

More Algorithms for Arbitrary-Precision Sampling

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[Peter Occil](#)

Abstract: This page contains additional algorithms for arbitrary-precision sampling of continuous distributions, Bernoulli factory algorithms (biased-coin to biased-coin algorithms), and algorithms to simulate irrational probabilities. They supplement my pages on Bernoulli factory algorithms and partially-sampled random numbers.

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1 Introduction

This page contains additional algorithms for arbitrary-precision sampling of continuous distributions, Bernoulli factory algorithms (biased-coin to biased-coin algorithms), and algorithms to simulate irrational probabilities. These samplers are designed to not rely on floating-point arithmetic. They may depend on algorithms given in the following pages:

- [Partially-Sampled Random Numbers for Accurate Sampling of Continuous Distributions](#)
- [Bernoulli Factory Algorithms](#)

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3 Bernoulli Factories and Irrational Probability Simulation

In the methods below, λ is the unknown probability of heads of the coin involved in the Bernoulli Factory problem.

3.1 Certain Numbers Based on the Golden Ratio

The following algorithm given by Fishman and Miller (2013)⁽¹⁾ finds the continued fraction expansion of certain numbers described as—

- $G(m, \ell) = (m + \sqrt{m^2 + 4 * \ell})/2$
or $(m - \sqrt{m^2 + 4 * \ell})/2$,

whichever results in a real number greater than 1, where m is a positive integer and ℓ is either 1 or -1 . In this case, $G(1, 1)$ is the golden ratio.

First, define the following operations:

- **Get the previous and next Fibonacci-based number given k , m , and ℓ :**
 1. If k is 0 or less, return an error.
 2. Set $g0$ to 0, $g1$ to 1, x to 0, and y to 0.
 3. Do the following k times: Set y to $m * g1 + \ell * g0$, then set x to $g0$, then set $g0$ to $g1$, then set $g1$ to y .
 4. Return x and y , in that order.
- **Get the partial denominator given pos , k , m , and ℓ** (this partial denominator is part of the continued fraction expansion found by Fishman and Miller):
 1. **Get the previous and next Fibonacci-based number given k , m , and ℓ** , call them p and n , respectively.
 2. If ℓ is 1 and k is odd, return $p + n$.
 3. If ℓ is -1 and pos is 0, return $n - p - 1$.
 4. If ℓ is 1 and pos is 0, return $(n + p) - 1$.
 5. If ℓ is -1 and pos is even, return $n - p - 2$. (The paper had an error here; the correction given here was verified by Miller via personal communication.)
 6. If ℓ is 1 and pos is even, return $(n + p) - 2$.
 7. Return 1.

An application of the continued fraction algorithm is the following algorithm that generates 1 with probability $G(m, \ell)^{-k}$ and 0 otherwise, where k is an integer that is 1 or greater (see "Continued Fractions" in my page on Bernoulli factory algorithms). The algorithm starts with $pos = 0$, then the following steps are taken:

1. **Get the partial denominator given pos , k , m , and ℓ** , call it kp .
2. With probability $kp/(1 + kp)$, return a number that is 1 with probability $1/kp$ and 0 otherwise.
3. Run this algorithm recursively, but with $pos = pos + 1$. If the algorithm returns 1, return 0. Otherwise, go to step 2.

3.2 Ratio of Lower Gamma Functions ($\gamma(m, x)/\gamma(m, 1)$).

1. Set ret to the result of **kthsmallest** with the two parameters m and m . (Thus, ret is distributed as $u^{1/m}$ where u is a uniform(0, 1) random number; although **kthsmallest** accepts only integers, this formula works for any m greater than 0.)
2. Set k to 1, then set u to point to the same value as ret .
3. Generate a uniform(0, 1) random number v .
4. If v is less than u : Set u to v , then add 1 to k , then go to step 3.
5. If k is odd, return a number that is 1 if ret is less than x and 0 otherwise. (If ret is implemented as a uniform partially-sampled random number (PSRN), this comparison should be done via **URandLessThanReal**.) If k is even, go to step 1.

Derivation: See Formula 1 in the section "[Probabilities Arising from Certain Permutations](#)", where:

- $ECDF(x)$ is the uniform(0,1) distribution's cumulative distribution function, namely x if x is in $[0, 1]$, 0 if x is less than 0, and 1 otherwise.
- $DPDF(x)$ is the probability density function for the maximum of m uniform(0,1) random numbers, namely $m \cdot x^{m-1}$ if x is in $[0, 1]$, and 0 otherwise.

3.3 Derivative (slope) of $\arctan(\lambda)$

This algorithm involves the series expansion of this function ($1 - \lambda^2 + \lambda^4 - \dots$) and involves the general martingale algorithm.

1. Set u to 1, set w to 1, set ℓ to 0, and set n to 1.
2. Generate a uniform(0, 1) random number ret .
3. (Loop.) If w is not 0, flip the input coin and multiply w by the result of the flip. Do this step again.
4. If n is even, set u to $\ell + w$. Otherwise, set ℓ to $u - w$.
5. If ret is less than (or equal to) ℓ , return 1. If ret is less than u , go to the next step. If neither is the case, return 0. (If ret is a uniform PSRN, these comparisons should be done via the **URandLessThanReal algorithm**, which is described in my [article on PSRNs](#).)
6. Add 1 to n and go to step 3.

3.4 $\cosh(\lambda) - 1$

This algorithm involves an application of the general martingale algorithm to the Taylor series for $\cosh(\lambda) - 1$, which is $\lambda^2/(2!) + \lambda^4/(4!) + \dots$. See (Łatuszyński et al. 2009/2011,

algorithm 3)⁽²⁾.

1. Set u to 0, set w to 1, set ℓ to 0, and set n to 1.
2. Generate a uniform(0, 1) random number ret .
3. If w is not 0, flip the input coin and multiply w by the result of the flip. Do this step again.
4. If w is 0, set u to ℓ and go to step 6. (The estimate λ^{n*2} is 0, so no more terms are added and we use ℓ as the final estimate for $\cosh(\lambda)-1$.)
5. Let m be $(n*2)$, let α be $1/(m!)$ (a term of the Taylor series), and let err be $2/((m+1)!)$ (the error term). Add α to ℓ , then set u to $\ell + err$.
6. If ret is less than (or equal to) ℓ , return 1. If ret is less than u , go to the next step. If neither is the case, return 0. (If ret is a uniform PSRN, these comparisons should be done via the **URandLessThanReal algorithm**, which is described in my [article on PSRNs](#).)
7. Add 1 to n and go to step 3.

In this algorithm, the error term, which follows from *Taylor's theorem*, has a numerator of 2 because 2 is higher than the maximum value that the function's slope, slope-of-slope, etc. functions can achieve anywhere in the interval $[0, 1]$.

3.5 $\sinh(\lambda)/2$

This algorithm involves an application of the general martingale algorithm to the Taylor series for $\sinh(\lambda)/2$, which is $\lambda^1/(1!*2) + \lambda^3/(3!*2) + \dots$, or as used here, $\lambda*(1/2 + \lambda^2/(3!*2) + \lambda^4/(5!*2) + \dots)$.

1. Flip the input coin. If it returns 0, return 0.
2. Set u to 0, set w to 1, set ℓ to $1/2$ (the first term is added already), and set n to 1.
3. Generate a uniform(0, 1) random number ret .
4. If w is not 0, flip the input coin and multiply w by the result of the flip. Do this step again.
5. If w is 0, set u to ℓ and go to step 7. (No more terms are added here.)
6. Let m be $(n*2+1)$, let α be $1/(m!*2)$ (a term of the Taylor series), and let err be $1/((m+1)!)$ (the error term). Add α to ℓ , then set u to $\ell + err$.
7. If ret is less than (or equal to) ℓ , return 1. If ret is less than u , go to the next step. If neither is the case, return 0. (If ret is a uniform PSRN, these comparisons should be done via the **URandLessThanReal algorithm**, which is described in my [article on PSRNs](#).)
8. Add 1 to n and go to step 4.

3.6 $\tanh(\lambda)$

There are two algorithms.

The first takes advantage of the so-called Lambert's continued fraction for $\tanh(\cdot)$, as well as Bernoulli Factory algorithm 3 for continued fractions. The algorithm begins with k equal to 1. Then the following steps are taken.

1. If k is 1: With probability $1/2$, flip the input coin and return the result.
2. If k is greater than 1, then do the following with probability $k/(1+k)$:
 - Flip the input coin twice. If any of these flips returns 0, return 0. Otherwise, return a number that is 1 with probability $1/k$ and 0 otherwise.
3. Run this algorithm recursively, but with $k = k + 2$. If the result is 1, return 0. Otherwise, go to step 1.

The second algorithm involves an alternating series expansion of $\tanh(\cdot)$ and involves the general martingale algorithm.

