# Notes 9: Numerics

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### Introduction

This document is the ninth of a set of notes, this document focusing on numerical questions, random number generation, and linear algebra. The notes are not meant to be particularly complete in terms of useful functions (Google and LLMs can now provide that quite well), but rather to introduce the language and consider key programming concepts in the context of Julia.

Given that, the document heavily relies on demos, with interpretation in some cases left to the reader.

### Random number generation

#### The seed

As usual, it's best to set the random number seed.

```
using Random
Random.seed!(1234);
                      # Seed number 1234
println(rand(3))
[0.32597672886359486, 0.5490511363155669, 0.21858665481883066]
println(rand(5))
[0.8942454282009883, 0.35311164439921205, 0.39425536741585077, 0.9531246272848422, 0.7955469475347194
                     # Re-seed with same number - will give the same sequence of random numbers
Random.seed! (1234);
println(rand(2))
[0.32597672886359486, 0.5490511363155669]
println(rand(6))
rand has a variety of methods.
rand(1:10, 3)
3-element Vector{Int64}:
 5
 8
 6
rand(['a','b','c'], 5)
5-element Vector{Char}:
 'c': ASCII/Unicode U+0063 (category L1: Letter, lowercase)
 'a': ASCII/Unicode U+0061 (category Ll: Letter, lowercase)
 'a': ASCII/Unicode U+0061 (category L1: Letter, lowercase)
 'b': ASCII/Unicode U+0062 (category L1: Letter, lowercase)
```

#### Generators

The manual discusses the available generators.

'c': ASCII/Unicode U+0063 (category L1: Letter, lowercase)

The default RNG is Xoshiro256++, but there are details related to random number streams when working in parallel that I have not absorbed yet.

```
Random.default_rng();
Random.seed!(1234)
println(rand(3))
```

```
[0.32597672886359486, 0.5490511363155669, 0.21858665481883066]
```

```
rng = Random.Xoshiro(1234)
println(rand(rng, 3))
```

[0.32597672886359486, 0.5490511363155669, 0.21858665481883066]

The default used to be the Mersenne Twister (still the default in R and formerly the default in Python/numpy).

```
rng = Random.MersenneTwister(1234);
println(rand(rng, 3))
```

[0.5908446386657102, 0.7667970365022592, 0.5662374165061859]

#### Distributions

0.1894867660152757 0.8329474331914342

As with distributions in SciPy in Python, we first define the distribution of interest and then carry out distributional operations with it.

```
using Distributions
beta_dist = Beta(2, 5)
beta_samples = rand(beta_dist, 10)
10-element Vector{Float64}:
 0.4542940512832609
 0.27004146470644036
 0.3870672630509891
 0.21116529046475746
 0.5041249425689261
 0.2784321400052276
 0.6829933994934323
 0.1295297689806119
 0.43795698919772685
 0.39395847238977155
pdf.(beta_dist, beta_samples[1:5])
5-element Vector{Float64}:
 1.2086291289041646
 2.300088115277408
 1.6389245754521693
 2.4529451430316644
 0.9144258514892248
logpdf.(beta_dist, beta_samples[1:5])
5-element Vector{Float64}:
```

```
0.4940402800647119
  0.8972894018105007
 -0.08945889536877027
Floating point issues
Integer and floating point types
64 bit integers can represent -2^{63}\dots 2^{63}. 32 bit integers can represent -2^{31}\dots 2^{31}.
Similarly for 16 and 128 bit integers.
Values outside that range overflow.
xi::Int64 = 2^62
4611686018427387904
xi::Int64 = 2^70 # Overflows
xi::Int64 = 2^63 # Just overflows.
-9223372036854775808
yi::Int128 = 2^63 # Hmmm.
-9223372036854775808
yi::Int128 = Int128(2)^63
9223372036854775808
Int64(yi)
LoadError: InexactError: trunc(Int64, 9223372036854775808)
InexactError: trunc(Int64, 9223372036854775808)
Stacktrace:
 [1] throw_inexacterror(f::Symbol, ::Type{Int64}, val::Int128)
   @ Core ./boot.jl:634
 [2] checked_trunc_sint
   @ ./boot.jl:656 [inlined]
 [3] toInt64
   @ ./boot.jl:705 [inlined]
 [4] Int64(x::Int128)
   @ Core ./boot.jl:784
 [5] top-level scope
   @ In[19]:1
```

9223372036854775807

xi::Int64 = 2^63 - 1 # Shouldn't this overflow when calculating 2^63?

```
x = parse(BigInt, "123456789012345678901234567890")
```

1234567890123456789012345678901234567890

x+1

1234567890123456789012345678901234567891

There are 16, 32, and 64 bit floating point numbers, as well as BigFloats.

