Arithmetic Algorithms for Extended Precision Using Floating-Point Expansions

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Abstract—Many numerical problems require a higher computing precision than the one offered by standard floating-point (FP) formats. One common way of extending the precision is to represent numbers in a *multiple component* format. By using the so-called *floating-point expansions*, real numbers are represented as the unevaluated sum of standard machine precision FP numbers. This representation offers the simplicity of using directly available, hardware implemented and highly optimized, FP operations. It is used by multiple-precision libraries such as Bailey's QD or the analogue Graphics Processing Units (GPU) tuned version, GQD. In this article we briefly revisit algorithms for adding and multiplying FP expansions, then we introduce and prove new algorithms for normalizing, dividing and square rooting of FP expansions. The new method used for computing the reciprocal a^{-1} and the square root \sqrt{a} of a FP expansion *a* is based on an adapted Newton-Raphson iteration where the intermediate calculations are done using "truncated" operations (additions, multiplications) involving FP expansions. We give here a thorough error analysis showing that it allows very accurate computations. More precisely, after *q* iterations, the computed FP expansion $x = x_0 + \ldots + x_{2^q-1}$ satisfies, for the reciprocal algorithm, the relative error bound: $|(x - a^{-1})/a^{-1}| \le 2^{-2^q(p-3)-1}$ and, respectively, for the square root one: $|x - 1/\sqrt{a}| \le 2^{-2^q(p-3)-1}/\sqrt{a}$, where p > 2 is the precision of the FP representation used (p = 24 for single precision and p = 53 for double precision).

Index Terms—Floating-point arithmetic, floating-point expansions, high precision arithmetic, multiple-precision arithmetic, division, reciprocal, square root, Newton-Raphson iteration

1 INTRODUCTION

ANY numerical problems in dynamical systems or planetary orbit dynamics, such as the long-term stability of the solar system [1], finding sinks in the Henon Map [2], iterating the Lorenz attractor [3], etc., require higher precisions than the standard double precision (now called binary64 [4]). Quad or higher precision is rarely implemented in hardware, and the most common solution is to use software emulated higher precision libraries, also called arbitrary precision libraries. There are mainly two ways of representing numbers in higher precision. The first one is the multiple-digit representation: numbers are represented by a sequence of possibly high-radix digits coupled with a single exponent. An example is the representation used in GNU MPFR [5], an open-source C library, which, besides arbitrary precision, also provides correct rounding for each atomic operation. The second way is the *multiple*term representation in which a number is expressed as the unevaluated sum of several standard floating-point (FP) numbers. This sum is usually called a FP expansion. Bailey's library QD [6] uses this approach and supports double-

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double (DD) and quad-double (QD) computations, i.e., numbers are represented as the unevaluated sum of 2 or 4 standard double-precision FP numbers. The DD and QD formats and the operations implemented in that library are not compliant with the IEEE 754-2008 standard, and do not provide correctly rounded operations. However, this multiple-term representation offers the simplicity of using directly available and highly optimized hardware implemented FP operations. This makes most multiple-term algorithms straightforwardly portable to highly parallel architectures, such as GPUs. In consequence, there is a demand for algorithms for arithmetic operations with FP expansions, that are sufficiently simple yet efficient, and for which effective error bounds and thorough proofs are given. Several algorithms already exist for addition and multiplication [6], [7, Thm. 44, Chap. 14].

In this article we mainly focus on division (and hence, *reciprocal*) and square root, which are less studied in literature. For these algorithms we provide a thorough error analysis and effective error bounds. There are two classes of algorithms for performing division and square root: the socalled *digit-recurrence algorithms* [8], that generalize the paper-and-pencil method, and the algorithms based on the Newton-Raphson (NR) iteration [9], [10]. While the algorithms suggested so far for dividing expansions belong to the former class, here we will be interested in studying the possible use of the latter class: since its very fast, quadratic convergence is appealing when high precision is at stake.

Another contribution of this article is a new method for the renormalization of FP expansions. This operation ensures certain precision related requirements and is an important basic brick in most computations with FP expansions. Our renormalization procedure takes advantage of the computer's pipeline, so it is fast in practice. For the sake

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of completeness, we also briefly present a variant of addition and multiplication algorithms which we implemented, and for which we intend on providing a full error analysis in a future related article.

A preliminary version of our work concerning only the case of division was recently presented in [11].

The outline of the paper is the following: in Section 2 we recall some basic notions about FP expansions and the algorithms used for handling them. Then, in Section 3 we give the new renormalization algorithm along with the proof of correctness. In Section 4 we present methods for performing divisions, including existing algorithms based on long classical division on expansions (Section 4.1) and the Newton based method (Section 4.2), followed by the correctness proof, the error analysis and the complexity analysis. After that, in Section 5 we give a similar method for computing the square root of an expansion along with the complexity analysis of the algorithm. Finally, in Section 6 we assess the performance of our algorithms—in terms of number of FP operations and proven accuracy bounds.

2 FLOATING-POINT EXPANSIONS

A normal binary precision-p floating-point number is a number of the form

$$x = M_x \cdot 2^{e_x - p + 1},$$

with $2^{p-1} \leq |M_x| \leq 2^p - 1$, where M_x is an integer. The integer e_x is called the *exponent* of x, and $M_x \cdot 2^{-p+1}$ is called the *significand* of x. We denote accordingly to Goldberg's definition: $ulp(x) = 2^{e_x-p+1}$ [7, Chap. 2] (ulp is an acronym for *unit in the last place*). Another useful concept is that of *unit in the last significant place*: $uls(x) = ulp(x) \cdot 2^{z_x}$, where z_x is the number of trailing zeros at the end of M_x .

In order to ensure the uniqueness of the representation we need to set the first bit of the significand to 1 and adjust the exponent according to that. This is called a *normalized* representation. This is not possible if x is less than $2^{e_{min}}$, where e_{min} is the smallest allowed exponent. Such numbers are called *subnormal*, where the first bit of the significand is 0 and the exponent is the minimum representable one. The IEEE 754 standard specifies that an *underflow* exception is raised every time a subnormal number occurs and the operation is inexact.

A natural extension of the notion of DD or QD is the notion of *floating-point expansion*. The arithmetic on FP expansions was first developed by Priest [12], and in a slightly different way by Shewchuk [13]. If, starting from a set of FP inputs, we only perform exact operations, then the values we obtain are always equal to finite sums of FP numbers. Such finite sums are called expansions. A natural idea is to try to manipulate such expansions, for performing calculations that are either exact, either approximate yet very accurate.

Definition 2.1. A FP expansion u with n terms is the unevaluated sum of n FP numbers u_0, \ldots, u_{n-1} , in which all nonzero terms are ordered by magnitude (i.e., $u_i \neq 0 \Rightarrow |u_i| \ge |u_{i+1}|$). Each u_i is called a component of u.

The notion of expansion is "redundant" since a nonzero number always has more than one representation as a FP



Fig. 1. Nonoverlapping sequence by (a) Priest's scheme and (b) Shewchuk's scheme [17].

expansion. To make the concept useful in practice and easy to manipulate, we must introduce a constraint on the components: the u_i 's cannot "overlap". The notion of overlapping varies depending on the authors. We give here two very different definitions, using the above-introduced notation.

- **Definition 2.2.** (Nonoverlapping FP numbers) Assuming x and y are normal numbers with representations $M_x \cdot 2^{e_x-p+1}$ and $M_y \cdot 2^{e_y-p+1}$ (with $2^{p-1} \leq |M_x|, |M_y| \leq 2^p - 1$), they are \mathcal{P} -nonoverlapping (that is, nonoverlapping according to Priest's definition [14]) if $|e_y - e_x| \geq p$.
- **Definition 2.3.** An expansion is *P*-nonoverlapping (that is, nonoverlapping according to Priest's definition [14]) if all of its components are mutually *P*-nonoverlapping.

A visual representation of Definition 2.3, inspired from [17] can be seen in Fig. 1a. Shewchuk [13] weakens this into *nonzero-overlapping* sequences as shown in Fig. 1b (also inspired from [17]):

Definition 2.4. An expansion $u_0, u_1, \ldots, u_{n-1}$ is S-nonoverlapping (that is, nonoverlapping according to Shewchuk's definition [13]) if for all 0 < i < n, we have $e_{u_{i-1}} - e_{u_i} \ge p - z_{u_{i-1}}$.

In general, a \mathcal{P} -nonoverlapping expansion carries more information than an S-nonoverlapping one with the same number of components. In the worst case, in radix 2, an S-nonoverlapping expansion with 53 components may not contain more information than one double-precision FP number; it suffices to put one bit of information into every component.

When Priest first started developing the FP expansion arithmetic, he considered that all the computations were done in faithful FP arithmetic (see [14]), since round-tonearest rounding mode was not so common. More recently, a slightly stronger sense of nonoverlapping was introduced in 2001 by Hida et al. [6]:

- **Definition 2.5.** An expansion $u_0, u_1, \ldots, u_{n-1}$ is \mathcal{B} -nonoverlapping (that is, nonoverlapping according to Bailey's definition [6]) if for all 0 < i < n, we have $|u_i| \le \frac{1}{2} \operatorname{ulp}(u_{i-1})$.
- **Remark 2.6.** For \mathcal{P} -nonoverlapping expansions we have $|u_i| \leq (2^p 1)/2^p \mathrm{ulp}(u_{i-1})$ and for \mathcal{S} -nonoverlapping expansions $|u_i| \leq (2^p 1)/2^p \mathrm{uls}(u_{i-1})$.

Even though we presented here three different types of nonoverlapping, in what follows we will focus only on the



Fig. 2. VecSum with n terms. Each 2Sum box performs Algorithm 1, the sum is outputted to the left and the error downwards.

