A GENERALIZATION OF THE WIENER RATIONAL BASIS
FUNCTIONS ON INFINITE INTERVALS
PART I – DERIVATION AND PROPERTIES

AKIL C. NARAYAN AND JAN S. HESTHAVEN

Abstract. We formulate and derive a generalization of an orthogonal rational-
function basis for spectral expansions over the infinite or semi-infinite inter-
val. The original functions, first presented by Wiener [30], are a mapping and
weighting of the Fourier basis to the infinite interval. By identifying the Fourier
series as a biorthogonal composition of Jacobi polynomials/functions, we are
able to define generalized Fourier series which, appropriately mapped to the
whole real line and weighted, form a generalization of Wiener’s basis functions.
It is known that the original Wiener rational functions inherit sparse Galerkin
matrices for differentiation, and can utilize the fast Fourier transform (FFT)
for computation of the modal coefficients. We show that the generalized ba-
sis sets also have a sparse differentiation matrix and we discuss connection
problems, which are necessary theoretical developments for application of the
FFT.

1. Introduction

The approximation of a function by a finite sum of basis functions has long been
a hallmark tool in numerical analysis. Over the finite interval much is known about
expansion properties and periodic Fourier expansions or polynomial expansions are
well-studied. On infinite intervals there are complications due to the unbounded
domain on which approximation is necessary. Nevertheless many basis sets have
been successfully investigated in this case: the Hermite functions provide a suit-
able method for approximation when it can be assumed that the function decays
exponentially; for functions that do not decay exponentially, the so-called mapped
Chebyshev rational functions can fill the void and open up the possibility for utiliz-
ing the fast Fourier transform (FFT); additionally, a Fourier basis mapped to the
real line has been explored and provides an additional method for function approx-
imation over the infinite interval. This last basis set serves as an inspiration for the
family of basis sets proposed in this paper.

Despite the available methods for function approximation over the infinite inter-
val, there are shortcomings. The Hermite functions/polynomials do not admit an
FFT exploitation and have problems approximating functions that do not decay
exponentially (which is to say, most functions). However, the solutions to differ-
etial equations have been relatively successful by Hermite approximation and in
some cases give superior approximations when compared to a Chebyshev (mapped

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or truncated) approximation [5]. The Whittaker cardinal interpolant functions [29], or Sinc functions, provide a remarkably simple method to approximate a function with known equispaced evaluations. The drawback is a relatively small class of functions for which such an expansion is complete. However, the case of applying Sinc methods has led to a great number of applications [21]. The Chebyshev rational functions [4], [9] are robust with respect to the deficiencies of the Hermite and Sinc bases, but they have some disadvantages compared with the generalized Wiener basis we will derive.

On the semi-infinite interval Laguerre polynomial/function expansions are the classical approximation technique [1], but these techniques suffer from the same problems as Hermite expansions. An alternative technique involves mapping Jacobi polynomials to the infinite interval [7]. This mapping technique makes it possible to accurately approximate algebraically-decaying functions on the semi-infinite interval, but introduces some computational issues for the solution to differential equations. The generalized Wiener basis can be employed on the semi-infinite interval; this results in a basis set that is also a mapped Jacobi polynomial method. However, the Wiener mapping is very different from that presented in the literature, and therefore acts as a competitor to these existing techniques.

Our generalized basis is inspired by a collection of orthogonal and complete functions originally proposed by Wiener [30]. He introduces the functions

\[ \phi_n(x) = \frac{(1 - i x)^n}{\sqrt{\pi (1 + i x)^{n+1}}}, \quad n \in \mathbb{N}_0 \]

as Fourier transforms of the Laguerre functions. He furthermore shows that these functions are orthogonal under the \( L^2 \) conjugate inner product. Higgins [19] expands this result by presenting the functions \( \phi_n \) along with their complex conjugates as a complete system in \( L^2 \). Following this, others have followed up on these functions by applying them to the solution of differential equations [13], [11]. We note that the functions \( \phi_n(x) \) presented above have magnitude that decays like \( \frac{1}{x^s} \) as \( |x| \to \infty \). We will generalize the above functions so that they have decay \( \frac{1}{x^s} \) for any \( s > \frac{1}{2} \). The ability to choose the rate of decay of the basis set is an advantage if such information is present about the nature of the function to be approximated or the differential equation to be solved (e.g. [20], [22]). Furthermore, we will show that this basis admits sparse Galerkin matrices and that the fast Fourier Transform can be used in certain cases to evaluate and manipulate the series.

This paper is concerned with the derivation and theoretical properties of the generalized Wiener rational function basis. Computational considerations, numerical examples, and comparisons with existing basis sets are presented in a second part. The outline of this paper is as follows. In Section 2 we formulate and derive the basis, which is heavily based upon a generalized Fourier series. Section 3 follows with some properties of the basis functions based on their close relationship to the canonical Fourier basis, and Section 4 concerns the properties that can be derived from the relation to Jacobi polynomials. In Section 5 we discuss how the Wiener basis set may be used to approximate functions on the semi-infinite interval. Finally, we briefly present mapped Jacobi polynomials as an alternative method in Section 6 and summarize and present an outlook in Section 7 for Part II, dealing with numerical issues.
2. Derivation of the basis

We begin by stating the major goals and the path we will take in accomplishing those goals. We seek a collection of $L^2$-orthogonal and complete basis functions whose domain is the entire real line. In addition, we desire the ability to specify a parameter $s > \frac{1}{2}$ that will denote the polynomial decay at $\pm \infty$ of each of the the basis functions.

Drawing inspiration from Wiener and his orthogonal basis functions, we seek a collection of functions $\phi_k^{(s)}(x)$ for $x \in \mathbb{R}$ and $k \in \mathbb{Z}$ such that $\{\phi_k^{(s)}\}_{k \in \mathbb{Z}}$ is a complete, orthogonal system for any valid $s$. Our method relies on the observation that the functions (1.1) are weighted maps of the canonical Fourier basis $e^{i\theta}$ for $\theta \in [0, 2\pi]$ (see e.g. [9], [28]). We will first generalize the Fourier basis on $[0, 2\pi]$ so that it will have the properties we desire on the infinite interval; we will then map the generalized Fourier basis to the real line and weight it accordingly to achieve the desired rate of decay.

2.1. Notation and setup. We shall reserve the variables $x, z, \theta, \text{and} r$ as independent variables on certain domains and list the domains and transformations in Table 1. The variable $r \in [-1, 1]$ is the standard interval over which the Jacobi polynomials are defined. The interval $\theta \in [0, \pi]$ is the image of the $r$ interval under the map $\theta = \arccos r$. The variable $z \in \mathbb{T}^+$ is the upper-half of the unit circle in the complex plane, and $x \in [0, \infty]$ is the positive half of the extended real line.

In much of what follows we will mix notation and write expressions both in terms of e.g. $r$ and $\theta$. It should then be understood that $r = r(\theta)$ and/or $\theta = \theta(r)$. Furthermore, we will extend the domains of $\theta, z, \text{and} x$ to be $[-\pi, \pi], \mathbb{T}, \text{and} \mathbb{R}$, respectively later in the paper.

We denote $L^2(A, B; w) = L^2_w(A, B)$ the space of square integrable functions $f : A \to B$ under the weight $w$. We endow $L^2_w(A, B)$ with the conjugate bilinear inner product; the notation for this inner product is $\langle \cdot, \cdot \rangle_w$. The omission of $w$ indicates the unit weight measure. The norm on this space will be denoted $\|\cdot\|_w$. The following weight functions will be used extensively in this article:

<table>
<thead>
<tr>
<th>$x$</th>
<th>$z$</th>
<th>$\theta$</th>
<th>$r$</th>
</tr>
</thead>
<tbody>
<tr>
<td>$x \in [0, \infty]$</td>
<td>$z = -\frac{x-i}{x+i}$</td>
<td>$\theta = 2\arctan(x)$</td>
<td>$r = \frac{1-x^2}{1+x^2}$</td>
</tr>
<tr>
<td>$z \in \mathbb{T}^+$</td>
<td>$\theta = \text{arg } z$</td>
<td>$r = \frac{1}{2}(z + \bar{z})$</td>
<td></td>
</tr>
<tr>
<td>$\theta = \tan \left(\frac{\theta}{2}\right)$</td>
<td>$z = e^{i\theta}$</td>
<td>$\theta \in [0, \pi]$</td>
<td>$r = \cos \theta$</td>
</tr>
<tr>
<td>$r \in [-1, 1]$</td>
<td>$z = e^{i\arccos r}$</td>
<td>$\theta = \arccos r$</td>
<td>$r \in [-1, 1]$</td>
</tr>
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Table 1. Isomorphic transforms between different domains.
\[ w_{\alpha}(r) = (1-r)^{\alpha} \]

\[ w_{\beta}(\theta) = w_{\alpha}(r(\theta)) = (1 + \cos \theta)^{\beta} (1 - \cos \theta)^{\delta} \]

\[ w_{\gamma}(x) = w_{\delta}(\theta(x)) = (1 + \cos \theta)^{\gamma} (1 - \cos \theta)^{\delta} \]

In addition, we will make use of a phase-shifted square root of \( w_{s,t}(x) \) and \( w_{\gamma,\delta}(\theta) \), which we define as:

\[ \sqrt[\gamma]{w_{s,t}(x)} := \sqrt{w_{s,t}(x)} \exp \left[ \frac{i(s + t)}{2} (\pi - \theta(x)) \right] = 2^{s + t} x^{t} \]

\[ \sqrt[\gamma]{w_{\gamma,\delta}(\theta)} = \sqrt[w_{\gamma,\delta}(x(\theta))]{\gamma} \frac{\sin \left( \frac{\theta}{2} \right)}{\cos \left( \frac{\theta}{2} \right)} \exp \left[ \frac{i(\gamma + \delta)}{2} (\pi - \theta) \right] \]

2.2. Jacobi polynomials. The classical Jacobi polynomials \( P_{n}^{(\alpha,\beta)} \) are a family of orthogonal polynomials [27] that have been used extensively in many applications due to their ability to approximate general classes of functions. They are a class of polynomials that encompass the Chebyshev, Legendre, and Gegenbauer/ultraspheric polynomials. These polynomials will form the building blocks for our generalization.

The Jacobi differential equation is

\[ (1 - r^2) p'' + [\beta - \alpha - (\alpha + \beta + 2)r] p' + n(n + \alpha + \beta + 1) p = 0, \quad r \in [-1, 1], \]

and for \( \alpha, \beta > -1, n \in \mathbb{N}_0 \) the only polynomial solution \( p = P_{n}^{(\alpha,\beta)}(x) \) is a polynomial of degree \( n \). The restriction \( \alpha, \beta > -1 \) is necessary to ensure integrability of the weight and thus existence of an \( L^2 \)-constant function solution. The family of polynomials \( \{ P_{n}^{(\alpha,\beta)}(x) \}_{n=0}^{\infty} \) is complete and orthogonal in \( L^2([-1, 1], \mathbb{R}; w_{r}^{(\alpha,\beta)}) \).

We denote \( h_{n}^{(\alpha,\beta)} = \| P_{n}^{(\alpha,\beta)} \|_{w_{r}^{(\alpha,\beta)}}^2 \), and define the normalized polynomials as

\[ \tilde{P}_{n}^{(\alpha,\beta)}(r) = \frac{P_{n}^{(\alpha,\beta)}(r)}{\sqrt{h_{n}^{(\alpha,\beta)}}}. \]

The orthonormal Jacobi polynomials \( \tilde{P}_{n}^{(\alpha,\beta)} \) will be integral in the derivation of the Wiener rational function basis on the real line. In addition, we require a minor generalization of Jacobi polynomials: we perform a change of the dependent variable in (2.3) to obtain:

**Lemma 2.1.** (Jacobi Functions) The Jacobi functions defined as

\[ P_{n}^{(\alpha,\beta,a,b)}(r) = (1 - r)^{a} (1 + r)^{b} P_{n}^{(\alpha,\beta)}(r) \]

satisfy the following properties:
(1) \( \{ P_n^{(\alpha, \beta, a, b)}(r) \}_{n \in \mathbb{N}_0} \) are orthogonal and complete in \( L^2([-1, 1], \mathbb{R}; w_r^{(\alpha-2a, \beta-2b)}) \).

(2) The \( P_n^{(\alpha, \beta, a, b)}(r) \) are eigenfunctions \( \rho_n(r) \) of the Sturm-Liouville problem

\[
\frac{d}{dr} [p(r) \rho'(r)] + q(r) \rho(r) - \lambda_n w(r) \rho(r) = 0,
\]

which is defined by the parameters

\[
p(r) = (1 - r)^{\alpha+1-2a}(1 + r)^{\beta+1-2b},
\]

\[
q(r) = \left[ a(\alpha - a)(1 - r)^{-2} + b(\beta - b)(1 + r)^{-2} \right] \times \frac{1}{(1 - r)^{\alpha+1-2a}(1 + r)^{\beta+1-2b}},
\]

\[
w(r) = (1 - r)^{\alpha-2a}(1 + r)^{\beta-2b},
\]

\[
\lambda_n = n(n + \alpha + \beta + 1) - 2ab + a(\beta + 1) + b(\alpha + 1)
\]

The proof is mathematically simple but algebraically tedious and we omit it. We shall actually only require the result of Lemma 2.1 for \( a = b = \frac{1}{2} \). Many of the results in this paper require the use of numerous recurrence relations involving Jacobi polynomials; these relations are given in Appendix A, equations (A.1)-(A.8).

The idea behind the formation of the Jacobi functions introduced in Lemma 2.1 is not novel and has already found use in the literature. In [17] the 'generalized Jacobi polynomials/functions' are denoted \( j_n^{(\alpha, \beta)} \), and are defined for all \( \alpha, \beta \in \mathbb{R} \) as

\[
j_n^{(\alpha, \beta)} \propto \begin{cases} 
P_{n_1}^{(-\alpha,-\beta,-\alpha,-\beta)}, & \alpha \leq -1 \text{ and } \beta \leq -1, \\
P_{n_1}^{(-\alpha,\beta,-\alpha,0)}, & \alpha \leq -1 \text{ and } \beta > -1, \\
P_{n_1}^{(\alpha,\beta,0,-\beta)}, & \alpha > -1 \text{ and } \beta \leq -1, \\
P_{n_1}^{(\alpha,\beta,0,0)}, & \text{else,}
\end{cases}
\]

where the index \( n_1 \) is defined as

\[
n_1 = \begin{cases} 
n - \lfloor -\alpha \rfloor - \lfloor -\beta \rfloor, & \alpha \leq -1 \text{ and } \beta \leq -1, \\
n - \lfloor -\alpha \rfloor, & \alpha \leq -1 \text{ and } \beta > -1, \\
n - \lfloor -\beta \rfloor, & \alpha > -1 \text{ and } \beta \leq -1, \\
n, & \text{else,}
\end{cases}
\]

and the integer floor function is denoted \( \lfloor \cdot \rfloor \). These functions are only defined for certain values of \( n \) but [17] presents significant approximation theory using them. They are advantageous for solving high-order differential equations with boundary conditions via a global spectral expansion.

