p}(\xi)w_i}, \quad (9)$$

where $\{N_{i,p}\}$ is a set of B-Spline basis functions and $\{w_i\}$ is a set of positive NURBS weights. By appropriate weight selection both polynomials and circular

arcs can be described. NURBS curves are defined as follows:

$$\mathbf{C}(\xi) = \sum_{i=1}^n R_i^p(\xi) \mathbf{B}_i. \quad (10)$$

Similarly, rational surfaces and volumes are defined in terms of rational basis functions as:

$$R_{i,j}^{p,q}(\xi, \eta) = \frac{N_{i,p}(\xi) M_{j,q}(\eta) w_{i,j}}{\sum_{i=1}^n \sum_{j=1}^m N_{i,p}(\xi) M_{j,q}(\eta) w_{i,j}}, \quad (11)$$

$$R_{i,j,k}^{p,q}(\xi, \eta, \zeta) = \frac{N_{i,p}(\xi) M_{j,q}(\eta) L_{k,r}(\zeta) w_{i,j,k}}{\sum_{i=1}^n \sum_{j=1}^m \sum_{k=1}^l N_{i,p}(\xi) M_{j,q}(\eta) L_{k,r}(\zeta) w_{i,j,k}}. \quad (12)$$

If the weights are all equal, NURBS basis functions will reduce to their B-Spline counterparts and the corresponding curve becomes a non-rational polynomial again. The geometry of an element in IGA can be expressed as:

$$\mathbf{x}^e(\tilde{\xi}) = \sum_{a=1}^{n_{en}} \mathbf{P}_a^e R_a^e(\tilde{\xi}), \quad (13)$$

where $n_{en} = (p+1)^{d_p}$ and d_p is the spatial dimension. The field $\mathbf{u}(\mathbf{x})$ is similarly expressed as:

$$\mathbf{u}(\tilde{\xi}) = \sum_{a=1}^{n_{en}} \mathbf{d}_a^e R_a^e(\tilde{\xi}), \quad (14)$$

where $\mathbf{d}_a^e R_a^e$ is the set of control variables. As a consequence of the parametric definition of basis functions in IGA, it is necessary to consider two mapping for integration

$$\begin{aligned} \int_{\Omega} f(\mathbf{x}) d\Omega &= \sum_{e=1}^{n_{el}} \int_{\Omega^e} f(\mathbf{x}) d\Omega = \sum_{e=1}^{n_{el}} \int_{\tilde{\Omega}} f(\mathbf{x}(\tilde{\xi})) |J_{\tilde{\xi}}| d\tilde{\Omega} \\ &= \sum_{e=1}^{n_{el}} \int_{\tilde{\Omega}} f(\mathbf{x}(\tilde{\phi}^e(\tilde{\xi}))) |J_{\tilde{\xi}}| |J_{\tilde{\xi}}| d\tilde{\Omega}, \end{aligned} \quad (15)$$

where $\tilde{\Omega}$ and $\hat{\Omega}$ are representing the parent and parameter spaces respectively, and Ω is the physical space.

3. A one-dimensional toy scattering problem

Now, let us come back to our physical problem by considering the simple one-dimensional Neumann scattering problem solved in [5]:

$$\begin{aligned} \partial_x^2 u + k^2 u &= 0, & \text{in } \Omega_b &= (0, 1), \\ \partial_x u &= ik, & \text{at } \Gamma &= \{0\}, \\ \partial_x u - ik u &= 0, & \text{at } \Sigma &= \{1\}. \end{aligned} \quad (16)$$

The variational formulation is given as:

$$\int_{\Omega_b} \{\partial_x u \partial_x \bar{v} - k^2 u \bar{v}\} dx - ik(u\bar{v})(1) = -(g\bar{v})(0), \quad \forall v \in H^1(\Omega_b). \quad (17)$$

The solution of the one-dimensional BVP (16) is: $u^{\text{ex}}(x) = e^{ikx}$, which represents the scattering of an incident plane wave by the left half space. An exact transparent boundary condition is considered for the fictitious boundary at $\Sigma = \{1\}$ using the Dirichlet-to-Neumann operator $\Lambda = ik$ on Σ , resulting in the transparent boundary condition $\partial_x u = \Lambda u$.

It is shown that the Finite Element approximated solution of the BVP described in system (16) suffers from the pollution error [5, 48, 51, 69]. The pollution error has a direct relation with the wavenumber k and increases with frequency. It is the polynomial basis used in conventional FEM which is inadequate to represent the wavefield. Since the main difference of the IGA with the conventional FEM is in the selection of the basis function, it is interesting to study the performance of IGA specially in high frequencies and its pollution error. We divide the computational bounded domain Ω_b into N_h uniform non-zero knot spans (elements) so that each segment length is $1/h$. We define the density of discretization as the number of elements per wavelength and denote it with $n_\lambda = \lambda/h$. A Matlab code was prepared to solve the above example in IGA context. The stiffness matrix for a single element is calculated and assembled into the global matrix following these steps:

- Set $K = 0$
- loop over number of elements
- Loop over Gaussian points $\tilde{\xi}_i, \tilde{w}_i$
 1. find the parametric coordinate $\xi = \tilde{\phi}^e(\tilde{\xi}_j)$
 2. compute the basis function $R_a^e(a = 1, \dots, p+1)$ at point ξ
 3. define vector $\mathbf{R} = [R_1^e, R_2^e, \dots, R_{p+1}^e]$
 4. compute the basis function derivatives $R_{a,\xi}^e(a = 1, \dots, p+1)$ at point ξ
 5. define vector $\mathbf{R}_\xi = [R_{1,\xi}^e, R_{2,\xi}^e, \dots, R_{p+1,\xi}^e]$
 6. compute $|J_\xi| = \|\mathbf{R}_\xi \mathbf{P}\|$
 7. compute $|J_{\tilde{\xi}}| = 0.5(\xi_{i+1} - \xi_i)$
 8. compute shape function derivatives $\mathbf{R}_x = J_\xi^{-1} \mathbf{R}_\xi^T$
 9. Set $\mathbf{k}^e = \mathbf{k}^e + \mathbf{R}_x \mathbf{R}_x^T |J_\xi| |J_{\tilde{\xi}}| \tilde{w}_j - k^2 \mathbf{R} \mathbf{R}^T |J_\xi| |J_{\tilde{\xi}}| \tilde{w}_j$
 10. End of both loops

then, the Neumann boundary condition is applied and the resulting linear system is solved to find the estimated solution in IGA. we plot the approximate solution u_h for $k = 40$ as well as the exact solution in Fig.3. The density of discretization for IGA approximations is $n_\lambda = 10$. As expected, IGA ($p = 1$) also suffers from pollution error. B-Spline functions generated with $p = 0$ and