First, define the following operation:

- **Get the m^{th} Bernoulli number:**
 1. If m is 0, 1, 2, 3, or 4, return 1, $-1/2$, $1/6$, 0, or $-1/30$, respectively. Otherwise, if m is odd, return 0.
 2. Set i to 2 and v to $1 - (m+1)/2$.
 3. While i is less than m :
 1. **Get the i^{th} Bernoulli number**, call it b . Add $b \cdot \text{choose}(m+1, i)$ to v .⁽³⁾
 2. Add 2 to i .
 4. Return $-v/(m+1)$.

The algorithm is then as follows:

1. Flip the input coin. If it returns 0, return 0.
2. Set u to 1, set w to 1, set ℓ to 0, and set n to 1.
3. Generate a uniform(0, 1) random number ret .
4. (Loop.) If w is not 0, flip the input coin. If the flip returns 0, set w to 0. Do this step again.
5. (Calculate the next term of the alternating series for $\tanh(\cdot)$) Let m be $2 \cdot (n+1)$. **Get the m^{th} Bernoulli number**, call it b . Let t be $\text{abs}(b) \cdot 2^{m \cdot (2^m - 1)} / (m!)$.
6. If n is even, set u to $\ell + w \cdot t$. Otherwise, set ℓ to $u - w \cdot t$.
7. If ret is less than (or equal to) ℓ , return 1. If ret is less than u , go to the next step. If neither is the case, return 0. (If ret is a uniform PSRN, these comparisons should be done via the **URandLessThanReal algorithm**, which is described in my [article on PSRNs](#).)
8. Add 1 to n and go to step 4.

3.7 Certain Piecewise Linear Functions

Let $f(\lambda)$ be a function of the form $\min(\lambda \cdot \text{mult}, 1 - \varepsilon)$. (This is a piecewise linear function with two pieces: a rising linear part and a constant part.) This section describes how to calculate the Bernstein coefficients for polynomials that converge from above and below to f , based on Thomas and Blanchet (2012)⁽⁴⁾. These polynomials can then be used to generate heads with probability $f(\lambda)$ using the algorithms given in "[General Factory Functions](#)". In this section, **fbelow(n, k)** and **fabove(n, k)** are the k^{th} Bernstein coefficients (with k starting at 0) of the lower and upper polynomials, respectively, of degree n .

The code in the **appendix** uses the computer algebra library SymPy to calculate a list of parameters for a sequence of polynomials converging from above. The method to do so is called `calc_linear_func(eps, mult, count)`, where `eps` is ε , `mult` = mult , and `count` is the number of polynomials to generate. Each item returned by `calc_linear_func` is a list of two items: the degree of the polynomial, and a *Y parameter*. The procedure to calculate the required polynomials is then logically as follows (as written, it runs very slowly, though):

1. Set i to 1.
2. Run `calc_linear_func(eps, mult, i)` and get the degree and *Y parameter* for the last listed item, call them n and y , respectively.
3. Set x to $-((y - (1 - \varepsilon))/\varepsilon)^5 / \text{mult} + y / \text{mult}$. (This exact formula doesn't appear in the Thomas and Blanchet paper; rather it comes from the [supplemental source code](#) uploaded by A. C. Thomas at my request.

4. For degree n , **fbelow**(n, k) is $\min((k/n)*mult, 1-\varepsilon)$, and **fabove**(n, k) is $\min((k/n)*y/x, y)$. (**fbelow** matches f because f is *concave* in the interval $[0, 1]$, which roughly means that its rate of growth there never goes up.)
5. Add 1 to i and go to step 2.

It would be interesting to find general formulas to find the appropriate polynomials (degrees and Y parameters) given only the values for $mult$ and ε , rather than find them "the hard way" via `calc_linear_func`. For this procedure, the degrees and Y parameters can be upper bounds, as long as the sequence of degrees is monotonically increasing and the sequence of Y parameters is nonincreasing.

Note: In Nacu and Peres (2005)⁽⁵⁾, the following polynomial sequences were suggested to simulate $\min(\lambda^2, 1 - 2\varepsilon)$, provided $\varepsilon < 1/8$, where n is a power of 2. However, with these sequences, an extraordinary number of input coin flips is required to simulate this function each time.

- **fbelow**(n, k) = $\min((k/n)^2, 1 - 2\varepsilon)$.
- **fabove**(n, k) = $\min((k/n)^2, 1 - 2\varepsilon) + (\max(0, k/n + 3\varepsilon - 1/2)/(\varepsilon/(1-\sqrt{2}/2))) * \sqrt{2/n} + (72 * \max(0, k/n - 1/9)/(1 - \exp(-2\varepsilon))) * \exp(-2\varepsilon * n)$.

My own algorithm for $\min(\lambda, 1/2)$ is as follows. See the [appendix](#) for the derivation of this algorithm.

1. With probability $1/2$, flip the input coin and return the result.
2. (Random walk.) Generate unbiased random bits until more zeros than ones are generated this way for the first time. Then set m to $(n-1)/2+1$, where n is the number of bits generated this way.
3. (Build a degree- m^2 polynomial equivalent to $(4\lambda*(1-\lambda))^m/2$.) Let z be $(4^m/2)/\text{choose}(m^2, m)$. Define a polynomial of degree m^2 whose $(m^2)+1$ Bernstein coefficients are all zero except the m^{th} coefficient (starting at 0), whose value is z . Elevate the degree of this polynomial enough times so that all its coefficients are 1 or less (degree elevation increases the polynomial's degree without changing its shape or position; see the derivation in the appendix). Let d be the new polynomial's degree.
4. (Simulate the polynomial, whose degree is d (Goyal and Sigman 2012)⁽⁶⁾.) Flip the input coin d times and set h to the number of ones generated this way. Let a be the h^{th} Bernstein coefficient (starting at 0) of the new polynomial. With probability a , return 1. Otherwise, return 0.

I suspected that the required degree d would be $\text{floor}(m^2/3)+1$, as described in the appendix. With help from the [MathOverflow community](#), steps 3 and 4 of the algorithm can be described more efficiently as follows:

- (3.) Let r be $\text{floor}(m^2/3)+1$, and let d be m^2+r .
- (4.) (Simulate the polynomial, whose degree is d .) Flip the input coin d times and set h to the number of ones generated this way. Let a be $(1/2) * 2^{m^2} * \text{choose}(r, h-m) / \text{choose}(d, h)$ (the polynomial's h^{th} Bernstein coefficient starting at 0; the first term is $1/2$ because the polynomial being simulated has the value $1/2$ at the point $1/2$). With probability a , return 1. Otherwise, return 0.

3.8 Non-Negative Factories

The Bernoulli factory approach can be extended in two ways to produce random numbers

beyond the interval $[0, 1]$. Both algorithms use a different *oracle* (black box) from coins that output heads or tails.

Algorithm 1. Say we have an oracle that produces random numbers in the interval $[a, b]$, and these numbers have an unknown mean of μ . The goal is now to produce non-negative random numbers that average to $f(\mu)$. This is possible if and only if f , in the interval $[a, b]$

- is continuous everywhere,
- is bounded from below,
- does not go to 0 exponentially fast in value, and
- returns 0 everywhere or returns a value 0 or greater at each of the points a and b and a value greater than 0 at each other point

(Jacob and Thiery 2015)⁽⁷⁾. (Here, a and b are both rational numbers and may be less than 0.)

In the algorithm below, let κ be greater than the maximum value of f in the interval $[a, b]$, and let $g(\lambda) = f(a + (b-a)\lambda)/\kappa$.

1. Create a λ input coin that does the following: "Generate a random number from the oracle, call it x . With probability $(x-a)/(b-a)$ (see note below), return 1. Otherwise, return 0."
2. Run a Bernoulli factory algorithm for $g(\lambda)$, using the λ input coin. Then return κ times the result.

Note: The check "With probability $(x-a)/(b-a)$ " is exact if the oracle produces only rational numbers. If the oracle can produce irrational numbers (such as numbers that follow a beta distribution or another continuous distribution), then the code for the oracle should use uniform [partially-sampled random numbers \(PSRNs\)](#). In that case, the check can be implemented as follows. Let x be a uniform PSRN representing a number generated by the oracle. Set y to **RandUniformFromReal**($b-a$), then the check succeeds if **URandLess**(y , **UniformAddRational**($x, -a$)) returns 1, and fails otherwise.

