

# Searching Worst Cases of a One-Variable Function Using Lattice Reduction

Damien Stehlé, Vincent Lefèvre, and Paul Zimmermann

**Abstract**—We propose a new algorithm to find worst cases for the correct rounding of a mathematical function of one variable. We first reduce this problem to the *real small value problem*—i.e., for polynomials with real coefficients. Then, we show that this second problem can be solved efficiently by extending Coppersmith’s work on the *integer small value problem*—for polynomials with integer coefficients—using lattice reduction. For floating-point numbers with a mantissa less than  $N$  and a polynomial approximation of degree  $d$ , our algorithm finds all worst cases at distance less than  $N^{\frac{d}{2d+1}}$  from a machine number in time  $O(N^{\frac{d}{2d+1}+\epsilon})$ . For  $d = 2$ , a detailed study improves on the  $O(N^{2/3+\epsilon})$  complexity from Lefèvre’s algorithm to  $O(N^{4/7+\epsilon})$ . For larger  $d$ , our algorithm can be used to check that there exist no worst cases at distance less than  $N^{-k}$  in time  $O(N^{1/2+\epsilon})$ .

**Index Terms**—Computer arithmetic, multiple precision arithmetic, special function approximations.

## 1 INTRODUCTION

THE IEEE-754 standard for binary floating-point arithmetic [11] requires that all four basic arithmetic operations (+, −, ×, ÷) and the square root are correctly rounded. For a given function, floating-point inputs for which it is difficult to guarantee correct rounding, called *worst cases*, are numbers for which the exact result—as computed in infinite precision—is near a machine number or near the middle of two consecutive machine numbers. This is the famous “Table Maker’s Dilemma” problem (TMD for short). Several authors [12], [13] have shown that, for the class of algebraic functions, such worst cases cannot be too close to a machine number or the middle of two consecutive machine numbers. Such bounds enable us to design some efficient algorithms that guarantee correct rounding for division and square root and less efficient algorithms for other algebraic functions.

However, for transcendental functions, number theoretic bounds—when they exist—are not sharp enough, which makes correct rounding much harder to implement. Defour et al. proposed in [7] to introduce different levels of quality for transcendental functions. However, an efficient implementation of Levels 1 and 2—correct rounding in the whole domain where the function is mathematically defined for Level 2 and in a subdomain for Level 1—requires first solving the TMD problem, i.e., to determine the worst cases for the given function in the given floating-point format.

Systematic work on the TMD was done by Lefèvre and Muller [15], who published worst cases for many elementary functions in double precision, over the full range for some functions. Alas, their approach is too expensive to deal with quadruple precision, which is included in the

revision of the IEEE-754 standard. Thus, the only possible approaches for higher precisions are either to guess a reasonable bound on the precision required for the hardest to round cases and to write a library computing up to that precision or to write a generic multiple-precision library. For instance, Ziv’s MathLib library does the former, where the guessed bound is 768 bits for double precision [20].

Having an efficient algorithm to find the hardest to round cases would help to replace guessed bounds—which are usually overestimated—by sharper and rigorous bounds. It would thus enable the design of very efficient libraries with correct rounding [6]. Then, there would no longer be a good reason to exclude those functions from the correct rounding requirements of the IEEE-754 standard.

Exhaustive search methods consist of finding the hardest to round cases of a given function in a given floating-point format. They give the best possible bound, but are very time-consuming. Moreover, a search for a given precision gives little knowledge for another precision. We propose here a new algorithm belonging to that class. It naturally extends Lefèvre’s algorithm [14] and is based on Coppersmith’s ideas.

Previous related work was done by Elkies, who gave in [8] a new algorithm using lattice reduction to find all rational points of small height near a plane curve; for example, his record:

$$5,853,886,516,781,223^3 - 447,884,928,428,402,042,307,918^2 \\ = 1,641,843$$

corresponds to a worst case of the function  $x^{3/2}$  for a 53-bit input and a 79-bit output; his other example:

$$2,220,422,932^3 - 283,059,965^3 - 2,218,888,517^3 = 30$$

corresponds to a worst case of  $(x^3 + y^3)^{1/3}$  in 32-bit arithmetic. Gonnet [9] also used lattice reduction to find worst cases, however, his approach seems equivalent to Lefèvre’s algorithm.

