Improved Halton sequences and discrepancy bounds

Henri Faure and Christiane Lemieux

Abstract. For about fifteen years, the surprising success of quasi-Monte Carlo methods in finance has been raising questions that challenge our understanding of these methods. At the origin are numerical experiments performed with so-called GSobol' and GFauve sequences by J. Traub and his team at Columbia University, following the pioneering work of S. Tezuka in 1993 on generalizations of Niederreiter \((t, s)\)−sequences, especially with \(t = 0\) (Faure sequences). Then in the early 2000, another breakthrough was achieved by E. Atanassov, who found clever generalizations of Halton sequences by means of permutations that are even asymptotically better than Niederreiter-Xing sequences in high dimensions. Unfortunately, detailed investigations of these GHalton sequences, together with numerical experiments, show that this good asymptotic behavior is obtained at the expense of remaining terms and is not sensitive to different choices of permutations of Atanassov. As the theory fails, the reasons of the success of GHalton, as well as GFauve, must be sought elsewhere, for instance in specific selections of good scramblings by means of tailor-made permutations. In this paper, we report on our assertions above and we give some tracks to tentatively remove a part of the mystery of QMC success.

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AMS classification. Low discrepancy sequences, Quasi-Monte Carlo methods, Computational finance.

1. Introduction

This paper aims to provide some new contributions to the thorny problem of high dimensional integration. It is well known that the number of nodes \(N\) required to approximate an \(s\)-dimensional integral with accuracy \(\varepsilon\) is exponential in \(s\) [17, 27]. For instance, using a cartesian product of trapezoidal rules gives an error in \(O(N^{-\varepsilon/s})\), hence needing \(N = 10^s\) points if \(\varepsilon = 10^{-2}\). This phenomenon is named the “curse of dimensionality” and for more than fifty years it has been the subject of countless tentatives to overcome its evil spell.

In short, these tentatives can be divided in two families, both being still unsatisfactory: (1) Monte Carlo (MC) methods seem to give a solution with an approximation in \(O(1/\sqrt{N})\), but it is only a probabilistic error bound still too slow and (2) quasi-Monte Carlo (QMC) methods do not have this disadvantage, but the deterministic bounds in

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that they provide, although asymptotically good, require a number of points practically too large to be meaningful. Nevertheless special sets of nodes used in QMC methods, among the so-called low discrepancy point sets to be discussed in Section 2, give very good results in finance—far better than MC methods—where integration problems are often in several hundreds of dimensions, see [16], [20] and [27, Section 2] where a very interesting “little history” of investigations on that dilemma at Columbia University is given. How can we explain that? Leaving aside the functional aspect of this problem based on weighted spaces and tractability, as surveyed in [27], we have taken in [9] and [14] another direction with the deepening of number theoretical properties of two main families of low discrepancy sequences, namely generalized Halton and Faure sequences. In the present study, we report some surprising results of experiments confirming that our knowledge on irregularities of distribution of multidimensional sequences is still poor in spite of remarkable results obtained in the last three decades. Its main lack is that theoretical bounds ignore the fundamental importance of permutations in the behavior of scrambled multidimensional sequences. Even for the modified Halton sequences studied by Atanassov [1], we will see that good asymptotic bounds depending on a clever choice of permutations are in fact deceiving.

On the other hand, in one dimension there are precise studies [4, 7, 8] that yield families of permutations with a tight control on the discrepancy of the resulting scrambled sequences. Using such scrambled one-dimensional sequences as coordinates of generalized Halton or generalized Faure sequences gives very good results in experiments with high dimensional integrals [9, 14], as good as Sobol’ sequences with well-chosen direction numbers. Moreover, it is easy to choose which coordinates have to be scrambled or not in relation with the presumed importance of such coordinates for the problem at hand. In this way, its effective dimension could be checked experimentally by comparing the results for different, more or less numerous, scrambled coordinates.

The paper is organized as follows. Bounds for the discrepancy of Halton sequences are recalled in Section 3, with the required background definitions given in Section 2. Similar bounds but for the class of digital \((t, s)\)-sequences are given in Section 4, and allow us, in Section 5, to establish comparisons between these sequences and Halton sequences with respect to the asymptotic behavior of their discrepancy bounds. In Section 6, we focus instead on the behavior of the bounds in the non-asymptotic regime, thus unveiling the huge gap between the behavior in these two regimes, in particular with respect to the prominent example of modified Halton sequences derived from the remarkable theoretical work of Atanassov [1]. We provide in Section 7 a discussion of digit permutations and their effect on the discrepancy in these two regimes, and then conclude the paper in Section 8.
2. Definitions

2.1. Halton and Faure sequences

Halton sequences are $s$-dimensional sequences, with values in the hypercube $I^s = [0, 1)^s$. They are obtained using one-dimensional van der Corput sequences in base $b$ for each coordinate, but in order to have good distribution properties, it is necessary to choose pairwise coprime bases, the simplest choice being to take the first $s$ primes. So, let us first recall the definition of a van der Corput sequence in base $b$, denoted $S_b$:

$$S_b(n) = \sum_{r=0}^{\infty} \frac{a_r(n)}{b^r + 1}, \quad (2.1)$$

where $a_r(n)$ is the $r$-th digit of the $b$-adic expansion of

$$n - 1 = \sum_{r=0}^{\infty} a_r(n) b^r.$$

An $s$-dimensional Halton sequence [11] $X_1, X_2, \ldots$ in $I^s$ is defined as

$$X_n = (S_{b_1}(n), \ldots, S_{b_s}(n)), \quad (2.2)$$

where the $b_j$'s, for $j = 1, \ldots, s$, are pairwise coprime.