Example: Suppose an oracle produces random numbers in the interval $[3, 13]$ with mean μ , and we seek to use the oracle to produce non-negative random numbers with mean $f(\mu) = -319/100 + \mu*103/50 - \mu^2*11/100$, which is a polynomial with Bernstein coefficients $[2, 9, 5]$ in the given interval. Then since 8 is greater than the maximum of f in that interval, $g(\lambda)$ is a degree-2 polynomial with Bernstein coefficients $[2/8, 9/8, 5/8]$ in the interval $[0, 1]$. g can't be simulated as is, though, but by increasing g 's degree to 3 we get the Bernstein coefficients $[1/4, 5/6, 23/24, 5/8]$, which are all less than 1 so we can proceed with the following algorithm (see "[Certain Polynomials in Bernstein Form](#)"):

1. Set *heads* to 0.
2. Generate three random numbers from the oracle (which must produce random numbers in the interval $[3, 13]$). For each number x : With probability $(x-3)/(10-3)$, add 1 to *heads*.
3. Depending on *heads*, return 8 (that is, 1 times the upper bound) with the given probability, or 0 otherwise: *heads*=0 \rightarrow probability 1/4; 1 \rightarrow 5/6; 2 \rightarrow 23/24; 3 \rightarrow 5/8.

Algorithm 2. This algorithm takes an oracle and produces non-negative random numbers that average to the mean of $f(X)$, where X is a number produced by the oracle.

The algorithm appears in the appendix, however, because it requires applying an arbitrary function (here, f) to a potentially irrational number.

4 General Arbitrary-Precision Samplers

4.1 Uniform Distribution Inside N-Dimensional Shapes

The following is a general way to describe an arbitrary-precision sampler for generating a point uniformly at random inside a geometric shape located entirely in the hypercube $[0, d1] \times [0, d2] \times \dots \times [0, dN]$ in N -dimensional space, where $d1, \dots, dN$ are integers greater than 0. The algorithm will generally work if the shape is reasonably defined; the technical requirements are that the shape must have a zero-volume boundary and a nonzero finite volume, and must assign zero probability to any zero-volume subset of it (such as a set of individual points).

The sampler's description has the following skeleton.

1. Generate N empty uniform partially-sampled random numbers (PSRNs), with a positive sign, an integer part of 0, and an empty fractional part. Call the PSRNs $p1, p2, \dots, pN$.
2. Set S to $base$, where $base$ is the base of digits to be stored by the PSRNs (such as 2 for binary or 10 for decimal). Then set N coordinates to 0, call the coordinates $c1, c2, \dots, cN$. Then set d to 1. Then, for each coordinate ($c1, \dots, cN$), set that coordinate to an integer in $0, dX$, chosen uniformly at random, where dX is the corresponding dimension's size.
3. For each coordinate ($c1, \dots, cN$), multiply that coordinate by $base$ and add a digit chosen uniformly at random to that coordinate.
4. This step uses a function known as **InShape**, which takes the coordinates of a box and returns one of three values: *YES* if the box is entirely inside the shape; *NO* if the box is entirely outside the shape; and *MAYBE* if the box is partly inside and partly outside the shape, or if the function is unsure. **InShape**, as well as the divisions of the coordinates by S , should be implemented using rational arithmetic. Instead of dividing those coordinates this way, an implementation can pass S as a separate parameter to **InShape**. See the [appendix for further implementation notes. In this step, run **InShape** using the current box, whose coordinates in this case are $((c1/S, c2/S, \dots, cN/S), ((c1+1)/S, (c2+1)/S, \dots, (cN+1)/S))$.
5. If the result of **InShape** is *YES*, then the current box was accepted. If the box is accepted this way, then at this point, $c1, c2$, etc., will each store the d digits of a coordinate in the shape, expressed as a number in the interval $[0, 1]$, or more precisely, a range of numbers. (For example, if $base$ is 10, d is 3, and $c1$ is 342, then the first coordinate is 0.342, or more precisely, a number in the interval $[0.342, 0.343]$.) In this case, do the following:
 1. For each coordinate ($c1, \dots, cN$), transfer that coordinate's least significant digits to the corresponding PSRN's fractional part. The variable d tells how many digits to transfer to each PSRN this way. Then, for each coordinate ($c1, \dots, cN$), set the corresponding PSRN's integer part to $\text{floor}(cX/base^d)$, where cX is that coordinate. (For example, if $base$ is 10, d is 3, and $c1$ is 7342, set $p1$'s fractional part to $[3, 4, 2]$ and $p1$'s integer part to 7.)
 2. For each PSRN ($p1, \dots, pN$), optionally fill that PSRN with uniform random digits as necessary to give its fractional part the desired number of digits

(similarly to **FillGeometricBag**).

3. For each PSRN, optionally do the following: With probability $1/2$, set that PSRN's sign to negative. (This will result in a symmetric shape in the corresponding dimension. This step can be done for some PSRNs and not others.)
4. Return the PSRNs p_1, \dots, p_N , in that order.
6. If the result of **InShape** is *NO*, then the current box lies outside the shape and is rejected. In this case, go to step 2.
7. If the result of **InShape** is *MAYBE*, it is not known whether the current box lies fully inside the shape, so multiply S by $base$, then add 1 to d , then go to step 3.

Notes:

- See (Li and El Gamal 2016)⁽⁸⁾ and (Oberhoff 2018)⁽⁹⁾ for related work on encoding random points uniformly distributed in a shape.
- Rejection sampling on a shape is subject to the "curse of dimensionality", since typical shapes of high dimension will tend to cover much less volume than their bounding boxes, so that it would take a lot of time on average to accept a high-dimensional box. Moreover, the more area the shape takes up in the bounding box, the higher the acceptance rate.
- Devroye (1986, chapter 8, section 3)⁽¹⁰⁾ describes grid-based methods to optimize random point generation. In this case, the space is divided into a grid of boxes each with size $1/base^k$ in all dimensions; the result of **InShape** is calculated for each such box and that box labeled with the result; all boxes labeled *NO* are discarded; and the algorithm is modified by adding the following after step 2: "2a. Choose a precalculated box uniformly at random, then set c_1, \dots, c_N to that box's coordinates, then set d to k and set S to $base^k$. If a box labeled *YES* was chosen, follow the substeps in step 5. If a box labeled *MAYBE* was chosen, multiply S by $base$ and add 1 to d ." (For example, if $base$ is 10, k is 1, N is 2, and $d_1 = d_2 = 1$, the space could be divided into a 10×10 grid, made up of 100 boxes each of size $(1/10) \times (1/10)$. Then, **InShape** is precalculated for the box with coordinates $((0, 0), (1, 1))$, the box $((0, 1), (1, 2))$, and so on [the boxes' coordinates are stored as just given, but **InShape** instead uses those coordinates divided by $base^k$, or 10^1 in this case], each such box is labeled with the result, and boxes labeled *NO* are discarded. Finally the algorithm above is modified as just given.)
- Besides a grid, another useful data structure is a *mapped regular paving* (Harlow et al. 2012)⁽¹¹⁾, which can be described as a binary tree with nodes each consisting of zero or two child nodes and a marking value. Start with a box that entirely covers the desired shape. Calculate **InShape** for the box. If it returns *YES* or *NO* then mark the box with *YES* or *NO*, respectively; otherwise it returns *MAYBE*, so divide the box along its first widest coordinate into two sub-boxes, set the parent box's children to those sub-boxes, then repeat this process for each sub-box (or if the nesting level is too deep, instead mark each sub-box with *MAYBE*). Then, to generate a random point (with a base-2 fractional part), start from the root, then: (1) If the box is marked *YES*, return a uniform random point between the given coordinates using the **RandUniformInRange** algorithm; or (2) if the box is marked *NO*, start over from the root; or (3) if the box is marked *MAYBE*, get the two child boxes bisected from the box, choose one of them with equal probability (e.g., choose the left child if an unbiased random bit is 0, or the right child otherwise), mark the chosen child with the result of **InShape** for that child, and repeat this process

with that child; or (4) the box has two child boxes, so choose one of them with equal probability and repeat this process with that child.

Examples:

- The following example generates a point inside a quarter diamond (centered at $(0, \dots, 0)$, "radius" k where k is an integer greater than 0): Let $d1, \dots, dN$ be k . Let **InShape** return *YES* if $((c1+1) + \dots + (cN+1)) < S*k$; *NO* if $(c1 + \dots + cN) > S*k$; and *MAYBE* otherwise. For a full diamond, step 5.3 in the algorithm is done for all N dimensions.
- The following example generates a point inside a quarter hypersphere (centered at $(0, \dots, 0)$, radius k where k is an integer greater than 0): Let $d1, \dots, dN$ be k . Let **InShape** return *YES* if $((c1+1)^2 + \dots + (cN+1)^2) < (S*k)^2$; *NO* if $(c1^2 + \dots + cN^2) > (S*k)^2$; and *MAYBE* otherwise. For a full hypersphere with radius 1, step 5.3 in the algorithm is done for all N dimensions. In the case of a 2-dimensional circle, this algorithm thus adapts the well-known rejection technique of generating X and Y coordinates until $X^2+Y^2 < 1$ (e.g., (Devroye 1986, p. 230 et seq.)⁽¹⁰⁾).
- The following example generates a point inside a quarter *astroid* (centered at $(0, \dots, 0)$, radius k where k is an integer greater than 0): Let $d1, \dots, dN$ be k . Let **InShape** return *YES* if $((sk-c1-1)^2 + \dots + (sk-cN-1)^2) > sk^2$; *NO* if $((sk-c1)^2 + \dots + (sk-cN)^2) < sk^2$; and *MAYBE* otherwise, where $sk = S*k$. For a full astroid, step 5.3 in the algorithm is done for all N dimensions.

