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Arb is a C library for arbitrary-precision floating-point ball arithmetic, developed by Fredrik Johansson (fredrik.johansson@gmail.com). It supports real and complex numbers, polynomials, power series, matrices, and evaluation of many transcendental functions. All is done with automatic, rigorous error bounds.

The git repository is https://github.com/fredrik-johansson/arb/

The documentation website is http://fredrikj.net/arb/
2.1 Feature overview

Ball arithmetic, also known as mid-rad interval arithmetic, is an extension of floating-point arithmetic in which an error bound is attached to each variable. This allows doing rigorous computations over the real numbers, while avoiding the overhead of traditional (inf-sup) interval arithmetic at high precision, and eliminating much of the need for time-consuming and bug-prone manual error analysis associated with standard floating-point arithmetic. (See for example [Hoe2009].)

Other implementations of ball arithmetic include iRRAM and Mathemagix. In contrast to those systems, Arb is more focused on low-level arithmetic and computation of transcendental functions needed for number theory. Arb also differs in some technical aspects of the implementation.

Arb 2.x contains:

- A module (`arf`) for correctly rounded arbitrary-precision floating-point arithmetic. Arb’s floating-point numbers have a few special features, such as arbitrary-size exponents (useful for combinatorics and asymptotics) and dynamic allocation (facilitating implementation of hybrid integer/floating-point and mixed-precision algorithms).
- A module (`mag`) for representing magnitudes (error bounds) more efficiently than with an arbitrary-precision floating-point type.
- A module (`arb`) for real ball arithmetic, where a ball is implemented as an `arf` midpoint and a `mag` radius.
- A module (`acb`) for complex numbers in rectangular form, represented as pairs real balls.
- Functions for fast high-precision evaluation of various mathematical constants and special functions, implemented using ball arithmetic with rigorous error bounds.
- Modules (`arb_poly`, `acb_poly`) for polynomials or power series over the real and complex numbers, implemented using balls as coefficients, with asymptotically fast polynomial multiplication and many other operations.
- Modules (`arb_mat`, `acb_mat`) for matrices over the real and complex numbers, implemented using balls as coefficients. At the moment, only rudimentary linear algebra operations are provided.

Arb 1.x used a different set of numerical base types (`fmpr`, `fmprb` and `fmpcb`). These types had a slightly simpler internal representation, but generally had worse performance. Almost all methods for the Arb 1.x types have now been ported to faster equivalents for the Arb 2.x types. The last version to include both the Arb 1.x and Arb 2.x types and methods was Arb 2.2. As of Arb 2.3, only a small set of `fmpr` and `fmprb` methods are left for fallback and testing purposes.

Planned features include more transcendental functions and more extensive polynomial and matrix functionality, as well as further optimizations.

Arb uses GMP / MPIR and FLINT for the underlying integer arithmetic and other functions. The code conventions borrow from FLINT, and the project might get merged back into FLINT when the code stabilizes in the future. Arb also uses MPFR for testing purposes and for evaluation of some functions.
2.2 Setup

2.2.1 Download

Tarballs of released versions can be downloaded from https://github.com/fredrik-johansson/arb/releases

Alternatively, you can simply install Arb from a git checkout of https://github.com/fredrik-johansson/arb/. The master branch is generally safe to use (the test suite should pass at all times), and recommended for keeping up with the latest changes.

2.2.2 Dependencies

Arb has the following dependencies:

- Either MPIR (http://www.mpir.org) 2.6.0 or later, or GMP (http://www.gmplib.org) 5.1.0 or later. If MPIR is used instead of GMP, it must be compiled with the `--enable-gmpcompat` option.
- MPFR (http://www.mpfr.org) 3.0.0 or later.
- FLINT (http://www.flintlib.org) version 2.4 or later. You may also use a git checkout of https://github.com/fredrik-johansson/flint2

2.2.3 Installation as part of FLINT

With a sufficiently new version of FLINT, Arb can be compiled as a FLINT extension package.

Simply put the Arb source directory somewhere, say /path/to/arb. Then go to the FLINT source directory and build FLINT using:

```
./configure --extensions=/path/to/arb <other options>
make
make check (optional)
make install
```

This is convenient, as Arb does not need to be configured or linked separately. Arb becomes part of the compiled FLINT library, and the Arb header files will be installed along with the other FLINT header files.

2.2.4 Standalone installation

To compile, test and install Arb from source as a standalone library, first install FLINT. Then go to the Arb source directory and run:

```
./configure <options>
make
make check (optional)
make install
```

If GMP/MPIR, MPFR or FLINT is installed in some other location than the default path /usr/local, pass `--with-gmp=...`, `--with-mpfr=...` or `--with-flint=...` with the correct path to configure (type `./configure --help` to show more options).
### 2.2.5 Running code

Here is an example program to get started using Arb:

```c
#include "arb.h"

int main()
{
    arb_t x;
    arb_init(x);
    arb_const_pi(x, 50 * 3.33);
    arb_printn(x, 50, 0); flint_printf("\n");
    flint_printf("Computed with arb-%s\n", arb_version);
    arb_clear(x);
}
```

Compile it with:

```
gcc -larb test.c
```

or (if Arb is built as part of FLINT):

```
gcc -lflint test.c
```

If the Arb/FLINT header and library files are not in a standard location (/usr/local on most systems), you may also have to pass options such as:

```
-I/path/to/arb -I/path/to/flint -L/path/to/flint -L/path/to/arb
```

to gcc. Finally, to run the program, make sure that the linker can find the FLINT (and Arb) libraries. If they are installed in a nonstandard location, you can for example add this path to the LD_LIBRARY_PATH environment variable.

The output of the example program should be something like the following:

```
[3.1415926535897932384626433832795028841971693993751 +/- 6.28e-50]
Computed with arb-2.4.0
```

### 2.3 Potential issues

#### 2.3.1 Interface changes

Most of the core API should be stable at this point, and significant compatibility-breaking changes will be specified in the release notes.

In general, Arb does not distinguish between “private” and “public” parts of the API. The implementation is meant to be transparent by design. All methods are intended to be fully documented and tested (exceptions to this are mainly due to lack of time on part of the author). The user should use common sense to determine whether a function is concerned with implementation details, making it likely to change as the implementation changes in the future. The interface of `arb_add()` is probably not going to change in the next version, but `_arb_get_mpn_fixed_mod_pi4()` just might.

#### 2.3.2 Correctness

Except where otherwise specified, Arb is designed to produce provably correct error bounds. The code has been written carefully, and the library is extensively tested. However, like any complex mathematical software, Arb is
virtually certain to contain bugs, so the usual precautions are advised:

- Perform sanity checks on the output (check known mathematical relations; recompute to another precision and compare)
- Compare against other mathematical software
- Read the source code to verify that it does what it is supposed to do

All bug reports are highly welcome!

### 2.3.3 Aliasing

As a rule, Arb allows aliasing of operands. For example, in the function call `arb_add(z, x, y, prec)`, which performs \( z \leftarrow x + y \), any two (or all three) of the variables \( x, y \) and \( z \) are allowed to be the same. Exceptions to this rule are documented explicitly.

The general rule that input and output variables can be aliased with each other only applies to variables of the same type (ignoring `const` qualifiers on input variables – a special case is that `arb_srcptr` is considered the `const` version of `arb_ptr`). This is a natural extension of the so-called strict aliasing rule in C.

For example, in `arb_poly_evaluate()` which evaluates \( y = f(x) \) for a polynomial \( f \), the output variable \( y \) is not allowed to be a pointer to one of the coefficients of \( f \) (but aliasing between \( x \) and \( y \) or between \( x \) and the coefficients of \( f \) is allowed). This also applies to `_arb_poly_evaluate()`: for the purposes of aliasing, `arb_srcptr` (the type of the coefficient array within \( f \)) and `arb_t` (the type of \( x \)) are not considered to be the same type, and therefore must not be aliased with each other, even though an `arb_ptr/arb_srcptr` variable pointing to a length 1 array would otherwise be interchangeable with an `arb_t/const arb_t`.

Moreover, in functions that allow aliasing between an input array and an output array, the arrays must either be identical or completely disjoint, never partially overlapping.

There are natural exceptions to these aliasing restrictions, which may used internally without being documented explicitly. However, third party code should avoid relying on such exceptions.

An important caveat applies to aliasing of input variables. Identical pointers are understood to give permission for algebraic simplification. This assumption is made to improve performance. For example, the call `arb_mul(z, x, x, prec)` sets \( z \) to a ball enclosing the set

\[
\{ t^2 : t \in x \}
\]

and not the (generally larger) set

\[
\{ tu : t \in x, u \in x \}.
\]

If the user knows that two values \( x \) and \( y \) both lie in the interval \([-1, 1]\) and wants to compute an enclosure for \( f(x, y) \), then it would be a mistake to create an `arb_t` variable \( x \) enclosing \([-1, 1]\) and reusing the same variable for \( y \), calling \( f(x, x) \). Instead, the user has to create a distinct variable \( y \) also enclosing \([-1, 1]\).

Algebraic simplification is not guaranteed to occur. For example, `arb_add(z, x, x, prec)` and `arb_sub(z, x, x, prec)` currently do not implement this optimization. It is better to use `arb_mul_2exp_si(z, x, 1)` and `arb_zero(z)`, respectively.

### 2.3.4 Integer overflow

Machine-size integers are used for precisions, sizes of integers in bits, lengths of polynomials, and similar quantities that relate to sizes in memory. Very few checks are performed to verify that such quantities do not overflow. Precisions and lengths exceeding a small fraction of `LONG_MAX`, say \( 2^{24} \sim 10^7 \) on 32-bit systems, should be regarded as resulting in undefined behavior. On 64-bit systems this should generally not be an issue, since most calculations will
exhaust the available memory (or the user’s patience waiting for the computation to complete) long before running into integer overflows. However, the user needs to be wary of unintentionally passing input parameters of order \( \text{LONG\_MAX} \) or negative parameters where positive parameters are expected, for example due to a runaway loop that repeatedly increases the precision.

This caveat does not apply to exponents of floating-point numbers, which are represented as arbitrary-precision integers, nor to integers used as numerical scalars (e.g. \texttt{arb\_mul\_si()}). However, it still applies to conversions and operations where the result is requested exactly and sizes become an issue. For example, trying to convert the floating-point number \( 2^{2^{100}} \) to an integer could result in anything from a silent wrong value to thrashing followed by a crash, and it is the user’s responsibility not to attempt such a thing.

### 2.3.5 Thread safety and caches

Arb should be fully threadsafe, provided that both MPFR and FLINT have been built in threadsafe mode. Use \texttt{flint\_set\_num\_threads()} to set the number of threads that Arb is allowed to use internally for single computations (this is currently only exploited by a handful of operations). Please note that thread safety is only tested minimally, and extra caution when developing multithreaded code is therefore recommended.

Arb may cache some data (such as the value of \( \pi \) and Bernoulli numbers) to speed up various computations. In threadsafe mode, caches use thread-local storage. There is currently no way to save memory and avoid recomputation by having several threads share the same cache. Caches can be freed by calling the \texttt{flint\_cleanup()} function. To avoid memory leaks, the user should call \texttt{flint\_cleanup()} when exiting a thread. It is also recommended to call \texttt{flint\_cleanup()} when exiting the main program (this should result in a clean output when running Valgrind, and can help catching memory issues).

There does not seem to be an obvious way to make sure that \texttt{flint\_cleanup()} is called when exiting a thread using OpenMP. A possible solution to this problem is to use OpenMP sections, or to use C++ and create a thread-local object whose destructor invokes \texttt{flint\_cleanup()}.

### 2.3.6 Use of hardware floating-point arithmetic

Arb uses hardware floating-point arithmetic (the \texttt{double} type in C) in two different ways.

Firstly, \texttt{double} arithmetic as well as transcendental \texttt{libm} functions (such as \texttt{exp}, \texttt{log}) are used to select parameters heuristically in various algorithms. Such heuristic use of approximate arithmetic does not affect correctness: when any error bounds depend on the parameters, the error bounds are evaluated separately using rigorous methods. At worst, flaws in the floating-point arithmetic on a particular machine could cause an algorithm to become inefficient due to inefficient parameters being selected.

Secondly, \texttt{double} arithmetic is used internally for some rigorous error bound calculations. To guarantee correctness, we make the following assumptions. With the stated exceptions, these should hold on all commonly used platforms.

- A \texttt{double} uses the standard IEEE 754 format (with a 53-bit significand, 11-bit exponent, encoding of infinities and NaNs, etc.)
- We assume that the compiler does not perform “unsafe” floating-point optimizations, such as reordering of operations. Unsafe optimizations are disabled by default in most modern C compilers, including GCC and Clang. The exception appears to be the Intel C++ compiler, which does some unsafe optimizations by default. These must be disabled by the user.
- We do not assume that floating-point operations are correctly rounded (a counterexample is the x87 FPU), or that rounding is done in any particular direction (the rounding mode may have been changed by the user). We assume that any floating-point operation is done with at most 1.1 ulp error.
- We do not assume that underflow or overflow behaves in a particular way (we only use doubles that fit in the regular exponent range, or explicit infinities).
• We do not use transcendental \texttt{libm} functions, since these can have errors of several ulps, and there is unfortunately no way to get guaranteed bounds. However, we do use functions such as \texttt{ldexp} and \texttt{sqrt}, which we assume to be correctly implemented.

2.4 History and changes

For more details, view the commit log in the git repository \url{https://github.com/fredrik-johansson/arb}

• 2015-12-31 - version 2.8.1
  – Fixed 32-bit test failure for the Laguerre function.
  – Made the Laguerre function indeterminate at negative integer orders, to be consistent with the test code.

• 2015-12-29 - version 2.8.0
  – Compatibility and build system
    * Windows64 support (contributed by Bill Hart).
    * Fixed a bug that broke basic arithmetic on targets where FLINT uses fallback code instead of assembly code, such as PPC64 (contributed by Jeroen Demeyer).
    * Fixed configure to use EXTRA\_SHARED\_FLAGS/LDFLAGS, and other build system fixes (contributed by Tommy Hofmann, Bill Hart).
    * Added soname versioning (contributed by Julien Puydt).
    * Fixed test code on MinGW (contributed by Hrvoje Abraham).
    * Miscellaneous fixes to simplify interfacing Arb from Julia.
  – Arithmetic and elementary functions
    * Fixed \texttt{arf\_get\_d} to handle underflow/overflow correctly and to support round-to-nearest.
    * Added more complex inverse hyperbolic functions (\texttt{acb\_asin}, \texttt{acb\_acos}, \texttt{acb\_asinh}, \texttt{acb\_acosh}, \texttt{acb\_atanh}).
    * Added \texttt{arb\_contains\_int} and \texttt{acb\_contains\_int} for testing whether an interval contains any integer.
    * Added \texttt{acb\_quadratic\_roots\_fmpz}.
    * Improved \texttt{arb\_sinh} to use a more accurate formula for \( x < 0 \).
    * Added sinc function (\texttt{arb\_sinc}) (contributed by Alex Griffing).
    * Fixed bug in \texttt{arb\_exp} affecting convergence for huge input.
    * Faster implementation of \texttt{arb\_div\_2expm1\_ui}.
    * Added \texttt{mag\_root}, \texttt{mag\_geom\_series}.
    * Improved and added test code for \texttt{arb\_add\_error} functions.
    * Changed \texttt{arb\_pow} and \texttt{acb\_pow} to make \texttt{pow(0,positive) = 0} instead of nan.
    * Improved \texttt{acb\_sqrt} to return finite output for finite input straddling the branch cut.
    * Improved \texttt{arb\_set\_interval\_arf} so that \([ inf, inf ] = inf\) instead of an infinite interval.
    * Added computation of Bell numbers (\texttt{arb\_bell\_fmpz}).
    * Added \texttt{arb\_power\_sum\_vec} for computing power sums using Bernoulli numbers.
    * Added computation of the Fujiwara root bound for \texttt{acb\_poly}.
- Added code to identify all the real roots of a real polynomial (acb_poly_validate_real_roots).
- Added several convenient assignment functions, including arb_set_d, acb_set_d, acb_set_d_d, acb_set_fmpz_fmpz (contributed by Ricky Farr).
- Added many accessor functions (_arb/acb_vec_entry_ptr, arb_get_mid/rad_arb, acb_real/imag_ptr, arb_mid/rad_ptr, acb_get_real/imag).
- Added missing functions acb_add_si, acb_sub_si.
- Renamed arb_root to arb_root_ui (keeping alias) and added acb_root_ui.

- Special functions
  - Implemented the Gauss hypergeometric function 2F1 and its regularized version.
  - Fixed two bugs in acb_hypgeom_pfq_series_direct discovered while implementing 2F1. In rare cases, these could lead to incorrect values for functions depending on parameter derivatives of hypergeometric series.
    - The first bug involved incorrect handling of negative integer parameters. The bug only affected 2F1 and higher functions; it did not affect correctness of any previously implemented functions that relied on acb_hypgeom_pfq_series_direct (such as Bessel Y and K functions of integer order).
    - The second bug involved a too small bound being computed for the sum of a geometric series. The geometric series bound is nearly tight for 2F1, and the incorrect version caused immediate test failures for that function. Theoretically, this bug affected correctness of some previously-implemented functions that relied on acb_hypgeom_pfq_series_direct (such as Bessel Y and K functions of integer order), but since the geometric bound is not as tight in those cases, those functions were still reliable in practice (no failing test case has been found).
  - Implemented Airy functions and their derivatives (acb_hypgeom_airy).
  - Implemented the confluent hypergeometric function 0F1 (acb_hypgeom_0f1).
  - Implemented associated Legendre functions P and Q.
  - Implemented Chebyshev, Jacobi, Gegenbauer, Laguerre, Hermite functions.
  - Implemented spherical harmonics.
  - Added function for computing Bessel J and Y functions simultaneously.
  - Added rising factorials for non-integer n (arb_rising, acb_rising).
  - Made rising factorials use gamma function for large integer n.
  - Faster algorithm for theta constants and Dedekind eta function at very high precision.
  - Fixed erf to give finite values instead of +/-inf for big imaginary input.
  - Improved acb_zeta (and arb_zeta) to automatically use fast code for integer zeta values.
  - Added double factorial (arb_doublefac_ui).
  - Added code for generating Hilbert class polynomials (acb_modular_hilbert_class_poly).

- Matrices
  - Added faster matrix squaring (arb/acb_mat_sqr) (contributed by Alex Griffing).
  - Added matrix trace (arb/acb_mat_trace) (contributed by Alex Griffing).
  - Added arb/acb_mat_set_round_fmpz_mat, acb_mat_set(_round)_arb_mat (contributed by Tommy Hofmann).
* Added arb/acb_mat_transpose (contributed by Tommy Hofmann).
* Added comparison methods arb/acb_mat_eq/ne (contributed by Tommy Hofmann).

– Other
  * Added complex_plot example program.
  * Added Airy functions to real_roots example program.
  * Other minor patches were contributed by Alexander Kobel, Marc Mezzarobba, Julien Puydt.
  * Removed obsolete file config.h.

• 2015-07-14 - version 2.7.0
  – hypergeometric functions
    * implemented Bessel I and Y functions (acb_hypgeom_bessel_i, acb_hypgeom_bessel_y)
    * fixed bug in Bessel K function giving the wrong branch for negative real arguments
    * added code for evaluating complex hypergeometric series binary splitting
    * added code for evaluating complex hypergeometric series using fast multipoint evaluation
  – gamma related functions
    * implemented the Barnes G-function and its continuous logarithm (acb_barnes_g, acb_log_barnes_g)
    * implemented the generalized polygamma function (acb_polygamma)
    * implemented the reflection formula for the logarithmic gamma function (acb_lgamma, acb_poly_lgamma_series)
    * implemented the digamma function of power series (arb_poly_digamma_series, acb_poly_digamma_series)
    * improved acb_poly_zeta_series to produce exact zero imaginary parts in most cases when the result should be real-valued
    * made the real logarithmic gamma function (arb_lgamma, arb_poly_lgamma_series) abort more quickly for negative input
  – elementary functions
    * added arb_exp_expinv and acb_exp_expinv functions for simultaneously computing exp(x), exp(-x)
    * improved acb_tan, acb_tan_pi, acb_cot and acb_cot_pi for input with large imaginary parts
    * added complex hyperbolic functions (acb_sinh, acb_cosh, acb_sinh_cosh, acb_tanh, acb_coth)
    * added acb_log_sin_pi for computing the logarithmic sine function without branch cuts away from the real line
    * added arb_poly_cot_pi_series, acb_poly_cot_pi_series
    * added arf_root and improved speed of arb_root
    * tuned algorithm selection in arb_pow_fmpq
  – other
    * added documentation for arb and acb vector functions

• 2015-04-19 - version 2.6.0
  – special functions
    * added the Bessel K function
* added the confluent hypergeometric functions $M$ and $U$
* added exponential, trigonometric and logarithmic integrals $e_i$, $s_i$, $c_i$, $c_i$, $l_i$
* added the complete elliptic integral of the second kind $E$
* added support for computing hypergeometric functions with power series as parameters
* fixed special cases in Bessel $J$ function returning useless output
* fix precision of zeta function accidentally being capped at 7000 digits (bug in 2.5)
* special-cased real input in the gamma functions for complex types
* fixed exp of huge numbers outputting unnecessarily useless intervals
* fixed broken code in erf that sometimes gave useless output
* made selection of number of terms in hypergeometric series more robust

– polynomials and power series
  * added $\sin_{\pi}$, $\cos_{\pi}$ and $\sin_{\cos_{\pi}}$ for real and complex power series
  * speeded up series reciprocal and division for length = 2
  * added add$_{\text{si}}$ methods for polynomials
  * made inv$_{\text{series}}$ and div$_{\text{series}}$ with zero input produce indeterminates instead of aborting
  * added arb$_{\text{poly}}$, arb$_{\text{poly}}$, arb$_{\text{poly}}$

– basic functions
  * added comparison methods arb$_{\text{eq}}$, arb$_{\text{ne}}$, arb$_{\text{lt}}$, arb$_{\text{le}}$, arb$_{\text{gt}}$, arb$_{\text{ge}}$, acb$_{\text{eq}}$, acb$_{\text{ne}}$
  * added acb$_{\text{rel}}$accuracy$_{\text{bits}}$ and improved the real version
  * fixed precision of constants like $\pi$ behaving more nondeterministically than necessary
  * fixed arf$_{\text{get}}$, mag$_{\text{lower}}$(nan) to output 0 instead of inf

– other
  * removed call to fmpq_dedekind$\text{sum}$ which only exists in the git version of flint
  * fixed a test code bug that could cause crashes on some systems
  * added fix for static build on OS X (thanks Marcello Seri)
  * miscellaneous corrections to the documentation

• 2015-01-28 - version 2.5.0
  – string conversion
    * added arb$_{\text{set}}$, arb$_{\text{str}}$
    * added arb$_{\text{get}}$, arb$_{\text{print}}$n for pretty-printed rigorous decimal output
    * added helper functions for binary to decimal conversion
  – core arithmetic
    * improved speed of division when using GMP instead of MPIR
    * improved complex division with a small denominator
    * removed a little bit of overhead for complex squaring
  – special functions
• faster code for atan at very high precision, used instead of mpfr_atan
• optimized elementary functions slightly for small input
• added modified error functions erfc and erfi
• added the generalized exponential integral
• added the upper incomplete gamma function
• implemented the complete elliptic integral of the first kind
• implemented the arithmetic-geometric mean of complex numbers
• optimized arb_digamma for small integers
• made mag_log_ui, mag_binpow_uiui and mag_polylog_tail proper functions
• added pow, agm, erf, elliptic_k, elliptic_p as functions of complex power series
• added incomplete gamma function of complex power series
• improved code for bounding complex rising factorials (the old code could potentially have given wrong results in degenerate cases)
• added arb_sqrt1pm1, arb_atanh, arb_asinh, arb_atanh
• added arb_log1p, acb_log1p, acb_atan
• added arb_hurwitz_zeta
• improved parameter selection in the Hurwitz zeta function to try to avoid stalling when given enormous input
• optimized sqrt and rsqrt of power series when given a binomial as input
• made arb_bermoulli_ui(2^64-2) not crash
• fixed rgamma of negative integers returning indeterminate

– polynomials and matrices
  • added characteristic polynomial computation for real and complex matrices
  • added polynomial set_round methods
  • added is_real methods for more types
  • added more get_unique_fmpz methods
  • added code for generating Swinnerton-Dyer polynomials
  • improved error bounding in det() and exp() of complex matrices to recognize when the result is real-valued
  • changed polynomial divrem to return success/fail instead of aborting on divide by zero

– miscellaneous
  • added logo to documentation
  • made inlined functions build as part of the library
  • silenced a clang warning
  • made _acb_vec_sort_pretty a library function

• 2014-11-15 - version 2.4.0
  • arithmetic and core functions
• made evaluation of sin, cos and exp at medium precision faster using the sqrt trick
• optimized arb_sinh and arb_sinh_cosh
• optimized complex division with a small denominator
• optimized cubing of complex numbers
• added floor and ceil functions for the arf and arb types
• added acb_poly powering functions
• added acb_exp_pi_i
• added functions for evaluation of Chebyshev polynomials
• fixed arb_div to output nan for input containing nan
  – added a module acb_hypgeom for hypergeometric functions
    • evaluation of the generalized hypergeometric function in convergent cases
    • evaluation of confluent hypergeometric functions using asymptotic expansions
    • the Bessel function of the first kind for complex input
    • the error function for complex input
  – added a module acb_modular for modular forms and elliptic functions
    • support for working with modular transformations
    • mapping a point to the fundamental domain
    • evaluation of Jacobi theta functions and their series expansions
    • the Dedekind eta function
    • the j-invariant and the modular lambda and delta function
    • Eisenstein series
    • the Weierstrass elliptic function and its series expansion
  – miscellaneous
    • fixed mag_print printing a too large exponent
    • fixed printd methods to use a fallback instead of aborting when printing numbers too large for MPFR
    • added version number string (arb_version)
    • various additions to the documentation

• 2014-09-25 - version 2.3.0
  – removed most of the legacy (Arb 1.x) modules
  – updated build scripts, hopefully fixing various issues
  – new implementations of arb_sin, arb_cos, arb_sin_cos, arb_atan, arb_log, arb_exp, arb_expm1, much faster up to a few thousand bits
  – ported the bit-burst code for high-precision exponentials to the arb type
  – speeded up arb_log_ui_from_prev
  – added mag_exp, mag_expm1, mag_exp_tail, mag_pow_fmpz
  – improved various mag functions
– added arb_get/set_interval_mpfr, arb_get_interval_arf, and improved arb_set_interval_arf
– improved arf_get_fmpz
– prettier printing of complex numbers with negative imaginary part
– changed some frequently-used functions from inline to non-inline to reduce code size

• 2014-08-01 - version 2.2.0
– added functions for computing polylogarithms and order expansions of polylogarithms, with support for real and complex s, z
– added a missing cast affecting C++ compatibility
– generalized powsum functions to allow a geometric factor
– improved powsum functions slightly when the exponent is an integer
– faster arb_log_ui_from_prev
– added mag_sqrt and mag_rsqrt functions
– fixed various minor bugs and added missing tests and documentation entries

• 2014-06-20 - version 2.1.0
– ported most of the remaining functions to the new arb/acb types, including:
  * elementary functions (log, atan, etc.)
  * hypergeometric series summation
  * the gamma function
  * the Riemann zeta function and related functions
  * Bernoulli numbers
  * the partition function
  * the calculus modules (rigorous real root isolation, rigorous numerical integration of complex-valued functions)
  * example programs
– added several missing utility functions to the arf and mag modules

• 2014-05-27 - version 2.0.0
– new modules mag, arf, arb, arb_poly, arb_mat, acb, acb_poly, acb_mat for higher-performance ball arithmetic
– poly_roots2 and hilbert_matrix2 example programs
– vector dot product and norm functions (contributed by Abhinav Baid)

• 2014-05-03 - version 1.1.0
– faster and more accurate error bounds for polynomial multiplication (error bounds are now always as good as with classical multiplication, and multiplying high-degree polynomials with approximately equal coefficients now has proper quasilinear complexity)
– faster and much less memory-hungry exponentials at very high precision
– improved the partition function to support n bigger than a single word, and enabled the possibility to use two threads for the computation
– fixed a bug in floating-point arithmetic that caused a too small bound for the rounding error to be reported when the result of an inexact operation was rounded up to a power of two (this bug did not affect the correctness of ball arithmetic, because operations on ball midpoints always round down)
– minor optimizations to floating-point arithmetic
– improved argument reduction of the digamma function and short series expansions of the rising factorial
– removed the holonomic module for now, as it did not really do anything very useful

• 2013-12-21 - version 1.0.0
– new example programs directory
  * poly_roots example program
  * real_roots example program
  * pi_digits example program
  * hilbert_matrix example program
  * keiper_li example program
– new fmpcrb_calc module for calculus with real functions
  * bisection-based root isolation
  * asymptotically fast Newton root refinement
– new fmpcb_calc module for calculus with complex functions
  * numerical integration using Taylor series
– scalar functions
  * simplified fmpcrb_const_euler using published error bound
  * added fmpcrb_inv
  * fmpcrb_trim, fmpcb_trim
  * added fmpcb_rsqrt (complex reciprocal square root)
  * fixed bug in fmpcrb_sqrtpos with nonfinite input
  * slightly improved fmpcrb powering code
  * added various functions for bounding fmprbs by powers of two
  * added fmpr_is_int
– polynomials and power series
  * implemented scaling to speed up blockwise multiplication
  * slightly faster basecase power series exponentials
  * improved sin/cos/tan/exp for short power series
  * added complex sqrt_series, rsqrt_series
  * implemented the Riemann-Siegel Z and theta functions for real power series
  * added fmpcrb_poly_pow_series, fmpcrb_poly_pow_ui and related methods
  * fmpcrb/fmpcb_poly_contains_fmpz_poly
  * faster composition by monomials
  * implemented Borel transform and binomial transform for real power series
– matrices
  * implemented matrix exponentials
  * multithreaded fmpcb_mat_mul
  * added matrix infinity norm functions
  * added some more matrix-scalar functions
  * added matrix contains and overlaps methods
– zeta function evaluation
  * multithreaded power sum evaluation
  * faster parameter selection when computing many derivatives
  * implemented binary splitting to speed up computing many derivatives
– miscellaneous
  * corrections for C++ compatibility (contributed by Jonathan Bober)
  * several minor bugfixes and test code enhancements
  * 2013-08-07 - version 0.7
  * floating-point and ball functions
    * documented, added test code, and fixed bugs for various operations involving a ball containing an infinity or NaN
    * added reciprocal square root functions (fmpc_rsqrt, fmpcb_rsqrt) based on mpfr_rec_sqrt
    * faster high-precision division by not computing an explicit remainder
    * slightly faster computation of pi by using new reciprocal square root and division code
    * added an mpfr function for approximate division to speed up certain radius operations
    * added fmpc_set_d for conversion from double
    * allow use of doubles to optionally compute the partition function faster but without an error bound
    * bypass mpfr overflow when computing the exponential function to extremely high precision (approximately 1 billion digits)
    * made fmpc_exp faster for large numbers at extremely high precision by skipping the log(2) removal
    * made fmpcb_lgamma faster at high precision by speeding up the argument reduction branch computation
    * added fmpc_asin, fmpc_acos
    * added various other utility functions to the fmpc module
    * added a function for computing the Glaisher constant
    * optimized evaluation of the Riemann zeta function at high precision
– polynomials and power series
  * made squaring of polynomials faster than generic multiplication
  * implemented power series reversion (various algorithms) for the fmpc_poly type
  * added many fmpc_poly utility functions (shifting, truncating, setting/getting coefficients, etc.)
  * improved power series division when either operand is short
* improved power series logarithm when the input is short
* improved power series exponential to use the basecase algorithm for short input regardless of the output size
* added power series square root and reciprocal square root
* added atan, tan, sin, cos, sin_cos, asin, acos fmprb_poly power series functions
* added Newton iteration macros to simplify various functions
* added gamma functions of real and complex power series ([fm-prb/fmpcb]_poly_[gamma/rgamma/lgamma]_series)
* added wrappers for computing the Hurwitz zeta function of a power series ([fm-prb/fmpcb]_poly_zeta_series)
* implemented sieving and other optimizations to improve performance for evaluating the zeta function of a short power series
* improved power series composition when the inner series is linear
* added many fmpcb_poly versions of nearly all fmprb_poly functions
* improved speed and stability of series composition/reversion by balancing the power table exponents

  – other
  * added support for freeing all cached data by calling flint_cleanup()
  * introduced fmprb_ptr, fmprb_srcptr, fmpcb_ptr, fmpcb_srcptr typedefs for cleaner function signatures
  * various bug fixes and general cleanup

• 2013-05-31 - version 0.6
  – made fast polynomial multiplication over the reals numerically stable by using a blockwise algorithm
  – disabled default use of the Gauss formula for multiplication of complex polynomials, to improve numerical stability
  – added division and remainder for complex polynomials
  – added fast multipoint evaluation and interpolation for complex polynomials
  – added missing fmprb_poly_sub and fmpcb_poly_sub functions
  – faster exponentials (fmprb_exp and dependent functions) at low precision, using precomputation
  – rewrote fnpr_add and fnpr_sub using mpn level code, improving efficiency at low precision
  – ported the partition function implementation from flint (using ball arithmetic in all steps of the calculation to guarantee correctness)
  – ported algorithm for computing the cosine minimal polynomial from flint (using ball arithmetic to guarantee correctness)
  – support using gmp instead of mpir
  – only use thread-local storage when enabled in flint
  – slightly faster error bounding for the zeta function
  – added some other helper functions

• 2013-03-28 - version 0.5
  – arithmetic and elementary functions
* added \texttt{fmpr\_get\_fmpz}, \texttt{fmpr\_get\_si}
* fixed accuracy problem with \texttt{fmprb\_div\_2expm1}
* special-cased squaring of complex numbers
* added various \texttt{fmpcb} convenience functions (\texttt{addmul\_ui}, etc)
* optimized \texttt{fmpr\_cmp\_2exp\_si} and \texttt{fmpr\_cmpabs\_2exp\_si}, and added test code for comparison functions
* added \texttt{fmprb\_atan2}, also fixing a bug in \texttt{fmpcb\_arg}
* added \texttt{fmprb\_sin\_pi}, \texttt{cos\_pi}, \texttt{sin\_cos\_pi} etc.
* added \texttt{fmprb\_sin\_pi\_fmpq} (etc.) using algebraic methods for fast evaluation of roots of unity
* faster \texttt{fmprb\_poly\_evaluate} and \texttt{evaluate\_fmpcb} using rectangular splitting
* added \texttt{fmprb\_poly\_evaluate2}, \texttt{evaluate2\_fmpcb} for simultaneously evaluating the derivative
* added \texttt{fmprb\_poly\_root} polishing code using near-optimal Newton steps (experimental)
* added \texttt{fmpr\_root}, \texttt{fmprb\_root} (currently based on MPFR)
* added \texttt{fmpr\_min}, \texttt{fmpr\_max}
* added \texttt{fmprb\_set\_interval\_fmpr}, \texttt{fmprb\_union}
* added \texttt{fmpr\_bits}, \texttt{fmprb\_bits}, \texttt{fmpcb\_bits} for obtaining the mantissa width
* added \texttt{fmprb\_hypot}
* added complex square roots
* improved \texttt{fmprb\_log} to slightly improve speed, and properly support huge arguments
* fixed \texttt{exp}, \texttt{cosh}, \texttt{sinh} to work with huge arguments
* added \texttt{fmprb\_expm1}
* fixed \texttt{sin}, \texttt{cos}, \texttt{atan} to work with huge arguments
* improved \texttt{fmprb\_pow} and \texttt{fmpcb\_pow}, including automatic detection of small integer and half-integer exponents
* added many more elementary functions: \texttt{fmprb\_tan/cot/tanh/coth}, \texttt{fmpcb\_tan/cot}, and pi versions
* added \texttt{fmprb\_const\_e}, \texttt{const\_log2}, \texttt{const\_log10}, \texttt{const\_catalan}
* fixed ball containment/overlap checking to work operate efficiently and correctly with huge exponents
* strengthened test code for many core operations

- special functions
  * reorganized zeta function related code
  * faster evaluation of the Riemann zeta function via sieving
  * documented and improved efficiency of the zeta constant binary splitting code
  * calculate error bound in Borwein’s algorithm with \texttt{fmprs} instead of using doubles
  * optimized divisions in zeta evaluation via the Euler product
  * use functional equation for Riemann zeta function of a negative argument
  * compute single Bernoulli numbers using ball arithmetic instead of relying on the floating-point code in \texttt{flint}
• initial code for evaluating the gamma function using its Taylor series
• much faster rising factorials at high precision, using difference polynomials
• much faster gamma function at high precision
• added complex gamma function, log gamma function, and other versions
• added fmprb_agm (real arithmetic-geometric mean)
• added fmprb_gamma_fmpq, supporting rapid computation of gamma(p/q) for q = 1, 2, 3, 4, 6
• added real and complex digamma function
• fixed unnecessary recomputation of Bernoulli numbers
• optimized computation of Euler’s constant, and added proper error bounds
• avoid reliance on doubles in the hypergeometric series tail bound
• cleaned up factorials and binomials, computing factorials via gamma
  – other
  • added an fmpz_extras module to collect various internal fmpz helper functions
  • fixed detection of flint header files
  • fixed various other small bugs

• 2013-01-26 - version 0.4
  – much faster fnmpr_mul, fnmprb_mul and set_round, resulting in general speed improvements
  – code for computing the complex Hurwitz zeta function with derivatives
  – fixed and documented error bounds for hypergeometric series
  – better algorithm for series evaluation of the gamma function at a rational point
  – much faster generation of Bernoulli numbers
  – complex log, exp, pow, trigonometric functions (currently based on MPFR)
  – complex nth roots via Newton iteration
  – added code for arithmetic on fmpcb_polys
  – code for computing Khinchin’s constant
  – code for rising factorials of polynomials or power series
  – faster sin_cos
  – better div_2expm1
  – many other new helper functions
  – improved thread safety
  – more test code for core operations

• 2012-11-07 - version 0.3
  – converted documentation to sphinx
  – new module fmpcb for ball interval arithmetic over the complex numbers
    • conversions, utility functions and arithmetic operations
  – new module fmpcb_mat for matrices over the complex numbers
• conversions, utility functions and arithmetic operations
  • multiplication, LU decomposition, solving, inverse and determinant
  – new module fmpcb_poly for polynomials over the complex numbers
    • root isolation for complex polynomials
  – new module fmpz_holonomic for functions/sequences defined by linear differential/difference equations with polynomial coefficients
    • functions for creating various special sequences and functions
    • some closure properties for sequences
    • Taylor series expansion for differential equations
    • computing the nth entry of a sequence using binary splitting
    • computing the nth entry mod p using fast multipoint evaluation
  – generic binary splitting code with automatic error bounding is now used for evaluating hypergeometric series
  – matrix powering
  – various other helper functions

• 2012-09-29 - version 0.2
  – code for computing the gamma function (Karatsuba, Stirling’s series)
  – rising factorials
  – fast exp_series using Newton iteration
  – improved multiplication of small polynomials by using classical multiplication
  – implemented error propagation for square roots
  – polynomial division (Newton-based)
  – polynomial evaluation (Horner) and composition (divide-and-conquer)
  – product trees, fast multipoint evaluation and interpolation (various algorithms)
  – power series composition (Horner, Brent-Kung)
  – added the fmprb_mat module for matrices of balls of real numbers
  – matrix multiplication
  – interval-aware LU decomposition, solving, inverse and determinant
  – many helper functions and small bugfixes

• 2012-09-14 - version 0.1

• 2012-08-05 - began simplified rewrite

• 2012-04-05 - experimental ball and polynomial code

### 2.5 Example programs

The examples directory (https://github.com/fredrik-johansson/arb/tree/master/examples) contains several complete C programs, which are documented below. Running:
make examples

will compile the programs and place the binaries in build/examples.

2.5.1 pi.c

This program computes π to an accuracy of roughly \( n \) decimal digits by calling the `arb_const_pi()` function with a working precision of roughly \( n \log_2(10) \) bits.

Sample output, computing π to one million digits:

```
> build/examples/pi 1000000
computing pi with a precision of 3321933 bits... cpu/wall(s): 0.58 0.586
virt/peak/res/peak(MB): 28.24 36.84 8.86 15.56
[3.14159265358979323846{...999959 digits...}42209010610577945815 +/- 3e-1000000]
```

The program prints an interval guaranteed to contain \( \pi \), and where all displayed digits are correct up to an error of plus or minus one unit in the last place (see `arb_printn()`). By default, only the first and last few digits are printed. Pass 0 as a second argument to print all digits (or pass \( m \) to print \( m + 1 \) leading and \( m \) trailing digits, as above with the default \( m = 20 \)).

2.5.2 hilbert_matrix.c

Given an input integer \( n \), this program accurately computes the determinant of the \( n \times n \) Hilbert matrix. Hilbert matrices are notoriously ill-conditioned: although the entries are close to unit magnitude, the determinant \( h_n \) decreases superexponentially (nearly as \( 1/4^{n^2} \)) as a function of \( n \). This program automatically doubles the working precision until the ball computed for \( h_n \) by `arb_mat_det()` does not contain zero.

Sample output:

```
> build/examples/hilbert_matrix 200
prec=20: 0 +/- 5.5777e-330
prec=40: 0 +/- 2.5785e-542
prec=80: 0 +/- 8.1169e-926
prec=160: 0 +/- 2.8538e-1924
prec=320: 0 +/- 6.3868e-4129
prec=640: 0 +/- 1.8545e-8826
prec=1280: 0 +/- 1.8545e-17758
prec=2560: 2.955454297e-23924 +/- 6.4586e-24044
success!
cpu/wall(s): 9.06 9.095
virt/peak/res/peak(MB): 55.52 55.52 35.50 35.50
```

2.5.3 keiper_li.c

Given an input integer \( n \), this program rigorously computes numerical values of the Keiper-Li coefficients \( \lambda_0, \ldots, \lambda_n \). The Keiper-Li coefficients have the property that \( \lambda_n > 0 \) for all \( n > 0 \) if and only if the Riemann hypothesis is true. This program was used for the record computations described in [Joh2013] (the paper describes the algorithm in some more detail).

The program takes the following parameters:

```
keiper_li n [-prec prec] [-threads num_threads] [-out out_file]
```

2.5. Example programs
The program prints the first and last few coefficients. It can optionally write all the computed data to a file. The working precision defaults to a value that should give all the coefficients to a few digits of accuracy, but can optionally be set higher (or lower). On a multicore system, using several threads results in faster execution.

Sample output:

```
> build/examples/keiper_1i 1000 -threads 2
zeta: cpu/wall(s): 0.4 0.244
ev/peak/res/peak(MB): 167.98 294.69 5.09 7.43
log: cpu/wall(s): 0.03 0.038
gamma: cpu/wall(s): 0.02 0.016
binomial transform: cpu/wall(s): 0.01 0.018
0: -0.69314718055994530941723212145817656807550013436026 +/- 6.5389e-347
1: 0.02395708966121033814310247906495291621932127152051 +/- 2.0924e-345
2: 0.04617286714023335192864234096033943387066108314123 +/- 1.674e-344
3: 0.069212973181082679304973488726010689942120263932 +/- 5.0219e-344
4: 0.092197619873060409647627872409439018065541673490213 +/- 2.0089e-343
5: 0.1151085429223549048662128109857276671349132303596 +/- 1.0044e-342
6: 0.1379266871372988290416713700341666356138966078654 +/- 6.0264e-342
7: 0.16063175965299421294040287257385366292282442046163 +/- 2.1092e-341
8: 0.183219459643382579081939331774721859849989089273432 +/- 8.4368e-341
9: 0.20565733870917046170289384721343304741236553410044 +/- 7.5931e-340
10: 0.2279336319315774369303405736844533074835942738 +/- 7.5931e-339
991: 0.31966179616133679283738996569486825626210430813341 +/- 2.461e-11
992: 0.3203766239254884035349986518332552033162909717288 +/- 9.5363e-11
993: 0.32109206129713238281165911633262682034375592414 +/- 1.8495e-10
994: 0.321807340188462110258826121503870112747188888893 +/- 3.5907e-10
995: 0.3225217392815185726928702951225314023773358152533 +/- 6.978e-10
996: 0.3232344858146283733322360941370912358283071281 +/- 1.3574e-09
997: 0.32394711488601452289542667580382034526509232475 +/- 2.6433e-09
998: 0.32465175910327008384143420432560514885689322209 +/- 5.1524e-09
999: 0.32535482758613821198125760520605269854493162101 +/- 1.0053e-08
1000: 0.326053161686466457406504694083223815804982041872 +/- 3.927e-08
virt/peak/res/peak(MB): 170.18 294.69 7.51 7.51
```

### 2.5.4 real_roots.c

This program isolates the roots of a function on the interval \((a, b)\) (where \(a\) and \(b\) are input as double-precision literals) using the routines in the `arb_calc` module. The program takes the following arguments:

```
real_roots function a b [-refine d] [-verbose] [-maxdepth n] [-maxeval n] [-maxfound n] [-prec n]
```

The following functions (specified by an integer code) are implemented:

- **0** - \(Z(x)\) (Riemann-Siegel \(Z\)-function)
- **1** - \(\sin(x)\)
- **2** - \(\sin(x^2)\)
- **3** - \(\sin(1/x)\)
- **4** - \(Ai(x)\) (Airy function)
- **5** - \(Ai'(x)\) (Airy function)
- **6** - \(Bi(x)\) (Airy function)
- **7** - \(Bi'(x)\) (Airy function)

The following options are available:
• `-refine d`: If provided, after isolating the roots, attempt to refine the roots to \( d \) digits of accuracy using a few bisection steps followed by Newton’s method with adaptive precision, and then print them.

• `-verbose`: Print more information.

• `-maxdepth n`: Stop searching after \( n \) recursive subdivisions.

• `-maxeval n`: Stop searching after approximately \( n \) function evaluations (the actual number evaluations will be a small multiple of this).

• `-maxfound n`: Stop searching after having found \( n \) isolated roots.

• `-prec n`: Working precision to use for the root isolation.

With `function 0`, the program isolates roots of the Riemann zeta function on the critical line, and guarantees that no roots are missed (there are more efficient ways to do this, but it is a nice example):

```
> build/examples/real_roots 0 0.0 50.0 -verbose
interval: [0, 50]
maxdepth = 30, maxeval = 100000, maxfound = 100000, low_prec = 30
found isolated root in: [14.111328125, 14.16015625]
found isolated root in: [20.99609375, 21.044921875]
found isolated root in: [25, 25.048828125]
found isolated root in: [30.419921875, 30.443359375]
found isolated root in: [32.91015625, 32.958984375]
found isolated root in: [37.548828125, 37.59765625]
found isolated root in: [40.91796875, 40.966796875]
found isolated root in: [43.310546875, 43.3349609375]
found isolated root in: [47.998046875, 48.0224609375]
found isolated root in: [49.755859375, 49.7802734375]
---------------------------------------------------------------
Found roots: 10
Subintervals possibly containing undetected roots: 0
Function evaluations: 3058
cpu/wall(s): 0.202 0.202
```

Find just one root and refine it to approximately 75 digits:

```
> build/examples/real_roots 0 0.0 50.0 -maxfound 1 -refine 75
interval: [0, 50]
maxdepth = 30, maxeval = 100000, maxfound = 1, low_prec = 30
refined root (0/8):
[14.134725141734693790457251983562470270784257115669243175685567460149963429809 +/- 2.57e-76]
---------------------------------------------------------------
Found roots: 1
Subintervals possibly containing undetected roots: 7
Function evaluations: 761
cpu/wall(s): 0.055 0.056
virt/peak/res/peak(MB): 26.12 26.14 2.75 2.75
```

Find the first few roots of an Airy function and refine them to 50 digits each:

```
> build/examples/real_roots 4 -10 0 -refine 50
interval: [-10, 0]
maxdepth = 30, maxeval = 100000, maxfound = 100000, low_prec = 30
refined root (0/6):
[-9.02265085334098038015819083988089256524677535156083 +/- 4.85e-52]
refined root (1/6):
```

2.5. Example programs
Find roots of \( \sin(x^2) \) on \((0, 100)\). The algorithm cannot isolate the root at \(x = 0\) (it is at the endpoint of the interval, and in any case a root of multiplicity higher than one). The failure is reported:

```
> build/examples/real_roots 2 0 100
interval: [0, 100]
maxdepth = 30, maxeval = 100000, maxfound = 100000, low_prec = 30
---------------------------------------------------------------
Found roots: 3183
Subintervals possibly containing undetected roots: 1
Function evaluations: 34058
cpu/wall(s): 0.032 0.032
```

This does not miss any roots:

```
> build/examples/real_roots 2 1 100
interval: [1, 100]
maxdepth = 30, maxeval = 100000, maxfound = 100000, low_prec = 30
---------------------------------------------------------------
Found roots: 3183
Subintervals possibly containing undetected roots: 0
Function evaluations: 34039
cpu/wall(s): 0.023 0.023
virt/peak/res/peak(MB): 26.32 26.37 2.01 2.01
```

Looking for roots of \( \sin(1/x) \) on \((0, 1)\), the algorithm finds many roots, but will never find all of them since there are infinitely many:

```
> build/examples/real_roots 3 0.0 1.0
interval: [0, 1]
maxdepth = 30, maxeval = 100000, maxfound = 100000, low_prec = 30
---------------------------------------------------------------
Found roots: 10198
Subintervals possibly containing undetected roots: 24695
Function evaluations: 202587
cpu/wall(s): 0.171 0.171
virt/peak/res/peak(MB): 28.39 30.38 4.05 4.05
```
Remark: the program always computes rigorous containing intervals for the roots, but the accuracy after refinement could be less than \( d \) digits.

### 2.5.5 poly\_roots.c

This program finds the complex roots of an integer polynomial by calling `acb\_poly\_find\_roots()` with increasing precision until the roots certainly have been isolated. The program takes the following arguments:

```bash
cpoly\_roots \[-refine d\] \[-print d\] <poly>
```

Isolates all the complex roots of a polynomial with integer coefficients. For convergence, the input polynomial is required to be squarefree.

If \(-\text{refine} d\) is passed, the roots are refined to an absolute tolerance better than \(10^{-d}\). By default, the roots are only computed to sufficient accuracy to isolate them. The refinement is not currently done efficiently.

If \(-\text{print} d\) is passed, the computed roots are printed to \(d\) decimals. By default, the roots are not printed.

The polynomial can be specified by passing the following as \(<\text{poly}>\):

- \(a <n>\): Easy polynomial \(1 + 2x + \ldots + (n+1)x^n\)
- \(t <n>\): Chebyshev polynomial \(T_n\)
- \(u <n>\): Chebyshev polynomial \(U_n\)
- \(p <n>\): Legendre polynomial \(P_n\)
- \(c <n>\): Cyclotomic polynomial \(\Phi_n\)
- \(s <n>\): Swinnerton-Dyer polynomial \(S_n\)
- \(b <n>\): Bernoulli polynomial \(B_n\)
- \(w <n>\): Wilkinson polynomial \(W_n\)
- \(e <n> <m>\): Taylor series of \(\exp(x)\) truncated to degree \(n\)
- \(m <n> <m>\): The Mignotte-like polynomial \(x^n + (100x+1)^m, n > m\)
- \(c0\ c1 \ldots cn\): \(c0 + c1 x + \ldots + cn x^n\) where all \(c:s\) are specified integers

This finds the roots of the Wilkinson polynomial with roots at the positive integers 1, 2, ..., 100:

```bash
> build/examples/poly\_roots -print 15 w 100
prec=53: 0 isolated roots | cpu/wall(s): 0.42 0.426
prec=106: 0 isolated roots | cpu/wall(s): 1.37 1.368
prec=212: 0 isolated roots | cpu/wall(s): 1.48 1.485
prec=424: 100 isolated roots | cpu/wall(s): 0.61 0.611
done!
(1 + 1.7285178043492e-125j) +/- (7.2e-122, 7.2e-122j)
(2 + 5.1605530263601e-122j) +/- (3.77e-118, 3.77e-118j)
(3 + -2.58115555871665e-118j) +/- (5.72e-115, 5.72e-115j)
(4 + 1.02141628524271e-115j) +/- (4.38e-112, 4.38e-112j)
(5 + 1.6132683409498e-113j) +/- (2.6e-109, 2.6e-109j)
...
(95 + 4.15294196875447e-62j) +/- (6.66e-59, 6.66e-59j)
(96 + 3.54502401922667e-64j) +/- (7.37e-60, 7.37e-60j)
(97 + -1.67755995325625e-65j) +/- (6.4e-61, 6.4e-61j)
(98 + 2.04638822352299e-65j) +/- (4e-62, 4e-62j)
(99 + -2.73425468022838e-66j) +/- (1.71e-63, 1.71e-63j)
(100 + -1.0095011302288e-68j) +/- (3.24e-65, 3.24e-65j)
cpu/wall(s): 3.88 3.893
```
This finds the roots of a Bernoulli polynomial which has both real and complex roots. Note that the program does not attempt to determine that the imaginary parts of the real roots really are zero (this could be done by verifying sign changes):

```
> build/examples/poly_roots -refine 100 -print 20 b 16
prec=53: 16 isolated roots | cpu/wall(s): 0 0.007
prec=106: 16 isolated roots | cpu/wall(s): 0 0.004
prec=212: 16 isolated roots | cpu/wall(s): 0 0.004
prec=424: 16 isolated roots | cpu/wall(s): 0 0.004
done!

(-0.94308706466055783383 + -5.512272663168484603e-128j) +/- (2.2e-125, 2.2e-125j)
(-0.75534059252067985752 + 1.937401283040249068e-128j) +/- (1.09e-125, 1.09e-125j)
(-0.24999757119077421009 + -4.5347924422246038692e-130j) +/- (3.6e-127, 3.6e-127j)
(0.24999757152512726002 + 4.219130761823281708e-129j) +/- (4.98e-127, 4.98e-127j)
(0.7500024284787273998 + 9.0360649917413170142e-128j) +/- (8.88e-126, 8.88e-126j)
(1.2499975711907742101 + 7.8804123808107088267e-127j) +/- (2.66e-124, 2.66e-124j)
(1.943070646605578338 + 3.303577342500953239e-125j) +/- (7.05e-123, 7.05e-123j)
(-0.9950933482925623279 + 0.44547958157103608805j) +/- (5.5e-125, 5.5e-125j)
(-0.9950933482925623279 + -0.44547958157103608805j) +/- (5.46e-125, 5.46e-125j)
(1.995093348292562328 + 0.44547958157103608805j) +/- (1.44e-122, 1.44e-122j)
(1.995093348292562328 + -0.44547958157103608805j) +/- (1.43e-122, 1.43e-122j)
(-0.92177327714429290564 + -1.0954360955079385542j) +/- (9.15e-123, 9.15e-123j)
(-0.92177327714429290564 + 1.0954360955079385542j) +/- (8.12e-123, 8.12e-123j)
cpu/wall(s): 0.02 0.02
```

### 2.5.6 complex_plot.c

This program plots one of the predefined functions over a complex interval 
\[ [x_a, x_b] + [y_a, y_b] i \] using domain coloring, at a resolution of \( xn \) times \( yn \) pixels.

The program takes the parameters:

```
complex_plot [-range xa xb ya yb] [-size xn yn] <func>
```

Defaults parameters are \([-10, 10] + [-10, 10] i\) and \(xn = yn = 512\).

The output is written to \(arbplot.ppm\). If you have ImageMagick, run convert arbplot.ppm arbplot.png to get a PNG.

Function codes \(<func>\) are:

- `gamma` - Gamma function
- `digamma` - Digamma function
- `lgamma` - Logarithmic gamma function
- `zeta` - Riemann zeta function
- `erf` - Error function
- `ai` - Airy function Ai
- `bi` - Airy function Bi
- `besselj` - Bessel function \( J_0 \)
- `bessely` - Bessel function \( Y_0 \)
• **besseli** - Bessel function $I_0$
• **besselk** - Bessel function $K_0$
• **modj** - Modular j-function
• **modeta** - Dedekind eta function
• **barnesg** - Barnes G-function
• **agm** - Arithmetic geometric mean

The function is just sampled at point values; no attempt is made to resolve small features by adaptive subsampling.

For example, the following plots the Riemann zeta function around a portion of the critical strip with imaginary part between 100 and 140:

```
> build/examples/complex_plot zeta -range -10 10 100 140 -size 256 512
```
3.1 mag.h – fixed-precision unsigned floating-point numbers for bounds

The `mag_t` type is an unsigned floating-point type with a fixed-precision mantissa (30 bits) and an arbitrary-precision exponent (represented as an `fmpz_t`), suited for representing and rigorously manipulating magnitude bounds efficiently. Operations always produce a strict upper or lower bound, but for performance reasons, no attempt is made to compute the best possible bound (in general, a result may a few ulps larger/smaller than the optimal value). The special values zero and positive infinity are supported (but not NaN). Applications requiring more flexibility (such as correct rounding, or higher precision) should use the `arf_t` type instead.

3.1.1 Types, macros and constants

`mag_struct`

A `mag_struct` holds a mantissa and an exponent. Special values are encoded by the mantissa being set to zero.

`mag_t`

A `mag_t` is defined as an array of length one of type `mag_struct`, permitting a `mag_t` to be passed by reference.

3.1.2 Memory management

void `mag_init (mag_t x)`

Initializes the variable `x` for use. Its value is set to zero.

void `mag_clear (mag_t x)`

Clears the variable `x`, freeing or recycling its allocated memory.

void `mag_init_set (mag_t x, const mag_t y)`

Initializes `x` and sets it to the value of `y`.

void `mag_swap (mag_t x, mag_t y)`

Swaps `x` and `y` efficiently.

void `mag_set (mag_t x, const mag_t y)`

Sets `x` to the value of `y`.

`mag_ptr_mag_vec_init (slong n)`

Allocates a vector of length `n`. All entries are set to zero.
void _mag_vec_clear (mag_ptr v, slong n)
    Clears a vector of length n.

### 3.1.3 Special values

void mag_zero (mag_t x)
    Sets x to zero.

void mag_one (mag_t x)
    Sets x to one.

void mag_inf (mag_t x)
    Sets x to positive infinity.

int mag_is_special (const mag_t x)
    Returns nonzero iff x is zero or positive infinity.

int mag_is_zero (const mag_t x)
    Returns nonzero iff x is zero.

int mag_is_inf (const mag_t x)
    Returns nonzero iff x is positive infinity.

int mag_is_finite (const mag_t x)
    Returns nonzero iff x is not positive infinity (since there is no NaN value, this function is exactly the negation of mag_is_inf()).

### 3.1.4 Comparisons

int mag_equal (const mag_t x, const mag_t y)
    Returns nonzero iff x and y have the same value.

int mag_cmp (const mag_t x, const mag_t y)
    Returns negative, zero, or positive, depending on whether x is smaller, equal, or larger than y.

int mag_cmp_2exp_si (const mag_t x, slong y)
    Returns negative, zero, or positive, depending on whether x is smaller, equal, or larger than $2^y$.

void mag_min (mag_t z, const mag_t x, const mag_t y)
void mag_max (mag_t z, const mag_t x, const mag_t y)
    Sets z respectively to the smaller or the larger of x and y.

