MULTIPLE-PRECISION CORRECTLY ROUNDED
NEWTON-COTES QUADRATURE

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Abstract. Numerical integration is an important operation for scientific computations. Although the different quadrature methods have been well studied from a mathematical point of view, the analysis of the actual error when performing the quadrature on a computer is often neglected. This step is however required for certified arithmetics.

We study the Newton-Cotes quadrature scheme in the context of multiple-precision arithmetic and give enough details on the algorithms and the error bounds to enable software developers to write a Newton-Cotes quadrature with bounded error.

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1. Introduction

Numerical integration is an operation that is common in mathematical software (\texttt{intnum} in Pari/GP [2], \texttt{quadrature} in MuPAD [7], \texttt{evalf(Int(...))} in Maple for example). At first glance this is a topic that seems well studied: several quadrature schemes with different convergence properties are known as well as strategies to use them (e.g. adaptive error analysis) or to combine them. When compared to the basic four operations (addition, subtraction, multiplication and division) one notices however the lack of correct rounding, which means the result is not always the floating-point number (see Def. 3.7) closest to the real value in the chosen direction.

Keywords and phrases. numerical integration, correct rounding, multiple-precision, Newton-Cotes.

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Let us illustrate the problem by an example, and compare the values returned respectively by Pari/GP and Maple for $I = \int_{10^6}^{10^6+\pi} (\sin(\cos(t)) - \cos(\sin(t))) \, dt$. On the one hand Pari/GP gives:

```
? \p19
? intnum(t = 10^6, 10^6+Pi, sin(cos(t)) - cos(sin(t)))
%1 = -1.810600390080270954
```

while Maple returns:

```
> Digits:=19: evalf(Int(sin(cos(t)) - cos(sin(t)), t=10^6..10^6 + Pi));
-1.810600390080269775
```

so clearly one software at least is wrong here.

The definition and normalization of rounding modes was a major advance in floating-point arithmetic with respect to the portability and reproducibility of computations. It is a challenging goal to extend this notion of correct rounding to more complex operations like numerical integration.

The error analysis of quadrature methods however is often limited to the mathematical error. In the context of correct rounding this is sadly not sufficient, because a precise bound on the total error is needed to be able to decide of the actual accuracy of the result. Previous works in the field include the study of the adaptive quadrature function of MuPAD [8] and dynamic error control of simple or multiple integrals [1, 6]. What differentiates our work is the careful study of the error term. It is often the case that the roundoff error is merely estimated or sometimes even dismissed by computing with a precision that is “good enough”. Instead, our goal is to give a rigorous formula bounding the total error made in the computation (both the mathematical and roundoff error) in order to be able to guarantee the final result. The use of multiple-precision arithmetic renews the questions that arise for the Newton-Cotes method; in particular higher orders are still up for consideration in this context. If the integral is not exactly representable then we can indeed provide a correctly rounded result, otherwise we can merely compute the value with an absolute error as small as desired.

In this paper $f : [a, b] \to \mathbb{R}$ is the $C^\infty$ function we want to integrate on a finite domain $[a, b]$, and $n$ is the number of evaluation points in the Newton-Cotes method. Let

$$I = \int_a^b f(x) \, dx$$

be the exact value of the integral.

The Newton-Cotes method uses equally-spaced evaluation points in the integration domain, commonly referred to as “abscissas” $x_0, x_1, \ldots, x_{n-1}$:

for $0 \leq i < n$, $x_i = a + ih$ where $h = \frac{b - a}{n - 1}$ is the step.
Since \( x_0 = a \) and \( x_{n-1} = b \), the bounds are used as abscissas and the method is said to be closed. The principle of the method is to approximate the function with the Lagrange interpolating polynomial with respect to the abscissas. The formulas below follow directly from this statement.

For \( i \in \{0, \ldots, n-1\} \), let \( l_i(x) = \prod_{j \neq i} \frac{(x-x_j)}{(x_i-x_j)} \) and \( w_i = \frac{1}{h} \int_a^b l_i(x) \, dx \). The approximated integral is then

\[
I_n = h \sum_{i=0}^{n-1} w_i f(x_i).
\]

The mathematical error \( E_n = I - I_n \) is of the form \( E_n = c_n h^{n+1} f^{(n)}(\zeta) \) for \( n \) even and \( E_n = c_n h^{n+2} f^{(n+1)}(\zeta) \) for \( n \) odd for some \( \zeta \in [a, b] \) (see [3], this will be detailed afterwards).

Firstly we describe the algorithms used in our implementation of the Newton-Cotes quadrature scheme. Then we establish some facts about the error due to the method as well as a few useful lemmas relevant to floating-point arithmetic. These results allow us to proceed to a thorough study of the error made when using the Newton-Cotes quadrature scheme on a computer using floating-point arithmetic. Then we state our main theorem (Th. 3.12). We conclude with some experiments and remarks about the quadrature scheme studied.

### 2. Algorithm for the Computation of the Newton-Cotes Coefficients

In the Newton-Cotes method we distinguish the computation of the coefficients from the quadrature itself, since the coefficients can be precomputed, and reused for several quadratures using the same number of points. For example the composition technique splits the initial integration interval in several parts and applies the same method on each part.

We describe here the algorithm for the computation of the coefficients. The full quadrature algorithm is explained in Section 3.5 together with a discussion about the error.

First we show that the coefficients do not depend on the integration interval. This is true for every linear quadrature scheme, even with non-equally spaced abscissas simply because of the linearity of the integral. We include however the proof only for the Newton-Cotes case because we want to derive the formula for the coefficients.