4.2 Building an Arbitrary-Precision Sampler

In many cases, if a continuous distribution—

- has a probability density function (PDF), or a function proportional to the PDF, with a known symbolic form,
- has a cumulative distribution function (CDF) with a known symbolic form,
- takes on only values 0 or greater, and
- has a PDF that has an infinite tail to the right, is bounded from above (that is, $PDF(0)$ is other than infinity), and decreases monotonically,

it may be possible to describe an arbitrary-precision sampler for that distribution. Such a description has the following skeleton.

1. With probability A , set *intval* to 0, then set *size* to 1, then go to step 4.
 - A is calculated as $(CDF(1) - CDF(0)) / (1 - CDF(0))$, where CDF is the distribution's CDF. This should be found analytically using a computer algebra system such as SymPy.
 - The symbolic form of A will help determine which Bernoulli factory algorithm, if any, will simulate the probability; if a Bernoulli factory exists, it should be used.
2. Set *intval* to 1 and set *size* to 1.
3. With probability $B(\text{size}, \text{intval})$, go to step 4. Otherwise, add *size* to *intval*, then multiply *size* by 2, then repeat this step.
 - This step chooses an interval beyond 1, and grows this interval by geometric steps, so that an appropriate interval is chosen with the correct probability.
 - The probability $B(\text{size}, \text{intval})$ is the probability that the interval is chosen given that the previous intervals weren't chosen, and is calculated as $(CDF(\text{size} + \text{intval}) - CDF(\text{intval})) / (1 - CDF(\text{intval}))$. This should be found analytically using a computer algebra system such as SymPy.
 - The symbolic form of B will help determine which Bernoulli factory algorithm, if

- any, will simulate the probability; if a Bernoulli factory exists, it should be used.
4. Generate an integer in the interval $[intval, intval + size)$ uniformly at random, call it i .
 5. Create a positive-sign zero-integer-part uniform PSRN, ret .
 6. Create an input coin that calls **SampleGeometricBag** on the PSRN ret . Run a Bernoulli factory algorithm that simulates the probability $C(i, \lambda)$, using the input coin (here, λ is the probability built up in ret via **SampleGeometricBag**, and lies in the interval $[0, 1]$). If the call returns 0, go to step 4.
 - The probability $C(i, \lambda)$ is calculated as $PDF(i + \lambda) / M$, where PDF is the distribution's PDF or a function proportional to the PDF, and should be found analytically using a computer algebra system such as SymPy.
 - In this formula, M is any convenient number in the interval $[PDF(intval), \max(1, PDF(intval))]$, and should be as low as feasible. M serves to ensure that C is as close as feasible to 1 (to improve acceptance rates), but no higher than 1. The choice of M can vary for each interval (each value of $intval$, which can only be 0, 1, or a power of 2). Any such choice for M preserves the algorithm's correctness because the PDF has to be monotonically decreasing and a new interval isn't chosen when λ is rejected.
 - The symbolic form of C will help determine which Bernoulli factory algorithm, if any, will simulate the probability; if a Bernoulli factory exists, it should be used.
 7. The PSRN ret was accepted, so set ret 's integer part to i , then optionally fill ret with uniform random digits as necessary to give its fractional part the desired number of digits (similarly to **FillGeometricBag**), then return ret .

Examples of algorithms that use this skeleton are the algorithm for the [ratio of two uniform random numbers](#), as well as the algorithms for the Rayleigh distribution and for the reciprocal of power of uniform, both given later.

Perhaps the most difficult part of describing an arbitrary-precision sampler with this skeleton is finding the appropriate Bernoulli factory for the probabilities A , B , and C , especially when these probabilities have a non-trivial symbolic form.

Note: The algorithm skeleton uses ideas similar to the inversion-rejection method described in (Devroye 1986, ch. 7, sec. 4.6)⁽¹⁰⁾; an exception is that instead of generating a uniform random number and comparing it to calculations of a CDF, this algorithm uses conditional probabilities of choosing a given piece, probabilities labeled A and B . This approach was taken so that the CDF of the distribution in question is never directly calculated in the course of the algorithm, which furthers the goal of sampling with arbitrary precision and without using floating-point arithmetic.

4.3 Mixtures

A *mixture* involves sampling one of several distributions, where each distribution has a separate probability of being sampled. In general, an arbitrary-precision sampler is possible if all of the following conditions are met:

- There is a finite number of distributions to choose from.
- The probability of sampling each distribution is a rational number, or it can be expressed as a function for which a [Bernoulli factory algorithm](#) exists.
- For each distribution, an arbitrary-precision sampler exists.

Example: One example of a mixture is two beta distributions, with separate parameters. One beta distribution is chosen with probability $\exp(-3)$ (a probability for which a Bernoulli factory algorithm exists) and the other is

chosen with the opposite probability. For the two beta distributions, an arbitrary-precision sampling algorithm exists (see my article on [partially-sampled random numbers \(PSRNs\)](#) for details).

5 Specific Arbitrary-Precision Samplers

5.1 Rayleigh Distribution

The following is an arbitrary-precision sampler for the Rayleigh distribution with parameter s , which is a rational number greater than 0.

1. Set k to 0, and set y to $2 * s * s$.
2. With probability $\exp(-(k * 2 + 1)/y)$, go to step 3. Otherwise, add 1 to k and repeat this step. (The probability check should be done with the **`exp(-x/y)` algorithm** in "[Bernoulli Factory Algorithms](#)", with $x/y = (k * 2 + 1)/y$.)
3. (Now we sample the piece located at $[k, k + 1)$.) Create a positive-sign zero-integer-part uniform PSRN, and create an input coin that returns the result of **`SampleGeometricBag`** on that uniform PSRN.
4. Set ky to $k * k / y$.
5. (At this point, we simulate $\exp(-U^2/y)$, $\exp(-k^2/y)$, $\exp(-U*k^2/y)$, as well as a scaled-down version of $U + k$, where U is the number built up by the uniform PSRN.) Call the **`exp(-x/y)` algorithm** with $x/y = ky$, then call the **`exp(-(\lambda^k * x))` algorithm** using the input coin from step 2, $x = 1/y$, and $k = 2$, then call the **`exp(-(\lambda^k * (x+m)))` algorithm** using the same input coin, $x+m = \text{floor}(k * 2 / y)$, and $k = 1$, then call the **`sub-algorithm`** given later with the uniform PSRN and $k = k$. If all of these calls return 1, the uniform PSRN was accepted. Otherwise, remove all digits from the uniform PSRN's fractional part and go to step 4.
6. If the uniform PSRN, call it *ret*, was accepted by step 5, set *ret*'s integer part to k , then optionally fill *ret* with uniform random digits as necessary to give its fractional part the desired number of digits (similarly to **`FillGeometricBag`**), and return *ret*.

The sub-algorithm below simulates a probability equal to $(U+k)/base^z$, where U is the number built by the uniform PSRN, *base* is the base (radix) of digits stored by that PSRN, k is an integer 0 or greater, and z is the number of significant digits in k (for this purpose, z is 0 if k is 0).

For base 2:

1. Set N to 0.
2. With probability 1/2, go to the next step. Otherwise, add 1 to N and repeat this step.
3. If N is less than z , return $\text{rem}(k / 2^{z-1-N}, 2)$. (Alternatively, shift k to the right, by $z - 1 - N$ bits, then return $k \text{ AND } 1$, where "AND" is a bitwise AND-operation.)
4. Subtract z from N . Then, if the item at position N in the uniform PSRN's fractional part (positions start at 0) is not set to a digit (e.g., 0 or 1 for base 2), set the item at that position to a digit chosen uniformly at random (e.g., either 0 or 1 for base 2), increasing the capacity of the uniform PSRN's fractional part as necessary.
5. Return the item at position N .

For bases other than 2, such as 10 for decimal, this can be implemented as follows (based on **`URandLess`**):

1. Set i to 0.
2. If i is less than z :
 1. Set da to $\text{rem}(k / 2^z - 1 - i, \text{base})$, and set db to a digit chosen uniformly at random (that is, an integer in the interval $[0, \text{base})$).
 2. Return 1 if da is less than db , or 0 if da is greater than db .
3. If i is z or greater:
 1. If the digit at position $(i - z)$ in the uniform PSRN's fractional part is not set, set the item at that position to a digit chosen uniformly at random (positions start at 0 where 0 is the most significant digit after the point, 1 is the next, etc.).
 2. Set da to the item at that position, and set db to a digit chosen uniformly at random (that is, an integer in the interval $[0, \text{base})$).
 3. Return 1 if da is less than db , or 0 if da is greater than db .
4. Add 1 to i and go to step 3.