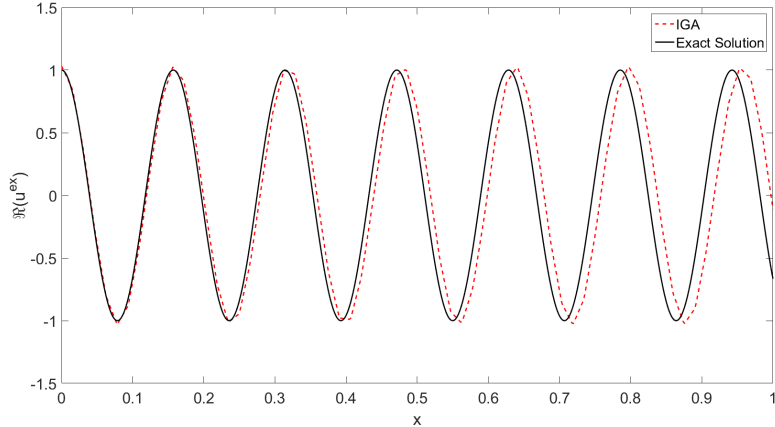


Figure 3: The pollution error is the numerical solutions $\Re(u_h)$ of IGA ($p = 1$), is visible when compared with the exact solution $\Re(u^{\text{ex}})$, for $k = 40$ and $n_\lambda = 10$.

$p = 1$ orders will result in the same piecewise constant and linear functions as standard FEM shape functions. The differences are found for cases using higher order shape functions. The p^{th} order B-Spline function has $p - 1$ continuous derivatives across the element boundaries. Each B-Spline basis is point-wise non-negative over the entire domain as a result, all of the entries of the mass matrix will be positive. The support of a B-Spline function of order p is always $p + 1$ knot spans. Therefore, higher order B-Spline function has support over much larger portion of the domain when compared to classical FEM. The total number of functions that any given function shares support with (including itself) is $2p + 1$ regardless of whether a FEM basis or a B-Splines is used. Hence, using higher order B-Spline basis functions will provide improved support compared to standard FEM without increasing the number of required shape functions. Increasing the order of Lagrangian polynomials will increase the amplitude of oscillations in FEM; this problem is eliminated in IGA as a result of non-negativity and non-interpolatory nature of B-Splines shape functions.

Next we investigate the effect of increasing the basis function order on the accuracy of the IGA approximation again for discretization density $n_\lambda = 10$. As expected increasing the order p reduces the error drastically as shown in Fig. 4. There is no visible pollution error in the IGA approximation if we keep all the parameters unchanged and increase the order of the basis function. One can reduce the error further by increasing the discretization density n_λ . For the numerical solution $f^{\text{calc}}(\mathbf{x})$, $\mathbf{x} \in \Omega_b$, we define the L_2 -error (in dimension d) as

$$\epsilon_2 = \left\{ \int_{\Omega_b} |f^{\text{calc}}(\mathbf{x}) - f^{\text{ex}}(\mathbf{x})|^2 d\mathbf{x} \right\}^{1/2}. \quad (18)$$

The evolution of the L_2 -error vs. the discretization density n_λ is reported in Fig. 5 for the wavenumber $k = 40$. For completeness, we also plot the curves relative

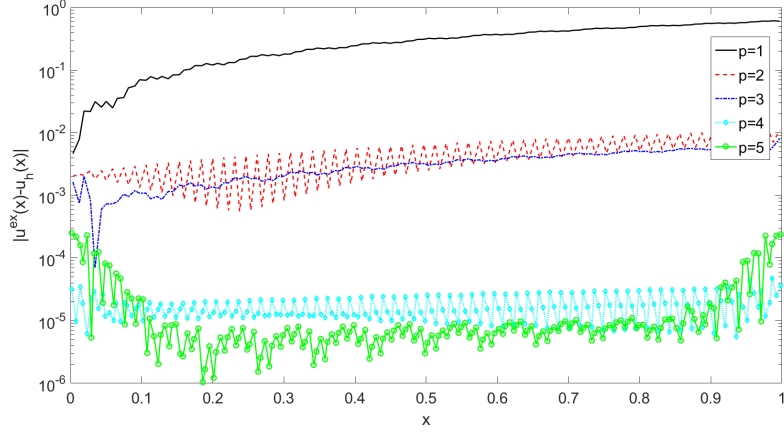


Figure 4: Absolute error $|u^{\text{ex}} - u_h|$ for $k = 40$, $n_\lambda = 10$ and various approximation orders $p = 1, \dots, 5$.

to the function x^{-p} , for $x = n_\lambda$. We can see that the error curves fits very well the polynomial curves. This lets to think that the pollution error is extremely weak, most particularly when p is large enough ($p \geq 3$) even for a small density n_λ . Next we present the evolution of the L_2 -error with the wavenumber k in Fig.

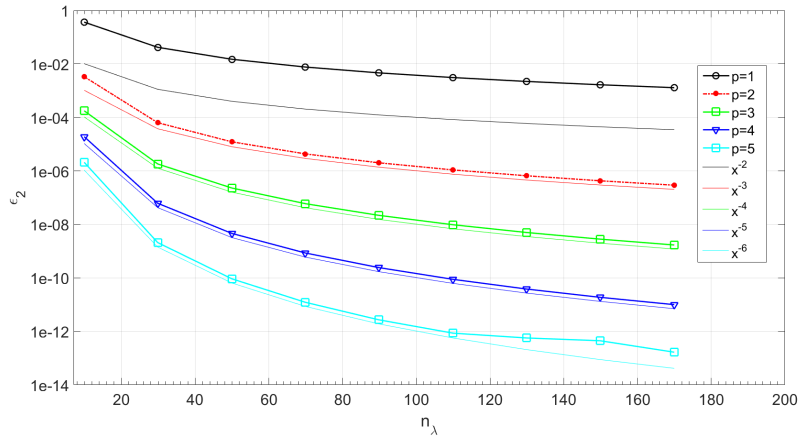


Figure 5: Evolution of the error ϵ_2 vs. the discretization density n_λ for $k = 40$ and $p = 1, \dots, 5$ (the slopes of the functions x^{-p} are plotted for comparison).

6 for a fixed density $n_\lambda = 20$, where for $p = 3$ and higher the pollution error is not visible and the error remains constant even for extremely high frequencies.

