

Accurately Computing $\log(1 - \exp(-|a|))$ Assessed by the **Rmpfr** package

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Abstract

In this note, we explain how $f(a) = \log(1 - e^{-a}) = \log(1 - \exp(-a))$ can be computed accurately, in a simple and optimal manner, building on the two related auxiliary functions **log1p(x)** ($= \log(1 + x)$) and **expm1(x)** ($= \exp(x) - 1 = e^x - 1$). The cutoff, a_0 , in use in R since 2004, is shown to be optimal both theoretically and empirically, using **Rmpfr** high precision arithmetic.

Keywords: Accuracy, Cancellation Error, R, MPFR, Rmpfr.

1. Introduction: Not **log()** nor **exp()**, but **log1p()** and **expm1()**

In applied mathematics, it has been known for a very long time that direct computation of $\log(1 + x)$ suffers from severe cancellation (in “ $1 + x$ ”) whenever $|x| \ll 1$, and for that reason, we have provided **log1p(x)** in R, since R version 1.0.0 (released, Feb. 29, 2000). Similarly, **log1p()** has been provided by C math libraries and has become part of C language standards around the same time, see, for example, [IEEE and Open Group \(2004\)](#).

Analogously, since R 1.5.0 (April 2002), the function **expm1(x)** computes $\exp(x) - 1 = e^x - 1$ accurately also for $|x| \ll 1$, where $e^x \approx 1$ is (partially) cancelled by “ -1 ”.

In both cases, a simple solution for small $|x|$ is to use a few terms of the Taylor series, as

$$\log1p(x) = \log(1 + x) = x - x^2/2 + x^3/3 - \dots, \text{ for } |x| < 1, \quad (1)$$

$$\expm1(x) = \exp(x) - 1 = x + x^2/2! + x^3/3! + \dots, \text{ for } |x| < 1, \quad (2)$$

and $n!$ denotes the factorial.

We have found, however, that in some situations, the use of **log1p()** and **expm1()** may not be sufficient to prevent loss of numerical accuracy. The topic of this note is to analyze the important case of computing $\log(1 - e^x) = \log(1 - \exp(x))$ for $x < 0$, computations needed in accurate computations of the beta, gamma, Weibull and logistic distributions, and even for the logit link function in logistic regression. For the beta and gamma distributions, see, for example, [DiDonato and Morris \(1992\)](#)¹, and further references mentioned in R’s **?pgamma** and **?pbeta** help pages. For the logistic distribution, $F_L(x) = \frac{e^x}{1+e^x}$, the inverse, aka quantile function is $q_L(p) = \text{logit}(p) := \log \frac{p}{1-p}$. If the argument p is provided on the log scale,

¹In the Fortran source, file “708”, also available as <http://www.netlib.org/toms/708>, the function **ALNREL()** computes **log1p()** and **REXP()** computes **expm1()**.

$\tilde{p} := \log p$, hence $\tilde{p} \leq 0$, we need

$$\text{qlogis}(\tilde{p}, \text{log.p} = \text{TRUE}) = q_L(e^{\tilde{p}}) = \text{logit}(e^{\tilde{p}}) = \log \frac{e^{\tilde{p}}}{1 - e^{\tilde{p}}} = \tilde{p} - \log(1 - e^{\tilde{p}}), \quad (3)$$

and the last term is exactly the topic of this note.

2. `log1p()` and `expm1()` for $\log(1 - \exp(x))$

Contrary to what one would expect, for computing $\log(1 - e^x) = \log(1 - \exp(x))$ for $x < 0$, neither

$$\log(1 - \exp(x)) = \log(-\text{expm1}(x)), \quad \text{nor} \quad (4)$$

$$\log(1 - \exp(x)) = \text{log1p}(-\exp(x)), \quad (5)$$

are uniformly sufficient for numerical evaluation. In (5), when x approaches 0, $\exp(x)$ approaches 1 and loses accuracy. In (4), when x is large, $\text{expm1}(x)$ approaches -1 and similarly loses accuracy. Because of this, we will propose to use a function `log1mexp(x)` which uses either `expm1` (4) or `log1p` (5), where appropriate. Already in R 1.9.0 ([R Development Core Team \(2004\)](#)), we have defined the macro `R_D_LExp(x)` to provide these two cases automatically².