Finally, we present two classical notational conventions that we will use briefly in the next section. The classical Jacobi polynomials that result from the cases \( \alpha = \beta = -\frac{1}{2} \) and \( \alpha = \beta = +\frac{1}{2} \) are the Chebyshev polynomials of the first and second kinds, respectively. Recalling the relation \( r = \cos \theta \), these polynomials are typically denoted \( T_n(r) \) and \( U_n(r) \) and they have a very special and concise
representation as trigonometric polynomials:
\[ \sqrt{\frac{2}{\pi}} \tilde{P}_{n}(-1/2,-1/2)(r) = T_{n}(r) = \cos{(n\theta)} = \cos{[n \arccos{(r)}]} \]
\[ \sqrt{\frac{2}{\pi}} \tilde{P}_{n}^{(1/2,1/2)}(r) = U_{n}(r) = \frac{\sin{[(n+1)\theta]}}{\sin{\theta}} = \frac{\sin{[(n+1) \arccos{(r)}]}}{\sin{[\arccos{(r)}]}}. \]

2.3. Generalizing the Fourier basis. In this section we will generalize the canonical Fourier basis given by
\[ \Psi_{k}(\theta) = e^{ik\theta}. \]

Our methodology is based upon the following dissection of the Fourier basis for \( k \neq 0 \):
\[ e^{ik\theta} = \cos{(k\theta)} + i \sin{(k\theta)} \]
\[ = \cos{(|k|\theta)} + i \, \text{sgn}(k) \sin{(|k|\theta)} \]
\[ = T_{|k|}(\cos{\theta}) + i \, \text{sgn}(k) \sin{\theta} U_{|k|-1}(\cos{\theta}) \]
\[ = \sqrt{\frac{2}{\pi}} \left[ \tilde{P}_{n}^{(-1/2,-1/2)}(\cos{\theta}) + i \, \text{sgn}(k) \sin{\theta} \tilde{P}_{n}^{(1/2,1/2)}(|k|-1)(\cos{\theta}) \right]. \]

We have broken down the Fourier basis into two components: the first component \((a)\) is even with respect to \( \theta \) as it is simply a polynomial in \( \cos{\theta} \). The second term \((b)\) is odd in \( \theta \) as it is a polynomial in \( \cos{\theta} \) (an even function) multiplied by the odd function \( \sin{\theta} \). This breakdown suggests that we can construct more general kinds of Fourier-type functions by augmenting the type of polynomials employed.

However, we cannot switch around polynomials with impunity; we want to retain orthogonality (at least with respect to some weight function). The separation into terms \((a)\) and \((b)\) above elucidates the biorthogonal decomposition of the Fourier basis. The \((a)\) functions are orthogonal with respect to each other, and with respect to the \((b)\) functions. In this case, the biorthogonality is manifested as an even-odd separation. Suppose we wish to generate a basis set orthogonal under the weight \( 1 + \cos{\theta} = 1 + r \). Naturally we can do this for basis \((a)\) by changing the second Jacobi class parameter from \( \beta = -\frac{1}{2} \) to \( \beta = \frac{1}{2} \). In order to do this for basis \((b)\), we use Lemma 2.1.

For \( \alpha, \beta > -1 \), we have the polynomials \( \tilde{P}_{n}^{(\alpha,\beta)} \) that are orthogonal in \( L^{2}\left([-1,1], \mathbb{R}; w_{r}^{(\alpha,\beta)}\right) \). By setting \( a = b = \frac{1}{2} \) in Lemma 2.1, we also observe that the Jacobi functions \( \tilde{P}_{n}^{(\alpha+1,\beta+1,1/2,1/2)}(1-r^{2})^{1/2} \tilde{P}_{n}^{(\alpha+1,\beta+1)} \) are orthogonal under the same weight. If we set \( \alpha = \beta = -\frac{1}{2} \), and add these two functions together with the appropriate scaling factors, then we exactly recover the Fourier basis by reversing the dissection steps above (i.e. by creating a biorthogonal construction). Of course, we are free to choose any values of \( (\alpha, \beta) \) that we desire in order to derive generalized trigonometric Fourier functions. In fact, this technique has already been used by Szegő [27] to determine orthogonal polynomials on the unit disk. Because the statement in [27] is merely a passing comment and is a markedly different result than what we desire, we present the following theorem:
Theorem 2.2. (cf. Szegő, [27]) For any \( \gamma > -\frac{1}{2} \), the functions
\( \Psi_k^{(\gamma)}(\theta) = \begin{cases} \frac{1}{\sqrt{2}} \tilde{P}_k^{(-1/2,\gamma-1/2)}(\cos \theta), & k = 0 \\ \frac{1}{2} \left[ \tilde{P}_k^{(-1/2,\gamma-1/2)}(\cos \theta) + i \text{sgn}(k) \sin(\theta) \tilde{P}_{|k|-1}^{(1/2,\gamma+1/2)}(\cos \theta) \right], & k \neq 0 \end{cases} \)
are complete and orthonormal in \( L^2\left([-\pi, \pi], \mathbb{C}; w_0^{(\gamma,0)}\right) \).

Proof. For orthonormality, it suffices to show
1. \( \left\langle \tilde{P}_k^{(-1/2,\gamma-1/2)}(\cos \theta), \tilde{P}_{|l|-1}^{(-1/2,\gamma-1/2)}(\cos \theta) \right\rangle_{w_0^{(\gamma,0)}} = 2\delta_{|k|,|l|} \)
2. \( \left\langle \sin \theta \tilde{P}_k^{(1/2,\gamma+1/2)}(\cos \theta), \sin \theta \tilde{P}_{|l|-1}^{(1/2,\gamma+1/2)}(\cos \theta) \right\rangle_{w_0^{(\gamma,0)}} = 2\delta_{|k|,|l|}, \) for \( k, l \neq 0 \).
3. \( \left\langle \tilde{P}_k^{(-1/2,\gamma-1/2)}(\cos \theta), \tilde{P}_{|l|-1}^{(1/2,\gamma+1/2)}(\cos \theta) \right\rangle_{w_0^{(\gamma,0)}} = 0, \) for \( l \neq 0 \).

The first property is a direct result of orthonormality of the normalized Jacobi polynomials \( \tilde{P} \) and the observation that on \([0, \pi]\), \( \left\langle f(\cos \theta), g(\cos \theta) \right\rangle_{w_0^{(\gamma,0)}} = \left\langle f(r), g(r) \right\rangle_{w_0^{(-1/2,\gamma-1/2)}} \). The second property is a result of the same observations as the first property along with the result of Lemma 2.1. The third property results from the fact that an odd function integrated over a symmetric interval is 0. Orthogonality then follows from an explicit calculation of \( \left\langle \Psi_k^{(\gamma)}, \Psi_l^{(\gamma)} \right\rangle_{w_0^{(\gamma,0)}} \) using the above three properties.

For completeness we note that any function \( f \in L^2 \) can be decomposed into an even \( f_e \) and an odd \( f_o \) part. That \( f_e \) is representable is clear from the fact that \( \tilde{P}_n^{(-1/2,\gamma-1/2)}(\cos \theta) \) is complete over \( \theta \in [0, \pi] \), which by symmetry implies completeness over all \( L^2 \)-even functions \( f_e \). Similarly, the collection of functions \( \sin \theta \tilde{P}_n^{(1/2,\gamma+1/2)}(\cos \theta) \) is complete over all \( L^2 \)-odd functions \( f_o \) by Lemma 2.1. Linearity and orthogonality of the even and odd parts yields the result. \(\square\)

Remark 2.3. Szegő [27] gives a more general result that involves orthogonality over the weight \( w_0^{(\gamma,\delta)} \) for \( \delta \neq 0 \). We do not require this level of generality; for \( \delta \neq 0 \) the weight function becomes zero at \( \theta = 0 \), which we will see does not help our cause. Indeed, it is possible to generalize Szegő’s result: he derived polynomials on the unit disk orthogonal with respect to \( w_0^{(\gamma,\delta)} \). By using Lemma 2.1 with \( a, b \) different from \( \frac{1}{2} \), we can derive non-polynomial basis sets that are orthogonal under a great variety of weights. These functions naturally may not be periodic on \( \theta \in [-\pi, \pi] \) if the quantity \( (1-r)^a(1+r)^b \) cannot be periodically extended in \( \theta \)-space to \( [-\pi, \pi] \).

We will refer to the functions (2.4) as either the generalized Fourier series, or the Szegő-Fourier functions. In the definition of the functions \( \Psi_k^{(\gamma)} \) it is desirable to use the \( L^2 \)-normalized versions of the Jacobi polynomials \( \tilde{P} \), rather than the standard polynomials \( P \). If the standard polynomials are used, then the norm of the Szegő-Fourier functions \( \Psi_k^{(\gamma)} \) depends on the rather unpleasant-looking sum \( h_{|k|}^{(-1/2,\gamma-1/2)} + h_{|k|-1}^{(1/2,\gamma+1/2)} \), and using this convention implies that \( \Psi_k^{(\gamma)} \) is not orthogonal to \( \Psi_{-k}^{(\gamma)} \).

We can also distribute the weight function onto the basis functions, which gives us orthogonality in the unweighted \( L^2 \)-norm:
Figure 1. Plots of the weighted Szegő-Fourier functions $\psi^{(2)}(\theta)$ for $k = 0, 1, 2, 3, \text{and} 4$. Real part (top) and imaginary part (bottom).

Corollary 2.1. For any $\gamma > -\frac{1}{2}$, the functions

$$
\psi^{(\gamma)}(\theta) = \begin{cases} \\
\frac{\sqrt{w^{(\gamma,0)}_\theta}}{\sqrt{2}} P_0(-1/2,\gamma-1/2)(\cos \theta), & k = 0 \\
\frac{\sqrt{w^{(\gamma,0)}_\theta}}{2} \left[ \tilde{P}_{|k|}^{(-1/2,\gamma-1/2)}(\cos \theta) + i \text{sgn}(k) \sin(\theta) \tilde{P}_{|k|-1}^{(1/2,\gamma+1/2)}(\cos \theta) \right], & k \neq 0 
\end{cases}
$$

are complete and orthonormal in $L^2([-\pi, \pi], C)$.

Due to the properties of $\sqrt{w^{(\gamma,0)}_\theta}$ given in (2.2), the functions $\psi^{(\gamma)}(\theta)$ decay like $(\cos \frac{\theta}{2})^\gamma$ at $\theta = \pm \pi$. This is exemplified in Figure 1 where we plot the real and imaginary parts of the functions for $\gamma = 2$. The even/odd behavior in $\theta$ for real/imaginary components depicted in the figure depends on the even/odd parity of $\gamma$. (There is no such characterization possible when $\gamma \notin \mathbb{N}_0$.) Clearly for $\gamma = 0$ we have $\Psi^{(0)}_k = \psi^{(0)}_k = \frac{1}{\sqrt{2\pi}} e^{ik\theta}$, the canonical Fourier basis.

2.4. Mapping to the real line. Having developed the necessary preliminaries on the finite interval, we now jump to the infinite line $x \in \mathbb{R}$ using the mappings introduced in Table 1. To facilitate the mapping, the following identities characterizing the mapping between $\theta$-space and $x$-space are useful:

$$
\cos \theta = \frac{1 - x^2}{1 + x^2}, \quad (1 - \cos \theta) = \frac{2x^2}{x^2 + 1}, \\
\sin \theta = \frac{2x}{x^2 + 1}, \quad (1 + \cos \theta) = \frac{2}{x^2 + 1}.
$$
Using these identities, we rewrite and relabel the functions $\Psi_k^{(s)}(\theta)$:

$$
\Phi_k^{(s)}(x) := \Psi_k^{(s-1)}(\theta(x))
$$

$$
= \begin{cases}
\frac{1}{\sqrt{2}} P_0^{(-1/2,s-3/2)} \left( \frac{1-x^2}{1+x^2} \right), & k = 0 \\
\frac{1}{2} \left[ P_k^{(-1/2,s-3/2)} \left( \frac{1-x^2}{1+x^2} \right) + 2i x \text{sgn}(k) P_{|k|-1}^{(1/2,s-1/2)} \left( \frac{1-x^2}{1+x^2} \right) \right], & k \neq 0
\end{cases}
$$

The above definition is valid for any $s > \frac{1}{2}$. $s = 1$ corresponds to a mapping of the canonical Fourier basis (i.e., $s = \gamma + 1$). These functions are orthogonal over the weight $w_x^{(s,0)}$. By following the route from Corollary 2.1 we can distribute the weight over the basis functions, and in this particular instance we choose the phase-shifted square root given in (2.1):

(2.5)

$$
\phi_k^{(s)} := \sqrt{w_x^{(s,0)}} \Phi_k^{(s)}(x)
$$

$$
= \begin{cases}
\frac{2(\frac{s}{2}-1)}{(s-1)} P_0^{(-1/2,s-3/2)} \left( \frac{1-x^2}{1+x^2} \right), & k = 0 \\
\frac{2(\frac{s}{2}-1)}{(s-1)} \left[ P_k^{(-1/2,s-3/2)} \left( \frac{1-x^2}{1+x^2} \right) + 2i x \text{sgn}(k) P_{|k|-1}^{(1/2,s-1/2)} \left( \frac{1-x^2}{1+x^2} \right) \right], & k \neq 0
\end{cases}
$$

The functions (2.5) are what we call the generalized Wiener rational functions. At present there is no clear reason why we have chosen to use $\sqrt{w_x^{(s,0)}}$ instead of the usual square root $\sqrt{w_x^{(s,0)}}$ to distribute the weight. However, the corollary following the coming proposition should provide part of the motivation.