 \mathcal{P} and \mathcal{B} -nonoverlapping expansions, since, in general, they provide more precision per given number of terms of a FP expansion.

2.1 Error Free Transforms

Most algorithms performing arithmetic operations on expansions are based on the so-called *error-free transforms* (EFT), that make it possible to compute both the result and the rounding error of a FP addition or multiplication. This implies that in general, each such EFT, applied to two FP numbers, still returns two FP numbers. Although these algorithms use native precision operations only, they keep track of all accumulated rounding errors, ensuring that no information is lost.

We present here two EFTs that we use as basic bricks for our work. Algorithm 2Sum (Algorithm 1) computes the exact sum of two FP numbers a and b and returns the result under the form s + e, where s is the result rounded to nearest and e is the rounding error. It requires only six native FP operations (*flops*), which it was proven to be optimal in [15], if we have no information on the ordering of a and b.

A	Algorithm 1. 2Sum (a, b) .					
	$s \leftarrow \operatorname{RN}(a+b)$					
	// RN stands for performing the operation in rounding to					
	nearest mode.					
	$t \leftarrow \operatorname{RN}(s-b)$					
	$e \leftarrow \text{RN}(\text{RN}(a-t) + \text{RN}(b - \text{RN}(s-t)))$					
	return (s, e) such that $s = RN(a + b)$ and $s + e = a + b$					

There exists a similar EFT, that performs the same addition using only three native precision FP operations. This one is called *Fast2Sum* [7] and it requires the exponent of *a* to be larger than or equal to that of *b*. This condition might be difficult to check, but of course, if $|a| \ge |b|$, it will be satisfied.

For multiplying two FP numbers there exist two algorithms: Dekker's product and 2MultFMA. They compute the product of two FP numbers *a* and *b* and return the exact result as π , the result rounded to nearest plus *e*, the rounding error. The first one requires 17 flops. The most expensive part of the algorithm is the computation of the error $e = a \cdot b - \pi$, but if a fused-multiply-add (FMA [7]) instruction, that takes only one *flop*, is available, it is easily computed. This gives Algorithm 2MultFMA (Algorithm 2), that takes only two *flops*. This algorithm works providing that the product $a \cdot b$ does not overflow and $e_a + e_b \ge$ $e_{min} + p - 1$, where e_a and e_b are the exponents of a and b and e_{min} is the minimum representable exponent. If the second condition is not satisfied, the product may not be representable as the exact sum of two FP numbers (e would be below the underflow threshold).



Fig. 3. Addition of FP expansions with n and m terms. The *Merge* box performs a classic algorithm for merging two arrays and the *Renormalize* box performs Algorithm 6.

Algorithm 2. 2MultFMA (a, b).

 $\pi \leftarrow \text{RN}(a \cdot b)$ // RN stands for performing the operation in rounding to nearest mode. $e \leftarrow \text{fma}(a, b, -\pi)$ **return** (π, e) such that $\pi = \text{RN}(a \cdot b)$ and $\pi + e = a \cdot b$

These EFT can be extended to work on several inputs by chaining, resulting in the so-called distillation algorithms [16]. From these we make use of an algorithm called *VecSum* by Ogita et al. [13], [17]. VecSum, presented in Fig. 2 and Algorithm 3, is simply a chain of 2Sum that performs an EFT on *n* FP numbers.

Algorithm 3. VecSum (x_0, \ldots, x_{n-1}) .
Input: $x_0,, x_{n-1}$ FP numbers. Output: $e_0 + \cdots + e_{n-1} = x_0 + \cdots + x_{n-1}$.
$s_{n-1} \leftarrow x_{n-1}$ for $i \leftarrow n-2$ to 0 do $(s_i, e_{i+1}) \leftarrow 2Sum(x_i, s_{i+1})$ end for $e_0 \leftarrow s_0$
return e_0, \ldots, e_{n-1}

2.2 Addition and Multiplication Algorithms for Expansions

An algorithm that performs the addition of two expansions a and b with n and m terms, respectively, will return a FP expansion with at most n + m terms. Similarly, for multiplication, the product can have at most 2nm terms [12]. So-called normalization algorithms are used to render the result nonoverlapping, and this also implies a potential reduction in the number of terms.

Many variants of algorithms that compute the sum and the product of two FP expansions have been presented in the literature [6], [12], [13], [16]. Here, we only briefly present the algorithms that we used in our actual implementation. The addition is based on the merge algorithm and the multiplication is a generalization of Bailey's algorithm for DD and QD [6] and it was first presented in [2].

The addition presented in Algorithm 4 and Fig. 3 is performed by merging the two FP expansions, a, with n and b, with m terms, respectively, and normalizing the resulted array for obtaining an approximation s on r terms of the sum a + b.

Subtraction is performed simply by negating the FP terms in *b*.



Fig. 4. Multiplication of FP expansions with k terms. Each 2MultFMA box performs Algorithm 2, the product is outputted downwards and the error to the right; the VecSum box performs Algorithm 3, in which the most significant component of the sum is outputted downwards; the circled + and * signs represent standard round-to-nearest FP addition and multiplication.

Input: FP expansions $a = a_0 + \dots + a_{n-1}$; $b = b_0 + \dots + b_{m-1}$; rlength of the result. Output: FP expansion $s = s_0 + \dots + s_{r-1}$. 1: $f[0:m+n-1] \leftarrow Merge(a[0:n-1], b[0:m-1])$ 2: $s[0:r-1] \leftarrow Renorm(f[0:m+n-1], r)$

3: **return** FP expansion $s = s_0 + \cdots + s_{r-1}$.

In Fig. 4 and Algorithm 5 we present the multiplication algorithm. Although we have implemented fully customized versions, for simplicity, we give here only the "*k input* - *k output*" variant of the algorithm. We consider two expansions *a* and *b*, each with *k* terms and we compute the *k* most significant components of the product $r = a \cdot b$. In the renormalization step (line 15 of Algorithm 5), we use an extra error correction term, so we perform our "error free transformation scheme" k + 1 times.

Here, we just give an intuitive explanation of the multiplication algorithm. Let $\varepsilon = \frac{1}{2} \operatorname{ulp}(r_0)$. Roughly speaking, if r_0 is of the order of Λ , then e_0 is of order $\mathcal{O}(\varepsilon \Lambda)$. So for the product $(p, e) = 2MultFMA(a_i, b_j)$, p is of order $\mathcal{O}(\varepsilon^n \Lambda)$ and e of order $\mathcal{O}(\varepsilon^{n+1}\Lambda)$, where n = i + j, and we consider only the terms for which $0 \le n \le k$. This implies that for each n we have n + 1 products to compute (line 4 of Algorithm 5). Next, we need to add all terms of the same order of magnitude. Beside the n + 1 products, we also have n^2 terms resulting from the previous iteration. This addition is performed using *VecSum* (Algorithm 3) to obtain r_n in line 6.

The remaining terms are concatenated with the errors from the n + 1 products, and the entire $e_0, \ldots, e_{(n+1)^2-1}$ array is used in the next iteration. The (k+1)-st component r_k is obtained by simple summation of all remaining errors with the simple products of order $\mathcal{O}(\varepsilon^k \Lambda)$. EFT are not needed in the last step since the errors are not reused.

Algorithm 5. Algorithm of multiplication of FP expansions with *k* terms.

Input: FP expansions $a = a_0 + \cdots + a_{k-1}$; $b = b_0 + \cdots + b_{k-1}$.						
Dutput: FP expansion $r = r_0 + \cdots + r_{k-1}$.						
1: $(r_0, e_0) \leftarrow 2MultFMA(a_0, b_0)$						
2 : for $n \leftarrow 1$ to $k - 1$ do						
3: for $i \leftarrow 0$ to n do						
4: $(p_i, \hat{e}_i) \leftarrow 2MultFMA(a_i, b_{n-i})$						
5: end for						
6: $r_n, e[0:n^2+n-1] \leftarrow VecSum(p[0:n], e[0:n^2-1])$						
7: $e[0:(n+1)^2-1] \leftarrow e[0:n^2+n-1], \hat{e}[0:n]$						
8: end for						
9: for $i \leftarrow 1$ to $k - 1$ do						
$10: r_k \leftarrow r_k + a_i \cdot b_{k-i}$						
1: end for						
12: for $i \leftarrow 0$ to $k^2 - 1$ do						
$3: r_k \leftarrow r_k + e_i$						
4: end for						
$15: r[0:k-1] \leftarrow Renormalize(r[0:k])$						
16: return FP expansion $r = r_0 + \cdots + r_{k-1}$.						

For the addition and multiplication algorithms presented in this section, we will provide an effective error analysis in a subsequent paper. An important step for this goal is to provide a thorough proof for the renormalization, which is used at the end of each of these two algorithms. So, in what follows we focus on our new algorithm for renormalization of expansions.

3 RENORMALIZATION ALGORITHM FOR EXPANSIONS

While several renormalization algorithms have been proposed in literature, Priest [12] algorithm seems to be the only one provided with a complete correctness proof. It has many conditional branches, which make it slow in practice, and has a worst case FP operation count of: R(n) = 20(n-1), for an input FP expansion with *n*-terms.

In an attempt to overcome the branching problem we developed a new algorithm (Algorithm 6), for which we provide a full correctness proof.

First, we need to define the concept of FP numbers that overlap by at most d digits.