**Proposition 2.1.** The coefficients of the Newton-Cotes methods do not depend on the integration interval, and are symmetric with respect to the middle of the interval.
Proof. We transform the expression of $w_i$ from Section 1:

$$w_i = \frac{1}{h} \int_a^b l_i(x)dx$$

$$= \int_0^{n-1} l_i(a + xh)dx$$

$$= \int_0^{n-1} \prod_{j \neq i} \frac{(a + xh - x_j)}{(x_i - x_j)}dx$$

$$= \int_0^{n-1} \prod_{j \neq i} \frac{(a + xh - (a + jh))}{h(i - j)}dx$$

$$= \int_0^{n-1} \prod_{j \neq i} \frac{(x - j)}{(i - j)}dx$$

$$= \frac{(-1)^{n-1-i}}{i!(n-1-i)!} \int_0^{n-1} \prod_{j \neq i} (x - j)dx.$$  

The variable change $x \mapsto n - 1 - x$ shows that $w_{n-1-i} = w_i$. □

Let $l^*_i(x) = \prod_{j \neq i} (x - j)$ and $L_i$ the antiderivative of $l^*_i$ such that $L_i(0) = 0$.

Let $u_i = \frac{(-1)^{n-1-i}}{l((n-1-i))} = (-1)^{n-1-i} \frac{n-i}{(n-1)!}$. Then we compute the weights as

$$w_i = u_i L_i(n-1).$$ (1)

From the formula one can notice the weights are rational. They can thus be computed exactly as $w_i = \frac{n_i}{d}$ by Algorithm 1 below. Note that with the $\delta$ factor in $l^*_i$ we ensure at each step that $L_i$ is an integer because in the polynomial integration only division by integers from 2 to $n - 1$ occur. Timings were done on

---

**Algorithm 1 Newton-Cotes coefficients**

1. $\delta \leftarrow \text{lcm}(2, 3, \ldots, n)$
2. $l^*_0 \leftarrow (x - 1)(x - 2) \ldots (x - (n - 1))\delta$
3. **for** $i \leftarrow 0$ to $n - 1$ **do**
4. $L_i \leftarrow \int_0^{n-1} l^*_i(x)dx$
5. $l_{i+1}^* \leftarrow \frac{x^{n-i-1} l^*_i}{x^{n-i}}$
6. $n_i \leftarrow (-1)^{n-1-i} \frac{n-1}{i} L_i$
7. **end for**
8. **return** $(n_0, n_1, \ldots, n_{\lfloor n/2 \rfloor}, d = \delta \cdot (n - 1)!)$
a Pentium-4 computer with MPFR [9] version 2.1.1 and are shown in Figure 1. We estimate the binary complexity of the algorithm as such:

- **line 1:** the size of $\delta$ is $O(n)$ [5] and it can be computed naively as $n$ steps of Euclidean's algorithm between the current least common multiple (lcm) of size $O(n)$ and the current integer of size $O(\log(n))$. The total complexity is thus $O(n^2 \log n)$.
- **line 2:** we build the polynomial product step by step where the coefficient of the current polynomial are of size $O(n \log n)$. The complexity is $O(n^3 \log^2 n)$.
- **line 4 and 5:** those lines are executed $n$ times and consist of $O(n)$ multiplications or divisions of integer of size $O(n \log n)$ by $O(\log n)$, thus a total complexity of $O(n^3 \log^2 n)$.
- **line 6:** the size of $L_i$ is $O(n \log^2 n)$, and the size of $\binom{n-1}{i}$ is $O(n)$, so this step costs $O(n^3 \log^2 n)$. We don’t compute $\binom{n-1}{i}$ at each step, instead we update from the value of the previous step (the cost of the update is negligible).

We get a final time complexity for Algorithm 1 of $O(n^3 \log^2 n)$.

### 3. Error bounds

When performing a numerical integration by means of a Newton-Cotes method, there are two sources of errors to consider: the mathematical error that comes from the method itself, and the roundoff error in the computation which depends on the way we implement the algorithm.
We first give bounds on the mathematical error with elementary proofs. Theorems~3.1 and 3.2 are standard and are included here for the convenience of the reader. A proof of Theorem 3.2 can be found e.g. in [3], pp. 286–287.

3.1. Bounds on the mathematical error

**Theorem 3.1.** For $n$ odd, the Newton-Cotes integration method on $[a, b]$ with $n$ points is exact for polynomials of degree $\leq n$. For $n$ even, it is exact for polynomials of degree $\leq n−1$.

**Proof.** For any $n$, the method is exact for polynomials of degree up to $n−1$ because in that case the Lagrange interpolating polynomial is $f$ exactly.

Let now $n$ be odd. The choice of the evaluation points for the method gives $x_i + x_{n−1−i} = a + b$. Let $g(x) = (x − x_0)(x − x_1)\ldots(x − x_{n−1})$.

$$g\left(\frac{a + b}{2} - x\right) = \prod_{i=0}^{n−1} \left(\frac{a + b}{2} - x - x_i\right)$$

$$= \prod_{i=0}^{n−1} \left(\frac{a + b}{2} - x - (a + b - x_{n−1−i})\right)$$

$$= \prod_{i=0}^{n−1} \left(-\frac{a + b}{2} + x + x_{n−1−i}\right)$$

$$= (-1)^n \prod_{i=0}^{n−1} \left(\frac{a + b}{2} + x - x_{n−1−i}\right)$$

$$= -g\left(\frac{a + b}{2} + x\right)$$

and then $\int_a^b g(x)dx = 0 = \sum_{i=0}^{n−1} w_i g(x_i)$ since $g(x_i) = 0$. The Newton-Cotes method is exact for polynomials of degree $n−1$ at most, and for $g$, which has degree $n$, hence by linearity is exact for all polynomials of degree at most $n$. □

3.2. Peano theorem

In this section we establish the expression of the mathematical error given in Section 1. For this the formalism of the Peano kernel of an integration method is a powerful tool.