5.2 Sum of Exponential Random Numbers

An arbitrary-precision sampler for the sum of n exponential random numbers (also known as the Erlang(n) or gamma(n) distribution) is doable via partially-sampled uniform random numbers, though it is obviously inefficient for large values of n .

1. Generate n exponential random numbers with a rate of 1 via the **ExpRand** or **ExpRand2** algorithm described in my article on [partially-sampled random numbers \(PSRNs\)](#). These numbers will be uniform PSRNs; this algorithm won't work for exponential PSRNs (e-rands), described in the same article, because the sum of two e-rands may follow a subtly wrong distribution. By contrast, generating exponential random numbers via rejection from the uniform distribution will allow unsampled digits to be sampled uniformly at random without deviating from the exponential distribution.
2. Generate the sum of the random numbers generated in step 1 by applying the [UniformAdd](#) algorithm given in another document.

5.3 Hyperbolic Secant Distribution

The following algorithm adapts the rejection algorithm from p. 472 in (Devroye 1986)⁽¹⁰⁾ for arbitrary-precision sampling.

1. Generate a uniform PSRN, call it ret , and turn it into an exponential random number with a rate of 1, using an algorithm that employs rejection from the uniform distribution.
2. Set ip to 1 plus ret 's integer part.
3. (The rest of the algorithm accepts ret with probability $1/(1+ret)$.) With probability $ip/(1+ip)$, generate a number that is 1 with probability $1/ip$ and 0 otherwise. If that number is 1, ret was accepted, in which case optionally fill it with uniform random digits as necessary to give its fractional part the desired number of digits (similarly to **FillGeometricBag**), then set ret 's sign to positive or negative with equal probability, then return ret .
4. Call **SampleGeometricBag** on ret 's fractional part (ignore ret 's integer part and sign). If the call returns 1, go to step 1. Otherwise, go to step 3.

5.4 Reciprocal of Power of Uniform

The following algorithm generates a PSRN of the form $1/U^{1/x}$, where U is a uniform random number in $[0, 1]$ and x is an integer greater than 0.

1. Set *intval* to 1 and set *size* to 1.
2. With probability $(4^x - 2^x)/4^x$, go to step 3. Otherwise, add *size* to *intval*, then multiply *size* by 2, then repeat this step.
3. Generate an integer in the interval [*intval*, *intval* + *size*) uniformly at random, call it *i*.
4. Create a positive-sign zero-integer-part uniform PSRN, *ret*.
5. Create an input coin that calls **SampleGeometricBag** on the PSRN *ret*. Call the **algorithm for $d^k / (c + \lambda)^k$** in "[Bernoulli Factory Algorithms](#)", using the input coin, where $d = \text{intval}$, $c = i$, and $k = x + 1$ (here, λ is the probability built up in *ret* via **SampleGeometricBag**, and lies in the interval [0, 1]). If the call returns 0, go to step 3.
6. The PSRN *ret* was accepted, so set *ret*'s integer part to *i*, then optionally fill *ret* with uniform random digits as necessary to give its fractional part the desired number of digits (similarly to **FillGeometricBag**), then return *ret*.

This algorithm uses the skeleton described earlier in "Building an Arbitrary-Precision Sampler". Here, the probabilities *A*, *B*, and *C* are as follows:

- $A = 0$, since the random number can't lie in the interval [0, 1).
- $B = (4^x - 2^x)/4^x$.
- $C = (x/(i + \lambda)^{x+1}) / M$. Ideally, *M* is either *x* if *intval* is 1, or x/intval^{x+1} otherwise. Thus, the ideal form for *C* is $\text{intval}^{x+1}/(i+\lambda)^{x+1}$.

5.5 Distribution of $U/(1-U)$

The following algorithm generates a PSRN of the form $U/(1-U)$, where *U* is a uniform random number in [0, 1].

1. With probability 1/2, set *intval* to 0, then set *size* to 1, then go to step 4.
2. Set *intval* to 1 and set *size* to 1.
3. With probability $\text{size}/(\text{size} + \text{intval} + 1)$, go to step 4. Otherwise, add *size* to *intval*, then multiply *size* by 2, then repeat this step.
4. Generate an integer in the interval [*intval*, *intval* + *size*) uniformly at random, call it *i*.
5. Create a positive-sign zero-integer-part uniform PSRN, *ret*.
6. Create an input coin that calls **SampleGeometricBag** on the PSRN *ret*. Call the **algorithm for $d^k / (c + \lambda)^k$** in "[Bernoulli Factory Algorithms](#)", using the input coin, where $d = \text{intval} + 1$, $c = i + 1$, and $k = 2$ (here, λ is the probability built up in *ret* via **SampleGeometricBag**, and lies in the interval [0, 1]). If the call returns 0, go to step 4.
7. The PSRN *ret* was accepted, so set *ret*'s integer part to *i*, then optionally fill *ret* with uniform random digits as necessary to give its fractional part the desired number of digits (similarly to **FillGeometricBag**), then return *ret*.

This algorithm uses the skeleton described earlier in "Building an Arbitrary-Precision Sampler". Here, the probabilities *A*, *B*, and *C* are as follows:

- $A = 1/2$.
- $B = \text{size}/(\text{size} + \text{intval} + 1)$.
- $C = (1/(i+\lambda+1)^2) / M$. Ideally, *M* is $1/(\text{intval}+1)^2$. Thus, the ideal form for *C* is $(\text{intval}+1)^2/(i+\lambda+1)^2$.

5.6 Arc-Cosine Distribution

Here we reimplement an example from Devroye's book *Non-Uniform Random Variate Generation* (Devroye 1986, pp. 128-129)⁽¹⁰⁾. The following arbitrary-precision sampler generates a random number from a distribution with the following cumulative distribution function (CDF): $1 - \cos(\pi x/2)$. The random number will be in the interval $[0, 1]$. Note that the result is the same as applying $\text{acos}(U)*2/\pi$, where U is a uniform $[0, 1]$ random number, as pointed out by Devroye. The algorithm follows.

1. Call the **kthsmallest** algorithm with $n = 2$ and $k = 2$, but without filling it with digits at the last step. Let *ret* be the result.
2. Set m to 1.
3. Call the **kthsmallest** algorithm with $n = 2$ and $k = 2$, but without filling it with digits at the last step. Let u be the result.
4. With probability $4/(4*m*m + 2*m)$, call the **URandLess** algorithm with parameters u and *ret* in that order, and if that call returns 1, call the **algorithm for $\pi/4$** , described in "[Bernoulli Factory Algorithms](#)", twice, and if both of these calls return 1, add 1 to m and go to step 3. (Here, we incorporate an erratum in the algorithm on page 129 of the book.)
5. If m is odd, optionally fill *ret* with uniform random digits as necessary to give its fractional part the desired number of digits (similarly to **FillGeometricBag**), and return *ret*.
6. If m is even, go to step 1.

And here is Python code that implements this algorithm. Note that it uses floating-point arithmetic only at the end, to convert the result to a convenient form, and that it relies on methods from *randomgen.py* and *bernoulli.py*.

```
def example_4_2_1(rg, bern, precision=53):
    while True:
        ret=rg.kthsmallest_psrn(2,2)
        k=1
        while True:
            u=rg.kthsmallest_psrn(2,2)
            kden=4*k*k+2*k # erratum incorporated
            if randomgen.urandless(rg,u, ret) and \
                rg.zero_or_one(4, kden)==1 and \
                bern.zero_or_one_pi_div_4()==1 and \
                bern.zero_or_one_pi_div_4()==1:
                k+=1
            elif (k&1)==1:
                return randomgen.urandfill(rg,ret,precision)/(1<<precision)
            else: break
```

5.7 Logistic Distribution

The following new algorithm generates a partially-sampled random number that follows the logistic distribution.

1. Set k to 0.
2. (Choose a 1-unit-wide piece of the logistic density.) Run the **algorithm for $(1+\exp(k))/(1+\exp(k+1))$** described in "[Bernoulli Factory Algorithms](#)". If the call returns 0, add 1 to k and repeat this step. Otherwise, go to step 3.
3. (The rest of the algorithm samples from the chosen piece.) Generate a uniform $(0, 1)$ random number, call it f .
4. (Steps 4 through 7 succeed with probability $\exp(-(f+k))/(1+\exp(-(f+k)))^2$.) With probability $1/2$, go to step 3.
5. Run the **algorithm for $\exp(-k/1)$** (described in "[Bernoulli Factory Algorithms](#)"),

- then **sample from the number f** (e.g., call **SampleGeometricBag** on f if f is implemented as a uniform PSRN). If any of these calls returns 0, go to step 4.
6. With probability $1/2$, accept f . If f is accepted this way, set f 's integer part to k , then optionally fill f with uniform random digits as necessary to give its fractional part the desired number of digits (similarly to **FillGeometricBag**), then set f 's sign to positive or negative with equal probability, then return f .
 7. Run the **algorithm for $\exp(-k/1)$** and **sample from the number f** (e.g., call **SampleGeometricBag** on f if f is implemented as a uniform PSRN). If both calls return 1, go to step 3. Otherwise, go to step 6.