To investigate the accuracy losses empirically, we make use of the R package **Rmpfr** for arbitrarily accurate numerical computation, and use the following simple functions:

```
R> library(Rmpfr)
R> t3.11e <- function(a)
  {
    c(def = log(1 - exp(-a)),
      expm1 = log(-expm1(-a)),
      log1p = log1p(-exp(-a)))
  }
R> ##' The relative Error of log1mexp computations:
R> relE.11e <- function(a, precBits = 1024) {
  stopifnot(is.numeric(a), length(a) == 1, precBits > 50)
  da <- t3.11e(a) ## double precision
  a. <- mpfr(a, precBits=precBits)
  ## high precision *and* using the correct case:
  mMa <- if(a <= log(2)) log(-expm1(-a.)) else log1p(-exp(-a.))
  structure(as.numeric(1 - da/mMa), names = names(da))
}
```

where the last one, `relE.11e()` computes the relative error of three different ways to compute $\log(1 - \exp(-a))$ for positive a (instead of computing $\log(1 - \exp(x))$ for negative x).

```
R> a.s <- 2^seq(-55, 10, length = 256)
R> ra.s <- t(sapply(a.s, relE.11e))
R> cbind(a.s, ra.s) # comparison of the three approaches
```

	a.s	def	expm1	log1p
[1,]	2.7756e-17	-Inf	-7.9755e-17	-Inf
[2,]	3.3119e-17	-Inf	-4.9076e-17	-Inf
[3,]	3.9520e-17	-Inf	-7.8704e-17	-Inf
[4,]	4.7157e-17	-Inf	-4.5998e-17	-Inf

²look for “ $\log(1 - \exp(x))$ ” in <http://svn.r-project.org/R/branches/R-1-9-patches/src/nmath/dpq.h>

```

[5,] 5.6271e-17 1.8162e-02 -7.3947e-17 1.8162e-02
[6,] 6.7145e-17 1.3504e-02 -4.4921e-17 1.3504e-02
[7,] 8.0121e-17 8.8009e-03 -1.2945e-17 8.8009e-03
.....
.....
[251,] 4.2329e+02 1.0000e+00 1.0000e+00 -3.3151e-17
[252,] 5.0509e+02 1.0000e+00 1.0000e+00 2.9261e-17
[253,] 6.0270e+02 1.0000e+00 1.0000e+00 1.7377e-17
[254,] 7.1917e+02 1.0000e+00 1.0000e+00 -4.7269e-12
[255,] 8.5816e+02 1.0000e+00 1.0000e+00 1.0000e+00
[256,] 1.0240e+03 1.0000e+00 1.0000e+00 1.0000e+00

```

This is revealing: Neither method, `log1p` or `expm1`, is uniformly good enough. Note that for large a , the relative errors evaluate to 1. This is because all three double precision methods give 0, *and* that is the best approximation in double precision (but not in higher `mpfr` precision), hence no problem at all, and we can restrict ourselves to smaller a (smaller than about 710, here).

What about really small a 's?

```

R> t3.11e(1e-20)

      def      expm1      log1p
      -Inf -46.052      -Inf

R> as.numeric(t3.11e(mpfr(1e-20, 256)))

[1] -46.052 -46.052 -46.052

```

so, indeed, the `expm1` method is absolutely needed here.

Figure 1 visualizes the relative errors of the three methods. Note that the default basically gives the maximum of the two methods' errors, whereas the final `log1mexp()` function will have (approximately) minimal error of the two.

```

R> matplot(a.s, abs(ra.s), type = "l", log = "xy",
           col=cc, lty=lt, lwd=ll, xlab = "a", ylab = "", axes=FALSE)
R> legend("top", leg, col=cc, lty=lt, lwd=ll, bty="n")
R> draw.machEps <- function(alpha.f = 1/3, col = adjustcolor("black", alpha.f)) {
  abline(h = .Machine$double.eps, col=col, lty=3)
  axis(4, at=.Machine$double.eps, label=quote(epsilon[c]), las=1, col.axis=col)
}
R> eaxis(1); eaxis(2); draw.machEps(0.4)