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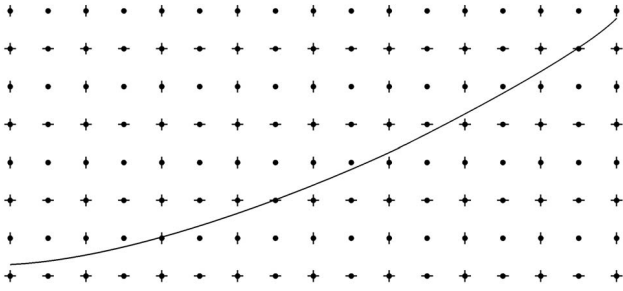


Fig. 1. A function graph and the grid of machine numbers. Worst cases correspond to grid points with a small vertical distance to the curve.

A preliminary version of this paper was presented at Arith 16 [18] and the technique described has also been used to break a cryptographic scheme proposed by Littlewood in the 1950s [19].

The paper is organized as follows: Section 2 explains in mathematical terms the problem we want to solve, recalls Lefèvre’s algorithm, and analyzes its complexity. Section 3 is a short survey on lattice reduction and Coppersmith’s work. Section 4 is the core of the paper: The new algorithm is described, proven, and analyzed. Section 5 explains in detail how the new algorithm can be tuned for the best practical choice of parameters. Section 6 presents experimental results for the  $2^x$  function, in double, double-extended, and quadruple precisions.

## 2 PRELIMINARIES

### 2.1 Definitions and Notations

In the following, we consider floating-point numbers with a mantissa of  $n$  bits. Let  $N = 2^n$ ; for instance,  $N = 2^{53}$  corresponds to double precision,  $N = 2^{64}$  corresponds to double-extended precision, and  $N = 2^{113}$  corresponds to quadruple precision. A worst case for a function  $f$  is a floating-point number  $x$  such that  $f(x)$  has  $m$  identical bits after the round bit. If all those  $m$  bits equal (respectively, differ from) the round bit,  $x$  is a worst case for the directed rounding modes (respectively, for rounding-to-nearest mode).

For the sake of simplicity, we consider directed rounding only (toward  $-\infty$ , toward  $+\infty$ , toward zero) since worst cases at precision  $n$  for all rounding modes are worst cases at precision  $n + 1$  for directed rounding. To find worst cases for directed rounding, we throw away the first  $n$  significant bits of the result mantissa and we look for runs of at least  $m$  zeros or ones in the following bits. Equivalently, a worst case of length  $m$  corresponds to  $|Nf(x) \bmod 1| < 2^{-m}$ , where  $x \bmod 1 := x - \lfloor x \rfloor$  denotes the “centered” fractional part (see Fig. 1).

We also assume that both argument  $x$  and result  $y = f(x)$  are normalized, i.e.,  $\frac{1}{2} \leq x, f(x) < 1$ . This is easy to achieve by multiplying  $x$  or  $f(x)$  by some fixed powers of 2, unless the exponent of  $f(x)$  varies a lot in the considered range. This excludes the case of numerically irregular functions like  $\sin x$  for large  $x$ . More generally, in the remainder of the paper, we assume that  $f^{(i)}(x) = O(1)$  for any  $i \geq 0$  and any  $x$  in the considered interval. This condition is verified in general, but it should be noted that

there are cases where it is not fulfilled (for example,  $\exp$  with large exponents).

Given a polynomial approximation  $P(t)$  to  $Nf(\frac{t}{N})$ —for example, a Taylor expansion—the TMD problem can be reduced to the following:

**Real Small Value Problem (Real SValP).** Given positive integers  $M$  and  $T$  and a polynomial  $P$  with real coefficients, find all integers  $|t| < T$  such that:

$$|P(t) \bmod 1| < \frac{1}{M}. \tag{1}$$

**Remark 1.** The mantissa bound  $N$  does not appear explicitly in the real SValP, however, the polynomial  $P(t)$  depends on  $N$  and so does the error made in the polynomial approximation.