More a bit later.

### Floating point precision

Let's consider how much precision we have with real-valued numbers because of limited floating point precision.

For 64-bit floats, 53 bits are used for precision, which translates to approximately 16 digits of accuracy in base 10, regardless of the magnitude of the number. How many digits of accuracy do we have with 32-bit and 16-bit floats?

"0.33325195312500000000"

BigFloat(1/3) # Hmmm.

 $\tt 0.333333333333333314829616256247390992939472198486328125$ 

```
BigFloat(1) / BigFloat(3)
```

```
BigFloat("0.3")
```

```
BigFloat("0.3", precision=500)
```

Computation with BigFloats will be slow, so you wouldn't want to do matrix operations with a matrix of them.

### Floating point details

With Float64, any number is stored as a base 2 number of the form:

$$(-1)^S \times 1.d \times 2^{e-1023} = (-1)^S \times 1.d_1d_2 \dots d_{52} \times 2^{e-1023}$$

where the computer uses base 2, b=2, (so  $d_i \in \{0,1\}$ ) because base-2 arithmetic is faster than base-10 arithmetic. The leading 1 normalizes the number; i.e., ensures there is a unique representation for a given computer number. This avoids representing any number in multiple ways, e.g., either  $1=1.0\times 2^0=0.1\times 2^1=0.01\times 2^2$ . For a double, we have 8 bytes=64 bits. Consider our representation as (S,d,e) where S is the sign. The leading 1 is the *hidden bit* and doesn't need to be stored because it is always present. In general e is represented using 11 bits ( $2^{11}=2048$ ), and the subtraction takes the place of having a sign bit for the exponent. (Note that in our discussion we'll just think of e in terms of its base 10 representation, although it is of course represented in base 2.) This leaves p=52=64-1-11 bits for d.

53 bits of precision in base 10 is about 16 digits. And 11 bits for the magnitude corresponds with when over/underflow occur. For 32-bit floats, we have both less precision and more easily over/underflow.

#### Overflow

Integer numbers can be represented exactly by Float64 up to  $2^{53}$ .

```
function pri(x)
    @sprintf("%.20i", x)
end

pri(2.0^52)

"4503599627370496"

pri(2.0^52 + 1)

"4503599627370497"

pri(2.0^53)

"9007199254740992"

pri(2.0^53 + 1)

"9007199254740992"

pri(2.0^53 + 2)

"9007199254740994"

pri(2.0^63)
```

```
pri(2.0^70) # No overflow here unlike Int64.
"1180591620717411303424"
pri(12345678123456781234.0) # Not exact.
"12345678123456780288"
Float64 overflow is not until \sim 2^{1023} \approx 10^{308}.
function prf(x)
  @sprintf("%.20f", x)
end
prf(2.0<sup>1022</sup>)
prf(2.0<sup>1023</sup>)
"8988465674311579538646525953945123668089884894711532863671504057886633790275048156635423866120376801
prf(2.0<sup>1024</sup>)
"Inf"
But Int64s that big do overflow.
pri(10<sup>307</sup>)
"000000000000000000000"
pri(10^310)
"000000000000000000000"
Implications for comparisons and calculations
prf(1-2/3)
"0.33333333333333337034"
prf(1/3)
"0.3333333333333331483"
prf(2/3-1/3)
"0.3333333333333331483"
prf(0.3 - 0.2)
"0.099999999999997780"
```

```
prf(0.1)

"0.10000000000000555"

0.3 - 0.2 == 0.1
```

#### false

Here's an example of catastrophic cancellation:

```
prf(123456781234.56)
prf(123456781234.00)
prf(123456781234.56 - 123456781234.00)
```

#### "0.55999755859375000000"

And here the precision is that of the larger magnitude number:

```
prf(1.0 + 1e-8)
"1.00000000999999993923"
```

```
prf(1.0 + 1e-17)
1.0 + 1e-17 == 1.0
```

true

### Avoid multiplying/dividing many numbers

```
using Distributions
normd = Normal(0, 1)
samples = rand(normd, 1000);
prod(pdf.(normd, samples)) # Log-likelihood/log-density

0.0
sum(logpdf.(normd, samples))
```

-1446.1968973565636

We see that a likelihood with only as many as 1000 terms underflows. (In other cases it could overflow.)

### Linear algebra

We'll see that linear algebra operations heavily exploit Julia's multiple dispatch system, calling the most appropriate method for different kinds of input matrices.