```

In Figure 2 below, we zoom into the region where all methods have about the same (good) accuracy. The region is the rectangle defined by the ranges of `a.` and `ra2`: In Figure 2 below, we zoom into the region where all methods have about the same (good) accuracy. The region is the rectangle defined by the ranges of `a.` and `ra2`:

```

R> a. <- (1:400)/256
R> ra <- t(sapply(a., relE.11e))
R> ra2 <- ra[,-1]

```

In addition to zooming in Figure 1, we want to smooth the two curves, using a method assuming approximately normal errors. Notice however that neither the original, nor the log-transformed values have approximately symmetric errors, so we use `MASS::boxcox()` to determine the “correct” power transformation,

```

R> da <- cbind(a = a., as.data.frame(ra2))
R> library(MASS)
R> bc1 <- boxcox(abs(expm1) ~ a, data = da, lambda = seq(0,1, by=.01), plotit=.plot.BC)

```

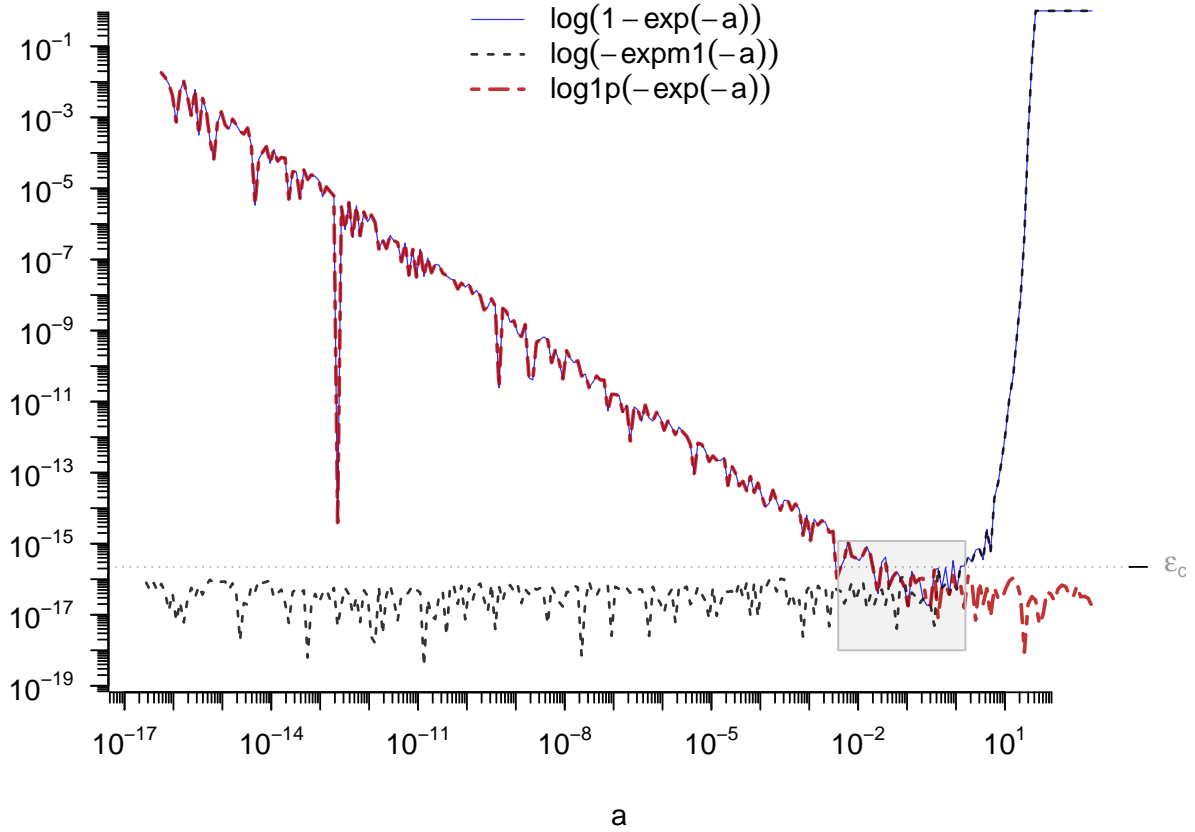


Figure 1: Absolute relative errors (with respect to 1024 bit **Rmpfr** computation) of the default, $\log(1 - e^{-a})$, and the two methods “**expm1**” $\log(-\text{expm1}(-a))$ and “**log1p**” $\log1p(-\exp(-a))$. Figure 2 will be a zoom into the gray rectangular region where all three curves are close.

```
R> bc2 <- boxcox(abs(log1p) ~ a, data = da, lambda = seq(0,1, by=.01), plotit=plot.BC)
R> c(with(bc1, x[which.max(y)]),
    with(bc2, x[which.max(y)]))## optimal powers
[1] 0.38 0.30
```

```
R> ## ==> taking ^ (1/3) :
R> s1 <- with(da, smooth.spline(a, abs(expm1)^(1/3), df = 9))
R> s2 <- with(da, smooth.spline(a, abs(log1p)^(1/3), df = 9))
```

and now plot a “zoom-in” of Figure 1. This already suggests that the cutoff, $a_0 = \log 2$ is empirically very close to optimal.

```
R> matplot(a., abs(ra2), type = "l", log = "y", # ylim = c(-1,1)*1e-12,
    col=cc[-1], lwd=ll[-1], lty=lt[-1],
    ylim = yl, xlab = "a", ylab = "", axes=FALSE)
R> legend("topright", leg[-1], col=cc[-1], lwd=ll[-1], lty=lt[-1], bty="n")
R> eaxis(1); eaxis(2); draw.machEps()
R> lines(a., predict(s1)$y ^ 3, col=cc[2], lwd=2)
R> lines(a., predict(s2)$y ^ 3, col=cc[3], lwd=2)
```

Why is it very plausible to take $a_0 := \log 2$ as approximately optimal cutoff? Already from Figure 2, empirically, an optimal cutoff a_0 is around 0.7. We propose to compute