**Proposition 2.1.** For any $s > \frac{1}{2}$, the functions $\Phi_k^{(s)}(x)$ are complete and orthonormal in $L^2 \left( \mathbb{R}, \mathbb{C}; w_x^{(s,0)} \right)$. The functions $\phi_k^{(s)}(x)$ are complete and orthonormal in $L^2 \left( \mathbb{R}, \mathbb{C} \right)$. Furthermore, the decay rate of these functions can be characterized as

$$
\lim_{|x| \to \infty} \left| x^t \phi_k^{(s)}(x) \right| < \infty, \quad t \leq s
$$

**Corollary 2.2.** Recalling the definition of Wiener’s original basis functions $\phi_n(x)$ in (1.1), the following relation holds:

$$
i \sqrt{2} \phi_n^{(1)}(x) = \phi_n(x), \quad n \in \mathbb{N}_0.
$$

We show plots of the functions $\phi_k^{(4)}$ in Figure 2. The conclusion of the corollary is easily seen if one makes the connection

$$
e^{it\theta} = \frac{i - x}{i + x},
$$

along with knowledge of the fact that $\Phi_k^{(1)}(x) = \psi_k^{(0)}(\theta) = \frac{1}{\sqrt{2\pi}} e^{ik\theta}$. We have thus shown that the orthogonal functions $\phi_k^{(s)}$ over the real line are a generalization of Wiener’s original basis set. Furthermore, $\phi_k^{(s)}$ decays like $x^{-s}$ while retaining
orthogonality under the same unit weight measure. When $s$ is an integer, the
functions are also purely rational: they are the division of one complex-valued
polynomial in $x$ by another. This connection was rather helpful in the nascent
stages of the computing when the calculation of a non-polynomial function required
significantly more computational investment, but now this property is probably
more aesthetic than functional. As a result, our use of the quantity $\sqrt{w_x(s,0)}$ is not
entirely necessary for purposes of evaluating the functions; it is equally valid to use
the traditional square root $\sqrt{w_x(s,0)}$.

By using the traditional square root, one sacrifice made is that the analogous
written form of Corollary 2.2 becomes less fortuitous and is complicated by $x$-
dependent phase-shift factors. The same observation is true of the weight $\sqrt{w_x^{(γ,0)}}$
used in the definition of the Szegö-Fourier functions $ψ_k^{(γ)}(θ)$ in Corollary 2.1. A
second reason to use the phase-shifted square root is that it can be written in the
following convenient form:

\begin{equation}
\sqrt{w_x^{(s,0)}} = \left[ \frac{i}{\sqrt{2}} (1 + e^{-iθ}) \right]^s.
\end{equation}

The utility of this expression will become clear when we consider the connection
problems in Section 4.

We have accomplished our goal of deriving basis functions satisfying tunable
decay rate while maintaining $L^2$-orthogonality. However, it is not clear that these
are superior or useful functions. We will now present some properties of the basis
and make the argument that these basis functions indeed are very useful for solving
problems in scientific computing.
3. Fourier-Derived Basis Properties

In this section we explore some of the desirable properties of the generalized Wiener basis set \( \{ \phi_k^{(s)} \}_{k \in \mathbb{Z}, s > \frac{1}{2}} \) based on their close relation to Fourier Series. The argument we make is that these functions inherit all the useful properties of the Fourier basis with the additional property that the decay rate \( s \) at \( |x| = \infty \) may be chosen. Many of these properties (e.g. the sparse modal differentiation matrix) rely on Jacobi polynomial properties covered in the next section. In particular, although application of the FFT is indeed a virtue of this basis, we will discuss it only in Part II, which focuses with computational issues.

3.1. Symmetry. The derivation of the basis functions automatically yields various simple properties. Note that due to the mapping, any property of the basis on the real line \( x \in \mathbb{R} \) also applies to the respective trigonometric interval \( \theta \in [-\pi, \pi] \). We omit the proof of these properties as they are elementary:

1. Index symmetry

\[
\Phi_k^{(s)}(x) = \Phi_{-k}^{(s)}(x)
\]

\[
|\Phi_k^{(s)}(x)| = |\Phi_{-k}^{(s)}(x)|
\]

\[
|\phi_k^{(s)}(x)| = |\phi_{-k}^{(s)}(x)|
\]

\[
\phi_k^{(1)}(x) = \phi_{-k-1}^{(1)}(x)
\]

2. Function symmetry

\[
\text{Re}\left\{ \Phi_k^{(s)}(x) \right\} = \text{Re}\left\{ \Phi_k^{(s)}(-x) \right\}
\]

\[
\text{Im}\left\{ \Phi_k^{(s)}(x) \right\} = -\text{Im}\left\{ \Phi_k^{(s)}(-x) \right\}
\]

\[
|\Phi_k^{(s)}(x)| = |\Phi_k^{(s)}(-x)|
\]

\[
|\phi_k^{(s)}(x)| = |\phi_k^{(s)}(-x)|
\]

3.2. Periodicity. Because trigonometric polynomials are periodic over \( \theta \in [-\pi, \pi] \), we cannot expect this condition to be violated on the infinite interval \( x \in \mathbb{R} \). From the viewpoint of expanding functions over the infinite interval \( \mathbb{R} \), the points \( x = \pm \infty \) are both unique points. However, because of the mapping, the basis functions view the points \( x = \pm \infty \) the same as they view the points \( \theta = \pm \pi \): i.e. they are the same point. This serves as a disadvantage if we wish to e.g. expand functions with different decay rates at \( \pm \infty \) because this is in effect non-smooth behavior of the function at a single point, which degrades the convergence rate of the approximation.

In particular it is known that although a Fourier series approximation will converge in the \( L^2 \) sense for an \( L^2 \) function, the rate of convergence is only algebraic if the function is non-periodic. Naturally, this deficiency will follow us to the infinite interval. Indeed, such observations have already been made [12]. Empirical studies
we have carried out show that the concern of periodicity is not paramount and frequently one can overlook it when comparing results to other basis expansions. Nonperiodic behavior is often manifested as algebraic decay at $x = \pm \infty$, where existing basis sets already have problems in approximation. In Part II we will present examples that explicitly illustrate this lack of fast convergence rate when the function to be expanded is not ‘periodic’ at $x = \pm \infty$.

![Figure 3. The linear fractional mapping that relates $x$ to $\theta$ has an illuminating representation when viewed as a transformation of the complex plane.](image)

Note that although it may seem a bit unnatural that periodicity is a condition at $x = \pm \infty$, in fact it is not surprising at all. One may consider our mapping as a rather unremarkable tangent map from $\theta$-space to $x$-space as written in Table 1. However, it is more deep than that: the functions $\Psi_k^{(\gamma)}(\theta)$ and $\psi_k^{(\gamma)}(\theta)$ are periodic basis sets for complex-valued functions on the unit circle $\mathbb{T}$. In other words, we may actually view these basis sets as functions of $z \in \mathbb{C}$. What looks like a tangent
mapping from $\theta$-space to $x$-space is actually a linear fractional map (a Möbius transformation) from the unit circle (the complex plane) to complexified $x$-space (the complex plane).

Linear fractional maps are structure-preserving maps of the complex plane; an illuminating way to consider this is by identifying the complex plane $\mathbb{C}$ with the Riemann Sphere (see Figure 3.2). Then the linear fractional map we’ve chosen merely takes one great circle (the unit circle in $z$-space) to another great circle (the real line in complexified $x$-space). Therefore, our approximation is nothing more than a rotation of functions on the Riemann Sphere (see Figure 3.2); the target space simply happens to correspond to the real line. From this point of view, periodicity at $|x| = \infty$ (analyticity at complexified $x = \infty$) is natural.

Nevertheless, this ‘natural’ periodicity can be problematic if we attempt to approximate a function that is not complex-analytic at $x = \infty$. In Part II we present examples of functions that are not analytic at $x = \infty$ and we will empirically analyze the impact of violating the assumption of periodicity.

4. Jacobi-Derived Basis Properties

The generalized Wiener functions are composed of Jacobi polynomials, and so it is reasonable to expect that we can use the properties of the Jacobi polynomials to perform certain tasks using the Wiener basis. Indeed, we can form recurrence relations, connection coefficients, a Gauss-like quadrature, and obtain an extremely useful sparsity result for the Galerkin stiffness matrix.

4.1. Recurrence Relations. Due to the strong dependence of the Szegő-Fourier functions on the Jacobi polynomials, they inherit six-term recurrence relations from the three-term recurrences for orthogonal polynomials.

\[
D_n^{(\gamma)} \Psi_{n+1}^{(\gamma)} = \left[A_n^{(\gamma)} e^{i\theta} - B_n^{(\gamma)} \right] \Psi_n^{(\gamma)} + \left[A_n^{(\gamma)} e^{-i\theta} - B_n^{(\gamma)} \right] \Psi_{-n}^{(\gamma)} +
C_n^{(\gamma)} \Psi_{n+1}^{(\gamma)} + C_n^{(\gamma)} \Psi_{n-1}^{(\gamma)},
\]

\[
\Psi_{n+1}^{(\gamma)} = \left[U_n^{(\gamma)} \cos \theta - V_n^{(\gamma)} \right] \Psi_n^{(\gamma)} + \left[U_n^{(\gamma)} \cos \theta - V_n^{(\gamma)} \right] \Psi_{-n}^{(\gamma)} +
W_n^{(\gamma)} \Psi_{n+1}^{(\gamma)} + W_n^{(\gamma)} \Psi_{n-1}^{(\gamma)},
\]

\[
\Psi_{n+1}^{(\gamma)} = \left[\tilde{U}_n^{(\gamma)} i \sin \theta - \tilde{V}_n^{(\gamma)} \right] \Psi_n^{(\gamma)} + \left[\tilde{U}_n^{(\gamma)} i \sin \theta - \tilde{V}_n^{(\gamma)} \right] \Psi_{-n}^{(\gamma)} +
\tilde{W}_n^{(\gamma)} \Psi_{n+1}^{(\gamma)} + \tilde{W}_n^{(\gamma)} \Psi_{n-1}^{(\gamma)},
\]

We give formulae for all the real-valued constants $A, B, C, D, U, V, W, \tilde{U}, \tilde{V}, \tilde{W}$ in Appendix A. Note that since the $\Psi_k^{(\gamma)}$ are not polynomials in $z = e^{i\theta}$, there is not a three-term recurrence as there would normally be for orthogonal polynomials on the unit disk (unless of course $\gamma = 0$). Although the above formulae are complex-valued six-term recurrence relations, they are no more difficult computationally than the pair of three-term recurrences necessary to generate $P_n^{(\alpha,\beta)}$ and $\tilde{P}_n^{(\alpha+1,\beta+1)}$ because $\Psi_n^{(\gamma)}$ is the complex conjugate of $\Psi_n^{(\gamma)}$ and therefore does not need to be generated independently. Direct use of any of the above six-term recurrences for generating the $\Psi_k^{(\gamma)}$ is just as expensive as forming $\Psi_k^{(\gamma)}$ by the even/odd synthesis of $\tilde{P}_n^{(\alpha,\beta)}$. 
and $\tilde{P}_n^{(\alpha+1,\beta+1)}$ in Theorem 2.2. However, using presumably existing routines for evaluating Jacobi polynomials and then synthesizing them is likely easier from an implementation view.

It is reassuring to note that simplifying the recurrence constants in the case $\gamma = 0$ yields, up to normalization, the trivial recurrence relations for the monomials on the unit disk $\Psi_{\gamma}(0) = z^k$:

$$
\begin{align*}
\Psi_{\gamma}(0)_{n+1} & = e^{i\theta}\Psi_{\gamma}(0)_{n}, \\
\Psi_{\gamma}(0)_{n+1} & = 2\cos \theta \Psi_{\gamma}(0)_{n} - \Psi_{\gamma}(0)_{n-1}, \\
\Psi_{\gamma}(0)_{n+1} & = 2i\sin \theta \Psi_{\gamma}(0)_{n} + \Psi_{\gamma}(0)_{n-1}.
\end{align*}
$$

Naturally, a recurrence relation for the unweighted $\Psi_{\gamma}(\theta)$ translates directly into one for the unweighted Wiener rational functions $\Phi_{\gamma}(x)$. The weighted functions $\psi_{\gamma}(\theta)$ and $\phi_{\gamma}(x)$ can be generated by first generating the unweighted functions and then multiplying by the phase-shifted square root $\sqrt{w}$.

4.2. Connection Problems. One advantage in using the generalized Wiener rational function basis is the ability to choose the parameter $s$, which indicates the rate of decay. In many applications, it may be useful to augment the basis functions mid-computation to suit the dynamics occurring at a particular time. In this case, one would like to be able to transfer from one basis to another while keeping the (finite-term) function expansion identical. We will also see in Part II that this problem also appears in an algorithm utilizing the FFT. In classical orthogonal polynomial theory, the problem of equating one expansion to another boils down to determining the connection coefficients. Before undertaking this task, we first outline the major tasks we wish to perform.

There are two main tasks on the infinite interval that require connections of some form:

1. Usage of the fast Fourier transform – transforming $N$ nodal evaluations into $N$ modal coefficients (or vice-versa) for an expansion in $\phi(s)$.
2. For a given expansion in $\phi(s)$ (i.e. a set of modal coefficients), translating this into a modal coefficient expansion in $\phi(S)$ for some $s \neq S$.