For a quadrature method $I : C^{\nu+1}([a, b]) \to \mathbb{R}$ the error $E : f \mapsto \int_a^b f(x)dx - I(f)$ can be seen as a linear function $C^{\nu+1}([a, b]) \to \mathbb{R}$. We have the following result:

**Theorem 3.2.** Define

$$K_\nu(t) = \frac{1}{\nu!} E[x \mapsto (x - t)^\nu]$$
Table 1. Newton-Cotes integration formulas for small \( n \). To simplify the notations we define \( f_i = f(x_i) \).

<table>
<thead>
<tr>
<th>( n )</th>
<th>Name</th>
<th>Formula</th>
<th>( c_n )</th>
</tr>
</thead>
<tbody>
<tr>
<td>2</td>
<td>trapezoidal rule</td>
<td>( I_2 = h \frac{f_0 + f_1}{2} )</td>
<td>( \frac{1}{12} )</td>
</tr>
<tr>
<td>3</td>
<td>Simpson’s rule</td>
<td>( I_3 = \frac{h}{3} (f_0 + 4f_1 + f_2) )</td>
<td>( \frac{1}{90} )</td>
</tr>
<tr>
<td>4</td>
<td>Simpson’s ( \frac{3}{8} ) rule</td>
<td>( I_4 = \frac{3h}{4} (f_0 + 3f_1 + 3f_2 + f_3) )</td>
<td>( \frac{3}{80} )</td>
</tr>
<tr>
<td>5</td>
<td>Boole’s rule</td>
<td>( I_5 = \frac{2}{45}h (7f_0 + 32f_1 + 12f_2 + 32f_3 + 7f_4) )</td>
<td>( \frac{8}{945} )</td>
</tr>
</tbody>
</table>

and

\[
(x - t)^\nu_+ = \begin{cases} 
(x - t)^\nu & \text{if } x > t \\
0 & \text{otherwise.}
\end{cases}
\]

If \( E[p] = 0 \) for all polynomials \( p \) of degree at most \( \nu \) then for \( f \in C^{\nu+1}([a, b]) \),

\[
E[f] = \int_a^b f^{(\nu+1)}(t)K_\nu(t)\,dt.
\]

\( K_\nu \) is called the Peano kernel of order \( \nu \) of \( E \).

**Proof.** Writing the Taylor series associated with \( f \) at origin \( a \):

\[
f(x) = p_\nu(x) + \int_a^x \frac{1}{\nu!}(x-t)^\nu f^{(\nu+1)}(t)\,dt
\]

\[
= p_\nu(x) + \int_a^b \frac{1}{\nu!}(x-t)^\nu f^{(\nu+1)}(t)\,dt,
\]

\[
E[f] = E \left[ \int_a^b \frac{1}{\nu!}(x-t)^\nu f^{(\nu+1)}(t)\,dt \right]
\]

\[
= \int_a^b E \left[ \frac{1}{\nu!}(x-t)^\nu \right] f^{(\nu+1)}(t)\,dt.
\]

This theorem links the mathematical error with the maximal degree of the polynomial the method integrates exactly (its maximal order). In Theorem 3.1 we proved that the order of an \( n \)-points Newton-Cotes method is at least \( n - 1 \) for \( n \) even and \( n \) for \( n \) odd. If we accept that for those orders the corresponding Peano kernel of the Newton-Cotes method does not change sign on \([a, b]\) (see [3], p. 289), then we have a method to compute the coefficient \( c_n \) given in Table 1. With the mean value theorem there exists \( \zeta \in [a, b] \) such that:

\[
E[f] = f^{(\nu+1)}(\zeta) \int_a^b K_\nu(t)\,dt \tag{2}
\]
and we obtain

\[ c_n := \begin{cases} 
\frac{1}{h^{n+1}} \int_a^b K_{n-1}(t) \, dt & \text{if } n \text{ is even,} \\
\frac{1}{h^{n+2}} \int_a^b K_n(t) \, dt & \text{if } n \text{ is odd.}
\end{cases} \]

For example for the 3-points method known as Simpson’s rule, we get

\[ E[f] = \int_a^b f(x) \, dx - \frac{b-a}{6} \left[ \left( f(a) + 4f \left( \frac{a+b}{2} \right) + f(b) \right) \right], \]

\[ K_3(t) = \frac{1}{6} E[x \mapsto (x-t)^3], \]

\[ 6K_3(t) = \int_a^b (x-t)^3 \, dx - \frac{b-a}{6} \left[ 4 \left( \frac{a+b}{2} - t \right)^3 + (b-t)^3 \right] \]

\[ = \int_t^b (x-t)^3 \, dx - \frac{b-a}{6} \left[ 4 \left( \frac{a+b}{2} - t \right)^3 + (b-t)^3 \right] \]

\[ = \begin{cases} 
\frac{(b-t)^4}{4} - \frac{b-a}{6} \left[ 4 \left( a+b \right)^3 - 4 \left( a+b \right)^3 + (b-t)^3 \right] & \text{if } t < \frac{a+b}{2} \\
\frac{(b-t)^4}{4} - \frac{b-a}{6} (b-t)^3 & \text{if } t \geq \frac{a+b}{2},
\end{cases} \]

\[ \int_a^b K_3(t) \, dt = \int_a^b \frac{(b-t)^4}{24} - \frac{b-a}{36} (b-t)^3 dt - \int_a^{\frac{a+b}{2}} \frac{(b-a)}{9} \left( \frac{a+b}{2} - t \right)^3 dt \]

\[ = \frac{(b-a)^5}{120} - \frac{(b-a)^5}{144} - \frac{(b-a)^5}{576} = \frac{1}{90} \left( \frac{b-a}{2} \right)^5 \]

and we find the value \( c_3 = -\frac{1}{90} \) given in the third row of Table 1.

