Rare-event simulation: Code demo 1 Patrick Laub February 12, 2020

1 Simulating (pseudo-) random numbers in Python

1.1 Setup

```
[1]: # numpy is the 'Numerical Python' package
     import numpy as np
     # Numpy's methods for pseudorandom number generation
     import numpy.random as rnd
     # scipy is the 'Scientific Python' package
     # We'll use this to get the gamma function
     from scipy.special import gamma
[2]: # Print out the versions of software I'm running
     import sys
     print("Python version:", sys.version)
     print("Numpy version:", np.__version__)
    Python version: 3.7.6 | packaged by conda-forge | (default, Jan 7 2020,
    21:00:34) [MSC v.1916 64 bit (AMD64)]
    Numpy version: 1.17.4
[3]: # Reminder that we need a relatively new version of numpy to make
     # use of the latest pseudorandom number generation algorithms.
     if int(np.__version__.split('.')[1]) < 17:</pre>
         raise RuntimeError("Need Numpy version >= 1.17")
```

1.2 Random numbers, seeds, accessing docs

Create a random number generator and call it rng.

```
[4]: rng = rnd.default_rng()
```

What kind of things can rng do? Let's look at the methods available to it using dir

```
[5]: print(dir(rng))
```

So it can simulate from a bunch of common distributions. That's nice. Let's try to generate a simulate standard uniform random variable:

```
[6]: print(rng.uniform())
```

0.78658288842436

```
[7]: rng.uniform()
```

[7]: 0.29599899160720367

Run that cell a few times, and you'll see it get different numbers.

Sometimes nice to have *same* random numbers. To do that, we set the 'seed' to be any fixed number.

```
[8]: print("First run")
  print(18*"-")

rng = rnd.default_rng(seed=1)
  print(rng.uniform())
  print(rng.uniform())
  print(rng.uniform())
```

```
First run
```

- 0.5118216247002567
- 0.9504636963259353
- 0.14415961271963373

```
[9]: print("Second run")
print(18*"-")
```

```
rng = rnd.default_rng(seed=1)
print(rng.uniform())
print(rng.uniform())
print(rng.uniform())
```

Second run

- 0.5118216247002567
- 0.9504636963259353
- 0.14415961271963373

Using rng = rnd.default_rng() and rng.uniform() is a pretty modern (≥ July 26, 2019). The old way was to run np.random.uniform() and the RNG was hidden away. Proper nerds (or those with trouble sleeping) can take a look at why the numpy developers moved away from this.

```
[10]: np.random.seed(1)
    print(np.random.uniform())

    np.random.seed(1)
    print(np.random.uniform())
```

- 0.417022004702574
- 0.417022004702574

Say we want to generate some other uniform variable, like Unif(a, b) with p.d.f.

$$f(x) = \begin{cases} \frac{1}{b-a} & \text{if } a \le x < b \\ 0 & \text{otherwise.} \end{cases}$$

Let's use help to look at the documentation for the uniform method to see if this is builtin.

[11]: help(rng.uniform)

Help on built-in function uniform:

uniform(...) method of numpy.random.generator.Generator instance uniform(low=0.0, high=1.0, size=None)

Draw samples from a uniform distribution.

Samples are uniformly distributed over the half-open interval ``[low, high)`` (includes low, but excludes high). In other words, any value within the given interval is equally likely to be drawn by `uniform`.

Parameters

low : float or array_like of floats, optional

Lower boundary of the output interval. All values generated will be greater than or equal to low. The default value is 0.

high : float or array_like of floats

Upper boundary of the output interval. All values generated will be less than high. The default value is 1.0.

size : int or tuple of ints, optional

Output shape. If the given shape is, e.g., ``(m, n, k)``, then
``m * n * k`` samples are drawn. If size is ``None`` (default),
a single value is returned if ``low`` and ``high`` are both scalars.
Otherwise, ``np.broadcast(low, high).size`` samples are drawn.

Returns

out : ndarray or scalar

Drawn samples from the parameterized uniform distribution.

See Also

integers : Discrete uniform distribution, yielding integers.

random : Floats uniformly distributed over ``[0, 1)``.

random : Alias for `random`.

Notes

The probability density function of the uniform distribution is

.. math::
$$p(x) = \frac{1}{b - a}$$

anywhere within the interval ``[a, b)``, and zero elsewhere.

