OPTIMAL QUADRATURE FOR HAAR WAVELET SPACES

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Abstract. This article considers the error of the scrambled equidistribution quadrature rules in the worst-case, random-case, and average-case settings. The underlying space of integrands is a Hilbert space of multidimensional Haar wavelet series, \( H_{\text{wav}} \). The asymptotic orders of the errors are derived for the case of the scrambled \((\Lambda, t, m, s)\)-nets and \((t, s)\)-sequences. These rules are shown to have the best asymptotic convergence rates for any random quadrature rule for the space of integrands \( H_{\text{wav}} \).

1. Introduction

Base-\(b\) scrambling quadrature, proposed by Owen [Owe95], is a hybrid method of Monte Carlo and quasi-Monte Carlo methods. One starts with a low discrepancy set or sequence and randomly permutes the digits of the points. When applied to a \((t, m, s)\)-net or a \((t, s)\)-sequence [Nie92, Lar98] the result is a new net or sequence with the same \(t\)-value. Since each scrambled point is uniformly distributed over the unit cube, the resulting scrambled quadrature rule is unbiased. Moreover, the quadrature error may be estimated by performing a number of replications of the scrambled rule. Furthermore, there is a central limit theorem for \((0, m, s)\)-nets as \(m\) tends to infinity [Loh01].

The purpose of this article is to show that scrambled net quadrature rules are optimal for spaces of functions that are multidimensional Haar wavelet series. This means that the asymptotic convergence rate for scrambled net quadrature is the highest possible for any quadrature rule. This result holds for worst-case, random-case and average-case error analyses, and the asymptotic convergence rates are obtained explicitly.

The integration problem studied here is integration over the \(s\)-dimensional unit cube:

\[
I(f) = \int_{[0,1]^s} f(x) \, dx.
\]
Quadrature rules to approximate this integral take the form:

\[ Q(f; P, \{ w_i \}) = \sum_{i=1}^{n} w_i f(x_i) \]

for some set of nodes \( P = \{ x_1, \ldots, x_n \} \subset [0,1]^d \) and some set of weights \( \{ w_i \} = \{ w_1, \ldots, w_n \} \). Quasi-Monte Carlo quadrature methods choose \( P \) to be a set of points evenly distributed over the integration domain and \( w_i = n^{-1} \) for all \( i \).

The quality of a quadrature rule can be assessed by a worst-case, random-case, or average-case error analysis [HW01]. Let \( \mathcal{H} \) denote some separable Hilbert space of measurable functions, and let \( B \) be the unit ball in \( \mathcal{H} \), i.e., \( B = \{ f \in \mathcal{H} : \| f \|_\mathcal{H} \leq 1 \} \). The quadrature error for a specific integrand in \( \mathcal{H} \) and a specific quadrature rule \( Q \) is given by \( \text{Err}(f; Q) = I(f) - Q(f; P, \{ w_i \}) \). Suppose that \( Q \) is random, i.e., the nodes, weights, and number of function evaluations are all chosen randomly. Specifically, let \( Q \) be chosen from some sample space, \( \mathcal{Q}_n \), equipped with some probability distribution, \( \mu \), where the average number of function evaluations is \( n \). (Deterministic quadrature rules are the case where \( \mathcal{Q}_n \) has a single element.) In the average case we assume that \( \mathcal{H} \) is equipped with a probability measure such that \( E \| f \|_\mathcal{H}^2 = 1 \).

The worst-case, random-case, and average-case error criteria for a given Hilbert space of integrands and type of rules are:

\[
\begin{align*}
(2a) & \quad \text{worst-case:} \quad e^w(\mathcal{H}; \mathcal{Q}_n) := \text{rms sup}_{Q \in \mathcal{Q}_n, f \in B} |\text{Err}(f; Q)|, \\
(2b) & \quad \text{random-case:} \quad e^r(\mathcal{H}; \mathcal{Q}_n) := \sup_{f \in B} \text{rms sup}_{Q \in \mathcal{Q}_n} |\text{Err}(f; Q)|, \\
(2c) & \quad \text{average-case:} \quad e^a(\mathcal{H}; \mathcal{Q}_n) := \text{rms rms sup}_{Q \in \mathcal{Q}_n, f \in \mathcal{H}} |\text{Err}(f; Q)|.
\end{align*}
\]

The operator rms means root mean square (or the respective expression with the upper integral if the function to which rms is applied is not measurable).

The optimal error criteria for a Hilbert space are defined as the infima of the above with respect to all possible quadrature rules:

\[
(3) \quad e^x(\mathcal{H}, n) := \inf_{\mathcal{Q}_n} e^x(\mathcal{H}; \mathcal{Q}_n), \quad x \in \{ w, r, a \}.
\]

A sequence of random quadrature rules \( (\mathcal{Q}_{nm})_{m=0,1,2,...} \) is said to be optimal if it has the same asymptotic order as the best possible quadrature rules. Specifically, one has worst-case, random-case, and average-case optimality, respectively, if there exists some nonzero constant \( C \) independent of \( n \) such that for all \( n = 1, 2, \ldots \)

\[
\min_{m \leq n} e^x(\mathcal{H}; \mathcal{Q}_{nm}) \leq C e^x(\mathcal{H}, n),
\]

where \( x = w, r, a \), respectively. It is possible for a sequence of quadrature rules to be optimal for one of the above criteria and not for the others.

This article is concerned with Haar wavelet spaces, \( \mathcal{H}_{\text{wav}} \), defined in Section 2, and quadrature rules based on scrambled \( (\lambda, t, m, s) \)-nets and \( (t, s) \)-sequences in base \( b \) and weights all chosen to be \( w_i = 1/n \). It is shown in Corollary 11 that scrambled net rules, \( \mathcal{Q}_{sc, nm} \), are optimal for integrating functions in \( \mathcal{H}_{\text{wav}} \) in the worst, random, and average cases. This is done by showing that for some constants \( C_{up} \) and \( C_{lo} \) and some function \( h^x(n) \),

\[
(4) \quad \min_{m \leq n} e^x(\mathcal{H}_{\text{wav}}; \mathcal{Q}_{sc, nm}) \leq C_{up} h^x(n), \quad n = 1, 2, \ldots,
\]
in Theorem 8, and
\[ C_{10} h^x(n) \leq e^x(\mathcal{H}_{\text{wav}}, n), \quad n = 1, 2, \ldots, \]
in Theorems 9 and 10. Here \( x \in \{ w, r, a \} \) again. The formula for \( h^x(n) \) is found explicitly, so the asymptotic convergence rate is known. Equation (4) can be derived by applying the results of [Hic96, HH99, HY01] for worst-case and average-case error and the results of [Owe97a, Owe97b, Owe98, Yue99, YM99] for random-case error. The problem of tractability (unbounded dimension) is not considered here but has been studied by [YH01b]. Obtaining lower bounds relies on constructing test functions (or systems of such, with a measure on them, in the random-case) that cannot be integrated too well by any quadrature rule. This approach is explained in [Nov88] and [Hei93].

There are a couple of reasons for considering the Hilbert space \( \mathcal{H}_{\text{wav}} \). First, the technical details of computing the upper bounds in (4) for scrambled nets are relatively easy for this space compared to other spaces. In other words, \( \mathcal{H}_{\text{wav}} \) seems to be the natural space for studying scrambled net quadrature. Second, it is known that for many other spaces, \( \mathcal{H} \), computing \( e^w(\mathcal{H}; \mathcal{Q}_{SC,n}) \) is equivalent to computing \( e^w(\mathcal{H}_{\text{wav}}; \mathcal{Q}_{SC,n}) \) [HY01]. A similar relationship holds for average-case error analysis, but unfortunately the situation is not so simple for random-case error analysis.