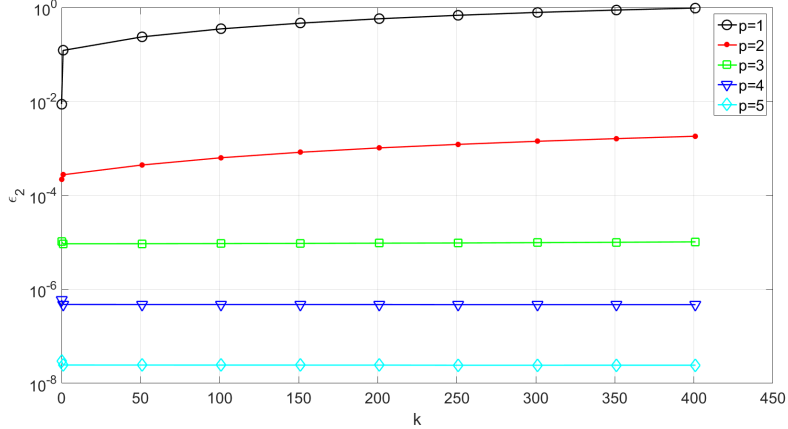


Figure 6: Evolution of the error ϵ_2 vs. the wavenumber k , for $n_\lambda = 20$ and various approximation orders $p = 1, \dots, 5$.

4. Two-dimensional examples

In this section, we evaluate the performance of IGA in solving two-dimensional acoustic problems. First, we consider a two-dimensional duct problem with rigid walls [47]. Next, we propose two examples of cylindrical disk scattering problems where we obtain, first the scattered field of a particular mode, and then the scattering of the disk subject to an incident plane wave. By separating the truncation error from the numerical basis approximation, we fairly investigate the performance of IGA in solving the two-dimensional sound-hard exterior scattering problem and numerically analyze the related pollution error.

4.1. The duct problem

The duct model is shown in Fig. 7 where $\Omega_b = [0, 2] \times [0, 1]$. We denote the outward boundary unit normal with v and assume that the lower and upper walls are rigid. We solve the Helmholtz problem stated in Eq. 19 for the acoustic pressure u

$$\begin{aligned}
 \Delta u + k^2 u &= 0, & \text{in } \Omega_b, \\
 \frac{\partial u}{\partial v} &= \cos(m\pi y), & \text{on } x = 0, \\
 \frac{\partial u}{\partial v} + iku &= 0, & \text{on } x = 2, \\
 \frac{\partial u}{\partial v} &= 0, & \text{on } y = 0, 1,
 \end{aligned} \tag{19}$$

where $m \in \mathbb{N}$ is the mode number. An inhomogeneous boundary condition is applied on the inlet boundary ($x = 0$) and an absorbing (and transparent for $m = 0$) boundary condition is set on the outlet boundary ($x = 2$). Since the boundaries at $y = 0, 1$ are assumed to be perfectly rigid, the normal derivative

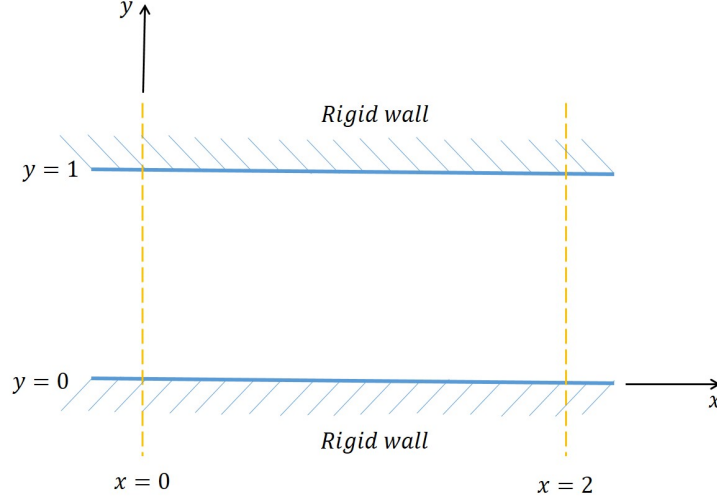


Figure 7: The duct domain.

of the acoustic pressure vanishes on these boundaries. The exact solution of problem (19) with ABC is as follows

$$u^{\text{ex}}(x, y) = \cos(m\pi y)(A_1 e^{-ik_x x} + A_1 e^{ik_x x}), \quad (20)$$

where $k_x = \sqrt{k^2 - (m\pi)^2}$ and the coefficients A_1 and A_2 are obtained from

$$\begin{pmatrix} k_x & -k_x \\ (k - k_x)e^{-2ik_x} & (k + k_x)e^{2ik_x} \end{pmatrix} \begin{pmatrix} A_1 \\ A_2 \end{pmatrix} = \begin{pmatrix} 1 \\ 0 \end{pmatrix}.$$

Solving this 2×2 linear system, we are able to get the expression of the solution to the duct problem with ABC. This is a very interesting point since we therefore can only focus the analysis on the finite-dimensional approximation without including the (continuous) truncation error. The cut-off frequency is $m_{\text{cut-off}} = k/\pi$. If the mode is such that $m \leq m_{\text{cut-off}}$, the solution is representing propagating modes and if $m > m_{\text{cut-off}}$ the solution corresponds to evanescent modes. The real parts of the exact and estimated IGA solutions are presented in Fig. 8a and 8b, respectively, for $k = 40$, $m = 2$ (propagative mode), $p = 3$, and $n_\lambda = 10$. The corresponding absolute error $|u^{\text{ex}} - u_h|$ is plotted in Fig. 9, the maximal value being of the order of 10^{-6} . For the propagative mode $m = 2$, the evolution of the L_2 -error with respect to the discretization density n_λ is shown in Fig. 10 for $k = 40$, and according to the wave number k in Fig. 11, where $n_\lambda = 10$. We can see that the error is very low, decreasing strongly with p . In addition, the error does not seem to depend on k as observed in Fig. 11, meaning that the pollution error is negligible.

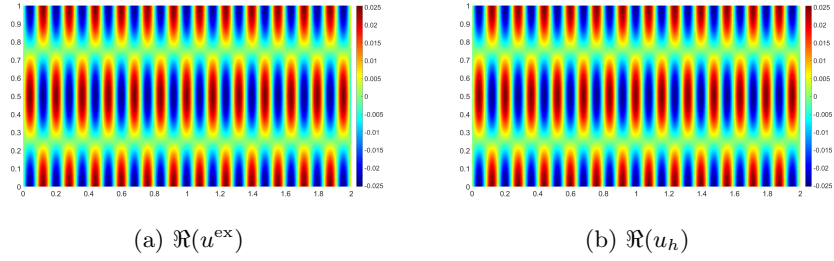


Figure 8: Comparing (a) the real parts of the exact solution u^{ex} and (b) the numerical solution u_h for $k = 40$, $m = 2$, $p = 3$, and $n_\lambda = 10$.