A one-dimensional generalized van der Corput sequence [4] is obtained first by choosing a sequence $\Sigma = (\sigma_r)_{r \geq 0}$ of permutations of $Z_b = \{0, 1, \ldots, b-1\}$; then, the $n$th term of the sequence is defined as

$$S^\Sigma_b(n) = \sum_{r=0}^{\infty} \frac{\sigma_r(a_r(n))}{b^r + 1}, \quad (2.3)$$

If the same permutation $\sigma$ is used for all digits, (i.e., if $\sigma_r = \sigma$ for all $r \geq 0$), then we use the notation $S^\sigma_b$ to denote $S^\Sigma_b$. The van der Corput sequence in base $b$ defined in (2.1) is obtained by taking $\sigma_r = id$ for all $r \geq 0$, where $id$ stands for the identity permutation over $Z_b$.

A generalized Halton sequence [6] is defined by choosing $s$ sequences of permutations $\Sigma_j = (\sigma_{j,r})_{r \geq 0}$, $j = 1, \ldots, s$, and then by defining the $n$th point $X_n \in I^s$ of the sequence as

$$X_n = (S^\Sigma_1(n), \ldots, S^\Sigma_s(n)), \quad (2.4)$$

where the $b_j$'s are pairwise coprime bases. In applications, these $b_j$'s are usually chosen as the first $s$ prime numbers. In this case, we denote the $j$th base as $p_j$.

Another way of enriching the van der Corput sequence is by applying a linear transformation to the digits $a_0(n), a_1(n), \ldots$ before outputting a number between 0 and 1.
We call this a linearly scrambled van der Corput sequence. For a prime base $b$, it is obtained by choosing an $\infty \times \infty$ matrix $C = (C_{r,l})$ with elements in $\mathbb{Z}_b$, and then defining the $n$th term of this sequence as

$$S_b^C(n) = \sum_{r=0}^{\infty} \sum_{l=0}^{\infty} C_{r+1,l+1}a_l(n)b^{-(r+1)}.$$

A generalized Faure sequence in base $b$ (where $b$ is prime) has its $n$th point $X_n \in I^s$ defined as

$$X_n = (S_b^{C_1}, \ldots, S_b^{C_s}), \quad n \geq 1,$$

where $C_j = A_j P_j$, with $P_j$ the (upper triangular) Pascal matrix in $\mathbb{Z}_b$ raised to the power $j-1$, and $A_j$ a nonsingular lower triangular matrix [24]. The original Faure sequences from [5] amounts to take $A_j$ as the identity matrix. We recall these definitions of $(0, s)$-sequences because we will point out some interesting connections with generalized Halton sequences in Section 7.

### 2.2. Discrepancy

Irregularities of distribution of sequences are measured with the concept of discrepancy. Various notions exist but, for short, here we only consider the so-called extreme discrepancy, which corresponds to the worst case error in the domain of complexity of multivariate problems. Let be given a point set $\mathcal{P}_N = \{X_1, \ldots, X_N\} \subseteq I^s$ and an interval $J$ of $I^s$ of the form $J = \prod_{j=1}^{s}(y_j, z_j)$, where $0 \leq y_j < z_j \leq 1$. Then the discrepancy function of $\mathcal{P}_N$ on $J$ is the difference

$$E(J; N) = A(J; N) - NV(J),$$

where $A(J; N) = \#\{n; 1 \leq n \leq N, X_n \in J\}$ is the number of points in $\mathcal{P}_N$ that fall in the subinterval $J$, and $V(J) = \prod_{j=1}^{s}(z_j - y_j)$ is the volume of $J$.

Let us denote by $J^*$ subintervals of the form $J^* = \prod_{j=1}^{s}[0, z_j)$, where $0 \leq z_j \leq 1$. Then, the star (extreme) discrepancy $D^*$ and the (extreme) discrepancy $D$ of $\mathcal{P}_N$ are defined by

$$D^*(\mathcal{P}_N) = \sup_{J^*}|E(J^*; N)| \quad \text{and} \quad D(\mathcal{P}_N) = \sup_{J}|E(J; N)|.$$

It is well known that $D^*(\mathcal{P}_N) \leq D(\mathcal{P}_N) \leq 2^s D^*(\mathcal{P}_N)$. For an infinite sequence $X$, we denote by $D(N, X)$ and $D^*(N, X)$ the discrepancies of its first $N$ points. Note that several authors have a $1/N$ factor when defining the above quantities.