3.3. Upper bound for the \( c_n \) coefficients

In order to be able to give an absolute bound on the mathematical error we need to bound the \( c_n \) coefficients. We detail here the bound and the proof for \( n \) even, and give only the result for \( n \) odd. Recall that for \( n \) even, the Newton-Cotes method is exact for all polynomials up to degree \( n - 1 \).

Take for \( f \) a monic degree-\( n \) polynomial \( p \) in equation (2) to get:

\[ E_n[p] = p^{(n)}(\zeta) \int_a^b K_{n-1}(t) \, dt \]

and then \( c_n = \frac{E_n[p]}{p^{(n)}(\zeta)} \), since \( p^{(n)}(\zeta) = n! \).
In particular for \( p(x) = (x - x_0)(x - x_1) \ldots (x - x_{n-1}) \), we have

\[
E_n[p] = \int_a^b p(x)dx - \sum_{i=0}^{n-1} w_i p(x_i) = \int_a^b p(x)dx
\]

so it is enough to bound \( \left| \int_a^b p(x)dx \right| \) in order to bound \( c_n \). We will use repeatedly the following simple lemma:

**Lemma 3.3.** For \((u, x, v) \in \mathbb{R}^3\) such that \( u \leq x \leq v \),

\[
|x - u||x - v| \leq \left( \frac{v - u}{2} \right)^2
\]

**Proof.**

\[
|x - u||x - v| = (x - u)(v - x)
= \left( x - \frac{u + v}{2} \right) \left( \frac{v - u}{2} - \left( x - \frac{u + v}{2} \right) \right)
= \left( \frac{v - u}{2} \right)^2 - \left( x - \frac{u + v}{v} \right)^2
\leq \left( \frac{v - u}{2} \right)^2.
\]

**Proposition 3.4.**

\[\forall x \in [a, b], \quad |p(x)| \leq \frac{h^n(n - 1)!}{4}.\]

**Proof.** Let \( x \in [a, b] \) such that \( p(x) \neq 0 \). Then there is \( i_0 \in \{0, \ldots, n-2\} \) such that \( x \in [x_{i_0}, x_{i_0+1}] \), and thus

\[
|x - x_{i_0}| |x - x_{i_0+1}| \leq \frac{h^2}{4}, \quad \text{(Lem. 3.3)}
\]

Then

\[
|p(x)| \leq \prod_{i=0}^{n-1} |x - x_i|
\leq \frac{h^2}{4} \prod_{i \neq \{i_0, i_{i_0+1}\}} |x - x_i|
\leq \frac{h^2}{4} \left[ \prod_{i > i_0+1} (i - i_0)h \right] \left[ \prod_{i < i_0} (i_0 + 1 - i)h \right],
\]

\[
\left[ \prod_{i > i_0+1} (i - i_0)h \right] \left[ \prod_{i < i_0} (i_0 + 1 - i)h \right] \leq (n - 1 - i_0)!(i_0 - 1)!h^{n-2}
\leq (n - 1)!h^{n-2},
\]

\[
|p(x)| \leq \frac{h^n(n - 1)!}{4}.
\]

\[\square\]
On the other hand we have \( b - a = (n - 1)h \), which yields

\[
|E_n[p]| = \left| \int_a^b p(x)dx \right| \leq \int_a^b |p(x)|dx \leq \frac{h^{n+1}(n-1)!}{4(n-1)}.
\]

With \( E_n[p] = c_n h^{n+1}n! \) we get

\[
|c_n| \leq \frac{n-1}{4n} \leq \frac{1}{4}.
\]

For \( n \) odd we take \( p(x) = (x-x_0)(x-x_1)\ldots(x-x_{n-1})(x - \frac{a+b}{2}) \); we have the evaluation points and the middle of the interval as zeroes of \( p \).

Similar computations give:

\[
|p(x)| \leq \frac{h^{n+1}(n-1)!}{8(n-1)}
\]

then

\[
\left| \int_a^b p(x)dx \right| \leq \frac{h^{n+2}(n-1)!}{8(n-1)^2}
\]

and with \( E_n = c_n h^{n+2}p^{(n+1)} \) we get

\[
|c_n| \leq \frac{(n-1)^2}{8n(n+1)} \leq \frac{1}{8}.
\]

3.4. ULP calculus

For the error analysis of Algorithm 2, we need a few useful lemmas concerning the “ulp\(^1\) calculus”, as well as some definitions. The floating-point numbers are represented with radix 2 (this could be generalized for any radix but radix 2 is simpler and is natural on computers). For this section, \( p \) is the working precision, and we assume that the exponent range is unbounded, which ensures that no underflow or overflow occur.

**Definition 3.5** (exponent). For a non-zero real number \( x \) we define \( E(x) := 1 + \lfloor \log_2 |x| \rfloor \), such that \( 2^{E(x)-1} \leq |x| < 2^{E(x)} \).

**Definition 3.6** (ulp). For a non-zero real number \( x \) we define \( \text{ulp}(x) := 2^{E(x)-p} \).

**Definition 3.7** (floating-point number). A real number \( x \) is a floating-point number with precision \( p \) if it can be written \( x = m \cdot 2^e \) where \( m,e \) are integers and \( |m| < 2^p \).