### 3.1.5 Input and output

void mag_print (const mag_t x)
    Prints x to standard output.

### 3.1.6 Random generation

void mag_randtest (mag_t x, flint_rand_t state, slong expbits)
    Sets x to a random finite value, with an exponent up to expbits bits large.

void mag_randtest_special (mag_t x, flint_rand_t state, slong expbits)
    Like mag_randtest(), but also sometimes sets x to infinity.
3.1.7 Conversions

void `mag_set_d` (mag_t y, double x)
void `mag_set_fmpr` (mag_t y, const fmp_t x)
void `mag_set_ui` (mag_t y, ulong x)
void `mag_set_fmpz` (mag_t y, const fmpz_t x)
Sets y to an upper bound for |x|.

void `mag_set_d_2exp_fmpz` (mag_t z, double x, const fmpz_t y)
void `mag_set_fmpz_2exp_fmpz` (mag_t z, const fmpz_t x, const fmpz_t y)
void `mag_set_ui_2exp_si` (mag_t z, ulong x, slong y)
Sets z to an upper bound for |x| × 2^y.

void `mag_get_fmpr` (fmp_t y, const mag_t x)
Sets y exactly to x.
void `mag_get_fmpq` (fmpq_t y, const mag_t x)
Sets y exactly to x. Assumes that no overflow occurs.

void `mag_set_ui_lower` (mag_t z, ulong x)
void `mag_set_fmpz_lower` (mag_t z, const fmpz_t x)
Sets y to a lower bound for |x|.

void `mag_set_fmpz_2exp_fmpz_lower` (mag_t z, const fmpz_t x, const fmpz_t y)
Sets z to a lower bound for |x| × 2^y.

3.1.8 Arithmetic

void `mag_mul_2exp_si` (mag_t z, const mag_t x, slong y)
void `mag_mul_2exp_fmpz` (mag_t z, const mag_t x, const fmpz_t y)
Sets z to x × 2^y. This operation is exact.

void `mag_mul` (mag_t z, const mag_t x, const mag_t y)
void `mag_mul_ui` (mag_t z, const mag_t x, ulong y)
void `mag_mul_fmpz` (mag_t z, const mag_t x, const fmpz_t y)
Sets z to an upper bound for xy.

void `mag_add` (mag_t z, const mag_t x, const mag_t y)
Sets z to an upper bound for x + y.

void `mag_addmul` (mag_t z, const mag_t x, const mag_t y)
Sets z to an upper bound for z + xy.

void `mag_add_2exp_fmpz` (mag_t z, const mag_t x, const fmpz_t e)
Sets z to an upper bound for x + 2^e.

void `mag_div` (mag_t z, const mag_t x, const mag_t y)
void `mag_div_ui` (mag_t z, const mag_t x, ulong y)
void `mag_div_fmpz` (mag_t z, const mag_t x, const fmpz_t y)
Sets z to an upper bound for x/y.

void `mag_mul_lower` (mag_t z, const mag_t x, const mag_t y)
3.1.9 Fast, unsafe arithmetic

The following methods assume that all inputs are finite and that all exponents (in all inputs as well as the final result) fit as fmpz inline values. They also assume that the output variables do not have promoted exponents, as they will be overwritten directly (thus leaking memory).

void **mag_fast_init_set** (mag_t x, const mag_t y)
Initialises x and sets it to the value of y.

void **mag_fast_zero** (mag_t x)
Sets x to zero.

int **mag_fast_is_zero** (const mag_t x)
Returns nonzero iff x to zero.

void **mag_fast_mul** (mag_t z, const mag_t x, const mag_t y)
Sets z to an upper bound for \( xy \).

void **mag_fast_addmul** (mag_t z, const mag_t x, const mag_t y)
Sets z to an upper bound for \( z + xy \).

void **mag_fast_add_2exp_si** (mag_t z, const mag_t x, slong e)
Sets z to an upper bound for \( x + 2^e \).

void **mag_fast_mul_2exp_si** (mag_t z, const mag_t x, slong e)
Sets z to an upper bound for \( x \cdot 2^e \).

3.1.10 Powers and logarithms

void **mag_pow_ui** (mag_t z, const mag_t x, ulong e)

void **mag_pow_fmpz** (mag_t z, const mag_t x, const fmpz_t e)
Sets z to an upper bound for \( x^e \). Requires \( e \geq 0 \).

void **mag_pow_ui_lower** (mag_t z, const mag_t x, ulong e)
Sets z to a lower bound for \( x^e \).

void **mag_sqrt** (mag_t z, const mag_t x)
Sets z to an upper bound for \( \sqrt{x} \).

void **mag_rsqrt** (mag_t z, const mag_t x)
Sets z to an upper bound for \( 1/\sqrt{x} \).

void **mag_hypot** (mag_t z, const mag_t x, const mag_t y)
Sets z to an upper bound for \( \sqrt{x^2 + y^2} \).

void **mag_root** (mag_t z, const mag_t x, ulong n)
Sets z to an upper bound for \( x^{1/n} \). We evaluate \( \exp(\log(1 + 2^{kn}x)/n)2^{-k} \), where \( k \) is chosen so that \( 2^{kn}x \approx 2^{30} \).
void **mag_log1p** *(mag_t z, const mag_t x)*  
Sets z to an upper bound for \( \log(1 + x) \). The bound is computed accurately for small x.

void **mag_log_ui** *(mag_t z, ulong n)*  
Sets z to an upper bound for \( \log(n) \).

void **mag_exp** *(mag_t z, const mag_t x)*  
Sets z to an upper bound for \( \exp(x) \).

void **mag_expm1** *(mag_t z, const mag_t x)*  
Sets z to an upper bound for \( \exp(x) - 1 \). The bound is computed accurately for small x.

void **mag_exp_tail** *(mag_t z, const mag_t x, ulong N)*  
Sets z to an upper bound for \( \sum_{k=N}^{\infty} x^k/k! \).

void **mag_binpow_uiui** *(mag_t z, ulong m, ulong n)*  
Sets z to an upper bound for \( (1 + 1/m)^n \).

void **mag_geom_series** *(mag_t res, const mag_t x, ulong N)*  
Sets res to an upper bound for \( \sum_{k=N}^{\infty} x^k \).

### 3.1.11 Special functions

void **mag_fac_ui** *(mag_t z, ulong n)*  
Sets z to an upper bound for \( n! \).

void **mag_rfac_ui** *(mag_t z, ulong n)*  
Sets z to an upper bound for \( 1/n! \).

void **mag_bernoulli_div_fac_ui** *(mag_t z, ulong n)*  
Sets z to an upper bound for \( |B_n|/n! \) where \( B_n \) denotes a Bernoulli number.

void **mag_polylog_tail** *(mag_t u, const mag_t z, slong s, ulong d, ulong N)*  
Sets u to an upper bound for

\[
\sum_{k=N}^{\infty} \frac{z^k \log^d(k)}{k^s}.
\]

Note: in applications where \( s \) in this formula may be real or complex, the user can simply substitute any convenient integer \( s' \) such that \( s' \leq \text{Re}(s) \).

Denote the terms by \( T(k) \). We pick a nonincreasing function \( U(k) \) such that

\[
\frac{T(k + 1)}{T(k)} = z \left( \frac{k}{k + 1} \right)^s \left( \frac{\log(k + 1)}{\log(k)} \right)^d \leq U(k).
\]

Then, as soon as \( U(N) < 1 \),

\[
\sum_{k=N}^{\infty} T(k) \leq T(N) \sum_{k=0}^{\infty} U(N)^k = \frac{T(N)}{1 - U(N)}.
\]

In particular, we take

\[
U(k) = z B(k, \max(0, -s)) B(k \log(k), d)
\]

where \( B(m, n) = (1 + 1/m)^n \). This follows from the bounds

\[
\left( \frac{k}{k + 1} \right)^s \leq \begin{cases} 1 & \text{if } s \geq 0 \\ (1 + 1/k)^{-s} & \text{if } s < 0. \end{cases}
\]
and

\[
\left( \frac{\log(k+1)}{\log(k)} \right)^d \leq \left( 1 + \frac{1}{k \log(k)} \right)^d.
\]

### 3.2 arf.h – arbitrary-precision floating-point numbers

A variable of type `arf_t` holds an arbitrary-precision binary floating-point number, i.e. a rational number of the form \(x \times 2^y\) where \(x, y \in \mathbb{Z}\) and \(x\) is odd; or one of the special values zero, plus infinity, minus infinity, or NaN (not-a-number).

The exponent of a finite and nonzero floating-point number can be defined in different ways: for example, as the component \(y\) above, or as the unique integer \(e\) such that \(x \times 2^y = m \times 2^e\) where \(1/2 \leq |m| < 1\). The internal representation of an `arf_t` stores the exponent in the latter format.

The conventions for special values largely follow those of the IEEE floating-point standard. At the moment, there is no support for negative zero, unsigned infinity, or a NaN with a payload, though some of these might be added in the future.

Except where otherwise noted, the output of an operation is the floating-point number obtained by taking the inputs as exact numbers, in principle carrying out the operation exactly, and rounding the resulting real number to the nearest representable floating-point number whose mantissa has at most the specified number of bits, in the specified direction of rounding. Some operations are always or optionally done exactly.

The `arf_t` type is almost identical semantically to the legacy `fmpr_t` type, but uses a more efficient internal representation. The most significant differences that the user has to be aware of are:

- The mantissa is no longer represented as a FLINT `fmpz`, and the internal exponent points to the top of the binary expansion of the mantissa instead of of the bottom. Code designed to manipulate components of an `fmpr_t` directly can be ported to the `arf_t` type by making use of `arf_get_fmpz_2exp()` and `arf_set_fmpz_2exp()`.

- Some `arf_t` functions return an `int` indicating whether a result is inexact, whereas the corresponding `fmpr_t` functions return an `slong` encoding the relative exponent of the error.

### 3.2.1 Types, macros and constants

**arf_struct**

**arf_t**

An `arf_struct` contains four words: an `fmpz` exponent (`exp`), a `size` field tracking the number of limbs used (one bit of this field is also used for the sign of the number), and two more words. The last two words hold the value directly if there are at most two limbs, and otherwise contain one `alloc` field (tracking the total number of allocated limbs, not all of which might be used) and a pointer to the actual limbs. Thus, up to 128 bits on a 64-bit machine and 64 bits on a 32-bit machine, no space outside of the `arf_struct` is used.

An `arf_t` is defined as an array of length one of type `arf_struct`, permitting an `arf_t` to be passed by reference.

**arf_rnd_t**

Specifies the rounding mode for the result of an approximate operation.

**ARF_RND_DOWN**

Specifies that the result of an operation should be rounded to the nearest representable number in the direction towards zero.
ARF_RND_UP
Specifies that the result of an operation should be rounded to the nearest representable number in the direction
away from zero.

ARF_RND_FLOOR
Specifies that the result of an operation should be rounded to the nearest representable number in the direction
towards minus infinity.

ARF_RND_CEIL
Specifies that the result of an operation should be rounded to the nearest representable number in the direction
towards plus infinity.

ARF_RND_NEAR
Specifies that the result of an operation should be rounded to the nearest representable number, rounding to an
odd mantissa if there is a tie between two values. Warning: this rounding mode is currently not implemented
(except for a few conversions functions where this stated explicitly).

ARF_PREC_EXACT
If passed as the precision parameter to a function, indicates that no rounding is to be performed. This must
only be used when it is known that the result of the operation can be represented exactly and fits in memory
(the typical use case is working with small integer values). Note that, for example, adding two numbers whose
exponents are far apart can easily produce an exact result that is far too large to store in memory.

3.2.2 Memory management

void arf_init (arf_t x)
Initializes the variable x for use. Its value is set to zero.

void arf_clear (arf_t x)
Clears the variable x, freeing or recycling its allocated memory.

3.2.3 Special values

void arf_zero (arf_t x)

void arf_one (arf_t x)

void arf_pos_inf (arf_t x)

void arf_neg_inf (arf_t x)

void arf_nan (arf_t x)
Sets x respectively to 0, 1, +∞, −∞, NaN.

int arf_is_zero (const arf_t x)

int arf_is_one (const arf_t x)

int arf_is_pos_inf (const arf_t x)

int arf_is_neg_inf (const arf_t x)

int arf_is_nan (const arf_t x)
Returns nonzero iff x respectively equals 0, 1, +∞, −∞, NaN.

int arf_is_inf (const arf_t x)
Returns nonzero iff x equals either +∞ or −∞.
int *arf_is_normal* (const *arf_t* x)

Returns nonzero iff *x* is a finite, nonzero floating-point value, i.e. not one of the special values 0, +∞, −∞, NaN.

int *arf_is_special* (const *arf_t* x)

Returns nonzero iff *x* is one of the special values 0, +∞, −∞, NaN, i.e. not a finite, nonzero floating-point value.

int *arf_is_finite* (*arf_t* x)

Returns nonzero iff *x* is a finite floating-point value, i.e. not one of the values +∞, −∞, NaN. (Note that this is not equivalent to the negation of *arf_is_inf*.)

### 3.2.4 Assignment, rounding and conversions

void *arf_set* (*arf_t* y, const *arf_t* x)

void *arf_set_mpz* (*arf_t* y, const *mpz_t* x)

void *arf_set_fmpz* (*arf_t* y, const *fmpz_t* x)

void *arf_set_ui* (*arf_t* y, ulong *x*)

void *arf_set_si* (*arf_t* y, slong *x*)

void *arf_set_mpfr* (*arf_t* y, const *mpfr_t* x)

void *arf_set_fmpr* (*arf_t* y, const *fmpr_t* x)

void *arf_set_d* (*arf_t* y, double *x*)

Sets *y* exactly to *x*.

void *arf_swap* (*arf_t* y, *arf_t* x)

Swaps *y* and *x* efficiently.

void *arf_init_set_ui* (*arf_t* y, ulong *x*)

Initialises *y* and sets it to *x* in a single operation.

int *arf_set_round* (*arf_t* y, const *arf_t* x, slong *prec*, *arf_rnd_t* rnd)

int *arf_set_round_si* (*arf_t* x, slong *v*, slong *prec*, *arf_rnd_t* rnd)

int *arf_set_round_ui* (*arf_t* x, ulong *v*, slong *prec*, *arf_rnd_t* rnd)

int *arf_set_round_mpz* (*arf_t* y, const *mpz_t* *x*, slong *prec*, *arf_rnd_t* rnd)

int *arf_set_round_fmpz* (*arf_t* y, const *fmpz_t* *x*, slong *prec*, *arf_rnd_t* rnd)

Sets *y* exactly to *x*, rounded to *prec* bits in the direction specified by *rnd*.

void *arf_set_si_2exp_si* (*arf_t* y, slong *m*, slong *e*)

void *arf_set_ui_2exp_si* (*arf_t* y, ulong *m*, slong *e*)

void *arf_set_fmpz_2exp* (*arf_t* y, const *fmpz_t* *m*, const *fmpz_t* *e*)

Sets *y* to $m \times 2^e$.

int *arf_set_round_fmpz_2exp* (*arf_t* y, const *fmpz_t* *x*, const *fmpz_t* *e*, slong *prec*, *arf_rnd_t* rnd)

Sets *y* to $x \times 2^e$, rounded to *prec* bits in the direction specified by *rnd*.

void *arf_get_fmpz_2exp* (*fmpz_t* *m*, *fmpz_t* *e*, const *arf_t* x)

Sets *m* and *e* to the unique integers such that $x = m \times 2^e$ and *m* is odd, provided that *x* is a nonzero finite fraction. If *x* is zero, both *m* and *e* are set to zero. If *x* is infinite or NaN, the result is undefined.
double \texttt{arf_get_d} (const \texttt{arf} \texttt{t} \texttt{x}, \texttt{arf_rnd_t} \texttt{rnd})

Returns \texttt{x} rounded to a double in the direction specified by \texttt{rnd}. This method supports rounding to nearest with \texttt{ARF_RND_NEAR}. It also rounds correctly when overflowing or underflowing the double exponent range (this was not the case in an earlier version).

void \texttt{arf_get_fmp} (\texttt{fmp} \texttt{y}, const \texttt{arf} \texttt{t} \texttt{x})

Sets \texttt{y} exactly to \texttt{x}.

\begin{verbatim}
int \texttt{arf_get_mp} (\texttt{mpfr} \texttt{y}, const \texttt{arf} \texttt{t} \texttt{x}, \texttt{mpfr_rnd_t} \texttt{rnd})

Sets the MPFR variable \texttt{y} to the value of \texttt{x}. If the precision of \texttt{x} is too small to allow \texttt{y} to be represented exactly, it is rounded in the specified MPFR rounding mode. The return value (-1, 0 or 1) indicates the direction of rounding, following the convention of the MPFR library.
\end{verbatim}

\begin{verbatim}
void \texttt{arf_get_fmp} (\texttt{fmpz} \texttt{z}, const \texttt{arf} \texttt{t} \texttt{x}, \texttt{arf_rnd_t} \texttt{rnd})

Sets \texttt{z} to \texttt{x} rounded to the nearest integer in the direction specified by \texttt{rnd}. If \texttt{rnd} is \texttt{ARF_RND_NEAR}, rounds to the nearest even integer in case of a tie. Aborts if \texttt{x} is infinite, NaN or if the exponent is unreasonably large.
\end{verbatim}

\begin{verbatim}
slong \texttt{arf_get_s} (const \texttt{arf} \texttt{t} \texttt{x}, \texttt{arf_rnd_t} \texttt{rnd})

Returns \texttt{x} rounded to the nearest integer in the direction specified by \texttt{rnd}. If \texttt{rnd} is \texttt{ARF_RND_NEAR}, rounds to the nearest even integer in case of a tie. Aborts if \texttt{x} is infinite, NaN, or the value is too large to fit in a slong.
\end{verbatim}

\begin{verbatim}
int \texttt{arf_get_fmp} (fmpz_t \texttt{y}, const \texttt{arf} \texttt{t} \texttt{x}, \texttt{arf_rnd_t} \texttt{rnd})

Sets \texttt{y} exactly to \texttt{x}.
\end{verbatim}

\begin{verbatim}
int \texttt{arf_equal} (const \texttt{arf} \texttt{t} \texttt{x}, const \texttt{arf} \texttt{t} \texttt{y})

Returns nonzero iff \texttt{x} and \texttt{y} are exactly equal. This function does not treat NaN specially, i.e. NaN compares as equal to itself.
\end{verbatim}

\begin{verbatim}
int \texttt{arf_equal_s} (const \texttt{arf} \texttt{t} \texttt{x}, slong \texttt{y})

Returns nonzero iff \texttt{x} and \texttt{y} are exactly equal. This function does not treat NaN specially, i.e. NaN compares as equal to itself.
\end{verbatim}

\begin{verbatim}
int \texttt{arf_cmp} (const \texttt{arf} \texttt{t} \texttt{x}, const \texttt{arf} \texttt{t} \texttt{y})

Returns negative, zero, or positive, depending on whether \texttt{x} is respectively smaller, equal, or greater compared to \texttt{y}. Comparison with NaN is undefined.
\end{verbatim}

\begin{verbatim}
int \texttt{arf_cmpabs} (const \texttt{arf} \texttt{t} \texttt{x}, const \texttt{arf} \texttt{t} \texttt{y})

Compares the absolute values of \texttt{x} and \texttt{y}.
\end{verbatim}

\begin{verbatim}
int \texttt{arf_cmpabs_ui} (const \texttt{arf} \texttt{t} \texttt{x}, ulong \texttt{y})

Compares the absolute values of \texttt{x} and \texttt{y}.
\end{verbatim}

\begin{verbatim}
int \texttt{arf_cmpabs_mag} (const \texttt{arf} \texttt{t} \texttt{x}, const \texttt{mag} \texttt{t} \texttt{y})

Compares the absolute values of \texttt{x} and \texttt{y}.
\end{verbatim}

\begin{verbatim}
int \texttt{arf_cmp_2exp_s} (const \texttt{arf} \texttt{t} \texttt{x}, slong \texttt{e})

Compares \texttt{x} (respectively its absolute value) with \texttt{2}^\texttt{e}.
\end{verbatim}

\begin{verbatim}
int \texttt{arf_sgn} (const \texttt{arf} \texttt{t} \texttt{x})

Returns -1, 0 or +1 according to the sign of \texttt{x}. The sign of NaN is undefined.
\end{verbatim}

\begin{verbatim}
void \texttt{arf_min} (\texttt{arf} \texttt{z}, const \texttt{arf} \texttt{t} \texttt{a}, const \texttt{arf} \texttt{t} \texttt{b})

\end{verbatim}

3.2.5 Comparisons and bounds
void **arf_max** (arf_t z, const arf_t a, const arf_t b)
Sets z respectively to the minimum and the maximum of a and b.

slong **arf_bits** (const arf_t x)
Returns the number of bits needed to represent the absolute value of the mantissa of x, i.e. the minimum precision sufficient to represent x exactly. Returns 0 if x is a special value.

int **arf_is_int** (const arf_t x)
Returns nonzero iff x is integer-valued.

int **arf_is_int_2exp_si** (const arf_t x, slong e)
Returns nonzero iff x equals \( n \cdot 2^e \) for some integer \( n \).

void **arf_abs_bound_lt_2exp_fmpz** (fmpz_t b, const arf_t x)
Sets b to the smallest integer such that \(|x| < 2^b\). If x is zero, infinity or NaN, the result is undefined.

void **arf_abs_bound_le_2exp_fmpz** (fmpz_t b, const arf_t x)
Sets b to the smallest integer such that \(|x| \leq 2^b\). If x is zero, infinity or NaN, the result is undefined.

slong **arf_abs_bound_lt_2exp_si** (const arf_t x)
Returns the smallest integer \( b \) such that \(|x| < 2^b\), clamping the result to lie between -ARF_PREC_EXACT and ARF_PREC_EXACT inclusive. If x is zero, -ARF_PREC_EXACT is returned, and if x is infinity or NaN, ARF_PREC_EXACT is returned.

### 3.2.6 Magnitude functions

void **arf_get_mag** (mag_t y, const arf_t x)
Sets y to an upper bound for the absolute value of x.

void **arf_get_mag_lower** (mag_t y, const arf_t x)
Sets y to a lower bound for the absolute value of x.

void **arf_set_mag** (arf_t y, const mag_t x)
Sets y to x.

void **mag_init_set_arf** (mag_t y, const arf_t x)
Initializes y and sets it to an upper bound for x.

void **mag_fast_init_set_arf** (mag_t y, const arf_t x)
Initializes y and sets it to an upper bound for x. Assumes that the exponent of y is small.

void **arf_mag_set_ulp** (mag_t z, const arf_t y, slong prec)
Sets z to the magnitude of the unit in the last place (ulp) of y at precision prec.

void **arf_mag_add_ulp** (mag_t z, const mag_t x, const arf_t y, slong prec)
Sets z to an upper bound for the sum of x and the magnitude of the unit in the last place (ulp) of y at precision prec.

void **arf_mag_fast_add_ulp** (mag_t z, const mag_t x, const arf_t y, slong prec)
Sets z to an upper bound for the sum of x and the magnitude of the unit in the last place (ulp) of y at precision prec. Assumes that all exponents are small.

### 3.2.7 Shallow assignment

void **arf_init_set_shallow** (arf_t z, const arf_t x)

void **arf_init_set_mag_shallow** (arf_t z, const mag_t x)
Initializes z to a shallow copy of x. A shallow copy just involves copying struct data (no heap allocation is performed).
The target variable \( z \) may not be cleared or modified in any way (it can only be used as constant input to functions), and may not be used after \( x \) has been cleared. Moreover, after \( x \) has been assigned shallowly to \( z \), no modification of \( x \) is permitted as long as \( z \) is in use.

```c
void arf_init_neg_shallow (arf_t z, const arf_t x)
```
Initializes \( z \) shallowly to the negation of \( x \).

### 3.2.8 Random number generation

```c
void arf_init_neg_mag_shallow (arf_t z, const mag_t x)
```

Initializes \( z \) shallowly to the negation of \( x \).

```c
void arf_randtest (arf_t x, flint_rand_t state, slong bits, slong mag_bits)
```
Generates a finite random number whose mantissa has precision at most \( \text{bits} \) and whose exponent has at most \( \text{mag_bits} \) bits. The values are distributed non-uniformly: special bit patterns are generated with high probability in order to allow the test code to exercise corner cases.

```c
void arf_randtest_not_zero (arf_t x, flint_rand_t state, slong bits, slong mag_bits)
```
Identical to \( \text{arf_randtest}() \), except that zero is never produced as an output.

```c
void arf_randtest_special (arf_t x, flint_rand_t state, slong bits, slong mag_bits)
```
Identical to \( \text{arf_randtest}() \), except that the output occasionally is set to an infinity or NaN.

### 3.2.9 Input and output

```c
void arf_debug (const arf_t x)
```
Prints information about the internal representation of \( x \).

```c
void arf_print (const arf_t x)
```
Prints \( x \) as an integer mantissa and exponent.

```c
void arf_printd (const arf_t y, slong d)
```
Prints \( x \) as a decimal floating-point number, rounding to \( d \) digits. This function is currently implemented using MPFR, and does not support large exponents.

### 3.2.10 Addition and multiplication

```c
void arf_abs (arf_t y, const arf_t x)
```
Sets \( y \) to the absolute value of \( x \).

```c
void arf_neg (arf_t y, const arf_t x)
```
Sets \( y = -x \) exactly.

```c
int arf_neg_round (arf_t y, const arf_t x, slong prec, arf_rnd_t rnd)
```
Sets \( y = -x \), rounded to \( \text{prec} \) bits in the direction specified by \( \text{rnd} \), returning nonzero iff the operation is inexact.

```c
void arf_mul_2exp_si (arf_t y, const arf_t x, slong e)
```
Sets \( y = x \times 2^e \) exactly.

```c
int arf_mul (arf_t z, const arf_t x, const arf_t y, const mpz_t prec, arf_rnd_t rnd)
```

```c
int arf_mul_ui (arf_t z, const arf_t x, ulong y, const mpz_t prec, arf_rnd_t rnd)
```

```c
int arf_mul_si (arf_t z, const arf_t x, slong y, const mpz_t prec, arf_rnd_t rnd)
```

```c
int arf_mul_mpz (arf_t z, const arf_t x, const mpz_t y, const mpz_t prec, arf_rnd_t rnd)
```
int \texttt{arf_mul_fmpz} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{const fmpz_t} \, y, \texttt{const slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

Sets \( z = x \times y \), rounded to \( \texttt{prec} \) bits in the direction specified by \( \texttt{rnd} \), returning nonzero iff the operation is inexact.

int \texttt{arf_add} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{arf_t} \, y, \texttt{const slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

int \texttt{arf_add_si} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{const slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

int \texttt{arf_add_ui} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{ulong} \, y, \texttt{const slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

int \texttt{arf_add_fmpz} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{arf_t} \, y, \texttt{const slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

int \texttt{arf_add} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{arf_t} \, y, \texttt{slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

int \texttt{arf_add_si} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

int \texttt{arf_add_ui} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{ulong} \, y, \texttt{slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

int \texttt{arf_add_fmpz} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{arf_t} \, y, \texttt{slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

Sets \( z = x + y \), rounded to \( \texttt{prec} \) bits in the direction specified by \( \texttt{rnd} \), returning nonzero iff the operation is inexact.

int \texttt{arf_add_fmpz_2exp} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{const fmpz_t} \, y, \texttt{const fmpz_t} \, e, \texttt{const slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

Sets \( z = x + y 2^e \), rounded to \( \texttt{prec} \) bits in the direction specified by \( \texttt{rnd} \), returning nonzero iff the operation is inexact.

int \texttt{arf_sub} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{arf_t} \, y, \texttt{const slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

int \texttt{arf_sub_si} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

int \texttt{arf_sub_ui} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{ulong} \, y, \texttt{slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

int \texttt{arf_sub_fmpz} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{arf_t} \, y, \texttt{slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

Sets \( z = x - y \), rounded to \( \texttt{prec} \) bits in the direction specified by \( \texttt{rnd} \), returning nonzero iff the operation is inexact.

int \texttt{arf_addmul} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{arf_t} \, y, \texttt{const slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

int \texttt{arf_addmul_ui} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{ulong} \, y, \texttt{const slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

int \texttt{arf_addmul_si} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{slong} \, y, \texttt{const slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

int \texttt{arf_addmul_fmpz} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{arf_t} \, y, \texttt{const slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

Sets \( z = z + x \times y \), rounded to \( \texttt{prec} \) bits in the direction specified by \( \texttt{rnd} \), returning nonzero iff the operation is inexact.

int \texttt{arf_submul} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{arf_t} \, y, \texttt{const slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

int \texttt{arf_submul_ui} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{ulong} \, y, \texttt{const slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

int \texttt{arf_submul_si} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{slong} \, y, \texttt{const slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

int \texttt{arf_submul_fmpz} (\texttt{arf_t} \, z, \texttt{arf_t} \, x, \texttt{arf_t} \, y, \texttt{const slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

Sets \( z = z - x \times y \), rounded to \( \texttt{prec} \) bits in the direction specified by \( \texttt{rnd} \), returning nonzero iff the operation is inexact.

### 3.2.11 Summation

int \texttt{arf_sum} (\texttt{arf_t} \, s, \texttt{arf_srcptr} \, \texttt{terms}, \texttt{const slong} \, \texttt{len}, \texttt{const slong} \, \texttt{prec}, \texttt{arf_rnd_t} \, \texttt{rnd})

Sets \( s \) to the sum of the array \( \texttt{terms} \) of length \( \texttt{len} \), rounded to \( \texttt{prec} \) bits in the direction specified by \( \texttt{rnd} \). The sum is computed as if done without any intermediate rounding error, with only a single rounding applied to the final result. Unlike repeated calls to \texttt{arf_add()} with infinite precision, this function does not overflow if the magnitudes of the terms are far apart. Warning: this function is implemented naively, and the running time is quadratic with respect to \( \texttt{len} \) in the worst case.
3.2.12 Division

int arf_div (arf_t z, const arf_t x, const arf_t y, slong prec, arf_rnd_t rnd)
int arf_div_ui (arf_t z, const arf_t x, ulong y, slong prec, arf_rnd_t rnd)
int arf_ui_div (arf_t z, ulong x, const arf_t y, slong prec, arf_rnd_t rnd)
int arf_div_si (arf_t z, const arf_t x, slong y, slong prec, arf_rnd_t rnd)
int arf_si_div (arf_t z, slong x, const arf_t y, slong prec, arf_rnd_t rnd)
int arf_div_fmpz (arf_t z, const arf_t x, const fmpz_t y, slong prec, arf_rnd_t rnd)
int arf_fmpz_div (arf_t x, const fmpz_t y, slong prec, arf_rnd_t rnd)
int arf_fmpz_div_fmpz (arf_t z, const fmpz_t x, const arf_t y, slong prec, arf_rnd_t rnd)

Sets \( z = x/y \), rounded to \( \text{prec} \) bits in the direction specified by \( \text{rnd} \), returning nonzero iff the operation is inexact. The result is NaN if \( y \) is zero.

3.2.13 Square roots

int arf_sqrt (arf_t z, const arf_t x, slong prec, arf_rnd_t rnd)
int arf_sqrt_ui (arf_t z, const arf_t x, ulong prec, arf_rnd_t rnd)
int arf_sqrt_fmpz (arf_t z, const arf_t x, const fmpz_t y, slong prec, arf_rnd_t rnd)
int arf_fmpz_div (arf_t x, const fmpz_t y, slong prec, arf_rnd_t rnd)
int arf_fmpz_div_fmpz (arf_t z, const fmpz_t x, const arf_t y, slong prec, arf_rnd_t rnd)

Sets \( z = \sqrt{x} \), rounded to \( \text{prec} \) bits in the direction specified by \( \text{rnd} \), returning nonzero iff the operation is inexact. The result is NaN if \( x \) is negative.

int arf_rsqrt (arf_t z, const arf_t x, slong prec, arf_rnd_t rnd)

Sets \( z = 1/\sqrt{x} \), rounded to \( \text{prec} \) bits in the direction specified by \( \text{rnd} \), returning nonzero iff the operation is inexact. The result is NaN if \( x \) is negative, and \( +\infty \) if \( x \) is zero.

int arf_root (arf_t z, const arf_t x, ulong k, slong prec, arf_rnd_t rnd)

Sets \( z = x^{1/k} \), rounded to \( \text{prec} \) bits in the direction specified by \( \text{rnd} \), returning nonzero iff the operation is inexact. The result is NaN if \( x \) is negative. Warning: this function is a wrapper around the MPFR root function. It gets slow and uses much memory for large \( k \).

3.2.14 Complex arithmetic

int arf_complex_mul (arf_t e, arf_t f, const arf_t a, const arf_t b, const arf_t c, const arf_t d, slong prec, arf_rnd_t rnd)
int arf_complex_mulFallback (arf_t e, arf_t f, const arf_t a, const arf_t b, const arf_t c, const arf_t d, slong prec, arf_rnd_t rnd)

Computes the complex product \( e + fi = (a + bi)(c + di) \), rounding both \( e \) and \( f \) correctly to \( \text{prec} \) bits in the direction specified by \( \text{rnd} \). The first bit in the return code indicates inexactness of \( e \), and the second bit indicates inexactness of \( f \).

If any of the components \( a, b, c, d \) is zero, two real multiplications and no additions are done. This convention is used even if any other part contains an infinity or NaN, and the behavior with infinite/NaN input is defined accordingly.

The fallback version is implemented naively, for testing purposes. No squaring optimization is implemented.

int arf_complex_sqr (arf_t e, arf_t f, const arf_t a, const arf_t b, slong prec, arf_rnd_t rnd)

Computes the complex square \( e + fi = (a + bi)^2 \). This function has identical semantics to arf_complex_mul() (with \( c = a, b = d \)), but is faster.

3.2. arf.h – arbitrary-precision floating-point numbers
3.3 arb.h – real numbers represented as floating-point balls

An \texttt{arb_t} represents a ball over the real numbers, that is, an interval \([m \pm r] \equiv [m-r, m+r]\) where the midpoint \(m\) and the radius \(r\) are (extended) real numbers and \(r\) is nonnegative (possibly infinite). The result of an (approximate) operation done on \texttt{arb_t} variables is a ball which contains the result of the (mathematically exact) operation applied to any choice of points in the input balls. In general, the output ball is not the smallest possible.

The precision parameter passed to each function roughly indicates the precision to which calculations on the midpoint are carried out (operations on the radius are always done using a fixed, small precision.)

For arithmetic operations, the precision parameter currently simply specifies the precision of the corresponding \texttt{arf_t} operation. In the future, the arithmetic might be made faster by incorporating sloppy rounding (typically equivalent to a loss of 1-2 bits of effective working precision) when the result is known to be inexact (while still propagating errors rigorously, of course). Arithmetic operations done on exact input with exactly representable output are always guaranteed to produce exact output.

For more complex operations, the precision parameter indicates a minimum working precision (algorithms might allocate extra internal precision to attempt to produce an output accurate to the requested number of bits, especially when the required precision can be estimated easily, but this is not generally required).

If the precision is increased and the inputs either are exact or are computed with increased accuracy as well, the output should converge proportionally, absent any bugs. The general intended strategy for using ball arithmetic is to add a few guard bits, and then repeat the calculation as necessary with an exponentially increasing number of guard bits (Ziv’s strategy) until the result is exact enough for one’s purposes (typically the first attempt will be successful).

The following balls with an infinite or NaN component are permitted, and may be returned as output from functions.

- The ball \([+\infty \pm c]\), where \(c\) is finite, represents the point at positive infinity. Such a ball can always be replaced by \([+\infty \pm 0]\) while preserving mathematical correctness (this is currently not done automatically by the library).
- The ball \([-\infty \pm c]\), where \(c\) is finite, represents the point at negative infinity. Such a ball can always be replaced by \([-\infty \pm 0]\) while preserving mathematical correctness (this is currently not done automatically by the library).
- The ball \([c \pm \infty]\), where \(c\) is finite or infinite, represents the whole extended real line \([-\infty, +\infty]\). Such a ball can always be replaced by \([0 \pm \infty]\) while preserving mathematical correctness (this is currently not done automatically by the library). Note that there is no way to represent a half-infinite interval such as \([0, \infty]\).
- The ball \([\text{NaN} \pm c]\), where \(c\) is finite or infinite, represents an indeterminate value (the value could be any extended real number, or it could represent a function being evaluated outside its domain of definition, for example where the result would be complex). Such an indeterminate ball can always be replaced by \([\text{NaN} \pm \infty]\) while preserving mathematical correctness (this is currently not done automatically by the library).

The \texttt{arb_t} type is almost identical semantically to the legacy \texttt{fmprb_t} type, but uses a more efficient internal representation. Whereas the midpoint and radius of an \texttt{fmprb_t} both have the same type, the \texttt{arb_t} type uses an \texttt{arf_t} for the midpoint and a \texttt{mag_t} for the radius. Code designed to manipulate the radius of an \texttt{fmprb_t} directly can be ported to the \texttt{arb_t} type by writing the radius to a temporary \texttt{arf_t} variable, manipulating that variable, and then converting back to the \texttt{mag_t} radius. Alternatively, \texttt{mag_t} methods can be used directly where available.

3.3.1 Types, macros and constants

\texttt{arb_struct}

\texttt{arb_t}

An \texttt{arb_struct} consists of an \texttt{arf_t} (the midpoint) and a \texttt{mag_t} (the radius). An \texttt{arb_t} is defined as an array of length one of type \texttt{arb_struct}, permitting an \texttt{arb_t} to be passed by reference.

\texttt{arb_ptr}

Alias for \texttt{arb_struct *}, used for vectors of numbers.
**arb_srcptr**
Alias for `const arb_struct *`, used for vectors of numbers when passed as constant input to functions.

**arb_midref** (x)
Macro returning a pointer to the midpoint of \( x \) as an `arf_t`.

**arb_radref** (x)
Macro returning a pointer to the radius of \( x \) as a `mag_t`.

### 3.3.2 Memory management

**void arb_init** (arb_t \( x \))
Initializes the variable \( x \) for use. Its midpoint and radius are both set to zero.

**void arb_clear** (arb_t \( x \))
Clears the variable \( x \), freeing or recycling its allocated memory.

**arb_ptr _arb_vec_init** (slong \( n \))
Returns a pointer to an array of \( n \) initialized `arb_struct` entries.

**void _arb_vec_clear** (arb_ptr \( v \), slong \( n \))
Clears an array of \( n \) initialized `arb_struct` entries.

**void arb_swap** (arb_t \( x \), arb_t \( y \))
Swaps \( x \) and \( y \) efficiently.

### 3.3.3 Assignment and rounding

**void arb_set_fmprb** (arb_t \( y \), const fmprb_t \( x \))

**void arb_get_fmprb** (fmprb_t \( y \), const arb_t \( x \))

**void arb_set** (arb_t \( y \), const arb_t \( x \))

**void arb_set_arf** (arb_t \( y \), const arf_t \( x \))

**void arb_set_si** (arb_t \( y \), slong \( x \))

**void arb_set_ui** (arb_t \( y \), ulong \( x \))

**void arb_set_d** (arb_t \( y \), double \( x \))

**void arb_set_fmpz** (arb_t \( y \), const fmpz_t \( x \))
Sets \( y \) to the value of \( x \) without rounding.

**void arb_set_fmpz_2exp** (arb_t \( y \), const fmpz_t \( x \), const fmpz_t \( e \))
Sets \( y \) to \( x \cdot 2^e \).

**void arb_set_round** (arb_t \( y \), const arb_t \( x \), slong \( prec \))

**void arb_set_round_fmpz** (arb_t \( y \), const fmpz_t \( x \), slong \( prec \))
Sets \( y \) to the value of \( x \), rounded to \( prec \) bits.

**void arb_set_round_fmpz_2exp** (arb_t \( y \), const fmpz_t \( x \), const fmpz_t \( e \), slong \( prec \))
Sets \( y \) to \( x \cdot 2^e \), rounded to \( prec \) bits.

**void arb_set_fmpq** (arb_t \( y \), const fmpq_t \( x \), slong \( prec \))
Sets \( y \) to the rational number \( x \), rounded to \( prec \) bits.
int **arb_set_str** (*arb* res, const char *inp, slong prec)
Sets res to the value specified by the human-readable string inp. The input may be a decimal floating-point literal, such as “25”, “0.001”, “7e+141” or “-31.4159e-1”, and may also consist of two such literals separated by the symbol “+/-” and optionally enclosed in brackets, e.g. “[3.25 +/- 0.0001]”, or simply “[+/- 10]” with an implicit zero midpoint. The output is rounded to prec bits, and if the binary-to-decimal conversion is inexact, the resulting error is added to the radius.

The symbols “inf” and “nan” are recognized (a nan midpoint results in an indeterminate interval, with infinite radius).

Returns 0 if successful and nonzero if unsuccessful. If unsuccessful, the result is set to an indeterminate interval.

char *arb_get_str (const *arb x, slong n, ulong flags)
Returns a nice human-readable representation of x, with at most n digits of the midpoint printed.

With default flags, the output can be parsed back with **arb_set_str()**, and this is guaranteed to produce an interval containing the original interval x.

By default, the output is rounded so that the value given for the midpoint is correct up to 1 ulp (unit in the last decimal place).

If **ARB_STR_MORE** is added to flags, more (possibly incorrect) digits may be printed.

If **ARB_STR_NO_RADIUS** is added to flags, the radius is not included in the output if at least 1 digit of the midpoint can be printed.

By adding a multiple m of **ARB_STR_CONDENSE** to flags, strings of more than three times m consecutive digits are condensed, only printing the leading and trailing m digits along with brackets indicating the number of digits omitted (useful when computing values to extremely high precision).

### 3.3.4 Assignment of special values

void **arb_zero** (*arb* f)
Sets x to zero.

void **arb_one** (*arb* f)
Sets x to the exact integer 1.

void **arb_pos_inf** (*arb* x)
Sets x to positive infinity, with a zero radius.

void **arb_neg_inf** (*arb* x)
Sets x to negative infinity, with a zero radius.

void **arb_zero_pm_inf** (*arb* x)
Sets x to [0 ±∞], representing the whole extended real line.

void **arb_indeterminate** (*arb* x)
Sets x to [NaN ±∞], representing an indeterminate result.

### 3.3.5 Input and output

void **arb_print** (const *arb* x)
Prints the internal representation of x.

void **arb_printd** (const *arb* x, slong digits)
Prints x in decimal. The printed value of the radius is not adjusted to compensate for the fact that the binary-to-decimal conversion of both the midpoint and the radius introduces additional error.
void **arb_printn** (const **arb_t** x, slong digits, ulong flags)
Prints a nice decimal representation of x. By default, the output is guaranteed to be correct to within one unit in the last digit. An error bound is also printed explicitly. See **arb_get_str()** for details.

### 3.3.6 Random number generation

void **arb_randtest** (**arb_t** x, flint_rand_t state, slong prec, slong mag_bits)
Generates a random ball. The midpoint and radius will both be finite.

void **arb_randtest_exact** (**arb_t** x, flint_rand_t state, slong prec, slong mag_bits)
Generates a random number with zero radius.

void **arb_randtest_precise** (**arb_t** x, flint_rand_t state, slong prec, slong mag_bits)
Generates a random number with magnitude of the midpoint around $2^{-\text{prec}}$.

void **arb_randtest_wide** (**arb_t** x, flint_rand_t state, slong prec, slong mag_bits)
Generates a random number with midpoint and radius chosen independently, possibly giving a very large interval.

void **arb_randtest_special** (**arb_t** x, flint_rand_t state, slong prec, slong mag_bits)
Generates a random interval, possibly having NaN or an infinity as the midpoint and possibly having an infinite radius.

void **arb_get_rand_fmpq** (**fmpq_t** q, flint_rand_t state, const **arb_t** x, slong bits)
Sets q to a random rational number from the interval represented by x. A denominator is chosen by multiplying the binary denominator of x by a random integer up to bits bits.

The outcome is undefined if the midpoint or radius of x is non-finite, or if the exponent of the midpoint or radius is so large or small that representing the endpoints as exact rational numbers would cause overflows.

### 3.3.7 Radius and interval operations

void **arb_get_mid_arb** (**arb_t** m, const **arb_t** x)
Sets m to the midpoint of x.

void **arb_get_rad_arb** (**arb_t** r, const **arb_t** x)
Sets r to the radius of x.

void **arb_add_error_arf** (**arb_t** x, const **arf_t** err)
void **arb_add_error_mag** (**arb_t** x, const **mag_t** err)
void **arb_add_error** (**arb_t** x, const **arb_t** err)
Adds the absolute value of err to the radius of x (the operation is done in-place).

void **arb_add_error_2exp_si** (**arb_t** x, slong e)
void **arb_add_error_2exp_fmpz** (**arb_t** x, const **fmpz_t** e)
Adds $2^e$ to the radius of x.

void **arb_union** (**arb_t** z, const **arb_t** x, const **arb_t** y, slong prec)
Sets z to a ball containing both x and y.

void **arb_get_abs_ubound_arf** (**arf_t** u, const **arb_t** x, slong prec)
Sets u to the upper bound for the absolute value of x, rounded up to prec bits. If x contains NaN, the result is NaN.

void **arb_get_abs_lbound_arf** (**arf_t** u, const **arb_t** x, slong prec)
Sets u to the lower bound for the absolute value of x, rounded down to prec bits. If x contains NaN, the result is NaN.
void \textbf{arb_get_mag} (mag_t z, const arb_t x)

Sets \(z\) to an upper bound for the absolute value of \(x\). If \(x\) contains NaN, the result is positive infinity.

void \textbf{arb_get_mag_lower} (mag_t z, const arb_t x)

Sets \(z\) to a lower bound for the absolute value of \(x\). If \(x\) contains NaN, the result is zero.

void \textbf{arb_get_mag_lower_nonnegative} (mag_t z, const arb_t x)

Sets \(z\) to a lower bound for the signed value of \(x\), or zero if \(x\) overlaps with the negative half-axis. If \(x\) contains NaN, the result is zero.

void \textbf{arb_get_interval_fmpz_2exp} (fmpz_t a, fmpz_t b, fmpz_t exp, const arb_t x)

Computes the exact interval represented by \(x\), in the form of an integer interval multiplied by a power of two, i.e. \(x = [a, b] \times 2^{\exp}\).

The outcome is undefined if the midpoint or radius of \(x\) is non-finite, or if the difference in magnitude between the midpoint and radius is so large that representing the endpoints exactly would cause overflows.

void \textbf{arb_set_interval_arf} (arb_t x, const arf_t a, const arf_t b, slong prec)

void \textbf{arb_set_interval_mpfr} (arb_t x, const mpfr_t a, const mpfr_t b, slong prec)

Sets \(x\) to a ball containing the interval \([a, b]\). We require that \(a \leq b\).

void \textbf{arb_get_interval_arf} (arf_t a, arf_t b, const arb_t x, slong prec)

void \textbf{arb_get_interval_mpfr} (mpfr_t a, mpfr_t b, const arb_t x)

Constructs an interval \([a, b]\) containing the ball \(x\). The MPFR version uses the precision of the output variables.

slong \textbf{arb_rel_error_bits} (const arb_t x)

Returns the effective relative error of \(x\) measured in bits, defined as the difference between the position of the top bit in the radius and the top bit in the midpoint, plus one. The result is clamped between plus/minus ARF\_PREC\_EXACT.

slong \textbf{arb_rel_accuracy_bits} (const arb_t x)

Returns the effective relative accuracy of \(x\) measured in bits, equal to the negative of the return value from \textbf{arb_rel_error_bits}().

slong \textbf{arb_bits} (const arb_t x)

Returns the number of bits needed to represent the absolute value of the mantissa of the midpoint of \(x\), i.e. the minimum precision sufficient to represent \(x\) exactly. Returns 0 if the midpoint of \(x\) is a special value.

void \textbf{arb_trim} (arb_t y, const arb_t x)

Sets \(y\) to a trimmed copy of \(x\): rounds \(x\) to a number of bits equal to the accuracy of \(x\) (as indicated by its radius), plus a few guard bits. The resulting ball is guaranteed to contain \(x\), but is more economical if \(x\) has less than full accuracy.

int \textbf{arb_get_unique_fmpz} (fmpz_t z, const arb_t x)

If \(x\) contains a unique integer, sets \(z\) to that value and returns nonzero. Otherwise (if \(x\) represents no integers or more than one integer), returns zero.

void \textbf{arb_floor} (arb_t y, const arb_t x, slong prec)

void \textbf{arb ceil} (arb_t y, const arb_t x, slong prec)

Sets \(y\) to a ball containing \([x]\) and \([x]\) respectively, with the midpoint of \(y\) rounded to at most \(prec\) bits.

void \textbf{arb_get_fmpz_mid_rad_10exp} (fmpz_t mid, fmpz_t rad, fmpz_t exp, const arb_t x, slong n)

Assuming that \(x\) is finite and not exactly zero, computes integers \(mid, rad, exp\) such that \(x \in [m - r, m + r] \times 10^{e}\) and such that the larger out of \(mid\) and \(rad\) has at least \(n\) digits plus a few guard digits. If \(x\) is infinite or exactly zero, the outputs are all set to zero.
### 3.3.8 Comparisons

```c
int arb_is_zero (const arb_t x)
Returns nonzero iff the midpoint and radius of x are both zero.
```

```c
int arb_is_nonzero (const arb_t x)
Returns nonzero iff zero is not contained in the interval represented by x.
```

```c
int arb_is_one (const arb_t x)
Returns nonzero iff x is exactly 1.
```

```c
int arb_is_finite (const arb_t x)
Returns nonzero iff the midpoint and radius of x are both finite floating-point numbers, i.e. not infinities or NaN.
```

```c
int arb_is_exact (const arb_t x)
Returns nonzero iff the radius of x is zero.
```

```c
int arb_is_int (const arb_t x)
Returns nonzero iff x is an exact integer.
```

```c
int arb_equal (const arb_t x, const arb_t y)
Returns nonzero iff x and y are equal as balls, i.e. have both the same midpoint and radius.
```

Note that this is not the same thing as testing whether both x and y certainly represent the same real number, unless either x or y is exact (and neither contains NaN). To test whether both operands might represent the same mathematical quantity, use `arb_overlaps()` or `arb_contains()`, depending on the circumstance.

```c
int arb_is_positive (const arb_t x)
int arb_is_nonnegative (const arb_t x)
int arb_is_negative (const arb_t x)
int arb_is_nonpositive (const arb_t x)
```

```c
int arb_overlaps (const arb_t x, const arb_t y)
Returns nonzero iff x and y have some point in common. If either x or y contains NaN, this function always returns nonzero (as a NaN could be anything, it could in particular contain any number that is included in the other operand).
```

```c
int arb_contains_arf (const arb_t x, const arf_t y)
int arb_contains_fmpq (const arb_t x, const fmpq_t y)
int arb_contains_fmpz (const arb_t x, const fmpz_t y)
int arb_contains_si (const arb_t x, slong y)
int arb_contains_mpfr (const arb_t x, const mpfr_t y)
int arb_contains (const arb_t x, const arb_t y)
Returns nonzero iff the given number (or ball) y is contained in the interval represented by x.
```

If x is contains NaN, this function always returns nonzero (as it could represent anything, and in particular could represent all the points included in y). If y contains NaN and x does not, it always returns zero.

```c
int arb_contains_int (const arb_t x)
Returns nonzero iff the interval represented by x contains an integer.
```

```c
int arb_contains_zero (const arb_t x)
int arb_contains_negative (const arb_t x)
```
int arb_contains_nonpositive (const arb_t x)
int arb_contains_positive (const arb_t x)
int arb_contains_nonnegative (const arb_t x)
  Returns nonzero iff there is any point p in the interval represented by x satisfying, respectively, p = 0, p < 0, p ≤ 0, p > 0, p ≥ 0. If x contains NaN, returns nonzero.
int arb_eq (const arb_t x, const arb_t y)
int arb_ne (const arb_t x, const arb_t y)
int arb_lt (const arb_t x, const arb_t y)
int arb_le (const arb_t x, const arb_t y)
int arb_gt (const arb_t x, const arb_t y)
int arb_ge (const arb_t x, const arb_t y)
  Respectively performs the comparison x = y, x ≠ y, x < y, x ≤ y, x > y, x ≥ y in a mathematically meaningful way. If the comparison t (op) u holds for all t ∈ x and all u ∈ y, returns 1. Otherwise, returns 0.
The balls x and y are viewed as subintervals of the extended real line. Note that balls that are formally different can compare as equal under this definition: for example, [−∞ ± 3] = [−∞ ± 0]. Also [−∞] ≤ [∞ ± ∞]. The output is always 0 if either input has NaN as midpoint.

3.3.9 Arithmetic

void arb_neg (arb_t y, const arb_t x)
void arb_neg_round (arb_t y, const arb_t x, slong prec)
  Sets y to the negation of x.
void arb_abs (arb_t y, const arb_t x)
  Sets y to the absolute value of x. No attempt is made to improve the interval represented by x if it contains zero.
void arb_add (arb_t z, const arb_t x, const arb_t y, slong prec)
void arb_add_arf (arb_t z, const arb_t x, const arf_t y, slong prec)
void arb_add_ui (arb_t z, const arb_t x, ulong y, slong prec)
void arb_add_si (arb_t z, const arb_t x, slong y, slong prec)
void arb_add_fmpz (arb_t z, const arb_t x, const fmpz_t y, slong prec)
  Sets z = x + y, rounded to prec bits. The precision can be ARF_PREC_EXACT provided that the result fits in memory.
void arb_add_fmpz_2exp (arb_t z, const arb_t x, const fmpz_t m, const fmpz_t e, slong prec)
  Sets z = x + m · 2^e, rounded to prec bits. The precision can be ARF_PREC_EXACT provided that the result fits in memory.
void arb_sub (arb_t z, const arb_t x, const arb_t y, slong prec)
void arb_sub_arf (arb_t z, const arb_t x, const arf_t y, slong prec)
void arb_sub_ui (arb_t z, const arb_t x, ulong y, slong prec)
void arb_sub_si (arb_t z, const arb_t x, slong y, slong prec)
void arb_sub_fmpz (arb_t z, const arb_t x, const fmpz_t y, slong prec)
  Sets z = x − y, rounded to prec bits. The precision can be ARF_PREC_EXACT provided that the result fits in memory.
void \texttt{arb\_mul}\,(arb\_t z,\ const\ arb\_t x,\ const\ arb\_t y,\ slong\ prec) \\
void \texttt{arb\_mul\_arf}\,(arb\_t z,\ const\ arb\_t x,\ const\ arf\_t y,\ slong\ prec) \\
void \texttt{arb\_mul\_si}\,(arb\_t z,\ const\ arb\_t x,\ slong\ y,\ slong\ prec) \\
void \texttt{arb\_mul\_ui}\,(arb\_t z,\ const\ arb\_t x,\ ulong\ y,\ slong\ prec) \\
void \texttt{arb\_mul\_fmpz}\,(arb\_t z,\ const\ arb\_t x,\ const\ fmpz\_t y,\ slong\ prec) \\
\hspace{1em}Sets\ z = x \cdot y,\ rounded\ to\ prec\ bits.\ The\ precision\ can\ be\ \texttt{ARF\_PREC\_EXACT}\ provided\ that\ the\ result\ fits\ in\ memory. \\
void \texttt{arb\_mul\_2exp\_si}\,(arb\_t y,\ const\ arb\_t x,\ slong\ e) \\
\hspace{1em}Sets\ y\ to\ x\ multiplied\ by\ 2^e. \\
void \texttt{arb\_addmul}\,(arb\_t z,\ const\ arb\_t x,\ const\ arb\_t y,\ slong\ prec) \\
void \texttt{arb\_addmul\_arf}\,(arb\_t z,\ const\ arb\_t x,\ const\ arf\_t y,\ slong\ prec) \\
void \texttt{arb\_addmul\_si}\,(arb\_t z,\ const\ arb\_t x,\ slong\ y,\ slong\ prec) \\
void \texttt{arb\_addmul\_ui}\,(arb\_t z,\ const\ arb\_t x,\ ulong\ y,\ slong\ prec) \\
void \texttt{arb\_addmul\_fmpz}\,(arb\_t z,\ const\ arb\_t x,\ const\ fmpz\_t y,\ slong\ prec) \\
\hspace{1em}Sets\ z = z + x \cdot y,\ rounded\ to\ prec\ bits.\ The\ precision\ can\ be\ \texttt{ARF\_PREC\_EXACT}\ provided\ that\ the\ result\ fits\ in\ memory. \\
void \texttt{arb\_submul}\,(arb\_t z,\ const\ arb\_t x,\ const\ arb\_t y,\ slong\ prec) \\
void \texttt{arb\_submul\_arf}\,(arb\_t z,\ const\ arb\_t x,\ const\ arf\_t y,\ slong\ prec) \\
void \texttt{arb\_submul\_si}\,(arb\_t z,\ const\ arb\_t x,\ slong\ y,\ slong\ prec) \\
void \texttt{arb\_submul\_ui}\,(arb\_t z,\ const\ arb\_t x,\ ulong\ y,\ slong\ prec) \\
void \texttt{arb\_submul\_fmpz}\,(arb\_t z,\ const\ arb\_t x,\ const\ fmpz\_t y,\ slong\ prec) \\
\hspace{1em}Sets\ z = z - x \cdot y,\ rounded\ to\ prec\ bits.\ The\ precision\ can\ be\ \texttt{ARF\_PREC\_EXACT}\ provided\ that\ the\ result\ fits\ in\ memory. \\
void \texttt{arb\_inv}\,(arb\_t y,\ const\ arb\_t x,\ slong\ prec) \\
\hspace{1em}Sets\ z\ to\ 1/x. \\
void \texttt{arb\_div}\,(arb\_t z,\ const\ arb\_t x,\ const\ arb\_t y,\ slong\ prec) \\
void \texttt{arb\_div\_arf}\,(arb\_t z,\ const\ arb\_t x,\ const\ arf\_t y,\ slong\ prec) \\
void \texttt{arb\_div\_si}\,(arb\_t z,\ const\ arb\_t x,\ slong\ y,\ slong\ prec) \\
void \texttt{arb\_div\_ui}\,(arb\_t z,\ const\ arb\_t x,\ ulong\ y,\ slong\ prec) \\
void \texttt{arb\_fmpz\_div\_fmpz}\,(arb\_t z,\ const\ arb\_t x,\ const\ fmpz\_t y,\ slong\ prec) \\
void \texttt{arb\_fmpz\_div\_mpz}\,(arb\_t z,\ const\ arb\_t x,\ const\ mpz\_t y,\ slong\ prec) \\
\hspace{1em}Sets\ z = x/y,\ rounded\ to\ prec\ bits.\ If\ y\ contains\ zero,\ z\ is\ set\ to\ 0 \pm \infty.\ Otherwise,\ error\ propagation\ uses\ the\ rule \\
\begin{equation*}
\left|\frac{x}{y} - \frac{x + \xi_1 a}{y + \xi_2 b}\right| = \frac{|x|}{|y|} \frac{1}{|\xi_1 a + \xi_2 b|} \leq \frac{|x|}{|y|} + \frac{|y a|}{|y|(|y| - b)}
\end{equation*}
\hspace{1em}where -1 \leq \xi_1, \xi_2 \leq 1,\ and\ where\ the\ triangle\ inequality\ has\ been\ applied\ to\ the\ numerator\ and\ the\ reverse\ triangle\ inequality\ has\ been\ applied\ to\ the\ denominator. \\
void \texttt{arb\_div\_2exp\_ui}\,(arb\_t z,\ const\ arb\_t x,\ ulong\ n,\ slong\ prec) \\
\hspace{1em}Sets\ z = x/(2^n - 1),\ rounded\ to\ prec\ bits.
### 3.3.10 Powers and roots

void \textbf{arb_sqrt} \((\texttt{arb}_z, \texttt{const arb}_x, \texttt{slong prec})\)

void \textbf{arb_sqrt_arf} \((\texttt{arb}_z, \texttt{const arb}_x, \texttt{slong prec})\)

void \textbf{arb_sqrt_fmpz} \((\texttt{arb}_z, \texttt{const fmpz}_t x, \texttt{slong prec})\)

void \textbf{arb_sqrt_ui} \((\texttt{arb}_z, \texttt{ulong} x, \texttt{slong prec})\)

Sets \(z\) to the square root of \(x\), rounded to \(\texttt{prec}\) bits.