In Part II where we outline computational considerations, we will address the above tasks. However, for now it suffices to note that these two tasks can be reduced to the following three connection problems in $\theta$-space:

1. The $\Psi(\gamma)$-$\Phi(\Gamma)$ connection (a necessary ingredient for all connection-like tasks)
2. The $\Psi(\gamma)$-$\psi(\gamma)$ connection (a generalization of the FFT task)
3. The $\psi(\gamma)$-$\phi(\Gamma)$ connection (identical to modification of $s$)

In Sections 4.2.1-4.2.3, we will tackle each of these problems. Note that modification of any of the following finite-interval algorithms for the infinite interval is trivial: the relations $\Psi_{\gamma}(\theta) \equiv \Phi_{k}^{(s-1)}(x)$, $\psi_{\gamma}(\theta) \equiv \phi_{k}^{(s-1)}(x)$, and $\gamma := s - 1$ allows for us to easily employ the same operations, whether we want to do it in $\theta$-space or $x$-space.
4.2.1. The $\Psi^+\Psi$ Connection Problem. Suppose we have a function $f \in L^2\left([−\pi, \pi], \mathbb{C}; w^{(\gamma)}_\theta\right) \cap L^2\left([−\pi, \pi], \mathbb{C}; w^{(\Gamma)}_\theta\right)$ with a Fourier expansion for some $\gamma > -\frac{1}{2}$:

$$f(x) = \sum_{k \in \mathbb{Z}} \hat{f}_k^{(\gamma)} \Psi_k^{(\gamma)}.$$  

The goal is to determine a way to re-expand $f$ in a Fourier expansion for a different decay parameter $\Gamma$:

$$f(x) = \sum_{k \in \mathbb{Z}} \hat{f}_k^{(\Gamma)} \Psi_k^{(\Gamma)}.$$  

The shift $\Gamma - \gamma$ can take values in the interval $(-\frac{1}{2} - \gamma, \infty)$. Naturally, one may equate the two expansions and use orthogonality to relate one set of expansion coefficients to the other:

$$\hat{f}_k^{(\Gamma)} = \sum_{l \in \mathbb{Z}, |l| \geq |k|} \hat{f}_l^{(\gamma)} \lambda^{\Psi}_{k,l}.$$  

We can then define the connection coefficients

$$\lambda^{\Psi}_{k,l} = \left\langle \Psi_k^{(\gamma)}, \Psi_l^{(\Gamma)} \right\rangle_{w^{(\Gamma)}_\theta},$$  

where we have suppressed the dependence of $\lambda$ on $\gamma$ and $\Gamma$. Our task is to determine how to calculate these connection coefficients. Due to orthogonality, it is clear that

$$\lambda^{\Psi}_{k,l} \equiv 0, \quad |l| < |k|.$$  

This implies that the connection problem is solved via the relation

$$\hat{f}_k^{(\Gamma)} = \sum_{l \in \mathbb{Z}, |l| \geq |k|} \hat{f}_l^{(\gamma)} \lambda^{\Psi}_{k,l}.$$  

Relation (4.2) is still not attractive: we must perform an infinite number of operations for an exact connection. If we only have a finite expansion (say a total of $N$ modal coefficients), we must still perform $O(N^2)$ operations to capture all the information at our disposal. However, we will show that, for integer values of the shift $\Gamma - \gamma$, the connection problem can be solved inexpensively. To be precise, we will show that for $G \in \mathbb{N}$, (4.2) reduces to

$$\hat{f}_k^{(\gamma + G)} = \sum_{k \geq |l| \geq |k|} \hat{f}_l^{(\gamma)} \lambda^{\Psi}_{k,l}. $$  

That is, only $2(G + 1)$ operations per coefficient are necessary to solve the connection problem (independent of $k$, and of any truncation size $N$). We refer to the above collapse of the infinite connection problem (4.2) into the finite $N$-independent problem (4.3) as a sparse connection.

In order to relate one Fourier function to another, we first recall a result from [25] using (A.4) – (A.7) that states that the connection coefficients binding one Jacobi polynomial class to another are sparse in certain special circumstances.

**Lemma 4.1.** For any $\alpha, \beta > -1$ and any $A, B, \in \mathbb{N}$, the connection problem

$$f(r) = \sum_{n=0}^{\infty} \tilde{f}_n^{(\alpha,\beta)} P_n^{(\alpha,\beta)}(r) \quad \rightarrow \quad f(r) = \sum_{n=0}^{\infty} \tilde{f}_n^{(\alpha+A,\beta+B)} P_n^{(\alpha+A,\beta+B)}(r),$$  

where
can be solved exactly via the relation
\begin{equation}
\hat{f}_n(\alpha+A,\beta+B) = \sum_{m=0}^{A+B} \lambda_{n,n+m}^{P} \hat{f}_{n+m}^{(\alpha,\beta)}.
\end{equation}

In the above we have suppressed the dependence of \( \lambda^P \) on \( \alpha, \beta, A \), and \( B \), but in the sequel we shall occasionally refer to the above coefficients as \( \lambda_{n,m}^{P} (\alpha, \beta, A, B) \). The result (4.4) is not a trivial one; the upper limit for the sum on the right-hand side is \( \infty \) for a general connection problem. For the very special cases satisfying the lemma, the exact connection becomes finite. We have not shown how to obtain the Jacobi-Jacobi connection coefficients \( \lambda^{P} \). For this, one may use explicit formulae given in [23] or [2], or one may utilize the algorithm given in [25].

The above result can be expanded to apply to the Szegő-Fourier functions \( \Psi^{(\gamma)}_k(\theta) \) and the corresponding mapped functions \( \Phi^{(s)}_k(x) \).

**Proposition 4.1.** For any \( \gamma > -\frac{1}{2} \) and any \( G \in \mathbb{N} \), the connection problem
\begin{equation}
f(\theta) = \sum_{k=-\infty}^{\infty} \hat{f}_k^{(\gamma)} \Psi^{(\gamma)}_k(\theta) \quad \longrightarrow \quad f(\theta) = \sum_{k=-\infty}^{\infty} \hat{f}_k^{(\gamma+G)} \Psi^{(\gamma+G)}_k(\theta),
\end{equation}
can be solved exactly via the relation
\begin{equation}
\hat{f}_k^{(\gamma+G)} = \sum_{|l|\leq|k|} \lambda_{k,l}^{\Psi} \hat{f}_l^{(\gamma)} + \sum_{|l|>|k|\geq-G} \lambda_{k,l}^{\Psi} \hat{f}_l^{(\gamma)}.
\end{equation}

Note that (4.5) is exactly (4.3). By making the connection \( s-1 \leftrightarrow \gamma \), we recover \( \lambda_{k,l}^{\Phi} \equiv \lambda_{k,l}^{\Psi} \), where \( \Phi^{(s)}_k(x) \) are the maps of the Szegő-Fourier functions \( \Psi^{(\gamma)}_k \). We stress again that this result is nontrivial. This also yields the functional connection
\begin{equation}
\Psi^{(\gamma)}_m(\theta) = \begin{cases} \sum_{|k|\leq|m|} \lambda_{k,m}^{\Psi} \Psi^{(\gamma+G)}_k(\theta), & |m| \leq G \\ \sum_{m-\geq G} \lambda_{k,m}^{\Psi} \Psi^{(\gamma+G)}_k(\theta), & |m| > G, \end{cases}
\end{equation}
i.e. \( \Psi^{(\gamma)}_m(\theta) \) is a linear combination of at most \( 2G + 1 \) functions \( \Psi^{(\gamma+G)}_k \). Note that the Fourier relation (4.6) parallels (4.5) in exactly the same way that the Jacobi relations (A.4) - (A.5) parallel (A.6) - (A.7).

We now illustrate how to calculate the Szegő-Fourier connection coefficients \( \lambda^{\Psi} \) in Proposition 4.1 from the Jacobi coefficients \( \lambda^{P} \). In the following, we make use of the notation:
\begin{align*}
n &:= |k| - 1, \quad \alpha = -\frac{1}{2}, \quad \beta = \gamma - \frac{1}{2}.
\end{align*}

From the definition of \( \Psi^{(\gamma)}_k \) in (2.4) we have
\begin{align*}
\tilde{P}^{(\alpha,\beta)}_{n+1} &= \Psi^{(\gamma)}_k + \Psi^{(\gamma)}_{-k} \\
\tilde{P}^{(\alpha+1,\beta+1)}_{n} &= \Psi^{(\gamma)}_{|k|} - \Psi^{(\gamma)}_{-|k|} \\
\tilde{P}^{(\alpha,\beta)}_{0} &= \sqrt{2} \Psi^{(\gamma)}_{0}.
\end{align*}
Therefore, from the modes \( \hat{f}_k^{(\gamma)} \) we can derive two sets of Jacobi modes:

\[
\hat{\epsilon}_n^{(\alpha,\beta)} = \hat{f}_n^{(\gamma)} + \hat{f}_{-n}^{(\gamma)}, \quad n \geq 1,
\]
\[
\hat{\delta}_n^{(\alpha+1,\beta+1)} = \hat{f}_{n+1}^{(\gamma)} - \hat{f}_{n-1}^{(\gamma)}, \quad n \geq 0,
\]
\[
\hat{\epsilon}_0^{(\alpha,\beta)} = \sqrt{2} \hat{f}_0^{(\gamma)}.
\]

The Jacobi modes \( \hat{\epsilon}_n \) are modes in an expansion in polynomials \( \hat{P}_n^{(\alpha,\beta)} \) and the modes \( \hat{\delta}_n \) are for an expansion in \( \hat{P}_n^{(\alpha+1,\beta+1)} \). With these modes in hand, we can use the Jacobi connection coefficients to promote the coefficients using Proposition 4.1 when \( n \geq 0 \):

\[
\hat{\epsilon}_n^{(\alpha,\beta+G)} = \sum_{m=0}^{G} \lambda_P^{(\alpha,\beta)} \hat{\epsilon}_n^{(\alpha,\beta)}, \quad \text{where} \quad \lambda_P = \lambda_P^{(\alpha,\beta,0,G)},
\]
\[
\hat{\delta}_n^{(\alpha+1,\beta+G+1)} = \sum_{m=0}^{G} \lambda_P^{(\alpha+1,\beta+1)} \hat{\delta}_n^{(\alpha+1,\beta+1)}, \quad \text{where} \quad \lambda_P = \lambda_P^{(\alpha+1,\beta+1,0,G)}.
\]

Finally we redistribute the modes back into Szegö-Fourier form to yield what we desired:

\[
\hat{f}_n^{(\gamma+G)} = \frac{1}{\sqrt{2}} \left[ \hat{\epsilon}_n^{(\alpha,\beta+G)} + \hat{\delta}_n^{(\alpha+1,\beta+1+G)} \right], \quad n \geq 1
\]
\[
\hat{f}_{-n}^{(\gamma+G)} = \frac{1}{\sqrt{2}} \left[ \hat{\epsilon}_n^{(\alpha,\beta+G)} - \hat{\delta}_n^{(\alpha+1,\beta+1+G)} \right], \quad n \geq 1
\]
\[
\hat{f}_0^{(\gamma+G)} = \frac{\hat{\epsilon}_0^{(\alpha,\beta+G)}}{\sqrt{2}}.
\]

The whole procedure is illustrated graphically in Figure 4. We may explicitly write
the connections as:

\[
\hat{f}_k^{(\gamma+G)} = \sum_{m=0}^{G} \left[ \frac{1}{2} \left( P_{\gamma}^{(\alpha,\beta)} k\mid k\mid + m \pm \text{sgn}(k) P_{\gamma}^{(\alpha+1,\beta+1)} k\mid k\mid + m \right) \right] \hat{f}_k^{(\gamma)} + \frac{1}{2} \sum_{m=0}^{G} \left[ \frac{1}{2} \left( P_{\gamma}^{(\alpha,\beta)} k\mid k\mid + m - \text{sgn}(k) P_{\gamma}^{(\alpha+1,\beta+1)} k\mid k\mid + m \right) \right] \hat{f}_k^{(\gamma)} + \frac{1}{2} \sum_{m=0}^{G} \left[ \frac{1}{2} \left( P_{\gamma}^{(\alpha,\beta)} k\mid k\mid + m \pm \text{sgn}(k) P_{\gamma}^{(\alpha+1,\beta+1)} k\mid k\mid + m \right) \right] \hat{f}_k^{(\gamma)} + \frac{1}{2} \sum_{m=0}^{G} \left[ \frac{1}{2} \left( P_{\gamma}^{(\alpha,\beta)} k\mid k\mid + m - \text{sgn}(k) P_{\gamma}^{(\alpha+1,\beta+1)} k\mid k\mid + m \right) \right] \hat{f}_k^{(\gamma)}.
\]

Therefore we have an explicit expression for the Szegö-Fourier connection coefficients in (4.1):

\[
\lambda^\Psi_{k,\pm}(\gamma + m) = \left\{ \begin{array}{ll}
\frac{1}{2} \left[ P_{\gamma}^{(\alpha,\beta)} k\mid k\mid + m \pm \text{sgn}(k) P_{\gamma}^{(\alpha+1,\beta+1)} k\mid k\mid + m \right], & |k| \geq 1 \\
\frac{1}{2} P_{\gamma}^{(\alpha,\beta)} k\mid k\mid + m, & k = 0
\end{array} \right.
\]

Of course, owing to observation (4.1), the above equation restricts \(0 \leq m \leq G\). As mentioned, this connection relation is also valid for converting an expansion in the functions \(\Phi_k^{(s)}(x)\) to one in the functions \(\Phi_k^{(s+S)}(x)\) for \(S \in \mathbb{N}\) since the modes for these two expansions are the same. Let \(f(\theta)\) be given and define \(g(x) = f(\theta(x))\). Then for all \(\gamma > \frac{1}{2}\):

\[
\hat{f}_k^{\Psi(\gamma)} = \left\langle f, \Psi_k^{(\gamma)} \right\rangle_w^{(\gamma,0)} = \left\langle g, \Phi_k^{(\gamma+1)} \right\rangle_w^{(\gamma+1,0)} = \hat{g}_k^{\Phi(\gamma+1)}.
\]

This completes the \(\Psi\)-\(\Psi\) connection problem. The reverse connection problem (converting \(f_k^{\Psi, (\gamma+G)}\) modes to \(f_k^{\Psi, (\gamma)}\) modes) is solved by reversing the above procedure (all steps are invertible) and use of the fact that the forward Jacobi connection problem with integral separation is banded upper-triangular and thus the backward connection is \(O(N)\) calculable sequentially via back-substitution. See [25].

We have determined how to quickly and exactly accomplish the connection problems for the unweighted functions

\[
\sum_k \hat{f}_k^{\Psi, (\gamma)} \Phi_k^{\Psi(\gamma)}(\theta) \mapsto \sum_k \hat{f}_k^{\Psi, (\gamma+G)} \Phi_k^{\Psi(\gamma+G)}(\theta),
\]

\[
\sum_k g_k^{\Phi, (s)} \Phi_k^{\Phi(s)}(x) \mapsto \sum_k \hat{g}_k^{\Phi, (s+S)} \Phi_k^{\Phi(s+S)}(x),
\]

in \(O(N)\) time where \(N\) is the total number of modes when \(S, G \in \mathbb{Z}\). These connections can be performed by utilizing the connection coefficients in (4.7) along with the sparse connection result of Proposition 4.1. For \(S, G \not\in \mathbb{Z}\), there is no sparse connection result for the modes, and so while the connection coefficients \(\lambda^\Psi_{k,\pm}\) can still be calculated based on known connection coefficients for Jacobi polynomials, the coefficients do not terminate finitely, and it is more expensive (that is, more costly than \(O(N)\)) to change \(s\) or \(\gamma\).