A sequence satisfying $D^*(N, X) \in O((\log N)^*)$ is typically considered to be a low-discrepancy sequence (according to the present knowledge on irregularities of distribution). But the constant hidden in the $O$ notation needs to be made explicit.
to make comparisons possible across sequences. This is achieved in many papers with
an inequality of the form
\[ D^*(N, X) \leq c_s(\log N)^s + O((\log N)^{s-1}). \]
Implicit in that inequality is a bound of the form (for some integer \( N_0 \))
\[ D^*(N, X) \leq c_s(\log N)^s + d_s(\log N)^{s-1}, \text{ for all } N \geq N_0. \] (2.6)
Then a common approach to compare low-discrepancy sequences is to assume the
second term above can be neglected and to study the behavior of \( c_s \) as a function of \( s \),
i.e., people consider the asymptotic behavior of the sequence as \( N \) goes to infinity.
But in applications, the number \( N \) of sample points used is limited and seldom goes
beyond \( 10^5 \) or \( 10^6 \). On the other hand, for the second term to be neglected, we should
have, for example,
\[ d_s(\log N)^{s-1} \leq c_s(\log N)^s, \]
which holds if and only if \( \log N > (d_s/c_s) \), i.e., \( N \geq N_1 \equiv e^{d_s/c_s} \). It means that
\( N \) might have to be huge before \( c_s \) can be a meaningful way of comparing sequences,
making \( c_s \) completely useless as a quality measure for common ranges of experiments
in QMC methods. Hence, it is important to take into account the comparative values of
\( c_s \) and \( d_s \) if we want to give more realistic comparisons of low discrepancy sequences.
Moreover, in the Koksma-Hlawka inequality, for high dimensional numerical integra-
tion, dividing (2.6) by \( N \) does not help because the term \( (\log N)^s/N \) remains prominent,
regardless of how small the leading constant \( c_s \) is. Such a behavior was already
noted, especially by Morokoff and Caflisch [16] and Kocis and Whiten [12]. See also
Section 2 in the remarkable survey of Woźniakowski [27], who gives an account of the
origin of research on tractability problems in relation with the surprising performance
of QMC methods in finance.

3. Discrepancy bounds for Halton sequences

This subsection describes results that have been established since the introduction of
Halton sequences in 1960 and that deal with the discrepancy of generalized Halton
sequences. In order to avoid confusion, we will denote the original Halton sequences
as in (2.1) by \( H \), the generalized Halton sequences as in (2.4) by \( GH \) and the so-called
modified Halton sequences constructed by Atanassov as \( HA \) (see (3.7) below). Note
that these \( HA \) sequences are a special case of \( GH \) sequences.

\[ D^*(N, H) \leq \prod_{j=1}^{s} \frac{b_j - 1}{\ln b_j} (\log N)^s + O((\log N)^{s-1}). \]
Next, Faure [5, 3] further improved the bound to
\[ D^*(N, H) \leq \prod_{j=1}^{s} \frac{b_j - 1}{2 \ln b_j} (\log N)^s + O((\log N)^{s-1}), \] (3.1)
rewritten with another proof by Niederreiter [17, p. 29] as (with the same \(c_s\))
\[ D^*(N, H) < \prod_{j=1}^{s} \left( \frac{b_j - 1}{2 \log b_j} \log N + \frac{b_j + 1}{2} \right) + s. \] (3.2)

Next, Faure [6] remarked that his proof in [3] extends easily to \(GH\) sequences. But the major theoretical improvement goes back to Atanassov [1, Theorem 2.1], with a completely different proof using an argument of diophantine geometry:
\[ D^*(N, H) \leq \frac{1}{s!} \prod_{j=1}^{s} \left( \frac{(b_j - 1) \ln N}{2 \ln b_j} + s \right) + \sum_{k=0}^{s-1} \frac{b_{k+1}}{k!} \prod_{j=1}^{k} \left( \frac{b_j}{2 \ln b_j} + k \right) + u, \] (3.3)
where \(u\) is 0 when all numbers \(b_j\) are odd, and
\[ u = \frac{b_j}{2(s-1)!} \prod_{1 \leq i \leq s, i \neq j} \left( \frac{(b_i - 1) \ln N}{2 \ln b_i} + s - 1 \right) \]
if \(b_j\) is the even number among them. Therefore estimate (2.6) holds with constant
\[ c_s = \frac{1}{s!} \prod_{j=1}^{s} \frac{b_j - 1}{2 \ln b_j}. \] (3.4)

By making the constant \(c_s\) smaller by a factor \(s!\), it is now going to 0 as \(s\) goes to infinity, whereas previously it was tending to infinity super-exponentially! Moreover, as for the preceding bounds, it is easy to extend this last result to generalized Halton sequences \(GH\). Therefore, in the comparisons to be made in Sections 5 and 6, we use the value of \(c_s\) given by (3.4) for \(GH\) sequences and denote it \(c_s(GH)\).

But in the same paper, when the bases are distinct primes, Atanassov was able to make the constant \(c_s\) even smaller in two cases. In the following, we deal only with the most interesting of these cases for applications. We need a definition before we can state the result.

**Definition.** Let \(p_1, \ldots, p_s\) be \(s\) distinct primes. The integers \(k_1, \ldots, k_s\) are called **admissible** for them, if \(p_j \nmid k_j\) and for each set of integers \(b_1, \ldots, b_s, p_j \nmid b_j\), there exists a set of integers \(a_1, \ldots, a_s,\) satisfying the congruences
\[ k_j^{(a_j)} \prod_{1 \leq i \leq s, i \neq j} p_i^{a_i} \equiv b_j \pmod{p_j}, \quad j = 1, \ldots, s. \] (3.5)
Finding sets of admissible integers is not a difficult task and amounts to solving integer linear systems. Many choices are possible and lists have been provided, see for instance [2] and [25].