Definition 3.1. Consider an array of FP numbers: $x_0, x_1, \ldots, x_{n-1}$. According to Priest's [12] definition, they overlap by at most d digits $(0 \le d < p)$ if and only if $\forall i, 0 \le i \le n-2, \exists k_i, \delta_i$ such that:

$$2^{k_i} \le |x_i| < 2^{k_i+1},\tag{1}$$

$$2^{k_i - \delta_i} \le |x_{i+1}| \le 2^{k_i - \delta_i + 1},\tag{2}$$

$$\delta_i \ge p - d,\tag{3}$$

$$\delta_i + \delta_{i+1} \ge p - z_{i-1},\tag{4}$$

where z_{i-1} is the number of trailing zeros at the end of x_{i-1} and for $i = 0, z_{-1} := 0$.

Proposition 3.2. Let $x_0, x_1, \ldots, x_{n-1}$ be an array of FP numbers which overlap by at most d digits ($0 \le d < p$). The following properties hold:

$$\left|x_{j+1}\right| < 2^d \mathrm{ulp}(x_j),\tag{5}$$

$$ulp(x_{j+1}) \le 2^{d-p} ulp(x_j), \tag{6}$$

$$|x_{j+2} + x_{j+1}| \le (2^d + 2^{2d-p}) \operatorname{ulp}(x_j).$$
 (7)

Proof. We have $ulp(x_j) = 2^{k_j-p+1}$ and from (3) we get $|x_{j+1}| < 2^{k_j-\delta_j+1} < 2^{p-\delta_j}ulp(x_j) < 2^dulp(x_j)$. This proves that (5) holds for all $0 \le j < n-1$.

By applying (3) we get $ulp(x_{j+1}) = 2^{k_j - \delta_j - p + 1} \le 2^{d-p} ulp(x_j)$, which proves that (6) holds for all $0 \le j \le n-1$.

We have $|x_{j+1}| \leq 2^d \operatorname{ulp}(x_j)$ and $|x_{j+2}| \leq 2^d \operatorname{ulp}(x_{j+1}) \leq 2^{2d-p} \operatorname{ulp}(x_j)$ from which (7) follows.

The renormalization algorithm (Algorithm 6, illustrated in Fig. 5) is based on different layers of chained 2*Sum*. For the sake of simplicity, these are grouped in simpler layers based on *VecSum*. We will prove that our algorithm returns a \mathcal{P} -nonoverlapping sequence.



Fig. 5. Renormalization of FP expansions with n terms. The VecSum box performs Algorithm 3, the VecSum-ErrBranch box, Algorithm 7 and the VecSumErr box, Algorithm 8.

Proposition 3.3. Consider an array $x_0, x_1, \ldots, x_{n-1}$ of FP numbers that overlap by at most $d \le p-2$ digits and let m be an input parameter, with $1 \le m \le n-1$. Provided that no underflow / overflow occurs during the calculations, Algorithm 6 returns a "truncation" to m terms of a \mathcal{P} -nonoverlapping FP expansion $f = f_0 + \cdots + f_{n-1}$ such that $x_0 + \cdots + x_{n-1} = f$.

Algorithm 6. Renormalization algorithm

Input: FP expansion $x = x_0 + \cdots + x_{n-1}$ consisting of FP numbers that overlap by at most *d* digits, with $d \le p - 2$; *m* length of output FP expansion.

Output: FP expansion
$$f = f_0 + \dots + f_{m-1}$$
 with $f_{i+1} \le (\frac{1}{2} + 2^{-p+2} + 2^{-p}) \operatorname{ulp}(f_i)$, for all $0 \le i < m-1$.

1:
$$e[0:n-1] \leftarrow VecSum(x[0:n-1])$$

 $f^{(0)}[0, -1] \leftarrow VecSum(x[0:n-1])$

2:
$$f^{(0)}[0:m] \leftarrow VecSumErrBranch(e[0:n-1], m+1)$$

3: **for** $i \leftarrow 0$ to m - 2 **do**

4:

$$f^{(i+1)}[i:m] \leftarrow VecSumErr(f^{(i)}[i:m])$$

5: end for 6: return FP expansion $f = f_0^{(1)} + \dots + f_{m-2}^{(m-1)} + f_{m-1}^{(m-1)}$.

To prove this proposition, in what follows, we first prove several intermediate properties. The notations used in the proof $(s_i, e_i, \varepsilon_i, f_i, \rho_i \text{ and } g_i)$ are defined on the schematic drawings of the algorithms discussed. We also raise the important remark that at each step we prove that all the *2Sum* blocks can be replaced by *Fast2Sum* ones, but for simplicity of the proof we chose to present first the *2Sum* version.

First Level (Line 1, Algorithm 6)

It consists in applying Algorithm 3, *VecSum* (see also Fig. 2) on the input array, from where we obtain the array $e = (e_0, e_1, \ldots, e_{n-1})$.

Proposition 3.4. After applying the VecSum algorithm, the output array $e = (e_0, e_1, \ldots, e_{n-1})$ is S-nonoverlapping and may contain interleaving zeros.

Proof. Since $s_i = \text{RN}(x_i + s_{i+1})$, s_i is closer to $x_i + s_{i+1}$ than x_i . Hence $|(x_i + s_{i+1}) - s_i| \le |(x_i + s_{i+1}) - x_i|$, and so $|e_{i+1}| \le |s_{i+1}|$. Similarly, s_i is closer to $x_i + s_{i+1}$ than s_{i+1} , so $|e_{i+1}| \le |x_i|$. From (5) we get:

$$\begin{aligned} |x_{j+1}| + |x_{j+2}| + \dots &\leq \\ &\leq [2^d + 2^{2d-p} + 2^{3d-2p} + 2^{4d-3p} + \dots] \mathrm{ulp}(x_j) \quad (8) \\ &\leq 2^d \frac{2^p}{2^p - 1} \mathrm{ulp}(x_j). \end{aligned}$$

We know that $s_{j+1} = \text{RN}(x_{j+1} + \text{RN}(... + x_{n-1}))$ and by using a property given by Jeannerod and Rump in [18] we get:

$$|s_{j+1} - (x_{j+1} + \dots + x_{n-1})| \le (n - j - 2) \cdot 2^{-p} \cdot (|x_{j+1}| + \dots + |x_{n-1}|).$$
(9)

From (8) and (9) we have:

$$|s_{j+1}| \le 2^d \frac{2^p}{2^p - 1} (1 + (n - j - 2)2^{-p}) ulp(x_j).$$

It is easily seen that

$$2^{d} \frac{2^{p}}{2^{p} - 1} (1 + (n - j - 2)2^{-p}) \le 2^{p-1},$$
(10)

is satisfied for $p \ge 4$ and $n \le 16$, for $p \ge 5$ and $n \le 32$ and so on. This includes all practical cases, when $d \le p - 2$, so that $ulp(s_{j+1}) < ulp(x_j)$. Therefore x_j and s_{j+1} are multiples of $ulp(s_{j+1})$, thus $x_j + s_{j+1}$ is multiple of $ulp(s_{j+1})$, hence $RN(x_j + s_{j+1})$ is multiple of $ulp(s_{j+1})$ and $|e_{j+1}| = |x_j + s_{j+1} - RN(x_j + s_{j+1})|$ is multiple of $ulp(s_{j+1})$.

Also, by definition of 2Sum, we have $|e_{j+2}| \leq \frac{1}{2} \operatorname{ulp}(s_{j+1})$. Now, we can compare $|e_{j+1}|$ and $|e_{j+2}|$. Since $|e_{j+1}|$ is a multiple of $\operatorname{ulp}(s_{j+1})$, either $e_{j+1} = 0$ or e_{j+1} is larger than $2|e_{j+2}|$ and multiple of 2^k , such that $2^k > |e_{j+2}|$. This implies that the array $e = (e_0, e_1, \ldots, e_{n-1})$ is *S*-nonoverlapping and may have interleaving zeros.

Remark 3.5. Since we have $|s_{j+1}| \leq 2^{p-1} \operatorname{ulp}(x_j)$, for $d \leq p-2$ and $p \geq 4$ for n up to 16 and also $\operatorname{ulp}(x_j) \leq 2^{1-p} |x_j|$ we deduce that $|s_{j+1}| \leq |x_j|$. Hence, at this level we can use instead of *2Sum* basic blocks the *Fast2Sum* ones.

Second Level (Line 2, Algorithm 6)

It is applied on the array *e* obtained previously. This is also a chain of *2Sum*, but instead of starting from the least significant, we start from the most significant component. Also, instead of propagating the sums we propagate the errors. If however, the error after a *2Sum* block is zero, then we propagate the sum (this is shown in Fig. 6). In what follows we will refer to this algorithm by *VecSumErrBranch* (see Algorithm 7). The following property holds:

Proposition 3.6. Let an input array $e = (e_0, ..., e_{n-1})$ of S-nonoverlapping terms and $1 \le m \le n$ the required number of output terms. After appling VecSumErrBranch, the output array of $f = (f_0, ..., f_{m-1})$, with $0 \le m \le n-1$ satisfies $|f_{i+1}| \le ulp(f_i)$ for all $0 \le i < m-1$.



Fig. 6. *VecSumErrBranch* with n terms. Each *2Sum* box performs Algorithm 1, the sum is outputted downwards and the error to the right. If the error is zero, the sum is propagated to the right.

Algorithm 7. Second level of the renormalization algorithm - *VecSumErrBranch*

Input: *S*-*nonoverlapping* FP expansion $e = e_0 + \cdots + e_{n-1}$; *m* length of the output expansion.