**Remark 2.** If the fractional bits of the function behave randomly, we can expect  $\approx \frac{T}{M}$  worst cases. Therefore, we may assume  $T \ll M$  if we want only a few worst cases. The notation  $x \ll y$  is equivalent to  $x = O(y)$ .

### 2.2 Lefèvre’s Algorithm

Lefèvre’s algorithm [14], [15] works as follows: One considers linear approximations to the function  $f$  on small intervals. Those approximations are computed from higher order polynomial approximations on larger intervals, using an efficient scheme based on the “table of differences” method. On each small interval, worst cases are found using a modified version of the Euclidean algorithm, which gives a lower bound for  $|Nf(\frac{t}{N}) \bmod 1|$  on that interval.

Assume  $f(x_0 + x) = a_0 + a_1x + a_2x^2 + O(x^3)$  around  $x_0$ . Since we neglect terms of order two or more in  $Nf(\frac{t}{N})$ , we need  $|a_2 \frac{T^2}{N}| \ll \frac{1}{M}$  so that the error coming from the polynomial approximation does not exceed the distance  $\frac{1}{M}$ . Together with  $T \ll M$ , it follows  $T \ll N^{1/3}$ . Therefore, the complexity of Lefèvre’s algorithm is  $O(N^{2/3+\epsilon})$  since we have to consider  $\frac{N}{T} \approx N^{2/3}$  small intervals to check a complete mantissa range.

In practice, Lefèvre’s algorithm is expensive but still feasible for double precision ( $N^{2/3} \approx 4 \cdot 10^{10}$ ), near the limits of current processors for double-extended precision ( $N^{2/3} \approx 7 \cdot 10^{12}$ ), and out of reach for quadruple precision ( $N^{2/3} \approx 5 \cdot 10^{22}$ ).

## 3 LATTICE REDUCTION AND COPPERSMITH’S TECHNIQUE

In this section, we first state some basic facts about lattices—we refer to [17] for an introduction to that topic—and we explain Coppersmith’s technique, on which our algorithm is based.

### 3.1 Some Basic Facts in Lattice Reduction Theory

A lattice  $L$  is a discrete subgroup of  $\mathbb{R}^n$  or, equivalently, the set of all integer linear combinations of  $\ell \leq n$  linearly independent vectors  $\mathbf{b}_i$  over  $\mathbb{R}$ , that is:

$$L = \left\{ \sum_{i=1}^{\ell} x_i \mathbf{b}_i \mid x_i \in \mathbb{Z} \right\}.$$



**Algorithm SLZ.**
**Input:** a function  $f$ , positive integers  $N, T, M, d, \alpha$ 
**Output:** all  $t \in [-T, T]$  such that  $|Nf(\frac{t}{N}) \bmod 1| < 1/M$ 

1. Let  $P(t)$  be the Taylor expansion of  $Nf(\frac{t}{N})$  up to order  $d$ , and  $n = \frac{(\alpha+1)(d\alpha+2)}{2}$
2. Compute  $\varepsilon$  such that  $|P(t) - Nf(\frac{t}{N})| < \varepsilon$  for  $|t| \leq T$
3. Let  $M' = \lfloor \frac{1/2}{1/M+\varepsilon} \rfloor$ ,  $C = (d+1)M'$ , and<sup>3</sup>  $\tilde{P}(\tau) = \lfloor CP(T\tau) \rfloor$
4. Let  $\{e_1, \dots, e_n\} \leftarrow \{\tau^i v^j\}$  for  $0 \leq i + dj \leq d\alpha$
5. Let  $\{g_1, \dots, g_n\} \leftarrow \{(T\tau)^i (\tilde{P}(\tau) + (d+1)v)^j C^{\alpha-j}\}$  for  $0 \leq i + dj \leq d\alpha$
6. Form the  $n \times n$  integral matrix  $L$  where  $L_{k,l}$  is the coefficient of the monomial  $e_k$  in  $g_l$
7.  $V \leftarrow \text{LatticeReduce}(L)$
8. Let  $\mathbf{v}_1, \mathbf{v}_2$  be the two smallest vectors from  $V$ ,  $Q_i(T\tau, v)$  the corresponding polynomials for  $i = 1, 2$
9. If there exists  $(t, v) \in [-T, T] \times [-1, 1]$  with  $|Q_1(t, v)| \geq C^\alpha$  or  $|Q_2(t, v)| \geq C^\alpha$ , return(FAIL)
10.  $p(t) \leftarrow \text{Res}_v(Q_1(t, v), Q_2(t, v))$   
if  $p(t) = 0$  then return(FAIL)
11. For each  $t_0$  in  $\text{IntegerRoots}(p(t), [-T, T])$  do  
if  $|Nf(\frac{t_0}{N}) \bmod 1| < \frac{1}{M}$  then output  $t_0$ .