When ``high`` == ``low``, values of ``low`` will be returned. If ``high`` < ``low``, the results are officially undefined and may eventually raise an error, i.e. do not rely on this function to behave when passed arguments satisfying that inequality condition.

Examples

Draw samples from the distribution:

```
>>> s = np.random.default_rng().uniform(-1,0,1000)
```

All values are within the given interval:

```
>>> np.all(s >= -1)
True
>>> np.all(s < 0)
True</pre>
```

Display the histogram of the samples, along with the probability density function:

```
>>> import matplotlib.pyplot as plt
>>> count, bins, ignored = plt.hist(s, 15, density=True)
>>> plt.plot(bins, np.ones_like(bins), linewidth=2, color='r')
>>> plt.show()
```

So, let's simulate from Unif(-10, 10).

```
[12]: rng.uniform(low=-10, high=10)
```

[12]: 8.972988942744877

The uniform method has the optional arguments low, high, and size. If we simply use them in this order, we don't need to write low= and high=, but can directly write:

```
[13]: rng.uniform(-10, 10)
```

[13]: -3.763370959790291

You can use you own judgement on whether to include the names of the arguments or omit them.

We can simulate many uniforms at the same time and the result will be an array filled with i.i.d. variables.

```
[14]: rng.uniform(-10, 10, size=5)
```

```
[14]: array([-1.53347102, 6.55405188, -1.81601727, 0.99187375, -9.44881774])
```

Let's simulate a large number of uniforms, and compare some of empirical quantities against the theoretical quantities.

```
[15]: # The number of random variables to simulate.
R = 10^6
print(R)
```

12

```
[16]: # The number of random variables to simulate.
R = 10**6
print(R)
```

1000000

```
[17]: # Simulate a bunch of i.i.d. uniform variables
uniforms = rng.uniform(-10, 10, R)
```

Sample mean: -0.0004428776080335241

Theoretical mean: 0.0

[19]: # Print the sample variance of these observations and the theoretical → variance for this distribution print("Sample variance:", uniforms.var()) print("Theoretical variance:", (10 - -10)**2 / 12)

Sample variance: 33.34011458681296

Theoretical variance: 33.333333333333333

It certainly looks like we correctly simulated from the desired/target distribution; well, at the very least, we simulated from some distribution which has the same mean and variance as our target distribution.

Let's try to simulate from a slightly more complication distribution, the exponential distribution. I'll define $\mathsf{Exp}(\lambda)$ to have p.d.f.

$$f(x) = \begin{cases} \lambda e^{-\lambda x} & \text{if } x > 0\\ 0 & \text{otherwise.} \end{cases}$$

There is a rng.exponential method, though we should call help on it to find out how to give it a specific rate λ .

[20]: help(rng.exponential)

Help on built-in function exponential:

exponential(...) method of numpy.random.generator.Generator instance
 exponential(scale=1.0, size=None)

Draw samples from an exponential distribution.

Its probability density function is

```
.. math:: f(x; \frac{1}{\beta}) = \frac{1}{\beta} \exp(-\frac{x}{\beta}),
```

for ``x > 0`` and 0 elsewhere. :math:`\beta` is the scale parameter, which is the inverse of the rate parameter :math:`\lambda = 1/\beta`. The rate parameter is an alternative, widely used parameterization of the exponential distribution [3]_.

The exponential distribution is a continuous analogue of the

geometric distribution. It describes many common situations, such as the size of raindrops measured over many rainstorms [1]_, or the time between page requests to Wikipedia [2].

Parameters

scale : float or array_like of floats
 The scale parameter, :math:`\beta = 1/\lambda`. Must be
 non-negative.

size : int or tuple of ints, optional
 Output shape. If the given shape is, e.g., ``(m, n, k)``, then
 ``m * n * k`` samples are drawn. If size is ``None`` (default),
 a single value is returned if ``scale`` is a scalar. Otherwise,
 ``np.array(scale).size`` samples are drawn.

Returns

out : ndarray or scalar

Drawn samples from the parameterized exponential distribution.

References

- .. [1] Peyton Z. Peebles Jr., "Probability, Random Variables and Random Signal Principles", 4th ed, 2001, p. 57.
- .. [2] Wikipedia, "Poisson process",
 https://en.wikipedia.org/wiki/Poisson_process

So, this one only takes a scale parameter, so we'll have to set the scale to be $1/\lambda$. Let's try simulate from the $\mathsf{Exp}(5)$ distribution.