2. Function Spaces Spanned by Haar Wavelets

This section defines the multidimensional spaces of Haar wavelets that are used in [Owe97a] and elsewhere. Let \( b \) be an integer greater than one that will denote the base of the Haar wavelets and the \((t, s)\)-sequences. Define the univariate basic wavelet functions
\[ \psi_\gamma(x) = b^{1/2} 1_{[b\gamma]} - b^{-1/2} 1_{[0]} , \quad \gamma = 0, 1, \ldots, b - 1, \]
where \( 1_{\bullet} \) denotes the characteristic function, and \([x]\) denotes the floor function of \( x \) or greatest integer less than or equal to \( x \). For integers \( 0 \leq \kappa \) and \( 0 \leq \tau < b^k \) the dilated and translated versions of the above functions are the following univariate wavelets:
\[ \psi_{\kappa \tau \gamma}(x) = b^{k/2} \psi_\gamma(b^{k} x - \tau) = b^{(k-1)/2}(b1_{[b^{k+1} \tau]} - 1_{[b^{k} \tau]} - 1_{[b^{k} x]}). \]

For each subset \( u \) of the coordinate axes \( \{1, \ldots, s\} \), let \( |u| \) denote the cardinality of \( u \), and let \( \bar{u} \) denote the complement of \( u \), \( \{1, \ldots, s\} - u \). For each \( r \in u \) let \( \kappa_r, \tau_r \) and \( \gamma_r \) be integers with \( \kappa_r \geq 0, 0 \leq \tau_r < b^{\kappa_r} \) and \( 0 \leq \gamma_r < b \). Define the vectors \( \kappa = (\kappa_r)_{r \in u} \), \( \tau = (\tau_r)_{r \in u} \), and \( \gamma = (\gamma_r)_{r \in u} \). Let \( \psi_{u \kappa \tau \gamma} \) be a product over \( r \in u \) of the dilated and translated wavelets, i.e.,
\[ \psi_{u \kappa \tau \gamma}(x) := \prod_{r \in u} \psi_{\kappa_r \tau_r \gamma_r}(x_r) \]
\[ = b^{(|u| - |\bar{u}|)/2} \prod_{r \in u} (b1_{[b^{\kappa_r + 1} \tau_r]} - 1_{[b^{\kappa_r} \tau_r]} - 1_{[b^{\kappa_r + 1} \tau_r + \gamma_r]}), \]
where \( |\kappa| = \sum_{r \in u} \kappa_r \). For \( u = \emptyset \) we take by convention \( \psi_{u \kappa \tau \gamma}(x) = \psi_\emptyset(x) = 1 \). The wavelets defined above are not orthogonal nor linearly independent, but they
are nearly so. As observed in [Owe97a],

\begin{equation}
\sum_{\gamma_r=0}^{b-1} \psi_{u \kappa \tau \gamma}(x) = 0, \quad \forall r \in u, \forall u \neq \emptyset, \kappa, \tau, \gamma_{u-\{r\}}
\end{equation}

\begin{equation}
\int_{[0,1]^s} \psi_{u \kappa \tau \gamma}(x) \psi_{u' \kappa' \tau' \gamma'}(x) \, dx = \delta_{u u'} \delta_{\kappa \kappa'} \delta_{\tau \tau'} \prod_{r \in u} (\delta_{\gamma_r \gamma'_r} - b^{-1}),
\end{equation}

where \( \delta \) is the Kronecker delta function.

From [Owe97a] any function \( f \in L_2([0,1]^s) \) can be represented as

\begin{equation}
f = \sum_{u, \kappa, \tau, \gamma} \hat{f}_{u \kappa \tau \gamma} \psi_{u \kappa \tau \gamma},
\end{equation}

where the coefficients \( \hat{f}_{u \kappa \tau \gamma} \) are given by

\begin{equation}
\hat{f}_{u \kappa \tau \gamma} := \int_{[0,1]^s} f(x) \psi_{u \kappa \tau \gamma}(x) \, dx.
\end{equation}

The coefficients satisfy

\[
\sum_{u, \kappa, \tau, \gamma} \hat{f}_{u \kappa \tau \gamma}^2 < \infty
\]

and the series (8) converges in \( L_2([0,1]^s) \). Note that \( \hat{f}_0 \) is simply \( I(f) \), the integral of the function. Because the wavelets are not linearly independent, they do not form a basis, and the series expression in (8) for \( f \) is not unique. By (6) and (9) it follows that

\begin{equation}
\sum_{\gamma_r=0}^{b-1} \int_{[0,1]^s} f_{u \kappa \tau \gamma} = 0, \quad \forall r \in u, \forall u \neq \emptyset, \kappa, \tau, \gamma_{u-\{r\}}.
\end{equation}

This condition ensures uniqueness of the representation in (8) in the sense that whenever

\[
f = \sum_{u, \kappa, \tau, \gamma} c_{u \kappa \tau \gamma} \psi_{u \kappa \tau \gamma},
\]

where the \( c_{u \kappa \tau \gamma} \) satisfy

\[
\sum_{u, \kappa, \tau, \gamma} c_{u \kappa \tau \gamma}^2 < \infty
\]

and

\[
\sum_{\gamma_r=0}^{b-1} c_{u \kappa \tau \gamma} = 0, \quad \forall r \in u, \forall u \neq \emptyset, \kappa, \tau, \gamma_{u-\{r\}},
\]

it follows that

\[
c_{u \kappa \tau \gamma} = \hat{f}_{u \kappa \tau \gamma} \quad \forall u, \kappa, \tau, \gamma.
\]

The space of integrands, \( H_{\text{wav}} \), considered in this article consists of all wavelet series whose coefficients converge to zero quickly enough. Let \( \omega_{u \kappa} \) denote some positive scalars defined for all \( u \) and \( \kappa \), satisfying

\begin{equation}
\sup_{u \kappa} \omega_{u \kappa} < \infty.
\end{equation}

Define

\begin{equation}
H_{\text{wav}} = \left\{ f \in L_2([0,1]^s) : \sum_{u, \kappa, \tau, \gamma} \omega_{u \kappa}^{-1} \hat{f}_{u \kappa \tau \gamma}^2 < \infty \right\},
\end{equation}
endowed with the inner product
\[ \langle f, g \rangle_{\mathcal{H}_{\text{wav}}} = \sum_{u, m, \tau, \gamma} \omega_{u_1}^{-1} \hat{f}_{u \tau \gamma} \hat{g}_{u \tau \gamma}. \]

Furthermore, define the scaled wavelets \( \psi_{u \tau \gamma}^\omega(x) = \omega_{u_1}^{1/2} \psi_{u \tau \gamma}(x) \), and let \( \hat{f}_{u \tau \gamma}^\omega = \omega_{u_1}^{-1/2} \hat{f}_{u \tau \gamma} \). Then integrands in \( \mathcal{H}_{\text{wav}} \) can be represented as
\[ f = \sum_{u, m, \tau, \gamma} \hat{f}_{u \tau \gamma}^\omega \psi_{u \tau \gamma}^\omega = \hat{f}_\omega^T \psi_\omega, \]
where \( \hat{f}_\omega \) is the column vector of the coefficients \( \hat{f}_{u \tau \gamma}^\omega \), and \( \psi_\omega \) is the column vector of the basis functions \( \psi_{u \tau \gamma}^\omega \). The inner product and norm of \( \mathcal{H}_{\text{wav}} \) can be expressed in terms of the scalar product and \( L_2 \)-norm of the coefficient vectors as follows:
\[ \langle f, g \rangle_{\mathcal{H}_{\text{wav}}} = \left\langle \hat{f}_\omega, \hat{g}_\omega \right\rangle_2 = \hat{f}_\omega^T \hat{g}_\omega, \quad \|f\|_{\mathcal{H}_{\text{wav}}} = \|\hat{f}_\omega\|_2 = (\hat{f}_\omega^T \hat{f}_\omega)^{1/2}. \]