## 4.2 Correctness of the Algorithm

**Theorem 2.** *In case algorithm SLZ does not return FAIL, it behaves correctly, i.e., it outputs exactly all integers  $t \in [-T, T]$  such that  $|Nf(\frac{t}{N}) \bmod 1| < 1/M$ .*

**Proof.** Because of the final check in Step 11, we only have to verify that no worst case is missed. Suppose there is  $t_0 \in [-T, T]$  with  $|Nf(\frac{t_0}{N}) \bmod 1| < 1/M$ . From the definition of  $P$ ,  $|P(t_0) \bmod 1| < 1/M + \varepsilon \leq \frac{1}{2M'}$ . Since  $|CP(T\tau) - \tilde{P}(\tau)| \leq \frac{d+1}{2}$  for  $|\tau| \leq 1$ , by choosing  $\tau = t_0/T$ , we get  $|\tilde{P}(t_0/T) \bmod C| < d+1$ .

Whence  $\tilde{P}(t/T) + (d+1)v = 0 \bmod C$  has a root  $(t_0, v_0)$  with  $|t_0| \leq T$  and  $|v_0| < 1$ . Since  $Q_1(t, v)$  and  $Q_2(t, v)$  are integer linear combinations of the  $g_i$ s, then  $(t_0, v_0)$  is a common root of  $Q_1(t, v)$  and  $Q_2(t, v)$  modulo  $C^\alpha$  and even over the reals since  $|Q_1|, |Q_2| < C^\alpha$ . Thus,  $t_0$  is an integer root of  $\text{Res}_v(Q_1(t, v), Q_2(t, v))$  and will be found at Step 11.  $\square$

## 4.3 Choice of Parameters and Complexity Analysis

### 4.3.1 Coppersmith's Bound

Because of the use of Coppersmith's technique in our algorithm, to ensure the algorithm does not return FAIL at Step 9, the bound " $X^d Y \ll A$ " has to be verified. In our case,  $X$  corresponds to  $T$ ,  $Y$  to  $(d+1)$ , and  $A$  to  $C$ , so we get:

$$T \ll M^{1/d}.$$

### 4.3.2 Choice of the Degree $d$ with Respect to $T$

Let  $a_0, a_1, \dots$  be the Taylor coefficients of  $f$ . Since we neglect Taylor coefficients of degree  $d+1$  and greater, the error made in the approximation to  $Nf(\frac{t}{N})$  by  $P(t)$  is  $\approx a_{d+1}T^{d+1}N^{-d}$ . Since we are looking for worst cases with

$|P(t) \bmod 1| < 1/M$ , we want  $T^{d+1}N^{-d} \ll 1/M$ , i.e.,  $T^{d+1} \ll N^d/M$ .

### 4.3.3 Complexity Analysis

We have two bounds for  $T$ : The first one,  $T \ll M^{1/d}$ , comes from Coppersmith's method, the second one,  $T^{d+1} \ll N^d/M$ , comes from the accuracy of the Taylor expansion. For  $M \ll N^{\frac{d}{2d+1}}$ , Coppersmith's bound wins and implies  $T \ll M^{1/d}$ , whereas, for  $M \gg N^{\frac{d}{2d+1}}$ , Taylor's bound gives  $T^{d+1} \ll N^d/M$ . The largest bound for  $T$  is obtained for  $M \approx N^{\frac{d}{2d+1}}$ , with  $T \ll N^{\frac{d}{2d+1}}$ . For  $d = 1$ , we find the constraint  $T \ll N^{1/3}$  from Lefèvre's method; for  $d = 2$ , this gives  $T \ll N^{2/5}$  with  $M \approx N^{4/5}$ ; for  $d = 3$ , this gives  $T \ll N^{3/7}$  with  $M \approx N^{9/7}$ . With  $M \approx N^k$ , we get a best possible interval length  $T \approx N^{1/2-1/(8k)+o(1/k)}$ .