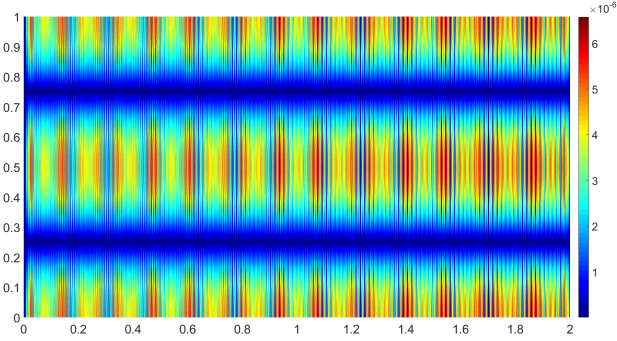


Figure 9: Absolute error $|u^{\text{ex}} - u_h|$ for $k = 40$, $m = 2$, $p = 3$ and $n_\lambda = 10$.

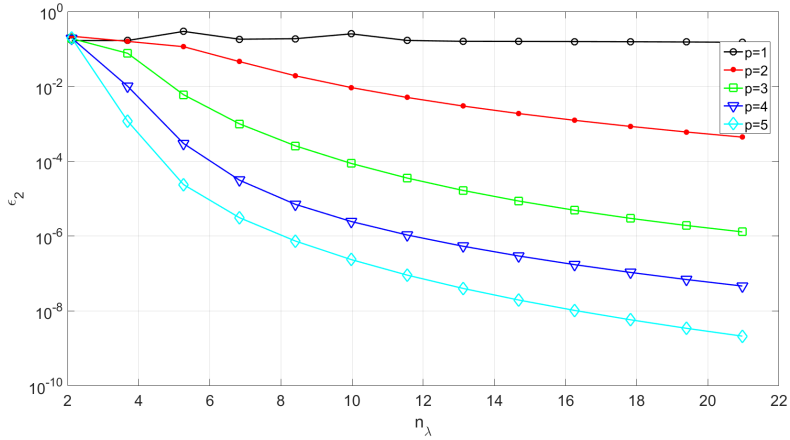


Figure 10: Evolution of the error ϵ_2 vs. the discretization density n_λ , for $k = 40$, $m = 2$ and $p = 1, \dots, 5$.

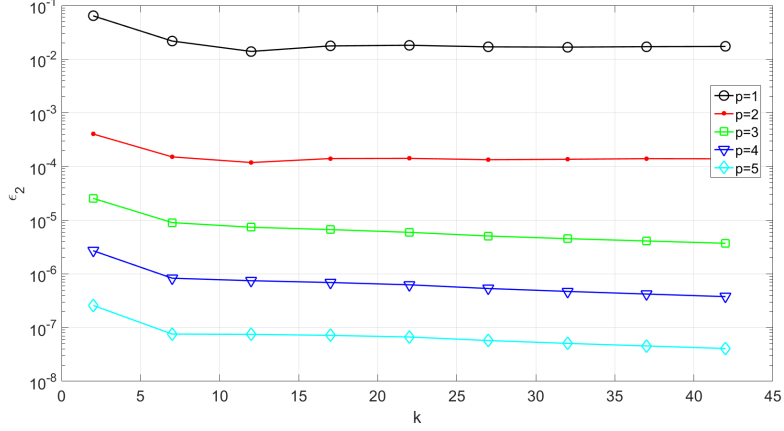


Figure 11: Evolution of the error ϵ_2 vs. the wavenumber k , for $m = 2$, $n_\lambda = 10$ and $p = 1, \dots, 5$.

4.2. The sound-hard circular cylinder problem

To further study the performance of IGA, we analyze the scattering by a 2D circular cylinder $\Omega^- = \mathcal{D}_0$ of radius R_0 centered at the origin, with boundary $\Gamma = \mathcal{C}_0$ (circle of radius R_0). The scatterer is surrounded by an outer fictitious circular boundary $\Sigma = \mathcal{C}_1$, again centered at the origin and with radius $R_1 > R_0$. Hence, the computational domain Ω_b is the annulus bounded between the inner \mathcal{C}_0 and outer \mathcal{C}_1 boundaries. On the exterior boundary, we set the symmetrical second-order Bayliss-Turkel ABC [2] given by

$$\partial_{\mathbf{n}_\Sigma} u + \left(-ik + \frac{\kappa}{2} - \frac{\kappa^2}{8(\kappa - ik)}\right)u - \frac{1}{2(\kappa - ik)}\partial_s^2 u = 0, \quad \text{on } \Sigma, \quad (21)$$

setting \mathbf{n}_Σ at the outwardly directed unit normal to Σ , $\partial_{\mathbf{n}_\Sigma} := \partial_r$ the normal derivative, $\kappa = 1/R_1$ is the curvature of $\Sigma := \mathcal{C}_1$ and $\partial_s^2 := R_1^{-2}\partial_\phi^2$ is the second-order curvilinear derivative on \mathcal{C}_1 . In the above expressions, the polar coordinate system is denoted by (r, ϕ) .

4.2.1. Scattering by a sound-hard circular cylinder : mode-by-mode analysis

We consider an incident wave with a fixed mode m which is described as

$$u_m^{\text{inc}}(\mathbf{x}) = J_m(kr)e^{im\phi}, \quad m \in \mathbb{Z}, \quad (22)$$

where J_m is the m -th order Bessel's function. The exact exterior modal solution u_m^{ex} to the truncated scattering problem of the inner sound-hard circular cylinder \mathcal{C}_0 in polar coordinate (r, ϕ) is given as

$$u_m^{\text{ex}}(\mathbf{x}) = \left(a_m H_m^{(1)}(kr) + b_m H_m^{(2)}(kr)\right)e^{im\phi}, \quad r \geq R_0, \quad m \in \mathbb{Z}. \quad (23)$$