We have not described the details of how this \(\Psi\)-\(\Psi\) connection problem relates to the two issues presented at the beginning of this section (i.e., using the FFT and modification of \(s\) for the weighted functions \(\phi^{(s)}(x)\)). The problem of using the FFT we will postpone until Part II, which describes computational issues. In
Section 4.2.3 we will describe a method for modification of the decay parameter \( s \), for which the connection process described in this section is an integral part.

4.2.2. The \( \Psi\)-\( \psi \) Connection Problem. We now consider the following problem: let \( f \in L^2 ([\pi, \pi], \mathbb{C}) \). We assume \( \gamma \geq 0 \) and consider two expansions:

\[
f(\theta) = \sum_{k \in \mathbb{Z}} \hat{f}_k^\Psi \Psi_k(\gamma)(\theta),
\]

\[
f(\theta) = \sum_{k \in \mathbb{Z}} \hat{f}_k^\psi \psi_k(\gamma)(\theta).
\]

The modal coefficients are defined in the following way:

\[
\hat{f}_k^\Psi = \left( f, \Psi_k(\gamma) \right)_{w(\gamma)}^w,
\]

\[
\hat{f}_k^\psi = \left( f, \psi_k(\gamma) \right)_{w(\gamma)}^w.
\]

We assume that the modal coefficients for the uppercase (unweighted function) expansion are known and that we wish to determine the lowercase modes \( \hat{f}_k^\psi \). From the definitions of the modal coefficients, it is clear that we can rewrite the lowercase modes as

\[
\hat{f}_k^\psi = \left( f, \psi_k(\gamma) \right)_{w(\gamma)}^w = \left( f \left[ \sqrt{2}i (1 + e^{-i\theta}) \right]^{-\gamma}, \Psi_k(\gamma) \right)_{w(\gamma)}^w.
\]

That is, the modal coefficients for the lowercase basis are identical to modal coefficients of a different function for the uppercase basis. To see how this helps us, we make a small digression; recall (2.6) and define

\[
g(\theta) := f \left[ \sqrt{w(\gamma)} \right]^{-\gamma} = f \times \left[ \frac{\sqrt{2}}{i (1 + e^{-i\theta})} \right]^{\gamma}.
\]

Suppose that \( \gamma = G \in \mathbb{N}_0 \) and that we can somehow find the modal coefficients \( \hat{g}_k^{\Psi,(0)} \) for \( \gamma = G \) due to the sparse connection. To see how we can find the modal coefficients \( \hat{g}_k^{\Psi,(0)} \), assume that we have the modal coefficients \( \hat{f}_k^{\Psi,(0)} \), assume that we have the modal coefficients \( \hat{f}_k^{\Psi,(0)} \). Then (4.8) implies that

\[
\sum_{m=0}^{G} \binom{G}{m} \hat{g}_{k+m}^{\Psi,(0)} = \hat{f}_k^{\Psi,(0)} \left( \frac{\sqrt{2}}{i} \right)^G.
\]

If we assume a finite expansion so that \( \hat{g}_k = 0 \) for \( |k| > 2N + 1 \), then we can solve (4.9) via back-substitution. Note that determining each coefficient costs \( O(G) \) operations, independent of \( N \); this is a similar operation count to the \( \Psi\)-\( \Psi \) connection cost.

Finally, we must obtain \( \hat{f}_k^{\Psi,(0)} \) from the given input \( \hat{f}_k^{\Psi,(G)} \). However, this is another \( \Psi\)-\( \Psi \) connection (albeit in reverse). Therefore, the three steps to take us from \( \hat{f}_k^{\Psi,(G)} \) modes to \( \hat{f}_k^{\psi,(G)} \) modes are
4.2.3. Modification Of $s$: The $\psi$-$\psi$ Connection. We have now developed the necessary tools for the modification of $s$, i.e., the $\psi$-$\psi$ connection problem. We assume that $G, F \in \mathbb{N}$ and that we know connection coefficients of some function $f \in L^2$ for an expansion in $\psi^{(F)}$, and wish to obtain the coefficients for a $\psi^{(G)}$ expansion. The whole procedure can be accomplished in three steps:

1. Compute $\hat{f}_{k}^{\psi,(0)}$ from $\hat{f}_{k}^{\psi,(G)}$, which is a (backward) $\Psi$-$\Psi$ connection
2. Compute $\hat{g}_{k}^{\psi,(0)}$ from $\hat{f}_{k}^{\psi,(0)}$ using (4.9).
3. Compute $f_{k}^{\psi,(G)} = g_{k}^{\psi,(G)}$ from $g_{k}^{\psi,(0)}$, a (forward) $\Psi$-$\Psi$ connection.

This is illustrated in Figure 5. For an expansion with $N$ modes, all three steps have $O(NG)$ cost asymptotically. The backward connection problem (determining $\hat{f}_{k}^{\psi,(G)}$ from $\hat{f}_{k}^{\psi,(G)}$) is also computable in $O(NG)$ operations, and is accomplished by reversing the above operations.

Note that if $\gamma \notin \mathbb{N}_0$ then all of these steps break down: the $\Psi$-$\Psi$ connection is not sparse, and (4.9) is not valid since $\gamma$ is not an integer in (4.8).

Performing these modal connections on the real line for expansions in $\Phi^{(s)}(x)$ and $\phi^{(s)}(x)$ is equivalent, except one must assign $\gamma := s - 1$ and then proceed as outlined above.

This particular connection problem is not necessarily useful explicitly since in many of our applications, we will have direct access to $\hat{f}_{k}^{\psi,(0)}$, but each of the pieces necessary for this computation are used extensively both in modification of the decay parameter $s$ and application of the FFT.
\( g = f \times \left[ \sqrt{w^{(3)}} \right]^{-1} \)

\( h = f \times \left[ \sqrt{w^{(5)}} \right]^{-1} \)

**Figure 6.** Flowchart of operations for modification of \( s \). The operator \(-=\) is the subtraction-assignment operator.

1. Obtain expansion coefficients for \( f \times \left[ \sqrt{w^{(F)}} \right]^{-1} \) in the \( \Psi^{(0)} \)
   \( (\Psi-\psi \text{ connection}) \)

2. Obtain expansion coefficients for \( f \times \left[ \sqrt{w^{(G)}} \right]^{-1} \) in the \( \Psi^{(0)} \)
   \( (\text{Fourier connection}) \)

3. Obtain the sought expansion coefficients of \( f \) in the \( \psi^{(G)} \)
   \( (\Psi-\psi \text{ connection}) \)

Step 2 is easily performed using a version of (4.9) by noting the relation between
\( f \times \left[ \sqrt{w^{(G)}} \right]^{-1} \) and \( f \times \left[ \sqrt{w^{(F)}} \right]^{-1} \) with knowledge of the canonical Fourier expansion coefficients \( (\Psi^{(0)}(\theta)) \). This is shown in Figure 6 for the special case \( F = 3, G = 5 \).

Note that this particular connection problem is very amenable to an FFT+collocation approach whereas the algorithm we have laid out is a ‘Galerkin’ approach. The problem with the collocation approach is that it requires \( \mathcal{O}(N \log N) \) operations with two FFT’s, whereas the above algorithm requires only \( \mathcal{O}(NG) \) steps.
As with the $\Psi$-$\psi$ connection of the previous section, if either the starting parameter $F$ or the target parameter $G$ are not integers, then this procedure cannot be used: the core of the fast algorithm is the ability to obtain the canonical Fourier modes, which cannot be done efficiently if the decay parameters are not integers.

### 4.3. Quadrature

We now turn to quadrature rules that will compute integrals over the real line. We adopt the following notation: the pair $\{r_n^{(\alpha, \beta)}(\gamma), \omega_n^{(\alpha, \beta)}(\gamma)\}_{n=1}^N$ denotes the $N$-point Gauss-quadrature for the Jacobi polynomial of class $(\alpha, \beta)$, i.e.,

$$
\int_{-1}^{1} f(r) w_r^{(\alpha, \beta)} \, dr = \sum_{n=1}^{N} f\left(r_n^{(\alpha, \beta)}(\gamma)\right) \omega_n^{(\alpha, \beta)}(\gamma), \quad \forall f \in B_{2N-1}^{N-1},
$$

where $B_{2N-1}^{N-1}$ is the space of polynomials of degree $2N-1$ or less. We suppress the dependence of $r_n^{(\alpha, \beta)}(\gamma)$ and $\omega_n^{(\alpha, \beta)}(\gamma)$ on $N$. We also denote $\{r_n^{(\alpha, \beta)\text{GR}}, \omega_n^{(\alpha, \beta)\text{GR}}\}_{n=1}^N$ as the $N$-point Gauss-Radau quadrature with the fixed node $r_1^{(\alpha, \beta)\text{GR}} \equiv 1$. We assume for clarity of presentation that the nodes are ordered by $n$, e.g., $r_{n-1}^{(\alpha, \beta)} < r_n^{(\alpha, \beta)}$.

With the goal that we wish to develop quadrature rules for the infinite line, we will take pains to develop quadrature rules in $\theta$-space that do not have nodes at $\theta = \pm \pi$, which map to $x = \pm \infty$. We use the Jacobi-Gauss quadrature rules as the building blocks for our generalized Fourier quadrature rules.

Suppose we wish to construct an $N$-point quadrature rule associated with the functions $\Psi_k^{(\gamma)}(\theta)$. If $N$ is even, then define

$$
\theta_n^{(\gamma)} = \begin{cases} 
- \arccos \left( r_n^{(-1/2, \gamma-1/2)} \right), & 1 \leq n \leq \frac{N}{2} \\
- \theta_{N+1-n}^{(\gamma)}, & \frac{N}{2} + 1 \leq n \leq N,
\end{cases}
$$

where $r_n^{(\alpha, \beta)}$ comes from an $\frac{N}{2}$-point quadrature rule, and

$$
\Omega_n^{(\gamma)} = \begin{cases} 
\omega_n^{(-1/2, \gamma-1/2)}, & 1 \leq n \leq \frac{N}{2} \\
\Omega_{N+1-n}^{(\gamma)}, & \frac{N}{2} + 1 \leq n \leq N,
\end{cases}
$$

and $\omega_n^{(\alpha, \beta)}$ comes from an $\frac{N}{2}$-point quadrature rule.

If $N$ is odd, then define

$$
\theta_n^{(\gamma)} = \begin{cases} 
- \arccos \left( r_n^{(-1/2, \gamma-1/2);\text{GR}} \right), & 1 \leq n \leq \frac{N+1}{2} \\
\theta_{N+1-n}^{(\gamma)}, & \frac{N+3}{2} \leq n \leq N,
\end{cases}
$$

where $r_n^{(\alpha, \beta);\text{GR}}$ comes from an $\frac{N+1}{2}$-point quadrature rule, and

$$
\Omega_n^{(\gamma)} = \begin{cases} 
\omega_n^{(-1/2, \gamma-1/2);\text{GR}}, & 1 \leq n \leq \frac{N+1}{2} \\
2 \omega_n^{(-1/2, \gamma-1/2);\text{GR}}, & n = \frac{N+1}{2} \\
\Omega_{N+1-n}^{(\gamma)}, & \frac{N+3}{2} \leq n \leq N.
\end{cases}
$$
Figure 7. Construction of Gauss-type quadrature for generalized Fourier functions. The new quadrature rules are symmetric combinations of Jacobi-Gauss-type quadrature rules. The constructions shown are accurate node locations for \( \gamma = 5 \).

For graphical descriptions of the above formulae, see Figure 7. We have used Jacobi-Gauss rules for \( N \) even and Jacobi-Gauss-Radau rules for \( N \) odd. By construction, when \( N \) is odd, \( \theta_{\frac{N+1}{2}}^{(\gamma)} = 0 \) due to the Gauss-Radau rule requirement that \( I_{\frac{\alpha,\beta}{N+1}}^{\text{GR}} \) = 1. The quadrature rules derived above have no nodes at \( \theta = \pm \pi \) (since there are no Jacobi-Gauss, or Jacobi-Gauss-Radau nodes at \( r = -1 \)) and are symmetric rules for any \( \gamma \). Thus they are always exact for any odd function. It is not difficult to show the following result:

**Proposition 4.2.** For \( N \) even, the \( N \)-point quadrature rule \( \{ \theta_n^{(\gamma)}, \Omega_n^{(\gamma)} \}_{n=1}^{N} \) satisfies

\[
\int_{-\pi}^{\pi} e^{ik\theta} w_0^{(\gamma,0)} d\theta = \sum_{n=1}^{N} e^{ik\theta_n^{(\gamma)}} \left( \theta_n^{(\gamma)} \right) \Omega_n^{(\gamma)}, \quad |k| \leq N - 1.
\]

When \( N \) is odd, the quadrature rule satisfies

\[
\int_{-\pi}^{\pi} e^{ik\theta} w_0^{(\gamma,0)} d\theta = \sum_{n=1}^{N} e^{ik\theta_n^{(\gamma)}} \left( \theta_n^{(\gamma)} \right) \Omega_n^{(\gamma)}, \quad |k| \leq N.
\]

The degeneracy in the quadrature rule for \( N \) even is exactly the same nature as the degeneracy in the canonical equispaced Fourier quadrature rule for an even number of grid points [18]. If \( \gamma = 0 \) the rule \( \{ \theta_n^{(0)}, \Omega_n^{(0)} \}_{n=1}^{N} \) is exactly the same as the equispaced Fourier quadrature rule, symmetric about \( \theta = 0 \). The quadrature rule \( \{ \theta_n^{(0)}, \Omega_n^{(0)} \}_{n=1}^{N} \) can be used to integrate against the weight function \( w_0^{(\gamma,0)} \) when \( \gamma \in \mathbb{N} \) since in this case the weight is itself a trigonometric polynomial.