3. Formulae for the Error Criteria Under Scrambled Quadrature

Now we turn to the problem of computing the error criteria defined in (2). The derivations in this section follow the arguments given by [HW01]. A key ingredient in this analysis is the \( \infty \times \infty \) matrix whose elements are the expectations of products of integration errors of any two wavelet functions by a randomized quadrature. Define
\[ \Lambda := E_{Q \in \mathcal{Q}_\omega} [\text{Err}(\psi_\omega; Q) \text{Err}(\psi_\omega; Q)^T]. \]

For the worst-case analysis one must assume that the quadrature rule is well-defined for any choice of sample points \( P \). This means that function evaluations must be bounded linear functionals on the Hilbert space \( \mathcal{H}_{\text{wav}} \). To ensure this it is assumed that
\[ \sum_{u, m, \tau, \gamma} \left| \psi_{u \tau \gamma}^\omega(x) \right|^2 = \sum_{u_1} (b - 1) |u_1| \omega_{u_1} < \infty, \]
which implies that the series defining the integrands in \( \mathcal{H}_{\text{wav}} \) is pointwise absolutely convergent.

For the random-case and average-case error analyses it is only necessary that the integrand be in \( L_2([0,1]^d) \), which is guaranteed by the definition of \( \mathcal{H}_{\text{wav}} \). For the average-case error analysis the integrand, \( f \), is assumed to be defined in terms of random series coefficients, \( \hat{f}_{u \tau \gamma}^\omega \). Specifically the vector of series coefficients, \( \hat{f}_\omega \), is assumed to have zero mean and covariance matrix \( \Sigma \). In this article this matrix is assumed to be block diagonal with blocks of the form
\[ \Sigma_{u \tau \gamma} = \sigma_{u_1} \otimes |u| \mathbf{H}_b, \]
where the \( b \times b \) matrix \( \mathbf{H}_b \) is defined as
\[ \mathbf{H}_b = \mathbf{I}_b - b^{-1} \mathbf{1}_b \mathbf{1}_b^T, \]
and \( \otimes |u| \mathbf{H}_b \) means taking the Kronecker product of \( \mathbf{H}_b \) with itself \( |u| \) times. This form of the covariance is consistent with condition (10). As noted in [HW01],
\[ E_{f \in \mathcal{H}_{\text{wav}}} \|f\|_{\mathcal{H}_{\text{wav}}}^2 = E_{f \in \mathcal{H}_{\text{wav}}} \langle f, f \rangle_{\mathcal{H}_{\text{wav}}} = E_{f \in \mathcal{H}_{\text{wav}}} \left\langle \hat{f}_\omega, \hat{f}_\omega \right\rangle_2 = \text{trace}(\Sigma). \]
In order to ensure that this quantity is unity, the positive scalars $\sigma_{un}$ must satisfy the following condition:

$$\sum_{un} (b - 1) |u| |p| \sigma_{un} = 1. $$

**Theorem 1.** Consider the case of random quadrature rules applied to integrands that take the form of wavelet series. The three error criteria defined in (2) are given by

$$e^{w}(\mathcal{H}_{\text{wav}}; Q_n) = \sqrt{\text{trace}(\Lambda)}, \text{ assuming (14)},$$

$$e^{r}(\mathcal{H}_{\text{wav}}; Q_n) = \sqrt{\rho(\Lambda)},$$

$$e^{s}(\mathcal{H}_{\text{wav}}; Q_n) = \sqrt{\text{trace}(\Lambda \Sigma)}, \text{ assuming (17)}.$$ 

**Proof.** To prove (18a) first define the vector $\xi_{Q,\omega} = \text{Err}(\psi_{\omega}; Q)$, and the function $\xi_{Q}(x) = \xi_{Q,\omega}^T \psi_{\omega}(x)$ for any quadrature rule $Q$. By assumption (14) it follows that the vector $\xi_{Q,\omega}$ has finite $\ell_2$-norm, and by (6) it follows that $\xi_{Q,\omega}$ satisfies (10). Thus, $\xi_{Q} \in \mathcal{H}_{\text{wav}}$. For any integrand $f = \tilde{f}_{\omega}^T \psi_{\omega} \in \mathcal{H}_{\text{wav}}$ its quadrature error may be expressed as the inner product of $f$ with $\xi_{Q}$:

$$\text{Err}(f; Q) = \tilde{f}_{\omega}^T \text{Err}(\psi_{\omega}; Q) = \langle f, \xi_{Q} \rangle_{\mathcal{H}_{\text{wav}}}. $$

Let $B_{\text{wav}}$ denote the unit ball in $\mathcal{H}_{\text{wav}}$. Then applying the Cauchy-Schwarz inequality yields a tight worst-case error bound for this quadrature rule:

$$\sup_{f \in B_{\text{wav}}} |\text{Err}(f; Q)| = \sup_{\|f\|_{\mathcal{H}_{\text{wav}}} \leq 1} |\langle f, \xi_{Q} \rangle_{\mathcal{H}_{\text{wav}}}| = \|\xi_{Q}\|_{\mathcal{H}_{\text{wav}}}. $$

The quantity $\|\xi_{Q}\|_{\mathcal{H}_{\text{wav}}}$ is often called the discrepancy. Note that the discrepancy may be expressed as the trace of a matrix:

$$\|\xi_{Q}\|_{\mathcal{H}_{\text{wav}}}^2 = \|\text{Err}(\psi_{\omega}; Q)\|^2 = \text{Err}(\psi_{\omega}; Q)^T \text{Err}(\psi_{\omega}; Q) = \text{trace}[\text{Err}(\psi_{\omega}; Q)^T \text{Err}(\psi_{\omega}; Q)^T]. $$

It follows from the definitions in (2a) and (13) that the worst-case error criterion is $e^{w}(\mathcal{H}_{\text{wav}}; Q_n) = \sqrt{\text{trace}(\Lambda)}$.

To prove (18b) it is required to compute the root mean square error in integrating an arbitrary integrand $f(x) = \tilde{f}_{\omega}^T \psi_{\omega}(x) \in \mathcal{H}_{\text{wav}}$. From the definition in (13) it follows that

$$\text{rms}_{Q \in \Omega_n} |\text{Err}(f; Q)| = \sqrt{\tilde{f}_{\omega}^T \Lambda \tilde{f}_{\omega}}. $$

The random-case error criterion is then given by

$$e^{r}(\mathcal{H}_{\text{wav}}; Q_n) = \sup_{f \in B_{\text{wav}}} \text{rms}_{Q \in \Omega_n} |\text{Err}(f; Q)| = \sup_{\|L\|_2 \leq 1} \sqrt{\tilde{f}_{\omega}^T \Lambda \tilde{f}_{\omega}}. $$

Without condition (10) this would be simply $\sqrt{\rho(\Lambda)}$, the square root of the spectral radius (largest absolute eigenvalue) of $\Lambda$.