If \(x = m \pm r\) where \(m \geq r \geq 0\), the propagated error is bounded by \(\sqrt{m} - \sqrt{m-r} = \sqrt{m}(1 - \sqrt{1-r/m}) \leq \sqrt{m}(r/m + (r/m)^2)/2\).

void \textbf{arb_sqrtpos} \((\texttt{arb}_z, \texttt{const arb}_x, \texttt{slong prec})\)

Sets \(z\) to the square root of \(x\), assuming that \(x\) represents a nonnegative number (i.e. discarding any negative numbers in the input interval).

void \textbf{arb_hypot} \((\texttt{arb}_z, \texttt{const arb}_x, \texttt{const arb}_y, \texttt{slong prec})\)

Sets \(z\) to \(\sqrt{x^2 + y^2}\).

void \textbf{arb_rsqrt} \((\texttt{arb}_z, \texttt{const arb}_x, \texttt{slong prec})\)

void \textbf{arb_rsqrt_ui} \((\texttt{arb}_z, \texttt{ulong} x, \texttt{slong prec})\)

Sets \(z\) to the reciprocal square root of \(x\), rounded to \(\texttt{prec}\) bits. At high precision, this is faster than computing a square root.

void \textbf{arb_sqrt1pm1} \((\texttt{arb}_z, \texttt{const arb}_x, \texttt{slong prec})\)

Sets \(z = 1 + x - 1\), computed accurately when \(x \approx 0\).

void \textbf{arb_root} \((\texttt{arb}_z, \texttt{const arb}_x, \texttt{ulong} k, \texttt{slong prec})\)

Sets \(z\) to the \(k\)-th root of \(x\), rounded to \(\texttt{prec}\) bits. This function selects between different algorithms. For large \(k\), it evaluates \(\exp\left(\log(x)/k\right)\). For small \(k\), it uses \texttt{arf_root()} at the midpoint and computes a propagated error bound as follows: if input interval is \([m-r, m+r]\) with \(r \leq m\), the error is largest at \(m-r\) where it satisfies

\[
m^{1/k} - (m-r)^{1/k} = m^{1/k}[1 - (1 - r/m)^{1/k}] \\
= m^{1/k}[1 - \exp(\log(1-r/m)/k)] \\
\leq m^{1/k} \min(1, -\log(1-r/m)/k) \\
= m^{1/k} \min(1, \log(1+r/(m-r))/k).
\]

This is evaluated using \texttt{mag_log1p()}.  

void \textbf{arb_root} \((\texttt{arb}_z, \texttt{const arb}_x, \texttt{ulong} k, \texttt{slong prec})\)

Alias for \texttt{arb_root_ui()}, provided for backwards compatibility.

void \textbf{arb_pow_fmpz\_binexp} \((\texttt{arb}_y, \texttt{const arb}_b, \texttt{const fmpz}_t e, \texttt{slong prec})\)

void \textbf{arb_pow_fmpz} \((\texttt{arb}_y, \texttt{const arb}_b, \texttt{const fmpz}_t e, \texttt{slong prec})\)

void \textbf{arb_pow_ui} \((\texttt{arb}_y, \texttt{const arb}_b, \texttt{ulong} e, \texttt{slong prec})\)

void \textbf{arb\_ui\_pow\_ui} \((\texttt{arb}_y, \texttt{ulong} b, \texttt{ulong} e, \texttt{slong prec})\)

void \textbf{arb\_si\_pow\_ui} \((\texttt{arb}_y, \texttt{long} b, \texttt{ulong} e, \texttt{slong prec})\)

Sets \(y = b^e\) using binary exponentiation (with an initial division if \(e < 0\)). Provided that \(b\) and \(e\) are small enough and the exponent is positive, the exact power can be computed by setting the precision to \texttt{ARF\_PREC\_EXACT}.

Note that these functions can get slow if the exponent is extremely large (in such cases \texttt{arb\_pow()} may be superior).

void \textbf{arb_pow\_fmpq} \((\texttt{arb}_y, \texttt{const arb}_x, \texttt{const fmpq}_t a, \texttt{slong prec})\)

Sets \(y = b^e\), computed as \(y = (b^{1/q})^q\) if the denominator of \(e = p/q\) is small, and generally as \(y = \exp(e \log b)\).
3.3.11 Exponentials and logarithms

void \texttt{arb\_pow} (\texttt{arb\_t \textbf{z}}, \texttt{const arb\_t \textbf{x}}, \texttt{const arb\_t \textbf{y}}, \texttt{slong \textit{prec}})

Sets \( z = x^y \), computed using binary exponentiation if \( y \) is a small exact integer, as \( z = (x^{1/2})^y \) if \( y \) is a small exact half-integer, and generally as \( z = \exp(y \log x) \).

At low to medium precision (up to about 4096 bits), \texttt{arb\_log\_arf()} uses table-based argument reduction and fast Taylor series evaluation via \_arb\_atan\_taylor\_rs(). At high precision, it falls back to MPFR.

The function \texttt{arb\_log()} simply calls \texttt{arb\_log\_arf()} with the midpoint as input, and separately adds the propagated error.

void \texttt{arb\_log\_ui\_from\_prev} (\texttt{arb\_t \textbf{z}}, \texttt{ulong \textbf{x}}, \texttt{slong \textit{prec}})

Computes \( \log(k_1) \), given \( \log(k_0) \) where \( k_0 < k_1 \). At high precision, this function uses the formula \( \log(k_1) = \log(k_0) + 2 \arctanh((k_1 - k_0)/(k_1 + k_0)) \), evaluating the inverse hyperbolic tangent using binary splitting (for best efficiency, \( k_0 \) should be large and \( k_1 - k_0 \) should be small). Otherwise, it ignores \( \log(k_0) \) and evaluates the logarithm the usual way.

void \texttt{arb\_log\_lp} (\texttt{arb\_t \textbf{z}}, \texttt{const arb\_t \textbf{x}}, \texttt{slong \textit{prec}})

Sets \( z = \log(1 + x) \), computed accurately when \( x \approx 0 \).

void \texttt{arb\_exp} (\texttt{arb\_t \textbf{z}}, \texttt{const arb\_t \textbf{x}}, \texttt{slong \textit{prec}})

Sets \( z = \exp(x) \). Error propagation is done using the following rule: assuming \( x = m \pm r \), the error is largest at \( m + r \), and we have \( \exp(m + r) - \exp(m) = \exp(m)(\exp(r) - 1) \leq r \exp(m + r) \).

void \texttt{arb\_exp\_m1} (\texttt{arb\_t \textbf{z}}, \texttt{const arb\_t \textbf{x}}, \texttt{slong \textit{prec}})

Sets \( z = \exp(x) - 1 \), computed accurately when \( x \approx 0 \).

void \texttt{arb\_exp\_invexp} (\texttt{arb\_t \textbf{z}}, \texttt{arb\_t \textbf{w}}, \texttt{const arb\_t \textbf{x}}, \texttt{slong \textit{prec}})

Sets \( z = \exp(x) \) and \( w = \exp(-x) \). The second exponential is computed from the first using a division, but propagated error bounds are computed separately.

3.3.12 Trigonometric functions

void \texttt{arb\_sin} (\texttt{arb\_t \textbf{s}}, \texttt{const arb\_t \textbf{x}}, \texttt{slong \textit{prec}})

void \texttt{arb\_cos} (\texttt{arb\_t \textbf{c}}, \texttt{const arb\_t \textbf{x}}, \texttt{slong \textit{prec}})

void \texttt{arb\_sin\_cos} (\texttt{arb\_t \textbf{s}}, \texttt{arb\_t \textbf{c}}, \texttt{const arb\_t \textbf{x}}, \texttt{slong \textit{prec}})

Sets \( s = \sin(x) \), \( c = \cos(x) \). Error propagation uses the rule \( |\sin(m \pm r) - \sin(m)| \leq \min(r, 2) \).

void \texttt{arb\_sin\_pi} (\texttt{arb\_t \textbf{s}}, \texttt{const arb\_t \textbf{x}}, \texttt{slong \textit{prec}})

void \texttt{arb\_cos\_pi} (\texttt{arb\_t \textbf{c}}, \texttt{const arb\_t \textbf{x}}, \texttt{slong \textit{prec}})

void \texttt{arb\_sin\_cos\_pi} (\texttt{arb\_t \textbf{s}}, \texttt{arb\_t \textbf{c}}, \texttt{const arb\_t \textbf{x}}, \texttt{slong \textit{prec}})

Sets \( s = \sin(\pi x) \), \( c = \cos(\pi x) \).
void \texttt{arb\_tan} (\texttt{arb\_t} y, const \texttt{arb\_t} x, slong \texttt{prec})
   
   Sets $y = \tan(x) = \sin(x) / \cos(y)$.

void \texttt{arb\_cot} (\texttt{arb\_t} y, const \texttt{arb\_t} x, slong \texttt{prec})
   
   Sets $y = \cot(x) = \cos(x) / \sin(y)$.

void \texttt{arb\_sin\_cos\_pi\_fmpq} (\texttt{arb\_t} s, \texttt{arb\_t} c, const \texttt{fmpq\_t} x, slong \texttt{prec})
   
   Sets $s = \sin(\pi x)$, $c = \cos(\pi x)$ where $x$ is a rational number (whose numerator and denominator are assumed to be reduced). We first use trigonometric symmetries to reduce the argument to the octant $[0, 1/4]$. Then we either multiply by a numerical approximation of $\pi$ and evaluate the trigonometric function the usual way, or we use algebraic methods, depending on which is estimated to be faster. Since the argument has been reduced to the first octant, the first of these two methods gives full accuracy even if the original argument is close to some root other the origin.

void \texttt{arb\_tan\_pi} (\texttt{arb\_t} y, const \texttt{arb\_t} x, slong \texttt{prec})
   
   Sets $y = \tan(\pi x)$.

void \texttt{arb\_cot\_pi} (\texttt{arb\_t} y, const \texttt{arb\_t} x, slong \texttt{prec})
   
   Sets $y = \cot(\pi x)$.

void \texttt{arb\_sinc} (\texttt{arb\_t} z, const \texttt{arb\_t} x, slong \texttt{prec})
   
   Sets $z = \text{sinc}(x) = \sin(x)/x$.

### 3.3.13 Inverse trigonometric functions

void \texttt{arb\_atan\_arf} (\texttt{arb\_t} z, const \texttt{arf\_t} x, slong \texttt{prec})
   
   The function \texttt{arb\_atan\_arf()} uses lookup tables if possible, and otherwise falls back to \texttt{arb\_atan\_arf\_bb()}. At low to medium precision (up to about 4096 bits), \texttt{arb\_atan\_arf()} uses table-based argument reduction and fast Taylor series evaluation via \_\texttt{arb\_atan\_taylor\_rs()}. At high precision, it falls back to MPFR. The function \texttt{arb\_atan()} simply calls \texttt{arb\_atan\_arf()} with the midpoint as input, and separately adds the propagated error.

void \texttt{arb\_atan2} (\texttt{arb\_t} z, const \texttt{arb\_t} b, const \texttt{arb\_t} a, slong \texttt{prec})
   
   Sets $r$ to an the argument (phase) of the complex number $a + bi$, with the branch cut discontinuity on $(-\infty, 0]$. We define \texttt{atan2}(0, 0) = 0, and for $a < 0$, \texttt{atan2}(0, a) = \pi$.

void \texttt{arb\_asin} (\texttt{arb\_t} z, const \texttt{arb\_t} x, slong \texttt{prec})
   
   Sets $z = \text{asin}(x) = \text{atan}(x/\sqrt{1 - x^2})$. If $x$ is not contained in the domain $[-1, 1]$, the result is an indeterminate interval.

void \texttt{arb\_acos} (\texttt{arb\_t} z, const \texttt{arb\_t} x, slong \texttt{prec})
   
   Sets $z = \text{acos}(x) = \pi/2 - \text{asin}(x)$. If $x$ is not contained in the domain $[-1, 1]$, the result is an indeterminate interval.

### 3.3.14 Hyperbolic functions

void \texttt{arb\_sinh} (\texttt{arb\_t} s, const \texttt{arb\_t} x, slong \texttt{prec})
   
void \texttt{arb\_cosh} (\texttt{arb\_t} c, const \texttt{arb\_t} x, slong \texttt{prec})
void arb_sinh_cosh (arb_t s, arb_t c, const arb_t x, slong prec)
    Sets \( s = \sinh(x) \), \( c = \cosh(x) \). If the midpoint of \( x \) is close to zero and the hyperbolic sine is to be computed, evaluates \((e^{2x} \pm 1)/(2e^x)\) via arb_expm1() to avoid loss of accuracy. Otherwise evaluates \((e^x \pm e^{-x})/2\).

void arb_tanh (arb_t y, const arb_t x, slong prec)
    Sets \( y = \tanh(x) = \sinh(x)/\cosh(x) \), evaluated via arb_expm1() as \( \tanh(x) = (e^{2x} - 1)/(e^{2x} + 1) \) if \( |x| \) is small, and as \( \tanh(\pm x) = 1 - 2e^{\mp 2x}/(1 + e^{\mp 2x}) \) if \( |x| \) is large.

void arb_coth (arb_t y, const arb_t x, slong prec)
    Sets \( y = \coth(x) = \cosh(x)/\sinh(x) \), evaluated using the same strategy as arb_tanh().

3.3.15 Inverse hyperbolic functions

void arb_atanh (arb_t z, const arb_t x, slong prec)
    Sets \( z = \operatorname{atanh}(x) \).

void arb_asinh (arb_t z, const arb_t x, slong prec)
    Sets \( z = \operatorname{asinh}(x) \).

void arb_acosh (arb_t z, const arb_t x, slong prec)
    Sets \( z = \operatorname{acosh}(x) \). If \( x < 1 \), the result is an indeterminate interval.

3.3.16 Constants

The following functions cache the computed values to speed up repeated calls at the same or lower precision. For further implementation details, see Algorithms for mathematical constants.

void arb_const_pi (arb_t z, slong prec)
    Computes \( \pi \).

void arb_const_sqrt_pi (arb_t z, slong prec)
    Computes \( \sqrt{\pi} \).

void arb_const_log_sqrt2pi (arb_t z, slong prec)
    Computes \( \log(\sqrt{2\pi}) \).

void arb_const_log2 (arb_t z, slong prec)
    Computes \( \log(2) \).

void arb_const_log10 (arb_t z, slong prec)
    Computes \( \log(10) \).

void arb_const_euler (arb_t z, slong prec)
    Computes Euler's constant \( \gamma = \lim_{k \to \infty} (H_k - \log k) \) where \( H_k = 1 + 1/2 + \ldots + 1/k \).

void arb_const_catalan (arb_t z, slong prec)
    Computes Catalan's constant \( C = \sum_{n=0}^{\infty} (-1)^n/(2n + 1)^2 \).

void arb_const_e (arb_t z, slong prec)
    Computes \( e = \exp(1) \).

void arb_const_khinchin (arb_t z, slong prec)
    Computes Khinchin's constant \( K_0 \).

void arb_const_glaisher (arb_t z, slong prec)
    Computes the Glaisher-Kinkelin constant \( A = \exp(1/12 - \zeta'(-1)) \).

void arb_const_apery (arb_t z, slong prec)
    Computes Apery's constant \( \zeta(3) \).
3.3.17 Gamma function and factorials

void \texttt{arb\_rising\_ui\_bs} (\texttt{arb\_t z}, const \texttt{arb\_t x}, ulong \texttt{n}, slong \texttt{prec})
void \texttt{arb\_rising\_ui\_rs} (\texttt{arb\_t z}, const \texttt{arb\_t x}, ulong \texttt{n}, ulong \texttt{step}, slong \texttt{prec})
void \texttt{arb\_rising\_ui\_rec} (\texttt{arb\_t z}, const \texttt{arb\_t x}, ulong \texttt{n}, slong \texttt{prec})
void \texttt{arb\_rising\_ui} (\texttt{arb\_t z}, const \texttt{arb\_t x}, ulong \texttt{n}, slong \texttt{prec})

void \texttt{arb\_rising} (\texttt{arb\_t z}, const \texttt{arb\_t x}, const \texttt{arb\_t n}, slong \texttt{prec})

Computes the rising factorial \( z = x(x+1)(x+2) \cdots (x+n-1) \).

The bs version uses binary splitting. The rs version uses rectangular splitting. The rec version uses either bs or rs depending on the input. The default version uses the gamma function unless \( n \) is a small integer.

The rs version takes an optional step parameter for tuning purposes (to use the default step length, pass zero).

void \texttt{arb\_rising\_fmpq\_ui} (\texttt{arb\_t z}, const \texttt{fmpq\_t x}, ulong \texttt{n}, slong \texttt{prec})

Computes the rising factorial \( z = x(x+1)(x+2) \cdots (x+n-1) \) using binary splitting. If the denominator or numerator of \( x \) is large compared to \texttt{prec}, it is more efficient to convert \( x \) to an approximation and use \texttt{arb\_rising\_ui}().

void \texttt{arb\_rising2\_ui\_bs} (\texttt{arb\_t u}, \texttt{arb\_t v}, const \texttt{arb\_t x}, ulong \texttt{n}, slong \texttt{prec})
void \texttt{arb\_rising2\_ui\_rs} (\texttt{arb\_t u}, \texttt{arb\_t v}, const \texttt{arb\_t x}, ulong \texttt{n}, ulong \texttt{step}, slong \texttt{prec})

Letting \( u(x) = x(x+1)(x+2) \cdots (x+n-1) \), simultaneously compute \( u(x) \) and \( v(x) = u'(x) \), respectively using binary splitting, rectangular splitting (with optional nonzero step length \texttt{step} to override the default choice), and an automatic algorithm choice.

void \texttt{arb\_fac\_ui} (\texttt{arb\_t z}, ulong \texttt{n}, slong \texttt{prec})
Computes the factorial \( z = n! \) via the gamma function.

void \texttt{arb\_doublefac\_ui} (\texttt{arb\_t z}, ulong \texttt{n}, slong \texttt{prec})
Computes the double factorial \( z = n!! \) via the gamma function.

void \texttt{arb\_bin\_ui} (\texttt{arb\_t z}, const \texttt{arb\_t n}, ulong \texttt{k}, slong \texttt{prec})
Computes the binomial coefficient \( z = \binom{n}{k} \), via the rising factorial as \( \binom{n}{k} = (n-k+1)_k/k! \).

void \texttt{arb\_gamma} (\texttt{arb\_t z}, const \texttt{arb\_t x}, slong \texttt{prec})

void \texttt{arb\_gamma\_fmpq} (\texttt{arb\_t z}, const \texttt{fmpq\_t x}, slong \texttt{prec})
void \texttt{arb\_gamma\_fmpz} (\texttt{arb\_t z}, const \texttt{fmpz\_t x}, slong \texttt{prec})
Computes the gamma function \( z = \Gamma(x) \).

void \texttt{arb\_lgamma} (\texttt{arb\_t z}, const \texttt{arb\_t x}, slong \texttt{prec})
Computes the logarithmic gamma function \( z = \log \Gamma(x) \). The complex branch structure is assumed, so if \( x \leq 0 \), the result is an indeterminate interval.

void \texttt{arb\_rgamma} (\texttt{arb\_t z}, const \texttt{arb\_t x}, slong \texttt{prec})
Computes the reciprocal gamma function \( z = 1/\Gamma(x) \), avoiding division by zero at the poles of the gamma function.

void \texttt{arb\_digamma} (\texttt{arb\_t y}, const \texttt{arb\_t x}, slong \texttt{prec})
Computes the digamma function \( z = \psi(x) = (\log \Gamma(x))' = \Gamma'(x)/\Gamma(x) \).
3.3. arb.h – real numbers represented as floating-point balls

### 3.3.18 Zeta function

**void arb_zeta_ui_vec_borwein (arb_ptr z, ulong start, slong num, ulong step, slong prec)**  
Evaluates $\zeta(s)$ at num consecutive integers $s$ beginning with start and proceeding in increments of step. Uses Borwein’s formula ([Bor2000], [GS2003]), implemented to support fast multi-evaluation (but also works well for a single $s$).

Requires start $\geq 2$. For efficiency, the largest $s$ should be at most about as large as prec. Arguments approaching LONG_MAX will cause overflows. One should therefore only use this function for $s$ up to about prec, and then switch to the Euler product.

The algorithm for single $s$ is basically identical to the one used in MPFR (see [MPFR2012] for a detailed description). In particular, we evaluate the sum backwards to avoid storing more than one $d_k$ coefficient, and use integer arithmetic throughout since it is convenient and the terms turn out to be slightly larger than $2^{\text{prec}}$. The only numerical error in the main loop comes from the division by $k^s$, which adds less than 1 unit of error per term. For fast multi-evaluation, we repeatedly divide by $k^{\text{step}}$. Each division reduces the input error and adds at most 1 unit of additional rounding error, so by induction, the error per term is always smaller than 2 units.

**void arb_zeta_ui_asympt (arb_t x, ulong s, slong prec)**  
Assuming $s \geq 2$, approximates $\zeta(s)$ by $1 + 2^{-s}$ along with a correct error bound. We use the following bounds: for $s > b$, $\zeta(s) - 1 < 2^{-b}$, and generally, $\zeta(s) - (1 + 2^{-s}) < 2^{1 - [s/2]}$.

**void arb_zeta_ui_euler_product (arb_t z, ulong s, slong prec)**  
Computes $\zeta(s)$ using the Euler product. This is fast only if $s$ is large compared to the precision.

Writing $P(a, b) = \prod_{a \leq p \leq b}(1 - p^{-s})$, we have $1/\zeta(s) = P(a, M)P(M + 1, \infty)$.

To bound the error caused by truncating the product at $M$, we write $P(M + 1, \infty) = 1 - \epsilon(s, M)$. Since $0 < P(a, M) \leq 1$, the absolute error for $\zeta(s)$ is bounded by $\epsilon(s, M)$.

According to the analysis in [Fil1992], it holds for all $s \geq 6$ and $M \geq 1$ that $1/\zeta(s) - 1 - \epsilon(s, M) \equiv 2M^{1-s}/(s^2/2 - 1)$. Thus, we have $1/(1 - \epsilon(s, M)) - 1 \leq f(s, M)$, and expanding the geometric series allows us to conclude that $\epsilon(M) \leq f(s, M)$.

**void arb_zeta_ui_bernoulli (arb_t x, ulong s, slong prec)**  
Computes $\zeta(s)$ for even $s$ via the corresponding Bernoulli number.

**void arb_zeta_ui_borwein_bsplit (arb_t x, ulong s, slong prec)**  
Computes $\zeta(s)$ for arbitrary $s \geq 2$ using a binary splitting implementation of Borwein’s algorithm. This has quasilinear complexity with respect to the precision (assuming that $s$ is fixed).

**void arb_zeta_ui_vec (arb_ptr x, ulong start, slong num, slong prec)**

**void arb_zeta_ui_vec_even (arb_ptr x, ulong start, slong num, slong prec)**

**void arb_zeta_ui_vec_odd (arb_ptr x, ulong start, slong num, slong prec)**

Computes $\zeta(s)$ at num consecutive integers (respectively num even or num odd integers) beginning with $s = \text{start} \geq 2$, automatically choosing an appropriate algorithm.

**void arb_zeta (arb_t x, const arb_t s, slong prec)**

Sets $z$ to the value of the Riemann zeta function $\zeta(s)$.

For computing derivatives with respect to $s$, use arb_poly_zeta_series().

**void hurwitz_zeta (arb_t x, const arb_t s, const arb_t a, slong prec)**

Sets $z$ to the value of the Hurwitz zeta function $\zeta(s, a)$. 
For computing derivatives with respect to $s$, use \texttt{arb\_poly\_zeta\_series()}. 

### 3.3.19 Bernoulli numbers

void \texttt{arb\_bernoulli\_ui (arb\_t b, ulong n, slong prec)}

Sets $b$ to the numerical value of the Bernoulli number $B_n$ accurate to $prec$ bits, computed by a division of the exact fraction if $B_n$ is in the global cache or the exact numerator roughly is larger than $prec$ bits, and using \texttt{arb\_bernoulli\_ui\_zeta()} otherwise. This function reads $B_n$ from the global cache if the number is already cached, but does not automatically extend the cache by itself.

void \texttt{arb\_bernoulli\_ui\_zeta (arb\_t b, ulong n, slong prec)}

Sets $b$ to the numerical value of $B_n$ accurate to $prec$ bits, computed using the formula $B_{2n} = \frac{(-1)^{n+1}(2(2n))!\zeta(2n)}{(2\pi)^n}$.

To avoid potential infinite recursion, we explicitly call the Euler product implementation of the zeta function. We therefore assume that the precision is small enough and $n$ large enough for the Euler product to converge rapidly (otherwise this function will effectively hang).

void \texttt{arb\_power\_sum\_vec (arb\_ptr res, const arb\_t a, const arb\_t b, slong len, slong prec)}

For $n$ from 0 to $len - 1$, sets entry $n$ in the output vector $res$ to 

$$S_n(a, b) = \frac{1}{n+1} (B_{n+1}(b) - B_{n+1}(a))$$

where $B_n(x)$ is a Bernoulli polynomial. If $a$ and $b$ are integers and $b \geq a$, this is equivalent to

$$S_n(a, b) = \sum_{k=a}^{b-1} k^n.$$ 

The computation uses the generating function for Bernoulli polynomials.

### 3.3.20 Polylogarithms

void \texttt{arb\_polylog (arb\_t w, const arb\_t s, const arb\_t z, slong prec)}

void \texttt{arb\_polylog\_si (arb\_t w, slong s, const arb\_t z, slong prec)}

Sets $w$ to the polylogarithm $\text{Li}_s(z)$.

### 3.3.21 Other special functions

void \texttt{arb\_fib\_fmpz (arb\_t z, const fmpz\_t n, slong prec)}

void \texttt{arb\_fib\_ui (arb\_t z, ulong n, slong prec)}

Computes the Fibonacci number $F_n$. Uses the binary squaring algorithm described in \cite{Tak2000}. Provided that $n$ is small enough, an exact Fibonacci number can be computed by setting the precision to \texttt{ARF\_PREC\_EXACT}.

void \texttt{arb\_agm (arb\_t z, const arb\_t x, const arb\_t y, slong prec)}

Sets $z$ to the arithmetic-geometric mean of $x$ and $y$.

void \texttt{arb\_chebyshev\_t\_ui (arb\_t a, ulong n, const arb\_t x, slong prec)}

void \texttt{arb\_chebyshev\_u\_ui (arb\_t a, ulong n, const arb\_t x, slong prec)}

Evaluates the Chebyshev polynomial of the first kind $a = T_n(x)$ or the Chebyshev polynomial of the second kind $a = U_n(x)$.

void \texttt{arb\_chebyshev\_t2\_ui (arb\_t a, arb\_t b, ulong n, const arb\_t x, slong prec)}
3.3. arb.h – real numbers represented as floating-point balls

### 3.3.22 Internals for computing elementary functions

- Function: `void arb_chebyshev_u2_ui (arb_t a, arb_t b, ulong n, const arb_t x, slong prec)`
  - Simultaneously evaluates \( a = T_n(x), b = T_{n-1}(x) \) or \( a = U_n(x), b = U_{n-1}(x) \). Aliasing between \( a, b \) and \( x \) is not permitted.

- Function: `void arb_bell_sum_bsplit (arb_t res, const fmpz_t n, const fmpz_t a, const fmpz_t b, const fmpz_t mmag, slong prec)`
  - Helper functions for Bell numbers, evaluating the sum \( \sum_{k=0}^{b-1} k^n/k! \). If \( mmag \) is non-NULL, it may be used to indicate that the target error tolerance should be \( 2^{mmag-prec} \).

- Function: `void arb_bell_sum_taylor (arb_t res, const fmpz_t n, const fmpz_t a, const fmpz_t b, const fmpz_t_mmag, slong prec)`
  - Computes an approximation of \( y = \sum_{k=0}^{N-1} x^{2k+1}/(2k + 1) \) (if \( alternating \) is 0) or \( y = \sum_{k=0}^{N-1} (-1)^k x^{2k+1}/(2k + 1) \) (if \( alternating \) is 1). Used internally for computing arctangents and logarithms. The naive version uses the forward recurrence, and the rs version uses a division-avoiding rectangular splitting scheme.
  - Requires \( N \leq 255, 0 \leq x \leq 1/16, \) and \( x_n \) positive. The input \( x \) and output \( y \) are fixed-point numbers with \( x_n \) fractional limbs. A bound for the ulp error is written to \( error \).

- Function: `void arb_exp_taylor_naive (mp_ptr y, mp_limb_t * error, mp_srcptr x, mp_size_t xn, ulong N)`
  - Computes an approximation of \( y = \sum_{k=0}^{N-1} x^k/k! \). Used internally for computing exponentials. The naive version uses the forward recurrence, and the rs version uses a division-avoiding rectangular splitting scheme.
  - Requires \( N \leq 287, 0 \leq x \leq 1/16, \) and \( x_n \) positive. The input \( x \) is a fixed-point number with \( x_n \) fractional limbs, and the output \( y \) is a fixed-point number with \( x_n \) fractional limbs plus one extra limb for the integer part of the result.
  - A bound for the ulp error is written to \( error \).

- Function: `void arb_sin_cos_taylor_naive (mp_ptr ysin, mp_ptr ycos, mp_limb_t * error, mp_srcptr x, mp_size_t xn, ulong N)`
  - Computes approximations of \( y_s = \sum_{k=0}^{N-1} (-1)^k x^{2k+1}/(2k + 1)! \) and \( y_c = \sum_{k=0}^{N-1} (-1)^k x^{2k}/(2k)! \). Used internally for computing sines and cosines. The naive version uses the forward recurrence, and the rs version uses a division-avoiding rectangular splitting scheme.
  - Requires \( N \leq 143, 0 \leq x \leq 1/16, \) and \( x_n \) positive. The input \( x \) and outputs \( ysin, ycos \) are fixed-point numbers with \( x_n \) fractional limbs. A bound for the ulp error is written to \( error \).
  - If \( sinonly \) is 1, only the sine is computed; if \( sinonly \) is 0 both the sine and cosine are computed. To compute sin and cos, \( alternating \) should be 1. If \( alternating \) is 0, the hyperbolic sine is computed (this is currently only intended to be used together with \( sinonly \)).
int _arb_get_mpn_fixed_mod_log2 (mp_ptr w, fmpz_t q, mp_limb_t * error, const arf_t x, mp_size_t wn)
Attempts to write \( w = x - q \log(2) \) with \( 0 \leq w < \log(2) \), where \( w \) is a fixed-point number with \( wn \) limbs and ulp error \( \text{error} \). Returns success.

int _arb_get_mpn_fixed_mod_pi4 (mp_ptr w, fmpz_t q, int * octant, mp_limb_t * error, const arf_t x, mp_size_t wn)
Attempts to write \( w = |x| - q\pi/4 \) with \( 0 \leq w < \pi/4 \), where \( w \) is a fixed-point number with \( wn \) limbs and ulp error \( \text{error} \). Returns success.

The value of \( q \mod 8 \) is written to \( \text{octant} \). The output variable \( q \) can be NULL, in which case the full value of \( q \) is not stored.

void _arb_atan_sum_bs_powtab (arf_t z, const arf_t x, slong prec)

Computes the arctangent of \( x \). Initially, the argument-halving formula
\[
\frac{x}{1 + \sqrt{1 + x^2}}
\]
is applied up to 8 times to get a small argument. Then a version of the bit-burst algorithm is used. The functional equation
\[
\frac{x}{1 + \sqrt{1 + x^2}} = \arctan(p/q) + \arctan(w), \quad w = \frac{qx - p}{px + q}, \quad p = \lfloor qx \rfloor
\]
is applied repeatedly instead of integrating a differential equation for the arctangent, as this appears to be more efficient.

### 3.3.23 Vector functions

- **void** `_arb_vec_zero`(arb_ptr vec, slong n)
  
  Sets all entries in vec to zero.
- **int** `_arb_vec_is_zero`(arb_srcptr vec, slong len)
  
  Returns nonzero iff all entries in vec are zero.
- **int** `_arb_vec_is_finite`(arb_srcptr x, slong len)
  
  Returns nonzero iff all entries in x certainly are finite.
- **void** `_arb_vec_set`(arb_ptr res, arb_srcptr vec, slong len)
  
  Sets res to a copy of vec.
- **void** `_arb_vec_set_round`(arb_ptr res, arb_srcptr vec, slong len, slong prec)
  
  Sets res to a copy of vec, rounding each entry to prec bits.
- **void** `_arb_vec_swap`(arb_ptr vec1, arb_ptr vec2, slong len)
  
  Swaps the entries of vec1 and vec2.
- **void** `_arb_vec_neg`(arb_ptr B, arb_srcptr A, slong n)
- **void** `_arb_vec_sub`(arb_ptr C, arb_srcptr A, arb_srcptr B, slong n, slong prec)
- **void** `_arb_vec_add`(arb_ptr C, arb_srcptr A, arb_srcptr B, slong n, slong prec)
- **void** `_arb_vec_scalar_mul`(arb_ptr res, arb_srcptr vec, slong len, const arb_t c, slong prec)
- **void** `_arb_vec_scalar_div`(arb_ptr res, arb_srcptr vec, slong len, const arb_t c, slong prec)
- **void** `_arb_vec_scalar_mul_fmpz`(arb_ptr res, arb_srcptr vec, slong len, const fmpz_t c, slong prec)
- **void** `_arb_vec_scalar_mul_2exp_si`(arb_ptr res, arb_srcptr src, slong len, slong c)
- **void** `_arb_vec_scalar_addmul`(arb_ptr res, arb_srcptr vec, slong len, const arb_t c, slong prec)
  
  Performs the respective scalar operation elementwise.
- **void** `_arb_vec_dot`(arb_t res, arb_srcptr vec1, arb_srcptr vec2, slong len2, slong prec)
  
  Sets res to the dot product of vec1 and vec2.
- **void** `_arb_vec_norm`(arb_t res, arb_srcptr vec, slong len, slong prec)
  
  Sets res to the dot product of vec with itself.
- **void** `_arb_vec_get_mag`(mag_t bound, arb_srcptr vec, slong len, slong prec)
  
  Sets bound to an upper bound for the entries in vec.
- **ulong** `_arb_vec_bits`(arb_srcptr x, slong len)
  
  Returns the maximum of arb_bits() for all entries in vec.
- **void** `_arb_vec_set_powers`(arb_ptr xs, const arb_t x, slong len, const arb_t c, slong prec)
  
  Sets xs to the powers 1, x, x^2, ..., x^{len-1}.
- **void** `_arb_vec_add_error_arf_vec`(arb_ptr res, arf_srcptr err, slong len)
- **void** `_arb_vec_add_error_mag_vec`(arb_ptr res, mag_srcptr err, slong len)
  
  Adds the magnitude of each entry in err to the radius of the corresponding entry in res.
- **void** `_arb_vec_indeterminate`(arb_ptr vec, slong len)
  
  Applies arb_indeterminate() elementwise.
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void _arb_vec_trim (arb_ptr res, arb_srcptr vec, slong len)
  Applies arb_trim() elementwise.

int _arb_vec_get_unique_fmpz_vec (fmpz * res, arb_srcptr vec, slong len)
  Calls arb_get_unique_fmpz() elementwise and returns nonzero if all entries can be rounded uniquely to integers. If any entry in vec cannot be rounded uniquely to an integer, returns zero.

3.4 arb_poly.h – polynomials over the real numbers

An arb_poly_t represents a polynomial over the real numbers, implemented as an array of coefficients of type arb_struct.

Most functions are provided in two versions: an underscore method which operates directly on pre-allocated arrays of coefficients and generally has some restrictions (such as requiring the lengths to be nonzero and not supporting aliasing of the input and output arrays), and a non-underscore method which performs automatic memory management and handles degenerate cases.

3.4.1 Types, macros and constants

arb_poly_struct

arb_poly_t
  Contains a pointer to an array of coefficients (coeffs), the used length (length), and the allocated size of the array (alloc).

  An arb_poly_t is defined as an array of length one of type arb_poly_struct, permitting an arb_poly_t to be passed by reference.

3.4.2 Memory management

void arb_poly_init (arb_poly_t poly)
  Initializes the polynomial for use, setting it to the zero polynomial.

void arb_poly_clear (arb_poly_t poly)
  Clears the polynomial, deallocating all coefficients and the coefficient array.

void arb_poly_fit_length (arb_poly_t poly, slong len)
  Makes sure that the coefficient array of the polynomial contains at least len initialized coefficients.

void _arb_poly_set_length (arb_poly_t poly, slong len)
  Directly changes the length of the polynomial, without allocating or deallocating coefficients. The value should not exceed the allocation length.

void _arb_poly_normalise (arb_poly_t poly)
  Strips any trailing coefficients which are identical to zero.

3.4.3 Basic manipulation

void arb_poly_zero (arb_poly_t poly)

void arb_poly_one (arb_poly_t poly)
  Sets poly to the constant 0 respectively 1.

void arb_poly_set (arb_poly_t dest, const arb_poly_t src)
  Sets dest to a copy of src.
void \texttt{arb\_poly\_set\_round} \((\texttt{arb\_poly\_t} \texttt{dest}, \texttt{const arb\_poly\_t src, slong prec})\)
Sets \texttt{dest} to a copy of \texttt{src}, rounded to \texttt{prec} bits.

void \texttt{arb\_poly\_set\_coeff\_si} \((\texttt{arb\_poly\_t poly, slong n, slong c})\)
Sets the coefficient with index \texttt{n} in \texttt{poly} to the value \texttt{c}. We require that \texttt{n} is nonnegative.

void \texttt{arb\_poly\_set\_coeff\_arb} \((\texttt{arb\_poly\_t poly, slong n, const arb\_t c})\)
Sets \texttt{v} to the value of the coefficient with index \texttt{n} in \texttt{poly}. We require that \texttt{n} is nonnegative.

void \texttt{arb\_poly\_get\_coeff\_ptr} \((\texttt{poly, n})\)
Given \texttt{n} \geq 0, returns a pointer to coefficient \texttt{n} of \texttt{poly}, or \texttt{NULL} if \texttt{n} exceeds the length of \texttt{poly}.

void \texttt{arb\_poly\_shift\_right} \((\texttt{arb\_ptr res, arb\_srcptr poly, slong len, slong n})\)
Sets \texttt{res} to \texttt{poly} divided by \(x^n\), throwing away the lower coefficients. We require that \texttt{n} is nonnegative.

void \texttt{arb\_poly\_shift\_left} \((\texttt{arb\_ptr res, arb\_srcptr poly, slong len, slong n})\)
Sets \texttt{res} to \texttt{poly} multiplied by \(x^n\). We require that \texttt{n} is nonnegative.

void \texttt{arb\_poly\_truncate} \((\texttt{arb\_poly\_t poly, slong n})\)
Truncates \texttt{poly} to have length at most \texttt{n}, i.e. degree strictly smaller than \texttt{n}.

\texttt{slong} \texttt{arb\_poly\_length} \((\texttt{const arb\_poly\_t poly})\)
Returns the length of \texttt{poly}, i.e. zero if \texttt{poly} is identically zero, and otherwise one more than the index of the highest term that is not identically zero.

\texttt{slong} \texttt{arb\_poly\_degree} \((\texttt{const arb\_poly\_t poly})\)
Returns the degree of \texttt{poly}, defined as one less than its length. Note that if one or several leading coefficients are balls containing zero, this value can be larger than the true degree of the exact polynomial represented by \texttt{poly}, so the return value of this function is effectively an upper bound.

### 3.4.4 Conversions

void \texttt{arb\_poly\_set\_fmpz\_poly} \((\texttt{arb\_poly\_t poly, const fmpz\_poly\_t src, slong prec})\)
void \texttt{arb\_poly\_set\_fmpq\_poly} \((\texttt{arb\_poly\_t poly, const fmpq\_poly\_t src, slong prec})\)
void \texttt{arb\_poly\_set\_si} \((\texttt{arb\_poly\_t poly, slong src})\)
Sets \texttt{poly} to \texttt{src}, rounding the coefficients to \texttt{prec} bits.

### 3.4.5 Input and output

void \texttt{arb\_poly\_printd} \((\texttt{const arb\_poly\_t poly, slong digits})\)
Prints the polynomial as an array of coefficients, printing each coefficient using \texttt{arb\_printd}.

### 3.4.6 Random generation

void \texttt{arb\_poly\_randtest} \((\texttt{arb\_poly\_t poly, flint\_rand\_t state, slong len, slong prec, slong mag\_bits})\)
Creates a random polynomial with length at most \texttt{len}.
3.4.7 Comparisons

```c
int arb_poly_contains (const arb_poly_t poly1, const arb_poly_t poly2)
int arb_poly_contains_fmpz_poly (const arb_poly_t poly1, const fmpz_poly_t poly2)
int arb_poly_contains_fmpq_poly (const arb_poly_t poly1, const fmpq_poly_t poly2)
```

Returns nonzero iff `poly1` contains `poly2`.

```c
int arb_poly_equal (const arb_poly_t A, const arb_poly_t B)
```

Returns nonzero iff `A` and `B` are equal as polynomial balls, i.e. all coefficients have equal midpoint and radius.

```c
int _arb_poly_overlaps (arb_srcptr poly1, slong len1, arb_srcptr poly2, slong len2)
int arb_poly_overlaps (const arb_poly_t poly1, const arb_poly_t poly2)
```

Returns nonzero iff `poly1` overlaps with `poly2`. The underscore function requires that `len1` is at least as large as `len2`.

```c
int arb_poly_get_unique_fmpz_poly (fmpz_poly_t z, const arb_poly_t x)
```

If `x` contains a unique integer polynomial, sets `z` to that value and returns nonzero. Otherwise (if `x` represents no integers or more than one integer), returns zero, possibly partially modifying `z`.

3.4.8 Bounds

```c
void _arb_poly_majorant (arb_ptr res, arb_srcptr poly, slong len, slong prec)
void arb_poly_majorant (arb_poly_t res, const arb_poly_t poly, slong prec)
```

Sets `res` to an exact real polynomial whose coefficients are upper bounds for the absolute values of the coefficients in `poly`, rounded to `prec` bits.

3.4.9 Arithmetic

```c
void _arb_poly_add (arb_ptr C, arb_srcptr A, slong lenA, arb_srcptr B, slong lenB, slong prec)
void arb_poly_add (arb_poly_t C, const arb_poly_t A, const arb_poly_t B, slong prec)
void arb_poly_add_si (arb_poly_t C, const arb_poly_t A, slong B, slong prec)
void _arb_poly_sub (arb_ptr C, arb_srcptr A, slong lenA, arb_srcptr B, slong lenB, slong prec)
void arb_poly_sub (arb_poly_t C, const arb_poly_t A, const arb_poly_t B, slong prec)
void arb_poly_sub (arb_poly_t C, const arb_poly_t A, slong B, slong prec)
void _arb_poly_neg (arb_poly_t C, const arb_poly_t A)
void arb_poly_neg (arb_poly_t C, const arb_poly_t A)
void arb_poly_scalar_mul_2exp_si (arb_poly_t C, const arb_poly_t A, slong c)
void _arb_poly_mullow_classical (arb_ptr C, arb_srcptr A, slong lenA, arb_srcptr B, slong lenB, slong n, slong prec)
void _arb_poly_mullow_block (arb_ptr C, arb_srcptr A, slong lenA, arb_srcptr B, slong lenB, slong n, slong prec)
```
void _arb_poly_mullow (arb_ptr C, arb_srcptr A, slong lenA, arb_srcptr B, slong lenB, slong n, slong prec)

Sets \( \{C, n\} \) to the product of \( \{A, \text{len}A\} \) and \( \{B, \text{len}B\} \), truncated to length \( n \). The output is not allowed to be aliased with either of the inputs. We require \( \text{len}A \geq \text{len}B \geq 0, n \geq 0, \text{len}A + \text{len}B - 1 \geq n \).

The classical version uses a plain loop. This has good numerical stability but gets slow for large \( n \).

The block version decomposes the product into several subproducts which are computed exactly over the integers.

It first attempts to find an integer \( c \) such that \( A(2^c x) \) and \( B(2^c x) \) have slowly varying coefficients, to reduce the number of blocks.

The scaling factor \( c \) is chosen in a quick, heuristic way by picking the first and last nonzero terms in each polynomial. If the indices in \( A \) are \( a_2, a_1 \) and the log-2 magnitudes are \( e_2, e_1 \), and the indices in \( B \) are \( b_2, b_1 \) with corresponding magnitudes \( f_2, f_1 \), then we compute \( c \) as the weighted arithmetic mean of the slopes, rounded to the nearest integer:

\[
c = \left\lfloor \frac{(e_2 - e_1) + (f_2 + f_1)}{(a_2 - a_1) + (b_2 - b_1)} + \frac{1}{2} \right\rfloor.
\]

This strategy is used because it is simple. It is not optimal in all cases, but will typically give good performance when multiplying two power series with a similar decay rate.

The default algorithm chooses the classical algorithm for short polynomials and the block algorithm for slong polynomials.

If the input pointers are identical (and the lengths are the same), they are assumed to represent the same polynomial, and its square is computed.

void arb_poly_mullow_classical (arb_poly_t C, const arb_poly_t A, const arb_poly_t B, slong n, slong prec)

void arb_poly_mullow_ztrunc (arb_poly_t C, const arb_poly_t A, const arb_poly_t B, slong n, slong prec)

void arb_poly_mullow_block (arb_poly_t C, const arb_poly_t A, const arb_poly_t B, slong n, slong prec)

void arb_poly_mullow (arb_poly_t C, const arb_poly_t A, const arb_poly_t B, slong n, slong prec)

Sets \( C \) to the product of \( A \) and \( B \), truncated to length \( n \). If the same variable is passed for \( A \) and \( B \), sets \( C \) to the square of \( A \) truncated to length \( n \).

void _arb_poly_mul (arb_ptr C, arb_srcptr A, slong lenA, arb_srcptr B, slong lenB, slong prec)

Sets \( \{C, \text{len}A + \text{len}B - 1\} \) to the product of \( \{A, \text{len}A\} \) and \( \{B, \text{len}B\} \). The output is not allowed to be aliased with either of the inputs. We require \( \text{len}A \geq \text{len}B > 0 \). This function is implemented as a simple wrapper for _arb_poly_mullow().

If the input pointers are identical (and the lengths are the same), they are assumed to represent the same polynomial, and its square is computed.

void arb_poly_mul (arb_poly_t C, const arb_poly_t A, const arb_poly_t B, slong prec)

Sets \( C \) to the product of \( A \) and \( B \). If the same variable is passed for \( A \) and \( B \), sets \( C \) to the square of \( A \).

void _arb_poly_inv_series (arb_ptr Q, arb_srcptr A, slong Alen, slong len, slong prec)

Sets \( \{Q, \text{len}\} \) to the power series inverse of \( \{A, \text{Alen}\} \). Uses Newton iteration.

void arb_poly_inv_series (arb_poly_t Q, const arb_poly_t A, slong n, slong prec)

Sets \( Q \) to the power series inverse of \( A \), truncated to length \( n \).

void _arb_poly_div_series (arb_ptr Q, arb_srcptr A, slong Alen, arb_srcptr B, slong Blen, slong n, slong prec)

Sets \( \{Q, n\} \) to the power series quotient of \( \{A, \text{Alen}\} \) by \( \{B, \text{Blen}\} \). Uses Newton iteration followed by multiplication.
3.4.10 Composition

void _arb_poly_compose_horner (arb_ptr res, arb_srcptr poly1, slong len1, arb_srcptr poly2, slong len2, slong prec)

Sets res to the composition \( h(x) = f(g(x)) \) where \( f \) is given by poly1 and \( g \) is given by poly2, respectively using Horner’s rule, divide-and-conquer, and an automatic choice between the two algorithms. The underscore methods do not support aliasing of the output with either input polynomial.

void _arb_poly_compose_series_horner (arb_ptr res, arb_srcptr poly1, slong len1, arb_srcptr poly2, slong len2, slong n, slong prec)

Sets res to the power series composition \( h(x) = f(g(x)) \) truncated to order \( O(x^n) \) where \( f \) is given by poly1 and \( g \) is given by poly2, respectively using Horner’s rule, the Brent-Kung baby step-giant step algorithm, and an automatic choice between the two algorithms. We require that the constant term in \( g(x) \) is exactly zero. The underscore methods do not support aliasing of the output with either input polynomial.
void \_arb\_poly\_revert\_series\_lagrange (arb\_ptr h, arb\_srcptr f, slong flen, slong n, slong prec)

void \_arb\_poly\_revert\_series\_lagrange (arb\_poly\_1 h, const arb\_poly\_1 f, slong n, slong prec)

void \_arb\_poly\_revert\_series\_newton (arb\_ptr h, arb\_srcptr f, slong flen, slong n, slong prec)

void \_arb\_poly\_revert\_series\_newton (arb\_poly\_1 h, const arb\_poly\_1 f, slong n, slong prec)

void \_arb\_poly\_revert\_series\_lagrange\_fast (arb\_ptr h, arb\_srcptr f, slong flen, slong n, slong prec)

void \_arb\_poly\_revert\_series\_lagrange\_fast (arb\_poly\_1 h, const arb\_poly\_1 f, slong n, slong prec)

void \_arb\_poly\_revert\_series (arb\_ptr h, arb\_srcptr f, slong flen, slong n, slong prec)

void \_arb\_poly\_revert\_series (arb\_poly\_1 h, const arb\_poly\_1 f, slong n, slong prec)

Sets \( h \) to the power series reversion of \( f \), i.e. the expansion of the compositional inverse function \( f^{-1}(x) \), truncated to order \( O(x^n) \), using respectively Lagrange inversion, Newton iteration, fast Lagrange inversion, and a default algorithm choice.

We require that the constant term in \( f \) is exactly zero and that the linear term is nonzero. The underscore methods assume that \( flen \) is at least 2, and do not support aliasing.

### 3.4.11 Evaluation

void \_arb\_poly\_evaluate\_horner (arb\_t y, arb\_srcptr f, slong len, const arb\_t x, slong prec)

void \_arb\_poly\_evaluate\_horner (arb\_t y, const arb\_poly\_1 f, const arb\_t x, slong prec)

void \_arb\_poly\_evaluate\_rectangular (arb\_t y, arb\_srcptr f, slong len, const arb\_t x, slong prec)

void \_arb\_poly\_evaluate\_rectangular (arb\_t y, const arb\_poly\_1 f, const arb\_t x, slong prec)

void \_arb\_poly\_evaluate (arb\_t y, arb\_srcptr f, slong len, const arb\_t x, slong prec)

void \_arb\_poly\_evaluate (arb\_t y, const arb\_poly\_1 f, const arb\_t x, slong prec)

Sets \( y = f(x) \), evaluated respectively using Horner’s rule, rectangular splitting, and an automatic algorithm choice.

void \_arb\_poly\_evaluate\_acb\_horner (acb\_t y, arb\_srcptr f, slong len, const acb\_t x, slong prec)

void \_arb\_poly\_evaluate\_acb\_horner (acb\_t y, const arb\_poly\_1 f, const acb\_t x, slong prec)

void \_arb\_poly\_evaluate\_acb\_rectangular (acb\_t y, arb\_srcptr f, slong len, const acb\_t x, slong prec)

void \_arb\_poly\_evaluate\_acb\_rectangular (acb\_t y, const arb\_poly\_1 f, const acb\_t x, slong prec)

void \_arb\_poly\_evaluate\_acb (acb\_t y, arb\_srcptr f, slong len, const acb\_t x, slong prec)

void \_arb\_poly\_evaluate\_acb (acb\_t y, const arb\_poly\_1 f, const acb\_t x, slong prec)

Sets \( y = f(x) \) where \( x \) is a complex number, evaluating the polynomial respectively using Horner’s rule, rectangular splitting, and an automatic algorithm choice.

void \_arb\_poly\_evaluate2\_horner (arb\_t y, arb\_t z, arb\_srcptr f, slong len, const arb\_t x, slong prec)

void \_arb\_poly\_evaluate2\_horner (arb\_t y, arb\_t z, const arb\_poly\_1 f, const arb\_t x, slong prec)

void \_arb\_poly\_evaluate2\_rectangular (arb\_t y, arb\_t z, arb\_srcptr f, slong len, const arb\_t x, slong prec)

void \_arb\_poly\_evaluate2\_rectangular (arb\_t y, arb\_t z, const arb\_poly\_1 f, const arb\_t x, slong prec)

void \_arb\_poly\_evaluate2 (arb\_t y, arb\_t z, arb\_srcptr f, slong len, const arb\_t x, slong prec)
void **arb_poly_evaluate2**(arb_t y, arb_t z, const arb_poly_t f, const arb_t x, slong prec)
Sets \( y = f(x), z = f'(x) \), evaluated respectively using Horner’s rule, rectangular splitting, and an automatic algorithm choice.

When Horner’s rule is used, the only advantage of evaluating the function and its derivative simultaneously is that one does not have to generate the derivative polynomial explicitly. With the rectangular splitting algorithm, the powers can be reused, making simultaneous evaluation slightly faster.

void **_arb_poly_evaluate2_acb_horner**(acb_t y, acb_t z, arb_srcptr f, slong len, const acb_t x, slong prec)
void **arb_poly_evaluate2_acb_horner**(acb_t y, acb_t z, const arb_poly_t f, const acb_t x, slong prec)

void **_arb_poly_evaluate2_acb_rectangular**(acb_t y, acb_t z, arb_srcptr f, slong len, const arb_t x, slong prec)
void **arb_poly_evaluate2_acb_rectangular**(acb_t y, acb_t z, const arb_poly_t f, const arb_t x, slong prec)

void **_arb_poly_evaluate2_acb**(acb_t y, acb_t z, arb_srcptr f, slong len, const arb_t x, slong prec)
void **arb_poly_evaluate2_acb**(acb_t y, acb_t z, const arb_poly_t f, const arb_t x, slong prec)

3.4.12 Product trees

void **_arb_poly_product_roots**(arb_ptr poly, arb_srcptr xs, slong n, slong prec)
void **arb_poly_product_roots**(arb_poly_t poly, arb_srcptr xs, slong n, slong prec)

Generates the polynomial \((x - x_0)(x - x_1) \cdots (x - x_{n-1})\).

**arb_ptr** **_arb_poly_tree_alloc**(slong len)
Returns an initialized data structure capable of representing a remainder tree (product tree) of len roots.

void **_arb_poly_tree_free**(arb_ptr* tree, slong len)
Deallocates a tree structure as allocated using _arb_poly_tree_alloc().

void **_arb_poly_tree_build**(arb_ptr* tree, arb_srcptr roots, slong len, slong prec)
Constructs a product tree from a given array of len roots. The tree structure must be pre-allocated to the specified length using _arb_poly_tree_alloc().

3.4.13 Multipoint evaluation

void **_arb_poly_evaluate_vec_iter**(arb_ptr ys, arb_srcptr poly, slong plen, arb_srcptr xs, slong n, slong prec)
void **arb_poly_evaluate_vec_iter**(arb_ptr ys, const arb_poly_t poly, arb_srcptr xs, slong n, slong prec)

Evaluates the polynomial simultaneously at n given points, calling _arb_poly_evaluate() repeatedly.

void **_arb_poly_evaluate_vec_fast_precomp**(arb_ptr vs, arb_srcptr poly, slong plen, arb_ptr* tree, slong len, slong prec)
void **arb_poly_evaluate_vec_fast**(arb_ptr vs, arb_srcptr poly, slong plen, arb_srcptr xs, slong n, slong prec)

void **arb_poly_evaluate_vec_fast**(arb_ptr vs, const arb_poly_t poly, arb_srcptr xs, slong n, slong prec)

Evaluates the polynomial simultaneously at n given points, using fast multipoint evaluation.
3.4.14 Interpolation

void _arb_poly_interpolate_newton (arb_ptr poly, arb_srcptr xs, arb_srcptr ys, slong n, slong prec)
void arb_poly_interpolate_newton (arb_poly_t poly, arb_srcptr xs, arb_srcptr ys, slong n, slong prec)

Recovers the unique polynomial of length at most \( n \) that interpolates the given \( x \) and \( y \) values. This implementation first interpolates in the Newton basis and then converts back to the monomial basis.

void _arb_poly_interpolate_barycentric (arb_ptr poly, arb_srcptr xs, arb_srcptr ys, slong n, slong prec)
void arb_poly_interpolate_barycentric (arb_poly_t poly, arb_srcptr xs, arb_srcptr ys, slong n, slong prec)

Recovers the unique polynomial of length at most \( n \) that interpolates the given \( x \) and \( y \) values. This implementation uses the barycentric form of Lagrange interpolation.

void _arb_poly_interpolation_weights (arb_ptr w, arb_ptr * tree, slong len, slong prec)
void arb_poly_interpolate_fast_precomp (arb_ptr poly, arb_srcptr ys, arb_ptr * tree, arb_srcptr weights, slong len, slong prec)
void _arb_poly_interpolate_fast (arb_ptr poly, arb_srcptr xs, arb_srcptr ys, slong n, slong prec)
void arb_poly_interpolate_fast (arb_poly_t poly, arb_srcptr xs, arb_srcptr ys, slong n, slong prec)

Recovers the unique polynomial of length at most \( n \) that interpolates the given \( x \) and \( y \) values, using fast Lagrange interpolation. The precomp function takes a precomputed product tree over the \( x \) values and a vector of interpolation weights as additional inputs.

3.4.15 Differentiation

void _arb_poly_derivative (arb_ptr res, arb_srcptr poly, slong len, slong prec)
Sets \( \{res, len - 1\} \) to the derivative of \( \{poly, len\} \). Allows aliasing of the input and output.
void arb_poly_derivative (arb_poly_t res, const arb_poly_t poly, slong prec)
Sets \( res \) to the derivative of \( poly \).

void _arb_poly_integral (arb_ptr res, arb_srcptr poly, slong len, slong prec)
Sets \( \{res, len\} \) to the integral of \( \{poly, len - 1\} \). Allows aliasing of the input and output.
void arb_poly_integral (arb_poly_t res, const arb_poly_t poly, slong prec)
Sets \( res \) to the integral of \( poly \).