In order to determine a quadrature rule to integrate the weighted functions \( \psi_k^{(\gamma)}(\theta) \), we can augment the weights \( \Omega_n^{(\gamma)} \) to contain information about the weight function. This can be summed up in the following result:
Corollary 4.1. The even $N$-point quadrature rule \( \left\{ \theta_n^{(\gamma)}, \omega_n^{(\gamma)} \right\}_{n=1}^N \), where \( \omega_n^{(\gamma)} := w_\theta^{(-\gamma,0)} \left( \theta_n^{(\gamma)} \right) \Omega_n^{(\gamma)} \) satisfies

\[
\int_{-\pi}^{\pi} \psi_k^{(\gamma)} \psi_l^{(\gamma)} d\theta = \sum_{n=1}^N \psi_k^{(\gamma)} \left( \theta_n^{(\gamma)} \right) \psi_l^{(\gamma)} \left( \theta_n^{(\gamma)} \right) \omega_n^{(\gamma)}, \quad |k| + |l| \leq N - 1
\]

Multiplying \( \Omega_n^{(\gamma)} \) by the inverse of the weight \( w_\theta^{(-\gamma,0)} \) is mathematically not a problem since none of the \( \theta_n^{(\gamma)} \) are equal to \( \pm \pi \), where the weight \( w_\theta^{(-\gamma,0)} \) is singular. Note that since the functions \( \Phi_k^{(s)}(x) \) are just a mapping of the functions \( \Psi_k^{(\gamma)}(\theta) \), the quadrature rule \( \left\{ x \left( \theta_n^{(s-1)} \right), \Omega_n^{(s-1)} \right\}_{n=1}^N \), which has nodal values over \( \mathbb{R} \), can be used to integrate the functions \( \Phi_k^{(s)}(x) \) over the real line. Similarly, the rule \( \left\{ x \left( \theta_n^{(s-1)} \right), \omega_n^{(s-1)} \right\}_{n=1}^N \) can be used to integrate Galerkin products of the generalized Wiener functions \( \phi_k^{(s)}(x) \) over the real line.

For various \( \gamma/s \) we graphically depict the location of the quadrature nodes for \( N = 21 \) in Figure 8 on the unit circle \( z \in \mathbb{T} \) and on the real line. Note that as we increase \( \gamma \) the quadrature nodes become more and more concentrated towards \( z = 1 \) (\( \theta = 0 \)). On the real line, this manifests itself as higher concentration near \( x = 0 \) which, although rectifiable via an affine mapping, is suboptimal if one wishes to resolve functions away from \( x = 0 \). Note that the tendency of Jacobi-Gauss nodes to become more equidistant on \([-1, 1]\) as \( \beta \) (i.e. \( \gamma \) or \( s \)) is increased [18] also suggests that these generalized quadrature rules for large \( \gamma \) or \( s \) will not be as good as the the ones for smaller \( \gamma \) or \( s \) since equidistant nodes are bad for finite-interval polynomial interpolation. In addition, when \( \gamma = 0 \), we can use these (equidistant) quadrature nodes to employ the FFT for modal-nodal transformations.

4.4. The Stiffness Matrix. In many applications to differential equations it is necessary to express the derivative of a basis function as a linear combination of basis functions. We devote this section to this endeavor. We define entries of the stiffness matrix as

\[
S_{k,l}^\phi = \left\langle \phi_k^{(s)}(x), \frac{d}{dx} \phi_l^{(s)}(x) \right\rangle.
\]

For the generalized Wiener rational functions, the following result can be proven:

Theorem 4.2. Let \( S^\phi \) denote the \( N \times N \) stiffness matrix for the weighted Wiener rational functions \( \phi_k^{(s)} \). \( S^\phi \) satisfies the following properties for any \( s > \frac{1}{2} \):

1. \( S^\phi \) is skew-Hermitian, i.e. \( S^\phi_{k,l} = -S^\phi_{l,k} \)
2. \( S^\phi \) is sparse with entries only on the super-, sub-, and main sinister and dexter diagonals: define

\[
k^\prime := \text{sgn}(k)(|k| - 1) = k - \text{sgn}(k), \quad k^\wedge := \text{sgn}(k)(|k| + 1) = k + \text{sgn}(k).
\]

Then

\[
\frac{d\phi_k^{(s)}(x)}{dx} = \sum_{l \in \{\pm k^\prime, \pm k, \pm k^\wedge\}} \phi_l^{(s)}(x),
\]
for some purely imaginary constants $\tau^{(s)}_{k,l}$. In other words,

$$S_{k,l}^\phi = 0, \quad l \notin \{\pm k^\vee, \pm k, \pm k^\wedge\}.$$

(3) The spectral radius of $S^\phi$ satisfies

$$\rho(S^\phi) \leq N + 5s.$$

The proof of Theorem 4.2 is quite tedious, so we only sketch the main points. Details are given in Appendix B.

**Proof.** Property 1 can easily be deduced by using integration by parts and noting that the functions $\phi_k^{(s)}(x)$ decay to zero as $|x| \to \infty$.

Property 2 is a highly nontrivial result that is provable using several properties of Jacobi Polynomials. We refer the reader to [24]. Most of the calculations are straightforward once a list of Jacobi Polynomial properties has been compiled. However, there are some difficulties whose resolutions rely on a couple of fortuitous properties: first, that $\frac{d}{d\theta}(\theta) = 1 + \cos \theta$, i.e. that the map we have chosen to take $\theta \to x$ has a Jacobian with a particular form. Second, that

$$\frac{d}{d\theta} \left[ (\sin \theta) \tilde{P}_{n}^{(\alpha+1, \beta+1)} (\cos \theta) \right]$$

is a sparse combination of $\tilde{P}_{n}^{(\alpha+1, \beta+1)} (\cos \theta)$, which is not a trivial result; we show this by using brute-force calculation with the compiled list of Jacobi Polynomial properties.

Property 3 can be derived from the second property. The key ingredient is Gerschgorin’s Theorem. Using the explicit entries for the constants $\tau^{(s)}_{k,l}$ given in
Theorem B.3 of Appendix B we can show that for each \(k\) satisfying \(|k| \geq 2\) the following crude bounds hold:

\[
|\tau_{k,k}| \leq n + 2s,
\]

\[
|\tau_{k,-k}| \leq |\tau_{k,k^\vee}| + |\tau_{k,-k^\vee}| + |\tau_{k,k^\wedge}| + |\tau_{k,-k^\wedge}| \leq n + 3s + 2,
\]

where \(n := |k| - 1\). Gershgorin’s Theorem can now be used to define a region in the complex plane in which all the eigenvalues lie. By the above properties, this region has distance from the origin at most \(2n + 5s + 2\). Once we consider the necessary relationship between \(n\), \(k\), and \(N\), the result is proven. (It is interesting, but not necessary, to note that the eigenvalues all lie on the imaginary axis due to the skew-Hermitian property of \(S\).)

\[\square\]

Remark 4.3. While the \(O(N)\) maximum eigenvalue does depend on \(s\), the proportionality factor is empirically around 2, not 5 as given in the theorem. See Table 2.

The sparsity pattern we have derived for the derivatives of these functions (property 2 of the above theorem) is illustrated in Figure 9. Note that the unweighted functions \(\Phi^{(s)}(x)\) also have a similar sparsity result; see Lemma B.2. However, the Fourier functions \(\Psi^{(\gamma)}(\theta)\) and \(\psi^{(\gamma)}(\theta)\) do not have sparse stiffness matrices (unless \(\gamma = 0\)). In addition, numerical values for the maximum eigenvalues of the stiffness matrix (property 3) are given in Table 2. The sparsity of the stiffness matrix is important for fast computations of derivatives for spectral methods for solving PDEs, and the \(O(N)\) maximum eigenvalue of the stiffness matrix indicates that we can take a relatively large timesteps for time-dependent problems. Finally, the skew-symmetry of the stiffness matrix easily leads to energy conservation for the Galerkin discretization of hyperbolic conservation laws.

5. THE SEMI-INFINITE INTERVAL

The generalized Wiener basis functions we have derived can be used for function expansions on the infinite line. In order to address expansions on semi-infinite
Figure 9. Sparsity plots for stiffness matrices of the weighted Wiener rational functions $\phi_k^{(s)}$. The sparsity patterns are representative of property 2 in Theorem 4.2 for $s = 1$ (left) and all $s \neq 1$ (right). The $s = 1$ sparsity pattern has been derived previously [13], and the expressions for the $\tau_{k,l}$ in Appendix B with $s = 1$ reduce to the pattern above.

interval, we can instead use either the even or odd Jacobi polynomial basis sets that make up the Fourier functions constructed in Section 2.3.

The Jacobi functions from Lemma 2.1 can be mapped and weighted in a procedure identical to the construction of the Wiener basis. The result is the collection of functions

$$
\rho_n^{(s)} = \sqrt{w_x^{(s)}(x)} \psi_n^{(s)}(x)
= \left( \frac{2}{x^2 + 1} \right)^{s/2} \tilde{P}_n^{(-1/2,s-3/2)} \left( \frac{1-x^2}{1+x^2} \right), \quad n \in \mathbb{N}_0
$$

These functions are a direct mapping and weighting of the Jacobi polynomials. Because of this, they are orthonormal and complete in $L^2(\mathbb{R}^+, \mathbb{R})$. Mapping techniques for classical functions are not novel and we discuss existing methods in Section 6. The classical competitor for spectral expansions on semi-infinite intervals is the set of Laguerre functions (weighted Laguerre polynomials). A comparison between the Laguerre functions and the functions defined in (5.1) will be made in Part II, and in Section 6 a different mapping transformaing Jacobi polynomials to the semi-infinite line will be addressed.
We make use of the regular square root function $\sqrt{w_x(s)}$ in (5.1) instead of the phase-shifted version $\sqrt{w_x^{(s)}}$ because there is no need to have complex-valued functions. The phase-shifted square root was a convenient choice for the Wiener functions on the infinite line: its compact Fourier series representation (2.6) enabled fast connections (Section 4.2) and sparse differentiation matrices (Section 4.4). By using the real-valued square root in (5.1) we sacrifice these two properties. However, the FFT can still be used for the evaluation of modal coefficients if $s$ is an integer.

The caveat in using these functions for expansions on the semi-infinite interval is the fact that they all have zero-valued odd derivatives at $x = 0$. This parallels the same property at $\theta = 0$ for a cosine series on $\theta \in [0, \pi]$. Alternative mappings of the Jacobi polynomials to the semi-infinite line do not exhibit this restriction, but those mappings also preserve the $O(N^2)$ time-stepping restriction for nodal-based polynomial solvers of time-dependent partial differential equations using explicit time-integration on finite intervals. In constrast, the functions (5.1) only have an $O(N)$ time-step restriction, similar to the time-step restriction for a finite-interval cosine basis expansion.

The restriction of the Wiener functions to the semi-infinite interval as defined in (5.1) comes both with advantages and sacrifices. These functions are purely weighted maps of Jacobi polynomials and are therefore easy to implement. Some of the attractive features of the Wiener rational basis functions on infinite intervals are lost (e.g. sparse stiffness matrices). However, these functions have properties that are advantageous when compared with existing mapping techniques (Section 6). A numerical comparison between those mapping techniques, the functions (5.1), and the Laguerre functions will be made in Part II.

6. Alternative Methods

Before concluding this article with a summary of the derived properties of the generalized Wiener basis, we first summarize existing results on the topic of mapping Jacobi polynomials from the finite interval to the infinite interval. This method is very closely related to our strategy of mapping a generalized Fourier series from the canonical finite Fourier interval to the real line. Numerical studies comparing these methods are presented in Part II, but it is appropriate to acknowledge these functions here, and to discuss how they relate to the Wiener rational function basis.

6.1. The Infinite Interval. The main idea for our generalization of Wiener’s original rational basis is using a ‘well-behaved’ mapping to transform functions on a finite interval to those on an infinite interval. This basic idea is classical [15]. Indeed one of the more popular mappings that has gained momentum in the literature are the so-called ‘mapped Chebyshev’ functions/polynomials.

In order to further generalize the mapped Chebyshev functions, we will briefly restate their derivation from our point of view. We begin with the Jacobi polynomials $P_n^{(\alpha,\beta)}(r)$ on $r \in [-1, 1]$. Mapping via $r = \cos \theta$ to $\theta \in [0, \pi]$ yields trigonometric polynomials. We now ‘stretch’ the domain to $\Theta \in [-\pi, \pi]$ via the affine mapping $\Theta = 2\theta - \pi$. Finally, we utilize the usual linear fractional map $e^{i\Theta} = \frac{e^{i\theta} - 1}{i + e^{i\theta}}$ (i.e. rotation of the Riemann Sphere) to yield functions on the real line $x \in \mathbb{R}$. For all
s, t > \frac{1}{2}, this results in the functions PB_{n}^{(s,t)}(x), defined as

$$PB_{n}^{(s,t)}(x) = \tilde{P}_{n}^{((2s-3)/2,(2t-3)/2)} \left( \frac{x}{\sqrt{1+x^2}} \right),$$

orthonormal on the real line under the weight

$$w_{PB}^{(s,t)} = \left[ 1 - \frac{x}{\sqrt{1+x^2}} \right]^{(2s-3)/2} \left[ 1 + \frac{x}{\sqrt{1+x^2}} \right]^{(2t-3)/2},$$

and the weighted functions

$$pb_{n}^{(s,t)} = \sqrt{w_{PB}^{(s,t)}} PB_{n}^{(s,t)},$$

are orthonormal under the unweighted inner product. When s = t = 1, the functions PB_{n}^{(s,t)} coincide with the mapped Chebyshev polynomials TB_{n}(x) introduced in [4] and subsequently developed in [8] and [10], although the original idea of applying spectral expansions over finite intervals to solving problems over infinite intervals seems to come from [16]. In any case, the mapped Jacobi functions pb_{n}^{(s,t)} decay like $\frac{1}{|x|}$ for $x \to -\infty$ and $\frac{1}{|x|^t}$ for $x \to +\infty$. The advantage of these functions is that the decay can be different as $|x| \to \infty$. Also, others have already explored some convergence theory in function spaces [3] and applications to differential equations [31] for the Chebyshev case $s = t = 1$. In Part II when we present numerical examples, we will use the basis set pb_{n}^{(s,t)} with $s = t = 1$, i.e. the Chebyshev case.