$$f(a) = \log(1 - e^{-a}) = \log(1 - \exp(-a)), \quad a > 0, \quad (6)$$

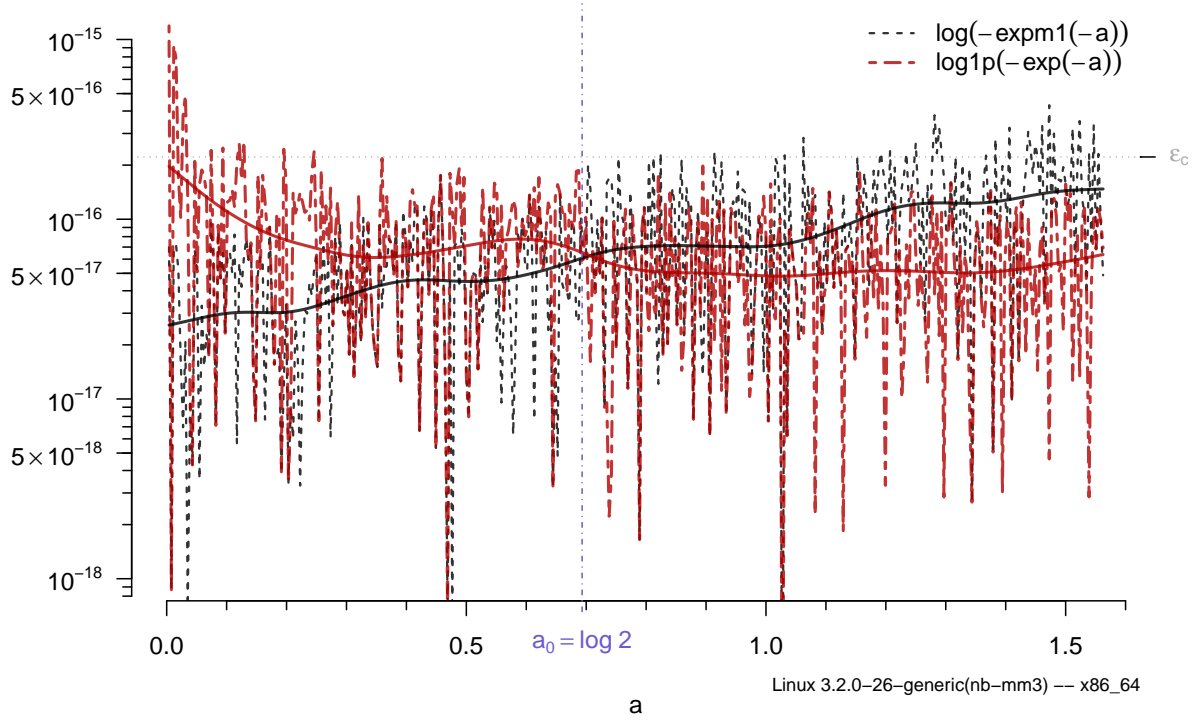


Figure 2: A “zoom in” of Figure 1 showing the region where the two basic methods, “`expm1`” and “`log1p`” switch their optimality with respect to their relative errors. Both have small relative errors in this region, typically below $\varepsilon_c := \text{Machine\$double.eps} = 2^{-52} \approx 2.22 \cdot 10^{-16}$. The smoothed curves indicate crossover close to $a = a_0 := \log 2$.

by a new method or function `log1mexp(a)`. It needs a cutoff a_0 between choosing `expm1` for $0 < a \leq a_0$ and `log1p` for $a > a_0$, i.e.,

$$f(a) = \text{log1mexp}(a) := \begin{cases} \log(-\text{expm1}(-a)) & 0 < a \leq a_0 \quad (:= \log 2 \approx 0.693) \\ \log1p(-\exp(-a)) & a > a_0. \end{cases} \quad (7)$$

The mathematical argument for choosing a_0 is quite simple, at least informally: In which situations does $1 - e^{-a}$ lose bits (binary digits) *entirely independently* of the computational algorithm? Well, as soon as it “spends” bits just to store its closeness to 1. And that is as soon as $e^{-a} < \frac{1}{2} = 2^{-1}$, because then, at least one bit cancels. This however is equivalent to $-a < \log(2^{-1}) = -\log(2)$ or $a > \log 2 =: a_0$.

3. Computation of $\log(1+\exp(x))$

Related to $\text{log1mexp}(a) = \log(1 - e^{-a})$ is the log survival function of the logistic distribution $\log(1 - F_L(x)) = \log \frac{1}{1+e^x} = -\log(1 + e^x)$,

$$g(x) := \log(1 + e^x) = \log1p(e^x), \quad (8)$$