3.4.16 Transforms

void _arb_poly_borel_transform (arb_ptr res, arb_srcptr poly, slong len, slong prec)
void arb_poly_borel_transform (arb_poly_t res, const arb_poly_t poly, slong prec)
Computes the Borel transform of the input polynomial, mapping \( \sum_k a_k x^k \) to \( \sum_k (a_k / k!) x^k \). The underscore method allows aliasing.

void _arb_poly_inv_borel_transform (arb_ptr res, arb_srcptr poly, slong len, slong prec)
void arb_poly_inv_borel_transform (arb_poly_t res, const arb_poly_t poly, slong prec)
Computes the inverse Borel transform of the input polynomial, mapping \( \sum_k a_k x^k \) to \( \sum_k a_k k! x^k \). The underscore method allows aliasing.

void _arb_poly_binomial_transform_basecase (arb_ptr b, arb_srcptr a, slong alen, slong len, slong prec)
3.4.17 Powers and elementary functions

void _arb_poly_pow_ui_trunc_binexp (arb_ptr res, arb_srcptr f, slong flen, ulong exp, slong len, slong prec)
Sets {res, len} to \( f(x)^\text{exp} \), truncated to length \( \text{len} \). Requires that \( \text{len} \) is no longer than the length of the power as computed without truncation (i.e. no zero-padding is performed). Does not support aliasing of the input and output, and requires that \( \text{flen} \) and \( \text{len} \) are positive. Uses binary exponentiation.

void _arb_poly_pow_ui_trunc_binexp (arb_poly_t res, const arb_poly_t poly, ulong exp, slong len, slong prec)
Sets \( \text{res} \) to \( \text{poly} \) raised to the power \( \text{exp} \), truncated to length \( \text{len} \). Uses binary exponentiation.

void _arb_poly_pow_ui (arb_ptr res, arb_srcptr f, slong flen, ulong exp, slong prec)
Sets \( \text{res} \) to \( f(x)^\text{exp} \). Does not support aliasing of the input and output, and requires that \( \text{flen} \) is positive.

void _arb_poly_pow_ui (arb_poly_t res, const arb_poly_t poly, ulong exp, slong prec)
Sets \( \text{res} \) to \( \text{poly} \) raised to the power \( \text{exp} \).

void _arb_poly_pow_series (arb_ptr h, arb_srcptr f, slong flen, arb_srcptr g, slong glen, slong len, slong prec)
Sets \( \{h, \text{len}\} \) to the power series \( f(x)^g(x) = \exp(g(x) \log f(x)) \), truncated to length \( \text{len} \). This function detects special cases such as \( g \) being an exact small integer or \( \pm 1/2 \), and computes such powers more efficiently. This function does not support aliasing of the output with either of the input operands. It requires that all lengths are positive, and assumes that \( \text{flen} \) and \( \text{glen} \) do not exceed \( \text{len} \).

void _arb_poly_pow_series (arb_poly_t h, const arb_poly_t f, const arb_poly_t g, slong len, slong prec)
Sets \( h \) to the power series \( f(x)^g(x) = \exp(g(x) \log f(x)) \), truncated to length \( \text{len} \). This function detects special cases such as \( g \) being an exact small integer or \( \pm 1/2 \), and computes such powers more efficiently.

void _arb_poly_pow_series (arb_ptr h, arb_srcptr f, slong flen, const arb_t g, slong len, slong prec)
Sets \( \{h, \text{len}\} \) to the power series \( f(x)^g = \exp(g \log f(x)) \), truncated to length \( \text{len} \). This function detects special
cases such as $g$ being an exact small integer or $\pm 1/2$, and computes such powers more efficiently. This function does not support aliasing of the output with either of the input operands. It requires that all lengths are positive, and assumes that $flen$ does not exceed $len$.

```c
void arb_poly_pow_arb_series (arb_poly_t res, const arb_poly_t f, const arb_poly_t g, slong flen, slong prec)
```

Sets $h$ to the power series $f(x)^g = \exp(g \log f(x))$ truncated to length $len$.

```c
void _arb_poly_sqrt_series (arb_ptr g, arb_srcptr h, slong hlen, slong n, slong prec)
```

Sets $g$ to the power series square root of $h$, truncated to length $n$. Uses division-free Newton iteration for the reciprocal square root, followed by a multiplication.

The underscore method does not support aliasing of the input and output arrays. It requires that $hlen$ and $n$ are greater than zero.

```c
void _arb_poly_rsqrt_series (arb_ptr g, arb_srcptr h, slong hlen, slong n, slong prec)
```

Sets $g$ to the reciprocal power series square root of $h$, truncated to length $n$. Uses division-free Newton iteration.

The underscore method does not support aliasing of the input and output arrays. It requires that $hlen$ and $n$ are greater than zero.

```c
void _arb_poly_log_series (arb_ptr res, arb_srcptr f, slong flen, slong n, slong prec)
```

Sets $res$ to the power series logarithm of $f$, truncated to length $n$. Uses the formula $\log(f(x)) = \int f'(x)/f(x)dx$, adding the logarithm of the constant term in $f$ as the constant of integration.

The underscore method supports aliasing of the input and output arrays. It requires that $flen$ and $n$ are greater than zero.

```c
void _arb_poly_atan_series (arb_ptr res, arb_srcptr f, slong flen, slong n, slong prec)
```

Sets $res$ respectively to the power series inverse tangent, inverse sine and inverse cosine of $f$, truncated to length $n$.

Uses the formulas

$$
\tan^{-1}(f(x)) = \int \frac{f'(x)}{1 + f(x)^2}dx,
$$

$$
\sin^{-1}(f(x)) = \int \frac{f'(x)}{(1 - f(x)^2)^{1/2}}dx,
$$

$$
\cos^{-1}(f(x)) = -\int \frac{f'(x)}{(1 - f(x)^2)^{1/2}}dx,
$$

adding the inverse function of the constant term in $f$ as the constant of integration.

The underscore method supports aliasing of the input and output arrays. They require that $flen$ and $n$ are greater than zero.

```c
void _arb_poly_exp_series_basecase (arb_ptr f, arb_srcptr h, slong hlen, slong n, slong prec)
```

To compute $\exp(h)$, which is more efficient when $h$ is an integer.

```c
void arb_poly_exp_series_basecase (arb_poly_t f, const arb_poly_t h, slong n, slong prec)
```

To compute $\exp(h)$.
void \_arb\_poly\_exp\_series (arb\_ptr f, arb\_srcptr h, slong hlen, slong n, slong prec)  

void arb\_poly\_exp\_series (arb\_poly\_t f, const arb\_poly\_t h, slong n, slong prec)  

Sets $f$ to the power series exponential of $h$, truncated to length $n$.

The basecase version uses a simple recurrence for the coefficients, requiring $O(nm)$ operations where $m$ is the length of $h$.

The main implementation uses Newton iteration, starting from a small number of terms given by the basecase algorithm. The complexity is $O(M(n))$. Redundant operations in the Newton iteration are avoided by using the scheme described in [HZ2004].

The underscore methods support aliasing and allow the input to be shorter than the output, but require the lengths to be nonzero.

void \_arb\_poly\_sin\_cos\_series\_basecase (arb\_ptr s, arb\_ptr c, arb\_srcptr h, slong hlen, slong n, slong prec, int times\_pi)  

void arb\_poly\_sin\_cos\_series\_basecase (arb\_poly\_t s, arb\_poly\_t c, const arb\_poly\_t h, slong n, slong prec, int times\_pi)  

void \_arb\_poly\_sin\_cos\_series\_tangent (arb\_ptr s, arb\_ptr c, arb\_srcptr h, slong hlen, slong n, slong prec, int times\_pi)  

void arb\_poly\_sin\_cos\_series\_tangent (arb\_poly\_t s, arb\_poly\_t c, const arb\_poly\_t h, slong n, slong prec, int times\_pi)  

void \_arb\_poly\_sin\_cos\_series (arb\_ptr s, arb\_ptr c, arb\_srcptr h, slong hlen, slong n, slong prec)  

void arb\_poly\_sin\_cos\_series (arb\_poly\_t s, arb\_poly\_t c, const arb\_poly\_t h, slong n, slong prec)  

Sets $s$ and $c$ to the power series sine and cosine of $h$, computed simultaneously.

The basecase version uses a simple recurrence for the coefficients, requiring $O(nm)$ operations where $m$ is the length of $h$.

The tangent version uses the tangent half-angle formulas to compute the sine and cosine via \_arb\_poly\_tan\_series(). This requires $O(M(n))$ operations. When $h = h_0 + h_1$ where the constant term $h_0$ is nonzero, the evaluation is done as $\sin(h_0 + h_1) = \cos(h_0)\sin(h_1) + \sin(h_0)\cos(h_1)$, $\cos(h_0 + h_1) = \cos(h_0)\cos(h_1) - \sin(h_0)\sin(h_1)$, to improve accuracy and avoid dividing by zero at the poles of the tangent function.

The default version automatically selects between the basecase and tangent algorithms depending on the input.

The basecase and tangent versions take a flag times\_pi specifying that the input is to be multiplied by $\pi$.

The underscore methods support aliasing and require the lengths to be nonzero.

void \_arb\_poly\_sin\_series (arb\_ptr s, arb\_srcptr h, slong hlen, slong n, slong prec)  

void arb\_poly\_sin\_series (arb\_poly\_t s, const arb\_poly\_t h, slong n, slong prec)  

void \_arb\_poly\_cos\_series (arb\_ptr c, arb\_srcptr h, slong hlen, slong n, slong prec)  

void arb\_poly\_cos\_series (arb\_poly\_t c, const arb\_poly\_t h, slong n, slong prec)  

Respectively evaluates the power series sine or cosine. These functions simply wrap \_arb\_poly\_sin\_cos\_series(). The underscore methods support aliasing and require the lengths to be nonzero.

void \_arb\_poly\_tan\_series (arb\_ptr g, arb\_srcptr h, slong hlen, slong len, slong prec)  

void arb\_poly\_tan\_series (arb\_poly\_t g, const arb\_poly\_t h, slong n, slong prec)  

Sets $g$ to the power series tangent of $h$.

For small $n$ takes the quotient of the sine and cosine as computed using the basecase algorithm. For large $n$, uses Newton iteration to invert the inverse tangent series. The complexity is $O(M(n))$.  

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The underscore version does not support aliasing, and requires the lengths to be nonzero.

```c
void _arb_poly_sin_cos_pi_series (arb_ptr s, arb_ptr c, arb_srcptr h, slong hlen, slong n, slong prec)
void _arb_poly_sin_pi_series (arb_ptr s, arb_srcptr h, slong hlen, slong n, slong prec)
void _arb_poly_cos_pi_series (arb_ptr c, arb_srcptr h, slong hlen, slong n, slong prec)
void _arb_poly_cot_pi_series (arb_ptr c, arb_srcptr h, slong hlen, slong n, slong prec)
void _arb_poly_sin_cos_pi_series (arb_poly_t s, arb_poly_t c, const arb_poly_t h, slong n, slong prec)
void _arb_poly_sin_pi_series (arb_poly_t s, const arb_poly_t h, slong n, slong prec)
void _arb_poly_cos_pi_series (arb_poly_t c, const arb_poly_t h, slong n, slong prec)
void _arb_poly_cot_pi_series (arb_poly_t c, const arb_poly_t h, slong n, slong prec)
void _arb_poly_rising_ui_series (arb_poly_t res, const arb_poly_t h, slong n, slong prec)
```

3.4.18 Gamma function and factorials

```c
void _arb_poly_gamma_series (arb_ptr res, arb_srcptr h, slong hlen, slong n, slong prec)
void _arb_poly_gamma_series (arb_poly_t res, const arb_poly_t h, slong n, slong prec)
void _arb_poly_rgamma_series (arb_poly_t res, const arb_poly_t h, slong n, slong prec)
void _arb_poly_lgamma_series (arb_poly_t res, const arb_poly_t h, slong n, slong prec)
void _arb_poly_lgamma_series (arb_poly_t res, arb_ptr h, slong n, slong prec)
void _arb_poly_rgamma_series (arb_ptr res, const arb_poly_t h, slong n, slong prec)
void _arb_poly_rgamma_series (arb_ptr res, arb_srcptr h, slong n, slong prec)
void _arb_poly_gamma_series (arb_poly_t res, const arb_poly_t h, slong n, slong prec)
void _arb_poly_rising_ui_series (arb_poly_t res, const arb_poly_t h, slong n, slong prec)
```

3.4.19 Zeta function

```c
void _arb_poly_zeta_series (arb_poly_t res, const arb_poly_t s, const arb_t a, int deflate, slong n, slong prec)
```

To evaluate the usual Riemann zeta function, set \( \alpha = 1 \).
If `deflate` is nonzero, evaluates \( \zeta(s, a) + 1/(1-s) \), which is well-defined as a limit when the constant term of \( s \) is 1. In particular, expanding \( \zeta(s, a) + 1/(1-s) \) with \( s = 1 + x \) gives the Stieltjes constants

\[
\sum_{k=0}^{n-1} \frac{(-1)^k}{k!} \gamma_k(a) x^k.
\]

If \( a = 1 \), this implementation uses the reflection formula if the midpoint of the constant term of \( s \) is negative.

```c
void _arb_poly_riemann_siegel_theta_series (arb_ptr res, arb_srcptr h, slong hlen, slong n, slong prec)

void arb_poly_riemann_siegel_theta_series (arb_poly_t res, const arb_poly_t h, slong n, slong prec)
```

Sets \( res \) to the series expansion of the Riemann-Siegel theta function

\[
\theta(h) = \arg \left( \Gamma \left( \frac{2ih + 1}{4} \right) \right) - \frac{\log \pi h}{2}
\]

where the argument of the gamma function is chosen continuously as the imaginary part of the log gamma function.

The underscore method does not support aliasing of the input and output arrays, and requires that the lengths are greater than zero.

```c
void _arb_poly_riemann_siegel_z_series (arb_ptr res, arb_srcptr h, slong hlen, slong n, slong prec)

void arb_poly_riemann_siegel_z_series (arb_poly_t res, const arb_poly_t h, slong n, slong prec)
```

Sets \( res \) to the series expansion of the Riemann-Siegel Z-function

\[
Z(h) = e^{i\theta(h)} \zeta(1/2 + ih).
\]

The zeros of the Z-function on the real line precisely correspond to the imaginary parts of the zeros of the Riemann zeta function on the critical line.

The underscore method supports aliasing of the input and output arrays, and requires that the lengths are greater than zero.

### 3.4.20 Root-finding

```c
void _arb_poly_root_bound_fujiwara (mag_t bound, arb_srcptr poly, slong len)

void arb_poly_root_bound_fujiwara (mag_t bound, arb_poly_t poly)
```

Sets \( bound \) to an upper bound for the magnitude of all the complex roots of \( poly \). Uses Fujiwara’s bound

\[
2 \max \left\{ \left| \frac{a_{n-1}}{a_n} \right|, \left| \frac{a_{n-2}}{a_n} \right|^{1/2}, \ldots, \left| \frac{a_1}{a_n} \right|^{1/(n-1)}, \left| \frac{a_0}{2a_n} \right|^{1/n} \right\}
\]

where \( a_0, \ldots, a_n \) are the coefficients of \( poly \).

```c
void _arb_poly_newton_convergence_factor (arf_t convergence_factor, arb_srcptr poly, slong len, const arb_t convergence_interval, slong prec)
```

Given an interval \( I \) specified by \( convergence_interval \), evaluates a bound for \( C = \sup_{t, u \in I} \frac{1}{2} \left| f''(t) \right| / \left| f'(u) \right| \), where \( f \) is the polynomial defined by the coefficients \( \{ poly, len \} \). The bound is obtained by evaluating \( f'(I) \) and \( f''(I) \) directly. If \( f \) has large coefficients, \( I \) must be extremely precise in order to get a finite factor.
Arb Documentation, Release 2.8.1

int _arb_poly_newton_step (arb_t xnew, arb_srcptr poly, slong len, const arb_t x, const arb_t convergence_interval, const arf_t convergence_factor, slong prec)

Performs a single step with Newton's method.

The input consists of the polynomial \( f \) specified by the coefficients \( \{poly, len\} \), an interval \( x = [m - r, m + r] \) known to contain a single root of \( f \), an interval \( I (\text{convergence_interval}) \) containing \( x \) with an associated bound \( (\text{convergence_factor}) \) for \( C = \sup_{t,u \in I} \frac{1}{2} \frac{|f''(t)|}{|f'(u)|} \), and a working precision \( prec \).

The Newton update consists of setting \( x' = [m' - r', m' + r'] \) where \( m' = m - \frac{f(m)}{f'(m)} \) and \( r' = Cr^2 \). The expression \( m - \frac{f(m)}{f'(m)} \) is evaluated using ball arithmetic at a working precision of \( prec \) bits, and the rounding error during this evaluation is accounted for in the output. We now check that \( x' \in I \) and \( m' < m \). If both conditions are satisfied, we set \( xnew \) to \( x' \) and return nonzero. If either condition fails, we set \( xnew \) to \( x \) and return zero, indicating that no progress was made.

void _arb_poly_newton_refine_root (arb_t r, arb_srcptr poly, slong len, const arb_t start, const arb_t convergence_interval, const arf_t convergence_factor, slong eval_extra_prec, slong prec)

Refines a precise estimate of a polynomial root to high precision by performing several Newton steps, using nearly optimally chosen doubling precision steps.

The inputs are defined as for \_arb_poly_newton_step, except for the precision parameters: \( prec \) is the target accuracy and \( eval_extra_prec \) is the estimated number of guard bits that need to be added to evaluate the polynomial accurately close to the root (typically, if the polynomial has large coefficients of alternating signs, this needs to be approximately the bit size of the coefficients).

3.4.21 Other special polynomials

void _arb_poly_swinnerton_dyer_ui (arb_ptr poly, ulong n, slong trunc, slong prec)
void arb_poly_swinnerton_dyer_ui (arb_poly_t poly, ulong n, slong prec)

Computes the Swinnerton-Dyer polynomial \( S_n \), which has degree \( 2^n \) and is the rational minimal polynomial of the sum of the square roots of the first \( n \) prime numbers.

If \( prec \) is set to zero, a precision is chosen automatically such that sizeof arb_poly_get_unique_fmpz_poly() should be successful. Otherwise a working precision of \( prec \) bits is used.

The underscore version accepts an additional \( trunc \) parameter. Even when computing a truncated polynomial, the array \( poly \) must have room for \( 2^n + 1 \) coefficients, used as temporary space.

3.5 arb_mat.h – matrices over the real numbers

An \( \text{arb_mat_t} \) represents a dense matrix over the real numbers, implemented as an array of entries of type \( \text{arb_struct} \).

The dimension (number of rows and columns) of a matrix is fixed at initialization, and the user must ensure that inputs and outputs to an operation have compatible dimensions. The number of rows or columns in a matrix can be zero.

3.5.1 Types, macros and constants

\text{arb_mat_struct}

\text{arb_mat_t}

Contains a pointer to a flat array of the entries (entries), an array of pointers to the start of each row (rows), and the number of rows (r) and columns (c).
An `arb_mat_t` is defined as an array of length one of type `arb_mat_struct`, permitting an `arb_mat_t` to be passed by reference.

`arb_mat_entry(mat, i, j)`
Macro giving a pointer to the entry at row `i` and column `j`.

`arb_mat_nrows(mat)`
Returns the number of rows of the matrix.

`arb_mat_ncols(mat)`
Returns the number of columns of the matrix.

### 3.5.2 Memory management

void `arb_mat_init(arb_mat_t mat, slong r, slong c)`
Initializes the matrix, setting it to the zero matrix with `r` rows and `c` columns.

void `arb_mat_clear(arb_mat_t mat)`
Clears the matrix, deallocating all entries.

### 3.5.3 Conversions

void `arb_mat_set(arb_mat_t dest, const arb_mat_t src)`

void `arb_mat_set_fmpz_mat(arb_mat_t dest, const fmpz_mat_t src)`

void `arb_mat_set_round_fmpz_mat(arb_mat_t dest, const fmpz_mat_t src, slong prec)`

void `arb_mat_set_fmpq_mat(arb_mat_t dest, const fmpq_mat_t src, slong prec)`
Sets `dest` to `src`. The operands must have identical dimensions.

### 3.5.4 Random generation

void `arb_mat_randtest(arb_mat_t mat, flint_rand_t state, slong prec, slong mag_bits)`
Sets `mat` to a random matrix with up to `prec` bits of precision and with exponents of width up to `mag_bits`.

### 3.5.5 Input and output

void `arb_mat_printd(const arb_mat_t mat, slong digits)`
Prints each entry in the matrix with the specified number of decimal digits.

### 3.5.6 Comparisons

int `arb_mat_equal(const arb_mat_t mat1, const arb_mat_t mat2)`
Returns nonzero iff the matrices have the same dimensions and identical entries.

int `arb_mat_overlaps(const arb_mat_t mat1, const arb_mat_t mat2)`
Returns nonzero iff the matrices have the same dimensions and each entry in `mat1` overlaps with the corresponding entry in `mat2`.

int `arb_mat_contains(const arb_mat_t mat1, const arb_mat_t mat2)`

int `arb_mat_contains_fmpz_mat(const arb_mat_t mat1, const fmpz_mat_t mat2)`
int \texttt{arb\_mat\_contains\_fmpq\_mat} (\texttt{const arb\_mat\_t mat1, const fmpq\_mat\_t mat2})

Returns nonzero iff the matrices have the same dimensions and each entry in \texttt{mat2} is contained in the corresponding entry in \texttt{mat1}.

int \texttt{arb\_mat\_eq} (\texttt{const arb\_mat\_t mat1, const arb\_mat\_t mat2})

Returns nonzero iff \texttt{mat1} and \texttt{mat2} certainly represent the same matrix.

int \texttt{arb\_mat\_ne} (\texttt{const arb\_mat\_t mat1, const arb\_mat\_t mat2})

Returns nonzero iff \texttt{mat1} and \texttt{mat2} certainly do not represent the same matrix.

3.5.7 Special matrices

void \texttt{arb\_mat\_zero} (\texttt{arb\_mat\_t mat})

Sets all entries in \texttt{mat} to zero.

void \texttt{arb\_mat\_one} (\texttt{arb\_mat\_t mat})

Sets the entries on the main diagonal to ones, and all other entries to zero.

3.5.8 Transpose

void \texttt{arb\_mat\_transpose} (\texttt{arb\_mat\_t dest, const arb\_mat\_t src})

Sets \texttt{dest} to the exact transpose \texttt{src}. The operands must have compatible dimensions. Aliasing is allowed.

3.5.9 Norms

void \texttt{arb\_mat\_bound\_inf\_norm} (\texttt{mag\_t b, const arb\_mat\_t A})

Sets \texttt{b} to an upper bound for the infinity norm (i.e. the largest absolute value row sum) of \texttt{A}.

3.5.10 Arithmetic

void \texttt{arb\_mat\_neg} (\texttt{arb\_mat\_t dest, const arb\_mat\_t src})

Sets \texttt{dest} to the exact negation of \texttt{src}. The operands must have the same dimensions.

void \texttt{arb\_mat\_add} (\texttt{arb\_mat\_t res, const arb\_mat\_t mat1, const arb\_mat\_t mat2, slong prec})

Sets \texttt{res} to the sum of \texttt{mat1} and \texttt{mat2}. The operands must have the same dimensions.

void \texttt{arb\_mat\_sub} (\texttt{arb\_mat\_t res, const arb\_mat\_t mat1, const arb\_mat\_t mat2, slong prec})

Sets \texttt{res} to the difference of \texttt{mat1} and \texttt{mat2}. The operands must have the same dimensions.

void \texttt{arb\_mat\_mul\_classical} (\texttt{arb\_mat\_t C, const arb\_mat\_t A, const arb\_mat\_t B, slong prec})

void \texttt{arb\_mat\_mul\_threaded} (\texttt{arb\_mat\_t C, const arb\_mat\_t A, const arb\_mat\_t B, slong prec})

void \texttt{arb\_mat\_mul} (\texttt{arb\_mat\_t res, const arb\_mat\_t mat1, const arb\_mat\_t mat2, slong prec})

Sets \texttt{res} to the matrix product of \texttt{mat1} and \texttt{mat2}. The operands must have compatible dimensions for matrix multiplication.

The \texttt{threaded} version splits the computation over the number of threads returned by \texttt{flint\_get\_num\_threads()}. The default version automatically calls the \texttt{threaded} version if the matrices are sufficiently large and more than one thread can be used.

void \texttt{arb\_mat\_sqr\_classical} (\texttt{arb\_mat\_t B, const arb\_mat\_t A, slong prec})

void \texttt{arb\_mat\_sqr} (\texttt{arb\_mat\_t res, const arb\_mat\_t mat, slong prec})

Sets \texttt{res} to the matrix square of \texttt{mat}. The operands must both be square with the same dimensions.
void *arb_mat_pow_ui (arb_mat_t res, const arb_mat_t mat, ulong exp, slong prec)

Sets res to mat raised to the power exp. Requires that mat is a square matrix.

3.5.11 Scalar arithmetic

void *arb_mat_scalar_mul_2exp_si (arb_mat_t B, const arb_mat_t A, slong c)

Sets B to A multiplied by \(2^c\).

void *arb_mat_scalar_addmul_si (arb_mat_t B, const arb_mat_t A, slong c, slong prec)

void *arb_mat_scalar_addmul_fmpz (arb_mat_t B, const arb_mat_t A, const fmpz_t c, slong prec)

void *arb_mat_scalar_addmul_arb (arb_mat_t B, const arb_mat_t A, const arb_t c, slong prec)

Sets B to \(B + A \times c\).

void *arb_mat_scalar_mul_si (arb_mat_t B, const arb_mat_t A, slong c, slong prec)

void *arb_mat_scalar_mul_fmpz (arb_mat_t B, const arb_mat_t A, const fmpz_t c, slong prec)

void *arb_mat_scalar_mul_arb (arb_mat_t B, const arb_mat_t A, const arb_t c, slong prec)

Sets B to \(A \times c\).

void *arb_mat_scalar_div_si (arb_mat_t B, const arb_mat_t A, slong c, slong prec)

void *arb_mat_scalar_div_fmpz (arb_mat_t B, const arb_mat_t A, const fmpz_t c, slong prec)

void *arb_mat_scalar_div_arb (arb_mat_t B, const arb_mat_t A, const arb_t c, slong prec)

Sets B to \(A/c\).

3.5.12 Gaussian elimination and solving

int *arb_mat_lu (slong * perm, arb_mat_t LU, const arb_mat_t A, slong prec)

Given an \(n \times n\) matrix \(A\), computes an LU decomposition \(PLU = A\) using Gaussian elimination with partial pivoting. The input and output matrices can be the same, performing the decomposition in-place.

Entry \(i\) in the permutation vector perm is set to the row index in the input matrix corresponding to row \(i\) in the output matrix.

The algorithm succeeds and returns nonzero if it can find \(n\) invertible (i.e. not containing zero) pivot entries. This guarantees that the matrix is invertible.

The algorithm fails and returns zero, leaving the entries in \(P\) and \(LU\) undefined, if it cannot find \(n\) invertible pivot elements. In this case, either the matrix is singular, the input matrix was computed to insufficient precision, or the LU decomposition was attempted at insufficient precision.

void *arb_mat_solve_lu_precomp (arb_mat_t X, const slong * perm, const arb_mat_t LU, const arb_mat_t B, slong prec)

Solves \(AX = B\) given the precomputed nonsingular LU decomposition \(A = PLU\). The matrices \(X\) and \(B\) are allowed to be aliased with each other, but \(X\) is not allowed to be aliased with \(LU\).

int *arb_mat_solve (arb_mat_t X, const arb_mat_t A, const arb_mat_t B, slong prec)

Solves \(AX = B\) where \(A\) is a nonsingular \(n \times n\) matrix and \(X\) and \(B\) are \(n \times m\) matrices, using LU decomposition.

If \(m > 0\) and \(A\) cannot be inverted numerically (indicating either that \(A\) is singular or that the precision is insufficient), the values in the output matrix are left undefined and zero is returned. A nonzero return value guarantees that \(A\) is invertible and that the exact solution matrix is contained in the output.
int arb_mat_inv(const arb_mat_t X, const arb_mat_t A, slong prec)

Sets \( X = A^{-1} \) where \( A \) is a square matrix, computed by solving the system \( AX = I \).

If \( A \) cannot be inverted numerically (indicating either that \( A \) is singular or that the precision is insufficient), the values in the output matrix are left undefined and zero is returned. A nonzero return value guarantees that the matrix is invertible and that the exact inverse is contained in the output.

void arb_mat_det(const arb_mat_t A, slong prec)

Computes the determinant of the matrix, using Gaussian elimination with partial pivoting. If at some point an invertible pivot element cannot be found, the elimination is stopped and the magnitude of the determinant of the remaining submatrix is bounded using Hadamard’s inequality.

3.5.13 Characteristic polynomial

void _arb_mat_charpoly(const arb_mat_t mat, slong prec)

Sets \( cp \) to the characteristic polynomial of \( mat \) which must be a square matrix. If the matrix has \( n \) rows, the underscore method requires space for \( n + 1 \) output coefficients. Employs a division-free algorithm using \( O(n^4) \) operations.

3.5.14 Special functions

void arb_mat_exp(const arb_mat_t B, const arb_mat_t A, slong prec)

Sets \( B \) to the exponential of the matrix \( A \), defined by the Taylor series

\[
\exp(A) = \sum_{k=0}^{\infty} \frac{A^k}{k!}.
\]

The function is evaluated as \( \exp(A/2^r)^{2^r} \), where \( r \) is chosen to give rapid convergence. The series is evaluated using rectangular splitting.

The elementwise error when truncating the Taylor series after \( N \) terms is bounded by the error in the infinity norm, for which we have

\[
\left\| \exp(2^{-r}A) - \sum_{k=0}^{N-1} \frac{(2^{-r}A)^k}{k!} \right\|_{\infty} \leq \sum_{k=N}^{\infty} \frac{(2^{-r}A)^k}{k!} \leq \sum_{k=N}^{\infty} \frac{(2^{-r}\|A\|)k}{k!}.
\]

We bound the sum on the right using \( \text{mag\_exp\_tail()} \).

void arb_mat_trace(const arb_mat_t mat, slong prec)

Sets \( \text{trace} \) to the trace of the matrix, i.e. the sum of entries on the main diagonal of \( mat \). The matrix is required to be square.

3.6 arb_calc.h – calculus with real-valued functions

This module provides functions for operations of calculus over the real numbers (intended to include root-finding, optimization, integration, and so on). It is planned that the module will include two types of algorithms:

- Interval algorithms that give provably correct results. An example would be numerical integration on an interval by dividing the interval into small balls and evaluating the function on each ball, giving rigorous upper and lower bounds.
• Conventional numerical algorithms that use heuristics to estimate the accuracy of a result, without guaranteeing
that it is correct. An example would be numerical integration based on pointwise evaluation, where the error is
estimated by comparing the results with two different sets of evaluation points. Ball arithmetic then still tracks
the accuracy of the function evaluations.

Any algorithms of the second kind will be clearly marked as such.

### 3.6.1 Types, macros and constants

**arb_calc_func_t**

Typedef for a pointer to a function with signature:

```c
int func(arb_ptr out, const arb_t inp, void * param, slong order, slong prec)
```

implementing a univariate real function $f(x)$. When called, `func` should write to `out` the first order coefficients
in the Taylor series expansion of $f(x)$ at the point `inp`, evaluated at a precision of `prec` bits. The `param` argument
may be used to pass through additional parameters to the function. The return value is reserved for future use as an
error code. It can be assumed that `out` and `inp` are not aliased and that `order` is positive.

**ARB_CALC_SUCCESS**
Return value indicating that an operation is successful.

**ARB_CALC_IMPRECISE_INPUT**
Return value indicating that the input to a function probably needs to be computed more accurately.

**ARB_CALC_NO_CONVERGENCE**
Return value indicating that an algorithm has failed to convergence, possibly due to the problem not having a
solution, the algorithm not being applicable, or the precision being insufficient.

### 3.6.2 Debugging

**int arb_calc_verbose**
If set, enables printing information about the calculation to standard output.

### 3.6.3 Subdivision-based root finding

**arf_interval_struct**

**arf_interval_t**
An `arf_interval_struct` consists of a pair of `arf_struct`, representing an interval used for
subdivision-based root-finding. An `arf_interval_t` is defined as an array of length one of type
`arf_interval_struct`, permitting an `arf_interval_t` to be passed by reference.

**arf_interval_ptr**
Alias for `arf_interval_struct *`, used for vectors of intervals.

**arf_interval_scptr**
Alias for `const arf_interval_struct *`, used for vectors of intervals.

```c
void arf_interval_init (arf_interval_t v)
void arf_interval_clear (arf_interval_t v)
arf_interval_ptr arf_interval_vec_init (slong n)
void _arf_interval_vec_clear (arf_interval_ptr v, slong n)
void arf_interval_set (arf_interval_t v, const arf_interval_t u)
```
void \texttt{arf\_interval\_swap} (\texttt{arf\_interval\_t v, arf\_interval\_t u})

void \texttt{arf\_interval\_get\_arb} (\texttt{arb\_t x, const arf\_interval\_t v, slong prec})

void \texttt{arf\_interval\_printd} (const \texttt{arf\_interval\_t v, slong n})

Helper functions for endpoint-based intervals.

\textbf{slong \texttt{arb\_calc\_isolate\_roots}} (\texttt{arf\_interval\_ptr * found, int ** flags, arb\_calc\_func\_t func, void * param, const arf\_interval\_t interval, slong maxdepth, slong maxeval, slong maxfound, slong prec})

Rigorously isolates single roots of a real analytic function on the interior of an interval.

This routine writes an array of \textit{n} interesting subintervals of \textit{interval} to \textit{found} and corresponding flags to \textit{flags}, returning the integer \textit{n}. The output has the following properties:

- The function has no roots on \textit{interval} outside of the output subintervals.
- Subintervals are sorted in increasing order (with no overlap except possibly starting and ending with the same point).
- Subintervals with a flag of 1 contain exactly one (single) root.
- Subintervals with any other flag may or may not contain roots.

If no flags other than 1 occur, all roots of the function on \textit{interval} have been isolated. If there are output subintervals on which the existence or nonexistence of roots could not be determined, the user may attempt further searches on those subintervals (possibly with increased precision and/or increased bounds for the breaking criteria). Note that roots of multiplicity higher than one and roots located exactly at endpoints cannot be isolated by the algorithm.

The following breaking criteria are implemented:

- At most \textit{maxdepth} recursive subdivisions are attempted. The smallest details that can be distinguished are therefore about \(2^{-\text{maxdepth}}\) times the width of \textit{interval}. A typical, reasonable value might be between 20 and 50.
- If the total number of tested subintervals exceeds \textit{maxeval}, the algorithm is terminated and any untested subintervals are added to the output. The total number of calls to \textit{func} is thereby restricted to a small multiple of \textit{maxeval} (the actual count can be slightly higher depending on implementation details). A typical, reasonable value might be between 100 and 100000.
- The algorithm terminates if \textit{maxfound} roots have been isolated. In particular, setting \textit{maxfound} to 1 can be used to locate just one root of the function even if there are numerous roots. To try to find all roots, \texttt{LONG\_MAX} may be passed.

The argument \textit{prec} denotes the precision used to evaluate the function. It is possibly also used for some other arithmetic operations performed internally by the algorithm. Note that it probably does not make sense for \textit{maxdepth} to exceed \textit{prec}.

Warning: it is assumed that subdivision points of \textit{interval} can be represented exactly as floating-point numbers in memory. Do not pass \(1 \pm 2^{-100}\) as input.

\textbf{int \texttt{arb\_calc\_refine\_root\_bisect}} (\texttt{arf\_interval\_t r, arb\_calc\_func\_t func, void * param, const arf\_interval\_t start, slong iter, slong prec})

Given an interval \textit{start} known to contain a single root of \textit{func}, refines it using \textit{iter} bisection steps. The algorithm can return a failure code if the sign of the function at an evaluation point is ambiguous. The output \textit{r} is set to a valid isolating interval (possibly just \textit{start}) even if the algorithm fails.
3.6.4 Newton-based root finding

void **arb_calc_newton_conv_factor** (arb_t conv_factor, arb_calc_func_t func, void *param, const
arb_t conv_region, slong prec)

Given an interval \( I \) specified by \( \text{conv} \_\text{region} \), evaluates a bound for \( C = \sup_{t,u \in I} \frac{1}{2} |f''(t)|/|f'(u)| \), where \( f \) is the function specified by \( \text{func} \) and \( \text{param} \). The bound is obtained by evaluating \( f'(I) \) and \( f''(I) \) directly. If \( f \) is ill-conditioned, \( I \) may need to be extremely precise in order to get an effective, finite bound for \( C \).

int **arb_calc_newton_step** (arb_t xnew, arb_calc_func_t func, void *param, const
arb_t x, const
arb_t conv_region, const arb_t conv_factor, slong prec)

Performs a single step with an interval version of Newton’s method. The input consists of the function \( f \) specified by \( \text{func} \) and \( \text{param} \), a ball \( x = [m - r, m + r] \) known to contain a single root of \( f \), a ball \( I \) (\( \text{conv} \_\text{region} \)) containing \( x \) with an associated bound (\( \text{conv} \_\text{factor} \)) for \( C = \sup_{t,u \in I} \frac{1}{2} |f''(t)|/|f'(u)| \), and a working precision \( \text{prec} \).

The Newton update consists of setting \( x' = [m' - r', m' + r'] \) where \( m' = m - f(m)/f'(m) \) and \( r' = Cr^2 \). The expression \( m - f(m)/f'(m) \) is evaluated using ball arithmetic at a working precision of \( \text{prec} \) bits, and the rounding error during this evaluation is accounted for in the output. We now check that \( x' \in I \) and \( r' < r \). If both conditions are satisfied, we set \( x\_\text{new} \) to \( x' \) and return **ARB_CALC_SUCCESS**. If either condition fails, we set \( x\_\text{new} \) to \( x \) and return **ARB_CALC_NO_CONVERGENCE**, indicating that no progress is made.

int **arb_calc_refine_root_newton** (arb_t r, arb_calc_func_t func, void *param, const
arb_t start, const
arb_t conv_region, const arb_t conv_factor,
slong eval_extra_prec, slong prec)

Refines a precise estimate of a single root of a function to high precision by performing several Newton steps, using nearly optimally chosen doubling precision steps.

The inputs are defined as for **arb_calc_newton_step**, except for the precision parameters: \( \text{prec} \) is the target accuracy and \( \text{eval} \_\text{extra} \_\text{prec} \) is the estimated number of guard bits that need to be added to evaluate the function accurately close to the root (for example, if the function is a polynomial with large coefficients of alternating signs and Horner’s rule is used to evaluate it, the extra precision should typically be approximately the bit size of the coefficients).

This function returns **ARB_CALC_SUCCESS** if all attempted Newton steps are successful (note that this does not guarantee that the computed root is accurate to \( \text{prec} \) bits, which has to be verified by the user), only that it is more accurate than the starting ball.

On failure, **ARB_CALC_IMPRECISE_INPUT** or **ARB_CALC_NO_CONVERGENCE** may be returned. In this case, \( r \) is set to a ball for the root which is valid but likely does have full accuracy (it can possibly just be equal to the starting ball).

3.7 acb.h – complex numbers

An **acb_t** represents a complex number with error bounds. An **acb_t** consists of a pair of real number balls of type **arb_struct**, representing the real and imaginary part with separate error bounds.

An **acb_t** thus represents a rectangle \([m_1 - r_1, m_1 + r_1] + [m_2 - r_2, m_2 + r_2]i\) in the complex plane. This is used instead of a disk or square representation (consisting of a complex floating-point midpoint with a single radius), since it allows implementing many operations more conveniently by splitting into ball operations on the real and imaginary parts. It also allows tracking when complex numbers have an exact (for example exactly zero) real part and an inexact imaginary part, or vice versa.

The interface for the **acb_t** type is slightly less developed than that for the **arb_t** type. In many cases, the user can easily perform missing operations by directly manipulating the real and imaginary parts.
3.7.1 Types, macros and constants

**acb_struct**

**acb_t**

An *acb_struct* consists of a pair of *arb_struct*s. An *acb_t* is defined as an array of length one of type *acb_struct*, permitting an *acb_t* to be passed by reference.

**acb_ptr**

Alias for *acb_struct* *, used for vectors of numbers.

**acb_srcptr**

Alias for *const acb_struct* *, used for vectors of numbers when passed as constant input to functions.

**acb_realref** (x)

Macro returning a pointer to the real part of x as an *arb_t*.

**acb_imagref** (x)

Macro returning a pointer to the imaginary part of x as an *arb_t*.

3.7.2 Memory management

void **acb_init** (*acb_t* x)

Initializes the variable x for use, and sets its value to zero.

void **acb_clear** (*acb_t* x)

Clears the variable x, freeing or recycling its allocated memory.

**acb_ptracb_vec_init** (slong n)

Returns a pointer to an array of n initialized *acb_struct*s.

void **acb_vec_clear** (*acb_ptr v*, slong n)

Clears an array of n initialized *acb_struct*s.

3.7.3 Basic manipulation

int **acb_is_zero** (const *acb_t* z)

Returns nonzero iff z is zero.

int **acb_is_one** (const *acb_t* z)

Returns nonzero iff z is exactly 1.

int **acb_is_finite** (const *acb_t* z)

Returns nonzero iff z certainly is finite.

int **acb_is_exact** (const *acb_t* z)

Returns nonzero iff z is exact.

int **acb_is_int** (const *acb_t* z)

Returns nonzero iff z is an exact integer.

void **acb_zero** (*acb_t* z)

void **acb_one** (*acb_t* z)

void **acb_onei** (*acb_t* z)

Sets z respectively to 0, 1, \( i = \sqrt{-1} \).

void **acb_set** (*acb_t* z, const *acb_t* x)

void **acb_set_ui** (*acb_t* z, slong x)
void \texttt{acb\_set\_si} (\texttt{acb\_t z}, \texttt{slong x})
void \texttt{acb\_set\_d} (\texttt{acb\_t z}, \texttt{double x})
void \texttt{acb\_set\_fmpz} (\texttt{acb\_t z}, \texttt{const fmpz\_t x})
void \texttt{acb\_set\_arb} (\texttt{acb\_t z}, \texttt{const arb\_t c})

Sets \( z \) to the value of \( x \).

void \texttt{acb\_set\_si\_si} (\texttt{acb\_t z}, \texttt{slong x}, \texttt{slong y})
void \texttt{acb\_set\_d\_d} (\texttt{acb\_t z}, \texttt{double x}, \texttt{double y})
void \texttt{acb\_set\_fmpz\_fmpz} (\texttt{acb\_t z}, \texttt{const fmpz\_t x}, \texttt{const fmpz\_t y})
void \texttt{acb\_set\_arb\_arb} (\texttt{acb\_t z}, \texttt{const arb\_t x}, \texttt{const arb\_t y})

Sets the real and imaginary part of \( z \) to the values \( x \) and \( y \) respectively.

void \texttt{acb\_set\_fmpq} (\texttt{acb\_t z}, \texttt{const fmpq\_t x}, \texttt{slong prec})
void \texttt{acb\_set\_round} (\texttt{acb\_t z}, \texttt{const \_acb\_t x}, \texttt{slong prec})
void \texttt{acb\_set\_round\_fmpz} (\texttt{acb\_t z}, \texttt{const fmpz\_t x}, \texttt{slong prec})
void \texttt{acb\_set\_round\_arb} (\texttt{acb\_t z}, \texttt{const arb\_t x}, \texttt{slong prec})

Sets \( z \) to \( x \), rounded to \( \texttt{prec} \) bits.

void \texttt{acb\_swap} (\texttt{acb\_t z}, \texttt{acb\_t x})
Swaps \( z \) and \( x \) efficiently.

void \texttt{acb\_add\_error\_mag} (\texttt{acb\_t x}, \texttt{const mag\_t err})

Adds \( \texttt{err} \) to the error bounds of both the real and imaginary parts of \( x \), modifying \( x \) in-place.

### 3.7.4 Input and output

void \texttt{acb\_print} (\texttt{const \_acb\_t x})
Prints the internal representation of \( x \).

void \texttt{acb\_printd} (\texttt{const \_acb\_t x}, \texttt{slong digits})
Prints \( x \) in decimal. The printed value of the radius is not adjusted to compensate for the fact that the binary-to-decimal conversion of both the midpoint and the radius introduces additional error.

### 3.7.5 Random number generation

void \texttt{acb\_randtest} (\texttt{acb\_t z}, \texttt{flint\_rand\_t state}, \texttt{slong prec}, \texttt{slong mag\_bits})
Generates a random complex number by generating separate random real and imaginary parts.

void \texttt{acb\_randtest\_special} (\texttt{acb\_t z}, \texttt{flint\_rand\_t state}, \texttt{slong prec}, \texttt{slong mag\_bits})
Generates a random complex number by generating separate random real and imaginary parts. Also generates NaNs and infinities.

void \texttt{acb\_randtest\_precise} (\texttt{acb\_t z}, \texttt{flint\_rand\_t state}, \texttt{slong prec}, \texttt{slong mag\_bits})
Generates a random complex number with precise real and imaginary parts.

void \texttt{acb\_randtest\_param} (\texttt{acb\_t z}, \texttt{flint\_rand\_t state}, \texttt{slong prec}, \texttt{slong mag\_bits})
Generates a random complex number, with very high probability of generating integers and half-integers.
3.7.6 Precision and comparisons

int acb_equal (const acb_t x, const acb_t y)
Returns nonzero iff x and y are identical as sets, i.e. if the real and imaginary parts are equal as balls.

Note that this is not the same thing as testing whether both x and y certainly represent the same complex number, unless either x or y is exact (and neither contains NaN). To test whether both operands might represent the same mathematical quantity, use acb_overlaps() or acb_contains(), depending on the circumstance.

int acb_eq (const acb_t x, const acb_t y)
Returns nonzero iff x and y are certainly equal, as determined by testing that arb_eq() holds for both the real and imaginary parts.

int acb_ne (const acb_t x, const acb_t y)
Returns nonzero iff x and y are certainly not equal, as determined by testing that arb_ne() holds for either the real or imaginary parts.

int acb_overlaps (const acb_t x, const acb_t y)
Returns nonzero iff x and y have some point in common.

void acb_get_abs_ubound_arf (arf_t u, const acb_t z, slong prec)
Sets u to an upper bound for the absolute value of z, computed using a working precision of prec bits.

void acb_get_abs_lbound_arf (arf_t u, const acb_t z, slong prec)
Sets u to a lower bound for the absolute value of z, computed using a working precision of prec bits.

void acb_get_rad_ubound_arf (arf_t u, const acb_t z, slong prec)
Sets u to an upper bound for the error radius of z (the value is currently not computed tightly).

void acb_get_mag (mag_t u, const acb_t x)
Sets u to an upper bound for the absolute value of x.

void acb_get_mag_lower (mag_t u, const acb_t x)
Sets u to a lower bound for the absolute value of x.

int acb_contains_fmpq (const acb_t x, const fmpq_t y)

int acb_contains_fmpz (const acb_t x, const fmpz_t y)

int acb_contains (const acb_t x, const acb_t y)
Returns nonzero iff y is contained in x.

int acb_contains_zero (const acb_t x)
Returns nonzero iff zero is contained in x.

int acb_contains_int (const acb_t x)
Returns nonzero iff the complex interval represented by x contains an integer.

slong acb_rel_error_bits (const acb_t x)
Returns the effective relative error of x measured in bits. This is computed as if calling arb_rel_error_bits() on the real ball whose midpoint is the larger out of the real and imaginary midpoints of x, and whose radius is the larger out of the real and imaginary radiuses of x.

slong acb_rel_accuracy_bits (const acb_t x)
Returns the effective relative accuracy of x measured in bits, equal to the negative of the return value from acb_rel_error_bits().

slong acb_bits (const acb_t x)
Returns the maximum of arb_bits applied to the real and imaginary parts of x, i.e. the minimum precision sufficient to represent x exactly.

void acb_indeterminate (acb_t x)
Sets x to \([NaN \pm \infty] + [NaN \pm \infty]i\), representing an indeterminate result.
void **acb_trim** (acb_t y, const acb_t x)
Sets y to a a copy of x with both the real and imaginary parts trimmed (see *arb_trim()*)

int **acb_is_real** (const acb_t x)
Returns nonzero iff the imaginary part of x is zero. It does not test whether the real part of x also is finite.

int **acb_get_unique_fmpz** (fmpz_t z, const acb_t x)
If x contains a unique integer, sets z to that value and returns nonzero. Otherwise (if x represents no integers or more than one integer), returns zero.

3.7.7 Complex parts

void **acb_get_real** (arb_t re, const acb_t z)
Sets re to the real part of z.

void **acb_get_imag** (arb_t im, const acb_t z)
Sets im to the imaginary part of z.

void **acb_arg** (arb_t r, const acb_t z, slong prec)
Sets r to a real interval containing the complex argument (phase) of z. We define the complex argument have a discontinuity on (−∞, 0], with the special value arg(0) = 0, and arg(a + 0i) = π for a < 0. Equivalently, if z = a + bi, the argument is given by atan2(b, a) (see *arb_atan2()*).

void **acb_abs** (arb_t r, const acb_t z, slong prec)
Sets r to the absolute value of z.

3.7.8 Arithmetic

void **acb_neg** (acb_t z, const acb_t x)
Sets z to the negation of x.

void **acb_conj** (acb_t z, const acb_t x)
Sets z to the complex conjugate of x.

void **acb_add_ui** (acb_t z, const acb_t x, ulong y, slong prec)
void **acb_add_si** (acb_t z, const acb_t x, slong y, slong prec)
void **acb_add_fmpz** (acb_t z, const acb_t x, const fmpz_t y, slong prec)
void **acb_add_arb** (acb_t z, const acb_t x, const arb_t y, slong prec)
void **acb_add** (acb_t z, const acb_t x, const acb_t y, slong prec)
Sets z to the sum of x and y.

void **acb_sub_ui** (acb_t z, const acb_t x, ulong y, slong prec)
void **acb_sub_si** (acb_t z, const acb_t x, slong y, slong prec)
void **acb_sub_fmpz** (acb_t z, const acb_t x, const fmpz_t y, slong prec)
void **acb_sub_arb** (acb_t z, const acb_t x, const arb_t y, slong prec)
void **acb_sub** (acb_t z, const acb_t x, const acb_t y, slong prec)
Sets z to the difference of x and y.

void **acb_mul_onei** (acb_t z, const acb_t x)
Sets z to x multiplied by the imaginary unit.

void **acb_div_onei** (acb_t z, const acb_t x)
Sets z to x divided by the imaginary unit.
void \texttt{acb\_mul\_ui} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{ulong} y, \texttt{slong} \texttt{prec})\)
void \texttt{acb\_mul\_si} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{slong} y, \texttt{slong} \texttt{prec})\)
void \texttt{acb\_mul\_fmpz} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{fmpz\_t} y, \texttt{slong} \texttt{prec})\)
void \texttt{acb\_mul\_arb} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{arb\_t} y, \texttt{slong} \texttt{prec})\)

Sets \(z\) to the product of \(x\) and \(y\).

void \texttt{acb\_mul} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{acb\_t} y, \texttt{slong} \texttt{prec})\)

Sets \(z\) to the product of \(x\) and \(y\). If at least one part of \(x\) or \(y\) is zero, the operations is reduced to two real multiplications. If \(x\) and \(y\) are the same pointers, they are assumed to represent the same mathematical quantity and the squaring formula is used.

void \texttt{acb\_mul\_2exp\_si} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{slong} e)\)
void \texttt{acb\_mul\_2exp\_fmpz} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{fmpz\_t} e)\)

Sets \(z\) to \(x\) multiplied by \(2^e\), without rounding.

void \texttt{acb\_cube} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{slong} \texttt{prec})\)

Sets \(z\) to \(x\) cubed, computed efficiently using two real squarings, two real multiplications, and scalar operations.

void \texttt{acb\_addmul} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{acb\_t} y, \texttt{slong} \texttt{prec})\)
void \texttt{acb\_addmul\_ui} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{ulong} y, \texttt{slong} \texttt{prec})\)
void \texttt{acb\_addmul\_si} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{slong} y, \texttt{slong} \texttt{prec})\)
void \texttt{acb\_addmul\_fmpz} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{fmpz\_t} y, \texttt{slong} \texttt{prec})\)
void \texttt{acb\_addmul\_arb} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{arb\_t} y, \texttt{slong} \texttt{prec})\)

Sets \(z\) to \(z\) plus the product of \(x\) and \(y\).

void \texttt{acb\_submul} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{acb\_t} y, \texttt{slong} \texttt{prec})\)
void \texttt{acb\_submul\_ui} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{ulong} y, \texttt{slong} \texttt{prec})\)
void \texttt{acb\_submul\_si} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{slong} y, \texttt{slong} \texttt{prec})\)
void \texttt{acb\_submul\_fmpz} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{fmpz\_t} y, \texttt{slong} \texttt{prec})\)
void \texttt{acb\_submul\_arb} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{arb\_t} y, \texttt{slong} \texttt{prec})\)

Sets \(z\) to \(z\) minus the product of \(x\) and \(y\).

void \texttt{acb\_inv} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{slong} \texttt{prec})\)

Sets \(z\) to the multiplicative inverse of \(x\).

void \texttt{acb\_div\_ui} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{ulong} y, \texttt{slong} \texttt{prec})\)
void \texttt{acb\_div\_si} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{slong} y, \texttt{slong} \texttt{prec})\)
void \texttt{acb\_div\_fmpz} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{fmpz\_t} y, \texttt{slong} \texttt{prec})\)
void \texttt{acb\_div\_arb} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{arb\_t} y, \texttt{slong} \texttt{prec})\)
void \texttt{acb\_div} \((\texttt{acb\_t} z, \texttt{acb\_t} x, \texttt{acb\_t} y, \texttt{slong} \texttt{prec})\)

Sets \(z\) to the quotient of \(x\) and \(y\).

### 3.7.9 Mathematical constants

void \texttt{acb\_const\_pi} \((\texttt{acb\_t} y, \texttt{slong} \texttt{prec})\)

Sets \(y\) to the constant \(\pi\).
3.7.10 Powers and roots
void \texttt{acb\_sqrt} (\texttt{acb\_t} \texttt{r}, const \texttt{acb\_t} \texttt{z}, slong \texttt{prec})
Sets \( r \) to the square root of \( z \). If either the real or imaginary part is exactly zero, only a single real square root is needed. Generally, we use the formula \( \sqrt{a + bi} = u/2 + ib/u, u = \sqrt{2(|a + bi| + a)} \), requiring two real square root extractions.

void \texttt{acb\_rsgqt} (\texttt{acb\_t} \texttt{r}, const \texttt{acb\_t} \texttt{z}, slong \texttt{prec})
Sets \( r \) to the reciprocal square root of \( z \). If either the real or imaginary part is exactly zero, only a single real reciprocal square root is needed. Generally, we use the formula \( 1/\sqrt{a + bi} = (|a + r|) / v, r = |a + bi|, v = \sqrt{|a + bi + r|^2} \), requiring one real square root and one real reciprocal square root.

void \texttt{acb\_quadratic\_roots\_fmpz} (\texttt{acb\_t} \texttt{r1}, \texttt{acb\_t} \texttt{r2}, const \texttt{fmpz\_t} \texttt{a}, const \texttt{fmpz\_t} \texttt{b}, const \texttt{fmpz\_t} \texttt{c}, slong \texttt{prec})
Sets \( r1 \) and \( r2 \) to the roots of the quadratic polynomial \( ax^2 + bx + c \). Requires that \( a \) is nonzero. This function is implemented so that both roots are computed accurately even when direct use of the quadratic formula would lose accuracy.

void \texttt{acb\_root\_ui} (\texttt{acb\_t} \texttt{r}, const \texttt{acb\_t} \texttt{z}, ulong \texttt{k}, slong \texttt{prec})
Sets \( r \) to the principal \( k \)-th root of \( z \).

void \texttt{acb\_pow\_fmpz} (\texttt{acb\_t} \texttt{y}, const \texttt{acb\_t} \texttt{b}, const \texttt{fmpz\_t} \texttt{e}, slong \texttt{prec})
void \texttt{acb\_pow\_ui} (\texttt{acb\_t} \texttt{y}, const \texttt{acb\_t} \texttt{b}, ulong \texttt{e}, slong \texttt{prec})
void \texttt{acb\_pow\_si} (\texttt{acb\_t} \texttt{y}, const \texttt{acb\_t} \texttt{b}, slong \texttt{e}, slong \texttt{prec})
Sets \( y = b^e \) using binary exponentiation (with an initial division if \( e < 0 \)). Note that these functions can get slow if the exponent is extremely large (in such cases \texttt{acb\_pow()} may be superior).

void \texttt{acb\_pow\_arb} (\texttt{acb\_t} \texttt{z}, const \texttt{acb\_t} \texttt{x}, const \texttt{arb\_t} \texttt{y}, slong \texttt{prec})
void \texttt{acb\_pow\_arb} (\texttt{acb\_t} \texttt{z}, const \texttt{acb\_t} \texttt{x}, const \texttt{arb\_t} \texttt{y}, slong \texttt{prec})
Sets \( z = x^y \), computed using binary exponentiation if \( y \) if a small exact integer, as \( z = (x^{1/2})^{2^y} \) if \( y \) is a small exact half-integer, and generally as \( z = \exp(y \log x) \).

3.7.11 Exponentials and logarithms
void \texttt{acb\_exp} (\texttt{acb\_t} \texttt{y}, const \texttt{acb\_t} \texttt{z}, slong \texttt{prec})
Sets \( y \) to the exponential function of \( z \), computed as \( \exp(a + bi) = \exp(a) (\cos(b) + \sin(b)i) \).

void \texttt{acb\_exp\_pi\_i} (\texttt{acb\_t} \texttt{y}, const \texttt{acb\_t} \texttt{z}, slong \texttt{prec})
Sets \( y \) to \( \exp(\pi i z) \).

void \texttt{acb\_exp\_invexp} (\texttt{acb\_t} \texttt{s}, \texttt{acb\_t} \texttt{t}, const \texttt{acb\_t} \texttt{z}, slong \texttt{prec})
Sets \( v = \exp(z) \) and \( w = \exp(-z) \).

void \texttt{acb\_log} (\texttt{acb\_t} \texttt{y}, const \texttt{acb\_t} \texttt{z}, slong \texttt{prec})
Sets \( y \) to the principal branch of the natural logarithm of \( z \), computed as \( \log(a + bi) = \frac{1}{2} \log(a^2 + b^2) + i \arg(a + bi) \).

void \texttt{acb\_log\_lp} (\texttt{acb\_t} \texttt{z}, const \texttt{acb\_t} \texttt{x}, slong \texttt{prec})
Sets \( z = \log(1 + x) \), computed accurately when \( x \approx 0 \).