[21]: $\lambda = 5$ rng.exponential(scale=1/ λ)

[21]: 0.3757960365292656

Fun fact: To get λ as a variable name, just type \Lambda then hit TAB. Try \sigma and some others.

Fun fact 2: Can use **sigma** as a variable name, but not **Lambda** as it has special significance in Python (for "lambda functions" = "anonymous functions").

Let's simulate a bunch of i.i.d. exponentials and check that their sample mean matches the theoretical value of 0.2 = 1/5.

[22]: exponentials = rng.exponential($1/\lambda$, R)

```
[23]: print("Sample mean:", exponentials.mean())
print("Theoretical mean:", 1/λ)
```

Sample mean: 0.20022323810176906

Theoretical mean: 0.2

The **exponential** function just specifies a **scale** argument, but sometimes the distribution we want to simulate from won't offer us this. There's an easy workaround though, just simulate from the default/unscaled distribution and multiply all the variables by our scale value. E.g. to simulate from the same exponential distribution this way:

```
[24]: exponentials = (1/λ) * rng.exponential(size=R)
print("Sample mean:", exponentials.mean())
print("Theoretical mean:", 1/λ)
```

Sample mean: 0.2000826239295432

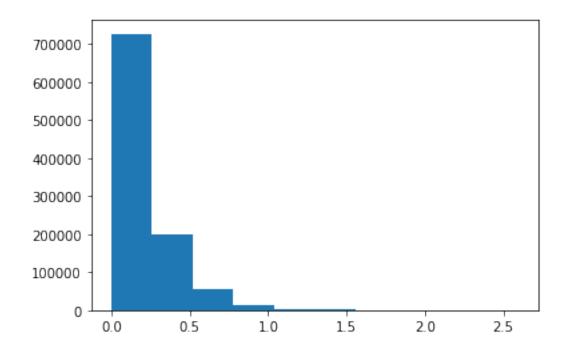
Theoretical mean: 0.2

1.3 Basic plotting

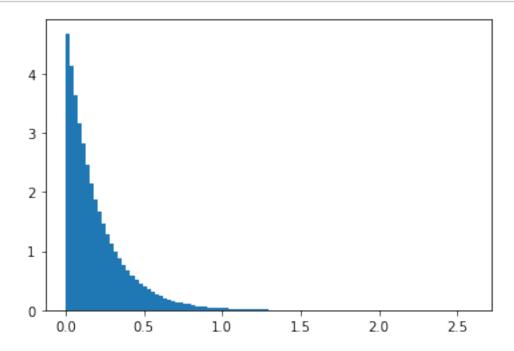
```
[25]: # Import the plotting library matplotlib
import matplotlib.pyplot as plt
```

```
[26]: plt.hist(exponentials)
```

```
[26]: (array([7.26228e+05, 1.98673e+05, 5.44270e+04, 1.48950e+04, 4.22500e+03, 1.12700e+03, 3.18000e+02, 6.80000e+01, 3.10000e+01, 8.00000e+00]), array([1.65267759e-07, 2.59168929e-01, 5.18337693e-01, 7.77506457e-01, 1.03667522e+00, 1.29584398e+00, 1.55501275e+00, 1.81418151e+00, 2.07335028e+00, 2.33251904e+00, 2.59168780e+00]), <a list of 10 Patch objects>)
```

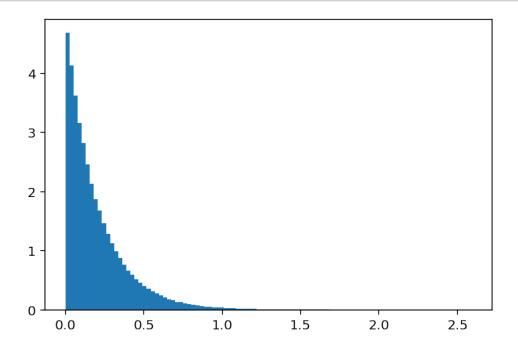


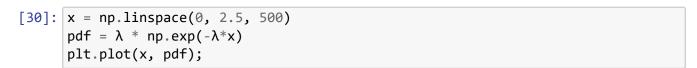
[27]: plt.hist(exponentials, bins=100, density=True);