In fact, $e^{r}(\mathcal{H}_{\text{wav}}; Q_n) = \sqrt{\rho(\Lambda)}$ even with condition (10). To prove this it is shown that the eigenvector, $\psi_{1}$, of $\Lambda$ corresponding to the maximum eigenvalue satisfies constraint (10). By definition $\Lambda$ is a symmetric, non-negative definite matrix. Without loss of generality assume that the largest eigenvalue is positive,
For any fixed $u, \kappa, \tau, r \in u$, and $\gamma_{u^{-\{1\}}}$, let $w$ be the vector whose $\mathbf{u} \kappa \tau \gamma$ elements are all one for $\gamma_r = 0, \ldots, b - 1$ and zero otherwise. By the definition of $\mathbf{A}$ and (6) it follows that this $w$ is an eigenvector of $\mathbf{A}$ with zero eigenvalue, i.e., $\mathbf{A}w = 0$. Since $\mathbf{A}$ is symmetric, $w^T \mathbf{v} = 0$, which implies that $\mathbf{v}$ satisfies (10).

For the average-case error criterion defined in (18c), note that the order of the two root mean squares can be exchanged. Starting from (19) and applying the standard formulas for expectations of quadratic forms of random variables yields

$$e^a(\mathcal{H}_{\text{wav}}; Q_n) = \sqrt{\frac{1}{E_{f \in \mathcal{H}_{\text{wav}}} \left[ \text{rms}(f; Q) \right]^2} = \sqrt{E_{f \in \mathcal{H}_{\text{wav}}} \left[ \hat{f}_{\mathbf{w}}^T \mathbf{A}^c \hat{f}_{\mathbf{w}} \right]}}$$

$$= \sqrt{\text{trace}(\mathbf{A} \Sigma)}.$$ 

The formulas for the worst-case and average-case errors are quite similar, and in fact, in a certain sense equivalent. This equivalence in a general setting is well-known (e.g., see [Rit00]). It is now applied to the problem at hand.

**Lemma 2.** Consider the case of random quadrature rules applied to integrands that take the form of wavelet series. Let $\mathcal{H}_{\text{wav}}$ be the space defined in (12) in terms of the parameters $\omega_{uk}$, and let $\tilde{\mathcal{H}}_{\text{wav}}$ be another such space defined in terms of the parameters $\tilde{\omega}_{uk} = \omega_{uu}\kappa\tau_{uk}$, where the $\kappa_{ur}$ are the parameters used to define $\Sigma$ in (15). For any kind of randomized quadrature rule $Q_n$ it follows that

$$e^w(\mathcal{H}_{\text{wav}}; Q_n) = e^a(\mathcal{H}_{\text{wav}}; Q_n),$$

provided (17) holds.

**Proof.** Given the vector of wavelets, $\psi_{\mathbf{w}}$ that defines the space $\mathcal{H}_{\text{wav}}$, and the vector $\psi_{\mathbf{w}}$ that defines the space $\tilde{\mathcal{H}}_{\text{wav}}$, it follows that $\psi_{\mathbf{w}}^{\mathbf{u} \kappa \tau \gamma} = \sigma_{\mathbf{u} \kappa \tau \gamma}^{1/2} \psi_{\mathbf{w}}^{\mathbf{u} \kappa \tau \gamma}$. Note that $\Sigma$ may be written as $\Sigma = \Sigma_{1/2}^1 \Sigma_{1/2}^1$, where $\Sigma_{1/2}^1$ is block diagonal with blocks $\Sigma_{u \kappa \tau \gamma}^{1/2} = \sigma_{\mathbf{u} \kappa \tau \gamma}^{1/2}[\otimes^u u \mathbf{H}_0]$. By condition (6) it follows that $[\otimes^u u \mathbf{H}_0]$ acting on the $\mathbf{u} \kappa \tau \gamma$ block of $\psi_{\mathbf{w}}$, leaves it unchanged. Thus, $\psi_{\mathbf{w}} = \Sigma_{1/2}^1 \psi_{\mathbf{w}}$.

Let $\tilde{\mathbf{A}}$ be the matrix associated with $\mathcal{H}_{\text{wav}}$ and $\tilde{\mathbf{A}}$ be the matrix associated with $\tilde{\mathcal{H}}_{\text{wav}}$. By the definition of these matrices, Theorem (1), and the discussion above,

$$\tilde{\mathbf{A}} = E_{Q \in Q_n} \left[ \text{Err}(\psi_{\mathbf{w}}; Q) \text{Err}(\psi_{\mathbf{w}}; Q)^T \right]$$

$$= E_{Q \in Q_n} \left[ \Sigma_{1/2}^1 \text{Err}(\psi_{\mathbf{w}}; Q) \text{Err}(\psi_{\mathbf{w}}; Q)^T \Sigma_{1/2}^1 \right] = \Sigma_{1/2}^1 \mathbf{A} \Sigma_{1/2}^1.$$ 

It then follows that

$$e^w(\mathcal{H}_{\text{wav}}; Q_n) = \sqrt{\text{trace}(\tilde{\mathbf{A}})} = \sqrt{\text{trace}(\Sigma_{1/2}^1 \mathbf{A} \Sigma_{1/2}^1)} = \sqrt{\text{trace}(\mathbf{A} \Sigma)}$$

$$= e^a(\mathcal{H}_{\text{wav}}; Q_n).$$ 

Now we consider the specific randomization where the quadrature rules are scrambled as proposed by Owen. For details of this scrambling readers are referred to [Owe95, Owe99] and the references therein. Although this scrambling is typically applied to $(t, m, s)$-nets and $(t, s)$-sequences, we do not need to assume anything about the points being scrambled. Given any $P = \{z_1, \cdots, z_n\}$ as the initial point set (before scrambling) to be used for quasi-Monte Carlo quadrature
as defined in (1), with \( w_i = 1/n \), we denote by \( Q_{\infty}(P) \) the space of scrambled quadratures.

**Lemma 3.** Let \( P = \{z_1, \cdots, z_n\} \). Then \( \Lambda^{\infty} \), the matrix \( \Lambda \) defined in (13) for scrambled quadrature, is a block diagonal matrix whose \( b[u] \times b[u] \) blocks, \( \Lambda^{\infty}_{u \tau}, \) are of the form

\[
\Lambda^{\infty}_{u \tau} = n^{-1} \omega_{u \tau} \Gamma_{u \tau}(P)[ \otimes b[u] H_b ].
\]

The gain coefficients of the set \( P, \Gamma_{u \tau}(P) \), are defined in [Owe97a] as

\[
\Gamma_{u \tau}(P) = \frac{1}{n(b-1)^u} \sum_{i,j=1}^{n} \prod_{r \in u} (b[1]_{b^r +1 z_{i_r}} = b^r z_{i_r} + 1 | b^r z_{j_r}]) \Delta_{u \tau}(P),
\]

and \( \Gamma_0 = 0 \).

**Proof.** Recall from (7) that \( \int_{[0,1]^d} \psi_{u',\tau'}^\omega(x) \, dx = \omega \delta_{u',0} \). Denote the entries of \( \Lambda^{\infty} \) by \( \lambda_{u \tau \gamma, u' \tau' \gamma'} \). Then

\[
\lambda_{u \tau \gamma, u' \tau' \gamma'} = n^{-2} \sum_{i,j=1}^{n} E \in Q_{\infty}(P) \{ [\omega \delta_{u,0} - \psi_{u \tau \gamma}^\omega(x_i)] [\omega \delta_{u',0} - \psi_{u' \tau' \gamma'}(x_j)] \}
\]

\[
= n^{-2} (\omega_{u \tau \gamma} \omega_{u' \tau' \gamma'})^{1/2} \sum_{i,j=1}^{n} E \in Q_{\infty}(P) \{ [\delta_{u,0} - \psi_{u \tau \gamma}^\omega(x_i)] [\delta_{u',0} - \psi_{u' \tau' \gamma'}(x_j)] \}.
\]