The functions $H_m^{(1)}$ (respectively $H_m^{(2)}$) is the first-kind (respectively second-kind) Hankel function of order m . The Neumann boundary condition is applied

on \mathcal{C}_0 and the ABC on \mathcal{C}_1 , resulting in the following linear system of equations to obtain the two unknown coefficients a_m and b_m

$$a_m = -\frac{A_{22}^m J_m'(kR_0)}{D_m}, \quad b_m = -\frac{A_{21}^m J_m'(kR_0)}{D_m}, \quad (24)$$

where $D_m = A_{11}^m A_{22}^m - A_{21}^m A_{12}^m$ and

$$\begin{cases} A_{11}^m = H_m^{(1)}(kR_0), & A_{21}^m = kH_m^{(1)}(kR_1) - \mathcal{B}_m H_m^{(1)}(kR_1), \\ A_{12}^m = H_m^{(2)}(kR_0), & A_{22}^m = kH_m^{(2)}(kR_1) - \mathcal{B}_m H_m^{(2)}(kR_1), \\ \mathcal{B}_m = -\left(\alpha_m \frac{m^2}{R_1^2} + \beta_m\right), & \alpha_m = -\frac{1}{2ik} \left(1 + \frac{i}{kR_1}\right)^{-1}, \\ \beta_m = -ik + \frac{1}{2R_1} + \frac{1}{8iR_1(1+kR_1)}. \end{cases}$$

The notation $f' := \partial_r f$ designates the radial derivative of a given function $f(r)$. Quite similarly to the duct problem, the modes $m \in \mathbb{N}$ such that $|m| < kR_0$ are propagating. For $|m| > kR_0$, the modes are evanescent (and therefore are not visible in the far-field since they do not propagate). A special case corresponds to $|m| \approx kR_0$, which is related to grazing modes that are tangent to the scatterer (generating glancing rays).

The modal analysis is meaningful since we can understand the accuracy of the IGA approximations thanks to the spatial frequencies m and the dimensionless wavenumber kR_0 . In addition, this also helps in clarifying what can be expected for the case of the scattering of a plane wave by the disk since the exact solution is built as a modal series expansion of the elementary mode solutions. We fix $R_0 = 1$ and $R_1 = 2$. The IGA model was generated using four identical patches as shown in Fig. 12. We report the real parts of u_m^{ex} and u_h on Fig. 13 where the wavenumber is $k = 40$ and $m = 2$ (which is a propagative mode since $m = 2 \leq kR_0 = 40$). For IGA, the order of the method is $p = 3$ for a density of discretization points per wavelength $n_\lambda = 10$. We clearly see that the wavefield is accurately computed since the absolute error $|u_m^{\text{ex}} - u_h|$ is below 10^{-5} as it can



Figure 12: Outline of the patches.

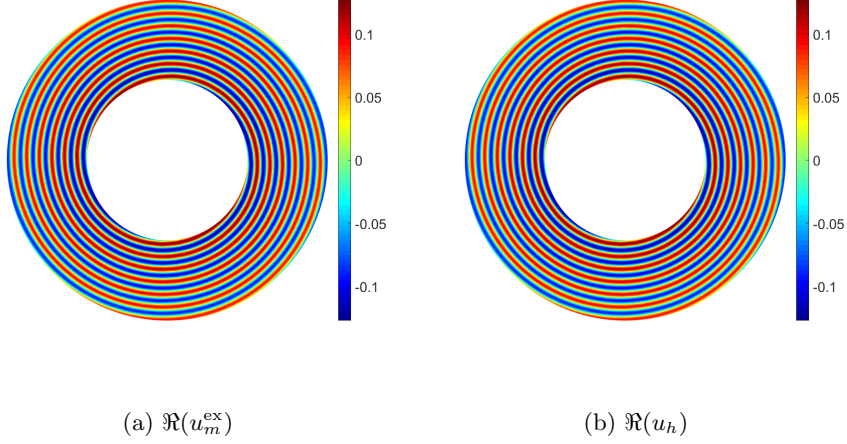


Figure 13: Comparing (a) the real parts of the exact solution u_m^{ex} and (b) the numerical IGA solution u_h for $k = 40$, $m = 2$, $p = 3$, and $n_\lambda = 10$.

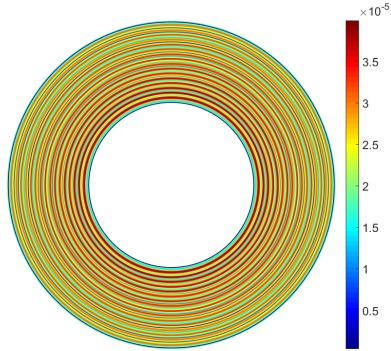


Figure 14: Absolute error $|u_m^{\text{ex}} - u_h|$ for $k = 40$, $m = 2$, $p = 3$ and $n_\lambda = 10$.

be deduced from Fig. 14. The convergence graph presented in Fig. 15 shows the effect of increasing the discretization density n_λ for various orders p of the underlying IGA basis function on the L_2 -error ϵ_2 for $k = 40$ and $m = 2$. We see that the error curves follow the slopes of the curves x^{-p} (for $x = n_\lambda$). The evolution of the error ϵ_2 with respect to the wavenumber k is plotted on Fig. 16, for $m = 2$, with various approximation orders p of IGA and $n_\lambda = 10$. We can observe that the accuracy does not really depend on k for a fixed density when $p \geq 2$. In fact, one question is to know how this error depends on m for a fixed k . To analyze this point, we report on Fig. 17 the evolution of the error ϵ_2

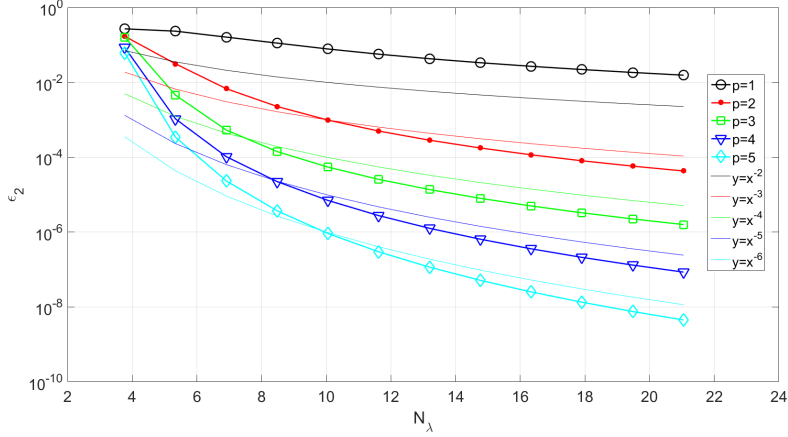


Figure 15: Evolution of the error ϵ_2 vs. the discretization density n_λ for $k = 40$, $m = 2$ and $p = 1, \dots, 5$.