With the help of admissible integers, Atanassov defines his modified Halton sequence $H_A$ as follows: For each coordinate $j$ ($1 \leq j \leq s$), let $T_j = (\tau_{j,r})_{r \geq 0}$ be the sequence of permutations of $\mathbb{Z}_{p_j}$ defined by
\[
\tau_{j,r}(x) = k_j^r x \pmod{p_j}, \quad r \geq 0,
\]
where $k_1, \ldots, k_s$ are admissible integers for the prime numbers $p_1, \ldots, p_s$. Then, the modified Halton sequence $H_A$ associated with admissible integers $k_1, \ldots, k_s$ is the generalized Halton sequence (as defined in (2.4)):
\[
H_A = (S_{T_{p_1}}, \ldots, S_{T_{p_s}}).
\]
(3.7)

This special choice of permutations $T_j$ leads to the best bounds found until now for generalized Halton sequences (where $K = \prod_{j=1}^{s} (p_j - 1)$):
\[
D^*(N, H_A) \leq \left( \frac{1}{s!} \prod_{j=1}^{s} \left( \frac{\log N}{K \log p_j} + s \right) \right) \cdot K^s.
\]
\[
\left( 1 + \sum_{j=1}^{s} \log p_j \prod_{j=1}^{s} \frac{p_j}{p_j - 1} \left( -1 + \prod_{j=1}^{s} (1 + \log p_j) \right) \right)
+ \sum_{k=0}^{s-1} \frac{p_k + 1}{k!} \prod_{j=1}^{k} \left( \frac{p_j}{2} \log N \log p_j + k \right)
+ \sum_{j=1}^{s} \frac{1}{(s - 1)!} \prod_{k=1}^{s} \left( \frac{p_k}{2} \log N \log p_k + s - 1 \right),
\]
(3.8)

The proof of this result is very long and difficult to follow. In the technical report [26], we checked carefully each step of the proof and provided three alternative proofs for an important intermediary result (Prop. 4.1 in [1]) whose proof contains a little inaccuracy in [1].

After rewriting, Atanassov took out the leading constant $c_s$ as in (2.6), everything else being rejected in the complementary term $O((\log N)^{s-1}) = d_s (\log N)^{s-1}$, thus getting:
\[
c_s(H_A) = \frac{1}{s!} \sum_{j=1}^{s} \log p_j \prod_{j=1}^{s} \frac{p_j (1 + \log p_j)}{(p_j - 1) \log p_j}.
\]
(3.9)

4. Discrepancy bounds for $(t, s)$—sequences

Before proceeding with comparisons between these various bounds for generalized Halton sequences, we briefly recall the main bounds corresponding to the other family...
of low discrepancy sequences, namely the \((t, s)-\)sequences. By doing so, we will be able to establish further comparisons in the next section, which hopefully will be useful to QMC practitioners.

For \((0, s)-\)sequences in prime base \(b \geq s\), in other words Faure sequences, denoted \(F\), we have an explicit bound deduced from [5, Section 6.4]: for all \(N \geq 1\)

\[
D^+(N, F) \leq \left(\frac{b-1}{2}\right)^s \frac{1}{(s-1)!} \sum_{m=0}^{n} (m + s)^{s-1},
\]

where the integer \(n\) satisfies \((\log N + \log 2)/\log b - 1 \leq n \leq (\log N + \log 2)/\log b\).

From this bound, we get after some calculations that the constant \(c_s\) in (2.6) for Faure sequences is

\[
c_s = \frac{1}{s!} \left(\frac{b-1}{2 \log b}\right)^s.
\]

The bound in (4.1) and constant \(c_s\) in (4.2) remain valid for generalized Faure sequences as defined in (2.5). In the following sections, we use that value of \(c_s\) given in (4.2) for \(c_s(GF)\).

For \((t, s)-\)sequences, e.g., Niederreiter sequences (and Sobol’ sequences in base 2), it is proved (see for instance [17]) that

\[
c_s = \frac{b^t b - 1}{s! \lfloor \frac{b}{2} \rfloor \left(\frac{b}{2\log b}\right)^s}.
\]

As for Faure sequences, this constant is also valid if the coordinates are scrambled by permutations of the digits. Note that in these constructions, the base \(b\) and the dimension \(s\) can be chosen independently from each other, but the quality parameter \(t\) depends on \(b\) and \(s\). While \(t \in O(s \log s)\) for the Sobol’ and Niederreiter sequences in a given base \(b\), \(t \in O(s)\) for the Niederreiter-Xing sequences studied in [18]. Note that Kritzer [13] recently improved the constants \(c_s\) in (4.2) and (4.3) by a factor \(1/2\) for odd \(b \geq 3\) and \(s \geq 2\), and by a factor \(1/3\) for \(b = 2\) and \(s \geq 5\) (a similar result holds for even \(b\)).

Finally, while (4.3) only gives the constant \(c_s\), Niederreiter also obtains explicit upper bounds for \((t, s)-\)sequences (see, for instance, [17, Theorem 4.12]). However, since these bounds are valid for integers \(N \geq b^t\), they are useless for usual ranges of samples in high dimensional integration, therefore we do not present them and refer the interested reader to [17].

5. Asymptotic comparisons

The results in this section come from computations done in [9]. We report them for sake of completeness and to enlighten our goal. They concern the constants \(c_s\) in the leading term of the upper bound (2.6). Such comparisons are a matter of Number
Theory competition, where the most tiny improvement in a direction can have a great importance, especially if people tried for a long time to progress in that direction. But regarding our present goal of practical implementations for MC and QMC users, asymptotic results that require astronomical and even cosmological numbers of trials to become relevant seem of little interest. Nevertheless, we think it is worthwhile to present them, because it can help clarify different points of view.

There are two possibilities for comparing the quantities \( c_s \) given in Sections 3 and 4: searching for asymptotic estimates as \( s \) grows to infinity or computing \( c_s \) for different values of the dimension \( s \).