For a real \( x \neq 0 \) and a working precision \( p \) we always have \( 2^{p-1}\text{ulp}(x) \leq |x| < 2^p\text{ulp}(x) \). If \( x \) is a floating-point number, then \( \text{ulp}(x) \) is the weight of the least significant bit — zero or not — in the \( p \)-bit mantissa of \( x \). For all real \( x \), \( \text{ulp}(x) \) is always greater than zero by definition.

\(^1\)Unit in the last place.
Lemma 3.8. If $c \neq 0$ and $x \neq 0$ then $c \cdot \text{ulp}(x) < 2 \cdot \text{ulp}(cx)$.

Proof. If $c < 0$ it is void. By definition of ulp(x) we have for all $c > 0$:

$$2^{p-1}\text{ulp}(x) \leq |x|$$

and

$$|cx| < 2^p\text{ulp}(cx)$$

so

$$c \cdot 2^{p-1}\text{ulp}(x) \leq |cx| < 2^p\text{ulp}(cx).$$

Lemma 3.9. Assuming no underflow (flush to zero) occurs then in all rounding modes for a non zero real $x$ we have:

$$\text{ulp}(x) \leq \text{ulp}(\circ(x)),$$

where $\circ(x)$ is the rounding of $x$ in the chosen mode with an unbounded exponent range.

Proof. We have $2^{E(x)-1} \leq |x| < 2^{E(x)}$ and $\text{ulp}(x) = 2^{E(x)-p}$. After rounding we get $2^{E(x)-1} \leq |\circ(x)| \leq 2^{E(x)}$ since $2^{E(x)}$ and $2^{E(x)-1}$ are exactly representable, therefore $\text{ulp}(\circ(x)) \geq 2^{E(x)-p} \geq \text{ulp}(x)$. \hfill \Box

Lemma 3.10. Let $x$ be a non-zero real and $\circ(x)$ its rounding to nearest on $p$ bits. Then $|x| \leq (1 + 2^{-p})|\circ(x)|$.

Proof. By definition of rounding to nearest we have

$$|x - \circ(x)| \leq \frac{1}{2}\text{ulp}(\circ(x)) \leq \frac{1}{2}2^{1-p}|\circ(x)|,$$

$$|x| \leq |\circ(x)| + 2^{-p}|\circ(x)|.$$ \hfill \Box

Lemma 3.11. Let $a$ and $b$ be two non-zero floating-point numbers of the same sign and precision $p$ then in all rounding modes

$$\text{ulp}(a) + \text{ulp}(b) \leq \frac{3}{2}\text{ulp}(\circ(a + b)).$$

Proof. It suffices to consider the case where $a$ and $b$ are positive. The definition of ulp gives:

$$2^{p-1}\text{ulp}(a) \leq a < 2^p\text{ulp}(a),$$

$$2^{p-1}\text{ulp}(b) \leq b < 2^p\text{ulp}(b)$$

thus

$$2^{p-1}[\text{ulp}(a) + \text{ulp}(b)] \leq a + b < 2^p[\text{ulp}(a) + \text{ulp}(b)].$$

If $\text{ulp}(a) = \text{ulp}(b)$ we get

$$2^p\text{ulp}(a) \leq a + b < 2^{p+1}\text{ulp}(a)$$

and therefore $\text{ulp}(\circ(a + b)) \geq \text{ulp}(a + b) \geq 2\text{ulp}(a) = \text{ulp}(a) + \text{ulp}(b)$ (Lem. 3.9) and the lemma holds.
Otherwise we can assume without loss of generality that \( \text{ulp}(a) > \text{ulp}(b) \), that is \( \text{ulp}(a) \geq 2 \cdot \text{ulp}(b) \). We deduce:

\[
\text{ulp}(a) + \text{ulp}(b) \leq \frac{3}{2} \text{ulp}(a),
\]

and together with \( \text{ulp}(\circ(a + b)) \geq \text{ulp}(a + b) \geq \text{ulp}(a) \) (Lem. 3.9) this concludes the proof. \( \square \)

**Example.** Let \( p = 4 \) and choose rounding to nearest: \( a = 1.010 \times 2^{-3} \), \( b = 0.1001 \times 2^{-4} \) in binary notation.

\[
a + b = 1.1101, \quad \circ(a + b) = 1.110,
\]

\[
\text{ulp}(a) + \text{ulp}(b) = 2^{-3} + 2^{-4} = \frac{3}{2} 2^{-3} = \frac{3}{2} \text{ulp}(\circ(a + b)).
\]

### 3.5. Roundoff errors

**Algorithm 2** Newton-Cotes integration

<table>
<thead>
<tr>
<th>Input: ( \hat{a}, \hat{b}, f, n, {n_i}, d. )</th>
</tr>
</thead>
<tbody>
<tr>
<td>Output: ( \hat{I}. )</td>
</tr>
<tr>
<td>1: for ( i \leftarrow 0 ) to ( n - 1 ) do</td>
</tr>
<tr>
<td>2: ( \hat{t} \leftarrow \circ(\hat{a} \times (n - i - 1)) )</td>
</tr>
<tr>
<td>3: ( \hat{u} \leftarrow \circ(\hat{b} \times i) )</td>
</tr>
<tr>
<td>4: ( \hat{v} \leftarrow \circ(\hat{t} + \hat{u}) )</td>
</tr>
<tr>
<td>5: ( \hat{x}_i \leftarrow \circ(\hat{v}/(n - 1)) )</td>
</tr>
<tr>
<td>6: ( \hat{z} \leftarrow \circ(f(\hat{x}_i)) )</td>
</tr>
<tr>
<td>7: ( \hat{y}_n \leftarrow \circ(\hat{z} \times n_i) )</td>
</tr>
<tr>
<td>8: end for</td>
</tr>
<tr>
<td>9: ( \hat{S} \leftarrow \text{sum}(\hat{y}_i, i = 0 \ldots n - 1) \quad \triangleright \text{with Demmel and Hida algorithm [4]} )</td>
</tr>
<tr>
<td>10: ( \hat{U} \leftarrow \circ(\hat{S}/d(n - 1)) )</td>
</tr>
<tr>
<td>11: ( \hat{D} \leftarrow \circ(\hat{b} - \hat{a}) )</td>
</tr>
<tr>
<td>12: return ( \circ(\hat{D}\hat{U}) )</td>
</tr>
</tbody>
</table>

In order to provide an error bound on the numerical result given by the Newton-Cotes method, we need to have a step-by-step look into Algorithm 2.