3.7.12 Trigonometric functions
void \texttt{acb\_sin} (\texttt{acb\_t} \texttt{s}, const \texttt{acb\_t} \texttt{z}, slong \texttt{prec})
void \texttt{acb\_cos} (\texttt{acb\_t} \texttt{c}, const \texttt{acb\_t} \texttt{z}, slong \texttt{prec})
void \texttt{acb\_sin\_cos} (\texttt{acb\_t} \textit{s}, \texttt{acb\_t} \textit{c}, const \texttt{acb\_t} \textit{z}, slong \textit{prec})

Sets \( s = \sin(z), c = \cos(z), \) evaluated as \( \sin(a + bi) = \sin(a) \cosh(b) + i \cos(a) \sinh(b), \cos(a + bi) = \cos(a) \cosh(b) - i \sin(a) \sinh(b) \).

void \texttt{acb\_tan} (\texttt{acb\_t} \textit{s}, const \texttt{acb\_t} \textit{z}, slong \textit{prec})

Sets \( s = \tan(z) = \sin(z) / \cos(z) \). For large imaginary parts, the function is evaluated in a numerically stable way as \( \pm i \) plus a decreasing exponential factor.

void \texttt{acb\_cot} (\texttt{acb\_t} \textit{s}, const \texttt{acb\_t} \textit{z}, slong \textit{prec})

Sets \( s = \cot(z) = \cos(z) / \sin(z) \). For large imaginary parts, the function is evaluated in a numerically stable way as \( \pm i \) plus a decreasing exponential factor.

void \texttt{acb\_sin\_pi} (\texttt{acb\_t} \textit{s}, const \texttt{acb\_t} \textit{z}, slong \textit{prec})

void \texttt{acb\_cos\_pi} (\texttt{acb\_t} \textit{s}, const \texttt{acb\_t} \textit{z}, slong \textit{prec})

void \texttt{acb\_sin\_cos\_pi} (\texttt{acb\_t} \textit{s}, \texttt{acb\_t} \textit{c}, const \texttt{acb\_t} \textit{z}, slong \textit{prec})

Sets \( s = \sin(\pi z), c = \cos(\pi z), \) evaluating the trigonometric factors of the real and imaginary part accurately via \texttt{arb\_sin\_cos\_pi}().

void \texttt{acb\_tan\_pi} (\texttt{acb\_t} \textit{s}, const \texttt{acb\_t} \textit{z}, slong \textit{prec})

Sets \( s = \tan(\pi z) \). Uses the same algorithm as \texttt{acb\_tan()}, but evaluates the sine and cosine accurately via \texttt{arb\_sin\_cos\_pi}().

void \texttt{acb\_cot\_pi} (\texttt{acb\_t} \textit{s}, const \texttt{acb\_t} \textit{z}, slong \textit{prec})

Sets \( s = \cot(\pi z) \). Uses the same algorithm as \texttt{acb\_cot()}, but evaluates the sine and cosine accurately via \texttt{arb\_sin\_cos\_pi}().

### 3.7.13 Inverse trigonometric functions

void \texttt{acb\_asin} (\texttt{acb\_t} \textit{res}, const \texttt{acb\_t} \textit{z}, slong \textit{prec})

Sets \( \textit{res} \) to \( \text{asin}(z) = -i \log(i z + \sqrt{1 - z^2}) \).

void \texttt{acb\_acos} (\texttt{acb\_t} \textit{res}, const \texttt{acb\_t} \textit{z}, slong \textit{prec})

Sets \( \textit{res} \) to \( \text{acos}(z) = \frac{1}{2} \pi - \text{asin}(z) \).

void \texttt{acb\_atan} (\texttt{acb\_t} \textit{res}, const \texttt{acb\_t} \textit{z}, slong \textit{prec})

Sets \( \textit{res} \) to \( \text{atan}(z) = \frac{1}{2} i(\log(1 - iz) - \log(1 + iz)) \).

### 3.7.14 Hyperbolic functions

void \texttt{acb\_sinh} (\texttt{acb\_t} \textit{s}, const \texttt{acb\_t} \textit{z}, slong \textit{prec})

void \texttt{acb\_cosh} (\texttt{acb\_t} \textit{c}, const \texttt{acb\_t} \textit{z}, slong \textit{prec})

void \texttt{acb\_sinh\_cosh} (\texttt{acb\_t} \textit{s}, \texttt{acb\_t} \textit{c}, const \texttt{acb\_t} \textit{z}, slong \textit{prec})

void \texttt{acb\_tanh} (\texttt{acb\_t} \textit{s}, const \texttt{acb\_t} \textit{z}, slong \textit{prec})

void \texttt{acb\_coth} (\texttt{acb\_t} \textit{s}, const \texttt{acb\_t} \textit{z}, slong \textit{prec})

Respectively computes \( \sinh(z) = -i \sin(iz), \cosh(z) = \cos(iz), \tanh(z) = -i \tan(iz), \coth(z) = i \cot(iz) \).

### 3.7.15 Inverse hyperbolic functions

void \texttt{acb\_asinh} (\texttt{acb\_t} \textit{res}, const \texttt{acb\_t} \textit{z}, slong \textit{prec})

Sets \( \textit{res} \) to \( \text{asinh}(z) = -i \text{asin}(iz) \).
void \texttt{acb\_acosh} (\texttt{acb\_t res}, const \texttt{acb\_t z}, slong prec)
\hspace{1em}Sets \texttt{res} to $\text{acosh}(z) = \log(z + \sqrt{z^2 + 1})$.

void \texttt{acb\_atanh} (\texttt{acb\_t res}, const \texttt{acb\_t z}, slong prec)
\hspace{1em}Sets \texttt{res} to $\text{atanh}(z) = -i \text{atan}(iz)$.

### 3.7.16 Rising factorials

void \texttt{acb\_rising\_ui\_bs} (\texttt{acb\_t z}, const \texttt{acb\_t x}, ulong \texttt{n}, slong prec)
\hspace{1em}Computes the rising factorial $z = x(x+1)(x+2)\cdots(x+n-1)$.

The \texttt{bs} version uses binary splitting. The \texttt{rs} version uses rectangular splitting. The \texttt{rec} version uses either \texttt{bs} or \texttt{rs} depending on the input. The default version uses the gamma function unless \texttt{n} is a small integer.

The \texttt{rs} version takes an optional \texttt{step} parameter for tuning purposes (to use the default step length, pass zero).

void \texttt{acb\_rising\_ui\_rec} (\texttt{acb\_t z}, const \texttt{acb\_t x}, ulong \texttt{n}, slong prec)
\hspace{1em}Using binary splitting, rectangular splitting (with optional nonzero step length \texttt{step} to override the default choice), and an automatic algorithm choice.

void \texttt{acb\_rising\_ui\_get\_mag} (\texttt{mag\_t bound}, const \texttt{acb\_t x}, ulong \texttt{n})
\hspace{1em}Computes an upper bound for the absolute value of the rising factorial $z = x(x+1)(x+2)\cdots(x+n-1)$. Not currently optimized for large \texttt{n}.

### 3.7.17 Gamma function

void \texttt{acb\_gamma} (\texttt{acb\_t y}, const \texttt{acb\_t x}, slong prec)
\hspace{1em}Computes the gamma function $y = \Gamma(x)$.

void \texttt{acb\_rgamma} (\texttt{acb\_t y}, const \texttt{acb\_t x}, slong prec)
\hspace{1em}Computes the reciprocal gamma function $y = 1/\Gamma(x)$, avoiding division by zero at the poles of the gamma function.

void \texttt{acb\_lgamma} (\texttt{acb\_t y}, const \texttt{acb\_t x}, slong prec)
\hspace{1em}Computes the logarithmic gamma function $y = \log \Gamma(x)$.

The branch cut of the logarithmic gamma function is placed on the negative half-axis, which means that $\log \Gamma(z) + \log z = \log \Gamma(z+1)$ holds for all $z$, whereas $\log \Gamma(z) \neq \log \Gamma(z)$ in general. In the left half plane, the reflection formula with correct branch structure is evaluated via \texttt{acb\_log\_sin\_pi()}.

void \texttt{acb\_digamma} (\texttt{acb\_t y}, const \texttt{acb\_t x}, slong prec)
\hspace{1em}Computes the digamma function $y = \psi(x) = (\log \Gamma(x))' = \Gamma'(x)/\Gamma(x)$.

void \texttt{acb\_log\_sin\_pi} (\texttt{acb\_t res}, const \texttt{acb\_t z}, slong prec)
\hspace{1em}Computes the logarithmic sine function defined by

$$S(z) = \log(\pi) - \log \Gamma(z) + \log \Gamma(1 - z)$$
which is equal to

\[ S(z) = \int_{1/2}^{z} \pi \cot(\pi t) dt \]

where the path of integration goes through the upper half plane if \( 0 < \arg(z) \leq \pi \) and through the lower half plane if \(-\pi < \arg(z) \leq 0\). Equivalently,

\[ S(z) = \log(\sin(\pi(z - n))) \pm n\pi i, \quad n = \lfloor \text{re}(z) \rfloor \]

where the negative sign is taken if \( 0 < \arg(z) \leq \pi \) and the positive sign is taken otherwise (if the interval \( \arg(z) \) does not certainly satisfy either condition, the union of both cases is computed). After subtracting \( n \), we have \( 0 \leq \text{re}(z) < 1 \). In this strip, we use \( S(z) = \log(\sin(\pi(z))) \) if the imaginary part of \( z \) is small. Otherwise, we use \( S(z) = i\pi(z - 1/2) + \log((1 + e^{-2i\pi z})/2) \) in the lower half-plane and the conjugated expression in the upper half-plane to avoid exponent overflow.

The function is evaluated at the midpoint and the propagated error is computed from \( S'(z) \) to get a continuous change when \( z \) is non-real and \( n \) spans more than one possible integer value.

```c
void acb_polygamma(acb_t z, const acb_t s, const acb_t z, slong prec)
Sets res to the value of the generalized polygamma function \( \psi(s, z) \).
If s is a nonnegative order, this is simply the s-order derivative of the digamma function. If \( s = 0 \), this function simply calls the digamma function internally. For integers \( s \geq 1 \), it calls the Hurwitz zeta function. Note that for small integers \( s \geq 1 \), it can be faster to use `acb_poly_digamma_series()` and read off the coefficients.

The generalization to other values of \( s \) is due to Espinosa and Moll [EM2004]:

\[ \psi(s, z) = \frac{\zeta'(s + 1, z) + (\gamma + \psi(-s))\zeta(s + 1, z)}{\Gamma(-s)} \]
```

```c
void acb_barnes_g(acb_t res, const acb_t z, slong prec)
void acb_log_barnes_g(acb_t res, const acb_t z, slong prec)
Computes Barnes G-function or the logarithmic Barnes G-function, respectively. The logarithmic version has branch cuts on the negative real axis and is continuous elsewhere in the complex plane, in analogy with the logarithmic gamma function. The functional equation

\[ \log G(z + 1) = \log \Gamma(z) + \log G(z). \]

holds for all \( z \).

For small integers, we directly use the recurrence relation \( G(z + 1) = \Gamma(z)G(z) \) together with the initial value \( G(1) = 1 \). For general \( z \), we use the formula

\[ \log G(z) = (z - 1) \log \Gamma(z) - \zeta'(-1, z) + \zeta'(-1). \]
```

### 3.7.18 Zeta function

```c
void acb_zeta(acb_t z, const acb_t s, slong prec)
Sets z to the value of the Riemann zeta function \( \zeta(s) \). Note: for computing derivatives with respect to \( s \), use `acb_poly_zeta_series()` or related methods.
```

```c
void acb_hurwitz_zeta(acb_t z, const acb_t s, const acb_t a, slong prec)
Sets z to the value of the Hurwitz zeta function \( \zeta(s, a) \). Note: for computing derivatives with respect to \( s \), use `acb_poly_zeta_series()` or related methods.
```
3.7.19 Polylogarithms

void \texttt{acb\_polylog}(\texttt{acb\_t w}, \texttt{const acb\_t s}, \texttt{const acb\_t z}, \texttt{slong prec})

void \texttt{acb\_polylog\_si}(\texttt{acb\_t w}, \texttt{slong s}, \texttt{const acb\_t z}, \texttt{slong prec})

Sets \( w \) to the polylogarithm \( \text{Li}_s(z) \).

3.7.20 Arithmetic-geometric mean

void \texttt{acb\_agm1}(\texttt{acb\_t m}, \texttt{const acb\_t z}, \texttt{slong prec})

Sets \( m \) to the arithmetic-geometric mean \( M(z) = \text{agm}(1, z) \), defined such that the function is continuous in the complex plane except for a branch cut along the negative half axis (where it is continuous from above). This corresponds to always choosing an “optimal” branch for the square root in the arithmetic-geometric mean iteration.

void \texttt{acb\_agm1\_cpx}(\texttt{acb\_ptr m}, \texttt{const acb\_t z}, \texttt{slong len}, \texttt{slong prec})

Sets the coefficients in the array \( m \) to the power series expansion of the arithmetic-geometric mean at the point \( z \) truncated to length \( len \), i.e. \( M(z + x) \in \mathbb{C}[[x]] \).

3.7.21 Other special functions

void \texttt{acb\_chebyshev\_t\_ui}(\texttt{acb\_t a}, \texttt{ulong n}, \texttt{const acb\_t x}, \texttt{slong prec})

void \texttt{acb\_chebyshev\_u\_ui}(\texttt{acb\_t a}, \texttt{ulong n}, \texttt{const acb\_t x}, \texttt{slong prec})

Evaluates the Chebyshev polynomial of the first kind \( a = T_n(x) \) or the Chebyshev polynomial of the second kind \( a = U_n(x) \).

void \texttt{acb\_chebyshev\_t2\_ui}(\texttt{acb\_t a}, \texttt{acb\_t b}, \texttt{ulong n}, \texttt{const acb\_t x}, \texttt{slong prec})

void \texttt{acb\_chebyshev\_u2\_ui}(\texttt{acb\_t a}, \texttt{acb\_t b}, \texttt{ulong n}, \texttt{const acb\_t x}, \texttt{slong prec})

Simultaneously evaluates \( a = T_n(x), b = T_{n-1}(x) \) or \( a = U_n(x), b = U_{n-1}(x) \). Aliasing between \( a, b \) and \( x \) is not permitted.

3.7.22 Vector functions

void \texttt{acb\_vec\_zero}(\texttt{acb\_ptr A}, \texttt{slong n})

Sets all entries in \( \text{vec} \) to zero.

int \texttt{acb\_vec\_is\_zero}(\texttt{acb\_srcptr vec}, \texttt{slong len})

Returns nonzero iff all entries in \( x \) are zero.

int \texttt{acb\_vec\_is\_real}(\texttt{acb\_srcptr v}, \texttt{slong len})

Returns nonzero iff all entries in \( x \) have zero imaginary part.

void \texttt{acb\_vec\_set}(\texttt{acb\_ptr res}, \texttt{acb\_srcptr vec}, \texttt{slong len})

Sets \( \text{res} \) to a copy of \( \text{vec} \).

void \texttt{acb\_vec\_set\_round}(\texttt{acb\_ptr res}, \texttt{acb\_srcptr vec}, \texttt{slong len}, \texttt{slong prec})

Sets \( \text{res} \) to a copy of \( \text{vec} \), rounding each entry to \( \text{prec} \) bits.

void \texttt{acb\_vec\_neg}(\texttt{acb\_ptr res}, \texttt{acb\_srcptr vec}, \texttt{slong len})

void \texttt{acb\_vec\_add}(\texttt{acb\_ptr res}, \texttt{acb\_srcptr vec1}, \texttt{acb\_srcptr vec2}, \texttt{slong len}, \texttt{slong prec})

void \texttt{acb\_vec\_sub}(\texttt{acb\_ptr res}, \texttt{acb\_srcptr vec1}, \texttt{acb\_srcptr vec2}, \texttt{slong len}, \texttt{slong prec})

void \texttt{acb\_vec\_scalar\_submul}(\texttt{acb\_ptr res}, \texttt{acb\_srcptr vec}, \texttt{slong len}, \texttt{const acb\_t c}, \texttt{slong prec})
void _acb_vec_scalar_addmul (acb_ptr res, acb_srcptr vec, slong len, const acb_t c, slong prec)
void _acb_vec_scalar_mul (acb_ptr res, acb_srcptr vec, slong len, const acb_t c, slong prec)
void _acb_vec_scalar_mul_ui (acb_ptr res, acb_srcptr vec, slong len, ulong c, slong prec)
void _acb_vec_scalar_mul_2exp_si (acb_ptr res, acb_srcptr vec, slong len, slong c)
void _acb_vec_scalar_mul_onei (acb_ptr res, acb_srcptr vec, slong len)
void _acb_vec_scalar_div_ui (acb_ptr res, acb_srcptr vec, slong len, ulong c, slong prec)
void _acb_vec_scalar_div (acb_ptr res, acb_srcptr vec, slong len, const acb_t c, slong prec)
void _acb_vec_scalar_mul_arb (acb_ptr res, acb_srcptr vec, slong len, const arb_t c, slong prec)
void _acb_vec_scalar_div_arb (acb_ptr res, acb_srcptr vec, slong len, const arb_t c, slong prec)
void _acb_vec_scalar_mul_fmpz (acb_ptr res, acb_srcptr vec, slong len, const fmpz_t c, slong prec)
void _acb_vec_scalar_div_fmpz (acb_ptr res, acb_srcptr vec, slong len, const fmpz_t c, slong prec)

Performs the respective scalar operation elementwise.

slong _acb_vec_bits (acb_srcptr vec, slong len)
Returns the maximum of arb_bits () for all entries in vec.
void _acb_vec_set_powers (acb_ptr xs, const acb_t x, slong len, slong prec)
Sets xs to the powers 1, x, x², . . . , x^{len-1}.
void _acb_vec_add_error_arf_vec (acb_ptr res, arf_srcptr err, slong len)
Add the magnitude of each entry in err to the radius of the corresponding entry in res.
void _acb_vec_add_error_mag_vec (acb_ptr res, mag_srcptr err, slong len)
void _acb_vec_indeterminate (acb_ptr vec, slong len)
Applies acb_indeterminate () elementwise.
void _acb_vec_trim (acb_ptr vec, slong len)
Applies acb_trim () elementwise.

int _acb_vec_get_unique_fmpz_vec (fmpz * res, acb_srcptr vec, slong len)
Calls acb_get_unique_fmpz () elementwise and returns nonzero if all entries can be rounded uniquely to integers. If any entry in vec cannot be rounded uniquely to an integer, returns zero.
void _acb_vec_sort_pretty (acb_ptr vec, slong len)
Sorts the vector of complex numbers based on the real and imaginary parts. This is intended to reveal structure when printing a set of complex numbers, not to apply an order relation in a rigorous way.

3.8 acb_poly.h – polynomials over the complex numbers

An acb_poly_t represents a polynomial over the complex numbers, implemented as an array of coefficients of type acb_struct.

Most functions are provided in two versions: an underscore method which operates directly on pre-allocated arrays of coefficients and generally has some restrictions (such as requiring the lengths to be nonzero and not supporting aliasing of the input and output arrays), and a non-underscore method which performs automatic memory management and handles degenerate cases.
### 3.8.1 Types, macros and constants

**acb_poly_struct**

**acb_poly_t**
Contains a pointer to an array of coefficients (coeffs), the used length (length), and the allocated size of the array (alloc).

An `acb_poly_t` is defined as an array of length one of type `acb_poly_struct`, permitting an `acb_poly_t` to be passed by reference.

### 3.8.2 Memory management

**void acb_poly_init (acb_poly_t poly)**
Initializes the polynomial for use, setting it to the zero polynomial.

**void acb_poly_clear (acb_poly_t poly)**
Clears the polynomial, deallocating all coefficients and the coefficient array.

**void acb_poly_fit_length (acb_poly_t poly, slong len)**
Makes sure that the coefficient array of the polynomial contains at least `len` initialized coefficients.

**void _acb_poly_set_length (acb_poly_t poly, slong len)**
Directly changes the length of the polynomial, without allocating or deallocating coefficients. The value should not exceed the allocation length.

**void _acb_poly_normalise (acb_poly_t poly)**
Strips any trailing coefficients which are identical to zero.

**void acb_poly_swap (acb_poly_t poly1, acb_poly_t poly2)**
Swaps `poly1` and `poly2` efficiently.

### 3.8.3 Basic properties and manipulation

**slong acb_poly_length (const acb_poly_t poly)**
Returns the length of `poly`, i.e. zero if `poly` is identically zero, and otherwise one more than the index of the highest term that is not identically zero.

**slong acb_poly_degree (const acb_poly_t poly)**
Returns the degree of `poly`, defined as one less than its length. Note that if one or several leading coefficients are balls containing zero, this value can be larger than the true degree of the exact polynomial represented by `poly`, so the return value of this function is effectively an upper bound.

**void acb_poly_zero (acb_poly_t poly)**
Sets `poly` to the zero polynomial.

**void acb_poly_one (acb_poly_t poly)**
Sets `poly` to the constant polynomial 1.

**void acb_poly_set (acb_poly_t dest, const acb_poly_t src)**
Sets `dest` to a copy of `src`.

**void acb_poly_set_round (acb_poly_t dest, const acb_poly_t src, slong prec)**
Sets `dest` to a copy of `src`, rounded to `prec` bits.

**void acb_poly_set_coeff_si (acb_poly_t poly, slong n, slong c)**
Sets the coefficient with index `n` in `poly` to the value `c`. We require that `n` is nonnegative.

**void acb_poly_set_coeff_acb (acb_poly_t poly, slong n, const acb_t c)**
Sets the coefficient with index `n` in `poly` to the value `c`. We require that `n` is nonnegative.
void \texttt{acb\_poly\_get\_coeff\_acb} (\texttt{acb\_t v}, \texttt{const acb\_poly\_t poly}, \texttt{long n})
\begin{itemize}
\item Sets \textit{v} to the value of the coefficient with index \textit{n} in \textit{poly}. We require that \textit{n} is nonnegative.
\end{itemize}

\texttt{acb\_poly\_get\_coeff\_ptr} (\textit{poly, n})
\begin{itemize}
\item Given \textit{n} \geq 0, returns a pointer to coefficient \textit{n} of \textit{poly}, or \texttt{NULL} if \textit{n} exceeds the length of \textit{poly}.
\end{itemize}

void \_\texttt{acb\_poly\_shift\_right} (\texttt{acb\_ptr res}, \texttt{acb\_srcptr poly}, \texttt{long len}, \texttt{long n})

void \texttt{acb\_poly\_shift\_right} (\texttt{acb\_poly\_t res}, \texttt{const acb\_poly\_t poly}, \texttt{long n})
\begin{itemize}
\item Sets \textit{res} to \textit{poly} divided by \textit{x}^{\textit{n}}, throwing away the lower coefficients. We require that \textit{n} is nonnegative.
\end{itemize}

void \_\texttt{acb\_poly\_shift\_left} (\texttt{acb\_ptr res}, \texttt{acb\_srcptr poly}, \texttt{long len}, \texttt{long n})

void \texttt{acb\_poly\_shift\_left} (\texttt{acb\_poly\_t res}, \texttt{const acb\_poly\_t poly}, \texttt{long n})
\begin{itemize}
\item Sets \textit{res} to \textit{poly} multiplied by \textit{x}^{\textit{n}}. We require that \textit{n} is nonnegative.
\end{itemize}

void \texttt{acb\_poly\_truncate} (\texttt{acb\_poly\_t poly}, \texttt{long n})
\begin{itemize}
\item Truncates \textit{poly} to have length at most \textit{n}, i.e. degree strictly smaller than \textit{n}.
\end{itemize}

3.8.4 Input and output

void \texttt{acb\_poly\_printd} (\texttt{const acb\_poly\_t poly}, \texttt{long digits})
\begin{itemize}
\item Prints the polynomial as an array of coefficients, printing each coefficient using \texttt{arb\_printd}.
\end{itemize}

3.8.5 Random generation

void \texttt{acb\_poly\_randtest} (\texttt{acb\_poly\_t poly}, \texttt{flint\_rand\_t state}, \texttt{long len}, \texttt{long prec}, \texttt{long mag\_bits})
\begin{itemize}
\item Creates a random polynomial with length at most \textit{len}.
\end{itemize}

3.8.6 Comparisons

int \texttt{acb\_poly\_equal} (\texttt{const acb\_poly\_t A}, \texttt{const acb\_poly\_t B})
\begin{itemize}
\item Returns nonzero iff \textit{A} and \textit{B} are identical as interval polynomials.
\end{itemize}

int \texttt{acb\_poly\_contains} (\texttt{const acb\_poly\_t poly1}, \texttt{const acb\_poly\_t poly2})

int \texttt{acb\_poly\_contains\_fmpz\_poly} (\texttt{const acb\_poly\_t poly1}, \texttt{const fmpz\_poly\_t poly2})
\begin{itemize}
\item Returns nonzero iff \textit{poly2} is contained in \textit{poly1}.
\end{itemize}

int \_\texttt{acb\_poly\_overlaps} (\texttt{acb\_srcptr poly1}, \texttt{long len1}, \texttt{acb\_srcptr poly2}, \texttt{long len2})

int \texttt{acb\_poly\_overlaps} (\texttt{const acb\_poly\_t poly1}, \texttt{const acb\_poly\_t poly2})
\begin{itemize}
\item Returns nonzero iff \textit{poly1} overlaps with \textit{poly2}. The underscore function requires that \textit{len1} is at least as large as \textit{len2}.
\end{itemize}

int \texttt{acb\_poly\_get\_unique\_fmpz\_poly} (\texttt{fmpz\_poly\_t z}, \texttt{const acb\_poly\_t x})
\begin{itemize}
\item If \textit{x} contains a unique integer polynomial, sets \textit{z} to that value and returns nonzero. Otherwise (if \textit{x} represents no integers or more than one integer), returns zero, possibly partially modifying \textit{z}.
\end{itemize}

int \texttt{acb\_poly\_is\_real} (\texttt{const acb\_poly\_t poly})
\begin{itemize}
\item Returns nonzero iff all coefficients in \textit{poly} have zero imaginary part.
3.8.7 Conversions

void \texttt{acb\_poly\_set\_fmpz\_poly} (\texttt{acb\_poly\_t poly}, \texttt{const fmpz\_poly\_t re}, \texttt{slong prec})

void \texttt{acb\_poly\_set2\_fmpz\_poly} (\texttt{acb\_poly\_t poly}, \texttt{const fmpz\_poly\_t re}, \texttt{const fmpz\_poly\_t im}, \texttt{slong prec})

void \texttt{acb\_poly\_set\_arb\_poly} (\texttt{acb\_poly\_t poly}, \texttt{const arb\_poly\_t re})

void \texttt{acb\_poly\_set2\_arb\_poly} (\texttt{acb\_poly\_t poly}, \texttt{const arb\_poly\_t re}, \texttt{const arb\_poly\_t im})

void \texttt{acb\_poly\_set\_fmpq\_poly} (\texttt{acb\_poly\_t poly}, \texttt{const fmpq\_poly\_t re}, \texttt{slong prec})

void \texttt{acb\_poly\_set2\_fmpq\_poly} (\texttt{acb\_poly\_t poly}, \texttt{const fmpq\_poly\_t re}, \texttt{const fmpq\_poly\_t im}, \texttt{slong prec})

Sets \textit{poly} to the given real part \textit{re} plus the imaginary part \textit{im}, both rounded to \textit{prec} bits.

void \texttt{acb\_poly\_set\_acb} (\texttt{acb\_poly\_t poly}, \texttt{slong src})

void \texttt{acb\_poly\_set\_si} (\texttt{acb\_poly\_t poly}, \texttt{slong src})

Sets \textit{poly} to \textit{src}.

3.8.8 Bounds

void \texttt{acb\_poly\_majorant} (\texttt{arb\_ptr res}, \texttt{acb\_srcptr poly}, \texttt{slong len}, \texttt{slong prec})

void \texttt{acb\_poly\_majorant} (\texttt{arb\_poly\_t res}, \texttt{acb\_poly\_t poly}, \texttt{slong prec})

Sets \textit{res} to an exact real polynomial whose coefficients are upper bounds for the absolute values of the coefficients in \textit{poly}, rounded to \textit{prec} bits.

3.8.9 Arithmetic

void \texttt{acb\_poly\_add} (\texttt{acb\_ptr C}, \texttt{acb\_srcptr A}, \texttt{slong lenA}, \texttt{acb\_srcptr B}, \texttt{slong lenB}, \texttt{slong prec})

Sets \{\textit{C}, max(lenA, lenB)\} to the sum of \{\textit{A}, lenA\} and \{\textit{B}, lenB\}. Allows aliasing of the input and output operands.

void \texttt{acb\_poly\_add} (\texttt{acb\_poly\_t C}, \texttt{const acb\_poly\_t A}, \texttt{const acb\_poly\_t B}, \texttt{slong prec})

void \texttt{acb\_poly\_add\_si} (\texttt{acb\_poly\_t C}, \texttt{const acb\_poly\_t A}, \texttt{slong B}, \texttt{slong prec})

Sets \textit{C} to the sum of \textit{A} and \textit{B}.

void \texttt{acb\_poly\_sub} (\texttt{acb\_ptr C}, \texttt{acb\_srcptr A}, \texttt{slong lenA}, \texttt{acb\_srcptr B}, \texttt{slong lenB}, \texttt{slong prec})

Sets \{\textit{C}, max(lenA, lenB)\} to the difference of \{\textit{A}, lenA\} and \{\textit{B}, lenB\}. Allows aliasing of the input and output operands.

void \texttt{acb\_poly\_sub} (\texttt{acb\_poly\_t C}, \texttt{const acb\_poly\_t A}, \texttt{const acb\_poly\_t B}, \texttt{slong prec})

Sets \textit{C} to the difference of \textit{A} and \textit{B}.

void \texttt{acb\_poly\_neg} (\texttt{acb\_poly\_t C}, \texttt{const acb\_poly\_t A})

Sets \textit{C} to the negation of \textit{A}.

void \texttt{acb\_poly\_scalar\_mul\_2exp\_si} (\texttt{acb\_poly\_t C}, \texttt{const acb\_poly\_t A}, \texttt{slong c})

Sets \textit{C} to \textit{A} multiplied by \textit{2}^\textit{c}.

void \texttt{acb\_poly\_mullow\_classical} (\texttt{acb\_ptr C}, \texttt{acb\_srcptr A}, \texttt{slong lenA}, \texttt{acb\_srcptr B}, \texttt{slong lenB}, \texttt{slong n}, \texttt{slong prec})

void \texttt{acb\_poly\_mullow\_transpose} (\texttt{acb\_ptr C}, \texttt{acb\_srcptr A}, \texttt{slong lenA}, \texttt{acb\_srcptr B}, \texttt{slong lenB}, \texttt{slong n}, \texttt{slong prec})

void \texttt{acb\_poly\_mullow\_transpose\_gauss} (\texttt{acb\_ptr C}, \texttt{acb\_srcptr A}, \texttt{slong lenA}, \texttt{acb\_srcptr B}, \texttt{slong lenB}, \texttt{slong n}, \texttt{slong prec})
void _acb_poly_mullow(acb_ptr C, acb_srcptr A, slong lenA, acb_srcptr B, slong lenB, slong n, slong prec)
Sets \( C \) to the product of \( A \), truncated to length \( n \). The output is not allowed to be aliased with either of the inputs. We require \( \text{lenA} \geq \text{lenB} > 0 \), \( n > 0 \), \( \text{lenA} + \text{lenB} - 1 \geq n \).

The \textit{classical} version uses a plain loop.

The \textit{transpose} version evaluates the product using four real polynomial multiplications (via \_arb_poly_mullow()).

The \textit{transpose\_gauss} version evaluates the product using three real polynomial multiplications. This is almost always faster than transpose, but has worse numerical stability when the coefficients vary in magnitude.

The default function _acb_poly_mullow() automatically switches been classical and transpose multiplication.

If the input pointers are identical (and the lengths are the same), they are assumed to represent the same polynomial, and its square is computed.

void acb_poly_mullow_classical(acb_poly_t C, const acb_poly_t A, const acb_poly_t B, slong n, slong prec)
void acb_poly_mullow_transpose(acb_poly_t C, const acb_poly_t A, const acb_poly_t B, slong n, slong prec)
void acb_poly_mullow_transpose_gauss(acb_poly_t C, const acb_poly_t A, const acb_poly_t B, slong n, slong prec)
void acb_poly_mullow(acb_poly_t C, const acb_poly_t A, const acb_poly_t B, slong n, slong prec)

Sets \( C \) to the product of \( A \) and \( B \), truncated to length \( n \). If the same variable is passed for \( A \) and \( B \), sets \( C \) to the square of \( A \) truncated to length \( n \).

void _acb_poly_mul(acb_ptr C, acb_srcptr A, slong lenA, acb_srcptr B, slong lenB, slong prec)
Sets \( C \), truncated to length \( n \). The output is not allowed to be aliased with either of the inputs. We require \( \text{lenA} \geq \text{lenB} > 0 \). This function is implemented as a simple wrapper for _acb_poly_mullow().

If the input pointers are identical (and the lengths are the same), they are assumed to represent the same polynomial, and its square is computed.

void acb_poly_mul(acb_poly_t C, const acb_poly_t A1, const acb_poly_t B2, slong prec)
Sets \( C \) to the product of \( A \) and \( B \). If the same variable is passed for \( A \) and \( B \), sets \( C \) to the square of \( A \) truncated to length \( n \).

void _acb_poly_inv_series(acb_ptr Qinv, acb_srcptr Q, slong Qlen, slong len, slong prec)
Sets \( Qinv \) to the power series inverse of \( Q \), truncated to length \( n \). Uses Newton iteration.

void acb_poly_inv_series(acb_poly_t Qinv, const acb_poly_t Q, slong n, slong prec)
Sets \( Qinv \) to the power series inverse of \( Q \).

void _acb_poly_div_series(acb_ptr Q, acb_srcptr A, slong Alen, acb_srcptr B, slong Blen, slong n, slong prec)
Sets \( Q \) to the power series quotient of \( A \), truncated to length \( n \). Uses Newton iteration followed by multiplication.

void acb_poly_div_series(acb_poly_t Q, const acb_poly_t A, const acb_poly_t B, slong n, slong prec)
Sets \( Q \) to the power series quotient \( A \) divided by \( B \), truncated to length \( n \).

void _acb_poly_div(acb_ptr Q, acb_srcptr A, slong lenA, acb_srcptr B, slong lenB, slong prec)
void _acb_poly_rem(acb_ptr R, acb_srcptr A, slong lenA, acb_srcptr B, slong lenB, slong prec)
void _acb_poly_divrem(acb_ptr Q, acb_ptr R, acb_srcptr A, slong lenA, acb_srcptr B, slong lenB, slong prec)
void `acb_poly_divrem` *(acb_poly_t Q, acb_poly_t R, const acb_poly_t A, const acb_poly_t B, slong prec)*

Performs polynomial division with remainder, computing a quotient Q and a remainder R such that \( A = BQ + R \). The implementation reverses the inputs and performs power series division.

If the leading coefficient of B contains zero (or if B is identically zero), returns 0 indicating failure without modifying the outputs. Otherwise returns nonzero.

void `_acb_poly_div_root` *(acb_ptr Q, acb_t R, acb_srcptr A, slong len, const acb_t c, slong prec)*

Divides A by the polynomial \( x - c \), computing the quotient Q as well as the remainder \( R = f(c) \).

### 3.8.10 Composition

void `_acb_poly_compose_horner` *(acb_ptr res, acb_srcptr poly1, slong len1, acb_srcptr poly2, slong len2, slong prec)*

void `acb_poly_compose_horner` *(acb_poly_t res, const acb_poly_t poly1, const acb_poly_t poly2, slong prec)*

void `_acb_poly_compose_divconquer` *(acb_ptr res, acb_srcptr poly1, slong len1, acb_srcptr poly2, slong len2, slong prec)*

void `acb_poly_compose_divconquer` *(acb_poly_t res, const acb_poly_t poly1, const acb_poly_t poly2, slong prec)*

void `acb_poly_compose` *(acb_ptr res, acb_srcptr poly1, slong len1, acb_srcptr poly2, slong len2, slong prec)*

void `acb_poly_compose` *(acb_poly_t res, const acb_poly_t poly1, const acb_poly_t poly2, slong prec)*

Sets res to the composition \( h(x) = f(g(x)) \) where \( f \) is given by poly1 and \( g \) is given by poly2, respectively using Horner’s rule, divide-and-conquer, and an automatic choice between the two algorithms. The underscore methods do not support aliasing of the output with either input polynomial.

void `_acb_poly_compose_series_horner` *(acb_ptr res, acb_srcptr poly1, slong len1, acb_srcptr poly2, slong len2, slong n, slong prec)*

void `acb_poly_compose_series_horner` *(acb_poly_t res, const acb_poly_t poly1, const acb_poly_t poly2, slong n, slong prec)*

void `_acb_poly_compose_series_brent_kung` *(acb_ptr res, acb_srcptr poly1, slong len1, acb_srcptr poly2, slong len2, slong n, slong prec)*

void `acb_poly_compose_series_brent_kung` *(acb_poly_t res, const acb_poly_t poly1, const acb_poly_t poly2, slong n, slong prec)*

void `_acb_poly_compose_series` *(acb_ptr res, acb_srcptr poly1, slong len1, acb_srcptr poly2, slong len2, slong n, slong prec)*

void `acb_poly_compose_series` *(acb_poly_t res, const acb_poly_t poly1, const acb_poly_t poly2, slong n, slong prec)*

Sets res to the power series composition \( h(x) = f(g(x)) \) truncated to order \( O(x^n) \) where \( f \) is given by poly1 and \( g \) is given by poly2, respectively using Horner’s rule, the Brent-Kung baby step-giant step algorithm, and an automatic choice between the two algorithms. We require that the constant term in \( g(x) \) is exactly zero. The underscore methods do not support aliasing of the output with either input polynomial.

void `_acb_poly_revert_series_lagrange` *(acb_ptr h, acb_srcptr f, slong flen, slong n, slong prec)*

void `acb_poly_revert_series_lagrange` *(acb_poly_t h, const acb_poly_t f, slong n, slong prec)*

void `_acb_poly_revert_series_newton` *(acb_ptr h, acb_srcptr f, slong flen, slong n, slong prec)*

void `acb_poly_revert_series_newton` *(acb_poly_t h, const acb_poly_t f, slong n, slong prec)*

void `_acb_poly_revert_series_lagrange_fast` *(acb_ptr h, acb_srcptr f, slong flen, slong n, slong prec)*
void `acb_poly_revert_series_lagrange_fast` (acb_poly_t h, const acb_poly_t f, slong n, slong prec)

void `_acb_poly_revert_series` (acb_ptr h, acb_srcptr f, slong flen, slong n, slong prec)

void `acb_poly_revert_series` (acb_poly_t h, const acb_poly_t f, slong n, slong prec)

Sets h to the power series reversion of f, i.e. the expansion of the compositional inverse function \( f^{-1}(x) \), truncated to order \( O(x^n) \), using respectively Lagrange inversion, Newton iteration, fast Lagrange inversion, and a default algorithm choice.

We require that the constant term in f is exactly zero and that the linear term is nonzero. The underscore methods assume that flen is at least 2, and do not support aliasing.

3.8.11 Evaluation

void `_acb_poly_evaluate_horner` (acb_t y, acb_srcptr f, slong len, const acb_t x, slong prec)

void `acb_poly_evaluate_horner` (acb_t y, const acb_poly_t f, const acb_t x, slong prec)

void `_acb_poly_evaluate_rectangular` (acb_t y, acb_srcptr f, slong len, const acb_t x, slong prec)

void `acb_poly_evaluate_rectangular` (acb_t y, const acb_poly_t f, const acb_t x, slong prec)

void `_acb_poly_evaluate` (acb_t y, acb_srcptr f, slong len, const acb_t x, slong prec)

void `acb_poly_evaluate` (acb_t y, const acb_poly_t f, const acb_t x, slong prec)

Sets \( y = f(x) \), evaluated respectively using Horner's rule, rectangular splitting, and an automatic algorithm choice.

void `_acb_poly_evaluate2_horner` (acb_t y, acb_t z, acb_srcptr f, slong len, const acb_t x, slong prec)

void `acb_poly_evaluate2_horner` (acb_t y, acb_t z, const acb_poly_t f, const acb_t x, slong prec)

void `_acb_poly_evaluate2_rectangular` (acb_t y, acb_t z, acb_srcptr f, slong len, const acb_t x, slong prec)

void `acb_poly_evaluate2_rectangular` (acb_t y, acb_t z, const acb_poly_t f, const acb_t x, slong prec)

void `_acb_poly_evaluate2` (acb_t y, acb_t z, acb_srcptr f, slong len, const acb_t x, slong prec)

void `acb_poly_evaluate2` (acb_t y, acb_t z, const acb_poly_t f, const acb_t x, slong prec)

Sets \( y = f(x) \), \( z = f'(x) \), evaluated respectively using Horner's rule, rectangular splitting, and an automatic algorithm choice.

When Horner's rule is used, the only advantage of evaluating the function and its derivative simultaneously is that one does not have to generate the derivative polynomial explicitly. With the rectangular splitting algorithm, the powers can be reused, making simultaneous evaluation slightly faster.

3.8.12 Product trees

void `_acb_poly_product_roots` (acb_ptr poly, acb_srcptr xs, slong n, slong prec)

void `acb_poly_product_roots` (acb_poly_t poly, acb_srcptr xs, slong n, slong prec)

Generates the polynomial \( (x - x_0)(x - x_1) \cdots (x - x_{n-1}) \).

`acb_ptr * _acb_poly_tree_alloc` (slong len)

Returns an initialized data structured capable of representing a remainder tree (product tree) of len roots.

void `_acb_poly_tree_free` (acb_ptr * tree, slong len)

Deallocates a tree structure as allocated using `_acb_poly_tree_alloc`.
void _acb_poly_tree_build (acb_ptr * tree, acb_srcptr roots, slong len, slong prec)
  Constructs a product tree from a given array of len roots. The tree structure must be pre-allocated to the specified length using _acb_poly_tree_alloc().

3.8.13 Multipoint evaluation

void _acb_poly_evaluate_vec_iter (acb_ptr ys, acb_srcptr poly, slong plen, acb_srcptr xs, slong n, slong prec)
void acb_poly_evaluate_vec_iter (acb_ptr ys, const acb_poly_t poly, acb_srcptr xs, slong n, slong prec)
  Evaluates the polynomial simultaneously at n given points, calling _acb_poly_evaluate() repeatedly.

void _acb_poly_evaluate_vec_fast_precomp (acb_ptr ys, acb_srcptr poly, slong plen, acb_ptr * tree, slong len, slong prec)
void _acb_poly_evaluate_vec_fast (acb_ptr ys, acb_srcptr poly, slong plen, acb_srcptr xs, slong n, slong prec)
  Evaluates the polynomial simultaneously at n given points, using fast multipoint evaluation.

3.8.14 Interpolation

void _acb_poly_interpolate_newton (acb_ptr poly, acb_srcptr xs, acb_srcptr ys, slong n, slong prec)
void acb_poly_interpolate_newton (acb_poly_t poly, acb_srcptr xs, acb_srcptr ys, slong n, slong prec)
  Recovers the unique polynomial of length at most n that interpolates the given x and y values. This implementation first interpolates in the Newton basis and then converts back to the monomial basis.

void _acb_poly_interpolate_barycentric (acb_ptr poly, acb_srcptr xs, acb_srcptr ys, slong n, slong prec)
void acb_poly_interpolate_barycentric (acb_poly_t poly, acb_srcptr xs, acb_srcptr ys, slong n, slong prec)
  Recovers the unique polynomial of length at most n that interpolates the given x and y values. This implementation uses the barycentric form of Lagrange interpolation.

void _acb_poly_interpolation_weights (acb_ptr w, acb_ptr * tree, slong len, slong prec)
void _acb_poly_interpolate_fast_precomp (acb_ptr poly, acb_srcptr ys, acb_ptr * tree, acb_srcptr weights, slong len, slong prec)
void _acb_poly_interpolate_fast (acb_ptr poly, acb_srcptr xs, acb_srcptr ys, slong len, slong prec)
void acb_poly_interpolate_fast (acb_poly_t poly, acb_srcptr xs, acb_srcptr ys, slong n, slong prec)
  Recovers the unique polynomial of length at most n that interpolates the given x and y values, using fast Lagrange interpolation. The precomp function takes a precomputed product tree over the x values and a vector of interpolation weights as additional inputs.

3.8.15 Differentiation

void _acb_poly_derivative (acb_ptr res, acb_srcptr poly, slong len, slong prec)
  Sets {res, len - 1} to the derivative of {poly, len}. Allows aliasing of the input and output.
void \texttt{acb\_poly\_derivative} (\texttt{acb\_poly\_t res, const acb\_poly\_t poly, slong prec})
    Sets \textit{res} to the derivative of \textit{poly}.

void \texttt{acbc\_poly\_integral} (\texttt{acb\_ptr res, acb\_srcptr poly, slong len, slong prec})
    Sets \{\textit{res}, \textit{len}\} to the integral of \{\textit{poly}, \textit{len} - 1\}. Allows aliasing of the input and output.

void \texttt{acb\_poly\_integral} (\texttt{acb\_poly\_t res, const acb\_poly\_t poly, slong prec})
    Sets \textit{res} to the integral of \textit{poly}.

\subsection*{3.8.16 Elementary functions}

void \texttt{acb\_poly\_pow\_ui\_trunc\_binexp} (\texttt{acb\_ptr res, acb\_srcptr f, slong flen, ulong exp, slong len, slong prec})
    Sets \{\textit{res}, \textit{len}\} to \{\textit{f}, \textit{flen}\} raised to the power \textit{exp}, truncated to length \textit{len}. Requires that \textit{len} is no longer than the length of the power as computed without truncation (i.e. no zero-padding is performed). Does not support aliasing of the input and output, and requires that \textit{flen} and \textit{len} are positive. Uses binary exponentiation.

void \texttt{acb\_poly\_pow\_ui\_trunc\_binexp} (\texttt{acb\_poly\_t res, const acb\_poly\_t poly, ulong exp, slong len, slong prec})
    Sets \textit{res} to \textit{poly} raised to the power \textit{exp}, truncated to length \textit{len}. Uses binary exponentiation.

void \texttt{acb\_poly\_pow\_ui} (\texttt{acb\_ptr res, acb\_srcptr f, slong flen, ulong exp, slong prec})
    Sets \textit{res} to \{\textit{f}, \textit{flen}\} raised to the power \textit{exp}. Does not support aliasing of the input and output, and requires that \textit{flen} is positive.

void \texttt{acb\_poly\_pow\_ui} (\texttt{acb\_poly\_t res, const acb\_poly\_t poly, ulong exp, slong prec})
    Sets \textit{res} to \textit{poly} raised to the power \textit{exp}.

void \texttt{acb\_poly\_pow\_series} (\texttt{acb\_ptr h, acb\_srcptr f, slong flen, acb\_srcptr g, slong glen, slong len, slong prec})
    Sets \{\textit{h}, \textit{len}\} to the power series \( f(x)^{g(x)} = \exp(g(x) \log f(x)) \) truncated to length \textit{len}. This function detects special cases such as \textit{g} being an exact small integer or \( \pm 1/2 \), and computes such powers more efficiently. This function does not support aliasing of the output with either of the input operands. It requires that all lengths are positive, and assumes that \textit{flen} and \textit{glen} do not exceed \textit{len}.

void \texttt{acb\_poly\_pow\_series} (\texttt{acb\_poly\_t h, const acb\_poly\_t f, const acb\_poly\_t g, slong len, slong prec})
    Sets \textit{h} to the power series \( f(x)^{g(x)} = \exp(g(x) \log f(x)) \) truncated to length \textit{len}. This function detects special cases such as \textit{g} being an exact small integer or \( \pm 1/2 \), and computes such powers more efficiently. This function does not support aliasing of the output with either of the input operands. It requires that all lengths are positive, and assumes that \textit{flen} does not exceed \textit{len}.

void \texttt{acb\_poly\_pow\_acb\_series} (\texttt{acb\_ptr h, acb\_srcptr f, slong flen, const acb\_t g, slong len, slong prec})
    Sets \{\textit{h}, \textit{len}\} to the power series \( f(x)^{g} = \exp(g \log f(x)) \) truncated to length \textit{len}. This function detects special cases such as \textit{g} being an exact small integer or \( \pm 1/2 \), and computes such powers more efficiently. This function does not support aliasing of the output with either of the input operands. It requires that all lengths are positive, and assumes that \textit{flen} does not exceed \textit{len}.

void \texttt{acb\_poly\_pow\_acb\_series} (\texttt{acb\_poly\_t h, const acb\_poly\_t f, const acb\_t g, slong len, slong prec})
    Sets \textit{h} to the power series \( f(x)^{g} = \exp(g \log f(x)) \) truncated to length \textit{len}.

void \texttt{acb\_poly\_sqrt\_series} (\texttt{acb\_ptr g, acb\_srcptr h, slong hlen, slong n, slong prec})

void \texttt{acb\_poly\_rsqrt\_series} (\texttt{acb\_ptr g, acb\_srcptr h, slong hlen, slong n, slong prec})
void \texttt{acb\_poly\_rsqrt\_series} (\texttt{acb\_poly\_t} \texttt{g}, \texttt{const} \texttt{acb\_poly\_t} \texttt{h}, \texttt{slong} \texttt{n}, \texttt{slong} \texttt{prec})

Sets \(g\) to the reciprocal power series square root of \(h\), truncated to length \(n\). Uses division-free Newton iteration.

The underscore method does not support aliasing of the input and output arrays. It requires that \(hlen\) and \(n\) are greater than zero.

void \texttt{acb\_poly\_log\_series} (\texttt{acb\_ptr} \texttt{res}, \texttt{acb\_srcptr} \texttt{f}, \texttt{slong} \texttt{flen}, \texttt{slong} \texttt{n}, \texttt{slong} \texttt{prec})

void \texttt{acb\_poly\_log\_series} (\texttt{acb\_poly\_t} \texttt{res}, \texttt{const} \texttt{acb\_poly\_t} \texttt{f}, \texttt{slong} \texttt{n}, \texttt{slong} \texttt{prec})

Sets \(res\) to the power series logarithm of \(f\), truncated to length \(n\). Uses the formula \(\log(f(x)) = \int f'(x)/f(x)dx\), adding the logarithm of the constant term in \(f\) as the constant of integration.

The underscore method supports aliasing of the input and output arrays. It requires that \(flen\) and \(n\) are greater than zero.

void \texttt{acb\_poly\_atan\_series} (\texttt{acb\_ptr} \texttt{res}, \texttt{acb\_srcptr} \texttt{f}, \texttt{slong} \texttt{flen}, \texttt{slong} \texttt{n}, \texttt{slong} \texttt{prec})

void \texttt{acb\_poly\_atan\_series} (\texttt{acb\_poly\_t} \texttt{res}, \texttt{const} \texttt{acb\_poly\_t} \texttt{f}, \texttt{slong} \texttt{n}, \texttt{slong} \texttt{prec})

Sets \(res\) the power series inverse tangent of \(f\), truncated to length \(n\).

Uses the formula

\[
\tan^{-1}(f(x)) = \int f'(x)/(1 + f(x)^2)dx,
\]

adding the function of the constant term in \(f\) as the constant of integration.

The underscore method supports aliasing of the input and output arrays. It requires that \(flen\) and \(n\) are greater than zero.

void \texttt{acb\_poly\_exp\_series\_basecase} (\texttt{acb\_ptr} \texttt{f}, \texttt{acb\_srcptr} \texttt{h}, \texttt{slong} \texttt{hlen}, \texttt{slong} \texttt{n}, \texttt{slong} \texttt{prec})

void \texttt{acb\_poly\_exp\_series\_basecase} (\texttt{acb\_poly\_t} \texttt{f}, \texttt{const} \texttt{acb\_poly\_t} \texttt{h}, \texttt{slong} \texttt{n}, \texttt{slong} \texttt{prec})

void \texttt{acb\_poly\_exp\_series} (\texttt{acb\_ptr} \texttt{f}, \texttt{acb\_srcptr} \texttt{h}, \texttt{slong} \texttt{hlen}, \texttt{slong} \texttt{n}, \texttt{slong} \texttt{prec})

void \texttt{acb\_poly\_exp\_series} (\texttt{acb\_poly\_t} \texttt{f}, \texttt{const} \texttt{acb\_poly\_t} \texttt{h}, \texttt{slong} \texttt{n}, \texttt{slong} \texttt{prec})

Sets \(f\) to the power series exponential of \(h\), truncated to length \(n\).

The basecase version uses a simple recurrence for the coefficients, requiring \(O(mn)\) operations where \(m\) is the length of \(h\).

The main implementation uses Newton iteration, starting from a small number of terms given by the basecase algorithm. The complexity is \(O(M(n))\). Redundant operations in the Newton iteration are avoided by using the scheme described in [HZZ2004].

The underscore methods support aliasing and allow the input to be shorter than the output, but require the lengths to be nonzero.

void \texttt{acb\_poly\_sin\_cos\_series\_basecase} (\texttt{acb\_ptr} \texttt{s}, \texttt{acb\_ptr} \texttt{c}, \texttt{acb\_srcptr} \texttt{h}, \texttt{slong} \texttt{hlen}, \texttt{slong} \texttt{n}, \texttt{slong} \texttt{prec}, \texttt{int} \texttt{times\_pi})

void \texttt{acb\_poly\_sin\_cos\_series\_basecase} (\texttt{acb\_poly\_t} \texttt{s}, \texttt{acb\_poly\_t} \texttt{c}, \texttt{const} \texttt{acb\_poly\_t} \texttt{h}, \texttt{slong} \texttt{n}, \texttt{slong} \texttt{prec}, \texttt{int} \texttt{times\_pi})

void \texttt{acb\_poly\_sin\_cos\_series\_tangent} (\texttt{acb\_ptr} \texttt{s}, \texttt{acb\_ptr} \texttt{c}, \texttt{acb\_srcptr} \texttt{h}, \texttt{slong} \texttt{hlen}, \texttt{slong} \texttt{n}, \texttt{slong} \texttt{prec}, \texttt{int} \texttt{times\_pi})

void \texttt{acb\_poly\_sin\_cos\_series\_tangent} (\texttt{acb\_poly\_t} \texttt{s}, \texttt{acb\_poly\_t} \texttt{c}, \texttt{const} \texttt{acb\_poly\_t} \texttt{h}, \texttt{slong} \texttt{n}, \texttt{slong} \texttt{prec}, \texttt{int} \texttt{times\_pi})

void \texttt{acb\_poly\_sin\_cos\_series} (\texttt{acb\_ptr} \texttt{s}, \texttt{acb\_ptr} \texttt{c}, \texttt{acb\_srcptr} \texttt{h}, \texttt{slong} \texttt{hlen}, \texttt{slong} \texttt{n}, \texttt{slong} \texttt{prec})

void \texttt{acb\_poly\_sin\_cos\_series} (\texttt{acb\_poly\_t} \texttt{s}, \texttt{acb\_poly\_t} \texttt{c}, \texttt{const} \texttt{acb\_poly\_t} \texttt{h}, \texttt{slong} \texttt{n}, \texttt{slong} \texttt{prec})

Sets \(s\) and \(c\) to the power series sine and cosine of \(h\), computed simultaneously.
The `basecase` version uses a simple recurrence for the coefficients, requiring $O(nm)$ operations where $m$ is the length of $h$.

The `tangent` version uses the tangent half-angle formulas to compute the sine and cosine via `_acb_poly_tan_series()`. This requires $O(M(n))$ operations. When $h = h_0 + h_1$ where the constant term $h_0$ is nonzero, the evaluation is done as $\sin(h_0 + h_1) = \cos(h_0)\sin(h_1) + \sin(h_0)\cos(h_1)$, $\cos(h_0 + h_1) = \cos(h_0)\cos(h_1) - \sin(h_0)\sin(h_1)$, to improve accuracy and avoid dividing by zero at the poles of the tangent function.

The default version automatically selects between the `basecase` and `tangent` algorithms depending on the input.

The `basecase` and `tangent` versions take a flag `times_pi` specifying that the input is to be multiplied by $\pi$.