In the first approach, it is easier to compare the logarithms, using Stirling formula and results from Number Theory. We refer to our paper [9] and Section B of its online appendix for details. The estimates are as follows, where \( c_s(\mathbf{GH}) \) is the value of the constant \( c_s \) given in (4.3), but in the special case of Niederreiter-Xing sequences, for which \( t \in O(s) \) as mentioned above:

\[
\begin{align*}
-s \log(1.5) &< \log(c_s(\mathbf{GH})) < -s \log s + O(1), \\
-s \log s &< \log(c_s(\mathbf{HA})) < -s \log s + O(s), \\
-s \log \log s &< \log(c_s(\mathbf{GF})) < -s \log \log s + O(s), \\
-s \log s &< \log(c_s(\mathbf{NX})) < -s \log s + O(s), \\
\end{align*}
\]

Therefore, from the point of view of asymptotic upper bounds, the best sequences are Halton-Atanassov and Niederreiter-Xing sequences, with order \(-s \log s\), followed by Faure \((0, s)\)–sequences, with order \(-s \log \log s\) and Halton sequences, with order \(-\log s\). Hence, this shows that modified Halton sequences \( \mathbf{HA} \) can have bounds on their discrepancy that are asymptotically as good as the Niederreiter-Xing sequences, which as mentioned before, are designed so that the quality parameter \( t \) behaves optimally as a function of \( s \). By constrast, the value of \( c_s \) given in (3.1)—so what was the best bound prior to the work of Atanassov—is such that \( \log c_s \in O(s \log s) \).

As a last comment for this first approach, we note that for Sobol' sequences, \( \log c_s \) is bounded from above by \( s \log \log s + O(s) \), whatever the direction numbers are. This fact is a first surprise since Sobol' sequences with carefully chosen direction numbers are among the best in numerical experiments.

Next, for the second approach, we give in Table 1 the values of \( c_s \) for Halton, Faure, Niederreiter-Xing and Halton-Atanassov sequences, for values of \( s \) ranging between 10 and 50. We have kept the values computed with \( D \) (instead of \( D' \)) in [9], hence the corrective factor \( 2^s \) for the leading constants in Table 1. The values \( t_s \) for the Niederreiter-Xing sequences were obtained using the database MinT [21]. That is, for each \( s \) we searched for the pair \((b, t)\) that was yielding the smallest value for \( c_s \) in (4.3) and we used the small improvement by Kritzer for (4.3).

These results show that for smaller values of \( s \), the Niederreiter-Xing sequences have the smallest constant, but around \( s = 30 \), the Halton-Atanassov sequences start to
Table 1. Values of constants in the discrepancy bounds for each sequence and different dimensions $s$.

<table>
<thead>
<tr>
<th>$s$</th>
<th>Halton $2^s c_s(GH)$</th>
<th>Faure $2^s c_s(GF)$</th>
<th>Nied-Xing $2^s c_s(NX)$</th>
<th>Halton-Atan $2^s c_s(HA)$</th>
</tr>
</thead>
<tbody>
<tr>
<td>10</td>
<td>2.18e-01</td>
<td>2.19e-01</td>
<td>4.46e-03</td>
<td>2.82e+00</td>
</tr>
<tr>
<td>20</td>
<td>2.85e-02</td>
<td>1.72e-02</td>
<td>1.75e-08</td>
<td>1.44e-07</td>
</tr>
<tr>
<td>30</td>
<td>3.53e-03</td>
<td>3.27e-05</td>
<td>1.40e-16</td>
<td>1.91e-17</td>
</tr>
<tr>
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<td>5.78e-04</td>
<td>1.20e-07</td>
<td>1.07e-24</td>
<td>6.17e-29</td>
</tr>
<tr>
<td>50</td>
<td>9.50e-05</td>
<td>1.19e-09</td>
<td>7.56e-35</td>
<td>1.34e-41</td>
</tr>
</tbody>
</table>

have a better constant, hence suggesting an advantage in favor of the Halton-Atanassov sequences inside the “$-s \log s$” order. Here is a second surprise since well chosen generalized Halton sequences and generalized Faure sequences (see Faure–Lemieux digital scramblings in [9] and [14]) perform as well as Sobol’ sequences in applications whereas Halton–Atanassov sequences as defined in [1] behave quite badly (see [25] and [9, Section 3 (2) and Appendix (Section H)].

### 6. Comparisons in the non-asymptotic regime

While we tried to understand the different steps of the proof of Theorem 2.3 [1, Section 4], which asserts that inequality (2.6) holds with leading constant (3.9), our attention focused on a quantity introduced in the final step, the constant $K = \prod_{j=1}^{s} (p_j - 1)$. This constant, which is already big (recall that $p_1, \ldots, p_s$ are distinct primes, hence at least the first $s$ primes), appears in the first line of (3.8) at the power $s$. Of course, it is compensated in the leading term because of the division by $K$, but a bit of rewriting shows that the term $Ks$, which is multiplied $s$ times, remains in the complementary term $d_s(\log N)^{s-1}$ of (2.6). This causes the bound (3.8) to be overall extremely large. This is a third surprise since nobody before paid attention to this complementary term when comparing the merits of low discrepancy sequences, neither for Halton nor for $(t, s)$–sequences.