5.8 Cauchy Distribution

Uses the skeleton for the uniform distribution inside N-dimensional shapes.

1. Generate two empty PSRNs, with a positive sign, an integer part of 0, and an empty fractional part. Call the PSRNs $p1$ and $p2$.
2. Set S to $base$, where $base$ is the base of digits to be stored by the PSRNs (such as 2 for binary or 10 for decimal). Then set $c1$ and $c2$ each to 0. Then set d to 1.
3. Multiply $c1$ and $c2$ each by $base$ and add a digit chosen uniformly at random to that coordinate.
4. If $((c1+1)^2 + (c2+1)^2) < S^2$, then do the following:
 1. Transfer $c1$'s least significant digits to $p1$'s fractional part, and transfer $c2$'s least significant digits to $p2$'s fractional part. The variable d tells how many digits to transfer to each PSRN this way. (For example, if $base$ is 10, d is 3, and $c1$ is 342, set $p1$'s fractional part to [3, 4, 2].)
 2. Run the **UniformDivision** algorithm (described in the article on PSRNs) on $p1$ and $p2$, in that order, then set the resulting PSRN's sign to positive or negative with equal probability, then return that PSRN.
5. If $(c1^2 + c2^2) > S^2$, then go to step 2.
6. Multiply S by $base$, then add 1 to d , then go to step 3.

5.9 Exponential Distribution with Rate $\ln(x)$

The following new algorithm generates a partially-sampled random number that follows the exponential distribution with rate $\ln(x)$. This is useful for generating a base- x logarithm of a uniform(0,1) random number. Here, x is a rational number that's greater than 1. In the algorithm, let b be $\text{floor}(\ln(x)/\ln(2))$.

1. (Samples the integer part of the random number.) Generate a random number that expresses the number of failed trials before the first success, where each trial succeeds with probability $1-1/x$. Set k to that random number. (This is also known as a "geometric random number", but this terminology is avoided because it has conflicting meanings in academic works. If x is a power of 2, this step can be implemented by generating blocks of b unbiased random bits until a **non-zero** block of bits is generated this way, then setting k to the number of **all-zero** blocks of bits generated this way.)
2. (The rest of the algorithm samples the fractional part.) Generate a uniform (0, 1) random number, call it f .
3. Create a μ input coin that does the following: "**Sample from the number f** (e.g., call **SampleGeometricBag** on f if f is implemented as a uniform PSRN), then run the **algorithm for $\ln(2)$** (described in "Bernoulli Factory Algorithms"). If both calls return 1, return 1. Otherwise, return 0." (This simulates the probability $\lambda = f \cdot \ln(2)$.) If x is not a power of 2, also create a ν input coin that does the following: "**Sample from the number f** , then run the **algorithm for $\ln(1 + y/z)$** (described in

- "Bernoulli Factory Algorithms") with $y/z = (x-2^b)/2^b$. If both calls return 1, return 1. Otherwise, return 0."
4. Run the **algorithm for exp(-λ)** (described in "Bernoulli Factory Algorithms") b times, using the μ input coin. If x is not a power of 2, run the same algorithm once, using the ν input coin. If all these calls return 1, accept f . If f is accepted this way, set f 's integer part to k , then optionally fill f with uniform random digits as necessary to give its fractional part the desired number of digits (similarly to **FillGeometricBag**), then return f .
 5. If f was not accepted by the previous step, go to step 2.

Note: A *bounded exponential* random number with rate $\ln(x)$ and bounded by m has a similar algorithm to this one. Step 1 is changed to read as follows: "Set k to a bounded-geometric($1-1/x$, m) random number (Bringmann and Friedrich 2013)⁽¹²⁾, or more simply, the lesser of m or the number of failed trials before the first success, where each trial succeeds with probability $1-1/x$. (If x is a power of 2, this can be implemented by generating blocks of b unbiased random bits until a **non-zero** block of bits or m blocks of bits are generated this way, whichever comes first, then setting k to the number of **all-zero** blocks of bits generated this way.) If k is m , return m (note that this m is a constant, not a uniform PSRN; if the algorithm would otherwise return a uniform PSRN, it can return something else in order to distinguish this constant from a uniform PSRN)." Additionally, instead of generating a uniform(0,1) random number in step 2, a uniform(0, μ) random number can be generated instead, such as a uniform PSRN generated via **RandUniformFromReal**, to implement an exponential distribution bounded by $m+\mu$ (where μ is a real number in the interval (0, 1)).

The following generator for the **rate ln(2)** is a special case of the previous algorithm and is useful for generating a base-2 logarithm of a uniform(0,1) random number. Unlike the similar algorithm of Ahrens and Dieter (1972)⁽¹³⁾, this one doesn't require a table of probability values.

1. (Samples the integer part of the random number. This will be geometrically distributed with parameter 1/2.) Generate unbiased random bits until a zero is generated this way. Set k to the number of ones generated this way.
2. (The rest of the algorithm samples the fractional part.) Generate a uniform (0, 1) random number, call it f .
3. Create an input coin that does the following: "**Sample from the number f** (e.g., call **SampleGeometricBag** on f if f is implemented as a uniform PSRN), then run the **algorithm for ln(2)** (described in "Bernoulli Factory Algorithms"). If both calls return 1, return 1. Otherwise, return 0." (This simulates the probability $\lambda = f \cdot \ln(2)$.)
4. Run the **algorithm for exp(-λ)** (described in "Bernoulli Factory Algorithms"), using the input coin from the previous step. If the call returns 1, accept f . If f is accepted this way, set f 's integer part to k , then optionally fill f with uniform random digits as necessary to give its fractional part the desired number of digits (similarly to **FillGeometricBag**), then return f .
5. If f was not accepted by the previous step, go to step 2.

5.10 Lindley Distribution and Lindley-Like Mixtures

A random number that follows the Lindley distribution (Lindley 1958)⁽¹⁴⁾ with parameter θ (a rational number greater than 0) can be generated as follows:

1. With probability $w = \theta/(1+\theta)$, generate an exponential random number with a rate of

- θ via **ExpRand** or **ExpRand2** (described in my article on PSRNs) and return that number.
- Otherwise, generate two exponential random numbers with a rate of θ via **ExpRand** or **ExpRand2**, then generate their sum by applying the **UniformAdd** algorithm, then return that sum.

For the Garima distribution (Shanker 2016)⁽¹⁵⁾, $w = (1+\theta)/(2+\theta)$.

For the i-Garima distribution (Singh and Das 2020)⁽¹⁶⁾, $w = (2+\theta)/(3+\theta)$.

For the mixture-of-weighted-exponential-and-weighted-gamma distribution in (Iqbal and Iqbal 2020)⁽¹⁷⁾, two exponential random numbers (rather than one) are generated in step 1, and three (rather than two) are generated in step 2.

6 Requests and Open Questions

- We would like to see new implementations of the following:
 - Algorithms that implement **InShape** for specific closed curves, specific closed surfaces, and specific signed distance functions. Recall that **InShape** determines whether a box lies inside, outside, or partly inside or outside a given curve or surface.
 - Descriptions of new arbitrary-precision algorithms that use the skeleton given in the section "Building an Arbitrary-Precision Sampler".
- The appendix contains implementation notes for **InShape**, which determines whether a box is outside or partially or fully inside a shape. However, practical implementations of **InShape** will generally only be able to evaluate a shape pointwise. What are necessary and/or sufficient conditions that allow an implementation to correctly classify a box just by evaluating the shape pointwise?
- Take a polynomial $f(\lambda)$ of even degree n of the form $\text{choose}(n, n/2) * \lambda^{n/2} * (1-\lambda)^{n/2} * k$, where k is greater than 1 (thus all f 's Bernstein coefficients are 0 except for the middle one, which equals k). Suppose $f(1/2)$ lies in the interval $(0, 1)$. If we do the degree elevation, described in the appendix, enough times (at least r times), then f 's Bernstein coefficients will all lie in $[0, 1]$. The question is: how many degree elevations are enough? A [MathOverflow answer](#) showed that r is at least $m = (n/f(1/2)^2)/(1-f(1/2)^2)$, but is it true that $\text{floor}(m)+1$ elevations are enough?

7 Notes

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8 Appendix

8.1 Ratio of Uniforms

The Cauchy sampler given earlier demonstrates the *ratio-of-uniforms* technique for sampling a distribution (Kinderman and Monahan 1977)⁽¹⁸⁾. It involves transforming the distribution's density function (PDF) into a compact shape. The ratio-of-uniforms method appears here in the appendix, particularly since it can involve calculating upper and lower bounds of transcendental functions which, while it's possible to achieve in rational arithmetic (Daumas et al., 2007)⁽¹⁹⁾, is less elegant than, say, the normal distribution sampler by Karney (2014)⁽²⁰⁾, which doesn't require calculating logarithms or other transcendental functions.