The underscore methods support aliasing and require the lengths to be nonzero.

```c
void _acb_poly_sin_series (acb_ptr s, acb_srcptr h, slong hlen, slong n, slong prec)
void acb_poly_sin_series (acb_poly_t s, const acb_poly_t h, slong n, slong prec)
void _acb_poly_cos_series (acb_ptr c, acb_srcptr h, slong hlen, slong n, slong prec)
void acb_poly_cos_series (acb_poly_t c, const acb_poly_t h, slong n, slong prec)

void _acb_poly_sin_cos_series (acb_ptr c, acb_ptr s, slong hlen, slong n, slong prec)
Respectively evaluates the power series sine or cosine. These functions simply wrap _acb_poly_sin_cos_series(). The underscore methods support aliasing and require the lengths to be nonzero.

void _acb_poly_tan_series (acb_ptr g, acb_srcptr h, slong hlen, slong len, slong prec)
void acb_poly_tan_series (acb_poly_t g, const acb_poly_t h, slong n, slong prec)

void _acb_poly_sin_cos_pi_series (acb_ptr s, acb_ptr c, acb_srcptr h, slong hlen, slong n, slong prec)
void acb_poly_sin_cos_pi_series (acb_poly_t s, acb_poly_t c, const acb_poly_t h, slong n, slong prec)
void _acb_poly_sin_pi_series (acb_ptr s, acb_srcptr h, slong hlen, slong n, slong prec)
void acb_poly_sin_pi_series (acb_poly_t s, const acb_poly_t h, slong n, slong prec)
void _acb_poly_cos_pi_series (acb_ptr c, acb_srcptr h, slong hlen, slong n, slong prec)
void acb_poly_cos_pi_series (acb_poly_t c, const acb_poly_t h, slong n, slong prec)
void _acb_poly_tan_pi_series (acb_ptr c, acb_srcptr h, slong hlen, slong n, slong prec)
void acb_poly_tan_pi_series (acb_poly_t c, const acb_poly_t h, slong n, slong prec)
void _acb_poly_sin_cos_pi_series (acb_ptr s, acb_ptr c, acb_srcptr h, slong hlen, slong n, slong prec)

Compute the respective trigonometric functions of the input multiplied by $\pi$.

### 3.8.17 Gamma function

```c
void _acb_poly_gamma_series (acb_ptr res, acb_srcptr h, slong hlen, slong n, slong prec)
void acb_poly_gamma_series (acb_poly_t res, const acb_poly_t h, slong n, slong prec)
void _acb_poly_rgamma_series (acb_ptr res, acb_srcptr h, slong hlen, slong n, slong prec)
void acb_poly_rgamma_series (acb_poly_t res, const acb_poly_t h, slong n, slong prec)
```
void _acb_poly_lgamma_series (acb_ptr res, acb_srcptr h, slong hlen, slong n, slong prec)
void acb_poly_lgamma_series (acb_poly_t res, const acb_poly_t h, slong n, slong prec)
void _acb_poly_digamma_series (acb_ptr res, acb_srcptr h, slong hlen, slong n, slong prec)
void acb_poly_digamma_series (acb_poly_t res, const acb_poly_t h, slong n, slong prec)

Sets res to the series expansion of $\Gamma(h(x))$, $1/\Gamma(h(x))$, or $\log \Gamma(h(x))$, $\psi(h(x))$, truncated to length n.

These functions first generate the Taylor series at the constant term of h, and then call _acb_poly_compose_series(). The Taylor coefficients are generated using Stirling’s series.

The underscore methods support aliasing of the input and output arrays, and require that hlen and n are greater than zero.

void _acb_poly_rising_ui_series (acb_ptr res, acb_srcptr f, slong flen, ulong r, slong trunc, slong prec)
void acb_poly_rising_ui_series (acb_poly_t res, const acb_poly_t f, ulong r, slong trunc, slong prec)

Sets res to the rising factorial $(f)(f+1)(f+2) \cdots (f+r-1)$, truncated to length trunc. The underscore method assumes that flen, r and trunc are at least 1, and does not support aliasing. Uses binary splitting.

### 3.8.18 Power sums

void _acb_poly_powsum_series_naive (acb_ptr z, const acb_t s, const acb_t a, const acb_t q, slong n, slong len, slong prec)
void _acb_poly_powsum_series_naive_threaded (acb_ptr z, const acb_t s, const acb_t a, const acb_t q, slong n, slong len, slong prec)

Computes

$$z = S(s, a, n) = \sum_{k=0}^{n-1} \frac{q^k}{(k+a)^{s+t}}$$

as a power series in $t$ truncated to length len. This function evaluates the sum naively term by term. The threaded version splits the computation over the number of threads returned by flint_get_num_threads().

void _acb_poly_powsum_one_series_sieved (acb_ptr z, const acb_t s, slong n, slong len, slong prec)

Computes

$$z = S(s, 1, n) \sum_{k=1}^{n} \frac{1}{k^s+t}$$

as a power series in $t$ truncated to length len. This function stores a table of powers that have already been calculated, computing $(ij)^r$ as $i^r j^r$ whenever $k = ij$ is composite. As a further optimization, it groups all even $k$ and evaluates the sum as a polynomial in $2^{(s+t)}$. This scheme requires about $n/\log n$ powers, $n/2$ multiplications, and temporary storage of $n/6$ power series. Due to the extra power series multiplications, it is only faster than the naive algorithm when len is small.

### 3.8.19 Zeta function

void _acb_poly_zeta_em_choose_param (mag_t bound, ulong * N, ulong * M, const acb_t s, const acb_t a, slong d, slong target, slong prec)

Chooses N and M for Euler-Maclaurin summation of the Hurwitz zeta function, using a default algorithm.

void _acb_poly_zeta_em_bound1 (mag_t bound, const acb_t s, const acb_t a, slong N, slong M, slong d, slong wp)
void _acb_poly_zeta_em_bound (arb_ptr vec, const acb_t s, const acb_t a, ulong N, ulong M, slong d, slong wp)
  Compute bounds for Euler-Maclaurin evaluation of the Hurwitz zeta function or its power series, using the formulas in [Joh2013].

void _acb_poly_zeta_em_tail_naive (acb_ptr z, const acb_t s, const acb_t Na, acb_srcptr Nasx, long M, slong len, slong prec)
  Evaluates the tail in the Euler-Maclaurin sum for the Hurwitz zeta function, respectively using the naive recurrence and binary splitting.

void _acb_poly_zeta_em_tail_bsplit (acb_ptr z, const acb_t s, const acb_t Na, acb_srcptr Nasx, slong M, slong len, slong prec)

void _acb_poly_zeta_em_sum (acb_ptr z, const acb_t s, const acb_t a, int deflate, ulong N, ulong M, slong d, slong prec)
  Evaluates the truncated Euler-Maclaurin sum of order N, M for the length-d truncated Taylor series of the Hurwitz zeta function $\zeta(s,a)$ at s, using a working precision of prec bits. With $a = 1$, this gives the usual Riemann zeta function.

If deflate is nonzero, $\zeta(s,a) - 1/(s - 1)$ is evaluated (which permits series expansion at $s = 1$).

void _acb_poly_zeta_cpx_series (acb_ptr z, const acb_t s, const acb_t a, int deflate, slong d, slong prec)
  Computes the series expansion of $\zeta(s + x, a)$ (or $\zeta(s + x, a) - 1/(s + x - 1)$ if deflate is nonzero) to order $d$. This function wraps _acb_poly_zeta_em_sum(), automatically choosing default values for $N, M$ using _acb_poly_zeta_em_choose_param() to target an absolute truncation error of $2^{-\text{prec}}$.

void _acb_poly_zeta_series (acb_ptr res, acb_srcptr h, slong hlen, const acb_t a, int deflate, slong len, slong prec)
  Sets res to the Hurwitz zeta function $\zeta(s,a)$ where $s$ a power series and $a$ is a constant, truncated to length $n$. To evaluate the usual Riemann zeta function, set $a = 1$.

If deflate is nonzero, evaluates $\zeta(s,a) + 1/(1-s)$, which is well-defined as a limit when the constant term of $s$ is 1. In particular, expanding $\zeta(s,a) + 1/(1-s)$ with $s = 1 + x$ gives the Stieltjes constants

$$\sum_{k=0}^{n-1} \frac{(-1)^k}{k!} \gamma_k(a)x^k.$$

If $a = 1$, this implementation uses the reflection formula if the midpoint of the constant term of $s$ is negative.

### 3.8.20 Other special functions

void _acb_poly_polylog_cpx_small (acb_ptr w, const acb_t s, const acb_t z, slong len, slong prec)

void _acb_poly_polylog_cpx_zeta (acb_ptr w, const acb_t s, const acb_t z, slong len, slong prec)

void _acb_poly_polylog_cpx (acb_ptr w, const acb_t s, const acb_t z, slong len, slong prec)
  Sets w to the Taylor series with respect to x of the polylogarithm $\text{Li}_{s+x}(z)$, where s and z are given complex constants. The output is computed to length len which must be positive. Aliasing between w and s or z is not permitted.

The small version uses the standard power series expansion with respect to z, convergent when $|z| < 1$. The zeta version evaluates the polylogarithm as a sum of two Hurwitz zeta functions. The default version automatically delegates to the small version when z is close to zero, and the zeta version otherwise. For further details, see Algorithms for polylogarithms.
void \_acb\_poly\_polylog\_series (acb\_ptr w, acb\_srcptr s, slong s\_len, const acb\_t z, slong l\_en, slong p\_rec)

void \acb\_poly\_polylog\_series (acb\_poly\_t w, const acb\_poly\_t s, const acb\_t z, slong l\_en, slong p\_rec)

Sets w to the polylogarithm \(Li_i(z)\) where s is a given power series, truncating the output to length len. The underscore method requires all lengths to be positive and supports aliasing between all inputs and outputs.

void \_acb\_poly\_erf\_series (acb\_ptr res, acb\_srcptr z, slong z\_len, slong n, slong p\_rec)

void \acb\_poly\_erf\_series (acb\_poly\_t res, const acb\_poly\_t z, slong n, slong p\_rec)

Sets res to the error function of the power series z, truncated to length n.

void \_acb\_poly\_upper\_gamma\_series (acb\_ptr res, acb\_t s, acb\_srcptr z, slong z\_len, slong n, slong p\_rec)

void \acb\_poly\_upper\_gamma\_series (acb\_poly\_t res, const acb\_t s, const acb\_poly\_t z, slong n, slong p\_rec)

Sets res to the upper incomplete gamma function \(\Gamma(s, z)\) where s is a constant and z is a power series, truncated to length n.

void \_acb\_poly\_agm1\_series (acb\_ptr res, acb\_srcptr z, slong z\_len, slong len, slong p\_rec)

void \acb\_poly\_agm1\_series (acb\_poly\_t res, const acb\_poly\_t z, slong n, slong p\_rec)

Sets res to the arithmetic-geometric mean of 1 and the power series z, truncated to length n.

void \_acb\_poly\_elliptic\_k\_series (acb\_ptr res, acb\_srcptr z, slong z\_len, slong len, slong p\_rec)

void \acb\_poly\_elliptic\_k\_series (acb\_poly\_t res, const acb\_poly\_t z, slong n, slong p\_rec)

Sets res to the complete elliptic integral of the first kind of the power series z, truncated to length n.

void \_acb\_poly\_elliptic\_p\_series (acb\_ptr res, acb\_srcptr z, slong z\_len, const acb\_t tau, slong len, slong p\_rec)

void \acb\_poly\_elliptic\_p\_series (acb\_poly\_t res, const acb\_poly\_t z, const acb\_t tau, slong n, slong p\_rec)

Sets res to the Weierstrass elliptic function of the power series z, with periods 1 and tau, truncated to length n.

### 3.8.21 Root-finding

void \_acb\_poly\_root\_bound\_fujiwara (mag\_t bound, acb\_srcptr poly, slong len)

void \acb\_poly\_root\_bound\_fujiwara (mag\_t bound, acb\_poly\_t poly)

Sets bound to an upper bound for the magnitude of all the complex roots of poly. Uses Fujiwara’s bound

\[
2 \max \left\{ \left| \frac{a_{n-1}}{a_n} \right|, \left| \frac{a_{n-2}}{a_n} \right|^{1/2}, \ldots, \left| \frac{a_1}{a_n} \right|^{1/(n-1)}, \left| \frac{a_0}{2a_n} \right|^{1/n} \right\}
\]

where \(a_0, \ldots, a_n\) are the coefficients of poly.

void \_acb\_poly\_root\_inclusion (acb\_t r, const acb\_t m, acb\_srcptr poly, acb\_srcptr polyder, slong len, slong p\_rec)

Given any complex number m, and a nonconstant polynomial \(f\) and its derivative \(f'\), sets r to a complex interval centered on \(m\) that is guaranteed to contain at least one root of \(f\). Such an interval is obtained by taking a ball of radius \(|f(m)/f'(m)|n\) where \(n\) is the degree of \(f\). Proof: assume that the distance to the nearest root exceeds \(r = |f(m)/f'(m)|n\). Then

\[
\left| \frac{f'(m)}{f(m)} \right| = \left| \sum_i \frac{1}{m - \zeta_i} \right| \leq \sum_i \left| \frac{1}{|m - \zeta_i|} \right| < \frac{n}{r} = \left| \frac{f'(m)}{f(m)} \right|
\]

which is a contradiction (see [Kob2010]).
Given a list of approximate roots of the input polynomial, this function sets a rigorous bounding interval for each root, and determines which roots are isolated from all the other roots. It then rearranges the list of roots so that the isolated roots are at the front of the list, and returns the count of isolated roots.

If the return value equals the degree of the polynomial, then all roots have been found. If the return value is smaller, all the remaining output intervals are guaranteed to contain roots, but it is possible that not all of the polynomial’s roots are contained among them.

Refines the given roots simultaneously using a single iteration of the Durand-Kerner method. The radius of each root is set to an approximation of the correction, giving a rough estimate of its error (not a rigorous bound).

Attempts to compute all the roots of the given nonzero polynomial \( p(x) \) using a working precision of \( \text{prec} \) bits. If \( n \) denotes the degree of \( p(x) \), the function writes \( n \) approximate roots with rigorous error bounds to the preallocated array \( \text{roots} \), and returns the number of roots that are isolated.

The roots are computed numerically by performing several steps with the Durand-Kerner method and terminating if the estimated accuracy of the roots approaches the working precision or if the number of steps exceeds \( \text{maxiter} \), which can be set to zero in order to use a default value. Finally, the approximate roots are validated rigorously.

Initial values for the iteration can be provided as the array \( \text{initial} \). If \( \text{initial} \) is set to NULL, default values \((0.4 + 0.9i)^k\) are used.

The polynomial is assumed to be squarefree. If there are repeated roots, the iteration is likely to find them (with low numerical accuracy), but the error bounds will not converge as the precision increases.

Given a strictly real polynomial \( p(x) \) (of length \( \text{len} \)) and isolating intervals for all its complex roots, determines if all the real roots are separated from the non-real roots. If this function returns nonzero, every root enclosure that touches the real axis (as tested by applying \( \text{arb_contains_zero()} \) to the imaginary part) corresponds to a real root (its imaginary part can be set to zero), and every other root enclosure corresponds to a non-real root (with known sign for the imaginary part).

If this function returns zero, then the signs of the imaginary parts are not known for certain, based on the accuracy of the inputs and the working precision \( \text{prec} \).

An \texttt{acb_mat_t} represents a dense matrix over the complex numbers, implemented as an array of entries of type \texttt{acb_struct}.

The dimension (number of rows and columns) of a matrix is fixed at initialization, and the user must ensure that inputs and outputs to an operation have compatible dimensions. The number of rows or columns in a matrix can be zero.
3.9.1 Types, macros and constants

acb_mat_struct

acb_mat_t
Contains a pointer to a flat array of the entries (entries), an array of pointers to the start of each row (rows), and the number of rows (r) and columns (c).

An acb_mat_t is defined as an array of length one of type acb_mat_struct, permitting an acb_mat_t to be passed by reference.

acb_mat_entry(mat, i, j)
Macro giving a pointer to the entry at row i and column j.

acb_mat_nrows(mat)
Returns the number of rows of the matrix.

acb_mat_ncols(mat)
Returns the number of columns of the matrix.

3.9.2 Memory management

void acb_mat_init(acb_mat_t mat, slong r, slong c)
Initializes the matrix, setting it to the zero matrix with r rows and c columns.

void acb_mat_clear(acb_mat_t mat)
Clears the matrix, deallocating all entries.

3.9.3 Conversions

void acb_mat_set(acb_mat_t dest, const acb_mat_t src)

void acb_mat_set_fmpz_mat(acb_mat_t dest, const fmpz_mat_t src)

void acb_mat_set_round_fmpz_mat(acb_mat_t dest, const fmpz_mat_t src, slong prec)

void acb_mat_set_fmpq_mat(acb_mat_t dest, const fmpq_mat_t src, slong prec)

void acb_mat_set_arb_mat(acb_mat_t dest, const arb_mat_t src)

void acb_mat_set_round_arb_mat(acb_mat_t dest, const arb_mat_t src, slong prec)
Sets dest to src. The operands must have identical dimensions.

3.9.4 Random generation

void acb_mat_randtest(acb_mat_t mat, flint_rand_t state, slong prec, slong mag_bits)
Sets mat to a random matrix with up to prec bits of precision and with exponents of width up to mag_bits.

3.9.5 Input and output

void acb_mat_printd(const acb_mat_t mat, slong digits)
Prints each entry in the matrix with the specified number of decimal digits.
3.9.6 Comparisons

int acb_mat_equal (const acb_mat_t mat1, const acb_mat_t mat2)
Returns nonzero iff the matrices have the same dimensions and identical entries.

int acb_mat_overlaps (const acb_mat_t mat1, const acb_mat_t mat2)
Returns nonzero iff the matrices have the same dimensions and each entry in mat1 overlaps with the corresponding entry in mat2.

int acb_mat_contains (const acb_mat_t mat1, const acb_mat_t mat2)
int acb_mat_contains_fmpz_mat (const acb_mat_t mat1, const fmpz_mat_t mat2)
int acb_mat_contains_fmpq_mat (const acb_mat_t mat1, const fmpq_mat_t mat2)
Returns nonzero iff the matrices have the same dimensions and each entry in mat2 is contained in the corresponding entry in mat1.

int acb_mat_eq (const acb_mat_t mat1, const acb_mat_t mat2)
Returns nonzero iff mat1 and mat2 certainly represent the same matrix.

int acb_mat_ne (const acb_mat_t mat1, const acb_mat_t mat2)
Returns nonzero iff mat1 and mat2 certainly do not represent the same matrix.

int acb_mat_is_real (const acb_mat_t mat)
Returns nonzero iff all entries in mat have zero imaginary part.

3.9.7 Special matrices

void acb_mat_zero (acb_mat_t mat)
Sets all entries in mat to zero.

void acb_mat_one (acb_mat_t mat)
Sets the entries on the main diagonal to ones, and all other entries to zero.

3.9.8 Transpose

void acb_mat_transpose (acb_mat_t dest, const acb_mat_t src)
Sets dest to the exact transpose src. The operands must have compatible dimensions. Aliasing is allowed.

3.9.9 Norms

void acb_mat_bound_inf_norm (mag_t b, const acb_mat_t A)
Sets b to an upper bound for the infinity norm (i.e. the largest absolute value row sum) of A.

3.9.10 Arithmetic

void acb_mat_neg (acb_mat_t dest, const acb_mat_t src)
Sets dest to the exact negation of src. The operands must have the same dimensions.

void acb_mat_add (acb_mat_t res, const acb_mat_t mat1, const acb_mat_t mat2, slong prec)
Sets res to the sum of mat1 and mat2. The operands must have the same dimensions.

void acb_mat_sub (acb_mat_t res, const acb_mat_t mat1, const acb_mat_t mat2, slong prec)
Sets res to the difference of mat1 and mat2. The operands must have the same dimensions.
void \texttt{acb_mat_mul} \((acb\_mat\_t \texttt{res}, \text{const } \texttt{acb\_mat\_t mat1}, \text{const } \texttt{acb\_mat\_t mat2}, \text{slong prec})\)

Sets \texttt{res} to the matrix product of \texttt{mat1} and \texttt{mat2}. The operands must have compatible dimensions for matrix multiplication.

void \texttt{acb_mat_sqr} \((acb\_mat\_t \texttt{res}, \text{const } \texttt{acb\_mat\_t mat}, \text{slong prec})\)

Sets \texttt{res} to the matrix square of \texttt{mat}. The operands must both be square with the same dimensions.

void \texttt{acb_mat_pow_ui} \((acb\_mat\_t \texttt{res}, \text{const } \texttt{acb\_mat\_t mat}, \text{ulong exp}, \text{slong prec})\)

Sets \texttt{res} to \texttt{mat} raised to the power \texttt{exp}. Requires that \texttt{mat} is a square matrix.

3.9.11 Scalar arithmetic

void \texttt{acb_mat_scalar_mul_2exp_si} \((acb\_mat\_t B, \text{const } \texttt{acb\_mat\_t A}, \text{slong c})\)

Sets \texttt{B} to \texttt{A} multiplied by \(2^c\).

void \texttt{acb_mat_scalar_addmul_si} \((acb\_mat\_t B, \text{const } \texttt{acb\_mat\_t A}, \text{long c, slong prec})\)

void \texttt{acb_mat_scalar_addmul_fmpz} \((acb\_mat\_t B, \text{const } \texttt{acb\_mat\_t A}, \text{fmpz_t c, slong prec})\)

void \texttt{acb_mat_scalar_addmul_arb} \((acb\_mat\_t B, \text{const } \texttt{acb\_mat\_t A}, \text{arb_t c, slong prec})\)

void \texttt{acb_mat_scalar_addmul_acb} \((acb\_mat\_t B, \text{const } \texttt{acb\_mat\_t A}, \text{acb_t c, slong prec})\)

Sets \texttt{B} to \(B + A \times c\).

void \texttt{acb_mat_scalar_mul_si} \((acb\_mat\_t B, \text{const } \texttt{acb\_mat\_t A}, \text{slong c, slong prec})\)

void \texttt{acb_mat_scalar_mul_fmpz} \((acb\_mat\_t B, \text{const } \texttt{acb\_mat\_t A}, \text{fmpz_t c, slong prec})\)

void \texttt{acb_mat_scalar_mul_arb} \((acb\_mat\_t B, \text{const } \texttt{acb\_mat\_t A}, \text{arb_t c, slong prec})\)

void \texttt{acb_mat_scalar_mul_acb} \((acb\_mat\_t B, \text{const } \texttt{acb\_mat\_t A}, \text{acb_t c, slong prec})\)

Sets \texttt{B} to \(A \times c\).

void \texttt{acb_mat_scalar_div_si} \((acb\_mat\_t B, \text{const } \texttt{acb\_mat\_t A}, \text{slong c, slong prec})\)

void \texttt{acb_mat_scalar_div_fmpz} \((acb\_mat\_t B, \text{const } \texttt{acb\_mat\_t A}, \text{fmpz_t c, slong prec})\)

void \texttt{acb_mat_scalar_div_arb} \((acb\_mat\_t B, \text{const } \texttt{acb\_mat\_t A}, \text{arb_t c, slong prec})\)

void \texttt{acb_mat_scalar_div_acb} \((acb\_mat\_t B, \text{const } \texttt{acb\_mat\_t A}, \text{acb_t c, slong prec})\)

Sets \texttt{B} to \(A/c\).

3.9.12 Gaussian elimination and solving

int \texttt{acb_mat_lu} \((\text{slong } \ast \texttt{perm}, \texttt{acb\_mat\_t LU, const } \texttt{acb\_mat\_t A}, \text{slong prec})\)

Given an \(n \times n\) matrix \(A\), computes an LU decomposition \(P LU = A\) using Gaussian elimination with partial pivoting. The input and output matrices can be the same, performing the decomposition in-place.

Entry \(i\) in the permutation vector \texttt{perm} is set to the row index in the input matrix corresponding to row \(i\) in the output matrix.

The algorithm succeeds and returns nonzero if it can find \(n\) invertible (i.e. not containing zero) pivot entries. This guarantees that the matrix is invertible.

The algorithm fails and returns zero, leaving the entries in \(P\) and \(LU\) undefined, if it cannot find \(n\) invertible pivot elements. In this case, either the matrix is singular, the input matrix was computed to insufficient precision, or the LU decomposition was attempted at insufficient precision.
void **acb_mat_solve_lu_precomp**(acb_mat_t X, const slong * perm, const acb_mat_t LU, const acb_mat_t B, slong prec)

Solves $AX = B$ given the precomputed nonsingular LU decomposition $A = PLU$. The matrices $X$ and $B$ are allowed to be aliased with each other, but $X$ is not allowed to be aliased with $LU$.

int **acb_mat_solve**(acb_mat_t X, const acb_mat_t A, const acb_mat_t B, slong prec)

Solves $AX = B$ where $A$ is a nonsingular $n \times n$ matrix and $X$ and $B$ are $n \times m$ matrices, using LU decomposition.

If $m > 0$ and $A$ cannot be inverted numerically (indicating either that $A$ is singular or that the precision is insufficient), the values in the output matrix are left undefined and zero is returned. A nonzero return value guarantees that $A$ is invertible and that the exact solution matrix is contained in the output.

int **acb_mat_inv**(acb_mat_t X, const acb_mat_t A, slong prec)

Sets $X = A^{-1}$ where $A$ is a square matrix, computed by solving the system $AX = I$.

If $A$ cannot be inverted numerically (indicating either that $A$ is singular or that the precision is insufficient), the values in the output matrix are left undefined and zero is returned. A nonzero return value guarantees that the matrix is invertible and that the exact inverse is contained in the output.

void **acb_mat_det**(acb_t det, const acb_mat_t A, slong prec)

Computes the determinant of the matrix, using Gaussian elimination with partial pivoting. If at some point an invertible pivot element cannot be found, the elimination is stopped and the magnitude of the determinant of the remaining submatrix is bounded using Hadamard’s inequality.

### 3.9.13 Characteristic polynomial

void **acb_mat_charpoly**(acb_ptr cp, const acb_mat_t mat, slong prec)

void **acb_mat_charpoly**(acb_poly_t cp, const acb_mat_t mat, slong prec)

Sets $cp$ to the characteristic polynomial of $mat$ which must be a square matrix. If the matrix has $n$ rows, the underscore method requires space for $n + 1$ output coefficients. Employs a division-free algorithm using $O(n^4)$ operations.

### 3.9.14 Special functions

void **acb_mat_exp**(acb_mat_t B, const acb_mat_t A, slong prec)

Sets $B$ to the exponential of the matrix $A$, defined by the Taylor series

$$\exp(A) = \sum_{k=0}^{\infty} \frac{A^k}{k!}.$$  

The function is evaluated as $\exp(A/2^r)^2^r$, where $r$ is chosen to give rapid convergence of the Taylor series. The series is evaluated using rectangular splitting. Error bounds are computed as for **arb_mat_exp**().

void **acb_mat_trace**(acb_t trace, const acb_mat_t mat, slong prec)

Sets $trace$ to the trace of the matrix, i.e. the sum of entries on the main diagonal of $mat$. The matrix is required to be square.

### 3.10 acb_calc.h – calculus with complex-valued functions

This module provides functions for operations of calculus over the complex numbers (intended to include root-finding, integration, and so on).
3.10.1 Types, macros and constants

//arb_calclnc_func_t
Typedef for a pointer to a function with signature:

```c
int func(acb_ptr out, const acb_t inp, void * param, slong order, slong prec)
```

implementing a univariate complex function \( f(x) \). When called, `func` should write to `out` the first order coefficients in the Taylor series expansion of \( f(x) \) at the point `inp`, evaluated at a precision of `prec` bits. The `param` argument may be used to pass through additional parameters to the function. The return value is reserved for future use as an error code. It can be assumed that `out` and `inp` are not aliased and that `order` is positive.

3.10.2 Bounds

//acb_calc_cauchy_bound
Sets `bound` to a ball containing the value of the integral

\[
C(x, r) = \frac{1}{2\pi r} \oint_{|z-x|=r} |f(z)|dz = \int_0^1 |f(x + re^{2\pi it})|dt
\]

where \( f \) is specified by (\( \text{func}, \text{param} \)) and \( r \) is given by `radius`. The integral is computed using a simple step sum. The integration range is subdivided until the order of magnitude of \( b \) can be determined (i.e. its error bound is smaller than its midpoint), or until the step length has been cut in half `maxdepth` times. This function is currently implemented completely naively, and repeatedly subdivides the whole integration range instead of performing adaptive subdivisions.

3.10.3 Integration

//acb_calc_integrate_taylor
Computes the integral

\[
I = \int_a^b f(t)dt
\]

where \( f \) is specified by (\( \text{func}, \text{param} \)), following a straight-line path between the complex numbers \( a \) and \( b \) which both must be finite.

The integral is approximated by piecewise centered Taylor polynomials. Rigorous truncation error bounds are calculated using the Cauchy integral formula. More precisely, if the Taylor series of \( f \) centered at the point \( m \) is \( f(m + x) = \sum_{n=0}^{\infty} a_n x^n \), then

\[
\int f(m + x) = \left( \sum_{n=0}^{N-1} a_n \frac{x^{n+1}}{n+1} \right) + \left( \sum_{n=N}^{\infty} a_n \frac{x^{n+1}}{n+1} \right).
\]

For sufficiently small \( x \), the second series converges and its absolute value is bounded by

\[
\sum_{n=N}^{\infty} \frac{C(m, R) |x|^{n+1}}{R^n (n+1)^N} = \frac{C(m, R) R x}{(R - x)(N + 1)} \left( \frac{x}{R} \right)^N.
\]

It is required that any singularities of \( f \) are isolated from the path of integration by a distance strictly greater than the positive value `outer_radius` (which is the integration radius used for the Cauchy bound). Taylor series
step lengths are chosen so as not to exceed $inner_radius$, which must be strictly smaller than $outer_radius$ for convergence. A smaller $inner_radius$ gives more rapid convergence of each Taylor series but means that more series might have to be used. A reasonable choice might be to set $inner_radius$ to half the value of $outer_radius$, giving roughly one accurate bit per term.

The truncation point of each Taylor series is chosen so that the absolute truncation error is roughly $2^{-p}$ where $p$ is given by $accuracy\_goal$ (in the future, this might change to a relative accuracy). Arithmetic operations and function evaluations are performed at a precision of $prec$ bits. Note that due to accumulation of numerical errors, both values may have to be set higher (and the endpoints may have to be computed more accurately) to achieve a desired accuracy.

This function chooses the evaluation points uniformly rather than implementing adaptive subdivision.

### 3.11 acb_hypgeom.h – hypergeometric functions in the complex numbers

The generalized hypergeometric function is formally defined by

$$pF_q(a_1, \ldots, a_p; b_1, \ldots, b_q; z) = \sum_{k=0}^{\infty} \frac{(a_1)_k \cdots (a_p)_k}{(b_1)_k \cdots (b_q)_k} \frac{z^k}{k!}. $$

It can be interpreted using analytic continuation or regularization when the sum does not converge. In a looser sense, we understand “hypergeometric functions” to be linear combinations of generalized hypergeometric functions with prefactors that are products of exponentials, powers, and gamma functions.

#### 3.11.1 Convergent series

In this section, we define

$$T(k) = \frac{\prod_{i=0}^{p-1} (a_i)_k}{\prod_{i=0}^{q-1} (b_i)_k} z^k$$

and

$$pH_q(a_0, \ldots, a_{p-1}; b_0 \ldots b_{q-1}; z) = p+1 F_q(a_0, \ldots, a_{p-1}, 1; b_0 \ldots b_{q-1}; z) = \sum_{k=0}^{\infty} T(k)$$

For the conventional generalized hypergeometric function $pF_q$, compute $pH_{q+1}$ with the explicit parameter $b_q = 1$, or remove a 1 from the $a_i$ parameters if there is one.

**void acb_hypgeom_pfq_bound_factor** (mag_t C, acb_srcptr a, slong p, acb_srcptr b, slong q, const acb_t z, ulong n)

Computes a factor $C$ such that

$$\sum_{k=n}^{\infty} T(k) \leq C |T(n)|.$$  

We check that $Re(b + n) > 0$ for all lower parameters $b$. If this does not hold, $C$ is set to infinity. Otherwise, we cancel out pairs of parameters $a$ and $b$ against each other. We have

$$\left| \frac{a + k}{b + k} \right| = \left| \frac{1 + \frac{a - b}{b + k}}{\frac{b}{b + k}} \right| \leq 1 + \left| \frac{a - b}{b + k} \right|$$
and
\[ \left| \frac{1}{b+k} \right| \leq \left| \frac{1}{b+n} \right| \]

for all \( k \geq n \). This gives us a constant \( D \) such that \( T(k+1) \leq DT(k) \) for all \( k \geq n \). If \( D \geq 1 \), we set \( C \) to infinity. Otherwise, we take \( C = \sum_{k=0}^{\infty} D^k = (1 - D)^{-1} \).

As currently implemented, the bound becomes infinite when \( n \) is too small, even if the series converges.

\texttt{slong \textbf{acb_hypgeom_pfq_choose_n}(acb_srcptr \ a, slong \ p, \ \textbf{acb_srcptr} \ b, slong \ q, \ \textbf{const} \ \textbf{acb_t} \ z, \ \textbf{slong} \ \textbf{prec})}

Heuristically attempts to choose a number of terms \( n \) to sum of a hypergeometric series at a working precision of \( \text{prec} \) bits.

Uses double precision arithmetic internally. As currently implemented, it can fail to produce a good result if the parameters are extremely large or extremely close to nonpositive integers.

Numerical cancellation is assumed to be significant, so truncation is done when the current term is \( \text{prec} \) bits smaller than the largest encountered term.

This function will also attempt to pick a reasonable truncation point for divergent series.

\texttt{void \textbf{acb_hypgeom_pfq_sum_forward}(acb_t \ s, \ \textbf{acb_t} \ t, \ \textbf{acb_srcptr} \ a, \ \textbf{slong} \ p, \ \textbf{acb_srcptr} \ b, \ \textbf{slong} \ q, \ \textbf{const} \ \textbf{acb_t} \ z, \ \textbf{slong} \ \textbf{prec})}

\texttt{void \textbf{acb_hypgeom_pfq_sum_rs}(acb_t \ s, \ \textbf{acb_t} \ t, \ \textbf{acb_srcptr} \ a, \ \textbf{slong} \ p, \ \textbf{acb_srcptr} \ b, \ \textbf{slong} \ q, \ \textbf{const} \ \textbf{acb_t} \ z, \ \textbf{slong} \ \textbf{prec})}

\texttt{void \textbf{acb_hypgeom_pfq_sum_bs}(acb_t \ s, \ \textbf{acb_t} \ t, \ \textbf{acb_srcptr} \ a, \ \textbf{slong} \ p, \ \textbf{acb_srcptr} \ b, \ \textbf{slong} \ q, \ \textbf{const} \ \textbf{acb_t} \ z, \ \textbf{slong} \ \textbf{prec})}

\texttt{void \textbf{acb_hypgeom_pfq_sum_fme}(acb_t \ s, \ \textbf{acb_t} \ t, \ \textbf{acb_srcptr} \ a, \ \textbf{slong} \ p, \ \textbf{acb_srcptr} \ b, \ \textbf{slong} \ q, \ \textbf{const} \ \textbf{acb_t} \ z, \ \textbf{slong} \ \textbf{prec})}

\texttt{void \textbf{acb_hypgeom_pfq_sum}(acb_t \ s, \ \textbf{acb_t} \ t, \ \textbf{acb_srcptr} \ a, \ \textbf{slong} \ p, \ \textbf{acb_srcptr} \ b, \ \textbf{slong} \ q, \ \textbf{const} \ \textbf{acb_t} \ z, \ \textbf{slong} \ \textbf{prec})}

Computes \( s = \sum_{k=0}^{n-1} T(k) \) and \( t = T(n) \). Does not allow aliasing between input and output variables. We require \( n \geq 0 \).

The \textit{forward} version computes the sum using forward recurrence.

The \textit{bs} version computes the sum using binary splitting.

The \textit{rs} version computes the sum in reverse order using rectangular splitting. It only computes a magnitude bound for the value of \( t \).

The \textit{fme} version uses fast multipoint evaluation.

The default version automatically chooses an algorithm depending on the inputs.

\texttt{void \textbf{acb_hypgeom_pfq_sum_bs_invz}(acb_t \ s, \ \textbf{acb_t} \ t, \ \textbf{acb_srcptr} \ a, \ \textbf{slong} \ p, \ \textbf{acb_srcptr} \ b, \ \textbf{slong} \ q, \ \textbf{const} \ \textbf{acb_t} \ w, \ \textbf{slong} \ n, \ \textbf{slong} \ \textbf{prec})}

\texttt{void \textbf{acb_hypgeom_pfq_sum_invz}(acb_t \ s, \ \textbf{acb_t} \ t, \ \textbf{acb_srcptr} \ a, \ \textbf{slong} \ p, \ \textbf{acb_srcptr} \ b, \ \textbf{slong} \ q, \ \textbf{const} \ \textbf{acb_t} \ z, \ \textbf{const} \ \textbf{acb_t} \ w, \ \textbf{slong} \ n, \ \textbf{slong} \ \textbf{prec})}

Like \texttt{acb_hypgeom_pfq_sum()}, but taking advantage of \( w = 1/z \) possibly having few bits.

\texttt{void \textbf{acb_hypgeom_pfq_direct}(acb_t \ \textbf{res}, \ \textbf{acb_srcptr} \ a, \ \textbf{slong} \ p, \ \textbf{acb_srcptr} \ b, \ \textbf{slong} \ q, \ \textbf{const} \ \textbf{acb_t} \ z, \ \textbf{slong} \ n, \ \textbf{slong} \ \textbf{prec})}

Computes
\[ pH_q(z) = \sum_{k=0}^{\infty} T(k) = \sum_{k=0}^{n-1} T(k) + \varepsilon \]
directly from the defining series, including a rigorous bound for the truncation error $\varepsilon$ in the output.

If $n < 0$, this function chooses a number of terms automatically using `acb_hypgeom_pfq_choose_n()`. void `acb_hypgeom_pfq_series_direct` (acb_poly_t res, const acb_poly_struct * a, slong p, const acb_poly_struct * b, slong q, const acb_poly_t z, int regularized, slong n, slong len, slong prec)

Computes $_pH_q(z)$ directly using the defining series, given parameters and argument that are power series. The result is a power series of length $\text{len}$.

An error bound is computed automatically as a function of the number of terms $n$. If $n < 0$, the number of terms is chosen automatically.

If `regularized` is set, the regularized hypergeometric function is computed instead.

### 3.11.2 Asymptotic series

Let $U(a, b, z)$ denote the confluent hypergeometric function of the second kind with the principal branch cut, and let $U^* = z^a U(a, b, z)$. For all $z \neq 0$ and $b \notin \mathbb{Z}$ (but valid for all $b$ as a limit), we have (DLMF 13.2.42)

$$U(a, b, z) = \frac{\Gamma(1-b)}{\Gamma(a-b+1)} M(a, b, z) + \frac{\Gamma(b-1)}{\Gamma(a)} z^{1-b} M(a-b+1, 2-b, z).$$

Moreover, for all $z \neq 0$ we have

$$\frac{1}{\Gamma(b)} \binom{a}{b} = \frac{(-z)^{-a}}{\Gamma(b-a)} U^*(a, b, z) + \frac{z^{a-b} e^z}{\Gamma(a)} U^*(b-a, b, -z)$$

which is equivalent to DLMF 13.2.41 (but simpler in form).

We have the asymptotic expansion

$$U^*(a, b, z) \sim 2F_{0}(a, a-b+1, -1/z)$$

where $2F_{0}(a, b, z)$ denotes a formal hypergeometric series, i.e.

$$U^*(a, b, z) = \sum_{k=0}^{n-1} \frac{(a)_k (a-b+1)_k}{k! (-z)^k} + \varepsilon_n(z).$$

The error term $\varepsilon_n(z)$ is bounded according to DLMF 13.7. A case distinction is made depending on whether $z$ lies in one of three regions which we index by $R$. Our formula for the error bound increases with the value of $R$, so we can always choose the larger out of two indices if $z$ lies in the union of two regions.

Let $r = |b - 2a|$. If $\text{Re}(z) \geq r$, set $R = 1$. Otherwise, if $\text{Im}(z) \geq r$ or $\text{Re}(z) > 0 \land |z| \geq r$, set $R = 2$. Otherwise, if $|z| \geq 2r$, set $R = 3$. Otherwise, the bound is infinite. If the bound is finite, we have

$$|\varepsilon_n(z)| \leq 2\alpha C_n \left| \frac{(a)_n (a-b+1)_n}{n! z^n} \right| \exp(2\alpha \rho C_1/|z|)$$
in terms of the following auxiliary quantities

\[
\sigma = \left| (b - 2a)/z \right|
\]

\[
C_n = \begin{cases} 
1 & \text{if } R = 1 \\
\chi(n) & \text{if } R = 2 \\
(\chi(n) + \rho \nu^2 n) \nu^n & \text{if } R = 3 
\end{cases}
\]

\[
\nu = \left( \frac{1}{2} + \frac{1}{2} \sqrt{1 - 4\sigma^2} \right)^{-1/2} \leq 1 + 2\sigma^2 
\]

\[
\chi(n) = \sqrt{\pi} \Gamma\left(\frac{1}{2}n + 1\right)/\Gamma\left(\frac{1}{2}n + \frac{1}{2}\right) 
\]

\[
\sigma' = \begin{cases} 
\sigma & \text{if } R \neq 3 \\
\nu \sigma & \text{if } R = 3 
\end{cases}
\]

\[
\alpha = (1 - \sigma')^{-1} 
\]

\[
\rho = \frac{1}{2}\left[2a^2 - 2ab + b\right] + \sigma'(1 + \frac{1}{4}\sigma')\left(1 - \sigma'\right)^{-2} 
\]

void \texttt{acb\_hypgeom\_u\_asymp} (\texttt{acb\_t res}, \texttt{const acb\_t a}, \texttt{const acb\_t b}, \texttt{const acb\_t z}, \texttt{slong n}, \texttt{slong prec})

Sets \(U^*(a, b, z)\) computed using \(n\) terms of the asymptotic series, with a rigorous bound for the error included in the output. We require \(n \geq 0\).

int \texttt{acb\_hypgeom\_u\_use\_asymp} (\texttt{const acb\_t z}, \texttt{slong prec})

Heuristically determines whether the asymptotic series can be used to evaluate \(U(a, b, z)\) to \(\text{prec}\) accurate bits (assuming that \(a\) and \(b\) are small).

### 3.11.3 Confluent hypergeometric functions

void \texttt{acb\_hypgeom\_u\_1f1\_series} (\texttt{acb\_poly\_t res}, \texttt{const acb\_poly\_t a}, \texttt{const acb\_poly\_t b}, \texttt{const acb\_poly\_t z}, \texttt{slong len}, \texttt{slong prec})

Computes \(U(a, b, z)\) as a power series truncated to length \(\text{len}\), given \(a, b, z \in \mathbb{C}[[x]]\). If \(b[0] \in \mathbb{Z}\), it computes one extra derivative and removes the singularity (it is then assumed that \(b[1] \neq 0\)). As currently implemented, the output is indeterminate if \(b\) is nonexact and contains an integer.

void \texttt{acb\_hypgeom\_u\_1f1} (\texttt{acb\_t res}, \texttt{const acb\_t a}, \texttt{const acb\_t b}, \texttt{const acb\_t z}, \texttt{slong prec})

Computes \(U(a, b, z)\) as a sum of two convergent hypergeometric series. If \(b \in \mathbb{Z}\), it computes the limit value via \texttt{acb\_hypgeom\_u\_1f1\_series}(). As currently implemented, the output is indeterminate if \(b\) is nonexact and contains an integer.

void \texttt{acb\_hypgeom\_u} (\texttt{acb\_t res}, \texttt{const acb\_t a}, \texttt{const acb\_t b}, \texttt{const acb\_t z}, \texttt{slong prec})

Computes \(U(a, b, z)\) using an automatic algorithm choice. The function \texttt{acb\_hypgeom\_u\_asymp}() is used if \(a\) or \(a - b + 1\) is a nonpositive integer (in which case the asymptotic series terminates), or if \(z\) is sufficiently large. Otherwise \texttt{acb\_hypgeom\_u\_1f1()} is used.

void \texttt{acb\_hypgeom\_m\_asymp} (\texttt{acb\_t res}, \texttt{const acb\_t a}, \texttt{const acb\_t b}, \texttt{const acb\_t z}, \texttt{int regularized}, \texttt{slong prec})

void \texttt{acb\_hypgeom\_m\_1f1} (\texttt{acb\_t res}, \texttt{const acb\_t a}, \texttt{const acb\_t b}, \texttt{const acb\_t z}, \texttt{int regularized}, \texttt{slong prec})

void \texttt{acb\_hypgeom\_m} (\texttt{acb\_t res}, \texttt{const acb\_t a}, \texttt{const acb\_t b}, \texttt{const acb\_t z}, \texttt{int regularized}, \texttt{slong prec})

Computes the confluent hypergeometric function \(M(a, b, z) = _1F_1(a, b, z)\), or \(M(a, b, z) = \frac{1}{\Gamma(\alpha)}F_1(a, b, z)\) if \texttt{regularized} is set.

void \texttt{acb\_hypgeom\_0f1\_asymp} (\texttt{acb\_t res}, \texttt{const acb\_t a}, \texttt{const acb\_t z}, \texttt{int regularized}, \texttt{slong prec})

void \texttt{acb\_hypgeom\_0f1\_direct} (\texttt{acb\_t res}, \texttt{const acb\_t a}, \texttt{const acb\_t z}, \texttt{int regularized}, \texttt{slong prec})
void \texttt{acb\_hypgeom\_0f1} (\texttt{acb\_t res, const acb\_t a, const acb\_t z, int regularized, slong prec})  
Computes the confluent hypergeometric function $\, _0F_1(a,z)$, or $\frac{1}{\Gamma(a)}$ if \texttt{regularized} is set, using asymptotic expansions, direct summation, or an automatic algorithm choice. The \texttt{asymp} version uses the asymptotic expansions of Bessel functions, together with the connection formulas  
\[ \frac{\, _0F_1(a,z)}{\Gamma(a)} = (-z)^{(1-a)/2}J_{a-1}(2\sqrt{-z}) = z^{(1-a)/2}I_{a-1}(2\sqrt{z}). \]

The Bessel-$J$ function is used in the left half-plane and the Bessel-$I$ function is used in the right half-plane, to avoid loss of accuracy due to evaluating the square root on the branch cut.

### 3.11.4 The error function

void \texttt{acb\_hypgeom\_erf\_1fla} (\texttt{acb\_t res, const acb\_t z, slong prec})  
void \texttt{acb\_hypgeom\_erf\_1flb} (\texttt{acb\_t res, const acb\_t z, slong prec})  
void \texttt{acb\_hypgeom\_erf\_asymp} (\texttt{acb\_t res, const acb\_t z, slong prec, slong prec2})  
void \texttt{acb\_hypgeom\_erf} (\texttt{acb\_t res, const acb\_t z, slong prec})  
Computes the error function respectively using
\[ \text{erf}(z) = \frac{2z}{\sqrt{\pi}} F_1(\frac{1}{2}, \frac{3}{2}, -z^2) \]
\[ \text{erf}(z) = \frac{2ze^{-z^2}}{\sqrt{\pi}} F_1(1, \frac{3}{2}, z^2) \]
\[ \text{erf}(z) = \frac{z}{\sqrt{z^2}} \left( 1 - \frac{e^{-z^2}}{\sqrt{\pi}} U\left(\frac{1}{2}, \frac{1}{2}, z^2\right) \right). \]

and an automatic algorithm choice. The \texttt{asymp} version takes a second precision to use for the $U$ term.

void \texttt{acb\_hypgeom\_erfc} (\texttt{acb\_t res, const acb\_t z, slong prec})  
Computes the complementary error function \texttt{erfc}(z) = 1-\texttt{erf}(z). This function avoids catastrophic cancellation for large positive $z$.

void \texttt{acb\_hypgeom\_erfi} (\texttt{acb\_t res, const acb\_t z, slong prec})  
Computes the imaginary error function \texttt{erfi}(z) = -i\texttt{erf}(iz). This is a trivial wrapper of \texttt{acb\_hypgeom\_erf}.

### 3.11.5 Bessel functions

void \texttt{acb\_hypgeom\_bessel\_j\_asymp} (\texttt{acb\_t res, const acb\_t nu, const acb\_t z, slong prec})  
Computes the Bessel function of the first kind via \texttt{acb\_hypgeom\_u\_asymp}. For all complex $\nu, z$, we have
\[ J_\nu(z) = \frac{z^\nu}{2^\nu e^{i\nu} \Gamma(\nu+1)} F_1(\nu + \frac{1}{2}, 2\nu + 1, 2iz) = A_+ B_+ + A_- B_- \]
where
\[ A_\pm = z^{\nu}(z^2)^{-\frac{1}{2}} \nu - i z (\mp i z)^{\frac{1}{2}} \nu^{2} (2\pi)^{-1/2} = (\pm iz)^{-1/2} \nu z^{\nu}(2\pi)^{-1/2} \]
\[ B_\pm = e^{\mp iz} U^{*}(\nu + \frac{1}{2}, 2\nu + 1, \mp 2iz). \]

Nicer representations of the factors $A_\pm$ can be given depending conditionally on the parameters. If $\nu + \frac{1}{2} = n \in \mathbb{Z}$, we have $A_\pm = (\pm i)^n (2\pi z)^{-1/2}$. And if $\text{Re}(z) > 0$, we have $A_\pm = \exp(\mp i [(2\nu + 1)/4]\pi) (2\pi z)^{-1/2}$.  

3.11. acb\_hypgeom.h – hypergeometric functions in the complex numbers

Arb Documentation, Release 2.8.1
void **acb_hypgeom_bessel_j_0f1** (*acb_t res*, const *acb_t nu*, const *acb_t z*, slong *prec*)

Computes the Bessel function of the first kind from

\[ J_{\nu}(z) = \frac{1}{\Gamma(\nu + 1)} \left( \frac{z}{2} \right)^{\nu} {}_0F_1 \left( \nu + 1, -\frac{z^2}{4} \right) \].

void **acb_hypgeom_bessel_j** (*acb_t res*, const *acb_t nu*, const *acb_t z*, slong *prec*)

Computes the Bessel function of the first kind \( J_{\nu}(z) \) using an automatic algorithm choice.

void **acb_hypgeom_bessel_y** (*acb_t res*, const *acb_t nu*, const *acb_t z*, slong *prec*)

Computes the Bessel function of the second kind \( Y_{\nu}(z) \) from the formula

\[ Y_{\nu}(z) = \frac{\cos(\nu \pi) J_{\nu}(z) - J_{-\nu}(z)}{\sin(\nu \pi)} \]

unless \( \nu = n \) is an integer in which case the limit value

\[ Y_n(z) = -\frac{2}{\pi} \left( i^n K_n(iz) + [\log(iz) - \log(z)] J_n(z) \right) \]

is computed. As currently implemented, the output is indeterminate if \( \nu \) is nonexact and contains an integer.

void **acb_hypgeom_bessel_jy** (*acb_t res1*, *acb_t res2*, const *acb_t nu*, const *acb_t z*, slong *prec*)

Sets res1 to \( J_{\nu}(z) \) and res2 to \( Y_{\nu}(z) \), computed simultaneously. From these values, the user can easily construct the Bessel functions of the third kind (Hankel functions) \( H^{(1)}_{\nu}(z), H^{(2)}_{\nu}(z) = J_{\nu}(z) \pm iY_{\nu}(z) \).

void **acb_hypgeom_bessel_i_0f1** (*acb_t res*, const *acb_t nu*, const *acb_t z*, slong *prec*)

Computes the modified Bessel function of the first kind \( I_{\nu}(z) = z^{\nu}e^{-\nu}J_{\nu}(iz) \) respectively using asymptotic series (see **acb_hypgeom_bessel_j_0f1**(1)), the convergent series

\[ I_{\nu}(z) = \frac{1}{\Gamma(\nu + 1)} \left( \frac{z}{2} \right)^{\nu} {}_0F_1 \left( \nu + 1, -\frac{z^2}{4} \right), \]

or an automatic algorithm choice.

void **acb_hypgeom_bessel_i** (*acb_t res*, const *acb_t nu*, const *acb_t z*, slong *prec*)

Computes the modified Bessel function of the second kind via **acb_hypgeom_u_0f1**(). For all \( \nu \) and all \( z \neq 0 \), we have

\[ K_{\nu}(z) = \left( \frac{2z}{\pi} \right)^{-1/2} e^{-z}U^{*}(\nu + \frac{1}{2}, 2\nu + 1, 2z). \]

void **acb_hypgeom_bessel_k_0f1_series** (*acb_poly_t res*, const *acb_poly_t nu*, const *acb_poly_t z*, slong *len*, slong *prec*)

Computes the modified Bessel function of the second kind \( K_{\nu}(z) \) as a power series truncated to length len, given \( \nu, z \in \mathbb{C}[[x]] \). Uses the formula

\[ K_{\nu}(z) = \frac{1}{2} \sin(\pi \nu) \left[ \left( \frac{z}{2} \right)^{-\nu} {}_0F_1 \left( 1 - \nu, \frac{z^2}{4} \right) - \left( \frac{z}{2} \right)^{\nu} {}_0F_1 \left( 1 + \nu, \frac{z^2}{4} \right) \right]. \]

If \( \nu[0] \in \mathbb{Z} \), it computes one extra derivative and removes the singularity (it is then assumed that \( \nu[1] \neq 0 \)). As currently implemented, the output is indeterminate if \( \nu[0] \) is nonexact and contains an integer.
void \texttt{acb\_hypgeom\_bessel\_k} (\texttt{acb\_t res}, const \texttt{acb\_t nu}, const \texttt{acb\_t z}, \texttt{slong prec})

Computes the modified Bessel function of the second kind from

\[
K_\nu(z) = \frac{1}{2} \left[ \left( \frac{z}{2} \right)^\nu \Gamma(\nu) F_1 \left( 1 - \nu, \frac{z^2}{4} \right) - \left( \frac{z}{2} \right)^\nu \frac{\pi}{\nu \sin(\pi \nu)} F_1 \left( \nu + 1, \frac{z^2}{4} \right) \right]
\]

if \( \nu \notin \mathbb{Z}. \) If \( \nu \in \mathbb{Z}, \) it computes the limit value via \texttt{acb\_hypgeom\_bessel\_k\_0f1\_series()}. As currently implemented, the output is indeterminate if \( \nu \) is nonexact and contains an integer.

void \texttt{acb\_hypgeom\_bessel\_k} (\texttt{acb\_t res}, const \texttt{acb\_t nu}, const \texttt{acb\_t z}, \texttt{slong prec})

Computes the modified Bessel function of the second kind \( K_\nu(z) \) using an automatic algorithm choice.

### 3.11.6 Airy functions

The Airy functions are linearly independent solutions of the differential equation \( y'' - zy = 0. \) All solutions are entire functions. The standard solutions are denoted \( \text{Ai}(z) \), \( \text{Bi}(z) \). For negative \( z \), both functions are oscillatory. For positive \( z \), the first function decreases exponentially while the second increases exponentially.

The Airy functions can be expressed in terms of Bessel functions of fractional order, but this is inconvenient since such formulas only hold piecewise (due to the Stokes phenomenon). Computation of the Airy functions can also be optimized more than Bessel functions in general. We therefore provide a dedicated interface for evaluating Airy functions.

The following functions optionally compute \( \text{Ai}(z), \text{Ai}'(z), \text{Bi}(z), \text{Bi}'(z) \) simultaneously. Any of the four functions can be omitted by passing \texttt{NULL} for the unwanted output variables. Note that higher derivatives of the Airy functions can be computed via recurrence relations.

void \texttt{acb\_hypgeom\_airy\_direct} (\texttt{acb\_t ai}, \texttt{acb\_t ai\_prime}, \texttt{acb\_t bi}, \texttt{acb\_t bi\_prime}, const \texttt{acb\_t z}, slong \( n \), \texttt{slong prec})

Computes the Airy functions using direct series expansions truncated at \( n \) terms. Error bounds are included in the output.

void \texttt{acb\_hypgeom\_airy\_asym}\texttt{p} (\texttt{acb\_t ai}, \texttt{acb\_t ai\_prime}, \texttt{acb\_t bi}, \texttt{acb\_t bi\_prime}, const \texttt{acb\_t z}, slong \( n \), \texttt{slong prec})

Computes the Airy functions using asymptotic expansions truncated at \( n \) terms. Error bounds, based on Olver (DLMF section 9.7), are included in the output. For \( \arg(z) < \pi \) and \( \zeta = (2/3)z^{3/2}, \) we have

\[
\text{Ai}(z) = \frac{e^{-\zeta}}{2 \sqrt{\pi} z^{1/4}} \left[ S_n(\zeta) + R_n(z) \right], \quad \text{Ai}'(z) = -\frac{z^{1/4} e^{-\zeta}}{2 \sqrt{\pi}} \left[ (S_n'(\zeta) + R_n'(z)) \right]
\]

\[
S_n(\zeta) = \sum_{k=0}^{n-1} (-1)^k \frac{u(k)}{\zeta^k}, \quad S_n'(\zeta) = \sum_{k=0}^{n-1} (-1)^k \frac{v(k)}{\zeta^k}
\]

\[
u(k) = \frac{(1/6)\zeta(5/6)_k}{2^k k!}, \quad v(k) = \frac{6k+1}{1-6k} u(k).
\]

Assuming that \( n \) is positive, the error terms are bounded by

\[
|R_n(z)| \leq C|u(n)||\zeta|^{-n}, \quad |R_n'(z)| \leq C|v(n)||\zeta|^{-n}
\]

where

\[
C = \begin{cases} 2 \exp(7/(36|\zeta|)) & |\arg(z)| \leq \pi/3 \\ 2 \chi(n) \exp(7\pi/72|\zeta|) & \pi/3 \leq |\arg(z)| \leq 2\pi/3 \\ 4\chi(n) \exp(7\pi/(36 \text{re}(\zeta))) |\cos(\arg(\zeta))|^{-n} & 2\pi/3 \leq |\arg(z)| < \pi. \end{cases}
\]

For computing \( \text{Bi} \) when \( z \) is roughly in the positive half-plane, we use the connection formulas

\[
\text{Bi}(z) = -i(2w^{+1} \text{Ai}(zw^{-2}) - \text{Ai}(z))
\]

\[
\text{Bi}(z) = +i(2w^{-1} \text{Ai}(zw^{+2}) - \text{Ai}(z))
\]

3.11. \texttt{acb\_hypgeom.h} – hypergeometric functions in the complex numbers
where \( w = \exp(\pi i/3) \). Combining roots of unity gives

\[
\text{Bi}(z) = \frac{1}{2\sqrt{\pi z}} \left[ 2X + iY \right] \\
\text{Bi}(z) = \frac{1}{2\sqrt{\pi z}} \left[ 2X - iY \right] \\
X = \exp(\mp \zeta)[S_n(-\zeta) + R_n(zw^2)], \quad Y = \exp(-\zeta)[S_n(\zeta) + R_n(z)]
\]

where the upper formula is valid for \(-\pi/3 < \arg(z) < \pi\) and the lower formula is valid for \(-\pi < \arg(z) < \pi/3\). We proceed analogously for the derivative of Bi.

In the negative half-plane, we use the connection formulas

\[
\text{Ai}(z) = e^{+\pi i/3} \text{Ai}(z_1) + e^{-\pi i/3} \text{Ai}(z_2) \\
\text{Bi}(z) = e^{-\pi i/6} \text{Ai}(z_1) + e^{+\pi i/6} \text{Ai}(z_2)
\]

where \( z_1 = -ze^{+\pi i/3} \), \( z_2 = -ze^{-\pi i/3} \). Provided that \(|\arg(-z)| < 2\pi/3\), we have \(|\arg(z_1)|, |\arg(z_2)| < \pi\), and thus the asymptotic expansion for Ai can be used. As before, we collect roots of unity to obtain

\[
\text{Ai}(z) = A_1[S_n(i\zeta) + R_n(z_1)] + A_2[S_n(-i\zeta) + R_n(z_2)] \\
\text{Bi}(z) = A_3[S_n(i\zeta) + R_n(z_1)] + A_4[S_n(-i\zeta) + R_n(z_2)]
\]

where \( \zeta = (2/3)(-z)^{3/2} \) and

\[
A_1 = \frac{\exp(-i(\zeta - \pi/4))}{2\sqrt{\pi(-z)^{1/4}}}, \quad A_2 = \frac{\exp(i(\zeta - \pi/4))}{2\sqrt{\pi(-z)^{1/4}}}, \quad A_3 = -iA_1, \quad A_4 = +iA_2.
\]

The differentiated formulas are analogous.

**void acb_hypgeomairybound** *(mag_t ai, mag_t aip, mag_t bi, mag_t bip, const acb_t z)*

Computes bounds for the Airy functions using first-order asymptotic expansions together with error bounds. This function uses some shortcuts to make it slightly faster than calling **acb_hypgeomairyasymp** with \( n = 1 \).

**void acb_hypgeomairy** *(acb_t ai, acb_t ai_prime, acb_t bi, acb_t bi_prime, const acb_t z, slong prec)*

Computes Airy functions using an automatic algorithm choice.

We use **acb_hypgeomairyasymp** whenever this gives full accuracy and **acb_hypgeomairydirec** otherwise.

In the latter case, we first use double precision arithmetic to determine an accurate estimate of the working precision needed to compute the Airy functions accurately for given \( z \). This estimate is obtained by comparing the leading-order asymptotic estimate of the Airy functions with the magnitude of the largest term in the power series. The estimate is generic in the sense that it does not take into account vanishing near the roots of the functions.

We subsequently evaluate the power series at the midpoint of \( z \) and bound the propagated error using derivatives. Derivatives are bounded using **acb_hypgeomairybound**.

### 3.11.7 Incomplete gamma functions

**void acb_hypgeomgammaupperasymp** *(acb_t res, const acb_t s, const acb_t z, int modified, slong prec)*

**void acb_hypgeomgammaupper1fla** *(acb_t res, const acb_t s, const acb_t z, int modified, slong prec)*

**void acb_hypgeomgammaupper1flb** *(acb_t res, const acb_t s, const acb_t z, int modified, slong prec)*
void **acb_hypgeom_gamma_upper_singular**(*acb_t* res, slong s, const *acb_t* z, int modified, slong prec)

void **acb_hypgeom_gamma_upper**(*acb_t* res, const *acb_t* s, const *acb_t* z, int modified, slong prec)

Computes the upper incomplete gamma function respectively using

\[
\Gamma(s, z) = e^{-z} U(1 - s, 1 - s, z) \\
\Gamma(s, z) = \Gamma(s) - \frac{z^s}{s} F_1(s, s + 1, -z) \\
\Gamma(s, z) = \Gamma(s) - \frac{z^s e^{-z}}{s} F_1(1, s + 1, z) \\
\Gamma(s, z) = \frac{(-1)^n}{n!}(\psi(n + 1) - \log(z)) + \frac{(-1)^n}{(n + 1)!} z^2 F_2(1, 1, 2 + n, -z) - z^{-n} \sum_{k=0}^{n-1} \frac{(-z)^k}{(k - n)k!}, \quad n = -s \in \mathbb{Z}_{\geq 0}
\]

and an automatic algorithm choice. The automatic version also handles other special input such as \(z = 0\) and \(s = 1, 2, 3\). The *singular* version evaluates the finite sum directly and therefore assumes that \(s\) is not too large. If *modified* is set, computes the exponential integral \(z^{-s}\Gamma(s, z) = E_1(z)\) instead.

### 3.11.8 Exponential and trigonometric integrals

The branch cut conventions of the following functions match Mathematica.

void **acb_hypgeom_expint** (*acb_t* res, const *acb_t* s, const *acb_t* z, slong prec)

Computes the generalized exponential integral \(E_s(z)\). This is a trivial wrapper of **acb_hypgeom_gamma_upper()**.

void **acb_hypgeom_ei_asym** (*acb_t* res, const *acb_t* z, slong prec)

void **acb_hypgeom_ei_2f2** (*acb_t* res, const *acb_t* z, slong prec)

void **acb_hypgeom_ei** (*acb_t* res, const *acb_t* z, slong prec)

Computes the exponential integral \(Ei(z)\), respectively using

\[
Ei(z) = -e^z U(1, 1, -z) - \log(-z) + \frac{1}{2} \left( \log(z) - \log \left( \frac{1}{z} \right) \right) \\
Ei(z) = z^2 F_2(1, 1; 2 + z) + \gamma + \frac{1}{2} \left( \log(z) - \log \left( \frac{1}{z} \right) \right)
\]

and an automatic algorithm choice.

void **acb_hypgeom_si_asym** (*acb_t* res, const *acb_t* z, slong prec)

void **acb_hypgeom_si_1f2** (*acb_t* res, const *acb_t* z, slong prec)

void **acb_hypgeom_si** (*acb_t* res, const *acb_t* z, slong prec)

Computes the sine integral \(Si(z)\), respectively using

\[
Si(z) = \frac{i}{2} \left[ e^{iz} U(1, 1, -iz) - e^{-iz} U(1, 1, iz) + \log(-iz) - \log(iz) \right] \\
Si(z) = z_1 F_2 \left( \frac{1}{2}; \frac{3}{2}; \frac{3}{2}; -\frac{z^2}{4} \right)
\]

and an automatic algorithm choice.

void **acb_hypgeom_ci_asym** (*acb_t* res, const *acb_t* z, slong prec)

void **acb_hypgeom_ci_2f3** (*acb_t* res, const *acb_t* z, slong prec)
void `acb_hypgeom_ci` (acb_t res, const acb_t z, slong prec)

Computes the cosine integral \( \text{Ci}(z) \), respectively using
\[
\text{Ci}(z) = \log(z) - \frac{1}{2} \left[ e^{iz}U(1, 1, -iz) + e^{-iz}U(1, 1, iz) + \log(-iz) + \log(iz) \right]
\]
\[
\text{Ci}(z) = -\frac{z^2}{4} F_3(1, 1; 2, 2, \frac{3}{2}; -\frac{z^2}{4}) + \log(z) + \gamma
\]

and an automatic algorithm choice.

void `acb_hypgeom_shi` (acb_t res, const acb_t z, slong prec)

Computes the hyperbolic sine integral \( \text{Shi}(z) = -i \text{Si}(iz) \). This is a trivial wrapper of `acb_hypgeom_si()`.

void `acb_hypgeom_chi_asymp` (acb_t res, const acb_t z, slong prec)

void `acb_hypgeom_chi_2f3` (acb_t res, const acb_t z, slong prec)

void `acb_hypgeom_chi` (acb_t res, const acb_t z, slong prec)

Computes the hyperbolic cosine integral \( \text{Chi}(z) \), respectively using
\[
\text{Chi}(z) = -\frac{1}{2} \left[ e^z U(1, 1, -z) + e^{-z} U(1, 1, z) + \log(-z) - \log(z) \right]
\]
\[
\text{Chi}(z) = \frac{z^2}{4} F_3(1, 1; 2, 2, \frac{3}{2}; \frac{z^2}{4}) + \log(z) + \gamma
\]

and an automatic algorithm choice.

void `acb_hypgeom_li` (acb_t res, const acb_t z, int offset, slong prec)

If offset is zero, computes the logarithmic integral \( \text{li}(z) = \text{Ei}(\log(z)) \).

If offset is nonzero, computes the offset logarithmic integral \( \text{Li}(z) = \text{li}(z) - \text{li}(2) \).

### 3.11.9 Gauss hypergeometric function

The following methods compute the Gauss hypergeometric function
\[
F(z) = _2F_1(a, b, c, z) = \sum_{k=0}^{\infty} \frac{(a)_k(b)_k}{(c)_k} \frac{z^k}{k!}
\]
or the regularized version \( F(z) = F(a, b, c, z) = _2F_1(a, b, c, z) / \Gamma(c) \) if the flag `regularized` is set.

void `acb_hypgeom_2f1_continuation` (acb_t res0, acb_t res1, const acb_t a, const acb_t b, const acb_t c, const acb_t z0, const acb_t z1, const acb_t f0, const acb_t f1, slong prec)

Given \( F(z_0), F'(z_0) \) in \( f0, f1 \), sets \( \text{res0} \) and \( \text{res1} \) to \( F(z_1), F'(z_1) \) by integrating the hypergeometric differential equation along a straight-line path. The evaluation points should be well-isolated from the singular points 0 and 1.

void `acb_hypgeom_2f1_series_direct` (acb_poly_t res, const acb_poly_t a, const acb_poly_t b, const acb_poly_t c, const acb_poly_t z, int `regularized`, slong `len`, slong prec)

Computes \( F(z) \) of the given power series truncated to length \( \text{len} \), using direct summation of the hypergeometric series.

void `acb_hypgeom_2f1_direct` (acb_t res, const acb_t a, const acb_t b, const acb_t c, const acb_t z, int `regularized`, slong `prec`)

Computes \( F(z) \) using direct summation of the hypergeometric series.

void `acb_hypgeom_2f1_transform` (acb_t res, const acb_t a, const acb_t b, const acb_t c, const acb_t z, int `regularized`, int `which`, slong `prec`)

---

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void \texttt{acb\_hypgeom\_2f1\_transform\_limit} \((acb\_t \text{ res}, \text{ const } acb\_t \text{ a}, \text{ const } acb\_t \text{ b}, \text{ const } acb\_t \text{ c}, \text{ const } acb\_t \text{ z}, \text{ int } \text{ regularized}, \text{ int } \text{ which}, \text{ long } \text{ prec})

Computes \( F(z) \) using an argument transformation determined by the flag \texttt{which}. Legal values are 1 for \( z/(z-1) \), 2 for \( 1/z \), 3 for \( 1/(1-z) \), 4 for \( 1-z \), and 5 for \( 1-1/z \).

The \texttt{limit} version assumes that \texttt{which} is not 1. If \texttt{which} is 2 or 3, it assumes that \( b-a \) represents an exact integer. If \texttt{which} is 4 or 5, it assumes that \( c-a-b \) represents an exact integer. In these cases, it computes the correct limit value.

void \texttt{acb\_hypgeom\_2f1\_corner} \((acb\_t \text{ res}, \text{ const } acb\_t \text{ a}, \text{ const } acb\_t \text{ b}, \text{ const } acb\_t \text{ c}, \text{ const } acb\_t \text{ z}, \text{ int } \text{ regularized}, \text{ long } \text{ prec})

Computes \( F(z) \) near the corner cases \( \exp(\pm \pi i \sqrt{3}) \) by analytic continuation.

\begin{verbatim}
int \texttt{acb\_hypgeom\_2f1\_choose} \((\text{ const } acb\_t \text{ z})

Chooses a method to compute the function based on the location of \( z \) in the complex plane. If the return value is 0, direct summation should be used. If the return value is 1 to 5, the transformation with this index in \texttt{acb\_hypgeom\_2f1\_transform()} should be used. If the return value is 6, the corner case algorithm should be used.

void \texttt{acb\_hypgeom\_2f1} \((acb\_t \text{ res}, \text{ const } acb\_t \text{ a}, \text{ const } acb\_t \text{ b}, \text{ const } acb\_t \text{ c}, \text{ const } acb\_t \text{ z}, \text{ int } \text{ regularized}, \text{ long } \text{ prec})

Computes \( F(z) \) (or \( F(z) \) if \texttt{regularized} is set) using an automatic algorithm choice.
\end{verbatim}

### 3.11.10 Orthogonal polynomials and functions

void \texttt{acb\_hypgeom\_chebyshev\_t} \((acb\_t \text{ res}, \text{ const } acb\_t \text{ n}, \text{ const } acb\_t \text{ z}, \text{ long } \text{ prec})

void \texttt{acb\_hypgeom\_chebyshev\_u} \((acb\_t \text{ res}, \text{ const } acb\_t \text{ n}, \text{ const } acb\_t \text{ z}, \text{ long } \text{ prec})

Computes the Chebyshev polynomial (or Chebyshev function) of first or second kind

\[
T_n(z) = 2 \, {}_2F_1\left( -n, n; \frac{1}{2}; \frac{1-z}{2} \right)
\]

\[
U_n(z) = (n+1) \, {}_2F_1\left( -n, n+2; \frac{3}{2}; \frac{1-z}{2} \right).
\]

The hypergeometric series definitions are only used for computation near the point 1. In general, trigonometric representations are used. For word-size integer \( n \), \texttt{acb\_chebyshev\_t\_ui()} and \texttt{acb\_chebyshev\_u\_ui()} are called.

void \texttt{acb\_hypgeom\_jacobi\_p} \((acb\_t \text{ res}, \text{ const } acb\_t \text{ n}, \text{ const } acb\_t \text{ a}, \text{ const } acb\_t \text{ b}, \text{ const } acb\_t \text{ z}, \text{ long } \text{ prec})

Computes the Jacobi polynomial (or Jacobi function)

\[
P_{n}^{(a,b)}(z) = \frac{(a+1)n}{\Gamma(n+1)} \, {}_2F_1\left( -a, a+b+1; n+1; \frac{1-z}{2} \right)
\]

For nonnegative integer \( n \), this is a polynomial in \( a, b \) and \( z \), even when the parameters are such that the hypergeometric series is undefined. In such cases, the polynomial is evaluated using direct methods.

void \texttt{acb\_hypgeom\_gegenbauer\_c} \((acb\_t \text{ res}, \text{ const } acb\_t \text{ n}, \text{ const } acb\_t \text{ m}, \text{ const } acb\_t \text{ z}, \text{ long } \text{ prec})

Computes the Gegenbauer polynomial (or Gegenbauer function)

\[
C_{n}^{m}(z) = \frac{(2m)_n}{\Gamma(n+1)} \, {}_2F_1\left( -2m, m+n; m+\frac{1}{2}; \frac{1-z}{2} \right).
\]

For nonnegative integer \( n \), this is a polynomial in \( m \) and \( z \), even when the parameters are such that the hypergeometric series is undefined. In such cases, the polynomial is evaluated using direct methods.

### 3.11. \texttt{acb\_hypgeom.h} – hypergeometric functions in the complex numbers
void \texttt{acb\_hypgeom\_laguerre\_l} (\texttt{acb\_t res}, const \texttt{acb\_t n}, const \texttt{acb\_t m}, const \texttt{acb\_t z}, \texttt{slong prec})

Computes the Laguerre polynomial (or Laguerre function)

\[ L^m_n(z) = \frac{(m + 1)_n}{\Gamma(n + 1)} F_1 (-n, m + 1, z). \]

For nonnegative integer \( n \), this is a polynomial in \( m \) and \( z \), even when the parameters are such that the hypergeometric series is undefined. In such cases, the polynomial is evaluated using direct methods.

There are at least two incompatible ways to define the Laguerre function when \( n \) is a negative integer. One possibility when \( m = 0 \) is to define \( L^m_n(z) = e^z L^m_{n-1}(-z) \). Another possibility is to cover this case with the recurrence relation \( L^m_{n-1}(z) + L^{m-1}_n(z) = L^n_m(z) \). Currently, we leave this case undefined (returning indeterminate).

void \texttt{acb\_hypgeom\_hermite\_h} (\texttt{acb\_t res}, const \texttt{acb\_t n}, const \texttt{acb\_t z}, \texttt{slong prec})

Computes the Hermite polynomial (or Hermite function)

\[ H_n(z) = 2^n \sqrt{\pi} \left( \frac{1}{\Gamma((1-n)/2)} F_1 \left( \frac{n}{2}, \frac{1}{2}, z^2 \right) - \frac{2z}{\Gamma(-n/2)} F_1 \left( \frac{1-n}{2}, \frac{3}{2}, z^2 \right) \right). \]

void \texttt{acb\_hypgeom\_legendre\_p} (\texttt{acb\_t res}, const \texttt{acb\_t n}, const \texttt{acb\_t m}, const \texttt{acb\_t z}, \texttt{int type}, \texttt{slong prec})

Sets \( \texttt{res} \) to the associated Legendre function of the first kind evaluated for degree \( n \), order \( m \), and argument \( z \). When \( m \) is zero, this reduces to the Legendre polynomial \( P_n(z) \).

Many different branch cut conventions appear in the literature. If \( \texttt{type} \) is 0, the version

\[ P^m_n(z) = \frac{(1 + z)^{m/2}}{(1 - z)^{m/2}} F \left( -n, n + 1, 1 - m, \frac{1 - z}{2} \right) \]

is computed, and if \( \texttt{type} \) is 1, the alternative version

\[ P^m_n(z) = \frac{(z + 1)^{m/2}}{(z - 1)^{m/2}} F \left( -n, n + 1, 1 - m, \frac{1 - z}{2} \right). \]

is computed. Type 0 and type 1 respectively correspond to type 2 and type 3 in \textit{Mathematica} and \textit{mpmath}.

void \texttt{acb\_hypgeom\_legendre\_q} (\texttt{acb\_t res}, const \texttt{acb\_t n}, const \texttt{acb\_t m}, const \texttt{acb\_t z}, \texttt{int type}, \texttt{slong prec})

Sets \( \texttt{res} \) to the associated Legendre function of the second kind evaluated for degree \( n \), order \( m \), and argument \( z \). When \( m \) is zero, this reduces to the Legendre function \( Q_n(z) \).

Many different branch cut conventions appear in the literature. If \( \texttt{type} \) is 0, the version

\[ Q^m_n(z) = \frac{\pi}{2 \sin(\pi m)} \left( \cos(\pi m) P^m_n(z) - \frac{\Gamma(1 + m + n)}{\Gamma(1 - m + n)} P^{-m}_n(z) \right) \]

is computed, and if \( \texttt{type} \) is 1, the alternative version

\[ Q^m_n(z) = \frac{\pi}{2 \sin(\pi m)} e^{\pi i m} \left( P^m_n(z) - \frac{\Gamma(1 + m + n)}{\Gamma(1 - m + n)} P^{-m}_n(z) \right) \]

is computed. Type 0 and type 1 respectively correspond to type 2 and type 3 in \textit{Mathematica} and \textit{mpmath}.

When \( m \) is an integer, either expression is interpreted as a limit. We make use of the connection formulas \([\text{WQ3a}], [\text{WQ3b}]\) and \([\text{WQ3c}]\) to allow computing the function even in the limiting case. (The formula \([\text{WQ3d}]\) would be useful, but is incorrect in the lower half plane.)

void \texttt{acb\_hypgeom\_legendre\_p\_uiui\_rec} (\texttt{acb\_t res}, \texttt{slong n}, \texttt{slong m}, const \texttt{acb\_t z}, \texttt{slong prec})

For nonnegative integer \( n \) and \( m \), uses recurrence relations to evaluate \((1 - z^2)^{-m/2} P^m_n(z)\) which is a polynomial in \( z \).
void \texttt{acb\_hypgeom\_spherical\_y} \( (acb\_t \text{ res, slong } n, \text{ slong } m, \text{ const } acb\_t \theta, \text{ const } acb\_t \phi, \text{ slong } \text{ prec}) \)

Computes the spherical harmonic of degree \( n \), order \( m \), latitude angle \( \theta \), and longitude angle \( \phi \), normalized such that

\[
Y_n^m(\theta, \phi) = \sqrt{\frac{2n + 1}{4\pi}} \frac{(n - m)!}{(n + m)!} e^{im\phi} P_n^m(\cos(\theta)).
\]

The definition is extended to negative \( m \) and \( n \) by symmetry. This function is a polynomial in \( \cos(\theta) \) and \( \sin(\theta) \). We evaluate it using \texttt{acb\_hypgeom\_legendre\_p\_uiui\_rec()}.

### 3.12 acb\_modular.h – modular forms in the complex numbers

This module provides methods for numerical evaluation of modular forms, Jacobi theta functions, and elliptic functions.

In the context of this module, \( \tau \) or \( \tau \) always denotes an element of the complex upper half-plane \( \mathbb{H} = \{ z \in \mathbb{C} : \text{Im}(z) > 0 \} \). We also often use the variable \( q \), variously defined as \( q = e^{2\pi i \tau} \) (usually in relation to modular forms) or \( q = e^{\pi i \tau} \) (usually in relation to theta functions) and satisfying \( |q| < 1 \). We will clarify the local meaning of \( q \) every time such a quantity appears as a function of \( \tau \).

As usual, the numerical functions in this module compute strict error bounds: if \( \tau \) is represented by an \texttt{acb\_t} whose content overlaps with the real line (or lies in the lower half-plane), and \( \tau \) is passed to a function defined only on \( \mathbb{H} \), then the output will have an infinite radius. The analogous behavior holds for functions requiring \( |q| < 1 \).

#### 3.12.1 The modular group

\texttt{psl2z\_struct}

\texttt{psl2z\_t}

Represents an element of the modular group \( \text{PSL}(2, \mathbb{Z}) \), namely an integer matrix

\[
\begin{pmatrix}
    a & b \\
    c & d
\end{pmatrix}
\]

with \( ad - bc = 1 \), and with signs canonicalized such that \( c \geq 0 \), and \( d > 0 \) if \( c = 0 \). The struct members \( a, b, c, d \) are of type \texttt{fmpz}.

void \texttt{psl2z\_init} \( (\text{psl2z\_t} \ g) \)

Initializes \( g \) and set it to the identity element.

void \texttt{psl2z\_clear} \( (\text{psl2z\_t} \ g) \)

Clears \( g \).

void \texttt{psl2z\_swap} \( (\text{psl2z\_t} \ f, \text{psl2z\_t} \ g) \)

Swaps \( f \) and \( g \) efficiently.

void \texttt{psl2z\_set} \( (\text{psl2z\_t} \ f, \text{const} \text{psl2z\_t} \ g) \)

Sets \( f \) to a copy of \( g \).

void \texttt{psl2z\_one} \( (\text{psl2z\_t} \ g) \)

Sets \( g \) to the identity element.

int \texttt{psl2z\_is\_one} \( (\text{const} \text{psl2z\_t} \ g) \)

Returns nonzero iff \( g \) is the identity element.

void \texttt{psl2z\_print} \( (\text{const} \text{psl2z\_t} \ g) \)

Prints \( g \) to standard output.
### 3.12.2 Modular transformations

- **void acb_modular_transform** (acb_t w, const psl2z_t g, const acb_t z, slong prec)
  
  Applies the modular transformation \( g \) to the complex number \( z \), evaluating
  
  \[
  w = g z = \frac{az + b}{cz + d}.
  \]

- **void acb_modular_fundamental_domain_approx_d** (psl2z_t g, double x, double y, double one_minus_eps)

  Attempts to determine a modular transformation \( g \) that maps the complex number \( x + yi \) to the fundamental domain or just slightly outside the fundamental domain, where the target tolerance (not a strict bound) is specified by \( \text{one\_minus\_eps} \).

  The inputs are assumed to be finite numbers, with \( y \) positive.

  Uses floating-point iteration, repeatedly applying either the transformation \( z \leftarrow z + b \) or \( z \leftarrow -1/z \). The iteration is terminated if \( |x| \leq 1/2 \) and \( x^2 + y^2 \geq 1 - \epsilon \) where \( 1 - \epsilon \) is passed as \( \text{one\_minus\_eps} \). It is also terminated if too many steps have been taken without convergence, or if the numbers end up too large or too small for the working precision.

  The algorithm can fail to produce a satisfactory transformation. The output \( g \) is always set to some correct modular transformation, but it is up to the user to verify a posteriori that \( g \) maps \( x + yi \) close enough to the fundamental domain.

- **void acb_modular_fundamental_domain_approx** (acb_t w, psl2z_t g, const acb_t z, const acb_t one_minus_eps, slong prec)

  Attempts to determine a modular transformation \( g \) that maps the complex number \( z \) to the fundamental domain or just slightly outside the fundamental domain, where the target tolerance (not a strict bound) is specified by \( \text{one\_minus\_eps} \). It also computes the transformed value \( w = g z \).

  This function first tries to use **acb_modular_fundamental_domain_approx_d()** and checks if the result is acceptable. If this fails, it calls **acb_modular_fundamental_domain_approx_arf()** with higher precision. Finally, \( w = g z \) is evaluated by a single application of \( g \).

  The algorithm can fail to produce a satisfactory transformation. The output \( g \) is always set to some correct modular transformation, but it is up to the user to verify a posteriori that \( w \) is close enough to the fundamental domain.

- **int acb_modular_is_in_fundamental_domain** (acb_t z, const arf_t tol, slong prec)

  Returns nonzero if it is certainly true that \( |z| \geq 1 - \epsilon \) and \( |\text{Re}(z)| \leq 1/2 + \epsilon \) where \( \epsilon \) is specified by \( \text{tol} \). Returns zero if this is false or cannot be determined.
### 3.12.3 Addition sequences

void **acb_modular_fill_addseq** (slong *tab, slong len)

Builds a near-optimal addition sequence for a sequence of integers which is assumed to be reasonably dense.

As input, the caller should set each entry in `tab` to −1 if that index is to be part of the addition sequence, and to 0 otherwise. On output, entry `i` in `tab` will either be zero (if the number is not part of the sequence), or a value `j` such that both `j` and `i − j` are also marked. The first two entries in `tab` are ignored (the number 1 is always assumed to be part of the sequence).

### 3.12.4 Jacobi theta functions

Unfortunately, there are many inconsistent notational variations for Jacobi theta functions in the literature. Unless otherwise noted, we use the functions

\[
\theta_1(z, \tau) = -i \sum_{n=-\infty}^{\infty} (-1)^n \exp((n + 1/2)^2 \tau + (2n + 1)z) = 2q_{1/4} \sum_{n=0}^{\infty} (-1)^n q^{n(1+1)} \sin((2n + 1)\pi z)
\]

where \(q = \exp(\pi i \tau)\) and \(q_{1/4} = \exp(\pi i \tau/4)\). Note that many authors write \(q_{1/4}\) as \(q^{1/4}\), but the principal fourth root \((q^{1/4} = \exp(\frac{1}{4} \log q)\) differs from \(q_{1/4}\) in general and some formulas are only correct if one reads \(q^{1/4} = \exp(\pi i \tau/4)\). To avoid confusion, we only write \(q^k\) when \(k\) is an integer.

void **acb_modular_theta_transform** (int *R, int *S, int *C, const psl2z_t i g)

We wish to write a theta function with quasiperiod \(\tau' = g \tau\), given some \(g = (a, b; c, d) \in \text{PSL}(2, \mathbb{Z})\). For \(i = 0, 1, 2, 3\), this function computes integers \(R_i\) and \(S_i\) \((R \text{ and } S \text{ should be arrays of length } 4)\) and \(C \in \{0, 1\}\) such that

\[
\theta_{i+1}(z, \tau) = \exp(\pi i R_i/4) \cdot A \cdot B \cdot \theta_{i+S_i}(z', \tau')
\]

where \(z' = z, A = B = 1\) if \(C = 0\), and

\[
z' = \frac{-z}{ct+d}, \quad A = \sqrt{\frac{i}{ct+d}}, \quad B = \exp\left(-\pi i c - \frac{z^2}{ct+d}\right)
\]

if \(C = 1\). Note that \(A\) is well-defined with the principal branch of the square root since \(A^2 = i/(ct+d)\) lies in the right half-plane.

Firstly, if \(c = 0\), we have \(\theta_i(z, \tau) = \exp(-\pi ib/4)\theta_i(z, \tau + b)\) for \(i = 1, 2\), whereas \(\theta_3\) and \(\theta_4\) remain unchanged when \(b\) is even and swap places with each other when \(b\) is odd. In this case we set \(C = 0\).

For an arbitrary \(g\) with \(c > 0\), we set \(C = 1\). The general transformations are given by Rademacher [Rad1973]. We need the function \(\theta_{m,n}(z, \tau)\) defined for \(m, n \in \mathbb{Z}\) by (beware of the typos in [Rad1973])

\[
\theta_{0,0}(z, \tau) = \theta_3(z, \tau), \quad \theta_{0,1}(z, \tau) = \theta_4(z, \tau)
\]

\[
\theta_{1,0}(z, \tau) = \theta_2(z, \tau), \quad \theta_{1,1}(z, \tau) = i\theta_1(z, \tau)
\]

\[
\theta_{m+2,n}(z, \tau) = (-1)^n \theta_{m,n}(z, \tau)
\]

\[
\theta_{m,n+2}(z, \tau) = \theta_{m,n}(z, \tau).
\]
Then we may write

\[
\begin{align*}
\theta_1(z, \tau) &= \varepsilon_1 \exp(\pi i R(a, b, c, d) + 1) / 4 \\
\theta_2(z, \tau) &= \varepsilon_2 \exp(\pi i R(a, b, c, d) + (5 + (2 - c)a) / 4) \\
\theta_3(z, \tau) &= \varepsilon_3 \exp(\pi i R(a, b, c, d) + (4 + (c - d - 2)(b - a)) / 4) \\
\theta_4(z, \tau) &= \varepsilon_4 \exp(\pi i R(a, b, c, d) + (3 - (2 + d)b) / 4)
\end{align*}
\]

where \( \varepsilon_i \) is an 8th root of unity. Specifically, if we denote the 24th root of unity in the transformation formula of the Dedekind eta function by \( \varepsilon(a, b, c, d) = \exp(\pi i R(a, b, c, d) / 12) \) (see [Rad1973] (Rademacher has the transformed/untransformed variables exchanged, and his \( \varepsilon \) differs from ours by a constant offset in the phase).

**void acb_modular_addseq_theta** (slong * exponents, slong * aindex, slong * bindex, slong num)

Constructs an addition sequence for the first \( \vert w \vert \) squares and triangular numbers interleaved (excluding zero), i.e. 1, 2, 4, 6, 9, 12, 16, 20, 25, 30 etc.

**void acb_modular_theta_sum** (acb_ptr theta1, acb_ptr theta2, acb_ptr theta3, acb_ptr theta4, const acb_t w, int w_is_unit, const acb_t q, slong len, slong prec)

Simultaneously computes the first \( \vert w \vert \) coefficients of each of the formal power series

\[
\begin{align*}
\theta_1(z + x, \tau) / q_1^4 &\in \mathbb{C}[[x]] \\
\theta_2(z + x, \tau) / q_1^4 &\in \mathbb{C}[[x]] \\
\theta_3(z + x, \tau) &\in \mathbb{C}[[x]] \\
\theta_4(z + x, \tau) &\in \mathbb{C}[[x]]
\end{align*}
\]

given \( w = \exp(\pi iz) \) and \( q = \exp(\pi i \tau / 4) \), by summing a finite truncation of the respective theta function series. In particular, with \( \vert w \vert \) equal to 1, computes the respective value of the theta function at the point \( z \). We require \( \vert w \vert \) to be positive. If \( w_is_unit \) is nonzero, \( w \) is assumed to lie on the unit circle, i.e. \( z \) is assumed to be real.

Note that the factor \( q_1^4 \) is removed from \( \theta_1 \) and \( \theta_2 \). To get the true theta function values, the user has to multiply this factor back. This convention avoids unnecessary computations, since the user can compute \( q_1^4 = \exp(\pi i \tau / 4) \) followed by \( q = (q_1^4)^{\lambda} \), and in many cases when computing products or quotients of theta functions, the factor \( q_1^4 \) can be eliminated entirely.

This function is intended for \( \vert q \vert \ll 1 \). It can be called with any \( q \), but will return useless intervals if convergence is not rapid. For general evaluation of theta functions, the user should only call this function after applying a suitable modular transformation.

We consider the sums together, alternatingly updating \((\theta_1, \theta_2)\) or \((\theta_3, \theta_4)\). For \( k = 0, 1, 2, \ldots \), the powers of \( q \) are \((k + 2)^2 / 4 = 1, 2, 4, 6, 9 \) etc. and the powers of \( w \) are \( \pm (k + 2) = \pm 2, \pm 3, \pm 4, \ldots \) etc. The scheme is illustrated by the following table:

\[
\begin{align*}
\theta_1, \theta_2 &\quad q^0 &\quad (w^1 \pm w^{-1}) \\
k = 0 &\quad \theta_3, \theta_4 &\quad q^4 &\quad (w^2 \pm w^{-2}) \\
k = 1 &\quad \theta_1, \theta_2 &\quad q^2 &\quad (w^3 \pm w^{-3}) \\
k = 2 &\quad \theta_3, \theta_4 &\quad q^4 &\quad (w^4 \pm w^{-4}) \\
k = 3 &\quad \theta_1, \theta_2 &\quad q^6 &\quad (w^5 \pm w^{-5}) \\
k = 4 &\quad \theta_3, \theta_4 &\quad q^9 &\quad (w^6 \pm w^{-6}) \\
k = 5 &\quad \theta_1, \theta_2 &\quad q^{12} &\quad (w^7 \pm w^{-7})
\end{align*}
\]
For some integer $N \geq 1$, the summation is stopped just before term $k = N$. Let $Q = |q|$, $W = \max(|w|, |w^{-1}|)$, $E = \lfloor (N + 2)^2/4 \rfloor$ and $F = \lfloor (N + 1)/2 \rfloor + 1$. The error of the zeroth derivative can be bounded as

$$2Q^E W^{N+2} \left[ 1 + Q^F W + Q^{2F} W^2 + \ldots \right] = \frac{2Q^E W^{N+2}}{1 - Q^F W}$$

provided that the denominator is positive (otherwise we set the error bound to infinity). When $len$ is greater than 1, consider the derivative of order $r$. The term of index $k$ and order $r$ picks up a factor of magnitude $(k + 2)^r$ from differentiation of $w^{k+2}$ (it also picks up a factor $\pi^r$, but we omit this until we rescale the coefficients at the end of the computation). Thus we have the error bound

$$2Q^E W^{N+2} (N + 2)^r \left[ 1 + Q^F W \frac{(N + 3)^r}{(N + 2)^r} + Q^{2F} W^2 \frac{(N + 4)^r}{(N + 2)^r} + \ldots \right]$$

which by the inequality $(1 + m/(N + 2))^r \leq \exp(mr/(N + 2))$ can be bounded as

$$\frac{2Q^E W^{N+2} (N + 2)^r}{1 - Q^F W \exp(r/(N + 2))}$$

again valid when the denominator is positive.

To actually evaluate the series, we write the even cosine terms as $w^{2n} + w^{-2n}$, the odd cosine terms as $w(w^{2n} + w^{-2n-2})$, and the sine terms as $w(w^{2n} - w^{-2n-2})$. This way we only need even powers of $w$ and $w^{-1}$. The implementation is not yet optimized for real $z$, in which case further work can be saved.

This function does not permit aliasing between input and output arguments.

void acb_modular_theta_const_sum_basecase (acb_t theta2, acb_t theta3, acb_t theta4, const acb_t q, slong N, slong prec)

void acb_modular_theta_const_sum_rs (acb_t theta2, acb_t theta3, acb_t theta4, const acb_t q, slong N, slong prec)

Computes the truncated theta constant sums $\theta_2 = \sum_{k(k+1)<N} q^{k(k+1)}$, $\theta_3 = \sum_{k^2<N} q^{k^2}$, $\theta_4 = \sum_{k^2<N} (-1)^k q^{k^2}$. The basecase version uses a minimal addition sequence. The rs version uses rectangular splitting.

void acb_modular_theta_const_sum (acb_t theta2, acb_t theta3, acb_t theta4, const acb_t q, slong prec)

Computes the respective theta constants by direct summation (without applying modular transformations). This function selects an appropriate $N$, calls either acb_modular_theta_const_sum_basecase() or acb_modular_theta_const_sum_rs() or depending on $N$, and adds a bound for the truncation error.

void acb_modular_theta_notransform (acb_t theta1, acb_t theta2, acb_t theta3, acb_t theta4, const acb_t z, const acb_t tau, slong prec)

Evaluates the Jacobi theta functions $\theta_i(z, \tau), i = 1, 2, 3, 4$ simultaneously. This function does not move $\tau$ to the fundamental domain. This is generally worse than acb_modular_theta(), but can be slightly better for moderate input.

void acb_modular_theta (acb_t theta1, acb_t theta2, acb_t theta3, acb_t theta4, const acb_t z, const acb_t tau, slong prec)

Evaluates the Jacobi theta functions $\theta_i(z, \tau), i = 1, 2, 3, 4$ simultaneously. This function moves $\tau$ to the fundamental domain before calling acb_modular_theta_sum().

### 3.12.5 The Dedekind eta function

void acb_modular_addseq_eta (slong * exponents, slong * aindex, slong * bindex, slong long num)

Constructs an addition sequence for the first $num$ generalized pentagonal numbers (excluding zero), i.e. 1, 2, 5, 7, 12, 15, 22, 26, 35, 40 etc.
void \texttt{acb\_modular\_eta\_sum}(acb\_t eta, const acb\_t q, slong prec)
  Evaluates the Dedekind eta function without the leading 24th root, i.e.
  \[
  \exp(-\pi\iota/12)\eta(\tau) = \sum_{n=-\infty}^{\infty} (-1)^n q^{(3n^2-n)/2}
  \]
given \(q = \exp(2\pi\iota\tau)\), by summing the defining series.

This function is intended for \(|q| \ll 1\). It can be called with any \(q\), but will return useless intervals if convergence is not rapid. For general evaluation of the eta function, the user should only call this function after applying a suitable modular transformation.

int \texttt{acb\_modular\_epsilon\_arg}(const psl2z\_t g)
  Given \(g = (a, b; c, d)\), computes an integer \(R\) such that \(\varepsilon(a, b, c, d) = \exp(\pi\iota R/12)\) is the 24th root of unity in the transformation formula for the Dedekind eta function,
  \[
  \eta\left(\frac{a\tau + b}{c\tau + d}\right) = \varepsilon(a, b, c, d)\sqrt{c\tau + d}\eta(\tau).
  \]

void \texttt{acb\_modular\_eta}(acb\_t r, const acb\_t tau, slong prec)
  Computes the Dedekind eta function \(\eta(\tau)\) given \(\tau\) in the upper half-plane. This function applies the functional equation to move \(\tau\) to the fundamental domain before calling \texttt{acb\_modular\_eta\_sum}().

### 3.12.6 Modular forms

void \texttt{acb\_modular\_j}(acb\_t r, const acb\_t tau, slong prec)
  Computes Klein’s j-invariant \(j(\tau)\) given \(\tau\) in the upper half-plane. The function is normalized so that \(j(i) = 1728\). We first move \(\tau\) to the fundamental domain, which does not change the value of the function. Then we use the formula \(j(\tau) = 32(\theta_2^6 + \theta_3^6 + \theta_4^6)^3 / (\theta_2 \theta_3 \theta_4)^6\) where \(\theta_i = \theta_i(0, \tau)\).

void \texttt{acb\_modular\_lambda}(acb\_t r, const acb\_t tau, slong prec)
  Computes the lambda function \(\lambda(\tau) = \theta_2^4(0, \tau) / \theta_3^4(0, \tau)\), which is invariant under modular transformations \((a, b; c, d)\) where \(a, d\) are odd and \(b, c\) are even.

void \texttt{acb\_modular\_delta}(acb\_t r, const acb\_t tau, slong prec)
  Computes the modular discriminant \(\Delta(\tau) = \eta(\tau)^{24}\), which transforms as
  \[
  \Delta\left(\frac{a\tau + b}{c\tau + d}\right) = (c\tau + d)^{12}\Delta(\tau).
  \]
  The modular discriminant is sometimes defined with an extra factor \((2\pi)^{12}\), which we omit in this implementation.

void \texttt{acb\_modular\_eisenstein}(acb\_ptr r, const acb\_t tau, slong len, slong prec)
  Computes simultaneously the first \(\textsc{len}\) entries in the sequence of Eisenstein series \(G_4(\tau), G_6(\tau), G_8(\tau), \ldots\), defined by
  \[
  G_{2k}(\tau) = \sum_{m^2 + n^2 \neq 0} \frac{1}{(m + n\tau)^{2k}}
  \]
  and satisfying
  \[
  G_{2k}\left(\frac{a\tau + b}{c\tau + d}\right) = (c\tau + d)^{2k}G_{2k}(\tau).
  \]
  We first evaluate \(G_4(\tau)\) and \(G_6(\tau)\) on the fundamental domain using theta functions, and then compute the Eisenstein series of higher index using a recurrence relation.
3.12.7 Elliptic functions

```c
void acb_modular_elliptic_p(acb_t wp, const acb_t z, const acb_t tau, slong prec)
```

Computes Weierstrass’s elliptic function

$$
\wp(z, \tau) = \frac{1}{z^2} + \sum_{n^2 + m^2 \neq 0} \left[ \frac{1}{(z + m + n\tau)^2} - \frac{1}{(m + n\tau)^2} \right]
$$

which satisfies \(\wp(z, \tau) = \wp(z + 1, \tau) = \wp(z + \tau, \tau)\). To evaluate the function efficiently, we use the formula

$$
\wp(z, \tau) = \frac{\pi^2}{2} \frac{\theta_2^4(0, \tau)}{\theta_4^4(0, \tau)} - \frac{\pi^2}{3} \left[ \theta_3^4(0, \tau) + \theta_4^4(0, \tau) \right].
$$

```c
void acb_modular_elliptic_p_zpx(acb_ptr wp, const acb_t z, const acb_t tau, slong len, slong prec)
```

Computes the formal power series \(\wp(z + x, \tau) \in \mathbb{C}[[x]]\), truncated to length \(\text{len}\). In particular, with \(\text{len} = 2\), simultaneously computes \(\wp(z, \tau), \wp'(z, \tau)\) which together generate the field of elliptic functions with periods 1 and \(\tau\).

3.12.8 Elliptic integrals

```c
void acb_modular_elliptic_k(acb_t w, const acb_t m, slong prec)
```

Computes the complete elliptic integral of the first kind \(K(m)\), using the arithmetic-geometric mean:

$$
K(m) = \frac{\pi}{2M(\sqrt{1 - m})}.
$$

```c
void acb_modular_elliptic_k_cpx(acb_ptr w, const acb_t m, slong len, slong prec)
```

Sets the coefficients in the array \(w\) to the power series expansion of the complete elliptic integral of the first kind at the point \(m\) truncated to length \(\text{len}\), i.e. \(K(m + x) \in \mathbb{C}[[x]]\).

```c
void acb_modular_elliptic_e(acb_t w, const acb_t m, slong prec)
```

Computes the complete elliptic integral of the second kind \(E(m)\), which is given by \(E(m) = (1 - m)(2mK'(m) + K(m))\) (where the prime denotes a derivative, not a complementary integral).

3.12.9 Class polynomials

```c
void acb_modular_hilbert_class_poly(fmpz_poly_t res, slong D)
```

Sets \(res\) to the Hilbert class polynomial of discriminant \(D\), defined as

$$
H_D(x) = \prod_{(a,b,c)} \left( x - j \left( \frac{-b + \sqrt{D}}{2a} \right) \right)
$$

where \((a, b, c)\) ranges over the primitive reduced positive definite binary quadratic forms of discriminant \(b^2 - 4ac = D\).

The Hilbert class polynomial is only defined if \(D < 0\) and \(D\) is congruent to 0 or 1 mod 4. If some other value of \(D\) is passed as input, \(res\) is set to the zero polynomial.

3.13 bernoulli.h – support for Bernoulli numbers

This module provides helper functions for exact or approximate calculation of the Bernoulli numbers, which are defined by the exponential generating function

$$
\frac{x}{e^x - 1} = \sum_{n=0}^{\infty} B_n \frac{x^n}{n!}.
$$
Efficient algorithms are implemented for both multi-evaluation and calculation of isolated Bernoulli numbers. A global (or thread-local) cache is also provided, to support fast repeated evaluation of various special functions that depend on the Bernoulli numbers (including the gamma function and the Riemann zeta function).

### 3.13.1 Generation of Bernoulli numbers

**bernoulli_rev_t**

An iterator object for generating a range of even-indexed Bernoulli numbers exactly in reverse order, i.e. computing the exact fractions $B_n, B_{n-2}, B_{n-4}, \ldots, B_0$. The Bernoulli numbers are generated from scratch, i.e. no caching is performed.

The Bernoulli numbers are computed by direct summation of the zeta series. This is made fast by storing a table of powers (as done by Bloemen et al. [http://remcobloemen.nl/2009/11/even-faster-zeta-calculation.html](http://remcobloemen.nl/2009/11/even-faster-zeta-calculation.html)). As an optimization, we only include the odd powers, and use fixed-point arithmetic.

The reverse iteration order is preferred for performance reasons, as the powers can be updated using multiplications instead of divisions, and we avoid having to periodically recompute terms to higher precision. To generate Bernoulli numbers in the forward direction without having to store all of them, one can split the desired range into smaller blocks and compute each block with a single reverse pass.

```c
void bernoulli_rev_init (bernoulli_rev_t iter, ulong n)
    Initializes the iterator iter. The first Bernoulli number to be generated by calling bernoulli_rev_next() is $B_n$. It is assumed that $n$ is even.

void bernoulli_rev_next (fmpz_t numer, fmpz_t denom, bernoulli_rev_t iter)
    Sets numer and denom to the exact, reduced numerator and denominator of the Bernoulli number $B_k$ and advances the state of iter so that the next invocation generates $B_{k-2}$.

void bernoulli_rev_clear (bernoulli_rev_t iter)
    Frees all memory allocated internally by iter.
```

### 3.13.2 Caching

**slong bernoulli_cache_num**

**fmpz * bernoulli_cache**

Cache of Bernoulli numbers. Uses thread-local storage if enabled in FLINT.

```c
void bernoulli_cache_compute (slong n)
    Makes sure that the Bernoulli numbers up to at least $B_{n-1}$ are cached. Calling flint_cleanup() frees the cache.
```

### 3.13.3 Bounding

**slong bernoulli_bound_2exp_si** (ulong n)

Returns an integer $b$ such that $|B_n| \leq 2^b$. Uses a lookup table for small $n$, and for larger $n$ uses the inequality $|B_n| < 4n!/(2\pi)^n < 4(n + 1)^{n+1}e^{-n}/(2\pi)^n$. Uses integer arithmetic throughout, with the bound for the logarithm being looked up from a table. If $|B_n| = 0$, returns LONG_MIN. Otherwise, the returned exponent $b$ is never more than one percent larger than the true magnitude.

This function is intended for use when $n$ small enough that one might comfortably compute $B_n$ exactly. It aborts if $n$ is so large that internal overflow occurs.

```c
void _bernoulli_fmpq_ui_zeta (fmpz_t num, fmpz_t den, ulong n)
    Sets num and den to the reduced numerator and denominator of the Bernoulli number $B_n$.
```
This function computes the denominator $d$ using von Staudt-Clausen theorem, numerically approximates $B_n$ using `arb_bernoulli_ui_zeta()`, and then rounds $dB_n$ to the correct numerator. If the working precision is insufficient to determine the numerator, the function prints a warning message and retries with increased precision (this should not be expected to happen).