This step is often neglected when doing numerical integration, where error analysis stops right after stating the well known bound for the mathematical error. In fact, the experiment illustrated in Figure 2 shows that much remains to be done to control the error on the result. For this section we denote by \( \hat{x} \) the value actually computed (i.e. with all roundoff errors) for a given “exact” value \( x \) that would be computed with an infinite precision from the beginning of the algorithm.

In Algorithm 2, we set \( p \) as the working precision expressed in the number of bits of the mantissa, the parameters \( \hat{a} \) and \( \hat{b} \) are assumed to be the rounded to nearest \( p \)-bit floating-point numbers of their exact counterpart \( a \) and \( b \). For the
error bound computation the algorithm needs additional parameters given by the user: an upper bound $M$ of $|f^{(n)}|$ on $[a, b]$ if $n$ is even, or an upper bound of $|f^{(n+1)}|$ if $n$ is odd; $m$ an upper bound of $|f'|$ on $[a, b]$.

In the rest of this section we will prove our main theorem:

**Theorem 3.12.** Let $\delta_{\hat{y}_i} = 6|n_i| \cdot m \cdot \text{ulp}(\hat{x}_i) + \frac{3}{2} \text{ulp}(\hat{y}_i)$ where $\hat{x}_i$ and $\hat{y}_i$ are defined in Algorithm 2. When computing the numerical quadrature of $f$ using Algorithm 2 with $a$ and $b$ of the same sign the total error on the result is bounded by:

$$E_{tot} = \left( \frac{45}{2} + 21 \cdot 2^{-p} \right) \text{ulp}(\hat{I}) + \frac{1}{2} |\hat{I}| \left( \text{ulp}(\hat{D}) + \text{ulp}(\hat{a}) + \text{ulp}(\hat{b}) \right) + 3(1 + 2^{-p})n \cdot \hat{D} \frac{\max(\delta_{\hat{y}_i})}{d(n-1)} + \begin{cases} 
\frac{1}{8} \left( \frac{b-a}{n-1} \right)^{n+2} M & \text{if } n \text{ is odd}, \\
\frac{1}{4} \left( \frac{b-a}{n-1} \right)^{n+1} M & \text{otherwise}.
\end{cases}$$

The algorithm can be analyzed in several steps:

1. The computation of the weights $w_i$, $i \in [0, n-1]$ of the method. For Newton-Cotes, those weights are rational and computed exactly: $w_i = \frac{n_i}{d}$ where $n_i, d \in \mathbb{Z}$, so no rounding error occurs at this step.
(2) The computation of $x_i$. This is done at line 5 of Algorithm 2:

$$
\hat{x}_i = \circ \left( \circ(\circ((n - 1 - i) \cdot \hat{a}) + \circ(i \cdot \hat{b})) \right).
$$

In order to simplify the notations we write $t = (n - i - 1) \alpha$, $u = i \cdot \beta$ and their inexact counterparts $\hat{t} = \circ((n - i - 1)\hat{a})$, $\hat{u} = \circ(i \hat{b})$. If $b = 0$ or $i = 0$ the error on $\hat{u}$ is zero. Otherwise the error estimation yields:

$$
|\circ(i \cdot \hat{b}) - i \beta| \leq \frac{1}{2} \text{ulp}(\circ(i \hat{b})) + \frac{i}{2} \text{ulp}(\hat{b})
$$

$$
\leq \frac{3}{2} \text{ulp}(\circ(i \hat{b})) = \frac{3}{2} \text{ulp}(\hat{u}). \quad \text{(Lems. 3.8 and 3.9)}
$$

Similarly if $a = 0$ or $n - i - 1 = 0$ the error on $\hat{t}$ is zero, and otherwise we get $|\hat{t} - t| \leq \frac{3}{4} \text{ulp}(\hat{t})$.

Since $a$ and $b$ have the same sign, $\hat{t}$ and $\hat{u}$ also have the same sign and we can use Lemma 3.11. This assumption is not restrictive in practice as we can split the integration interval at 0, and apply our error bound on each part to get the consolidated error bound on $[a, b]$. Moreover assume without loss of generality that $0 \leq a < b$, which gives $0 \leq \hat{u} \leq \hat{b}$; then:

$$
|\hat{v} - v| \leq \frac{1}{2} \text{ulp}(\hat{v}) + \frac{3}{2} (\text{ulp}(\hat{t}) + \text{ulp}(\hat{u}))
$$

$$
\leq \frac{11}{4} \text{ulp}(\hat{v}). \quad \text{(Lem. 3.11)}
$$

Taking into account the error coming from the division by $n - 1$ we get:

$$
\delta_{\hat{x}_i} = |x_i - \hat{x}_i| \leq \frac{1}{2} \text{ulp}(\hat{x}_i) + \frac{11}{4(n - 1)} \text{ulp}(\hat{v})
$$

$$
\leq \frac{11}{4} \text{ulp}(\hat{x}_i) + \frac{11}{2} \text{ulp}(\hat{x}_i) \quad \text{(Lems. 3.8 and 3.9)}
$$

$$
\leq 6 \cdot \text{ulp}(\hat{x}_i).
$$

(3) The computation of $f(x_i)$. We assume we have an implementation of $f$ with correct rounding, and we call the function $f$ requesting the rounding to nearest of the exact value with precision $p$. Such correctly rounded implementations of mathematical functions with arbitrary precision on the result can be found for example in MPFR [9] for non-trivial functions like exp, sin, arctan and numerous others.