This algorithm works for any univariate (one-variable) distribution as long as—

- for all x , $PDF(x) < \infty$ and $PDF(x)*x^2 < \infty$, where PDF is the distribution's PDF or a function proportional to the PDF,
- PDF is continuous almost everywhere, and
- either—
 - the distribution's ratio-of-uniforms shape (the transformed PDF) is covered entirely by the rectangle $[0, \text{ceil}(d1)] \times [0, \text{ceil}(d2)]$, where $d1$ is not less than the highest value of $x*\text{sqrt}(PDF(x))$ anywhere, and $d2$ is not less than the highest value of $\text{sqrt}(PDF(x))$ anywhere, or
 - half of that shape is covered this way and the shape is symmetric about the v -axis.

The algorithm follows.

1. Generate two empty PSRNs, with a positive sign, an integer part of 0, and an empty fractional part. Call the PSRNs $p1$ and $p2$.
2. Set S to $base$, where $base$ is the base of digits to be stored by the PSRNs (such as 2 for binary or 10 for decimal). Then set $c1$ to an integer in the interval $[0, d1)$, chosen uniformly at random, then set $c2$ to an integer in $[0, d2)$, chosen uniformly at random, then set d to 1.
3. Multiply $c1$ and $c2$ each by $base$ and add a digit chosen uniformly at random to that coordinate.
4. Run an **InShape** function that determines whether the transformed PDF is covered by the current box. In principle, this is the case when $z \leq 0$ everywhere in the box, where u lies in $[c1/S, (c1+1)/S]$, v lies in $[c2/S, (c2+1)/S]$, and z is $v^2 - PDF(u/v)$. **InShape** returns *YES* if the box is fully inside the transformed PDF, *NO* if the box is fully outside it, and *MAYBE* in any other case, or if evaluating z fails for a given box (e.g., because $\ln(0)$ would be calculated or v is 0). See the next section for implementation notes.
5. If **InShape** as described in step 4 returns *YES*, then do the following:
 1. Transfer $c1$'s least significant digits to $p1$'s fractional part, and transfer $c2$'s least significant digits to $p2$'s fractional part. The variable d tells how many digits to transfer to each PSRN this way. Then set $p1$'s integer part to $\text{floor}(c1/base^d)$ and $p2$'s integer part to $\text{floor}(c2/base^d)$. (For example, if $base$ is 10, d is 3, and $c1$ is 7342, set $p1$'s fractional part to [3, 4, 2] and $p1$'s integer part to 7.)
 2. Run the **UniformDivision** algorithm (described in the article on PSRNs) on $p1$ and $p2$, in that order.
 3. If the transformed PDF is symmetric about the v -axis, set the resulting PSRN's sign to positive or negative with equal probability. Otherwise, set the PSRN's sign to positive.
 4. Return the PSRN.
6. If **InShape** as described in step 4 returns *NO*, then go to step 2.
7. Multiply S by $base$, then add 1 to d , then go to step 3.

Examples:

1. For the normal distribution, PDF is proportional to $\exp(-x^2/2)$, so that z after a logarithmic transformation (see next section) becomes $4*\ln(v) + (u/v)^2$, and since the distribution's ratio-of-uniforms shape is symmetric about the v -axis, the return value's sign is positive or negative with equal probability.
2. For the standard lognormal distribution ([Gibrat's distribution](#)), $PDF(x)$ is proportional to $\exp(-(\ln(x))^2/2)/x$, so that z after a logarithmic

transformation becomes $2*\ln(v)-(-\ln(u/v)^2/2 - \ln(u/v))$, and the returned PSRN has a positive sign.

3. For the gamma distribution with shape parameter $a > 1$, $PDF(x)$ is proportional to $x^{a-1}*\exp(-x)$, so that z after a logarithmic transformation becomes $2*\ln(v)-(a-1)*\ln(u/v)-(u/v)$, or 0 if u or v is 0, and the returned PSRN has a positive sign.

8.2 Implementation Notes for Box/Shape Intersection

The "**Uniform Distribution Inside N-Dimensional Shapes**" algorithm uses a function called **InShape** to determine whether an axis-aligned box is either outside a shape, fully inside the shape, or partially inside the shape. The following are notes that will aid in developing a robust implementation of **InShape** for a particular shape, especially because the boxes being tested can be arbitrarily small.

1. **InShape**, as well as the divisions of the coordinates by S , should be implemented using rational arithmetic. Instead of dividing those coordinates this way, an implementation can pass S as a separate parameter to **InShape**.
2. If the shape is convex, and the point $(0, 0, \dots, 0)$ is on or inside that shape, **InShape** can return—
 - *YES* if all the box's corners are in the shape;
 - *NO* if none of the box's corners are in the shape and if the shape's boundary does not intersect with the box's boundary; and
 - *MAYBE* in any other case, or if the function is unsure.

In the case of two-dimensional shapes, the shape's corners are $(c1/S, c2/S)$, $((c1+1)/S, c2/S)$, $(c1, (c2+1)/S)$, and $((c1+1)/S, (c2+1)/S)$. However, checking for box/shape intersections this way is non-trivial to implement robustly, especially if interval arithmetic is not used.

3. If the shape is given as an inequality of the form $f(t1, \dots, tN) \leq 0$, **InShape** should use rational interval arithmetic (such as the one given in (Daumas et al., 2007)⁽¹⁹⁾), where the two bounds of each interval are rational numbers with arbitrary-precision numerators and denominators. Then, **InShape** should build one interval for each dimension of the box and evaluate f using those intervals⁽²¹⁾ with an accuracy that increases as S increases. Then, **InShape** can return—
 - *YES* if the interval result of f has an upper bound less than or equal to 0;
 - *NO* if the interval result of f has a lower bound greater than 0; and
 - *MAYBE* in any other case.

For example, if f is $(t1^2+t2^2-1)$, which describes a quarter disk, **InShape** should build two intervals, namely $t1 = [c1/S, (c1+1)/S]$ and $t2 = [c2/S, (c2+1)/S]$, and evaluate $f(t1, t2)$ using interval arithmetic.

One thing to point out, though: If f calls the $\exp(x)$ function where x can potentially have a high absolute value, say 10000 or higher, the \exp function can run a very long time in order to calculate proper bounds for the result, since the number of digits in $\exp(x)$ grows linearly with x . In this case, it may help to transform the inequality to its logarithmic version. For example, by applying $\ln(\cdot)$ to each side of the inequality $y^2 \leq \exp(-(x/y)^2/2)$, the inequality becomes $2*\ln(y) \leq -(x/y)^2/2$ and thus becomes $2*\ln(y) + (x/y)^2/2 \leq 0$ and thus becomes $4*\ln(y) + (x/y)^2 \leq 0$.

4. If the shape is such that every axis-aligned line segment that begins in one face of the hypercube and ends in another face crosses the shape at most once, ignoring the segment's endpoints (an example is an axis-aligned quarter of a circular disk where the disk's center is (0, 0)), then **InShape** can return—
 - *YES* if all the box's corners are in the shape;
 - *NO* if none of the box's corners are in the shape; and
 - *MAYBE* in any other case, or if the function is unsure.

If **InShape** uses rational interval arithmetic, it can build an interval per dimension *per corner*, evaluate the shape for each corner individually and with an accuracy that increases as S increases, and treat a corner as inside or outside the shape only if the result of the evaluation clearly indicates that. Using the example of a quarter disk, **InShape** can build eight intervals, namely an x - and y -interval for each of the four corners; evaluate $(x^2 + y^2 - 1)$ for each corner; and return *YES* only if all four results have upper bounds less than or equal to 0, *NO* only if all four results have lower bounds greater than 0, and *MAYBE* in any other case.

5. If **InShape** expresses a shape in the form of a [*signed distance function*](#), namely a function that describes the closest distance from any point in space to the shape's boundary, it can return—
 - *YES* if the signed distance (or an upper bound of such distance) at each of the box's corners, after dividing their coordinates by S , is less than or equal to $-\sigma$ (where σ is an upper bound for $\sqrt{N}/(S \cdot 2)$, such as $1/S$);
 - *NO* if the signed distance (or a lower bound of such distance) at each of the box's corners is greater than σ ; and
 - *MAYBE* in any other case, or if the function is unsure.
6. **InShape** implementations can also involve a shape's *implicit curve* or *algebraic curve* equation (for closed curves), its *implicit surface* equation (for closed surfaces), or its *signed distance field* (a quantized version of a signed distance function).
7. An **InShape** function can implement a set operation (such as a union, intersection, or difference) of several simpler shapes, each with its own **InShape** function. The final result depends on the shape operation (such as union or intersection) as well as the result returned by each component for a given box. The following are examples of set operations:
 - For unions, the final result is *YES* if any component returns *YES*; *NO* if all components return *NO*; and *MAYBE* otherwise.
 - For intersections, the final result is *YES* if all components return *YES*; *NO* if any component returns *NO*; and *MAYBE* otherwise.
 - For differences between two shapes, the final result is *YES* if the first shape returns *YES* and the second returns *NO*; *NO* if the first shape returns *NO* or if both shapes return *YES*; and *MAYBE* otherwise.
 - For the exclusive OR of two shapes, the final result is *YES* if one shape returns *YES* and the other returns *NO*; *NO* if both shapes return *NO* or both return *YES*; and *MAYBE* otherwise.