In order to have a concrete idea of the importance of the complementary term in comparison with the leading one, we have computed the bounds (3.2) and (3.3) for generalized Halton sequences $GH$ (with the first $s$ prime numbers, excluding prime 2 in (3.3)) and the bound (3.8) for modified Halton sequences $HA$, but without the two last lines (so as to lighten the computation) as in Atanasov’s paper [1, p. 31]. We report these results in Tables 2, 3, 4. As a matter of fact, since the Koksma-Hlawka inequality involves the ratio $D^*(N, X)/N$ and since numbers occuring in the results are very huge, we have actually computed the bounds of $D^*(N, X)/N$ and $c_s(\log N)^*/N$. And finally, since the bounds of $D^*(N, X)/N$ were still too large, we
decided to show their logarithm. The results have been rounded to two decimals.

To sum up, in the first column we give the value of $c_s$, in the second one the value of $\log \left( \frac{bd \left( \frac{D^*(N)}{N} \right)}{N} \right)$, in the third one the value of $c_s \left( \log N \right)^s / N$ and in the last one the ratio $c_s \left( \log N \right)^s / \text{bound} \left( \frac{D^*(N, X)}{N} \right)$. Hence, a small ratio means the complementary term $d_s \left( \log N \right)^s - 1$ in (2.6) is largely dominating the leading term $c_s \left( \log N \right)^s$.

In light of these results, it is obvious that a bound in which $c_s$ behaves extremely well does not necessarily mean that overall the bound is better. Namely, although modified Halton sequences $HA$ have the constant $c_s$ with the best behavior, we can see that when $s$ is small, $c_s(GH)$ is actually larger than $c_s(HA)$ for generalized Halton sequences; and when $s$ is large, the good behavior of $c_s(HA)$ would require astronomical values of $N$ in order to provide smaller discrepancy bounds. In addition, it should be pointed out that except for $GH$ when $s = 5$, all the above bounds are still bigger than 1, while from the definitions it is clear that $D^*(N, X)/N \leq 1 \ldots$ These experiments dramatically show that the good behavior of $c_s$ is obtained at the expense of $d_s$ for $HA$ sequences.

### 7. Primordial importance of permutations

A remarkable characteristic of bounds on the discrepancy of Halton and $(t, s)-$sequences listed in Sections 3–4 and compared in Sections 5–6 is that they are valid for their generalizations with sequences of permutations, whatever these sequences of
<table>
<thead>
<tr>
<th>$N$</th>
<th>sequence</th>
<th>$c_{10}$</th>
<th>$\log \left( \frac{bd^\star(N)}{N} \right)$</th>
<th>$c_{10} \log(N)^{10}/N$</th>
<th>$c_{10} \log(N)^{10}/bd^\star(N)$</th>
</tr>
</thead>
<tbody>
<tr>
<td>$10^4$</td>
<td>(3.2) GH</td>
<td>7.72e+02</td>
<td>22.34</td>
<td>3.39e+08</td>
<td>6.73e-02</td>
</tr>
<tr>
<td></td>
<td>(3.3) GH</td>
<td>1.29e-03</td>
<td>11.09</td>
<td>5.66e+02</td>
<td>8.68e-03</td>
</tr>
<tr>
<td></td>
<td>(3.8) HA</td>
<td>2.76e-03</td>
<td>222.53</td>
<td>1.21e+03</td>
<td>2.74e-94</td>
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<td>$10^5$</td>
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<td>7.72e+02</td>
<td>21.79</td>
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<tr>
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<td>5.27e+02</td>
<td>1.98e-02</td>
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<tr>
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<td>2.76e-03</td>
<td>220.23</td>
<td>1.13e+03</td>
<td>2.55e-93</td>
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<tr>
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<td>(3.2) GH</td>
<td>7.72e+02</td>
<td>20.97</td>
<td>1.96e+08</td>
<td>1.53e-01</td>
</tr>
<tr>
<td></td>
<td>(3.3) GH</td>
<td>1.29e-03</td>
<td>9.12</td>
<td>3.27e+02</td>
<td>3.57e-02</td>
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<tr>
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<td>(3.8) HA</td>
<td>2.76e-03</td>
<td>217.93</td>
<td>6.98e+02</td>
<td>1.58e-92</td>
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<td>$10^{10}$</td>
<td>(3.2) GH</td>
<td>7.72e+02</td>
<td>16.16</td>
<td>3.24e+06</td>
<td>3.10e-01</td>
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<tr>
<td></td>
<td>(3.3) GH</td>
<td>1.29e-03</td>
<td>3.76</td>
<td>5.40e+00</td>
<td>1.26e-01</td>
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<tr>
<td></td>
<td>(3.8) HA</td>
<td>2.76e-03</td>
<td>208.72</td>
<td>1.15e+01</td>
<td>2.61e-90</td>
</tr>
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Table 3. Computations of bounds (3.2), (3.3) and (3.8) in dimension 10.