Note that because all of these mapped types of polynomials and the generalized Wiener basis we have presented ultimately stem from Jacobi polynomials and mappings of similar character, all these basis sets are related in some fashion. To relate the mapped Jacobi functions to the generalized Wiener rational functions, we have

$$PB_{n}^{(s,s)}(x) \propto \text{Re} \left\{ \Phi_{n}^{(s)} \left( \frac{x + \sqrt{x^2 + 1} - 1}{x - \sqrt{x^2 + 1} + 1} \right) \right\}.$$ 

In Table 3 we relate the unweighted functions to the generalized Wiener rational basis, modulo multiplicative constants. In this article we make no observations about how mapped Jacobi polynomials compare to the Wiener basis set as a practical tool for function expansions. However, such a comparison will be a central theme in Part II.

### 6.2. The Semi-Infinite Interval

To perform spectral expansions on semi-infinite intervals, the only classical technique is the Laguerre polynomial/function method. However, mapping techniques can be used to transform finite-interval methods to semi-infinite interval methods.

As with Section 6.1, we explain the choice of mapping from our point of view as a mapping of the Riemann Sphere. The Jacobi polynomials are defined on $r \in [-1, 1]$. If we allow complex values of $r$, then we may consider using a linear fractional map to transform the Jacobi polynomial domain to the semi-infinite line. The ordered assignments $r = \{1, 0, -1\}$ to $x = \{0, 1, \infty\}$ specify the transformation uniquely as

$$x = \frac{1-r}{1+r} \quad r = \frac{1-x}{1+1}.$$ 

(6.1)

If necessary, one can also specify the relationship to $\theta$ and the cosine series on $[0, \pi]$. For details, see [7]. Our definition of the transformation differs only in orientation from that presented in [7]. We have chosen this orientation so that the Jacobi
Previous function  | Name/classification          | Interval     | Reference | Relation |
--- | --- | --- | --- | --- |
TB\(_n\)  | Chebyshev rational functions (1st)       | \(\mathbb{R}\) | [9], [4], [6]  | PB\(^{(1,1)}_n\) |
SB\(_n\)/UB\(_n\)  | Chebyshev rational functions (2nd)     | \(\mathbb{R}\) | [9], [6], [10]  | PB\(^{(2,2)}_n\) |
C\(_n\)/CC\(_n\)  | Christov functions (even)          | \([0, \infty)\) | [13], [9]  | Im\(\phi^{(1)}_n\) |
S\(_n\)/SC\(_n\)  | Christov functions (odd)           | \([0, \infty)\) | [13], [9]  | Re\(\phi^{(1)}_n\) |
CH\(_n\)  | Higgins functions (even)          | \([0, \infty)\) | [9]  | Re\(\Phi^{(1)}_n\) |
SH\(_n\)  | Higgins functions (odd)           | \([0, \infty)\) | [9]  | Im\(\Phi^{(1)}_n\) |
\(\rho_k\)  | (Complex) Higgins functions        | \([0, \infty)\) | [19], [13]  | \(\Phi^{(1)}_k\) |
\(\sigma_k\)  | (Complex) Wiener rational functions | \([0, \infty)\) | [30], [13]  | \(\phi^{(1)}_k\) |
TL\(_n\)  | Half-infinite Chebyshev rational functions  | \([0, \infty)\) | [7]  | PL\(^{(1/2)}_n\) |

**Table 3.** Relationship between orthogonal functions in previous work and the current bases presented.

The parameter \(\beta\) is assigned to the location \(x = \infty\) in order to mimic to the same assignment for the Wiener functions.

In the literature the maps of the Chebyshev polynomials under the transformation (6.1) are labeled TL\(_n\)(\(x\)). Adopting similar notation, we define

\[
PL^{(s)}_n(x) = \tilde{P}^{(-1/2, 2s - 2)}_n \left( \frac{1 - x}{1 + x} \right), \quad x \in [0, \infty],
\]

which are \(L^2\)-complete and orthonormal under the weight function

\[
w^{(s)}_{PL}(x) = \frac{1}{2\sqrt{x}} \left( \frac{2}{1 + x} \right)^{(2s)}.
\]

It is then possible to define the weighted functions

\[
pl^{(s)}(x) = \left( \frac{2}{1 + x} \right)^s PL^{(s)}_n(x)
\]

(6.2)

which are \(L^2\)-complete and orthonormal under the weighted \(L^2\) inner product

\[
\langle f, g \rangle_{w^{(s)}_{PL}} = \int_0^\infty f(x) g(x) \frac{1}{2\sqrt{x}} \, dx,
\]

for any \(s > \frac{1}{2}\). The \(pl^{(s)}_n\) are defined for \(x \in [0, \infty]\) and decay like \(x^s\) as \(x \to \infty\). A significant difference between the Wiener-type functions (both on the infinite and semi-infinite intervals) and the set defined in (6.2) is the fact that these functions are not orthogonal in the unweighted \(L^2\) inner product, but instead in the norm defined by the above inner product. This choice was made (as opposed to defining functions in the unweighted inner product) to ensure that integer values of \(s\) resulted in a Jacobi polynomial family that was amenable to usage of the FFT.

The main observation we make regarding this basis is that these functions are the result of a linear fractional map directly from the Jacobi domain; therefore,
they will inherit the $O(N^2)$ time-step restriction of nodal explicit time-integration methods for time-dependent partial differential equations. The same observation can be made about the functions defined in [7].

7. Conclusion

We have presented a collection of generalized Fourier series which, when mapped and weighted appropriately, generates a basis set on the infinite interval with a tunable rate of decay. For each rate of decay $s$ satisfying $s > \frac{1}{2}$ the resulting basis set $\phi^{(s)}_k$:

- is orthonormal and complete in $L^2(\mathbb{R}, \mathbb{C})$
- is characterized by $x^{-s}$ decay for $|x| \to \infty$
- can be generated via Jacobi polynomial recurrence relations
- has sparse connection properties that can be efficiently exploited via combinations of sparse Fourier and Jacobi connections
- has an $N \times N$ Galerkin stiffness/differentiation matrix that has at most $6N$ nonzero entries with $O(N)$ spectral radius
- is characterized by a ‘Gauss-like’ quadrature rule.

When $s \in \mathbb{N}$, the basis set is a rational function; we will show in Part II that in this case we can use the FFT for modal-nodal transformations. The case $s = 1$ corresponds to a mapping and weighting of the canonical Fourier series, as discovered by others previously. Due to the original presentation of the $s = 1$ basis by Wiener [30], we call the functions $\phi^{(s)}_k$ the generalized Wiener rational basis functions.

These basis functions have a similar flavor to directly mapped and weighted Jacobi polynomials (called $p\phi^{(s)}_n$ here). In Part II we will compare these basis sets and discuss advantages and disadvantages of each. In addition, we will also employ the Sinc and Hermite functions in test cases in an attempt to investigate a relatively broad class of spectral approximation methods. In contrast to [26] which reviews much of the theory present for expansions on the infinite interval, we concentrate on numerical issues, including application of the FFT. We will extend our investigation to the semi-infinite interval to compare the Laguerre polynomials/functions, the mapped Jacobi functions (denoted $p\ell^{(s)}_n$ here), and the restriction of the Wiener functions to the semi-infinite interval as given in Section 5.

We do not wish to claim that, on the infinite or semi-infinite intervals, genuinely global spectral expansions are truly superior to alternative numerical approximations; rather we wish to identify the generalized Wiener basis set as a novel competitor to existing global spectral expansions. Part II will follow up to show that the Wiener basis set is very competitive with existing expansions.

Appendix A. Recurrence Coefficients

In this appendix we compile various recurrence relations for the Jacobi/Szegő-Fourier/Wiener rational functions. We state the recurrences in terms of the Szegő-Fourier functions $\Psi^{(s)}_k$, but note that they all apply equally well to the unweighted Wiener rational functions as well. Note that we only list recurrences for $k \geq 0$; for $k < 0$, we may use the conjugation relation (3.1) to obtain $\Psi^{(s)}_{-|k|}$ at almost no additional cost. We first require a tour of some Jacobi polynomials recurrences:
\[
\sqrt{b_{n+1}} \tilde{P}_{n+1}^{(\alpha, \beta)} = \left[ r - a_n^{(\alpha, \beta)} \right] \tilde{P}_n^{(\alpha, \beta)} - \sqrt{b_n} \tilde{P}_{n-1}^{(\alpha, \beta)},
\]

(A.1)

\[
(1 - r^2) \tilde{P}_n^{(\alpha, \beta)} = \sum_{i=0}^{2} \varepsilon_{n,i}^{(\alpha, \beta)} \tilde{P}_{n+i}^{(\alpha-1, \beta-1)},
\]

(A.2)

\[
\tilde{P}_n^{(\alpha, \beta)} = \sum_{i=0}^{2} \eta_{n-i}^{(\alpha, \beta)} \tilde{P}_{n-i}^{(\alpha+1, \beta+1)},
\]

(A.3)

\[
(1 - r) \tilde{P}_n^{(\alpha, \beta)} = \mu_{n,0}^{(\alpha, \beta)} \tilde{P}_{n-1}^{(\alpha, \beta)} - \mu_{n,1}^{(\alpha, \beta)} \tilde{P}_{n+1}^{(\alpha, \beta)},
\]

(A.4)

\[
(1 + r) \tilde{P}_n^{(\alpha, \beta)} = \nu_{n,0}^{(\alpha, \beta)} \tilde{P}_{n-1}^{(\alpha, \beta-1)} + \nu_{n,-1}^{(\beta, \alpha)} \tilde{P}_{n-1}^{(\alpha+1, \beta)},
\]

(A.5)

\[
\tilde{P}_n^{(\alpha, \beta)} = \nu_{n,0}^{(\beta, \alpha)} \tilde{P}_{n+1}^{(\alpha, \beta)} + \nu_{n,-1}^{(\beta, \alpha)} \tilde{P}_{n-1}^{(\alpha, \beta+1)},
\]

(A.6)

\[
\frac{d}{dr} \tilde{P}_n^{(\alpha, \beta)} = \gamma_{n}^{(\alpha, \beta)} \tilde{P}_{n-1}^{(\alpha+1, \beta+1)},
\]

(A.7)

\[
\mu_{n,0}^{(\alpha, \beta)} = \sqrt{\frac{2(n + \alpha)(n + \alpha + \beta)}{(2n + \alpha + \beta)(2n + \alpha + \beta + 1)}},
\]

(A.9)

\[
\mu_{n,1}^{(\alpha, \beta)} = \sqrt{\frac{2(n + 1)(n + \beta + 1)}{(2n + \alpha + \beta + 1)(2n + \alpha + \beta + 2)}},
\]

(A.10)

\[
\nu_{n,0}^{(\alpha, \beta)} = \sqrt{\frac{2(n + \alpha + 1)(n + \alpha + \beta + 1)}{(2n + \alpha + \beta + 1)(2n + \alpha + \beta + 2)}},
\]

(A.11)

\[
\nu_{n,-1}^{(\alpha, \beta)} = \sqrt{\frac{2n(n + \beta)}{(2n + \alpha + \beta)(2n + \alpha + \beta + 1)}},
\]

(A.12)

\[
\nu_{n}^{(\alpha, \beta)} = \sqrt{n(n + \alpha + \beta + 1)}.
\]

(A.13)

where \(\mu_{n,0}^{(\alpha, \beta)}, \nu_{n,0}^{(\alpha, \beta)},\) and \(\gamma_{n}^{(\alpha, \beta)}\) in (A.4)-(A.8) are constants for which we take explicit formulae from [25]:
The three-term recurrence coefficients in (A.1) are given by [14]:

(A.14) \( a_n^{(\alpha, \beta)} = \begin{cases} \frac{\beta - \alpha}{\alpha + \beta + 2}, & n = 0, \\ \frac{2^{n+1} \Gamma(\alpha + 1) \Gamma(\beta + 1)}{\Gamma(\alpha + \beta + 2)}, & n > 0. \end{cases} \)

(A.15) \( b_n^{(\alpha, \beta)} = \begin{cases} \frac{2^{n+1} \Gamma(\alpha + 1) \Gamma(\beta + 1)}{\Gamma(\alpha + \beta + 2)}, & n = 0, \\ \frac{4(\alpha + 1)(\beta + 1)}{(\alpha + \beta + 2)^2}, & n = 1, \\ \frac{4(n + \alpha)(n + \beta)(n + \alpha + \beta)}{(2n + \alpha + \beta - 1)(2n + \alpha + \beta + 1)}, & n > 1. \end{cases} \)

The demotion recurrence coefficients in (A.2) can be obtained by determining the analogous relations for the monic orthogonal polynomials ([1], [27]) and then employing the appropriate normalizations:

(A.16) \( \varepsilon_{n,0}^{(\alpha, \beta)} = \begin{cases} 2^{\alpha \beta}, & n = 0, \\ \frac{1}{(\alpha + \beta + 2) \sqrt{\alpha + \beta}}, & n = 1, \\ \frac{2}{(n + \alpha)(n + \beta)(n + \alpha + \beta)} \frac{1}{(2n + \alpha + \beta - 1)(2n + \alpha + \beta + 1)}, & n > 1. \end{cases} \)

(A.17) \( \varepsilon_{n,1}^{(\alpha, \beta)} = \begin{cases} 2\sqrt{\alpha + \beta}, & n = 0, \\ \frac{2(\alpha - \beta)(n + \alpha + \beta)}{(2n + \alpha + \beta)(2n + \alpha + \beta + 2)}, & n > 0. \end{cases} \)

(A.18) \( \varepsilon_{n,2}^{(\alpha, \beta)} = \begin{cases} \frac{2(\alpha + 1)(\beta + 1)}{(\alpha + \beta + 1)(\alpha + \beta + 3)}, & n = 0, \\ \frac{2}{\alpha + \beta + 2} \sqrt{\frac{2(\alpha + 1)(\beta + 1)}{(\alpha + \beta + 1)(\alpha + \beta + 3)}}, & n > 0. \end{cases} \)

Finally, the promotion relation (A.3) coefficients can also be determined:

(A.19) \( \eta_{n,0}^{(\alpha, \beta)} = \varepsilon_{n,0}^{(\alpha + 1, \beta + 1)}, \)