```c
void _bernoulli_fmpq_ui(fmpz_t num, fmpz_t den, ulong n)
void bernoulli_fmpq_ui(fmpq_t b, ulong n)
```
Computes the Bernoulli number $B_n$ as an exact fraction, for an isolated integer $n$. This function reads $B_n$ from the global cache if the number is already cached, but does not automatically extend the cache by itself.

### 3.14 hypgeom.h – support for hypergeometric series

This module provides functions for high-precision evaluation of series of the form

$$
\sum_{k=0}^{n-1} \frac{A(k)}{B(k)} \prod_{j=1}^{k} \frac{P(j)}{Q(j)} z^k
$$

where $A, B, P, Q$ are polynomials. The present version only supports $A, B, P, Q \in \mathbb{Z}[k]$ (represented using the FLINT `fmpz_poly_t` type). This module also provides functions for high-precision evaluation of infinite series ($n \to \infty$), with automatic, rigorous error bounding.

Note that we can standardize to $A = B = 1$ by setting $\tilde{P}(k) = P(k)A(k)B(k-1), \tilde{Q}(k) = Q(k)A(k-1)B(k)$. However, separating out $A$ and $B$ is convenient and improves efficiency during evaluation.

#### 3.14.1 Strategy for error bounding

We wish to evaluate $S(z) = \sum_{k=0}^{\infty} T(k)z^k$ where $T(k)$ satisfies $T(0) = 1$ and

$$
T(k) = R(k)T(k-1) = \left( \frac{P(k)}{Q(k)} \right) T(k-1)
$$

for given polynomials

$$
P(k) = a_pk^p + a_{p-1}k^{p-1} + \ldots + a_0
$$

$$
Q(k) = b_qk^q + b_{q-1}k^{q-1} + \ldots + b_0.
$$

For convergence, we require $p < q$, or $p = q$ with $|z||a_p| < |b_q|$. We also assume that $P(k)$ and $Q(k)$ have no roots among the positive integers (if there are positive integer roots, the sum is either finite or undefined). With these conditions satisfied, our goal is to find a parameter $n \geq 0$ such that

$$
\left| \sum_{k=n}^{\infty} T(k)z^k \right| \leq 2^{-d}.
$$

We can rewrite the hypergeometric term ratio as

$$
zR(k) = z \frac{P(k)}{Q(k)} = z \left( \frac{a_p}{b_q} \right) \frac{1}{k^{q-p}} F(k)
$$

where

$$
F(k) = \frac{1 + \tilde{a}_1/k + \tilde{a}_2/k^2 + \ldots + \tilde{a}_q/k^p}{1 + b_1/k + b_2/k^2 + \ldots + b_q/k^q} = 1 + O(1/k)
$$
and where \( \tilde{a}_i = a_{p-i}/a_p \) and \( \tilde{b}_i = b_{q-i}/b_q \). Next, we define
\[
C = \max_{1 \leq i \leq p} |\tilde{a}_i|^{1/i}, \quad D = \max_{1 \leq i \leq q} |\tilde{b}_i|^{1/i}.
\]

Now, if \( k > C \), the magnitude of the numerator of \( F(k) \) is bounded from above by
\[
1 + \sum_{i=1}^{p} \left( \frac{C}{k} \right)^i \leq 1 + \frac{C}{k-C}
\]
and if \( k > 2D \), the magnitude of the denominator of \( F(k) \) is bounded from below by
\[
1 - \sum_{i=1}^{q} \left( \frac{D}{k} \right)^i \geq 1 + \frac{D}{D-k}.
\]

Putting the inequalities together gives the following bound, valid for \( k > K = \max(C, 2D) \):
\[
|F(k)| \leq \frac{k(k-D)}{(k-C)(k-2D)} = \left( 1 + \frac{C}{k-C} \right) \left( 1 + \frac{D}{k-2D} \right).
\]

Let \( r = q - p \) and \( \tilde{z} = |z a_p / b_q| \). Assuming \( k > \max(C, 2D, \tilde{z}^{1/r}) \), we have
\[
|z R(k)| \leq G(k) = \frac{\tilde{z} F(k)}{k^r}
\]
where \( G(k) \) is monotonically decreasing. Now we just need to find an \( n \) such that \( G(n) < 1 \) and for which \( |T(n)|/(1-G(n)) \leq 2^{-d} \). This can be done by computing a floating-point guess for \( n \) then trying successively larger values.

This strategy leaves room for some improvement. For example, if \( \tilde{b}_1 \) is positive and large, the bound \( B \) becomes very pessimistic (a larger positive \( \tilde{b}_1 \) causes faster convergence, not slower convergence).

### 3.14.2 Types, macros and constants

**hypgeom struct**

**hypgeom_t**

Stores polynomials \( A, B, P, Q \) and precomputed bounds, representing a fixed hypergeometric series.

### 3.14.3 Memory management

void **hypgeom_init** (hypgeom_t hyp)

void **hypgeom_clear** (hypgeom_t hyp)

### 3.14.4 Error bounding

slong **hypgeom_estimate_terms** (const mag_t z, int r, slong d)

Computes an approximation of the largest \( n \) such that \( |z|^n/(n!)^r = 2^{-d} \), giving a first-order estimate of the number of terms needed to approximate the sum of a hypergeometric series of weight \( r \geq 0 \) and argument \( z \) to an absolute precision of \( d \geq 0 \) bits. If \( r = 0 \), the direct solution of the equation is given by \( n = (\log(1-z) - d \log 2)/\log z \). If \( r > 0 \), using \( \log n! \approx n \log n - n \) gives an equation that can be solved in terms of the Lambert \( W \)-function as \( n = (d \log 2)/(r W(t)) \) where \( t = (d \log 2)/(e z^{1/r}) \).

The evaluation is done using double precision arithmetic. The function aborts if the computed value of \( n \) is greater than or equal to \( \text{LONG\_MAX}/2 \).
slong hypgeom_bound (mag_t error, int r, slong C, slong D, slong K, const mag_t TK, const mag_t z, slong prec)
  Computes a truncation parameter sufficient to achieve prec bits of absolute accuracy, according to the strategy described above. The input consists of r, C, D, K, precomputed bound for T(K), and \( \tilde{z} = z(a_p/b_q) \), such that for \( k > K \), the hypergeometric term ratio is bounded by
  \[
  \frac{\tilde{z}}{k(k-D)} \leq 2^{-\text{prec}}.
  \]
  Given this information, we compute a \( \varepsilon \) and an integer \( n \) such that \( |\sum_{k=n}^{\infty} T(k)| \leq \varepsilon \leq 2^{-\text{prec}} \). The output variable error is set to the value of \( \varepsilon \), and \( n \) is returned.

void hypgeom_precompute (hypgeom_t hyp)
Precomputes the bounds data \( C, D, K \) and an upper bound for \( T(K) \).

### 3.15.5 Summation

void fmprb_hypgeom_sum (fmprb_t P, fmprb_t Q, const hypgeom_t hyp, const slong n, slong prec)
Computes \( P/Q = \sum_{k=0}^{n-1} T(k) \) where \( T(k) \) is defined by hyp, using binary splitting and a working precision of prec bits.

void fmprb_hypgeom_infsum (fmprb_t P, fmprb_t Q, hypgeom_t hyp, slong tol, slong prec)
Computes \( P/Q = \sum_{k=0}^{\infty} T(k) \) where \( T(k) \) is defined by hyp, using binary splitting and working precision of prec bits. The number of terms is chosen automatically to bound the truncation error by at most \( 2^{-\text{tol}} \). The bound for the truncation error is included in the output as part of \( P \).

void arb_hypgeom_sum (arb_t P, arb_t Q, const hypgeom_t hyp, const slong n, slong prec)
Computes \( P/Q = \sum_{k=0}^{n-1} T(k) \) where \( T(k) \) is defined by hyp, using binary splitting and a working precision of prec bits.

void arb_hypgeom_infsum (arb_t P, arb_t Q, hypgeom_t hyp, slong tol, slong prec)
Computes \( P/Q = \sum_{k=0}^{\infty} T(k) \) where \( T(k) \) is defined by hyp, using binary splitting and working precision of prec bits. The number of terms is chosen automatically to bound the truncation error by at most \( 2^{-\text{tol}} \). The bound for the truncation error is included in the output as part of \( P \).

### 3.15 partitions.h – computation of the partition function

This module implements the asymptotically fast algorithm for evaluating the integer partition function \( p(n) \) described in [Joh2012]. The idea is to evaluate a truncation of the Hardy-Ramanujan-Rademacher series using tight precision estimates, and symbolically factoring the occurring exponential sums.

An implementation based on floating-point arithmetic can also be found in FLINT. That version relies on some numerical subroutines that have not been proved correct.

The implementation provided here uses ball arithmetic throughout to guarantee a correct error bound for the numerical approximation of \( p(n) \). Optionally, hardware double arithmetic can be used for low-precision terms. This gives a significant speedup for small (e.g. \( n < 10^6 \)).

void partitions_rademacher_bound (arf_t b, const fmpz_t n, ulong N)
Sets \( b \) to an upper bound for
  \[
  M(n, N) = \frac{44\pi^2}{225\sqrt{3}} N^{-1/2} + \frac{\pi\sqrt{2}}{75} \left( \frac{N}{n-1} \right)^{1/2} \sinh \left( \frac{\pi}{N} \sqrt{\frac{2n}{3}} \right).
  \]

This formula gives an upper bound for the truncation error in the Hardy-Ramanujan-Rademacher formula when the series is taken up to the term \( t(n, N) \) inclusive.
partitions_hrr_sum_arb

Evaluates the partial sum \( \sum_{k=N_0}^{N} t(n, k) \) of the Hardy-Ramanujan-Rademacher series.

If \( \text{use\_doubles} \) is nonzero, doubles and the system's standard library math functions are used to evaluate the smallest terms. This significantly speeds up evaluation for small \( n \) (e.g. \( n < 10^6 \)), and gives a small speed improvement for larger \( n \), but the result is not guaranteed to be correct. In practice, the error is estimated very conservatively, and unless the system's standard library is broken, use of doubles can be considered safe. Setting \( \text{use\_doubles} \) to zero gives a fully guaranteed bound.

void partitions_fmpz_fmpz

Computes the partition function \( p(n) \) using the Hardy-Ramanujan-Rademacher formula. This function computes a numerical ball containing \( p(n) \) and verifies that the ball contains a unique integer.

If \( n \) is sufficiently large and a number of threads greater than 1 has been selected with \( \text{flint\_set\_num\_threads()} \), the computation time will be reduced by using two threads.

See partitions_hrr_sum_arb() for an explanation of the \( \text{use\_doubles} \) option.

void partitions_fmpz_ui

Computes the partition function \( p(n) \) using the Hardy-Ramanujan-Rademacher formula. This function computes a numerical ball containing \( p(n) \) and verifies that the ball contains a unique integer.

void partitions_fmpz_ui_using_doubles

Computes the partition function \( p(n) \), enabling the use of doubles internally. This significantly speeds up evaluation for small \( n \) (e.g. \( n < 10^6 \)), but the error bounds are not certified (see remarks for partitions_hrr_sum_arb()).
4.1 Algorithms for mathematical constants

Most mathematical constants are evaluated using the generic hypergeometric summation code.

4.1.1 Pi

\( \pi \) is computed using the Chudnovsky series

\[
\frac{1}{\pi} = 12 \sum_{k=0}^{\infty} \frac{(-1)^k (6k)! (13591409 + 545140134k)}{(3k)! (k!)^3 640320^{3k+3/2}}
\]

which is hypergeometric and adds roughly 14 digits per term. Methods based on the arithmetic-geometric mean seem to be slower by a factor three in practice.

A small trick is to compute \( \frac{1}{\sqrt{640320}} \) instead of \( \sqrt{640320} \) at the end.

4.1.2 Logarithms of integers

We use the formulas

\[
\log(2) = \frac{3}{4} \sum_{k=0}^{\infty} \frac{(-1)^k (k!)^2}{2^k (2k + 1)!}
\]

\[
\log(10) = 46 \operatorname{atanh}(1/31) + 34 \operatorname{atanh}(1/49) + 20 \operatorname{atanh}(1/161)
\]

4.1.3 Euler’s constant

Euler’s constant \( \gamma \) is computed using the Brent-McMillan formula (\cite{BM1980}, \cite{MPFR2012})

\[
\gamma = \frac{S_0(2n) - K_0(2n)}{I_0(2n)} - \log(n)
\]

in which \( n \) is a free parameter and

\[
S_0(x) = \sum_{k=0}^{\infty} \frac{H_k}{(k!)^2} \left( \frac{x}{2} \right)^{2k}, \quad I_0(x) = \sum_{k=0}^{\infty} \frac{1}{(k!)^2} \left( \frac{x}{2} \right)^{2k}
\]

\[
2x I_0(x) K_0(x) \sim \sum_{k=0}^{\infty} \frac{[(2k)!]^3}{(k!)^4 8^{2k} x^{2k}}.
\]
All series are evaluated using binary splitting. The first two series are evaluated simultaneously, with the summation taken up to \( k = N - 1 \) inclusive where \( N \geq \alpha n + 1 \) and \( \alpha \approx 4.9706257595442318644 \) satisfies \( \alpha(\log \alpha - 1) = 3 \). The third series is taken up to \( k = 2n - 1 \) inclusive. With these parameters, it is shown in [BJ2013] that the error is bounded by \( 24e^{-8n} \).

### 4.1.4 Catalan’s constant

Catalan’s constant is computed using the hypergeometric series

\[
C = \sum_{k=0}^{\infty} \frac{(-1)^k 4^{4k+1} (40k^2 + 56k + 19) \left[(k+1)\right]^2(2k+2)!^3}{(k+1)^3(2k+1)(4k+4)!^2}
\]

### 4.1.5 Khinchin’s constant

Khinchin’s constant \( K_0 \) is computed using the formula

\[
\log K_0 = \frac{1}{\log 2} \left[ \log \left( \frac{k+1}{k} \right) - \sum_{k=2}^{N-1} \log \left( \frac{k-1}{k} \right) \right]
\]

where \( N \geq 2 \) is a free parameter that can be used for tuning [BBC1997]. If the infinite series is truncated after \( n = M \), the remainder is smaller in absolute value than

\[
\sum_{n=M+1}^{\infty} \zeta(2n, N) = \sum_{n=M+1}^{\infty} \sum_{k=0}^{\infty} (k+N)^{-2n} \leq \sum_{n=M+1}^{\infty} \left( N^{-2n} + \int_0^{\infty} (t+N)^{-2n} dt \right)
\]

Thus, for an error of at most \( 2^{-p} \) in the series, it is sufficient to choose \( M \geq p/(2 \log_2 N) \).

### 4.1.6 Glaisher’s constant

Glaisher’s constant \( A = \exp\left(\frac{1}{12} - \zeta'(-1)\right) \) is computed directly from this formula. We don’t use the reflection formula for the zeta function, as the arithmetic in Euler-Maclaurin summation is faster at \( s = -1 \) than at \( s = 2 \).

### 4.1.7 Apery’s constant

Apery’s constant \( \zeta(3) \) is computed using the hypergeometric series

\[
\zeta(3) = \frac{1}{64} \sum_{k=0}^{\infty} (-1)^k (205k^2 + 250k + 77) \frac{(k!)^{10}}{(2k+1)!5}
\]

### 4.2 Algorithms for gamma functions

#### 4.2.1 The Stirling series

In general, the gamma function is computed via the Stirling series

\[
\log \Gamma(z) = \left( z - \frac{1}{2} \right) \log z - z + \frac{\ln 2\pi}{2} + \sum_{k=1}^{n-1} \frac{B_{2k}}{2k(2k-1)z^{2k-1}} + R(n, z)
\]

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where ([Olv1997] pp. 293-295) the remainder term is exactly
\[ R_n(z) = \int_0^\infty \frac{B_{2n} - \hat{B}_{2n}(x)}{2n(x + z)^{2n}} \, dx. \]

To evaluate the gamma function of a power series argument, we substitute \( z \rightarrow z + t \in \mathbb{C}[[t]] \).

Using the bound for \( |x + z| \) given by [Olv1997] and the fact that the numerator of the integrand is bounded in absolute value by \( 2|B_{2n}| \), the remainder can be shown to satisfy the bound
\[ |\{t^k R_n(z + t)\}| \leq 2|B_{2n}| \frac{\Gamma(2n + k - 1)}{\Gamma(k + 1) \Gamma(2n + 1)} |z| (b/|z|)^{2n + k} \]
where \( b = 1/\cos(\arg(z)/2) \). Note that by trigonometric identities, assuming that \( z = x + yi \), we have \( b = \sqrt{1 + t^2} \) where
\[ t = \frac{y}{\sqrt{x^2 + y^2 + x}} = \frac{\sqrt{x^2 + y^2} - x}{y}. \]

To use the Stirling series at \( p \)-bit precision, we select parameters \( r, n \) such that the remainder \( R(n, z) \) approximately is bounded by \( 2^{-p} \). If \( |z| \) is too small for the Stirling series to give sufficient accuracy directly, we first translate to \( z + t \) using the formula \( \Gamma(z) = \Gamma(z + r)/(z(z + 1)(z + 2)\cdots(z + r - 1)) \).

To obtain a remainder smaller than \( 2^{-p} \), we must choose an \( r \) such that, in the real case, \( z + r > \beta p \), where \( \beta > \log(2)/(2\pi) \approx 0.11 \). In practice, a slightly larger factor \( \beta \approx 0.2 \) more closely balances \( n \) and \( r \). A much larger \( \beta \) (e.g., \( \beta = 1 \)) could be used to reduce the number of Bernoulli numbers that have to be precomputed, at the expense of slower repeated evaluation.

### 4.2.2 Rational arguments

We use efficient methods to compute \( y = \Gamma(p/q) \) where \( q \) is one of \( 1, 2, 3, 4, 6 \) and \( p \) is a small integer.

The cases \( \Gamma(1) = 1 \) and \( \Gamma(1/2) = \sqrt{\pi} \) are trivial. We reduce all remaining cases to \( \Gamma(1/3) \) or \( \Gamma(1/4) \) using the following relations:

\[
\begin{align*}
\Gamma(2/3) &= \frac{2\pi}{3^{1/2} \Gamma(1/3)}, \\
\Gamma(3/4) &= \frac{2^{1/2} \pi}{\Gamma(1/4)}, \\
\Gamma(1/6) &= \frac{\Gamma(1/3)^2}{(\pi/3)^{1/2} 2^{1/3}}, \\
\Gamma(5/6) &= \frac{2\pi (\pi/3)^{1/2} 2^{1/3}}{\Gamma(1/3)^2}.
\end{align*}
\]

We compute \( \Gamma(1/3) \) and \( \Gamma(1/4) \) rapidly to high precision using
\[
\begin{align*}
\Gamma(1/3) &= \left( \frac{12\pi^4}{\sqrt{10}} \sum_{k=0}^\infty \frac{(6k)!(-1)^k}{(k!)^3 (3k)! 3^{2k} 160^{3k}} \right)^{1/6}, \\
\Gamma(1/4) &= \sqrt{\frac{(2\pi)^{3/2}}{\text{agm}(1, \sqrt{2})}}.
\end{align*}
\]

An alternative formula which could be used for \( \Gamma(1/3) \) is
\[
\Gamma(1/3) = \frac{2^{4/9} \pi^{2/3}}{3^{1/12} \left( \text{agm} \left(1, \frac{1}{2} \sqrt{2 + \sqrt{3}}\right)\right)^{1/3}},
\]
but this appears to be slightly slower in practice.

### 4.2. Algorithms for gamma functions

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4.3 Algorithms for polylogarithms

The polylogarithm is defined for $s, z \in \mathbb{C}$ with $|z| < 1$ by

$$\operatorname{Li}_s(z) = \sum_{k=1}^{\infty} \frac{z^k}{k^s}$$

and for $|z| \geq 1$ by analytic continuation, except for the singular point $z = 1$.

4.3.1 Computation for small $z$

The power sum converges rapidly when $|z| \ll 1$. To compute the series expansion with respect to $s$, we substitute $s \rightarrow s + x \in \mathbb{C}[[x]]$ and obtain

$$\operatorname{Li}_{s+x}(z) = \sum_{d=0}^{\infty} x^d \frac{(-1)^d}{d!} \sum_{k=1}^{\infty} T(k)$$

where

$$T(k) = \frac{z^k \log^d(k)}{k^s}.$$ 

The remainder term $|\sum_{k=N}^{\infty} T(k)|$ is bounded via mag_polylog_tail().

4.3.2 Expansion for general $z$

For general complex $s, z$, we write the polylogarithm as a sum of two Hurwitz zeta functions

$$\operatorname{Li}_s(z) = \frac{\Gamma(v)}{(2\pi)^v} \left[ i^v \zeta \left( v, \frac{1}{2} + \frac{\log(-z)}{2\pi i} \right) + i^{-v} \zeta \left( v, \frac{1}{2} - \frac{\log(-z)}{2\pi i} \right) \right]$$

in which $s = 1 - v$. With the principal branch of $\log(-z)$, we obtain the conventional analytic continuation of the polylogarithm with a branch cut on $z \in (1, +\infty)$.

To compute the series expansion with respect to $v$, we substitute $v \rightarrow v + x \in \mathbb{C}[[x]]$ in this formula (at the end of the computation, we map $x \rightarrow -x$ to obtain the power series for $\operatorname{Li}_{s+x}(z)$). The right hand side becomes

$$\Gamma(v+x)[E_1 Z_1 + E_2 Z_2]$$

where $E_1 = (i/(2\pi))^{v+x}$, $Z_1 = \zeta(v+x, \ldots)$, $E_2 = (1/(2\pi i))^{v+x}$, $Z_2 = \zeta(v+x, \ldots)$.

When $v = 1$, the $Z_1$ and $Z_2$ terms become Laurent series with a leading $1/x$ term. In this case, we compute the deflated series $\tilde{Z}_1, \tilde{Z}_2 = \zeta(x, \ldots) - 1/x$. Then

$$E_1 Z_1 + E_2 Z_2 = (E_1 + E_2)/x + E_1 \tilde{Z}_1 + E_2 \tilde{Z}_2.$$ 

Note that $(E_1 + E_2)/x$ is a power series, since the constant term in $E_1 + E_2$ is zero when $v = 1$. So we simply compute one extra derivative of both $E_1$ and $E_2$, and shift them one step. When $v = 0, -1, -2, \ldots$, the $\Gamma(v+x)$ prefactor has a pole. In this case, we proceed analogously and formally multiply $x \Gamma(v+x)$ with $[E_1 Z_1 + E_2 Z_2]/x$.

Note that the formal cancellation only works when the order $s$ (or $v$) is an exact integer: it is not currently possible to use this method when $s$ is a small ball containing any of $0, 1, 2, \ldots$ (then the result becomes indeterminate).

The Hurwitz zeta method becomes inefficient when $|z| \rightarrow 0$ (it gives an indeterminate result when $z = 0$). This is not a problem since we just use the defining series for the polylogarithm in that region. It also becomes inefficient when $|z| \rightarrow \infty$, for which an asymptotic expansion would better.
5.1 fmpr.h – arbitrary-precision floating-point numbers

This type is now obsolete: use arf_t instead.

A variable of type fmpr_t holds an arbitrary-precision binary floating-point number, i.e. a rational number of the form $x \times 2^y$ where $x, y \in \mathbb{Z}$ and $x$ is odd; or one of the special values zero, plus infinity, minus infinity, or NaN (not-a-number).

The component $x$ is called the mantissa, and $y$ is called the exponent. Note that this is just one among many possible conventions: the mantissa (alternatively significand) is sometimes viewed as a fraction in the interval $[1/2, 1)$, with the exponent pointing to the position above the top bit rather than the position of the bottom bit, and with a separate sign.

The conventions for special values largely follow those of the IEEE floating-point standard. At the moment, there is no support for negative zero, unsigned infinity, or a NaN with a payload, though some these might be added in the future.

An fmpr number is exact and has no inherent “accuracy”. We use the term precision to denote either the target precision of an operation, or the bit size of a mantissa (which in general is unrelated to the “accuracy” of the number: for example, the floating-point value $1$ has a precision of $1$ bit in this sense and is simultaneously an infinitely accurate approximation of the integer $1$ and a 2-bit accurate approximation of $\sqrt{2} = 1.011010100\ldots_2$).

Except where otherwise noted, the output of an operation is the floating-point number obtained by taking the inputs as exact numbers, in principle carrying out the operation exactly, and rounding the resulting real number to the nearest representable floating-point number whose mantissa has at most the specified number of bits, in the specified direction of rounding. Some operations are always or optionally done exactly.

5.1.1 Types, macros and constants

fmpr_struct

An fmpr_struct holds a mantissa and an exponent. If the mantissa and exponent are sufficiently small, their values are stored as immediate values in the fmpr_struct; large values are represented by pointers to heap-allocated arbitrary-precision integers. Currently, both the mantissa and exponent are implemented using the FLINT fmpz type. Special values are currently encoded by the mantissa being set to zero.

fmpr_t

An fmpr_t is defined as an array of length one of type fmpr_struct, permitting an fmpr_t to be passed by reference.

fmpr_rnd_t

Specifies the rounding mode for the result of an approximate operation.
FMPR_RND_DOWN
   Specifies that the result of an operation should be rounded to the nearest representable number in the direction
towards zero.

FMPR_RND_UP
   Specifies that the result of an operation should be rounded to the nearest representable number in the direction
away from zero.

FMPR_RND_FLOOR
   Specifies that the result of an operation should be rounded to the nearest representable number in the direction
towards minus infinity.

FMPR_RND_CEIL
   Specifies that the result of an operation should be rounded to the nearest representable number in the direction
towards plus infinity.

FMPR_RND_NEAR
   Specifies that the result of an operation should be rounded to the nearest representable number, rounding to an
odd mantissa if there is a tie between two values. Warning: this rounding mode is currently not implemented
(except for a few conversions functions where this stated explicitly).

FMPR_PREC_EXACT
   If passed as the precision parameter to a function, indicates that no rounding is to be performed. This must only
be used when it is known that the result of the operation can be represented exactly and fits in memory (the
typical use case is working small integer values). Note that, for example, adding two numbers whose exponents
are far apart can easily produce an exact result that is far too large to store in memory.

5.1.2 Memory management

void fmpr_init (fmpr_t x)
   Initializes the variable x for use. Its value is set to zero.

void fmpr_clear (fmpr_t x)
   Clears the variable x, freeing or recycling its allocated memory.

5.1.3 Special values

void fmpr_zero (fmpr_t x)
void fmpr_one (fmpr_t x)
void fmpr_pos_inf (fmpr_t x)
void fmpr_neg_inf (fmpr_t x)
void fmpr_nan (fmpr_t x)
   Sets x respectively to 0, 1, +∞, −∞, NaN.

int fmpr_is_zero (const fmpr_t x)
int fmpr_is_one (const fmpr_t x)
int fmpr_is_pos_inf (const fmpr_t x)
int fmpr_is_neg_inf (const fmpr_t x)
int fmpr_is_nan (const fmpr_t x)
   Returns nonzero iff x respectively equals 0, 1, +∞, −∞, NaN.
int \texttt{fmpr\_is\_inf} (\texttt{const fmpr\_t x})

Returns nonzero iff \( x \) equals either \(+\infty\) or \(-\infty\).

int \texttt{fmpr\_is\_normal} (\texttt{const fmpr\_t x})

Returns nonzero iff \( x \) is a finite, nonzero floating-point value, i.e. not one of the special values \( 0, +\infty, -\infty, \text{NaN} \).

int \texttt{fmpr\_is\_special} (\texttt{const fmpr\_t x})

Returns nonzero iff \( x \) is one of the special values \( 0, +\infty, -\infty, \text{NaN} \), i.e. not a finite, nonzero floating-point value.

int \texttt{fmpr\_is\_finite} (\texttt{fmpr\_t x})

Returns nonzero iff \( x \) is a finite floating-point value, i.e. not one of the values \(+\infty, -\infty, \text{NaN}\). (Note that this is not equivalent to the negation of \texttt{fmpr\_is\_inf}().)

### 5.1.4 Assignment, rounding and conversions

\texttt{slong \_fmpr\_normalise} (\texttt{fmpz\_t man, fmpz\_t exp, slong prec, fmpr\_rnd\_t rnd})

Rounds the mantissa and exponent in-place.

\texttt{void \textbf{fmpr\_set}} (\texttt{fmpr\_t y, const fmpr\_t x})

Sets \( y \) to a copy of \( x \).

\texttt{void \textbf{fmpr\_swap}} (\texttt{fmpr\_t x, fmpr\_t y})

Swaps \( x \) and \( y \) efficiently.

\texttt{slong \textbf{fmpr\_set\_round}} (\texttt{fmpr\_t y, const fmpr\_t x, slong prec, fmpr\_rnd\_t rnd})

Sets \( y \) to a copy of \( x \) rounded in the direction specified by \texttt{rnd} to the number of bits specified by \texttt{prec}.

\texttt{slong \_fmpr\_set\_round\_mpn} (\texttt{slong * shift, fmpz\_t man, mp\_srcptr x, mp\_size\_t xn, int negative, slong prec, fmpr\_rnd\_t rnd})

Given an integer represented by a pointer \( x \) to a raw array of \( xn \) limbs (negated if \texttt{negative} is nonzero), sets \texttt{man} to the corresponding floating-point mantissa rounded to \texttt{prec} bits in direction \texttt{rnd}, sets \texttt{shift} to the exponent, and returns the error bound. We require that \( xn \) is positive and that the leading limb of \( x \) is nonzero.

\texttt{slong \textbf{fmpr\_set\_round\_ui\_2exp\_fmpz}} (\texttt{fmpr\_t z, mp\_limb\_t lo, const fmpz\_t exp, int negative, slong prec, fmpr\_rnd\_t rnd})

Sets \( z \) to the unsigned integer \( lo \) times two to the power \( exp \), negating the value if \texttt{negative} is nonzero, and rounding the result to \texttt{prec} bits in direction \texttt{rnd}.

\texttt{slong \textbf{fmpr\_set\_round\_uiui\_2exp\_fmpz}} (\texttt{fmpr\_t z, mp\_limb\_t hi, mp\_limb\_t lo, const fmpz\_t exp, int negative, slong prec, fmpr\_rnd\_t rnd})

Sets \( z \) to the unsigned two-limb integer \{\( hi, lo \)\} times two to the power \( exp \), negating the value if \texttt{negative} is nonzero, and rounding the result to \texttt{prec} bits in direction \texttt{rnd}.

\texttt{void \textbf{fmpr\_set\_error\_result}} (\texttt{fmpr\_t err, const fmpr\_t result, slong rret})

Given the return value \texttt{rret} and output variable \texttt{result} from a function performing a rounding (e.g. \texttt{fmpr\_set\_round} or \texttt{fmpr\_add}), sets \texttt{err} to a bound for the absolute error.

\texttt{void \textbf{fmpr\_add\_error\_result}} (\texttt{fmpr\_t err, const fmpr\_t err\_in, const fmpr\_t result, slong rret, slong prec, fmpr\_rnd\_t rnd})

Like \texttt{fmpr\_set\_error\_result}, but adds \texttt{err\_in} to the error.

\texttt{void \textbf{fmpr\_ulp}} (\texttt{fmpr\_t u, const fmpr\_t x, slong prec})

Sets \texttt{u} to the floating-point unit in the last place (ulp) of \( x \). The ulp is defined as in the MPFR documentation and satisfies \( 2^{-n}|x| \leq u \leq 2^{-n+1}|x| \) for any finite nonzero \( x \). If \( x \) is a special value, \texttt{u} is set to the absolute value of \( x \).
int fmpr_check_ulp (const fmpr_t x, slong r, slong prec)
    Assume that r is the return code and x is the floating-point result from a single floating-point rounding. Then this function returns nonzero iff x and r define an error of exactly 0 or 1 ulp. In other words, this function checks that fmpr_set_error_result() gives exactly 0 or 1 ulp as expected.

int fmpr_get_mpfr (mpfr_t x, const fmpr_t y, mpfr_rnd_t rnd)
    Sets the MPFR variable x to the value of y. If the precision of x is too small to allow y to be represented exactly, it is rounded in the specified MPFR rounding mode. The return value indicates the direction of rounding, following the standard convention of the MPFR library.

void fmpr_set_mpfr (fmpr_t x, const mpfr_t y)
    Sets x to the exact value of the MPFR variable y.

double fmpr_get_d (const fmpr_t x, mpfr_rnd_t rnd)
    Returns x rounded to a double in the direction specified by rnd.

void fmpr_set_d (fmpr_t x, double v)
    Sets x the the exact value of the argument v of type double.

void fmpr_set_ui (fmpr_t x, ulong c)

void fmpr_set_si (fmpr_t x, slong c)

void fmpr_set_fmpz (fmpr_t x, const fmpz_t c)
    Sets x exactly to the integer c.

void fmpr_get_fmpz (fmpz_t z, const fmpr_t x, fmpr_rnd_t rnd)
    Sets z to x rounded to the nearest integer in the direction specified by rnd. If rnd is FMPR_RND_NEAR, rounds to the nearest even integer in case of a tie. Aborts if x is infinite, NaN or if the exponent is unreasonably large.

slong fmpr_get_si (const fmpr_t x, fmpr_rnd_t rnd)
    Returns x rounded to the nearest integer in the direction specified by rnd. If rnd is FMPR_RND_NEAR, rounds to the nearest even integer in case of a tie. Aborts if x is infinite, NaN, or the value is too large to fit in an slong.

void fmpr_get_fmpq (fmpq_t y, const fmpr_t x)
    Sets y to the exact value of x. The result is undefined if x is not a finite fraction.

slong fmpr_set_fmpq (fmpr_t x, const fmpq_t y, long prec, fmpr_rnd_t rnd)
    Sets x to the value of y, rounded according to prec and rnd.

void fmpr_set_fmpz_2exp (fmpr_t x, const fmpz_t man, const fmpz_t exp)

void fmpr_set_si_2exp_si (fmpr_t x, slong man, slong exp)

void fmpr_set_ui_2exp_si (fmpr_t x, ulong man, slong exp)
    Sets x to man \times 2^\text{exp}.

slong fmpr_set_round_fmpz_2exp (fmpr_t x, const fmpz_t man, const fmpz_t exp, slong prec, fmpr_rnd_t rnd)
    Sets x to man \times 2^\text{exp}, rounded according to prec and rnd.

void fmpr_get_fmpz_2exp (fmpr_t x, const fmpz_t man, const fmpz_t exp)
    Sets man and exp to the unique integers such that x = man \times 2^\text{exp} and man is odd, provided that x is a nonzero finite fraction. If x is zero, both man and exp are set to zero. If x is infinite or NaN, the result is undefined.

int fmpr_get_fmpz_fixed_fmpz (fmpz_t y, const fmpr_t x, const fmpz_t e)

int fmpr_get_fmpz_fixed_si (fmpr_t y, const fmpr_t x, slong e)
    Converts x to a mantissa with predetermined exponent, i.e. computes an integer y such that y \times 2^e \approx x, truncating if necessary. Returns 0 if exact and 1 if truncation occurred.
5.1.5 Comparisons

int \texttt{fmpr\_equal} (const \texttt{fmpr\_t} x, const \texttt{fmpr\_t} y)
Returns nonzero iff \( x \) and \( y \) are exactly equal. This function does not treat NaN specially, i.e. NaN compares as equal to itself.

int \texttt{fmpr\_cmp} (const \texttt{fmpr\_t} x, const \texttt{fmpr\_t} y)
Returns negative, zero, or positive, depending on whether \( x \) is respectively smaller, equal, or greater compared to \( y \). Comparison with NaN is undefined.

int \texttt{fmpr\_cmpabs} (const \texttt{fmpr\_t} x, const \texttt{fmpr\_t} y)
Compares the absolute values of \( x \) and \( y \).

int \texttt{fmpr\_cmpabs\_ui} (const \texttt{fmpr\_t} x, ulong y)
Compares \( x \) (respectively its absolute value) with \( 2^e \).

int \texttt{fmpr\_sgn} (const \texttt{fmpr\_t} x)
Returns \(-1, 0 \) or \(+1 \) according to the sign of \( x \). The sign of NaN is undefined.

void \texttt{fmpr\_min} (\texttt{fmpr\_t} z, const \texttt{fmpr\_t} a, const \texttt{fmpr\_t} b)
Sets \( z \) respectively to the minimum and the maximum of \( a \) and \( b \).

slong \texttt{fmpr\_bits} (const \texttt{fmpr\_t} x)
Returns the number of bits needed to represent the absolute value of the mantissa of \( x \), i.e. the minimum precision sufficient to represent \( x \) exactly. Returns 0 if \( x \) is a special value.

int \texttt{fmpr\_is\_int} (const \texttt{fmpr\_t} x)
Returns nonzero iff \( x \) is integer-valued.

int \texttt{fmpr\_is\_int\_2exp\_si} (const \texttt{fmpr\_t} x, slong e)
Returns nonzero iff \( x \) equals \( n2^e \) for some integer \( n \).

void \texttt{fmpr\_abs\_bound\_le\_2exp\_fmpz} (fmpz_t b, const \texttt{fmpr\_t} x)
Sets \( b \) to the smallest integer such that \(|x| \leq 2^b \). If \( x \) is zero, infinity or NaN, the result is undefined.

void \texttt{fmpr\_abs\_bound\_lt\_2exp\_fmpz} (fmpz_t b, const \texttt{fmpr\_t} x)
Sets \( b \) to the smallest integer such that \(|x| < 2^b \). If \( x \) is zero, infinity or NaN, the result is undefined.

slong \texttt{fmpr\_abs\_bound\_lt\_2exp\_si} (const \texttt{fmpr\_t} x)
Returns the smallest integer \( b \) such that \(|x| < 2^b \), clamping the result to lie between \textit{FMPR\_PREC\_EXACT} and \textit{FMPR\_PREC\_EXACT} inclusive. If \( x \) is zero, \textit{-FMPR\_PREC\_EXACT} is returned, and if \( x \) is infinity or NaN, \textit{FMPR\_PREC\_EXACT} is returned.

5.1.6 Random number generation

void \texttt{fmpr\_randtest} (\texttt{fmpr\_t} x, flint\_rand\_t state, slong bits, slong mag\_bits)
Generates a finite random number whose mantissa has precision at most \( \text{bits} \) and whose exponent has at most \( \text{mag\_bits} \) bits. The values are distributed non-uniformly: special bit patterns are generated with high probability in order to allow the test code to exercise corner cases.

void \texttt{fmpr\_randtest\_not\_zero} (\texttt{fmpr\_t} x, flint\_rand\_t state, slong bits, slong mag\_bits)
Identical to \texttt{fmpr\_randtest}, except that zero is never produced as an output.

void \texttt{fmpr\_randtest\_special} (\texttt{fmpr\_t} x, flint\_rand\_t state, slong bits, slong mag\_bits)
Identical to \texttt{fmpr\_randtest}, except that the output occasionally is set to an infinity or NaN.
5.1.7 Input and output

void **fmpr_print** (const fmpr_t x)
Prints the mantissa and exponent of x as integers, precisely showing the internal representation.

void **fmpr_printd** (const fmpr_t x, slong digits)
Prints x as a decimal floating-point number, rounding to the specified number of digits. This function is currently implemented using MPFR, and does not support large exponents.

5.1.8 Arithmetic

void **fmpr_neg** (fmpr_t y, const fmpr_t x)
Sets y to the negation of x.

slong **fmpr_neg_round** (fmpr_t y, const fmpr_t x, slong prec, fmpr_rnd_t rnd)
Sets y to the negation of x, rounding the result.

void **fmpr_abs** (fmpr_t y, const fmpr_t x)
Sets y to the absolute value of x.

slong **fmpr_add** (fmpr_t z, const fmpr_t x, const fmpr_t y, slong prec, fmpr_rnd_t rnd)
Sets z = x + y, rounded according to prec and rnd. The precision can be FMPR_PREC_EXACT to perform an exact addition, provided that the result fits in memory.

slong **fmpr_add_eps** (fmpr_t z, const fmpr_t x, int sign, slong prec, fmpr_rnd_t rnd)
Sets z to the value that results by adding an infinitesimal quantity of the given sign to x, and rounding. The result is undefined if x is zero.

slong **fmpr_sub** (fmpr_t z, const fmpr_t x, const fmpr_t y, slong prec, fmpr_rnd_t rnd)
Sets z = x − y, rounded according to prec and rnd. The precision can be FMPR_PREC_EXACT to perform an exact addition, provided that the result fits in memory.

slong **fmpr_sum** (fmpr_t s, const fmpr_struct * terms, slong len, slong prec, fmpr_rnd_t rnd)
Sets s to the sum of the array terms of length len, rounded to prec bits in the direction rnd. The sum is computed as if done without any intermediate rounding error, with only a single rounding applied to the final result. Unlike repeated calls to fmpr_add, this function does not overflow if the magnitudes of the terms are far apart. Warning: this function is implemented naively, and the running time is quadratic with respect to len in the worst case.

slong **fmpr_mul** (fmpr_t z, const fmpr_t x, const fmpr_t y, slong prec, fmpr_rnd_t rnd)
slong **fmpr_mul_ui** (fmpr_t z, const fmpr_t x, ulong y, slong prec, fmpr_rnd_t rnd)
slong **fmpr_mul_si** (fmpr_t z, const fmpr_t x, slong y, slong prec, fmpr_rnd_t rnd)
slong **fmpr_mul_fmpz** (fmpr_t z, const fmpr_t x, const fmpz_t y, slong prec, fmpr_rnd_t rnd)
Sets z = x × y, rounded according to prec and rnd. The precision can be FMPR_PREC_EXACT to perform an exact multiplication, provided that the result fits in memory.

void **fmpr_mul_2exp_si** (fmpr_t y, const fmpr_t x, slong e)
void \texttt{fmpr\_mul\_2exp\_fmpz} (\texttt{fmpr\_t y}, \texttt{const fmpr\_t x}, \texttt{const fmpz\_t e})
\begin{itemize}
\item Sets \(y\) to \(x\) multiplied by \(2^e\) without rounding.
\end{itemize}

\texttt{slong fmpr\_div} (\texttt{fmpr\_t z}, \texttt{const fmpr\_t x}, \texttt{const fmpr\_t y}, \texttt{slong prec, fmpr\_rnd\_t rnd})

\texttt{slong fmpr\_div\_ui} (\texttt{fmpr\_t z}, \texttt{const fmpr\_t x}, \texttt{ulong y}, \texttt{slong prec, fmpr\_rnd\_t rnd})

\texttt{slong fmpr\_ui\_div} (\texttt{fmpr\_t z}, \texttt{ulong x}, \texttt{const fmpr\_t y}, \texttt{slong prec, fmpr\_rnd\_t rnd})

\texttt{slong fmpr\_div\_si} (\texttt{fmpr\_t z}, \texttt{const fmpr\_t x}, \texttt{slong y}, \texttt{slong prec, fmpr\_rnd\_t rnd})

\texttt{slong fmpr\_si\_div} (\texttt{fmpr\_t z}, \texttt{slong x}, \texttt{const fmpr\_t y}, \texttt{slong prec, fmpr\_rnd\_t rnd})

\texttt{slong fmpr\_div\_fmpz} (\texttt{fmpr\_t z}, \texttt{const fmpr\_t x}, \texttt{const fmpr\_t y}, \texttt{slong prec, fmpr\_rnd\_t rnd})

\texttt{slong fmpr\_fmpz\_div\_fmpz} (\texttt{fmpr\_t z}, \texttt{const fmpr\_t x}, \texttt{const fmpr\_t y}, \texttt{slong prec, fmpr\_rnd\_t rnd})
\begin{itemize}
\item Sets \(z = x/y\), rounded according to \texttt{prec} and \texttt{rnd}. If \(y\) is zero, \(z\) is set to NaN.
\end{itemize}

\texttt{void fmpr\_divapppr\_abs\_ubound} (\texttt{fmpr\_t z}, \texttt{const fmpr\_t x}, \texttt{const fmpr\_t y}, \texttt{slong prec})
\begin{itemize}
\item Sets \(z\) to an upper bound for \(|x|/|y|\), computed to a precision of approximately \texttt{prec} bits. The error can be a few ulp.
\end{itemize}

\texttt{slong fmpr\_addmul} (\texttt{fmpr\_t z}, \texttt{const fmpr\_t x}, \texttt{const fmpr\_t y}, \texttt{slong prec, fmpr\_rnd\_t rnd})

\texttt{slong fmpr\_addmul\_ui} (\texttt{fmpr\_t z}, \texttt{const fmpr\_t x}, \texttt{ulong y}, \texttt{slong prec, fmpr\_rnd\_t rnd})

\texttt{slong fmpr\_addmul\_si} (\texttt{fmpr\_t z}, \texttt{const fmpr\_t x}, \texttt{slong y}, \texttt{slong prec, fmpr\_rnd\_t rnd})

\texttt{slong fmpr\_addmul\_fmpz} (\texttt{fmpr\_t z}, \texttt{const fmpr\_t x}, \texttt{const fmpr\_t y}, \texttt{slong prec, fmpr\_rnd\_t rnd})
\begin{itemize}
\item Sets \(z = z + x \times y\), rounded according to \texttt{prec} and \texttt{rnd}. The intermediate multiplication is always performed without roundoff. The precision can be \texttt{FMPR\_PREC\_EXACT} to perform an exact addition, provided that the result fits in memory.
\end{itemize}

\texttt{slong fmpr\_submul} (\texttt{fmpr\_t z}, \texttt{const fmpr\_t x}, \texttt{const fmpr\_t y}, \texttt{slong prec, fmpr\_rnd\_t rnd})

\texttt{slong fmpr\_submul\_ui} (\texttt{fmpr\_t z}, \texttt{const fmpr\_t x}, \texttt{ulong y}, \texttt{slong prec, fmpr\_rnd\_t rnd})

\texttt{slong fmpr\_submul\_si} (\texttt{fmpr\_t z}, \texttt{const fmpr\_t x}, \texttt{slong y}, \texttt{slong prec, fmpr\_rnd\_t rnd})

\texttt{slong fmpr\_submul\_fmpz} (\texttt{fmpr\_t z}, \texttt{const fmpr\_t x}, \texttt{const fmpr\_t y}, \texttt{slong prec, fmpr\_rnd\_t rnd})
\begin{itemize}
\item Sets \(z = z - x \times y\), rounded according to \texttt{prec} and \texttt{rnd}. The intermediate multiplication is always performed without roundoff. The precision can be \texttt{FMPR\_PREC\_EXACT} to perform an exact subtraction, provided that the result fits in memory.
\end{itemize}

\texttt{slong fmpr\_sqrt} (\texttt{fmpr\_t y}, \texttt{const fmpr\_t x}, \texttt{slong prec, fmpr\_rnd\_t rnd})

\texttt{slong fmpr\_sqrt\_fmpz} (\texttt{fmpr\_t z}, \texttt{const fmpr\_t x}, \texttt{slong prec, fmpr\_rnd\_t rnd})
\begin{itemize}
\item Sets \(z\) to the square root of \(x\), rounded according to \texttt{prec} and \texttt{rnd}. The result is NaN if \(x\) is negative.
\end{itemize}

\texttt{slong fmpr\_rsqrt} (\texttt{fmpr\_t z}, \texttt{const fmpr\_t x}, \texttt{slong prec, fmpr\_rnd\_t rnd})
\begin{itemize}
\item Sets \(z\) to the reciprocal square root of \(x\), rounded according to \texttt{prec} and \texttt{rnd}. The result is NaN if \(x\) is negative. At high precision, this is faster than computing a square root.
\end{itemize}

\texttt{slong fmpr\_root} (\texttt{fmpr\_t z}, \texttt{const fmpr\_t x}, \texttt{ulong k}, \texttt{slong prec, fmpr\_rnd\_t rnd})
\begin{itemize}
\item Sets \(z\) to the \(k\)-th root of \(x\), rounded to \texttt{prec} bits in the direction \texttt{rnd}. Warning: this function wraps MPFR, and is currently only fast for small \(k\).
\end{itemize}

\texttt{void fmpr\_pow\_sloppy\_fmpz} (\texttt{fmpr\_t y}, \texttt{const fmpr\_t b}, \texttt{const fmpz\_t e}, \texttt{slong prec, fmpr\_rnd\_t rnd})

\texttt{void fmpr\_pow\_sloppy\_ui} (\texttt{fmpr\_t y}, \texttt{const fmpr\_t b}, \texttt{ulong e}, \texttt{slong prec, fmpr\_rnd\_t rnd})
void \texttt{fmp\_pow\_sloppy\_si}(\texttt{fmp\_t} \ y, \texttt{const fmp\_t} \ b, \texttt{ulong} \ e, \texttt{ulong} \ \texttt{prec}, \texttt{fmp\_rnd\_t} \ \texttt{rnd})

Sets \( y = b^e \), computed using without guaranteeing correct (optimal) rounding, but guaranteeing that the result is a correct upper or lower bound if the rounding is directional. Currently requires \( b \geq 0 \).

### 5.1.9 Special functions

**fmp\_log** (\texttt{fmp\_t} \ y, \texttt{const fmp\_t} \ x, \texttt{ulong} \ \texttt{prec}, \texttt{fmp\_rnd\_t} \ \texttt{rnd})

Sets \( y = \log(x) \), rounded according to \texttt{prec} and \texttt{rnd}. The result is NaN if \( x \) is negative. This function is currently implemented using MPFR and does not support large exponents.

**fmp\_log1p** (\texttt{fmp\_t} \ y, \texttt{const fmp\_t} \ x, \texttt{ulong} \ \texttt{prec}, \texttt{fmp\_rnd\_t} \ \texttt{rnd})

Sets \( y = \log(1 + x) \), rounded according to \texttt{prec} and \texttt{rnd}. This function computes an accurate value when \( x \) is small. The result is NaN if \( 1 + x \) is negative. This function is currently implemented using MPFR and does not support large exponents.

**fmp\_exp** (\texttt{fmp\_t} \ y, \texttt{const fmp\_t} \ x, \texttt{ulong} \ \texttt{prec}, \texttt{fmp\_rnd\_t} \ \texttt{rnd})

Sets \( y = \exp(x) \), rounded according to \texttt{prec} and \texttt{rnd}. This function is currently implemented using MPFR and does not support large exponents.

**fmp\_expm1** (\texttt{fmp\_t} \ y, \texttt{const fmp\_t} \ x, \texttt{ulong} \ \texttt{prec}, \texttt{fmp\_rnd\_t} \ \texttt{rnd})

Sets \( y = \exp(x) - 1 \), rounded according to \texttt{prec} and \texttt{rnd}. This function computes an accurate value when \( x \) is small. This function is currently implemented using MPFR and does not support large exponents.

### 5.2 fmp\_rb.h – real numbers represented as floating-point balls

This type is now obsolete: use \texttt{arb\_t} instead.

An \texttt{fmp\_rb\_t} represents a ball over the real numbers.

#### 5.2.1 Types, macros and constants

**fmp\_rb\_struct**

**fmp\_rb\_t**

An \texttt{fmp\_rb\_struct} consists of a pair of \texttt{fmp\_struct}s. An \texttt{fmp\_rb\_t} is defined as an array of length one of type \texttt{fmp\_rb\_struct}, permitting an \texttt{fmp\_rb\_t} to be passed by reference.

**fmp\_rb\_ptr**

Alias for \texttt{fmp\_rb\_struct} *, used for vectors of numbers.

**fmp\_rb\_srcptr**

Alias for \texttt{const fmp\_rb\_struct} *, used for vectors of numbers when passed as constant input to functions.

**FMPRB\_RAD\_PREC**

The precision used for operations on the radius. This is small enough to fit in a single word, currently 30 bits.

**fmp\_rb\_midref** (\texttt{x})

Macro returning a pointer to the midpoint of \texttt{x} as an \texttt{fmp\_t}.

**fmp\_rb\_radref** (\texttt{x})

Macro returning a pointer to the radius of \texttt{x} as an \texttt{fmp\_t}.
5.2.2 Memory management

void \texttt{fmprb\_init} (\texttt{fmprb\_t x})
\hspace{1cm}Initializes the variable \texttt{x} for use. Its midpoint and radius are both set to zero.

void \texttt{fmprb\_clear} (\texttt{fmprb\_t x})
\hspace{1cm}Clears the variable \texttt{x}, freeing or recycling its allocated memory.

\texttt{fmprb\_ptr \_fmprb\_vec\_init} (slong \texttt{n})
\hspace{1cm}Returns a pointer to an array of \texttt{n} initialized \texttt{fmprb\_struct}:s.

void \texttt{\_fmprb\_vec\_clear} (\texttt{fmprb\_ptr v}, slong \texttt{n})
\hspace{1cm}Clears an array of \texttt{n} initialized \texttt{fmprb\_struct}:s.

5.2.3 Assignment and rounding

void \texttt{fmprb\_set} (\texttt{fmprb\_t y}, const \texttt{fmprb\_t x})
\hspace{1cm}Sets \texttt{y} to a copy of \texttt{x}.

void \texttt{fmprb\_set\_round} (\texttt{fmprb\_t y}, const \texttt{fmprb\_t x}, slong \texttt{prec})
\hspace{1cm}Sets \texttt{y} to a copy of \texttt{x}, rounded to \texttt{prec} bits.

void \texttt{fmprb\_set\_fmp} (\texttt{fmprb\_t y}, const \texttt{fmprb\_t x})

void \texttt{fmprb\_set\_si} (\texttt{fmprb\_t y}, slong \texttt{x})

void \texttt{fmprb\_set\_ui} (\texttt{fmprb\_t y}, ulong \texttt{x})

void \texttt{fmprb\_set\_fmpz} (\texttt{fmprb\_t y}, const \texttt{fmpz\_t x})
\hspace{1cm}Sets \texttt{y} exactly to \texttt{x}.

void \texttt{fmprb\_set\_fmpq} (\texttt{fmprb\_t y}, const \texttt{fmpq\_t x}, slong \texttt{prec})
\hspace{1cm}Sets \texttt{y} to the rational number \texttt{x}, rounded to \texttt{prec} bits.

void \texttt{fmprb\_set\_fmpz\_2exp} (\texttt{fmprb\_t x}, const \texttt{fmpz\_t y}, const \texttt{fmpz\_t exp})
\hspace{1cm}Sets \texttt{x} to \texttt{y} multiplied by 2 raised to the power \texttt{exp}.

void \texttt{fmprb\_set\_round\_fmpz\_2exp} (\texttt{fmprb\_t y}, const \texttt{fmpz\_t x}, const \texttt{fmpz\_t exp}, slong \texttt{prec})
\hspace{1cm}Sets \texttt{x} to \texttt{y} multiplied by 2 raised to the power \texttt{exp}, rounding the result to \texttt{prec} bits.

5.2.4 Assignment of special values

void \texttt{fmprb\_zero} (\texttt{fmprb\_t x})
\hspace{1cm}Sets \texttt{x} to zero.

void \texttt{fmprb\_one} (\texttt{fmprb\_t x})
\hspace{1cm}Sets \texttt{x} to the exact integer 1.

void \texttt{fmprb\_pos\_inf} (\texttt{fmprb\_t x})
\hspace{1cm}Sets \texttt{x} to positive infinity, with a zero radius.

void \texttt{fmprb\_neg\_inf} (\texttt{fmprb\_t x})
\hspace{1cm}Sets \texttt{x} to negative infinity, with a zero radius.

void \texttt{fmprb\_zero\_pm\_inf} (\texttt{fmprb\_t x})
\hspace{1cm}Sets \texttt{x} to $[0 \pm \infty]$, representing the whole extended real line.

void \texttt{fmprb\_indeterminate} (\texttt{fmprb\_t x})
\hspace{1cm}Sets \texttt{x} to $[\text{NaN} \pm \infty]$, representing an indeterminate result.
5.2.5 Input and output

void `fmprb_print` (const `fmprb_t x`)
   Prints the internal representation of `x`.

void `fmprb_printd` (const `fmprb_t x`, slong `digits`)
   Prints `x` in decimal. The printed value of the radius is not adjusted to compensate for the fact that
   the binary-to-decimal conversion of both the midpoint and the radius introduces additional error.

5.2.6 Random number generation

void `fmprb_randtest` (`fmprb_t x`, flint_rand_t `state`, slong `prec`, slong `mag_bits`)
   Generates a random ball. The midpoint and radius will both be finite.

void `fmprb_randtest_exact` (`fmprb_t x`, flint_rand_t `state`, slong `prec`, slong `mag_bits`)
   Generates a random number with zero radius.

void `fmprb_randtest_precise` (`fmprb_t x`, flint_rand_t `state`, slong `prec`, slong `mag_bits`)
   Generates a random number with radius at most $2^{-\text{prec}}$ the magnitude of the midpoint.

void `fmprb_randtest_wide` (`fmprb_t x`, flint_rand_t `state`, slong `prec`, slong `mag_bits`)
   Generates a random number with midpoint and radius chosen independently, possibly giving a very large
   interval.

void `fmprb_randtest_special` (`fmprb_t x`, flint_rand_t `state`, slong `prec`, slong `mag_bits`)
   Generates a random interval, possibly having NaN or an infinity as the midpoint and possibly having an
   infinite radius.

void `fmprb_get_rand_fmpq` (``fmpq_t`` `q`, flint_rand_t `state`, const `fmprb_t x`, slong `bits`)
   Sets `q` to a random rational number from the interval represented by `x`. A denominator is chosen by
   multiplying the binary denominator of `x` by a random integer up to `bits` bits.
   
   The outcome is undefined if the midpoint or radius of `x` is non-finite, or if the exponent of the midpoint
   or radius is so large or small that representing the endpoints as exact rational numbers would cause
   overflows.

5.2.7 Radius and interval operations

void `fmprb_add_error_fmpr` (``fmprb_t`` `x`, const `fmpr_t err`)
   Adds `err`, which is assumed to be nonnegative, to the radius of `x`.

void `fmprb_add_error_2exp_si` (``fmprb_t`` `x`, slong `e`)
   Adds $2^e$ to the radius of `x`.

void `fmprb_add_error_2exp_fmpz` (``fmprb_t`` `x`, const `fmpz_t` `e`)
   Adds `2^e` to the radius of `x`.

void `fmprb_add_error` (``fmprb_t`` `x`, const `fmprb_t err`)
   Adds the supremum of `err`, which is assumed to be nonnegative, to the radius of `x`.

void `fmprb_union` (``fmprb_t`` `z`, const `fmprb_t x`, const `fmprb_t y`, slong `prec`)
   Sets `z` to a ball containing both `x` and `y`.

void `fmprb_get_abs_ubound_fmpr` (``fmpr_t`` `u`, const `fmprb_t x`, slong `prec`)
   Sets `u` to the upper bound of the absolute value of `x`, rounded up to `prec` bits. If `x` contains NaN, the
   result is NaN.

void `fmprb_get_abs_lbound_fmpr` (``fmpr_t`` `u`, const `fmprb_t x`, slong `prec`)
   Sets `u` to the lower bound of the absolute value of `x`, rounded down to `prec` bits. If `x` contains NaN, the
   result is NaN.
void \textbf{fmprb\_get\_interval\_fmpz\_2exp} (fmpz\_t \textit{a}, fmpz\_t \textit{b}, fmpz\_t \textit{exp}, const fmpr\_t \textit{x})

Computes the exact interval represented by \textit{x}, in the form of an integer interval multiplied by a power of two, i.e. \( x = [a, b] \times 2^{\textit{exp}} \).

The outcome is undefined if the midpoint or radius of \textit{x} is non-finite, or if the difference in magnitude between the midpoint and radius is so large that representing the endpoints exactly would cause overflows.

void \textbf{fmprb\_set\_interval\_fmpr} (fmpr\_t \textit{x}, const fmpr\_t \textit{a}, const fmpr\_t \textit{b}, slong \textit{prec})

Sets \textit{x} to a ball containing the interval \([a, b]\). We require that \( a \leq b \).

slong \textbf{fmprb\_rel\_error\_bits} (const fmpr\_t \textit{x})

Returns the effective relative error of \textit{x} measured in bits, defined as the difference between the position of the top bit in the radius and the top bit in the midpoint, plus one. The result is clamped between plus/minus \texttt{FMPR\_PREC\_EXACT}.

slong \textbf{fmprb\_rel\_accuracy\_bits} (const fmpr\_t \textit{x})

Returns the effective relative accuracy of \textit{x} measured in bits, equal to the negative of the return value from \textbf{fmprb\_rel\_error\_bits}.

slong \textbf{fmprb\_bits} (const fmpr\_t \textit{x})

Returns the number of bits needed to represent the absolute value of the mantissa of the midpoint of \textit{x}, i.e. the minimum precision sufficient to represent \textit{x} exactly. Returns 0 if the midpoint of \textit{x} is a special value.

void \textbf{fmprb\_trim} (fmpr\_t \textit{y}, const fmpr\_t \textit{x})

Sets \textit{y} to a trimmed copy of \textit{x}: rounds \textit{x} to a number of bits equal to the accuracy of \textit{x} (as indicated by its radius), plus a few guard bits. The resulting ball is guaranteed to contain \textit{x}, but is more economical if \textit{x} has less than full accuracy.

int \textbf{fmprb\_get\_unique\_fmpz} (fmpz\_t \textit{z}, const fmpr\_t \textit{x})

If \textit{x} contains a unique integer, sets \textit{z} to that value and returns nonzero. Otherwise (if \textit{x} represents no integers or more than one integer), returns zero.

5.2.8 Comparisons

int \textbf{fmprb\_is\_zero} (const fmpr\_t \textit{x})

Returns nonzero iff the midpoint and radius of \textit{x} are both zero.

int \textbf{fmprb\_is\_nonzero} (const fmpr\_t \textit{x})

Returns nonzero iff zero is not contained in the interval represented by \textit{x}.

int \textbf{fmprb\_is\_one} (const fmpr\_t \textit{x})

Returns nonzero iff \textit{x} is exactly 1.

int \textbf{fmprb\_is\_finite} (fmpr\_t \textit{x})

Returns nonzero iff the midpoint and radius of \textit{x} are both finite floating-point numbers, i.e. not infinities or NaN.

int \textbf{fmprb\_is\_exact} (const fmpr\_t \textit{x})

Returns nonzero iff the radius of \textit{x} is zero.

int \textbf{fmprb\_is\_int} (const fmpr\_t \textit{x})

Returns nonzero iff \textit{x} is an exact integer.

int \textbf{fmprb\_equal} (const fmpr\_t \textit{x}, const fmpr\_t \textit{y})

Returns nonzero iff \textit{x} and \textit{y} are equal as balls, i.e. have both the same midpoint and radius.

Note that this is not the same thing as testing whether both \textit{x} and \textit{y} certainly represent the same real number, unless either \textit{x} or \textit{y} is exact (and neither contains NaN). To test whether both operands \texttt{might} represent the same mathematical quantity, use \textbf{fmprb\_overlaps()} or \textbf{fmprb\_contains()}, depending on the circumstance.

int \textbf{fmprb\_is\_positive} (const fmpr\_t \textit{x})

5.2. fmprb\_h – real numbers represented as floating-point balls
int fmprb_is_nonnegative (const fmprb_t x)
int fmprb_is_negative (const fmprb_t x)
int fmprb_is_nonpositive (const fmprb_t x)
returns nonzero iff all points $p$ in the interval represented by $x$ satisfy, respectively, $p > 0$, $p \geq 0$, $p < 0$, $p \leq 0$. If $x$ contains NaN, returns zero.

int fmprb_overlaps (const fmprb_t x, const fmprb_t y)
returns nonzero iff $x$ and $y$ have some point in common. If either $x$ or $y$ contains NaN, this function always returns nonzero (as a NaN could be anything, it could in particular contain any number that is included in the other operand).

int fmprb_contains_fmpr (const fmprb_t x, const fmpfr_t y)
int fmprb_contains_fmpq (const fmprb_t x, const fmpq_t y)
int fmprb_contains_fmpz (const fmprb_t x, const fmpz_t y)
int fmprb_contains_si (const fmprb_t x, slong y)
int fmprb_contains_mpfr (const fmprb_t x, const mpfr_t y)
int fmprb_contains_zero (const fmprb_t x)
returns nonzero iff the given number (or ball) $y$ is contained in the interval represented by $x$.
If $x$ contains NaN, this function always returns nonzero (as it could represent anything, and in particular could represent all the points included in $y$). If $y$ contains NaN and $x$ does not, it always returns zero.

int fmprb_contains_negative (const fmprb_t x)
int fmprb_contains_nonpositive (const fmprb_t x)
int fmprb_contains_positive (const fmprb_t x)
int fmprb_contains_nonnegative (const fmprb_t x)
returns nonzero iff there is any point $p$ in the interval represented by $x$ satisfying, respectively, $p < 0$, $p \leq 0$, $p > 0$, $p \geq 0$. If $x$ contains NaN, returns nonzero.

5.2.9 Arithmetic

void fmprb_neg (fmprb_t y, const fmprb_t x)
sets $y$ to the negation of $x$.

void fmprb_abs (fmprb_t y, const fmprb_t x)
sets $y$ to the absolute value of $x$. No attempt is made to improve the interval represented by $x$ if it contains zero.

void fmprb_add (fmprb_t z, const fmprb_t x, const fmprb_t y, slong prec)
void fmprb_add_ui (fmprb_t z, const fmprb_t x, ulong y, slong prec)
void fmprb_add_si (fmprb_t z, const fmprb_t x, slong y, slong prec)
void fmprb_add_fmpz (fmprb_t z, const fmprb_t x, const fmpz_t y, slong prec)
void fmprb_add_fmpfr (fmprb_t z, const fmprb_t x, const mpfr_t y, slong prec)
sets $z = x + y$, rounded to $prec$ bits. The precision can be FMPR_PREC_EXACT provided that the result fits in memory.

void fmprb_sub (fmprb_t z, const fmprb_t x, const fmprb_t y, slong prec)
void fmprb_sub_ui (fmprb_t z, const fmprb_t x, const fmprb_t y, slong prec)
void `fmprb_sub_si` (const `fmprb_t` z, `slong` y, `slong` prec)

void `fmprb_sub_fmpz` (const `fmprb_t` z, const `fmpz_t` x, `const fmpz_t` y, `slong` prec)

Sets \( z = x - y \), rounded to `prec` bits. The precision can be `FMPR_PREC_EXACT` provided that the result fits in memory.

void `fmprb_mul` (const `fmprb_t` z, const `fmprb_t` x, const `fmprb_t` y, `slong` prec)

void `fmprb_mul_ui` (const `fmprb_t` z, const `fmprb_t` x, `ulong` y, `slong` prec)

void `fmprb_mul_si` (const `fmprb_t` z, const `fmprb_t` x, `slong` y, `slong` prec)

void `fmprb_mul_fmpz` (const `fmprb_t` z, const `fmprb_t` x, const `fmpz_t` y, `slong` prec)

void `fmprb_mul_2exp_fmpz` (const `fmprb_t` z, `ulong` e, const `fmpz_t` y, `slong` prec)

void `fmprb_mul_2exp_si` (const `fmprb_t` z, `ulong` e)

void `fmprb_addmul` (const `fmprb_t` z, const `fmprb_t` x, const `fmprb_t` y, `slong` prec)

void `fmprb_addmul_ui` (const `fmprb_t` z, const `fmprb_t` x, `ulong` y, `slong` prec)

void `fmprb_addmul_si` (const `fmprb_t` z, const `fmprb_t` x, `slong` y, `slong` prec)

void `fmprb_submul` (const `fmprb_t` z, const `fmprb_t` x, const `fmprb_t` y, `slong` prec)

void `fmprb_submul_ui` (const `fmprb_t` z, const `fmprb_t` x, `ulong` y, `slong` prec)

void `fmprb_submul_si` (const `fmprb_t` z, const `fmprb_t` x, `slong` y, `slong` prec)

void `fmprb_submul_fmpz` (const `fmprb_t` z, const `fmprb_t` x, const `fmpz_t` y, `slong` prec)

Sets \( z = x - y \), rounded to `prec` bits. The precision can be `FMPR_PREC_EXACT` provided that the result fits in memory.

---

5.2. `fmprb.h` – real numbers represented as floating-point balls

\[
\frac{x - x + \xi_1a}{y + \xi_2b} = \frac{x\xi_2b - y\xi_1a}{y(y + \xi_2b)} \leq \frac{|xb| + |ya|}{|y||y - b|}
\]

where \(-1 \leq \xi_1, \xi_2 \leq 1\), and where the triangle inequality has been applied to the numerator and the reverse triangle inequality has been applied to the denominator.
5.2.10 Powers and roots

void \texttt{fmprb\_sqrt} (\texttt{fmprb\_t z}, \texttt{const fmprb\_t x}, \texttt{slong prec})

void \texttt{fmprb\_sqrt\_ui} (\texttt{fmprb\_t z}, \texttt{ulong x}, \texttt{slong prec})

void \texttt{fmprb\_sqrt\_fmpz} (\texttt{fmprb\_t z}, \texttt{const fmpz\_t x}, \texttt{slong prec})

Sets \texttt{z} to the square root of \texttt{x}, rounded to \texttt{prec} bits. Error propagation is done using the following rule: assuming \( m > r \geq 0 \), the error is largest at \( m - r \), and we have \( \sqrt{m} - \sqrt{m - r} \leq r/(2\sqrt{m - r}) \).

void \texttt{fmprb\_sqrtpos} (\texttt{fmprb\_t z}, \texttt{const fmprb\_t x}, \texttt{slong prec})

Sets \texttt{z} to the square root of \texttt{x}, assuming that \texttt{x} represents a nonnegative number (i.e. discarding any negative numbers in the input interval), and producing an output interval not containing any negative numbers (unless the radius is infinite).

void \texttt{fmprb\_hypot} (\texttt{fmprb\_t z}, \texttt{const fmprb\_t x}, \texttt{const fmprb\_t y}, \texttt{slong prec})

Sets \texttt{z} to \( \sqrt{x^2 + y^2} \).

void \texttt{fmprb\_rsqrt} (\texttt{fmprb\_t z}, \texttt{const fmprb\_t x}, \texttt{slong prec})

void \texttt{fmprb\_rsqrt\_ui} (\texttt{fmprb\_t z}, \texttt{ulong x}, \texttt{slong prec})

Sets \texttt{z} to the reciprocal square root of \texttt{x}, rounded to \texttt{prec} bits. At high precision, this is faster than computing a square root.

void \texttt{fmprb\_root} (\texttt{fmprb\_t z}, \texttt{const fmprb\_t x}, \texttt{ulong k}, \texttt{slong prec})

Sets \texttt{z} to the \texttt{k}-th root of \texttt{x}, rounded to \texttt{prec} bits. As currently implemented, this function is only fast for small fixed \texttt{k}. For large \texttt{k} it is better to use \texttt{fmprb\_pow\_fmpq()} or \texttt{fmprb\_pow()}.  

void \texttt{fmprb\_agm} (\texttt{fmprb\_t z}, \texttt{const fmprb\_t x}, \texttt{const fmprb\_t y}, \texttt{slong prec})

Sets \texttt{z} to the arithmetic-geometric mean of \texttt{x} and \texttt{y}. 

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6.1 Credits and references

Arb is licensed GNU General Public License version 2, or any later version.

Fredrik Johansson is the main author. The project was started in 2012 as a numerical extension of FLINT, and the initial design was heavily based on FLINT 2.0 (with particular credit to Bill Hart and Sebastian Pancratz).

From 2012 to July 2014, Fredrik’s work on Arb was supported by Austrian Science Fund FWF Grant Y464-N18 (Fast Computer Algebra for Special Functions). During that period, he was a PhD student (and briefly a postdoc) at RISC, Johannes Kepler University, Linz, supervised by Manuel Kauers.

From September 2014 to October 2015, Fredrik was a postdoc in the LFANT project-team at INRIA Bordeaux and Institut de Mathématiques de Bordeaux, headed by Andreas Enge. During that period, Fredrik’s work on Arb was supported by ERC Starting Grant ANTICS 278537 (Algorithmic Number Theory in Computer Science) http://cordis.europa.eu/project/rcn/101288_en.html Since October 2015, Fredrik is a CR2 researcher in the LFANT team, funded by INRIA.

6.1.1 Contributors

Several people have contributed patches, bug reports, or substantial feedback. This list is probably incomplete.

- Bill Hart - build system, Windows 64 support, design of FLINT
- Sebastian Pancratz - divide-and-conquer polynomial composition algorithm (taken from FLINT)
- The MPFR development team - Arb includes two-limb multiplication code taken from MPFR
- Jonathan Bober - C++ compatibility fixes
- Yuri Matiyasevich - feedback about the zeta function and root-finding code
- Abhinav Baid - dot product and norm functions
- Ondřej Čertík - bug reports, feedback
- Andrew Booker - bug reports, feedback
- Francesco Biscani - C++ compatibility fixes, feedback
- Clemens Heuberger - work on Arb interface in Sage, feedback
- Marc Mezzarobba - work on Arb interface in Sage, bug reports, feedback
- Pascal Molin - feedback
- Ricky Farr - convenience functions, feedback
- Marcello Seri - fix for static builds on OS X
• Tommy Hofmann - matrix transpose, comparison, other utility methods, Julia interface
• Alexander Kobel - documentation and code cleanup patches
• Hrvoje Abraham - patches for MinGW compatibility
• Julien Puydt - soname versioning support
• Alex Griffing - sinc function, matrix trace, improved matrix squaring, miscellaneous patches
• Jeroen Demeyer - patch for major bug on PPC64

6.1.2 Software

The following software has been helpful in the development of Arb.

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• MPIR (Brian Gladman, Jason Moxham, William Hart and others), http://mpir.org
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• FLINT (William Hart, Sebastian Pancratz, Andy Novocin, Fredrik Johansson, David Harvey and others), http://flintlib.org
• Sage (William Stein and others), http://sagemath.org
• Pari/GP (The Pari group), http://pari.math.u-bordeaux.fr/
• SymPy (Ondřej Čertík, Aaron Meurer and others), http://sympy.org
• mpmath (Fredrik Johansson and others), http://mpmath.org
• Mathematica (Wolfram Research), http://www.wolfram.com/mathematica
• HolonomicFunctions (Christoph Koutschan), http://www.risc.jku.at/research/combinat/software/HolonomicFunctions/
• Sphinx (George Brandl and others), http://sphinx.pocoo.org
• CM (Andreas Enge), http://www.multiprecision.org/index.php?prog=cm

6.1.3 Citing Arb

If you wish to cite Arb in a scientific paper, the following reference can be used (you may also cite the manual or the website directly):


In BibTeX format:

@article{Johansson2013arb,
  title={Arb: a C library for ball arithmetic},
  author={F. Johansson},
  journal={ACM Communications in Computer Algebra},
  volume={47},
  number={4},
  pages={166--169},
  year={2013},
  publisher={ACM}
}
6.1.4 Bibliography
[WQ3b] http://functions.wolfram.com/07.12.27.0014.01


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