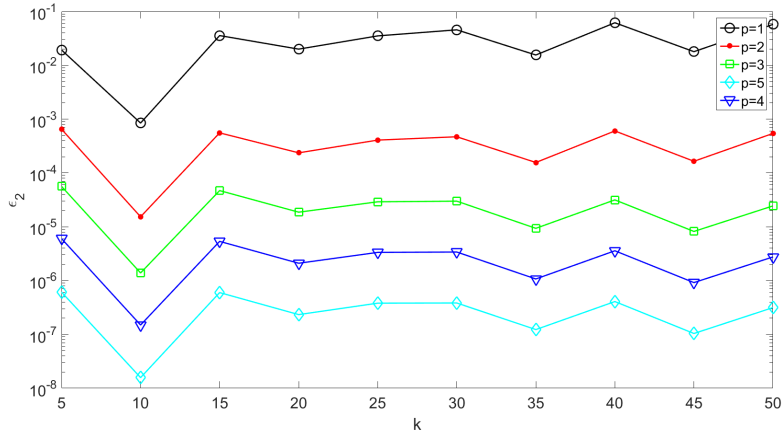


Figure 16: Evolution of the error ϵ_2 vs. the wavenumber k for $m = 2$, $n_\lambda = 10$ and $p = 1, \dots, 5$.

thanks to m and various values of p , for a fixed framework $k = 40$ and $n_\lambda = 10$. We can remark that the error almost does not depend on m , the value $m = 2$ corresponding more or less to the worst situation. When $|m| \geq kR_0$, the error starts decreasing strongly. Indeed, the solution is then evanescent and rather corresponds to the solution of a positive definite problem (similar to a shifted laplacian). Such solutions are easier to compute with high accuracy than for a propagative solution and do not participate to the far-field. From these remarks, and since the solution of the scattering of a plane wave by a disk is computed as the superposition of these modes solutions, we can expect that similar graphs to Fig. 16 could be obtained for the full plane wave problem (see subsection

4.2.2).

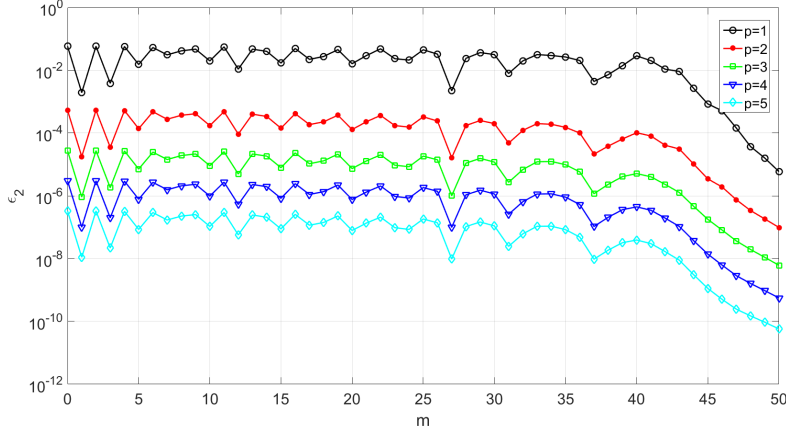


Figure 17: Evolution of the error ϵ_2 vs. the mode m , for $k = 40$, $n_\lambda = 10$ and $p = 1, \dots, 5$.

4.2.2. Scattering by a sound-hard circular cylinder: plane wave scattering

In the previous section, the scattering problem of a sound-hard circular cylinder was found for a given harmonics m . The solution for the full plane wave incidence can be constructed by superposition as a Fourier series expansion of these harmonics. In this section, we consider an incident plane wave $u^{\text{inc}}(\mathbf{x}) = e^{i\mathbf{k}\cdot\mathbf{x}}$, where \mathbf{d} is the incidence direction $\mathbf{d} = (\cos(\theta^{\text{inc}}), \sin(\theta^{\text{inc}}))^T$ and θ^{inc} is the scattering angle. Because of the symmetry of the problem, we fix the incidence direction to $\mathbf{d} = (1, 0)^T$ and the scatterer as the unit disk with $R_0 = 1$. Similar to the previous example, the second-order Bayliss-Turkel ABC (21) is placed on the circle with radius $R_1 = 2$. We consider the following exact solution [53] to analyze the pollution and approximation errors and to avoid the domain truncation error

$$u^{\text{ex}} = \sum_{m \in \mathbb{Z}} (i)^m u_m^{\text{ex}}, \quad (25)$$

where u_m^{ex} is given by Eq. (23). To get an accurate reference solution, we truncate the above series expansion by summing up on m from $-m^{\text{max}}$ to $m^{\text{max}} := [kR_0] + 30$ (where $[r]$ denotes the integer part of a real-valued positive number r). We report on Fig. 18 the real parts of u^{ex} (see Fig. 18a) and u_h (see Fig. 18b), where the wavenumber is $k = 40$. For IGA, the order of the method is $p = 3$ for a density of discretization points per wavelength $n_\lambda = 10$. We clearly see that the wavefield is accurately computed since the absolute error $|u_m^{\text{ex}} - u_h|$ is below 10^{-4} as it can be deduced from Fig. 19a.

The evolution of the error with respect to the discretization density n_λ for $p = 1, \dots, 5$, is shown in Fig. 20 where the wavenumber k is equal to 40. This shows that if p is larger than 3, then the ϵ_2 -error starts being small (e.g. less

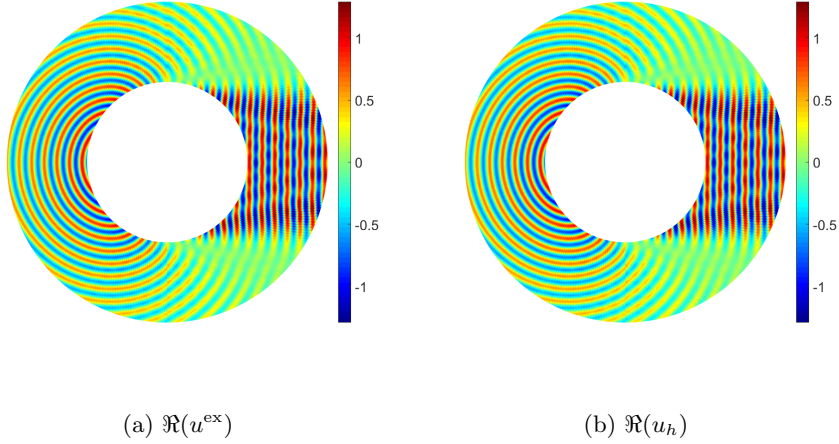


Figure 18: Comparing the real parts of (a) the exact solution u^{ex} and (b) the numerical IGA solution u_h for $k = 40$, $p = 3$, and $n_\lambda = 10$.