#### Arithmetic

```
A = rand(3, 3);
B = rand(3, 3);
```

```
A + B # Element-wise
A .+ B # Element-wise
A * B # Matrix multiplication
A .* B # Element-wise multiplication
```

### Solving systems of equations / inversion

```
using LinearAlgebra
A = rand(3, 3);
AtA = A'A;
            # A' * A
b = rand(3);
AtA \ b
         # Solve system of equations (AtA^{-1} b)
3-element Vector{Float64}:
 -5839.367752001243
   237.62992428763377
  1388.1477149501256
inv(AtA) * b # Not as efficient
3-element Vector{Float64}:
 -5839.367752001244
   237.62992428763377
  1388.1477149501254
eigvals(AtA)
3-element Vector{Float64}:
 1.0406188999502348e-5
 0.20302101952602597
 2.8527433608929575
det(AtA)
              # Be careful of over/underflow!
6.026919865436598e-6
logdet(AtA)
-12.019274479480647
tr(A)
1.3719983709419976
chol = cholesky(AtA)
Cholesky{Float64, Matrix{Float64}}
U factor:
```

```
3×3 UpperTriangular{Float64, Matrix{Float64}}:
 0.290417 1.25712
                       1.00661
           0.605775 -0.103239
                      0.0139545
chol \ b
                         # Automatically exploits triangularity.
3-element Vector{Float64}:
 -5839.3677520089905
   237.62992428794826
  1388.1477149519676
chol.U \ (chol.L \ b)
                         # Manual equivalent solution.
3-element Vector{Float64}:
 -5839.3677520089905
   237.62992428794826
  1388.1477149519676
typeof(chol)
                         # Special type of object
Cholesky{Float64, Matrix{Float64}}
typeof(chol.U)
                         # Special kind of matrix.
UpperTriangular{Float64, Matrix{Float64}}
This works in Julia but is having problems via Quarto, so I'll just paste in some timings.
using BenchmarkTools
n = 5000
A = rand(n, n);
AtA = A'A;
b = rand(n);
                       # Solve system of equations (AtA^{-1} b) via Gaussian elimination (LU)
Obtime AtA \ b;
```

In principle, the Cholesky approach should involve  $n^3/6$  calculations and the Gaussian elimination  $n^3/3$ , but we don't see a two-fold difference here in practice.

Obtime cholesky(AtA) \ b; # Best if matrix is positive definite.

Obtime chol = cholesky(AtA); chol.U \ (chol.L \ b); # As good but verbose.

### Spectral (eigen) decomposition

# 633.046 ms (6 allocations: 190.81 MiB)
@btime inv(AtA) \* b; # Not as efficient
# 2.538 s (8 allocations: 193.25 MiB)

# 546.957 ms (5 allocations: 190.77 MiB)

# 520.021 ms (3 allocations: 190.73 MiB)

```
\Gamma = eigvecs(AtA);
 = eigvals(AtA);
result = \Gamma * diagm() * \Gamma';
result[1:3,1:3]
3×3 Matrix{Float64}:
 0.0843421 0.365089 0.292337
 0.365089
            1.94732
                       1.20289
            1.20289
 0.292337
                       1.02412
AtA[1:3,1:3]
3×3 Matrix{Float64}:
 0.0843421 0.365089
                       0.292337
 0.365089
            1.94732
                       1.20289
 0.292337
            1.20289
                       1.02412
result == AtA
false
isapprox(result, AtA)
true
         AtA # \approx TAB
result
true
```

### Don't forget floating point issues

```
n=500
A = rand(n, n);
det(A'A)

Inf
logdet(A'A)
```

#### 1369.773541109337

Here's a positive definite matrix (mathematically) that has all real, positive eigenvalues. On a computer, it's not positive definite, and therefore not invertible/full rank.

```
xs = 0:99
# Compute distance matrix.
dists = abs.(xs .- xs') # Using broadcasting with ' (an "outer" operation).
# Create correlation matrix.
corr_matrix = exp.(-(dists/10).^2)
```

```
# Compute eigenvalues and get last 20
eigvals(corr_matrix)[1:20]
20-element Vector{Float64}:
 -5.050107193681261e-15
 -3.975288045003424e-15
 -3.0343039148117795e-15
 -2.3309913036821875e-15
 -1.5093549661154687e-15
 -1.3511621460923302e-15
 -1.2848549714671506e-15
 -1.0472866233731733e-15
 -8.644575073124488e-16
 -8.364136467422076e-16
 -7.989065965702987e-16
 -6.829697931404996e-16
 -6.091008178138255e-16
 -6.030845253775599e-16
 -5.700968077566563e-16
 -5.37906121536584e-16
 -4.615826595950015e-16
 -4.4365869197860045e-16
 -4.0894133652379844e-16
 -3.580464402267892e-16
chol = cholesky(corr_matrix);
LoadError: PosDefException: matrix is not positive definite; Cholesky factorization failed.
```

PosDefException: matrix is not positive definite; Cholesky factorization failed.