From [Owe97a, Lemmas 4-5] it is known that

\[
E \in Q_{\infty}(P) \left[ \psi_{u \tau \gamma}^\omega(x_i) \psi_{u' \tau' \gamma'}^\omega(x_j) \right] = \delta_{uu'} \delta_{\kappa \kappa'} \delta_{\tau \tau'} \Upsilon_{ij \tau \gamma \gamma'} \prod_{r \in u} (\delta_{r \gamma'}, b^{-1}),
\]

where

\[
\Upsilon_{ij \tau \gamma} = \frac{1}{(b-1)^{u[t]}} \prod_{r \in u} (b[1]_{b^r +1 z_{i_r}} = b^r z_{i_r} + 1 | b^r z_{j_r} = b^r z_{j_r}).
\]

It then follows from the definition of the gain coefficients in (21) that

\[
\lambda_{u \tau \gamma, u' \tau' \gamma'} = \begin{cases} 0, & \text{if } u \neq u' \text{ or } \kappa \neq \kappa' \text{ or } \tau \neq \tau', \\ n^{-1} \omega_{u \tau \gamma} \Gamma_{u \tau}(P) \prod_{r \in u} (\delta_{r \gamma'}, b^{-1}), & \text{otherwise}. \end{cases}
\]

This completes the proof of (20). \( \Box \)

Given the formula for \( \Lambda^{\infty} \) one can now apply Theorem 1 to compute the three error criteria for scrambled quadrature rules. These calculations are feasible because \( \Lambda^{\infty} \) is block diagonal. The simple structure of \( \Lambda^{\infty} \) results from the good match between the space of integrands, \( H_w \), and Owen’s scrambled quadrature rules, \( Q_{\infty}(P) \).

**Theorem 4.** Consider the case of Owen’s scrambled quadrature rules applied to integrands that take the form of wavelet series. The three error criteria defined in
(2) are given by

\[
e^w(\mathcal{H}_{\text{wav}}; Q_{\text{gc}}(P)) = n^{-1/2} \left[ \sum_{u \neq 0, \kappa} (b - 1) |u| \lambda_{u \kappa} \Gamma_{u \kappa}(P) \right]^{1/2},
\]
assuming (14),

\[
e^\tau(\mathcal{H}_{\text{wav}}; Q_{\text{gc}}(P)) = n^{-1/2} \left[ \max_{u, \kappa} \{\omega_{u \kappa} \Gamma_{u \kappa}(P)\} \right]^{1/2},
\]

\[
e^n(\mathcal{H}_{\text{wav}}; Q_{\text{gc}}(P)) = n^{-1/2} \left[ \sum_{u \neq 0, \kappa} (b - 1) |u| \lambda_{u \kappa} \sigma_{u \kappa} \omega_{u \kappa} \Gamma_{u \kappa}(P) \right]^{1/2},
\]
assuming (17).

**Proof.** To prove (22a) note from (18a) that \(e^w(\mathcal{H}_{\text{wav}}; Q_{\text{gc}}(P)) = \sqrt{\text{trace}(A^{sc})}\). This trace can then be computed by Lemma 3 in straightforward manner to give (22a):

\[
\text{trace}(A^{sc}) = \sum_{u \neq 0, \kappa} \text{trace}(A^{sc}_{u \kappa}) = \sum_{u \neq 0, \kappa} n^{-1} (b - 1) |u| \lambda_{u \kappa} \omega_{u \kappa} \Gamma_{u \kappa}.
\]

Formula (22b) follows directly from (18b) and Lemma 3. Formula (22c) is a consequence of (22a) and Lemma 2. \(\Box\)

4. **Asymptotic Behavior of Error Criteria for Nets**

The gain coefficients were defined to measure the performance of an arbitrary set against a simple random sample. Let \(Q_{\text{mc}, n}\) denote the sample space of the simple Monte Carlo quadrature, that is, the points in \(P \in Q_{\text{mc}, n}\) are independent and uniformly distributed over \([0, 1]^s\) random variables and the weights are \(w_i = 1/n\). The space \(Q_{\text{mc}, n}\) is invariant under scrambling. Moreover, \(E_{P \in Q_{\text{mc}, n}} \Gamma_{u \kappa}(P) = 1\) for all \(u, \kappa\). This leads to the following formulas in (23) for the error criteria for classical Monte Carlo quadrature. Note that these three error criteria all decay like proportionally to \(n^{-1/2}\), but the constants in front are largest for the worst-case and smallest for the average-case.

**Corollary 5.** Consider the case of Monte Carlo quadrature rules based on simple random samples applied to integrands that take the form of wavelet series. The three error criteria defined in (2) are:

\[
e^w(\mathcal{H}_{\text{wav}}; Q_{\text{mc}, n}) = n^{-1/2} \left[ \sum_{u \neq 0, \kappa} (b - 1) |u| \lambda_{u \kappa} \omega_{u \kappa} \right]^{1/2}, \text{ assuming (14)},
\]

\[
e^\tau(\mathcal{H}_{\text{wav}}; Q_{\text{mc}, n}) = n^{-1/2} \left[ \max_{u, \kappa} \{\omega_{u \kappa}\} \right]^{1/2},
\]

\[
e^n(\mathcal{H}_{\text{wav}}; Q_{\text{mc}, n}) = n^{-1/2} \left[ \sum_{u \neq 0, \kappa} (b - 1) |u| \lambda_{u \kappa} \sigma_{u \kappa} \omega_{u \kappa} \right]^{1/2},
\]
assuming (17).
In order for the error criteria to decay faster than \(O(n^{-1/2})\) the gain coefficients must become small or even zero, especially for larger \(\omega_{\text{un}}\). The gain coefficients for \((\lambda, 0, m, s)\)-nets and \((0, s)\)-sequences in base \(b\) can be computed precisely [Owe97a, HY01]. For general \((\lambda, t, m, s)\)-nets and \((t, s)\)-sequences with \(t > 0\) one can only obtain upper bounds on the gain coefficients [Owe97a, HY01]. In the case of digital nets and sequences, precise formulae for the gain coefficients can be obtained [NP01, YH01a], but these offer no improvement in the asymptotic decay rates over the upper bounds on the gain coefficients for general nets and sequences. The results mentioned in this paragraph are summarized in the following two lemmas.

**Lemma 6.** Let \(P\) be a \((\lambda, t, m, s)\)-net in base \(b\). Then the gain coefficients of \(P\) have the following upper bounds:

\[
\Gamma_{\text{un}}(P) = \begin{cases} 0, & |\kappa| < m - t - |u|, \\ b^t \left[ \frac{b+2^{2m+|\kappa|}|u|}{2(b-1)^{|u|}} \right], & m - t - |u| < |\kappa| \leq m - t, \\ b^t \left[ \frac{b+2^{2m+|\kappa|}|u|}{2(b-1)^{|u|}} \right], & |\kappa| > m - t. \end{cases}
\]

For the case of \(t = 0\), the gain coefficients can be computed exactly as

\[
\Gamma_{\text{un}}(P) = \Gamma_{\text{un}}^0(\lambda, m) = \frac{1}{(b-1)^{|u|}} \sum_{\alpha = 1}^{-m+|\kappa|+|u|+1} \binom{|u|}{\alpha} (-1)^{\alpha} [b|u|-\alpha - \lambda b^m-|\kappa|] \\
= \begin{cases} 0, & |\kappa| \leq m - |u|, \\ 1, & |\kappa| \geq m + 1. \end{cases}
\]

Let \(P\) contain the first \(n\) terms of a \((t, s)\)-sequence in base \(b\). Let \(n\) be expressed as