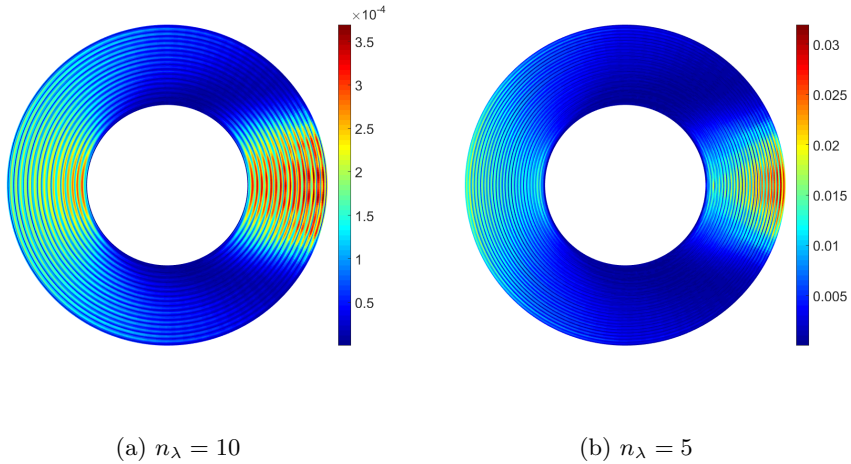


Figure 19: Absolute error $|u^{\text{ex}} - u_h|$ for $k = 40$ and $p = 3$. The discretization density is (a) $n_\lambda = 10$ and (b) $n_\lambda = 5$.

than 10^{-2}) even for small densities n_λ (typically $n_\lambda \geq 6$). In particular, the error gap between $p = 2$ and $p = 3$ is important. In addition, for $n_\lambda = 5$, we can see on Fig. 19b that the error remains acceptable but of the order of 2–3%, where the maximal error is located where the amplitude of the wavefield is larger and the solution oscillates.

The evolution of the error ϵ_2 according to the wavenumber k is depicted in

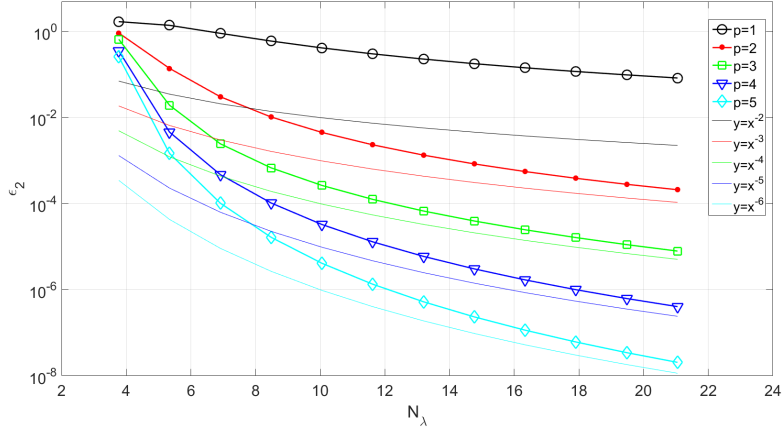


Figure 20: Evolution of the error ϵ_2 vs. the discretization density n_λ for $k = 40$ and $p = 1, \dots, 5$.

Fig. 21. For this test, we fix $n_\lambda = 10$ and $p = 3$ for IGA. As we can see, the error almost does not depend on k , which means that the pollution error is very small. Again, a very good accuracy is obtained for $p \geq 3$, the p -th order of the method being then visible for p larger than 3. The dependence with respect to k seems more present for the lowest orders IGA approximations, i.e. $p = 1$ and $p = 2$.

We consider now a much higher wavenumber, i.e. $k = 200$. We report on Fig. 22b the real part of the numerical IGA solution for $p = 3$ and the low density of discretization points per wavelength $n_\lambda = 5$. The solution can be compared with the exact solution available on Fig. 22a. We immediately see

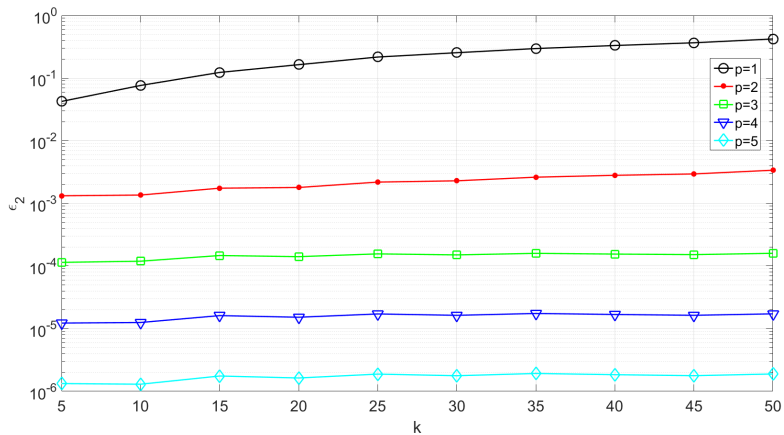


Figure 21: Evolution of the error ϵ_2 vs. the wavenumber k , for $n_\lambda = 10$ and $p = 1, \dots, 5$.