#### Stacktrace:

- [1] checkpositivedefinite
  - @ /system/linux/julia-1.10.4/share/julia/stdlib/v1.10/LinearAlgebra/src/factorization.jl:67 [inlinearAlgebra/src/factorization.jl:67 [inlinearAlgebra/src/factor
- [2] cholesky!(A::Hermitian{Float64, Matrix{Float64}}, ::NoPivot; check::Bool)
  - @ LinearAlgebra /system/linux/julia-1.10.4/share/julia/stdlib/v1.10/LinearAlgebra/src/cholesky.jl:
- [3] cholesky!
  - @ /system/linux/julia-1.10.4/share/julia/stdlib/v1.10/LinearAlgebra/src/cholesky.jl:267 [inlined]
- [4] cholesky!(A::Matrix{Float64}, ::NoPivot; check::Bool)
  - @ LinearAlgebra /system/linux/julia-1.10.4/share/julia/stdlib/v1.10/LinearAlgebra/src/cholesky.jl:
- [5] cholesky! (repeats 2 times)
  - @ /system/linux/julia-1.10.4/share/julia/stdlib/v1.10/LinearAlgebra/src/cholesky.jl:295 [inlined]
- [6] cholesky
  - @ /system/linux/julia-1.10.4/share/julia/stdlib/v1.10/LinearAlgebra/src/cholesky.jl:401 [inlined]
- [7] cholesky(A::Matrix{Float64})
  - @ LinearAlgebra /system/linux/julia-1.10.4/share/julia/stdlib/v1.10/LinearAlgebra/src/cholesky.jl:
- [8] top-level scope
  - @ In[73]:1

One should be able to use an approximate pivoted Cholesky that sets diagonal elements to zero corresponding to the rank deficiency. I'm having trouble seeing how to do that.

This does seem to work to solve the system of equations. We'd have to investigate to know what Julia is doing behind the scenes, but it's probably using a pivoted LU decomposition.

```
b = rand(100);
out = corr_matrix \ b;
```

#### **Smart factorization**

```
n = 500
A = rand(n, n);

typeof(factorize(A'A))

Cholesky{Float64, Matrix{Float64}}

typeof(factorize(A))  # I wouldn't have guessed that A is invertible!

LU{Float64, Matrix{Float64}, Vector{Int64}}

typeof(factorize(A[:,1:10]))
```

```
QRPivoted{Float64, Matrix{Float64}, Vector{Float64}, Vector{Int64}}
```

The orthogonal matrices generated by certain factorizations can most efficiently be worked with by having them treated as "matrix-backed, function-based linear operators".

```
QRresult = qr(A[:,1:10]);
typeof(QRresult.Q)
```

```
LinearAlgebra.QRCompactWYQ{Float64, Matrix{Float64}}, Matrix{Float64}}
```

```
QRresult.Q * rand(10);
```

#### Use known structure!

```
using BenchmarkTools

n = 5000
A = rand(n, n);
AtA = A'A;
b = rand(n);
chol = cholesky(AtA);
typeof(chol.U)
@btime chol.U \ b;
```

```
5.996 ms (3 allocations: 39.12 KiB)
```

```
U_dense = Matrix(chol.U);
Obtime U_dense \ b;
  14.779 ms (2 allocations: 39.11 KiB)
Obtime logdet(chol.U);
  104.481 s (2 allocations: 32 bytes)
Obtime logdet(U_dense);
  647.519 ms (5 allocations: 190.77 MiB)
Obtime sum(log.(diag(chol.U)));
  100.264 s (7 allocations: 78.27 KiB)
using SparseArrays
n = 5000;
A = Matrix(1.0I, n, n);
A[1,3] = 7.7;
A[5,9] = 2.3;
sA = sparse(A);
sizeof(A)
200000000
sizeof(sA) # Ok, so that's not helpful given the pointers involved.
40
# sA. TAB # What are the components of `sA`?
sizeof(sA.colptr) + sizeof(sA.nzval) + sizeof(sA.rowval)
120040
b = rand(n);
@btime A * b;
  9.429 ms (2 allocations: 39.11 KiB)
Obtime sA * b;
  20.277 s (2 allocations: 39.11 KiB)
```