<table>
<thead>
<tr>
<th>$N$</th>
<th>sequence</th>
<th>$c_{50}$</th>
<th>$\log \left( \frac{bd^\star(N)}{N} \right)$</th>
<th>$c_{50} \log(N)^{50}/N$</th>
<th>$c_{50} \log(N)^{50}/bd^\star(N)$</th>
</tr>
</thead>
<tbody>
<tr>
<td>$10^4$</td>
<td>(3.2) GH</td>
<td>2.57e+45</td>
<td>225.77</td>
<td>4.20e+89</td>
<td>3.72e-09</td>
</tr>
<tr>
<td></td>
<td>(3.3) GH</td>
<td>2.49e-18</td>
<td>91.34</td>
<td>4.07e+26</td>
<td>8.73e-14</td>
</tr>
<tr>
<td></td>
<td>(3.8) HA</td>
<td>1.19e-56</td>
<td>10521</td>
<td>1.95e-12</td>
<td>0</td>
</tr>
<tr>
<td>$10^5$</td>
<td>(3.2) GH</td>
<td>2.57e+45</td>
<td>231.32</td>
<td>2.94e+93</td>
<td>1.02e-07</td>
</tr>
<tr>
<td></td>
<td>(3.3) GH</td>
<td>2.49e-18</td>
<td>95.79</td>
<td>2.85e+30</td>
<td>7.17e-12</td>
</tr>
<tr>
<td></td>
<td>(3.8) HA</td>
<td>1.19e-56</td>
<td>10519</td>
<td>1.36e-08</td>
<td>0</td>
</tr>
<tr>
<td>$10^6$</td>
<td>(3.2) GH</td>
<td>2.57e+45</td>
<td>235.79</td>
<td>2.68e+96</td>
<td>1.06e-06</td>
</tr>
<tr>
<td></td>
<td>(3.3) GH</td>
<td>2.49e-18</td>
<td>99.37</td>
<td>2.60e+33</td>
<td>1.82e-10</td>
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<tr>
<td></td>
<td>(3.8) HA</td>
<td>1.19e-56</td>
<td>10517</td>
<td>1.24e-05</td>
<td>0</td>
</tr>
<tr>
<td>$10^{10}$</td>
<td>(3.2) GH</td>
<td>2.57e+45</td>
<td>247.08</td>
<td>3.31e+103</td>
<td>1.64e-04</td>
</tr>
<tr>
<td></td>
<td>(3.3) GH</td>
<td>2.49e-18</td>
<td>108.3</td>
<td>3.21e+40</td>
<td>3.07e-07</td>
</tr>
<tr>
<td></td>
<td>(3.8) HA</td>
<td>1.19e-56</td>
<td>10507</td>
<td>1.54e+02</td>
<td>0</td>
</tr>
</tbody>
</table>

Table 4. Computations of bounds (3.2), (3.3) and (3.8) in dimension 50.
permutations are, either the identity or sophisticated ones.

As we saw in Section 3, there is an exception to this, which are the Halton–Atanassov sequences tailored with sequences of permutations obtained from admissible integers. But in this case, there still exists many possibilities for the choice of admissible integers (see our comment after (3.5)) and once again many sequences of permutations give the same bound. But there is more: in the definition of sequences $T_j$ of permutations in [1], recalled in (3.6), the first permutation for each coordinate $\tau_{j,0}$ is equal to the identity $id$ and therefore, the first digit of each coordinate is never scrambled whatever the admissible integers are.

However, there is no need to begin with $r = 0$ to define the sequences of permutations $T_j$. In [2] Atanassov and Durchova suggest, “because it is intuitively better”, to change $\tau_{j,r}$ in (3.6) for $\tau_{j,r}(x) = k_j^{r+1}x + b_{j,r} \pmod{p_j}$ with $r \geq 0$ and with arbitrary integers $b_{j,r}$ which, moreover, does not change the bound (3.8). The proof of that assertion, outlined in [2], can be extended to permutations $\tau_{j,r}(x) = k_j^{r+a_j}x + b_{j,r} \pmod{p_j}$ with arbitrary integers $a_j$ and even in a more general setting involving matrices, as discussed in [10].

Our implementations of Halton–Atanassov sequences, named AD in [9], using power $r + 1$ (instead of $r$) perform well in various experiments and are comparable to our own scrambled Halton sequences, named FL in [9], themselves comparable to GSobol’ and GFaure sequences as shown in [14]. But on the contrary, Halton–Atanassov sequences are hardly distinguishable from original Halton sequences $H$ when using power $r$. This fact has already been noted by Vandewoestyne and Cools [25] and is quite prominent in examples shown in Figures 1 and 3 where “HApowr” stands for $HA$ sequences with power $r$ in (3.6). With such examples, we hope it is now clear for the reader that using the identity for the first digit only is the cause of a dramatic damage.

The same situation prevails for implementations of GFaure sequences. That is, generalized Faure sequences whose first digit is not permuted do not perform much better than the original Faure sequences. This is illustrated on Figures 2 and 3, where GF is obtained from the first construction in [14, Section 4] and GFunperm is the version of GF where the first digit of each coordinate is not permuted, while F is the original Faure sequence.

All these observations confirm the primordial importance of permutations in the use of low discrepancy sequences for QMC methods, in spite of the theory of irregularities of distribution which, until now, provides only upper bounds that are unable to discriminate the identity among other permutations. The only exception is dimension one for which several theoretical and practical results exist showing the gap between the identity and well chosen permutations. To end this section, we give below a short excerpt of these results (from [4] and [8]) to further emphasize the crucial role of permutations for readers of the MCM community.

We only consider sequences $S^r_\sigma := S^r_\sigma$ from (2.3) where $\sigma_r = \sigma$ for all $r \geq 0$ and
we give results for $D$ only because they are easier to state, but analog formulas also exist for $D^\sigma$. For such sequences, there exists effectively computable constants $\alpha_b^\sigma$, which here depend on the permutation $\sigma$, such that for all $N \geq 1$

$$D(N, S_b^\sigma) \leq \frac{\alpha_b^\sigma}{\log b} \log N + \alpha_b^\sigma + 2 \quad \text{and} \quad \limsup_{N \to \infty} \frac{D(N, S_b^\sigma)}{\log N} = \frac{\alpha_b^\sigma}{\log b}. \quad (7.1)$$

It is important to point out that these formulas directly give both the asymptotic and the non-asymptotic regimes.