Output: FP expansion $f = f_0 + \dots + f_{m-1}$ with $f_{j+1} \le ulp(f_j)$, $0 \le j < m-1$. 1: $j \leftarrow 0$ 2: $\varepsilon_0 = e_0$

- 3: for $i \leftarrow 0$ to n 2 do 4: $(f_i, \varepsilon_{i+1}) \leftarrow 2Sum(\varepsilon_i, e_{i+1})$
- 5: if $\varepsilon_{i+1} \neq 0$ then
- 6: **if** $j \ge m 1$ then
- 7: **return** FP expansion $f = f_0 + \dots + f_{m-1}$. //enough output terms

8: end if 9: $j \leftarrow j + 1$ 10: else 11: $\varepsilon_{i+1} \leftarrow f_j$ 12: end if 13: end for if $\varepsilon_{n-1} \neq 0$ and j < m then 14: 15: $f_i \leftarrow \varepsilon_{n-1}$ 16: end if

- 17: **return** FP expansion $f = f_0 + \cdots + f_{m-1}$.
- **Proof.** The case when *e* contains 1 or 2 elements is trivial. Consider now at least 3 elements. By definition of 2Sum, we have $|\varepsilon_1| \leq \frac{1}{2} \operatorname{ulp}(f_0)$ and by definition of *S*-nonoverlapping,

$$e_0 = E_0 \cdot 2^{k_0}$$
 with $|e_1| < 2^{k_0}$,
 $e_1 = E_1 \cdot 2^{k_1}$ with $|e_2| < 2^{k_1}$.

Hence, f_0 and ε_1 are both multiples of 2^{k_1} . Two possible cases may occur:

(*i*) $\varepsilon_1 = 0$. If we choose to propagate directly $\varepsilon_1 = 0$, then $f_1 = e_2$ and $\varepsilon_2 = 0$. This implies by induction that $f_i = e_{i+1}, \forall i \ge 1$. So, directly propagating the error poses a problem, since the whole remaining chain of 2Sum is executed without any change. So, as shown in Algorithm 7, line 11, when $\varepsilon_{i+1} = 0$ we propagate the sum f_j .

(*ii*) $\varepsilon_1 \neq 0$. Then $|e_2| < |\varepsilon_1|$ and $|\varepsilon_1 + e_2| < 2|\varepsilon_1|$, from where we get $|f_1| = |\text{RN}(\varepsilon_1 + e_2)| \le 2|\varepsilon_1| \le \text{ulp}(f_0)$.

We prove by induction the following statement: at step i > 0 of the loop in Algorithm 7, both f_{j-1} and ε_i are multiples of 2^{k_i} with $|e_{i+1}| < 2^{k_i}$. We proved above that i = 1 holds. Suppose now it holds for i and prove it for i + 1. Since f_{j-1} and ε_i are multiples of 2^{k_i} with $|e_{i+1}| < 2^{k_i}$ and $e_{i+1} = E_{i+1} \cdot 2^{k_{i+1}}$ with $|e_{i+2}| < 2^{k_{i+1}}$ (by definition of *S*-nonoverlapping), it follows that both f_j and ε_{i+1} are multiples of $2^{k_{i+1}}$ (by definition of *S*-monoverlapping).



Fig. 7. *VecSumErr* with m+1 terms. Each *2Sum* box performs Algorithm 1, sums are outputted downwards and errors to the right.

Finally, we prove the relation between f_j and f_{j-1} . If $\varepsilon_{i+1} = 0$, we propagate f_j , i.e., $\varepsilon_{i+1} = f_j$. Otherwise $|e_{i+1}| < |\varepsilon_i|$, so $|e_{i+1} + \varepsilon_i| < 2|\varepsilon_i|$ and finally $|f_j| = |\operatorname{RN}(e_{i+1} + \varepsilon_i)| \le 2|\varepsilon_i| \le \operatorname{ulp}(f_{j-1})$.

Remark 3.7. For this algorithm we can also use *Fast2Sum* instead of *2Sum*. We already showed that either $|e_{i+1}| < |\varepsilon_i|$, or $\varepsilon_i = 0$, in which case we replace $\varepsilon_i = f_{j-1}$, which is a multiple of 2^{k_i} with $|e_{i+1}| < 2^{k_i}$.

Third Level and Further (Lines 3-5, Algorithm 6)

On the previously obtained array we apply a similar chain of *2Sum*, starting from the most significant component and propagating the error. In these subsequent levels, no conditional branching is needed anymore (see Algorithm 8 and Fig. 7).

Algorithm 8. Third level of the renormalization algorithm - *VecSumErr*

Input: FP expansion $f = f_0 + \cdots + f_m$ with $|f_{i+1}| \le ulp(f_i)$, for all $0 \le i \le m - 1$. Output: FP expansion $g = g_0 + \cdots + g_m$ with $|g_1| \le (\frac{1}{2} + 2^{-p+2})$ $ulp(g_0)$ and $|g_{i+1}| \le ulp(g_i)$, for $0 < i \le m - 1$.

- 1: $\rho_0 = f_0$
- 2: for $i \leftarrow 0$ to m 1 do
- 3: $(g_i, \rho_{i+1}) \leftarrow 2Sum(\rho_i, f_{i+1})$
- 4: end for
- 5: $g_m \leftarrow \varepsilon_m$
- 6: **return** FP expansion $g = g_0 + \cdots + g_m$.

We prove the following property:

- **Proposition 3.8.** After applying Algorithm 8, VecSumErr on an input array $f = (f_0, \ldots, f_m)$, with $|f_{i+1}| \le ulp(f_i)$, for all $0 \le i \le m 1$, the output array $g = (g_0, \ldots, g_m)$ satisfies $|g_1| \le (\frac{1}{2} + 2^{-p+2})ulp(g_0)$ and $|g_{i+1}| \le ulp(g_i)$, for 0 < i < m 1.
- **Proof.** Since $|f_1| \leq \operatorname{ulp}(f_0)$ and $g_0 = \operatorname{RN}(f_0 + f_1)$ we have: $(1/2)\operatorname{ulp}(f_0) \leq \operatorname{ulp}(g_0) \leq 2\operatorname{ulp}(f_0)$, and $|\rho_1| \leq (1/2)$ $\operatorname{ulp}(g_0)$. We also have:

$$|f_1| \le ulp(f_0)$$
, which implies $ulp(f_1) \le 2^{-p+1} ulp(f_0)$,
 $|f_2| \le ulp(f_1) \le 2^{-p+2} ulp(g_0)$.

Hence:

$$|\rho_1 + f_2| \le (1/2 + 2^{-p+2}) \operatorname{ulp}(g_0).$$

Since $(1/2 + 2^{-p+2})ulp(g_0)$ is a FP number, we also have:

$$|g_1| = |\operatorname{RN}(\rho_1 + f_2)| \le (1/2 + 2^{-p+2}) \operatorname{ulp}(g_0).$$

This bound is very close to $(1/2)ulp(g_0)$ and it seems that in most practical cases, one actually has $|g_1| \leq \frac{1}{2}ulp(g_0)$. This implies that g_0 and g_1 are "almost" *B*-nonoverlapping and a simple computation shows that they are *P*-nonoverlapping as soon as $p \geq 3$, which occurs in all practical cases. As we iterate further, we get:

$$\begin{aligned} \left| \rho_{i+1} \right| &\leq (1/2) \mathrm{ulp}(g_i), \\ \left| f_{i+1} \right| &\leq \mathrm{ulp}(f_i), \text{ which implies } \mathrm{ulp}(f_{i+1}) \leq 2^{-p+1} \mathrm{ulp}(f_i). \end{aligned}$$

We know that ρ_i is multiple of $ulp(f_i)$ and from this we derive two cases: (i) $\rho_i = 0$, and as a consequence $\forall j \ge i, g_j = f_{j+1}$ and $g_m = 0$. In the second case (ii) we get:

$$|f_{i+1}| \le |\rho_i| \le (1/2) \operatorname{ulp}(g_{i-1}), |f_{i+1} + \rho_i| \le \operatorname{ulp}(g_{i-1}), |g_i| = |\operatorname{RN}(f_{i+1} + \rho_i)| \le \operatorname{ulp}(g_{i-1}).$$

Remark 3.9. For Algorithm 8 we can also use the faster algorithm, $Fast2Sum(\rho_i, f_{i+1})$, because we either have $\rho_i = 0$ or $|f_{i+1}| \le |\rho_i|$.

The above proposition shows that while we obtain a *nonoverlapping* condition for the first two elements of the resulting array g, for the others we don't strengthen the existing bound $|g_{i+1}| \leq \text{ulp}(g_i)$. There is an advantage however: if zeros appear in the summation process, they are pushed at the end; we don't use any branching. This suggests to continue applying a subsequent level of the same algorithm on the remaining elements, say g_1, \ldots, g_m . This is the idea of applying m - 1 levels of *VecSumErr* in lines 3-5, Algorithm 6. We are now able to prove Proposition 3.3.

Proof (of Proposition 3.3). Consider $m \ge 2$, otherwise the output reduces to only one term. The loop in lines 3-5 of Algorithm 6 is executed at least once. From Propositions 3.4, 3.6 and 3.8 we deduce that $|f_1^{(1)}| \le (1/2 + 2^{-p+2})$ $ulp(f_0^{(1)})$ and $|f_{i+1}^{(1)}| \le ulp(f_i^{(1)})$, for i > 0. If m = 2 then $f_0^{(1)}, f_1^{(1)}$ are outputted and the proposition is proven. Otherwise, $f_0^{(1)}$ is kept unchanged and another *VecSumErr* is applied to remaining $f_1^{(1)}, \ldots, f_m^{(1)}$. We have:

$$\begin{split} \left| f_1^{(1)} \right| &\leq \left(1/2 + 2^{-p+2} \right) \mathrm{ulp}(f_0^{(1)}), \\ \left| f_2^{(1)} \right| &\leq \mathrm{ulp}(f_1^{(1)}) \leq 2^{-p+1} \left(1/2 + 2^{-p+2} \right) \mathrm{ulp}(f_0^{(1)}), \\ &\leq 2^{-p+1} \mathrm{ulp}(f_0^{(1)}). \end{split}$$

Hence,

$$\begin{split} f_1^{(2)} &|= \Big| \mathrm{RN}(f_1^{(1)} + f_2^{(1)}) \Big|, \\ &\leq \big(1/2 + 2^{-p+2} + 2^{-p+1} \big) \mathrm{ulp}(f_0^{(1)}) \end{split}$$

Similarly,

$$\begin{split} \left| f_2^{(2)} \right| &\leq \left(1/2 + 2^{-p+2} \right) \mathrm{ulp}(f_1^{(2)}), \\ \left| f_3^{(2)} \right| &\leq \mathrm{ulp}(f_2^{(2)}) \leq 2^{-p+1} \big(1/2 + 2^{-p+2} \big) \mathrm{ulp}(f_1^{(2)}), \\ &\leq 2^{-p+1} \mathrm{ulp}(f_1^{(2)}). \end{split}$$



Fig. 8. Illustration of the effect of Algorithm 6. Expansion x is the input FP sequence, e is the sequence obtained after the first level and $f^{(0)}$ is the sequence obtained after the second level.

