

# Low discrepancy sequence

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- We wish approximate an integral  $\int_0^1 f(u)du$  using a total of  $n$  evaluations of the function  $f$ .
- Suppose that the function is smooth and I use points  $u_i = \frac{2i-1}{2n}$ ,  $i = 1, 2, \dots, n$ .
- Note :  $\int_0^{1/n} f(u)du = \frac{1}{n}f(u_1)$ .  
For a point in the interval  $u_1$

- How good is this approximation ? Using a Taylor series expansion around the point  $f(u_1)$ ,

$$f(x) = f(u_1) + f'(u_1)(x - u_1)$$

- and integrating both sides over the interval  $[0, \frac{1}{n}]$  we obtain

$$\int_0^{1/n} f(u) du \approx \frac{1}{n} f(u_1) + \frac{f'(u_1)}{4n^2}$$

- If the function  $f$  has a bounded first derivative this means that the integral over each subinterval

$$\int_{(j-1)/n}^{j/n} f(u) du = \frac{1}{n} f(u_1)$$

with an error that is less than

$$\text{constant} \times \frac{1}{n^2}$$

- Therefore the error in the sum over  $n$  such intervals is less than a constant  $\times \frac{1}{n}$ . How does this compare to a crude Monte Carlo integral ?
- We have seen that if we randomly select  $n$  uniform(0,1) points  $u_i$  and use the crude estimator

$$\frac{1}{n} \sum_{i=1}^n f(u_i)$$

- then the estimator has variance

$$\frac{\sigma^2}{n}$$

where

$$\sigma^2 = \text{var}(f(U_i))$$

and standard error

$$\frac{\sigma}{\sqrt{n}}$$

## 2 dimension

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- For large values of  $n$  notice that

$$constant \times \frac{1}{n} < \frac{\sigma}{\sqrt{n}}$$

. i.e. the error in the numerical integral is less than that in the monte carlo integral. **For large values of  $n$ , and for smooth functions  $f$  in one-dimension, numerical integration is better than Monte Carlo integration.**

- Suppose we now want to find an integral of the form

$$\int \int f(u_1, u_2) du_1 du_2$$

- where the integral is over the unit square. Again we wish to use  $n$  evaluation of the function. Suppose we use equally spaced points on a lattice in the two dimensional unit square and suppose  $n = m^2$ . If we define

$$u_i = \frac{2i-1}{2m}, \quad i = 1, 2, \dots, m$$

- the error in this approximation is less than or equal to

$$constant \times \frac{1}{m} = constant \times \frac{1}{\sqrt{n}}$$

- Note that this is same order of magnitude as the standard error of a Monte Carlo integral.
- For dimensions higher than 2, for example for evaluating an integral like

$$\int \int \int f(u_1, u_2, u_3) du_1 du_2 du_3$$

the Monte Carlo integral as measured by the order of the error term than is a numerical integral based on placement of points on a lattice.

- Equally spaced points on the line, or in space have the advantage that they fill holes efficiently (they get reasonably close to all points in the space). **One disadvantage is that I need to know in advance how many points ( $n$ ) are to be selected to that I can space them  $1/n$  apart.**  
**Is it possible to construct a sequence so that at least periodically the sequence consists of equally spaced points ?**

# Low Discrepancy Sequences

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- low-discrepancy: successive numbers are added in a position as far as possible from the other numbers.
- i.e. avoiding clustering
- the numbers generated sequentially fill in the larger “gaps” between the previous numbers of the sequence.
- In dimension 1, the van der Corput (1935) sequence in base 2, starts from zero, and is confined in the interval  $[0, 1)$ .



# Vander Corput Sequence, base b

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- The  $n$ 'th term of the van der Corput sequence, for base  $b$ , is generated as follows: - The decimal-base number  $n$  is expanded in the base  $b$ . For example,  $n = 4$  in base 2 is 100 ( $4 = 1 \times 4 + 0 \times 2 + 0 \times 1$ );
- The number in base  $b$  is reflected. In the example, 100 becomes 001;
- Map into interval  $[0,1)$ . 001 becomes 0.001 (binary decimal) corresponds to the decimal number  $1/8$ , that is  $1/8$  ( $= 0 \times (1/2) + 0 \times (1/4) + 1 \times (1/8)$ ).

# General Van der Corput

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- In general, for base  $b$  if

$$n = \sum_{j=0}^m a_j(n) b^j$$

- Van der corput base  $b(n) = \Phi_b(n) = \sum_{j=0}^m a_j(n) b^{-j-1}$

# Halton sequence

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Let  $p_1, p_2, \dots, p_m$  be pairwise prime integers. The Halton sequence is defined as the sequence of vectors :

$x_i := (\phi_{p_1}(i), \phi_{p_2}(i), \dots, \phi_{p_m}(i)), \quad i = 1, 2, \dots$  Usually one takes  $p_1, p_2, \dots, p_m$  to be the first  $m$  prime numbers.

<i>m</i>	repres base 2	first comp	repres base 3	second comp	repres base 5	third comp
1	1	$1/2$	1	$1/3$	1	$1/5$
2	10	$1/4$	2	$2/3$	2	$2/5$
3	11	$3/4$	10	$1/9$	3	$3/5$
4	100	$1/8$	11	$4/9$	4	$4/5$
5	101	$5/8$	12	$7/9$	10	$1/25$
6	110	$3/8$	20	$2/9$	11	$6/25$
7	111	$7/8$	21	$5/9$	12	$11/25$
9	1000	$1/16$	22	$8/9$	13	$16/25$
10	1001	$9/16$	100	$1/27$	14	$21/25$