\(n = \lambda_0 + \lambda_1 b + \cdots + \lambda_q b^q = \cdots 0\lambda_q\lambda_{q-1} \cdots \lambda_1\lambda_0\) (base \(b\)),

where \(q = \lfloor \log_b(n) \rfloor\) and where \(\lambda_m \in \{0, 1, \ldots, b-1\}\). Then the gain coefficients of \(P\) have the following upper bounds:

\[
\Gamma_{\text{un}}(P) \leq \frac{b^{h q+1} b^{-h m+1+t+|\kappa|+|u|+q+1}}{n b (b-1)} \left( \frac{b+1}{b-1} \right)^{|u|+1} \leq b^{t+1} \left( \frac{b+1}{b-1} \right)^{|u|+1}.
\]

For \(t = 0\), these gain coefficients can be computed precisely as

\[
\Gamma_{\text{un}}(P) = \frac{1}{n b (b-1)} \sum_{m=0}^{q+1} b^m C_m \Gamma_{\text{un}}^0(1, m),
\]

where \(C_0 = b\lambda_0(b-\lambda_0)\),

\[
C_m = \lambda_{m-1}(\lambda_{m-1} - 1) + b\lambda_m(b - \lambda_m) - 2\lambda_{m-1}\lambda_m \\
+ 2(\lambda_{m-1} - \lambda_m)b^{-m+1} \sum_{h=0}^{m-2} \lambda_h b^h, \quad m = 1, \ldots, q,
\]

and

\[
C_{q+1} = \lambda_q(\lambda_q - 1) + 2\lambda_q b^{-q} \sum_{h=0}^{q-1} \lambda_h b^h.
\]
Lemma 7. Let \( a > 0 \) be a scalar, and let \( u \) be a subset of \( \{1, \ldots, s\} \). If \( P \) is a \((\lambda, t, m, s)\)-net in base \( b \), then

\[
    n^{-1} \sum_{\kappa} b^{-a|\kappa|} \Gamma_{u\kappa}(P) = O(n^{-1-\alpha}[\log n]^{s-1}),
\]

where \( n = \lambda b^m \). If \( P \) consists of the first \( n \) points of a \((t, s)\)-sequence in base \( b \), then

\[
    n^{-1} \sum_{\kappa} b^{-a|\kappa|} \Gamma_{u\kappa}(P) = \begin{cases} O(n^{-1-\alpha}[\log n]^{s-1}), & 0 < a < 1, \\ O(n^{-\alpha}[\log n]^s), & a = 1, \\ O(n^{-2}[\log n]^{s-1}), & a > 1, \end{cases}
\]

with constants in the \( O \)-notation not depending on \( n \) and \( P \).

According to Theorem 1 the asymptotic convergence rates of the three error criteria depend, not only on the decay rates of the gain coefficients, which are given in the lemma above, but also on the decay rates of the the \( \omega_{u\kappa} \) (and \( \sigma_{u\kappa} \) for the average case). From here on it is assumed that

\begin{equation}
    (24) \quad \omega_{u\kappa} = \theta_0 \theta_1 |u|^\alpha b^{-2\alpha|\kappa|}, \quad \sigma_{u\kappa} = \left[ 1 + \frac{\eta(b-1)}{1-b^{-2\beta}} \right]^{-\beta} \eta^{|u|b^{-2\beta|\kappa|}},
\end{equation}

where \( \theta_0, \theta_1, \eta \) are positive numbers, and \( \alpha \) is non-negative. The summability condition in (14) for worst-case analysis is equivalent to \( \alpha > 1/2 \), while condition (11) is equivalent to \( \alpha \geq 0 \). Condition (17) for the average-case analysis is equivalent to \( \beta > 1/2 \). The constants \( \alpha \) and \( \beta \) determine the asymptotic rates of decay of the three error criteria.

Let \( Q_{sc,n} \) denote Owen’s scrambled quasi-Monte Carlo quadrature using a \((\lambda, t, m, s)\)-net with \( n = \lambda b^m \) points as the node set, where \( \lambda = 0, 1, \ldots, b-1 \) and \( m = 0, 1, \ldots \).

Theorem 8. Let the space \( H_{wav} \) be defined as in Section 2 with \( \omega_{u\kappa} \) and \( \sigma_{u\kappa} \) given as in (24). Then for scrambled \((\lambda, t, m, s)\)-nets in base \( b \) the three error criteria have the following asymptotic orders:

\begin{align}
    (25a) \quad e^w(H_{wav}; Q_{sc,n}) &= O(n^{-\alpha}[\log n]^{(s-1)/2}), \quad \alpha > 1/2, \\
    (25b) \quad e^r(H_{wav}; Q_{sc,n}) &= O(n^{-\alpha-1/2}), \quad \alpha \geq 0, \\
    (25c) \quad e^a(H_{wav}; Q_{sc,n}) &= O(n^{-\alpha-\beta}[\log n]^{(s-1)/2}), \quad \alpha \geq 0, \quad \beta > 1/2.
\end{align}

Proof. This theorem is obtained by applying the formulas in Lemma 7 to those of Theorem 1 under the assumption of (24). The key step is to note that the value of \( a \) in the term \( b^{-a|\kappa|} \) determines the asymptotic rate of decay.

The formulae given in Theorem 4 have been evaluated numerically for scrambled \((\lambda, 0, m, s)\)-nets and \((0, s)\)-sequences in base \( b \), where \( s = 1, \ldots, 9 \), and the base \( b \) equal to all prime powers between \( s \) and 9 inclusively. For the \((\lambda, 0, m, s)\)-net, the sample size is \( n = \lambda b^m \) with \( m \geq 0, 1 \leq \lambda < b \), while for the \((0, s)\)-sequence, \( n \) can be any integer. The \( \omega_{u\kappa} \) and \( \sigma_{u\kappa} \) are assumed to be as follows:

\begin{equation}
    (26) \quad \omega_{u\kappa} = (6\lambda)^{-|u|b^{-2|\kappa|}}, \quad \sigma_{u\kappa} = (b^{\lambda}-1)^{-1}(1-b^{-1})^{|u|b^{-2|\kappa|}}.
\end{equation}

Conditions (14) and (17) are automatically satisfied. Some of these calculations are shown in Figures 1 and 2. The three error criteria for simple Monte Carlo quadrature as given in Theorem are also shown for comparison. The error criteria
for scrambled net or sequence quadrature are never larger than the corresponding errors of the simple Monte Carlo quadrature rules. Moreover, as the dimension $s$ increases, $e^w(\mathcal{H}_{\text{wav}}; Q_{\text{sc},n})$ increases, $e^s(\mathcal{H}_{\text{wav}}; Q_{\text{sc},n})$ stays about the same, and $e^u(\mathcal{H}_{\text{wav}}; Q_{\text{sc},n})$ decreases.

5. LOWER BOUNDS FOR THE ERROR CRITERIA

In the previous section the asymptotic convergence rates of worst-case, random-case and average-case quadrature errors were determined for scrambled net quadrature applied to integrands in $\mathcal{H}_{\text{wav}}$. In this section it is shown that these convergence rates are optimal, i.e., they are the best possible for any random quadrature rule.

**Theorem 9.** Let the space $\mathcal{H}_{\text{wav}}$ be defined as in Section 2 with $\omega_{\text{sc}}$ and $\sigma_{\text{sc}}$ given as in (24). Then the worst-case error and average-case criteria for this space as defined in (3) have the lower bounds

$$
e^w(\mathcal{H}_{\text{wav}}, n) \geq C_w n^{-\alpha} \log n \right)^{s-1}/2, \quad \alpha > 1/2,$$

$$
e^s(\mathcal{H}_{\text{wav}}, n) \geq C_a n^{-\alpha - \beta} \log n \right)^{s-1}/2, \quad \alpha > 0, \beta > 1/2,$$

for some constants $C_w$ and $C_a$ that are independent of $n$.