So, $f_0^{(1)}, f_1^{(2)}, f_2^{(2)}$ are *nonoverlapping*. It follows by induction that after m-1 loop iterations the output $f_0^{(1)}, \ldots, f_{m-2}^{(m-1)}, f_{m-1}^{(m-1)}$ is a *P*-nonoverlapping expansion. Finally, when all n-1 terms are considered, after at most n-1 loop iterations we have: $x_0 + \cdots + x_{n-1} = f_0^{(1)} + \cdots + f_{n-2}^{(n-1)} + f_{n-1}^{(n-1)}$.

Fig. 8 gives an intuitive drawing showing the different constraints between the FP numbers before and after the first two levels of Algorithm 6. The notation is the same as in Fig. 5.

Remark 3.10. In the worst case, Algorithm 6 performs n - 1*Fast2Sum* calls in the first level and n - 2 *Fast2Sum* calls plus n - 1 comparisons in the second one. During the following m - 1 levels we perform m - i *Fast2Sum* calls, with $0 \le i < m - 2$. This accounts for a total of $R_{new}(n,m) = 7n + \frac{3}{2}m^2 + \frac{3}{2}m - 13$ FP operations.

Table 1 shows some effective values of the worst case FP operation count for Priest's renormalization algorithm [12] and Algorithm 6. One can see that for $n \leq 7$ our algorithm performs better or the same. Even though from values of n > 7 Algorithm 6 performs worse in terms of operation count than Priest's one, in practice, the last m - 1 levels will take advantage of the computers pipeline, because we do not need branching conditions anymore, which makes it faster in practice.

In what follows we denote by *AddRoundE* (x[0:n-1], y[0:m-1], r), an algorithm for expansions addition, which given two (\mathcal{P} - or \mathcal{B} -) nonoverlapping expansions, returns the r most significant terms of the exact normalized (\mathcal{P} - or \mathcal{B} -) nonoverlapping sum. If no request is made on the number of terms to be returned, then we denote simply by AddE(x[0:n-1], y[0:m-1]). Similarly, we denote by MulRoundE, MulE, SubRoundE, SubE, DivRoundE, RenormalizeE algorithms for multiplication, subtraction, division and normalization.

4 RECIPROCAL ALGORITHM

4.1 Algorithms Using Classical Long Division on Expansions

In reference [12], division is done using the classical long division algorithm (a variation of the paper-and-pencil method), which is recalled in Algorithm 9. Bailey's division algorithm [6] is similar. For instance, let $a = a_0 + a_1 + a_2 + a_3$ and $b = b_0 + b_1 + b_2 + b_3$ be QD numbers. First, one approximates the quotient $q_0 = a_0/b_0$, then computes the

TABLE 1 FP Operation Count for Algorithm 6 versus Priest's Renormalization Algorithm [12]

q	2	4	7	8	10	12	16
Alg. 6 Priest's alg. [12]	10 20	45 60	120 120	$\begin{array}{c} 151 \\ 140 \end{array}$	222 180	$305 \\ 220$	$507 \\ 300$

We consider that both algorithms compute n-1 terms in the output expansion.

remainder $r = a - q_0 b$ in quad-double. The next correction term is $q_1 = r_0/b_0$. Subsequent terms q_i are obtained by continuing this process. At each step when computing r, full quad-double multiplication and subtraction are performed since most of the bits will be canceled out when computing q_3 and q_4 , in Bailey's algorithm. A renormalization step is performed only at the end, on $q_0 + q_1 + q_2 + ...$ in order to ensure non-overlapping. No error bound is given in [6].

Note that in Algorithm 9 [12] a renormalization step is performed after each computation of $r = r - q_i b$. An error bound is given in [12]:

Algorithm 9. Priest's [12] division algorithm. We denote by f[0:...] and expansion f whose number of terms is not known in advance.

Input: FP expansion $a = a_0 + \cdots + a_{n-1}$; $b = b_0 + \cdots + b_{m-1}$; length of output quotient FP expansion *d*.

Output: FP expansion $q = q_0 + \dots$ with at most d terms s.t. $\left|\frac{q-a/b}{a/b}\right| < 2^{1-\lfloor (p-4)d/p \rfloor}$.

1: $q_0 = \mathrm{RN}(a_0/b_0)$ $r^{(0)}[0:n-1] \leftarrow a[0:n-1]$ 2: 3: **for** $i \leftarrow 1$ to d - 1 **do** 4: $f[0:\ldots] \leftarrow MulE(q_{i-1}, b[0:m-1])$ 5: $r^{(i)}[0:\ldots] \leftarrow RenormalizeE(SubE(r^{(i-1)}[0:\ldots], f[0:\ldots]))$ $q_i = \mathrm{RN}(r_0^{(i)}/b_0)$ 6: 7: end for 8: $q[0:\ldots] \leftarrow RenormalizeE(q[0:d-1])$

9: **return** FP expansion $q = q_0 + \ldots$

Proposition 4.1. [12] Consider two \mathcal{P} -nonoverlapping expansions: $a = a_0 + \cdots + a_{n-1}$ and $b = b_0 + \cdots + b_{m-1}$, Priest division algorithm [12] computes a quotient expansion $q = q_0 + \cdots + q_{d-1}$ s.t.

$$\left|\frac{q-a/b}{a/b}\right| < 2^{1-\lfloor (p-4)d/p\rfloor}.$$
(11)

Daumas and Finot [19] modify Priest's division algorithm by using only estimates of the most significant component of the remainder r_0 and storing the less significant components of the remainder and the terms $-q_ib$ unchanged in a set that is managed with a priority queue. While the asymptotic complexity of this algorithm is better, in practical simple cases Priest's algorithm is faster due to the control overhead of the priority queue [19]. The error bound obtained with Daumas' algorithm is (using the same notations as above):

$$\left|\frac{q-a/b}{a/b}\right| < 2^{-d(p-1)} \prod_{i=0}^{d-1} (4i+6).$$
(12)

4.2 Reciprocal of Expansions with an Adapted Newton-Raphson Iteration

The classical Newton-Raphson iteration for computing reciprocals [7, Chap. 2] [9], [10], is based on the general NR iteration for computing the roots of a given function *f*, which is:

$$x_{n+1} = x_n - \frac{f(x_n)}{f'(x_n)}.$$
(13)

When x_0 is close to a root α , $f'(\alpha) \neq 0$, the iteration converges quadratically. For computing 1/a we choose f(x) = 1/x - a, which gives

$$x_{n+1} = x_n(2 - ax_n). (14)$$

The iteration converges to 1/a for all $x_0 \in (0, 2/a)$. However, taking any point in (0, 2/a) as the starting point x_0 would be a poor choice. A much better choice is to choose x_0 equal to a FP number very close to 1/a. This only requires one FP division. The quadratic convergence of (14) is deduced from $x_{n+1} - 1/a = -a(x_n - 1/a)^2$. This iteration is *self-correcting* because rounding errors do not modify the limit value.

Algorithm 10. Truncated Newton iteration based algorithm for reciprocal of an FP expansion.

In	put: FP expansion $a = a_0 + \cdots + a_{2^k-1}$; length of output FP
	expansion 2^q .
0	utput: FP expansion $x = x_0 + \cdots + x_{2^{q-1}}$ s.t. $ x - \frac{1}{q} \leq$
	$\frac{2^{-2^q(p-3)-1}}{ a }$.
1:	$x_0 = \mathrm{RN}(1/a_0)$
2:	for $i \leftarrow 0$ to $q - 1$ do
3:	$\hat{v}[0:2^{i+1}-1] \leftarrow MulRoundE(x[0:2^{i}-1],a[0:2^{i+1}-1],2^{i+1})$
4:	$\hat{w}[0:2^{i+1}-1] \leftarrow SubRoundE(2, \hat{v}[0:2^{i+1}-1], 2^{i+1})$
5:	$x[0:2^{i+1}-1] \leftarrow MulRoundE(x[0:2^{i}-1], \hat{w}[0:2^{i+1}-1], 2^{i+1})$
6:	end for
7:	return FP expansion $x = x_0 + \cdots + x_{2^{q-1}}$.

While iteration (14) is well known, in Algorithm 10 we use an adaptation for computing reciprocals of FP expansions, with truncated operations involving FP expansions. Our algorithm works with both *B- and P-nonoverlapping* FP expansions. For the sake of clarity we consider first the case of *B-nonoverlapping* FP expansions, and then make the necessary adjustments for *P-nonoverlapping* expansions in Proposition 4.4.