With the already estimated error on $\hat{x}_i$ we have:

$$
|f(\hat{x}_i) - f(x_i)| = |f'(\theta_i)(\hat{x}_i - x_i)|, \quad \theta_i \in [\min(x_i, \hat{x}_i), \max(x_i, \hat{x}_i)]
$$

and with an upper bound on $f'$ we can bound this error absolutely. Let $\hat{f}_i = \circ(f(\hat{x}_i))$ be the floating-point number computed. At this step we
MULTIPLE-PRECISION CORRECTLY ROUNDED QUADRATURE

now have:

\[ |\hat{f}_i - f(x_i)| \leq |f'(\theta_i)(\hat{x}_i - x_i)| + \frac{1}{2} \text{ulp}(\hat{f}_i) \]
\[ \leq 6m \cdot \text{ulp}(\hat{x}_i) + \frac{1}{2} \text{ulp}(\hat{f}_i). \]

(4) The computation of the \( y_i = f(x_i) \cdot n_i \). The accumulated error so far is:

\[ |\hat{y}_i - y_i| \leq |n_i| \cdot |\hat{f}_i - f_i| + \frac{1}{2} \text{ulp}(\hat{y}_i) \]
\[ \leq 6|n_i| \cdot m \cdot \text{ulp}(\hat{x}_i) + \frac{3}{2} \text{ulp}(\hat{y}_i) = \delta_{\hat{y}_i}. \] (Lems. 3.8 and 3.9)

**Remark.** When bounding the error on \( \hat{x}_i, \hat{f}_i \) as well as \( \hat{y}_i \), the term with \( \text{ulp}(\hat{x}_i) \) vanishes if the error on \( \hat{x}_i \) is zero. One can easily show that with our assumption that no underflow occurs, if \( \hat{x}_i = 0 \) then the error on \( \hat{x}_i \) is zero (i.e. \( x_i = 0 \)) and the ill-defined quantity \( \text{ulp}(\hat{x}_i) \) vanishes. For the error bound we keep track of only \( \max(\delta_{\hat{y}_i}) \).

(5) Summation of the \( y_i \)'s: this is done with Demmel and Hida summation algorithm \([4]\), which guarantees an error of at most 1.5 ulp on the final result. This algorithm uses a larger working precision \( p' \approx p + \log_2(n) \).

Let \( S = \sum_{i=0}^{n-1} y_i \).

\[ |\hat{S} - S| \leq \frac{3}{2} \text{ulp}(\hat{S}) + n \cdot \max(\delta_{\hat{y}_i}). \]

(6) Division of \( S \) by \( d(n-1) \): \( U = \frac{S}{d(n-1)} \). The computation of \( d(n-1) \) is done with integer arithmetic and is exact. The error at this step is thus:

\[ |\hat{U} - U| \leq \frac{1}{2} \text{ulp}(\hat{U}) + \frac{3}{2d(n-1)} \text{ulp}(\hat{S}) + \frac{n}{d(n-1)} \max(\delta_{\hat{y}_i}) \]
\[ \leq \frac{7}{2} \text{ulp}(\hat{U}) + \frac{n}{d(n-1)} \max(\delta_{\hat{y}_i}). \] (Lems. 3.8 and 3.9)

(7) Multiplication by \( b-a \): \( I = (b-a)U \). We note \( D = b-a \) and \( \hat{D} = \circ(\hat{b}-\hat{a}) \).

\[ |\hat{D} - D| \leq \frac{1}{2} \left[ \text{ulp}(\hat{D}) + \text{ulp}(\hat{a}) + \text{ulp}(\hat{b}) \right]. \]

We have by hypothesis

\[ \hat{b} \geq b - \frac{1}{2} \text{ulp}(\hat{b}) \]
\[ \hat{a} \leq a + \frac{1}{2} \text{ulp}(\hat{a}) \]
where $\text{ulp}(0) = 0$ by convention and therefore

$$
\hat{b} - \hat{a} \geq b - a - \frac{1}{2} \left( \text{ulp}(\hat{a}) + \text{ulp}(\hat{b}) \right)
$$

$$
\geq b - a - \text{ulp}(\hat{b}) .
$$

On the other hand we know $\hat{b} \geq \hat{a}$ and we further discard the case $\hat{a} = \hat{b}$ because it is of no practical interest: in this case the current precision $p$ is not even sufficient to decide if $a = b$. We can however still compute an error bound with the knowledge of $m$, $\hat{a}$ and one call to $f(\hat{a})$.