\( \eta_{n,-1}^{(\alpha, \beta)} = \varepsilon_{n,-1}^{(\alpha + 1, \beta + 1)}, \)

\( \eta_{n,-2}^{(\alpha, \beta)} = -\varepsilon_{n,-2}^{(\alpha + 1, \beta + 1)}. \)

Of course, (A.2)-(A.3) are consequences of combining (A.4)-(A.7). Using the orthogonal polynomial three-term recurrence relation (A.1) we can show the following recurrence relation for the Szegö-Fourier functions \( \Psi_n^{(\gamma)}(\theta) \):

\( \Psi_{n+1}^{(\gamma)} = \left[ U_n^{(\gamma)} \cos \theta - V_n^{(\gamma)} \right] \Psi_n^{(\gamma)} + \left[ U_n^{(\gamma)} \cos \theta - V_n^{(\gamma)} \right] \Psi_{-n}^{(\gamma)} - W_n^{(\gamma)} \Psi_{n-1}^{(\gamma)} - W_n^{(\gamma)} \Psi_{-(n-1)}^{(\gamma)}. \)
In the following expressions, we make use of the following definitions: for a given \( \gamma > -\frac{1}{2} \),

\[
\alpha := -\frac{1}{2}, \quad \beta := \gamma - \frac{1}{2}.
\]

The recurrence constants are then given by

\[
U^{(\gamma)}_{\pm n} = \frac{1}{2} \left[ \sqrt{\frac{1}{b_n^{(\alpha,\beta)+1}}} \pm \sqrt{\frac{1}{b_n^{(\alpha+1,\beta)+1}}} \right],
\]

\[
V^{(\gamma)}_{\pm n} = \pm \frac{1}{2} \left[ \frac{\sigma_n^{(\alpha,\beta)}}{b_n^{(\alpha,\beta)+1}} \pm \frac{\sigma_{n-1}^{(\alpha+1,\beta+1)}}{b_n^{(\alpha+1,\beta+1)+1}} \right],
\]

\[
W^{(\gamma)}_{\pm n} = \pm \frac{1}{2} \left[ \sqrt{\frac{1}{b_n^{(\alpha,\beta)+1}}} \pm \sqrt{\frac{1}{b_n^{(\alpha+1,\beta)+1}}} \right].
\]

Using the promotion and demotion three-term recurrences (A.2-A.3) we also have the following recurrence relation:

\[(A.20) \quad \Psi^{(\gamma)}_{n+1} = \left[ \hat{U}^{(\gamma)}_{n} i \sin \theta - \hat{V}^{(\gamma)}_{n} \right] \Psi^{(\gamma)}_{n} + \left[ \hat{U}^{(\gamma)}_{-n} i \sin \theta - \hat{V}^{(\gamma)}_{-n} \right] \Psi^{(\gamma)}_{-n} - \]

\[
\hat{W}^{(\gamma)}_{n} \Psi^{(\gamma)}_{n-1} - \hat{W}^{(\gamma)}_{-n} \Psi^{(\gamma)}_{-(n-1)},
\]

where the recurrence constants are given by

\[
\hat{U}^{(\gamma)}_{\pm n} = \frac{1}{2} \left[ \frac{\sigma_n^{(\alpha,\beta)}}{b_n^{(\alpha,\beta)+1}} \pm \frac{\sigma_{n-1}^{(\alpha+1,\beta+1)}}{b_n^{(\alpha+1,\beta+1)+1}} \right],
\]

\[
\hat{V}^{(\gamma)}_{\pm n} = \pm \frac{1}{2} \left[ \frac{\sigma_n^{(\alpha+1,\beta+1)}}{b_n^{(\alpha+1,\beta)+1}} \pm \frac{\sigma_{n-1}^{(\alpha,\beta)}}{b_n^{(\alpha,\beta)+1}} \right],
\]

\[
\hat{W}^{(\gamma)}_{\pm n} = \frac{1}{2} \left[ \frac{\sigma_n^{(\alpha+1,\beta+1)}}{b_n^{(\alpha+1,\beta)+1}} \pm \frac{\sigma_{n-1}^{(\alpha,\beta)}}{b_n^{(\alpha,\beta)+1}} \right].
\]

Finally, putting these last two recurrences together yields

\[
D^{(\gamma)}_{n+1} \Psi^{(\gamma)}_{n+1} = \left[ A^{(\gamma)}_{n} e^{i\theta} - B^{(\gamma)}_{n} \right] \Psi^{(\gamma)}_{n} + \left[ A^{(\gamma)}_{-n} e^{-i\theta} - B^{(\gamma)}_{-n} \right] \Psi^{(\gamma)}_{-n} -
\]

\[
C^{(\gamma)}_{n} \Psi^{(\gamma)}_{n-1} + C^{(\gamma)}_{-n} \Psi^{(\gamma)}_{-(n-1)},
\]

\[
\Psi^{(\gamma)}_{n+1} = \left[ A^{(\gamma)}_{n} e^{i\theta} - B^{(\gamma)}_{n} \right] \Psi^{(\gamma)}_{n} + \left[ A^{(\gamma)}_{-n} e^{-i\theta} - B^{(\gamma)}_{-n} \right] \Psi^{(\gamma)}_{-n} +
\]

\[
C^{(\gamma)}_{n} \Psi^{(\gamma)}_{n-1} + C^{(\gamma)}_{-n} \Psi^{(\gamma)}_{-(n-1)},
\]
with the following values for the recurrence coefficients:

\[ D_{n}^{(\gamma)} = \begin{cases} 4\epsilon_{0,2}(\alpha,\beta)\sqrt{\gamma}, & n = 0, \\ 2\epsilon_{n,2}(\alpha,\beta)[\sqrt{n+\gamma} + \sqrt{n}], & n > 0. \end{cases} \]

\[ A_{\pm n}^{(\gamma)} = \begin{cases} \sqrt{2} \left[ \sqrt{\gamma + 1} \pm 1 \right], & n = 0, \\ \sqrt{n+\gamma + 1} \pm \sqrt{n+1}, & n > 0. \end{cases} \]

\[ B_{\pm n}^{(\gamma)} = \begin{cases} -\frac{\epsilon_{n,1}(\alpha,\beta)}{2\sqrt{(n+1)(n+\gamma-1)}} \left( \gamma A_{\pm n}^{(\gamma)} + \left[ 2\sqrt{n(n+\gamma)} - 1 \right] A_{\mp n}^{(\gamma)} \right), & n > 0. \end{cases} \]

\[ C_{\pm n}^{(\gamma)} = \begin{cases} 0, & n = 0, \\ -\frac{1}{A_{\pm (\gamma+1)}} \sqrt{\frac{(2\gamma+1)\gamma}{\gamma+2}}, & n = 1, \\ \frac{\gamma\epsilon_{n,0}(\alpha,\beta)}{\sqrt{(n+\gamma-1)(n+\gamma-1)}A_{n-1}^{(\gamma)}} \left[ \sqrt{(n+\gamma)^2 - 1} - \sqrt{n^2 - 1} \right], & n > 1. \end{cases} \]

**Appendix B. The Stiffness Matrix**

We assume the decay parameter \( s > \frac{1}{2} \) is given and we derive \( \alpha \) and \( \beta \) from the value \( \gamma := s - 1 \) as in Appendix A. Also, we define increments and decrements of the integer index \( k \in \mathbb{Z} \):

\[ \alpha := -\frac{1}{2}, \quad \beta := s - \frac{3}{2}, \]

\[ k^{\vee} = \text{sgn}(k) \left( |k| - 1 \right), \quad k^{\wedge} = \text{sgn}(k) \left( |k| + 1 \right), \quad n := |k| - 1. \]

We begin by noting the sparse representation of the product of \( \phi_k^{(s)} \) and \( \frac{1}{(x-\tau)^{-1}} \):

**Lemma B.1.** We have the representation:

\[ \frac{-s}{(x-\tau)^{-1}} \phi_k^{(s)} = \sum_{i \in \{\pm k^{\vee}, \pm k^{\wedge} \}} \chi_{k,i}^{(s)} \phi_i^{(s)}, \]

for some constants \( \chi_{k,i}^{(s)} \).

**Proof.** We first note that

\[ \frac{-s}{x-\tau} = -\frac{s}{2} \left[ \sin(\theta(x)) + i(1 + \cos(\theta(x))) \right], \]

after making the transformation to \( \theta(x) \). Then making the identification \( \Phi_k^{(s)} = \Psi_k^{(s-1)} \), we may use recurrence relations (A.19)-(A.20) to obtain the result. \( \square \)

A second more potent result is the sparsity result for the unweighted Wiener rational functions \( \Phi_k^{(s)}(x) \):
Lemma B.2. We have
\[
\frac{d\phi_k^{(s)}(x)}{dx} = \sum_{l \in \{\pm k^\vee, \pm k, \pm k^\wedge\}} \sigma_{k,l}^{(s)} \phi_l^{(s)},
\]
with
\[
\sigma_{k,\pm k^\vee}^{(s)} = i \text{sgn}(k) \frac{n+s}{2(2n+s)} \sqrt{\frac{(2n+1)(2n+2s-1)}{(2n+s-1)(2n+s+1)}} \times \\
\left[ \sqrt{(n+s-1)(n+s)} \pm \sqrt{n(n+1)} \right],
\]
\[
\sigma_{k,\pm k}^{(s)} = i \text{sgn}(k) \left\{ \begin{array}{ll}
\sqrt{(n+1)(n+s)}, & +k, \\
\sqrt{(n+1)(n+s)} - \frac{s(1-s)}{(2n+s)(2n+s+2)}, & -k.
\end{array} \right.
\]
\[
\sigma_{k,\pm k^\wedge}^{(s)} = \pm \frac{n+1}{n+s+1} \sigma_{k^\wedge, \pm k}.
\]

Proof. This result can be proven by brute-force calculation of the derivatives using (A.4)-(A.8) and the recurrence formula (A.1). Two critical steps are necessary: a highly nontrivial collapsing of a special arithmetic combination involving various constants in several Jacobi polynomial relations, and the very special form of the \( \theta \rightarrow x \) Jacobian for the mapping. Thus, the particular form of the mapping is critical in proving this result. \qed

Putting the two lemmas together, we have the desired sparsity result for the \( \phi_k^{(s)}(x) \) stiffness matrix:

Theorem B.3. The following equality holds for any \( s > \frac{1}{2} \):
\[
\frac{d\phi_k^{(s)}(x)}{dx} = \sum_{l \in \{\pm k^\vee, \pm k, \pm k^\wedge\}} \tau_{k,l}^{(s)} \phi_l^{(s)},
\]
where the constants \( \tau_{k,l}^{(s)} \) are equal to
\[
\tau_{k,\pm k^\vee} = i \sqrt{1 - \frac{s(s-2)}{(2n+s-1)(2n+s+1)}} \times \\
\left\{ \text{sgn}(k) \left( \sqrt{(n+s-1)(n+s)} \pm \sqrt{n(n+1)} \right) + \\
-\frac{s}{2(2n+s)} \left( \sqrt{(n+1)(n+s-1)} \pm \sqrt{n(n+s)} \right) \right\},
\]
\[
\tau_{k,k} = i \text{sgn}(k) \sqrt{(n+1)(n+s)} - \frac{is(s-1)^2}{2(2n+s)(2n+s+2)} - \frac{is}{2},
\]
\[
\tau_{k,-k} = \frac{is(s-1)}{2(2n+s)(2n+s+2)},
\]
\[
\tau_{k,\pm k^\wedge} = i \sqrt{1 - \frac{s(s-2)}{(2n+s-1)(2n+s+3)}} \times \\
\left\{ -\frac{s}{2n+s+2} \left[ \sqrt{(n+2)(n+s)} \pm \sqrt{(n+1)(n+s+1)} \right] + \\
\text{sgn}(k) \left[ \sqrt{(n+1)(n+2)} \pm \sqrt{(n+s)(n+s+1)} \right] \right\}.
\]
Clearly for $|k| = 1$ we have $\tau_{k, \pm k^\lor} = \tau_{k,0}$ so that, taking into account the different normalization constant in the definition of $\phi_k^{(s)}$ for $|k| = 1$ we have:

$$\tau_{k,0} = \frac{i}{2} \sqrt{\frac{2s-1}{2s+2}} \left\{ \text{sgn}(k) \sqrt{s} - 1 \right\}, \quad |k| = 1.$$  

And for $k = 0$:

$$\frac{d\phi_0^{(s)}}{dx} = -\frac{i}{2} \sqrt{s - \frac{1}{2}} \left[ \frac{1 + \sqrt{s}}{\sqrt{1+s}} \phi_{-1}^{(s)} + 2\sqrt{s - \frac{1}{2}} \phi_0^{(s)} + \frac{1 - \sqrt{s}}{\sqrt{1+s}} \phi_1^{(s)} \right].$$

Proof. The hard work was completed in Lemma B.2. The result of this theorem follows from a simple application of the product rule of differentiation to

$$\frac{d}{dx} \left[ \frac{2^{s/2}}{(x-i)^s} \phi_k^{(s)} \right],$$

with the appropriate use of Lemmas B.1 and B.2. \hfill $\Box$

With explicit entries for the stiffness matrix derived, we are now ready to give the proof the third property of Theorem 4.2, the spectral radius of the stiffness matrix.

Proof. We can crudely bound the entries of the stiffness matrix for $n := |k| - 1 > 1$

$$|\tau_{k, k^\lor}| \leq \frac{n}{2} + s,$$

$$|\tau_{k, -k^\lor}| \leq \frac{s}{2},$$

$$|\tau_{k, \pm k^\lor}| \leq \frac{n}{2} + s,$$

$$|\tau_{k, k}| \leq n + 2s,$$

$$|\tau_{k, -k}| \leq 1,$$

$$|\tau_{k, k^\land}| \leq \frac{n}{2} + s + \frac{1}{2},$$

$$|\tau_{k, -k^\land}| \leq \frac{s}{4}.$$

Thus we have:

$$|\tau_{k}| + |\tau_{k^\lor}| + |\tau_{-k^\lor}| + |\tau_{k^\land}| + |\tau_{-k^\land}| \leq 1 + \frac{n}{2} + s + \frac{s}{2} + \frac{n}{2} + s + \frac{1}{2} + \frac{s}{4} = n + 3s + 2.$$

An application of Gerschgorin’s Theorem then proves the result. \hfill $\Box$

References

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