(with a “+” instead of a “−”) which is easier to analyze and compute, its only problem being large x ’s where e^x overflows numerically. As $g(x) = \log(1 + e^x) = \log(e^x(e^{-x} + 1)) = x + \log(1 + e^{-x})$, we see from (1) that

$$g(x) = x + \log(1 + e^{-x}) = x + e^{-x} + \mathcal{O}((e^{-x})^2), \quad (9)$$

for $x \rightarrow \infty$. Using double precision arithmetic, a fast and accurate computational method is to use

$$g(x) = \log1pexp(x) := \begin{cases} \log1p(\exp(x)) & x < x_1 := 33.3, \\ x & x \geq x_1, \end{cases} \quad (10)$$

where x_1 can be replaced by a larger number such as 34 (and even 100, but not 800^3).

4. Conclusion

We have used high precision arithmetic (R package **Rmpfr**) to empirically verify that computing $f(a) = \log(1 - e^{-a})$ is accomplished best via equation (7). In passing, we have also shown that accurate computation of $g(x) = \log(1 + e^x)$ can be achieved via (10).

Session Information

```
R> toLatex(sessionInfo())
```

- R version 2.15.1 Patched (2012-07-06 r59741), x86_64-unknown-linux-gnu
- Locale: LC_CTYPE=de_CH.UTF-8, LC_NUMERIC=C, LC_TIME=en_US.UTF-8, LC_COLLATE=C, LC_MONETARY=en_US.UTF-8, LC_MESSAGES=C, LC_PAPER=C, LC_NAME=C, LC_ADDRESS=C, LC_TELEPHONE=C, LC_MEASUREMENT=de_CH.UTF-8, LC_IDENTIFICATION=C
- Base packages: base, datasets, grDevices, graphics, methods, stats, utils
- Other packages: MASS 7.3-19, Rmpfr 0.5-0, gmp 0.5-2, polynom 1.3-6, sfsmisc 1.0-20
- Loaded via a namespace (and not attached): tools 2.15.1

References

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³The R plot `curve(log1p(exp(x)) - x, 33.1, 33.5, n=2^10)` reveals a somewhat fuzzy cutoff x_1

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