**Proof.** This proof proceeds by constructing a function in $\mathcal{H}_{\text{wav}}$ that cannot be integrated very well by any quadrature rule. Let $n$ be any given positive integer. Choose an integer $m$ such that

$$b^{m-1} < 2n \leq b^m.$$  \hfill (29)

For each $s$-vector $\nu = (\nu_1, \ldots, \nu_s)$ of non-negative integers and each $s$-vector $\ell = (\ell_1, \ldots, \ell_s)$ of integers $0 \leq \ell_s < b^{\nu_s}$, define the elementary interval in base $b$ as

$$B_{\nu\ell} = \prod_{r=1}^s \left[ \frac{\ell_r}{b^{\nu_r}}, \frac{\ell_r + 1}{b^{\nu_r}} \right].$$  \hfill (30)

It is clear that the set of all $B_{\nu\ell}$ is a partition of $[0, 1]^s$ into $b^{\left| \nu \right|}$ elementary intervals of volume $b^{-\left| \nu \right|}$.

Let $P = \{x_1, \ldots, x_n \} \subset [0, 1]^s$ be any given set of points. We define the function $f_\nu$ for each $\nu$ with $\left| \nu \right| = m$ by

$$f_\nu(x) = \begin{cases} 1, & \text{for all } x \in B_{\nu\ell} \text{ with } B_{\nu\ell} \cap P = \emptyset, \\ 0, & \text{otherwise}. \end{cases}$$

Then $f_\nu \in \mathcal{H}_{\text{wav}}$ and

$$I(f_\nu) = \int_{[0,1]^s} f_\nu(x) \, dx \geq b^{-m}(b^m - n) \geq \frac{1}{2}$$

since $2n \leq b^m$ from (29). Now we define $f_0 = \sum_{|\nu|=m} f_\nu$. Then $f_0 \in \mathcal{H}_{\text{wav}}$, $f_0(x_i) = 0$, for $i = 1, \ldots, n$, and

$$\text{Err}(f_0, Q) = I(f_0) = \sum_{|\nu|=m} I(f_\nu) \geq \frac{1}{2} \left( m + s - 1 \right) \geq c_1 m^{s-1},$$

where $c_1$ is a positive constant.
Figure 1. Performance of scrambled \((\lambda, 0, m, s)\)-net quadrature for the space \(H_{\text{wav}}\) with the \(\omega_{\text{wav}}\) and \(\sigma_{\text{wav}}\) specified in (26). The three reference lines are for \(e^\omega(H_{\text{wav}}; Q_{\text{mc}, n})\) (solid), \(e^r(H_{\text{wav}}; Q_{\text{mc}, n})\) (dashed) and \(e^s(H_{\text{wav}}; Q_{\text{mc}, n})\) (dot-dashed), respectively.
Figure 2. Performance of scrambled \((0,s)\)-sequence quadrature for the space \(\mathcal{H}_{\text{wav}}\) with the \(\omega_{\text{nr}}\) and \(\sigma_{\text{nr}}\) specified in (26). The three reference lines are for \(e^w(\mathcal{H}_{\text{wav}}; Q_{\text{mc},n})\) (solid), \(e^f(\mathcal{H}_{\text{wav}}; Q_{\text{mc},n})\) (dashed) and \(e^k(\mathcal{H}_{\text{wav}}; Q_{\text{mc},n})\) (dot-dashed), respectively.
To complete the proof the norm of \( f_0 \) must be estimated. For any \(|u|\)-vector \( \kappa \) and any \( s \)-vector \( \nu \), the notation \( \kappa \leq \nu \) means that \( \kappa_r \leq \nu_r \) for all \( r \in u \). The wavelet series expansion for the function \( f_\nu \) defined above is a finite sum:

\[
f_\nu(x) = \sum_{\kappa \leq \nu} \hat{f}_{\nu,\kappa}\psi_{\nu,\kappa}(x), \quad \text{where} \quad \hat{f}_{\nu,\kappa} = \int_{[0,1]} f_\nu(x)\psi_{\nu,\kappa}(x) \, dx.
\]

Note that the set of support for \( \psi_{\nu,\kappa}\) is \( B(\kappa, 0) \tau \) by definition (30), where \( (\kappa, 0) \) is the \( s \)-vector obtained by padding \( \kappa \) with zeroes for all \( r \notin u \). Since \( \|f_\nu\|_\infty = 1 \) and \( \|\psi_{\nu,\kappa}\|_\infty = \epsilon b^{-|\kappa|} / (b-1)^{|u|} \), it follows that

\[
\left| \hat{f}_{\nu,\kappa}\right| = \left| \int_{B(\kappa, 0) \tau} f_\nu(x)\psi_{\nu,\kappa}(x) \, dx \right| \\
\leq b^{-|\kappa|} (b^{-|\kappa|} - 1)^{|u|} (b-1)^{|u|} = (b-1)^{|u|} b^{-|\kappa| + |u|} / 2.
\]

Based on the above upper bound, the square norm of \( f_0 \) can now be written as follows:

\[
\|f_0\|_{\mathcal{H}_{\text{wav}}}^2 = \langle f_0, f_0 \rangle_{\mathcal{H}_{\text{wav}}} = \sum_{|\nu|=|\nu'|=m} \langle f_\nu, f_\nu' \rangle_{\mathcal{H}_{\text{wav}}}
\]

\[
\leq \sum_{|\nu|=|\nu'|=m} \sum_{u \subseteq \{1, \ldots, s\}} \sum_{\kappa \leq \min\{\nu, \nu'\}} \omega_{|\kappa|}^{-1} \left| \hat{f}_{\nu,\kappa}\right| \left| \hat{f}_{\nu',\kappa}\right|
\]

\[
\leq \sum_{u \subseteq \{1, \ldots, s\}} \sum_{\nu, \nu' \geq \kappa} \theta_0^{-1} \theta_1^{-|u|} 2^{|\kappa|} (b-1)^{|u|}
\]

\[
\leq \theta_0^{-1} \max(1, \theta_1^{-s}) (b-1)^{2s} \sum_{p=0}^{m} b^{2\alpha p} \left[ \sum_{u \subseteq \{1, \ldots, s\}} \sum_{|\nu| = p} \sum_{|\nu'| = m} \right] 1.
\]

Given \( u \) and \( \kappa \), there are \( \binom{m - p + q - 1}{s - 1} \) ways to choose a non-negative integer \( s \)-vector \( \nu - (\kappa, 0) \) with cardinality \( m - p \). Moreover, there are \( \binom{s}{q} \) subsets \( u \) with \(|u| = q\), and for each of them \( \binom{p + q - 1}{q} \) ways to construct a \( \kappa \) with \(|\kappa| = p\). Since

\[
\sum_{q=0}^{s} \binom{s}{q} \binom{p + q - 1}{q - 1} \leq 2^s (p + s - 1)^{s-1} \leq 2^s (m + s - 1)^{s-1},
\]
the upper bound on $\|f_0\|_{H_{n_{\text{wv}}}}$ may be written as

$$
\|f_0\|_{H_{n_{\text{wv}}}}^2 \leq \theta_0^{-1} \max(1, \theta_1^{-s})(b - 1)^{2s}2^s(m + s - 1)^{s-1} \\
\times \sum_{p=0}^{m} (m - p + s - 1)^{2(s-1)} b^{2\alpha p} \\
\leq \theta_0 \max(1, \theta_1^{-s})(b - 1)^{2s}2^s(m + s - 1)^{s-1} b^{2\alpha m} \\
\times \sum_{q=0}^{m} (q + s - 1)^{2(s-1)} b^{-2\alpha q} \\
\leq c_2 m^{s-1} b^{2\alpha m}.
$$

for some $c_2$ depending on $b$, $s$, and $\alpha$, but not on $m$.