For large $b$, computing $\alpha_b^\sigma$ is a difficult task, but it has a good approximation by means of another quantity which is easy to compute: set $Z_b^\sigma = (\frac{\sigma(0)}{b}, \cdots, \frac{\sigma(b-1)}{b})$ and define $d_b^\sigma := \max_{1 \leq k \leq b} \max_{0 \leq h' < k < b} \left| E \left( \frac{h'}{b}, \frac{h}{b} \right) ; Z_b^\sigma \right|$, the so-called discrete discrepancy of the two-dimensional net $\{(\frac{k-1}{b}, Z_b^\sigma(k)) ; 1 \leq k \leq b\}$. Then $d_b^\sigma - 1 \leq \alpha_b^\sigma \leq d_b^\sigma$.

First of all, we recall formulas for the original van der Corput sequences $S_b^{id}$: for $b = 2$, $d_2^{id} = 1/2$ and $\alpha_2^{id} = 1/3$. Then for odd $b \geq 3$ (analogs exist for even $b$)

$$d_b^{id} = \frac{b - 1}{4} + \frac{b - 1}{4b} \quad \text{and} \quad \alpha_b^{id} = \frac{b - 1}{4}.$$

For instance, with the least prime base larger or equal to 360, we obtain $d_{367}^{id} = 33672/367 \approx 91.75$ and $\alpha_{367}^{id} = 366/4 = 91.5$.

Let us now introduce a new definition, which we restrict to a prime base $b$ for simplicity: a linear digit scrambling is a permutation $\pi$ of the set $Z_b$ of the form $\pi(k) = fk + g \pmod{b}$, where $f \neq 0$ and $g$ are given in $Z_b$. If $g = 0$, we obtain

**Figure 1.** 1000 first points of HApowr (left) and H (right), dimensions 39 and 40.
the so-called multipliers $f$ of our preceding papers [9], [14]. The additive factor $g$ is a translation usually called digital shift. In base 367, with one of the best multipliers, $f = 97$, we get ($\pi$ is identified to $f$) $d_{367}^{97} = 1088/367 \approx 2.96$ while $d_{367}^{id} \approx 91.75$, more than 30 times bigger!

Among all prime bases up to $b = 1301$, the best leading constant in (7.1), that is $\alpha_b^f / \log b$, is obtained with base 233 with $\alpha_{233}^{99} / \log 233 \approx 0.44$. Among the 212 prime numbers we have considered, few such constants are above 0.52, mainly for small $b$. Recall that $\alpha_2^{id} = 1/3$ and $\alpha_3^{id} = 1/2$ so that $\alpha_2^{id} / \log 2 \approx 0.48$ and $\alpha_3^{id} / \log 3 \approx 0.45$. Good multipliers in any base often give bounds close to these exact values, whereas without scrambling, we would have $\alpha_b^f / \log b = (b - 1)/4 \log b$, a quantity increasing to infinity with $b$.

8. Conclusion

In this paper we have reviewed discrepancy bounds for Halton sequences and their generalizations. We have provided numerical evidence demonstrating that the implied constant in these bounds—while being a useful tool to make comparisons in the asymptotic regime—could dramatically misrepresent the behavior of the bounds in the non-asymptotic regime. We also discussed the primordial importance of permutations, whose effect on the quality of a sequence is well understood theoretically in one dimension, but less so in higher dimensions.

In view of all this, we are seemingly left with no good answer to the question: how should one choose permutations and/or scrambling matrices to construct “good” generalized Halton or Faure sequences? On one hand, as we have seen in the previous sections, there is a huge gap between what appears to be a good sequence depending
on whether we consider the asymptotic or non-asymptotic regime; on the other hand, tight theoretical results are only known for one-dimensional sequences. In an attempt to best take into account these two realities, the approach we adopted in the two recent papers [9, 14] is to make use of the better knowledge for one-dimensional sequences as a primary criterion for building sequences, while also looking at the discrepancy of low-dimensional projections in the non-asymptotic regime as a secondary criterion. This approach automatically avoids using the identity to permute the first digit, because as we just saw the identity yields a large discrepancy in one dimension. In the case of GF sequences though, not only the identity should be avoided, but also using the same permutations for different coordinates should be avoided as much as possible, as this causes the corresponding projections to be almost the same as if the identity had been used on each of them. We are working to overcome such difficulties with the use of larger bases, around $2^s$, instead of the least prime base greater than $s$. We plan to report on this later in another paper.

To conclude the present study, we want to emphasize that like many people do, we think the superiority of QMC methods over MC for high dimensional integration, as in finance, comes from the fact that these problems involve only a reasonable set of “nevralgic” coordinates. Finding such sets (even roughly) could be helpful to remove a part of the mystery evoked by Woźniakowski in his paper [27] at the end of Sections 3, 4 and, last but not least, 10. The remarks on effective dimension given in the introduction suggest directions for future research. We will pursue our efforts in these

Figure 3. Absolute error for digital option problem (see [14] for more details) with $s = 75$: H and HApowr are indistinguishable.
directions in the near future.

Finally, recall that the great open conjecture in irregularities of distribution, i.e., the exact order for the extreme discrepancy of $s$-dimensional sequences is $(\log N)^s$, already starts with dimension $s = 2$. Hence it is quite hopeless to wait for theoretical results able to distinguish between permutations in dimensions greater than one and so empirical studies like [9] and [14] currently seem the only way to improve scramblings of quasi-random sequences.

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