4.3 Error Analysis of Algorithm 10

Let $a = a_0 + \cdots + a_{2^k-1}$ be a *B*-nonoverlapping FP expansion with 2^k terms and $q \ge 0$. We will prove that our algorithm returns an approximation $x = x_0 + \cdots + x_{2^q-1}$ of 1/a, in the form of a *B*-nonoverlapping FP expansion with 2^q terms, such that

$$|x - 1/a| \le 2^{-2^q(p-3)-1}/|a|.$$
 (15)

We will first prove the following proposition:

Proposition 4.2. Consider a \mathcal{B} -nonoverlapping expansion $u = u_0 + u_1 + \cdots + u_k$ with k normal binary FP terms of precision p. Denote $u^{(i)} = u_0 + u_1 + \cdots + u_i, i \ge 0$, i.e., "a truncation" of u to i + 1 terms. The following inequalities hold for $0 \le i \le k$:

$$|u_i| \le 2^{-ip} |u_0|, \tag{16}$$

$$\left|u - u^{(i)}\right| \le 2^{-ip} |u| \frac{\eta}{1 - \eta},$$
(17)

$$\left(1 - \frac{2^{-ip}\eta}{1 - \eta}\right)|u| \le \left|u^{(i)}\right| \le \left(1 + \frac{2^{-ip}\eta}{1 - \eta}\right)|u|, \tag{18}$$

$$1/u - 1/u_0| \le \eta/|u|, \tag{19}$$

where

$$\eta = \sum_{j=0}^{\infty} 2^{(-j-1)p} = \frac{2^{-p}}{1 - 2^{-p}}$$

Proof. By definition of a *B*-nonoverlapping expansion and since for any normal binary FP number u_i , $ulp(u_i) \le 2^{-p+1}|u_i|$ we have $|u_i| \le (1/2)ulp(u_{i-1}) \le 2^{-p}|u_{i-1}|$ and (16) follows by induction.

Consequently we have $|u - u_0| = |u_1 + u_2 + \dots + u_k|$ $\leq 2^{-p}|u_0| + 2^{-2p}|u_0| + \dots + 2^{-kp}|u_0| \leq |u_0|\eta$. One observes that u and u_0 have the same sign. A possible proof is by noticing that $1 - \eta > 0$ and $-|u_0|\eta \leq u - u_0 \leq |u_0|\eta$. Suppose $u_0 > 0$, then $-u_0\eta \leq u - u_0 \leq u_0\eta$, hence $u_0(1 - \eta) \leq u \leq u_0(1 + \eta)$ which implies u > 0. The case $u_0 < 0$ is similar. It follows that

$$|u|/(1+\eta) \le |u_0| \le |u|/(1-\eta).$$
⁽²⁰⁾

For (17) we use (20) together with:

$$\left|u-u^{(i)}\right| \leq \sum_{j=0}^{\infty} 2^{(-i-j-1)p} |u_0| \leq 2^{-ip} \eta |u_0|,$$

- and (18) is a simple consequence of (17). Similarly, (19) follows from $|1/u 1/u_0| = (1/|u|) \cdot |(u_0 u)/u_0| \le \eta/|u|$.
- **Proposition 4.3.** Provided that no underflow / overflow occurs during the calculations, Algorithm 10 is correct when run with B-nonoverlapping expansions.
- **Proof.** The input of the algorithm is a non-overlapping FP expansion $a = a_0 + a_1 + \cdots + a_{2^{k}-1}$, in which every term a_i is a normal binary FP number of precision p. Let $f_i = 2^{i+1} 1$ and $a^{(f_i)} = a_0 + a_1 + \cdots + a_{f_i}$ i.e., "a truncation" of a to $f_i + 1$ terms, with $0 \le i$.

For computing 1/a we use Newton iteration: $x_0 = \text{RN}(1/a_0), x_{i+1} = x_i(2 - a^{(f_i)}x_i)), i \ge 0$ by truncating each FP expansion operation as follows:

- let $v_i := (a^{(f_i)} \cdot x_i)$ be the exact product represented as a non-overlapping expansion on $2^{2(i+1)}$ terms, we compute $\hat{v}_i := v_i^{(2^i)}$ i.e., v_i "truncated to" 2^{i+1} terms;
- let w_i := 2 v̂_i be the exact result of the subtraction represented as a non-overlapping expansion on 2ⁱ⁺¹ + 1 terms, we compute ŵ_i := w_i^(2ⁱ) i.e., v_i "truncated to" 2ⁱ⁺¹ terms;
- let τ_i := x_i · ŵ_i be the exact product represented as a non-overlapping expansion on 2 · 2ⁱ(2ⁱ⁺¹) terms, we compute x_{i+1} := τ_i^(2ⁱ⁺¹-1) i.e., τ_i "truncated to" 2ⁱ⁺¹ terms.

Let us first prove a simple upper bound for the approximation error in x_0 :

$$\varepsilon_0 = |x_0 - 1/a| \le 2\eta/|a|.$$
 (21)

Since $x_0 = \text{RN}(1/a_0)$, then $|x_0 - 1/a_0| \le 2^{-p} |1/a_0|$, so $|x_0 - 1/a| \le 2^{-p} |1/a_0| + |1/a - 1/a_0| \le ((1 + \eta)2^{-p} + \eta)/|a| \le 2\eta/|a|$ (from (20)).

Let us deduce an upper bound for the approximation error in x at step i + 1, $\varepsilon_{i+1} = |x_{i+1} - 1/a|$. For this, we will use a chain of triangular inequalities that make the transition from our "truncated" Newton error to the "untruncated" one. Let $\gamma_i = 2^{-(2^{i+1}-1)p}\eta/(1-\eta)$. We have from Proposition 4.2, eq. (17):

$$|x_{i+1} - \tau_i| \le \gamma_i |x_i \cdot \hat{w}_i|,\tag{22}$$

$$|w_i - \hat{w}_i| \le \gamma_i |w_i| \le \gamma_i |2 - \hat{v}_i|, \tag{23}$$

$$|v_i - \hat{v}_i| \le \gamma_i |a^{(f_i)} \cdot x_i|, \qquad (24)$$

$$\left|a - a^{(f_i)}\right| \le \gamma_i |a|. \tag{25}$$

From (22) we have:

$$\begin{split} \varepsilon_{i+1} &\leq |x_{i+1} - \tau_i| + |\tau_i - 1/a| \\ &\leq \gamma_i |x_i \cdot \hat{w}_i| + |x_i \cdot \hat{w}_i - 1/a| \\ &\leq \gamma_i |x_i(w_i - \hat{w}_i)| + \gamma_i |x_i w_i| + |x_i \cdot \hat{w}_i - 1/a| \\ &\leq (1 + \gamma_i) |x_i| |w_i - \hat{w}_i| + \gamma_i |x_i w_i| \\ &+ |x_i \cdot w_i - 1/a|. \end{split}$$

Using (23) and (24):

$$\begin{split} \varepsilon_{i+1} &\leq |x_i \cdot w_i - 1/a| + ((1+\gamma_i)\gamma_i + \gamma_i)|x_iw_i| \\ &\leq |x_i \cdot (2-v_i) - 1/a| + |x_i| \cdot |(v_i - \hat{v}_i)| \\ &+ (\gamma_i(1+\gamma_i) + \gamma_i)|x_i|(|(2-v_i)| + |v_i - \hat{v}_i|) \\ &\leq |x_i \cdot (2-a^{(f_i)} \cdot x_i) - 1/a| \\ &+ \gamma_i(1+\gamma_i)^2 |x_i^2| |a^{(f_i)}| \\ &+ (\gamma_i(1+\gamma_i) + \gamma_i)|x_i(2-a^{(f_i)} \cdot x_i)|. \end{split}$$

By (25), we have:

$$\begin{aligned} \left| x_i \cdot (2 - a^{(f_i)} \cdot x_i) - 1/a \right| &\leq |a| |x_i - 1/a|^2 + \gamma_i |x_i|^2 |a|, \\ |x_i|^2 |a^{(f_i)}| &\leq (1 + \gamma_i) |x_i|^2 |a|, \end{aligned}$$

and

$$|x_i \cdot (2 - a^{(f_i)} \cdot x_i)| \le |a| |x_i - 1/a|^2 + \gamma_i |x_i|^2 |a| + 1/|a|.$$

Hence we have:

$$\varepsilon_{i+1} \le (1+\gamma_i)^2 |a| |x_i - 1/a|^2 + \gamma_i (1+\gamma_i)^2 (2+\gamma_i) |x_i^2| |a| + \gamma_i (2+\gamma_i) 1/|a|.$$
(26)

We now prove by induction that for all $i \ge 0$:

$$\varepsilon_i = |x_i - 1/a| \le 2^{-2^i(p-3)-1}/|a|.$$
 (27)

For i = 0, this holds from (21) and the fact that $\eta = 1/(2^p - 1) \le 2^{-p+1}$. For the induction step, we have from (26):

$$\varepsilon_{i+1} \leq (1+\gamma_i)^2 |a| |\varepsilon_i|^2 + \gamma_i (1+\gamma_i)^2 (2+\gamma_i) (1\pm\varepsilon_i |a|)^2 \frac{1}{|a|} + \gamma_i (2+\gamma_i) \frac{1}{|a|},$$
(28)

which implies

$$\frac{|a|\varepsilon_{i+1}}{2^{-2^{i+1}(p-3)}} \leq \frac{(1+\gamma_i)^2}{4} + \frac{(1+2^{-p+2})(2+\gamma_i)}{64} \\ \cdot (1+(1+\gamma_i)^2(1+2^{-2^i(p-3)-1})^2) \\ \leq 1/2.$$
(29)

This completes our proof.