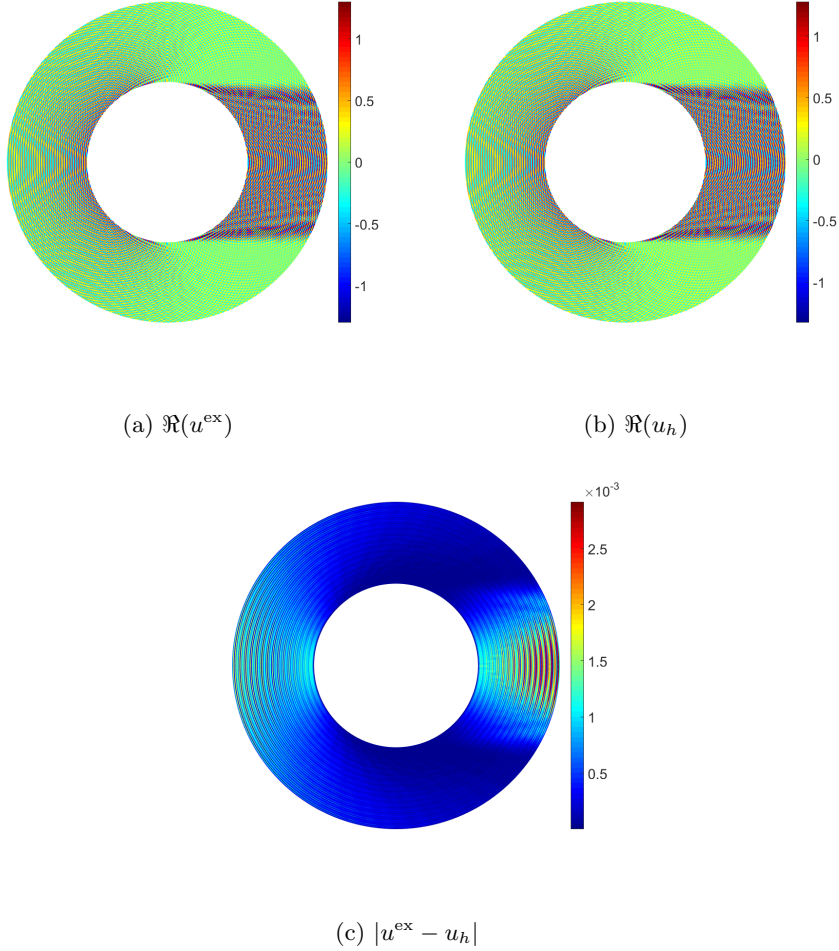


Figure 22: Comparing the real parts of (a) the exact solution u^{ex} and (b) the numerical IGA solution u_h , and (c) absolute error $|u^{\text{ex}} - u_h|$, for $k = 200$, $p = 3$, and $n_\lambda = 5$.

that the two solutions are the same. This is confirmed on Fig. 22c where we report the absolute error $|u^{\text{ex}} - u_h|$ between the two solutions. This error is of the order of 10^{-3} , which is very small.

Finally, we report now the results for a very high wavenumber $k = 500$ on Fig. 23. We provide the real parts of the exact solution u^{ex} (Fig. 23a) and IGA numerical solution u_h (Fig. 23b). For IGA, the order of the method is $p = 3$ for a density of discretization points per wavelength $n_\lambda = 5$. Even if there is no visible difference, plotting the absolute error on Fig. 23c shows an order of about 6%. This can be made much smaller by increasing the order of approximation to $p = 4$ by keeping $n_\lambda = 5$ as seen on Fig. 23d, leading to an

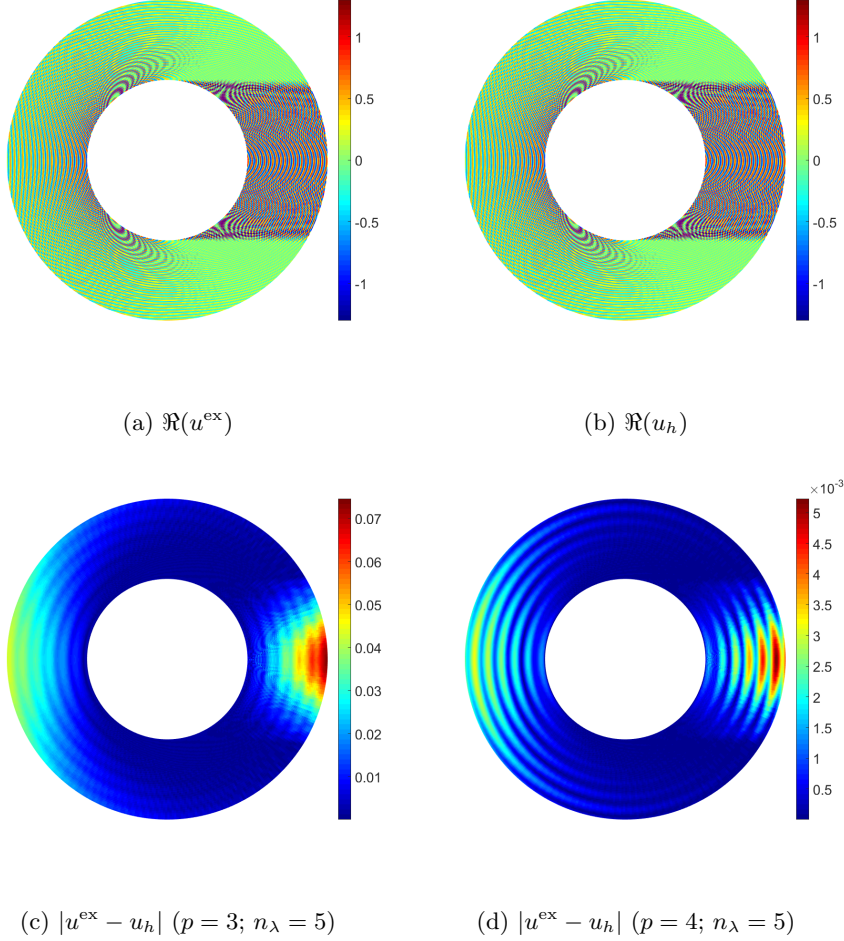


Figure 23: Comparing the real parts of (a) the exact solution u^{ex} and (b) the numerical IGA solution u_h for $k = 500$, $p = 3$, and $n_\lambda = 5$. Absolute error $|u^{\text{ex}} - u_h|$ for (c) $p = 3$; $n_\lambda = 5$, (d) $p = 4$; $n_\lambda = 5$.

error of about 10^{-3} .

5. Conclusions

The exterior Helmholtz problem was solved for one-, and two-dimensional problems using IGA. An absorbing boundary condition is applied on a fictitious boundary to truncate the infinite space. This truncation introduces an error in the numerical solution. To better study the performance of IGA, the truncation error was included in the analytical solution so that the pollution error is distinguishable. It is shown that IGA is less prone to pollution error even for large

wavenumbers. Also, the pollution error was considerably reduced by increasing the order of the basis functions which can be conveniently done in IGA, even for small densities of discretization points per wavelength. The smooth representation of the boundaries in IGA and the low pollution error makes IGA a suitable platform for scattering analysis, in particular for prospecting high frequency problems. For high wave numbers it will be interesting to investigate the use of different spaces for geometry and field approximations. Low order geometry description can be used with high order field approximations to investigate super and sub-geometric approaches, also known as GIFT [75].

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