### 4.3.4 Working Precision

In Step 1, we can use floating-point coefficients in the Taylor expansion  $P(t)$  instead of symbolic coefficients as long as it introduces no error in Step 3 while computing  $\tilde{P}(\tau)$ . Let  $a_i$  be the  $i$ th Taylor coefficient of  $f$ . Then, to get  $\tilde{P}(\tau)$  correct at Step 3, the error on  $CN(T/N)^i a_i$  must be less than  $1/2$ , thus the error on  $a_i$  must be less than  $1/(2CN)(N/T)^i$ . Since  $N \geq T$ , it thus suffices to compute  $a_i$  with  $\log_2(2CN) \leq 2n$  bits after the binary point.

**Remark 3.** When searching worst cases with  $M \ll N$ , degree-2 approximations suffice. Indeed,  $N^{1-d}T^d \ll N^{1-d}M$  since  $T^d \ll M$  (Coppersmith's bound) and, for  $d \geq 3$ ,  $N^{1-d}T^d \ll N^{2-d} \ll 1/N \ll 1/M$ . Thus, all Taylor terms of degree  $\geq 3$  give a negligible contribution to  $Nf(\frac{t}{N})$  and the largest value of  $T$  is  $\approx N^{2/5}$ , giving a complexity  $\approx N^{3/5}$  to search a whole range of  $N/2$  values. More generally, for  $M \ll N^k$ , degree- $2k$  approximations suffice, giving a complexity  $\approx N^{\frac{2k+1}{2k+1}}$ .

### 4.3.5 About the Size of the Coefficients of $\tilde{P}$

Here, we study the size of the coefficients of the polynomial  $\tilde{P}$  obtained at Step 3 of the algorithm and used in the lattice reduction.

**Theorem 3.** *Suppose that the derivatives of  $f$  satisfy  $|f^{(i)}(x)| = O(1)$  and that  $T$  is chosen to reach Taylor's bound, i.e.,  $T^{d+1} = \Theta(N^d/M)$ . Then, the degree- $d$  polynomial  $\tilde{P}(\tau) = \tilde{p}_0 + \tilde{p}_1\tau + \dots + \tilde{p}_d\tau^d \in \mathbb{Z}[\tau]$  computed at Step 3 of the algorithm satisfies  $\tilde{p}_i = O((N/T)^{d-i+1})$  for any  $1 \leq i \leq d$ .*

**Proof.** Let  $1 \leq i \leq d$ . Since  $P(t) = p_0 + p_1t + \dots + p_d t^d$  is the degree- $d$  Taylor expansion of  $Nf(\frac{t}{N})$ , we have  $p_i = O(N^{1-i})$ .  $\tilde{P}$  is defined by

$$\tilde{P}(\tau) = \tilde{p}_0 + \tilde{p}_1\tau + \dots + \tilde{p}_d\tau^d = \lfloor CP(T\tau) \rfloor.$$

Since  $C = \Theta(M)$ , we have  $\tilde{p}_i = O(MT^i N^{1-i})$ . Because  $T$  is chosen to reach Taylor's bound, we have  $M = \Theta(N^d T^{-d-1})$ , which concludes the proof.  $\square$

This result highlights the fact that the obtained instantiations of the integer SValP are very special in the sense that the high degree coefficients of the input polynomial are far smaller than expected. As we show in the next section, this weakness of the high degree coefficients can be used in order to improve Coppersmith's bound and, therefore, to obtain a better complexity bound.

3. The notation  $\lfloor CP(T\tau) \rfloor$  means that we round to the nearest integer each coefficient of  $CP(T\tau)$ . This gives an element of  $\mathbb{Z}[\tau]$ .





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