- **Proposition 4.4.** Algorithm 10 is correct when run with \mathcal{P} -nonoverlapping expansions.
- **Proof.** The previous analysis holds provided that we use Remark 2.6. This mainly implies the following changes $\eta' = \frac{2}{2^p - 3'} \gamma'_i = \left(\frac{2}{2^p - 1}\right)^{2^{i+1} - 1} \frac{\eta'}{1 - \eta'}$. With this change it is easy to verify that equations (21)-(28) hold as soon as p > 2. Note that for the induction step at i = 1, a tighter bound is needed for $\varepsilon'_0 \leq \frac{2^{-p}(1 + \eta') + \eta'}{|a|} \leq \frac{2\eta'}{|a|} \frac{3 - 2^{-p}}{4}$, but the rest of the proof is identical, safe for some tedious computations. □

4.4 Complexity Analysis for Reciprocal

Our algorithm has the feature of using "truncated" expansions, while some variants of *Add-RoundE* and *MulRoundE* compute the result fully and only then truncate. This is the case of Priest's algorithms, which are not tuned for obtaining "truncated" expansions on the fly—and thus penalize our algorithm—. On the other hand, the algorithms presented in Section 2.2, can take into account only the significant terms of the input expansions in order to compute the result. Even though these algorithms have not yet been proven to work properly, we obtained promising results in practice, so we will perform the complexity analysis based on them.

We present here the operation count of our algorithms, by taking ([7]) 6 FP operations for 2*Sum*, 3 for *Fast2Sum* and 2 for 2*MultFMA*. For the sake of simplicity, for multiplication, we will consider that the input expansions have the same number of terms.

- The renormalization (Algorithm 6) of an overlapping expansion x with n terms, requires $(2n-3) + \sum_{i=0}^{m-2} m - i$ Fast2Sum calls and n-1 comparisons. This accounts for $R_{new}(n,m) = 7n + \frac{3}{2}m^2 + \frac{3}{2}m - 13$ FP operations.
- The addition (Algorithm 4) of two *P*-nonoverlapping expansions requires n + m 1 comparisons and a renormalization $R_{new}(n + m, r)$. This accounts for $A(n, m, r) = \frac{3}{2}r^2 + \frac{3}{2}r + 8n + 8m 14$ FP operations.
- The multiplication (Algorithm 5) of two *P*-nonoverlapping expansions requires $\sum_{i=1}^{k} i$ 2MultFMA calls, $\sum_{i=1}^{k-1} (n^2 + n)$ 2Sum calls, k - 1 FP multiplications, followed by $k^2 + k - 2$ FP additions and a

renormalization $R_{new}(k+1,k)$ in the end. This accounts for $M(k) = 2k^3 + \frac{7}{2}k^2 + \frac{19}{2}k - 9$ FP operations.

- -> The special case of multiplying a FP expansion to a single FP value accounts for only $M1(k) = \frac{9}{2}k^2 + \frac{17}{2}k 7$.
- Using these algorithms for addition and multiplication of FP expansions, Priest's division (Algorithm 9) requires d divisions and $(d-1)(M1(k)) + \sum_{i=0}^{d-1} A(k+2k(i-1),k+2k(i-1),k+2k(i-1)) + R_{new}(d,d)$ function calls in the worst case. This accounts for $D(d,k) = 2d^3k^2 - 6d^2k^2 + \frac{32}{2}d^2k + \frac{3}{2}d^2 + 10dk^2 - \frac{52}{2}dk - \frac{23}{2}d - \frac{9}{2}k^2 - \frac{17}{2}k - 6$ FP operations.
- **Proposition 4.5.** Using for addition and multiplication of FP expansions the algorithms presented in Section 2.2, Algorithm 10 requires $\frac{32}{7}8^q + \frac{34}{3}4^q + 57 \cdot 2^q 24q \frac{1510}{21}$ FP operations.
- **Proof.** During the *i*th iteration the following operations with expansions are performed: two multiplications $M(2^{i+1})$; one addition $A(2^{i+1}, 1, 2^{i+1})$. Since *q* iterations are done, the total number of FP operations is: $\frac{32}{7}8^q + \frac{34}{3}4^q + 57 \cdot 2^q 24q \frac{1510}{21}$ FP operations.
- **Remark 4.6.** Division is simply performed with Algorithm 10 followed by a multiplication $M(2^q)$ where the numerator expansion has 2^q terms.

5 SQUARE ROOT ALGORITHMS

The families of algorithms most commonly used are exactly the same as for division, although, in the case of FP expansions the digit-recurrence algorithm is typically more complicated than for division. This is why a software implementation would be tedious. Moreover, Newton-Raphson based algorithms offer the advantage of assuring a quadratic convergence.

5.1 Square Root of Expansions with an Adapted Newton-Raphson Iteration

Starting from the general Newton-Raphson iteration (13), we can compute the square root in two different ways. We can look for the zeros of the function $f(x) = x^2 - a$ that leads to the so called "Heron iteration":

$$x_{n+1} = \frac{1}{2} \left(x_n + \frac{a}{x_n} \right).$$
(30)

If $x_0 > 0$, then x_n goes to \sqrt{a} . This iteration needs a division at each step, which counts as a major drawback.

To avoid performing a division at each step we can look for the positive root of the function $f(x) = 1/x^2 - a$. This gives the iteration

$$x_{n+1} = \frac{1}{2}x_n \left(3 - ax_n^2\right). \tag{31}$$

This iteration converges to $1/\sqrt{a}$, provided that $x_0 \in (0, \sqrt{3}/\sqrt{a})$. The result can be multiplied by *a* in order to get an approximation of \sqrt{a} . To obtain fast, quadratic, convergence, the first point x_0 must be a close approximation to $1/\sqrt{a}$. The division by 2 is done by multiplying each of the terms of the input expansion by 0.5, separately.

As in the case of the reciprocal (Section 4.2), in Algorithm 11 we use an adaption of iteration (31), using the truncated algorithms presented above.

Algorithm 11. Truncated "division-free" Newton iteration (31) based algorithm for reciprocal of the square root of an FP expansion. By "division-free" we mean that we do not need a division of FP expansions.

Input: FP expansion $a = a_0 + \cdots + a_{2^{k}-1}$; length of output FP expansion 2^q . **Output:** FP expansion $x = x_0 + \cdots + x_{2^q-1}$ s.t.

$$\left|x - 1/\sqrt{a}\right| \le 2^{-2^{q}(p-3)-1}/\sqrt{a}.$$
 (32)

1: $x_0 = \text{RN}(1/\sqrt{a_0})$ 2: **for** $i \leftarrow 0$ to q - 1 **do** 3: $\hat{v}[0: 2^{i+1} - 1] \leftarrow MulRoundE(x[0: 2^i - 1], a[0: 2^{i+1} - 1], 2^{i+1})$ 4: $\hat{w}[0: 2^{i+1} - 1] \leftarrow MulRoundE(x[0: 2^i - 1], \hat{v}[0: 2^{i+1} - 1], 2^{i+1})$ 5: $\hat{y}[0: 2^{i+1} - 1] \leftarrow SubRoundE(3, \hat{w}[0: 2^{i+1} - 1], 2^{i+1})$ 6: $\hat{z}[0: 2^{i+1} - 1] \leftarrow MulRoundE(x[0: 2^i - 1], \hat{y}[0: 2^{i+1} - 1], 2^{i+1})$ 7: $x[0: 2^{i+1} - 1] \leftarrow \hat{z}[0: 2^{i+1} - 1] * 0.5$ 8: **end for** 9: **return** FP expansion $x = x_0 + \dots + x_{2^q-1}$.

The error analysis for this algorithm follows the same principle as the one for the reciprocal algorithm. The detailed proof is given in Appendix A. We show that the relative error decreases after every loop of the algorithm, by taking into account the truncations performed after each operation. The strategy is to make the exact Newton iteration term and bound appear. We show that by the end of the *i*th iteration of the loop, $\varepsilon_i = |x^{(2^i)} - 1/\sqrt{a}| \le 2^{-2^i(p-3)-1}/\sqrt{a}$.

In his library, QD, Bailey also uses the Newton iteration for the square root computation. Although he uses the same function as we do, he uses the iteration under the form: $x_{i+1} = x_i + \frac{1}{2}x_i(1 - ax_i^2)$, which from a mathematical point of view is the same, but it requires a different implementation. Even though Bailey does not provide an error analysis for his algorithm, we managed to prove that the error bound is preserved when using this iteration (see Appendix A for the detailed proof).

Algorithm 12. Truncated "Heron iteration" (30) based algorithm for square root of an FP expansion.

Input: FP expansion $a = a_0 + \cdots + a_{2^k-1}$; length of output FP expansion 2^q .

Output: FP expansion $x = x_0 + \cdots + x_{2^q-1}$ s.t.

$$|x - \sqrt{a}| \le 3\sqrt{a} \cdot 2^{-2^q(p-3)-2}.$$
(33)

1: $x_0 = \text{RN}(\sqrt{a_0})$ 2: for $i \leftarrow 0$ to q - 1 do 3: $\hat{v}[0: 2^{i+1} - 1] \leftarrow DivRoundE(a[0: 2^{i+1} - 1], x[0: 2^i - 1], 2^{i+1})$ 4: $\hat{w}[0: 2^{i+1} - 1] \leftarrow AddRoundE(x[0: 2^i - 1], \hat{v}[0: 2^{i+1} - 1], 2^{i+1})$ 5: $x[0: 2^{i+1} - 1] \leftarrow \hat{w}[0: 2^{i+1} - 1] * 0.5$ 6: end for 7: return FP expansion $x = x_0 + \dots + x_{2^{q-1}}$.