So we may assume $\hat{b} > \hat{a}$ and we have $\hat{b} - \hat{a} \geq \max(\frac{1}{2}\text{ulp}(\hat{b}), b - a - \text{ulp}(\hat{b}))$ which gives $\text{ulp}(\hat{b}) \leq 2(\hat{b} - \hat{a})$; then

$$
D \leq \hat{b} - \hat{a} + \text{ulp}(\hat{b}) \leq 3(\hat{b} - \hat{a}) \leq 3(1 + 2^{-p})\hat{D} . \quad (\text{Lem. 3.10}) \quad (3)
$$

If we put all the results and bounds gathered so far, we can reach the following final error on $\hat{I} = \circ(\hat{D}\hat{U})$:

$$
|\hat{I} - I| \leq \frac{1}{2} \text{ulp}(\hat{I}) + |\hat{D}\hat{U} - D \cdot U|
$$

$$
\leq \frac{1}{2} \text{ulp}(\hat{I}) + |\hat{U}| \cdot |\hat{D} - D| + |D| \cdot |\hat{U} - U| 
$$

$$
\leq \frac{1}{2} \text{ulp}(\hat{I}) + |\hat{U}| \cdot |\hat{D} - D| + 3(1 + 2^{-p})|\hat{D}| \cdot |\hat{U} - U| \quad \text{[Inequality (3)]}
$$

$$
\leq \frac{1}{2} \text{ulp}(\hat{I}) + |\hat{U}| \cdot |\hat{D} - D|
$$

$$
+ 3(1 + 2^{-p})|\hat{D}| \left( \frac{7}{2} \text{ulp}(\hat{U}) + \frac{n}{d(n - 1)} \max(\delta_{\hat{y}_i}) \right)
$$

$$
\leq \left( \frac{43}{2} + 21 \cdot 2^{-p} \right) \text{ulp}(\hat{I}) + |\hat{U}| \cdot |\hat{D} - D|
$$

$$
+ 3(1 + 2^{-p})n \cdot \hat{D} \max(\delta_{\hat{y}_i}) \quad \text{(Lems. 3.8 and 3.9)}
$$

$$
\leq \left( \frac{45}{2} + 21 \cdot 2^{-p} \right) \text{ulp}(\hat{I}) + \frac{1}{2} |\hat{U}| \left( \text{ulp}(\hat{a}) + \text{ulp}(\hat{b}) \right)
$$

$$
+ 3(1 + 2^{-p})n \cdot \hat{D} \max(\delta_{\hat{y}_i}) . \quad \text{(Lems. 3.8 and 3.9)}
$$

This bound for the error is satisfactory for use in the algorithm, because it is made of quantities that we can compute before the algorithm is started ($p$, $n$), or which are naturally computed in the flow of the algorithm ($\hat{I}$, $\hat{U}$, $\hat{a}$, $\hat{D}$, $d$, $\delta_{\hat{y}_i}$).
For the final error bound we need to add a bound on the mathematical error:

\[
E_{\text{math}} \leq \begin{cases} 
\frac{1}{8} \frac{(b-a)}{(n-1)}^{n+2} M & \text{if } n \text{ is odd,} \\
\frac{1}{4} \frac{(b-a)}{(n-1)}^{n+1} M & \text{otherwise}
\end{cases}
\]  

which is easily computed as well. While computing the error bound we carefully choose for every operation the directed rounding mode which ensures that the computed bound is larger than the theoretical error bound.

4. Experiments

Algorithm 2 was implemented using the MPFR library [9]. In addition to the result of the integration, the program gives an error bound on the computed result split in two terms:

1. the mathematical error, whose expression is given in equation (4);
2. the “roundoffs” error \( E_{\text{roundoff}} = E_{\text{tot}} - E_{\text{math}} \).

For our experiments we choose a function and an integration domain where the exact value is known, so that we can measure precisely the actual error of the computation (denoted by \( E_{\text{meas}} \)). Figure 3 shows the different errors when computing the integral \( I = \int_{0}^{3} e^x \, dx \) with 113 bits of working precision, the number of evaluation points varying from 2 to 30.
Figure 4. Optimal values of the number $n$ of points for several working precisions (experimental data gathered with $\int_0^3 e^x \, dx$).

The mathematical error decreases rapidly but it appears clearly that it is well compensated by the roundoff errors as soon as more than about 15 evaluation points are used, for the considered function and parameters. The theoretical gain of increasing the order of the method is lost. Figure 4 gives the smallest value of the number of points for which the mathematical error is inferior to the roundoff errors, for different working precisions chosen. This is commonly interpreted as the optimal value of $n$ in the following sense: for higher values of $n$ the benefit of an higher order method is lost in the noise of the roundoff error, and for smaller values the accuracy on the evaluation of the function is not exploited to its fullest. Although the coefficients generating algorithm is slow for high values of $n$, no particular attempts were made to optimize it yet; this is motivated partly by the slow growth in Figure 4 (the other reasons being the numerical instability discussed below as well as the possibility to use composition).

The bound on the total error as given by the algorithm is somehow close to the measured error. In the experimental data we observe a maximal ratio of about 46,000 – which seems to be huge, but with a logarithmic scale it means we lost a mere 16 bits of precision by our estimation. In particular this means our algorithm is not too grossly pessimistic.

The numerical instability of the method when $n$ grows is not surprising, and not new either. The fact that negative coefficients appear in the formula as soon as $n \geq 8$ partly explains this fact which is demonstrated here. Considering the smoothness of the function chosen for the experiment, the instability is to be attributed to the method. If you read Figure 4 from the other point of view, the
required working precision for higher orders increases rapidly. Small values of \( n \) are therefore recommended for the Newton-Cotes quadrature method.

5. Conclusion and future work

The Newton-Cotes quadrature scheme is the simplest numerical quadrature method, which made it the natural candidate for a detailed study. We were able to provide a rigorous analysis of the method that is self-contained and covers every aspect that is relevant to an implementation, that is, the description of the algorithms and the establishment of proven error bounds.

However the Newton-Cotes family of quadrature methods were not a goal per se but rather a proof of concept that such a study of the error is feasible and indeed desirable. It is planned to perform the same kind of work with other quadrature schemes, notably the Gauss-Legendre methods (which have order \( 2n \) for \( n \) points and are numerically more stable). These analyses might serve as a mathematical foundation of a correctly rounded quadrature library.

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References