8.3 SymPy Code for Piecewise Linear Factory Functions

```
def bernstein_n(func, x, n, pt=None):
    # Bernstein operator.
```

```

# Create a polynomial that approximates func, which in turn uses
# the symbol x. The polynomial's degree is n and is evaluated
# at the point pt (or at x if not given).
if pt==None: pt=x
ret=0
v=[binomial(n,j) for j in range(n//2+1)]
for i in range(0, n+1):
    oldret=ret
    bino=v[i] if i<len(v) else v[n-i]
    ret+=func.subs(x,S(i)/n)*bino*pt**i*(1-pt)**(n-i)
    if pt!=x and ret==oldret and ret>0: break
return ret

def inflec(y,eps=S(2)/10,mult=2):
    # Calculate the inflection point (x) given y, eps, and mult.
    # The formula is not found in the paper by Thomas and
    # Blanchet 2012, but in
    # the supplemental source code uploaded by
    # A.C. Thomas.
    po=5 # Degree of y-to-x polynomial curve
    eps=S(eps)
    mult=S(mult)
    x=-((y-(1-eps))/eps)**po/mult + y/mult
    return x

def xfunc(y,sym,eps=S(2)/10,mult=2):
    # Calculate Bernstein "control polygon" given y,
    # eps, and mult.
    return Min(sym*y/inflec(y,eps,mult),y)

def calc_linear_func(eps=S(5)/10, mult=1, count=10):
    # Calculates the degrees and Y parameters
    # of a sequence of polynomials that converge
    # from above to min(x*mult, 1-eps).
    # eps must be in the interval (0, 1).
    # Default is 10 polynomials.
    polys=[]
    eps=S(eps)
    mult=S(mult)
    count=S(count)
    bs=20
    ypt=1-(eps/4)
    x=symbols('x')
    tfunc=Min(x*mult,1-eps)
    tfn=tfunc.subs(x,(1-eps)/mult).n()
    xpt=xfunc(ypt,x,eps=eps,mult=mult)
    bits=5
    #olddb=None
    i=0
    lastbxn = 1
    diffs=[]
    while i<count:
        bx=bernstein_n(xpt,x,bits,(1-eps)/mult)
        bxn=bx.n()
        if bxn > tfn and bxn < lastbxn:
            # Dominates target function
            #if oldbx!=None:
            #    diffs.append(bx)
            #    diffs.append(olddb-bx)
            #olddb=bx
            oldxpt=xpt
            lastbxn = bxn

```

```

    polys.append([bits,ypt])
    print("    [%d,%s]," % (bits,ypt))
    # Find y2 such that y2 < ypt and
    # bernstein_n(oldxpt,x,bits,inflec(y2, ...)) >= y2,
    # so that next Bernstein expansion will go
    # underneath the previous one
    while True:
        ypt=(ypt-(1-eps))/4
        xpt=inflec(ypt,eps=eps,mult=mult).n()
        bxs=bernstein_n(oldxpt,x,bits,xpt).n()
        if bxs>=ypt.n():
            break
        xpt=xfunc(ypt,x,eps=eps,mult=mult)
        bits+=20
        i+=1
    else:
        bits=int(bits*200/100)
    return polys

```

```
calc_linear_func(count=8)
```

8.4 Derivation of My Algorithm for $\min(\lambda, 1/2)$

The following explains how the algorithm is derived.

The function $\min(\lambda, 1/2)$ can be rewritten as $A + B$ where—

- $A = (1/2) * \lambda$, and
- $B = (1/2) * ((1 - \sqrt{1 - 4 * \lambda * (1 - \lambda)}) / 2) = (1/2) * \sum_{k=1, 2, \dots} g(k) * h_k(\lambda)$,

revealing that the function is a [convex combination](#), and B is itself a convex combination where—

- $g(k) = (2^k * k) / ((2^k - 1) * 2^{2^k})$, and
- $h_k(\lambda) = (4 * \lambda * (1 - \lambda))^k / 2 = (\lambda * (1 - \lambda))^k * 4^k / 2$

(see also Wästlund (1999)⁽²²⁾; Dale et al. (2015)⁽²³⁾). The right-hand side of h , which is the polynomial built in step 3 of the algorithm, is a polynomial of degree $k * 2$ with Bernstein coefficients—

- $z = (4^{v/2}) / \text{choose}(v * 2, v)$ at $v = k$, and
- 0 elsewhere.

Unfortunately, z is generally greater than 1, so that the polynomial can't be simulated, as is, using the Bernoulli factory algorithm for [polynomials in Bernstein form](#). Fortunately, the polynomial's degree can be elevated to bring the Bernstein coefficients to 1 or less (for degree elevation and other algorithms, see (Tsai and Farouki 2001)⁽²⁴⁾). Moreover, due to the special form of the Bernstein coefficients in this case, the degree elevation process can be greatly simplified. Given an even degree d as well as z (as defined above), the degree elevation is as follows:

1. Set r to $\text{floor}(d/3) + 1$. (This starting value is because when this routine finishes, r/d appears to converge to $1/3$ as d gets large, for the polynomial in question.) Let c be $\text{choose}(d, d/2)$.
2. Create a list of $d + r + 1$ Bernstein coefficients, all zeros.
3. For each integer i in the interval $[0, d + r]$:
 - If $d/2$ is in the interval $[\max(0, i - r), \min(d, i)]$, set the i^{th} Bernstein coefficient

- (starting at 0) to $z^*c*\text{choose}(r,i-d/2)^* / \text{choose}(d+r, i)$.
4. If all the Bernstein coefficients are 1 or less, return them. Otherwise, add $d/2$ to r and go to step 2.

8.5 Algorithm 2 for Non-Negative Factories

Algorithm 2. Say we have an *oracle* that produces random real numbers. The goal is now to produce non-negative random numbers that average to the mean of $f(X)$, where X is a number produced by the oracle. This is possible whenever the mean of $f(X)$ is not less than δ , where δ is a known number greater than 0. The algorithm to do so follows (see Lee et al. 2014)⁽²⁵⁾:

1. Let m be an upper bound of the maximum value of $\text{abs}(f(\mu))$ anywhere. Create a ν input coin that does the following: "Generate a random number from the oracle, call it x . With probability $\text{abs}(f(x))/m$, return a number that is 1 if $f(x) < 0$ and 0 otherwise. Otherwise, repeat this process."
2. Use one of the [linear Bernoulli factories](#) to simulate $2*\nu$ (2 times the ν coin's probability of heads), using the ν input coin, with $\epsilon = \delta/m$. If the factory returns 1, return 0. Otherwise, generate a random number from the oracle, call it ξ , and return $\text{abs}(f(\xi))$.

Example: An example from Lee et al. (2014)⁽²⁵⁾. Say the oracle produces uniform random numbers in $[0, 3*\pi]$, and let $f(\mu) = \sin(\mu)$. Then the mean of $f(X)$ is $2/(3*\pi)$, which is greater than 0 and found in SymPy by `sympy.stats.E(sin(sympy.stats.Uniform('U', 0, 3*pi)))`, so the algorithm can produce non-negative random numbers that average to that mean.

Notes:

1. Averaging to the mean of $f(X)$ (that is, $\mathbf{E}[f(X)]$ where $\mathbf{E}[\cdot]$ means expected or average value) is not the same as averaging to $f(\mu)$ where μ is the mean of the oracle's numbers (that is, $f(\mathbf{E}[X])$). For example, if X is 0 or 1 with equal probability, and $f(\lambda) = \exp(-\lambda)$, then $\mathbf{E}[f(X)] = \exp(0) + (\exp(-1) - \exp(0))*(1/2)$, and $f(\mathbf{E}[X]) = f(1/2) = \exp(-1/2)$.
2. (Lee et al. 2014, Corollary 4)⁽²⁵⁾: If $f(\mu)$ is known to return only values in the interval $[a, c]$, the mean of $f(X)$ is not less than δ , $\delta > b$, and δ and b are known numbers, then Algorithm 2 can be modified as follows:
 - Use $f(\mu) = f(\mu) - b$, and use $\delta = \delta - b$.
 - m is taken as $\max(b-a, c-b)$.
 - When Algorithm 2 finishes, add b to its return value.

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