TABLE 2 Error Bounds Values for Priest (11) versus Daumas (12) versus Our Analysis (15)

Prec, iteration	Eq. (11)	Eq. (12)	Eq. (15)	β
p = 53, q = 0	2	2^{-49}	2^{-51}	2^{-52}
p = 53, q = 1	1	2^{-98}	2^{-101}	2^{-104}
p = 53, q = 2	2^{-2}	2^{-195}	2^{-201}	2^{-208}
p = 53, q = 3	2^{-6}	2^{-387}	2^{-401}	2^{-416}
p = 53, q = 4	2^{-13}	2^{-764}	2^{-801}	2^{-833}
p = 24, q = 0	2	2^{-20}	2^{-22}	2^{-23}
p = 24, q = 1	1	2^{-40}	2^{-43}	2^{-46}
p = 24, q = 2	2^{-2}	2^{-79}	2^{-85}	2^{-92}
p = 24, q = 3	2^{-5}	2^{-155}	2^{-169}	*
p = 24, q = 4	2^{-12}	2^{-300}	2^{-337}	*

 β gives the largest obtained errors for Algorithm 10 using the standard FP formats double and single. *underflow occurs.

"Heron Iteration" Algorithm

The same type of proof as above can be applied for the algorithm using the "Heron iteration" (30) and the same type of truncations. In this case (Algorithm 12) we obtain a slightly larger error bound for both types of *nonoverlapping* FP expansions: $|x - \sqrt{a}| \leq 3\sqrt{a} \cdot 2^{-2^{q}(p-3)-2}$.

5.2 Complexity Analysis for Square Root

We will perform our operation count based on the addition and multiplication presented in Section 2.2, the same as in Section 4.4.

- **Proposition 5.1.** Using for addition, multiplication and division of FP expansions the algorithms previously presented, Algorithm 11 requires $\frac{48}{7}8^q + 16 \cdot 4^q + 78 \cdot 2^q 33q \frac{699}{7}$ FP operations.
- **Proof.** During the *i*th iteration, three multiplications $M(2^{i+1})$, one addition $A(2^{i+1}, 1, 2^{i+1})$ and one division by 2 are performed. Since *q* iterations are done, the total number of FP operations is: $\frac{48}{7}8^q + 16 \cdot 4^q + 78 \cdot 2^q 33q \frac{699}{7}$.
- **Remark 5.2.** We obtain the square root of an expansion by simply multiplying the result obtained from Algorithm 11 by the input expansion *a*. This means an additional $M(2^q)$, where 2^q is the number of terms in *a*.
- **Proposition 5.3.** Using for addition, multiplication and division of FP expansions the algorithms previously presented, Algorithm 12 requires $\frac{368}{49}8^q + \frac{196}{9}4^q + 170 \cdot 2^q 12q^2 \frac{2245}{21}q \frac{87445}{441}$ FP operations.
- **Proof.** One addition $A(2^{i+1}, 2^{i+1}, 2^{i+1})$, one division $D(2^{i+1})$ and a division by 2 are performed during each *i*th iteration. Since *q* iterations are done, the total number of FP operations is: $\frac{368}{49}8^q + \frac{196}{9}4^q + 170 \cdot 2^q - 12q^2 - \frac{2245}{21}q - \frac{87445}{441}$.

Based on these values, Algorithm 11 performs slightly better than Algorithm 12, and in the same time the obtained error bound is tighter.

6 COMPARISON AND DISCUSSION

In Table 2 we show values of the bounds provided by our error analysis, compared with those of Priest and Daumas for the reciprocal computation. Our algorithm performs

TABLE 3FP Operation Count for Priest versus Our Algorithm; $d = 2^q$ Terms Are Computed in the Quotient

d	2	4	8	16
Alg. 9 [12]	24	1,714	52,986	1,808,698
Alg. 10 + Alg. 5	140	795	4,693	31,601

better for the same number of terms in the computed quotient, say $d = 2^q$ in equations (11) and (12). Moreover, our algorithm provides a unified error bound with quadratic convergence independent of using underlying \mathcal{P} - or \mathcal{B} *nonoverlapping* expansions. In the last column of the same table we give the largest errors that we actually obtained through direct computation of the reciprocal using our algorithm. The given value represents the obtained value upper rounded to the immediate power of 2. For each table entry we performed one million random tests.

The complexity analysis shows that our algorithm performs better, for expansions with more than two terms, even if no error bound is requested (see Table 3 for some effective values of the worst case FP operation count).

Note that, for instance, to guarantee an error bound of $2^{-d(p-3)-1}$, Priest's algorithm (based on the bound given in Prop. 4.1) needs at least (dp - 3d + 2)p/(p - 4) terms, which entails a very poor complexity. This implies that Daumas' algorithm might be a good compromise in this case, provided that the priority queue used there can be efficiently implemented.

This plus the performance tests that we ran confirm our hypothesis that for higher precisions the Newton-Raphson iteration is preferable to classical division.

In the case of the square root, because no error bound is given for the *digit-recurrence* algorithm we can only compare between the errors that we obtain if using the two different types of Newton iteration available for computing the square root. The effective values of the bounds are given in Table 4. The bound provided for Algorithm 11 is only slightly tighter that the one for Algorithm 12. The same as for the reciprocal, in the last column we present the bounds obtained through direct computation using Algorithm 11.

In Table 5 we give some effective values of the worst case FP operation count for Algorithm 11 versus Algorithm 12 based on Section 5.2.

TABLE 4 Error Bounds Values for (44) versus (33)

Prec, iteration	Eq. (44)	Eq. (33)	β
p = 53, q = 0	2^{-51}	$3 \cdot 2^{-52}$	2^{-52}
p = 53, q = 1	2^{-101}	$3 \cdot 2^{-102}$	2^{-103}
p = 53, q = 2	2^{-201}	$3 \cdot 2^{-202}$	2^{-206}
p = 53, q = 3	2^{-401}	$3 \cdot 2^{-402}$	2^{-412}
p = 53, q = 4	2^{-801}	$3 \cdot 2^{-802}$	2^{-823}
p = 24, q = 0	2^{-22}	$3 \cdot 2^{-23}$	2^{-23}
p = 24, q = 1	2^{-43}	$3 \cdot 2^{-44}$	2^{-45}
p = 24, q = 2	2^{-85}	$3 \cdot 2^{-86}$	2^{-90}
p = 24, q = 3	2^{-169}	$3 \cdot 2^{-170}$	*
p = 24, q = 4	2^{-337}	$3 \cdot 2^{-338}$	*

 β gives the Largest Obtained Errors for Algorithm 11 using the Standard FP formats double and single. *underflow occurs.

TABLE 5FP Operation Count for Algorithm 11 versus Algorithm 12;2^q Terms Are Computed in the Quotient

\overline{q}	1	2	3	4
Alg. 11 + Alg. 5 Alg. 12	182 170	$1,054 \\ 1,049$	$6,275 \\ 5,972$	42,430 38,239

The algorithms presented in this article were implemented in the CAMPARY (CudA Multiple Precision ARithmetic librarY) software available at http://homepages.laas.fr/mmjoldes/campary. The library is implemented in CUDA—an extension of the C language developed by NVIDIA [20] for their GPUs. The algorithms presented are very suitable for the GPU: all basic operations $(+, -, *, /, \sqrt{)}$ conform to the IEEE 754-2008 standard for FP arithmetic for single and double precision; support for the four rounding modes is provided and dynamic rounding mode change is supported without any penalties. The fma instruction is supported for all devices with *Compute Capability* at least 2.0.

In the implementation we use templates for both the number of terms in the expansion and the native type for the terms. In other words, we allow static generation of any input-output precision combinations (e.g. add a double double with a quad-double and store the result on triple-double) and operations with types like single-single, quad-single, etc. are supported. All the functions are defined using __host___device__ specifiers, which allows for the library to be used on both CPU and GPU.

In Table 6 we give some GPU performance measurements for the reciprocal and square root algorithms implemented in CAMPARY compared to the GQD implementation. The tests were performed on a Tesla C2075 GPU, using CUDA 7.0 software architecture, using a single thread of execution. More extensive comparisons, on both CPU and GPU, can be consulted at http:// homepages.laas.fr/mmjoldes/campary.

As a future work we intend to generalize the theoretical analysis of DD and QD addition/multiplication algorithms and thus to be able to provide a full error analysis for these algorithms.

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TABLE 6 GPU Performance in MFlops/s for the Reciprocal, Division and Square Root Algorithms Implemented in CAMPARY versus GQD

		CAMPARY	QD		
d_i, d_o	Alg. 10	Alg. 10+Alg. 5	Alg. 11	Div.	Sqrt.
2, 2 4, 4 8, 8 16, 16 1, 2 2, 4 1, 4	$\begin{array}{c} 0.0501\\ 0.0114\\ 0.0025\\ 0.00031\\ 0.0672\\ 0.012\\ 0.0122\\ \end{array}$	0.0348 0.0083 0.0017 0.00023 0.0419 0.0086 0.0087	$\begin{array}{c} 0.027\\ 0.0063\\ 0.0013\\\\ 0.0338\\ 0.0068\\ 0.0074 \end{array}$	0.0632 0.0044 * 0.102 * 0.2564	0.0495 0.0015 * * * *
4, 2 2, 8 4, 8 4, 16 8, 16	$\begin{array}{c} 0.0501 \\ 0.0030 \\ 0.0027 \\ 0.0004 \\ 0.00039 \end{array}$	$\begin{array}{c} 0.0348 \\ 0.002 \\ 0.0018 \\ 0.00028 \\ 0.00027 \end{array}$	0.0268 0.0016 0.0014 	* * * *	* * * *

 d_i represents the input size (the numerator size for division) and d_o is the size of the computed result (the denominator and the quotient for division). -- error due to the GPU's limited stack size and local memory.

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