Letting $f_* = f_0/\|f_0\|_{H_{n_{\text{wv}}}}$, it follows that $f_*$ lies in the unit ball, $B_{\text{wv}}$. Moreover, from (29), (31) and (32) it follows that

$$
|\text{Err}(f_*, Q)| \geq \frac{c_1 m^{s-1} b^{2\alpha m}}{\sqrt{c_2 m^{s-1} b^{2\alpha m}}} = \frac{c_1}{\sqrt{c_2}} m^{(s-1)/2} b^{-\alpha m} \geq C_w n^{-\alpha} \log n)^{(s-1)/2}
$$

for some constant $C_w$. This implies (27). Formula (28) then follows from Lemma (2). \hfill \Box

**Theorem 10.** Let the space $H_{n_{\text{wv}}}$ be defined as in Section 2 with $\omega_{u\kappa}$ and $\sigma_{u\kappa}$ given as in (24). Then the random-case error criterion for this space as defined in (3) has the lower bound

$$
e^r(H_{n_{\text{wv}}}, n) \geq C_r n^{-\alpha-1/2}, \quad \alpha \geq 0,
$$

for some constant $C_r$ that is independent of $n$.

**Proof.** It is sufficient to consider the case with $s = 1$, since integration in the space $H_{n_{\text{wv}}}$ with $s = 1$ is no harder than integration in the respective $H_{n_{\text{wv}}}$ with $s > 1$ (the one dimensional $H_{n_{\text{wv}}}$ can be identified with the subspace of the multidimensional $H_{n_{\text{wv}}}$ consisting of functions depending only on the first variable, or equivalently, of functions $f$ with $f_{uK\gamma} = 0$ for all $u$ except $u = 0$ and $u = \{1\}$). By convention the weights $\omega_{u\kappa}$ in (24) for the one dimensional case are written as $\omega_0 = \theta_0$ and $\omega_{\kappa} = \theta_0 b^{-2\alpha \kappa}$, for $\kappa = 0, 1, 2, \ldots$. The square norm of the function in $H_{n_{\text{wv}}}$ with $s = 1$ is

$$
\|f\|^2 = \hat{f}_0^2 \omega_0^{-1} + \sum_{\kappa=0}^{\infty} b^{-1} \sum_{\tau=0}^{b-1} \sum_{\gamma=0}^{b^{-1}} \hat{f}_{K\tau\gamma}^2 \omega_{\kappa}^{-1},
$$

where

$$
\hat{f}_0 = \int_0^1 f(x) \, dx \quad \text{and} \quad \hat{f}_{K\tau\gamma} = \int_0^1 f(x) \psi_{K\tau\gamma}(x) \, dx.
$$

Let $n, m$ be taken as in (29) in the proof of the previous theorem. Again define basic intervals

$$
B_{m\ell} = \left[ \frac{\ell}{b^m}, \frac{\ell + 1}{b^m} \right), \quad \ell = 0, 1, \ldots, b^m - 1,
$$
and let \( g_t = 1_{B_{nt}}(x) \) be the indicator function of \( B_{nt} \). It follows that \( I(g_t) = b^{-m} \). Define

\[
g = \sum_{\ell=0}^{b^m-1} \varepsilon g_\ell, \quad \text{where } \varepsilon \in \{1, -1\}.
\]

The norm of this function can be bounded in a similar manner as was done in the proof of the previous theorem. First the series coefficients of \( g \) are bounded as follows: \( |\hat{g}_0| \leq 1 \) and

\[
|\hat{g}_{\kappa \gamma}| = \int_{\mathbb{T}_b} g(x) \psi_{\kappa \gamma}(x) \, dx \leq b^{(\kappa-1)/2} (b-1) b^{-\kappa} = b^{-(\kappa+1)/2} (b-1).
\]

This gives an upper bound on the norm of \( g \):

\[
\|g\|_{\mathcal{M}_{\text{wav}}}^2 = \omega_0^{-1} \sum_{\kappa=0}^{m} \sum_{\tau=0}^{b^m-1} \sum_{\gamma=0}^{b^m-1} \omega_\kappa^{-1} \theta_\tau^{-1} \theta_0^{-1} (b-1) \sum_{\kappa=0}^{m} b^{2\alpha\kappa} \leq c_3^2 b^{2\alpha m},
\]

for some constant \( c_3 \) that is independent of \( m \).

Now consider the following functions \( f_\ell = g_\ell / (c_3 b^{\alpha m}) \). These \( f_\ell \) have disjoint supports and satisfy

\[
I(f_\ell) \geq c_3 b^{-(\alpha+1)m}, \quad \left\| \sum_{\ell=0}^{b^m-1} \varepsilon f_\ell \right\| \leq 1 \text{ for any } \varepsilon \in \{1, -1\}.
\]

From [Nov88, Section 2.2.4, Proposition 1] it follows that

\[
e_\ell^*(\mathcal{H}_{\text{wav}}, n) \geq c_4 b^{-(\alpha+1)m} (b^m)^{1/2} = c_4 b^{-(\alpha+1/2)m} \geq C_r n^{-\alpha-1/2}.
\]

This completes the proof of (33). \( \square \)

Theorems 8, 9, and 10 may be combined to summarize the main results of this article, namely that scrambled net quadrature is optimal for Haar wavelet series in all three settings. We use the standard notation \( a_n \approx b_n \) for non-negative sequences \( a_n \) and \( b_n \) meaning that there are constants \( c_{0l}, c_{up} > 0 \) such that \( c_{0l} b_n \leq a_n \leq c_{up} b_n \).

**Corollary 11.** Let the space \( \mathcal{H}_{\text{wav}} \) be defined as in Section 2 with \( \omega_{\alpha, \kappa} \) and \( \sigma_{\alpha, \kappa} \) given as in (24). Let \( Q_{sc, n} \) denote Owen's scrambled quasi-Monte Carlo quadrature using a \((\lambda, t, m, s)\)-net with \( \lambda b^m \) points as the node set, where \( \lambda = 0, 1, \ldots, b-1 \) and \( m = 0, 1, \ldots \). Then,

\[
\min_{\lambda b^m \leq n} e^w(\mathcal{H}_{\text{wav}}, Q_{sc, \lambda b^m}) \approx e^w(\mathcal{H}_{\text{wav}}, n) \asymp n^{-\alpha} [\log n]^{(s-1)/2}, \quad \alpha > 1/2,
\]

\[
\min_{\lambda b^m \leq n} e^r(\mathcal{H}_{\text{wav}}, Q_{sc, \lambda b^m}) \approx e^r(\mathcal{H}_{\text{wav}}, n) \asymp n^{-\alpha-1/2}, \quad \alpha \geq 0,
\]

\[
\min_{\lambda b^m \leq n} e'^*(\mathcal{H}_{\text{wav}}, Q_{sc, \lambda b^m}) \approx e'^*(\mathcal{H}_{\text{wav}}, n) \asymp n^{-\alpha-\beta} [\log n]^{(s-1)/2}, \quad \alpha \geq 0, \ \beta > 1/2.
\]

**Remark.** Usually the gap between worst-case and random-case \( n \)-th minimal errors is not more than of order \( n^{1/2} \). This is the first time that the authors have encountered a situation where this gap is larger, namely, \( e^w(\mathcal{H}_{\text{wav}}, n) / e^r(\mathcal{H}_{\text{wav}}, n) \) is of the order \( n^{1/2} [\log n]^{(s-1)/2} \).
References


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