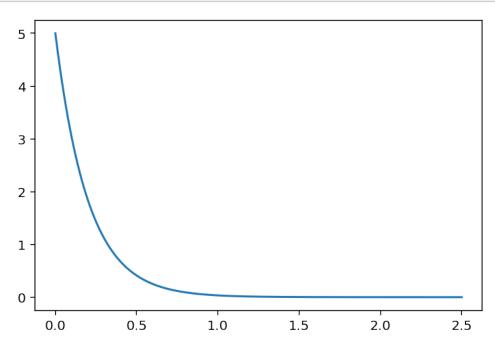


[28]: %config InlineBackend.figure_format = 'retina'

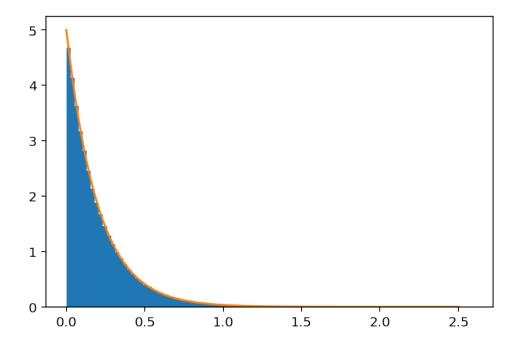
[29]: plt.hist(exponentials, bins=100, density=True);







```
[31]: plt.hist(exponentials, bins=100, density=True);
plt.plot(x, pdf); # Or plt.plot(x, pdf, 'r');
```



Try again with $R=10^5~{\sf Gamma}(2,3)$ random variables, where our definition of ${\sf Gamma}(r,m)$ has the p.d.f.

$$f(x) = \begin{cases} \frac{x^{r-1}e^{-\frac{x}{m}}}{\Gamma(r)m^r} & \text{if } x > 0\\ 0 & \text{otherwise.} \end{cases}$$

```
[32]: R = 10**5
r = 2
m = 3
gammas = rng.gamma(r, m, R)
```

```
[33]: x = np.linspace(0, 40, 500)
pdf = (x**(r-1) * np.exp(-x/m)) / (gamma(r) * m**r)

plt.hist(gammas, bins=100, density=True)
plt.plot(x, pdf);
```

```
[34]: R = 10**4
      # Mean vector and covariance matrix
      n = 2
      \mu = (1, 2)
      \sigma 2 = 2
      \rho = 0.8
      \Sigma = \sigma 2 * ((1-\rho) * np.eye(n) + \rho * np.ones(n))
      # Simulating the index value at T
      rng = rnd.default_rng()
      normals = rng.multivariate_normal(μ, Σ, size=R)
[35]: import pandas as pd
      df = pd.DataFrame(normals, columns=["x", "y"])
      df
[35]:
                    Χ
            -0.228046 1.382743
      1
            0.351114 1.827377
      2
            1.752873 3.419312
      3
            4.276756 5.734045
            2.827853 4.172706
      4
      9995 -2.081927 -0.750787
```

```
      9996
      2.892714
      2.892638

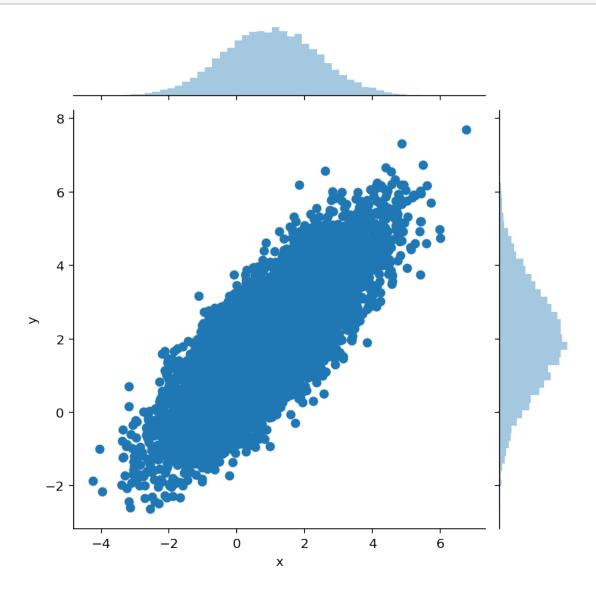
      9997
      1.445032
      2.041335

      9998
      1.598780
      2.429420

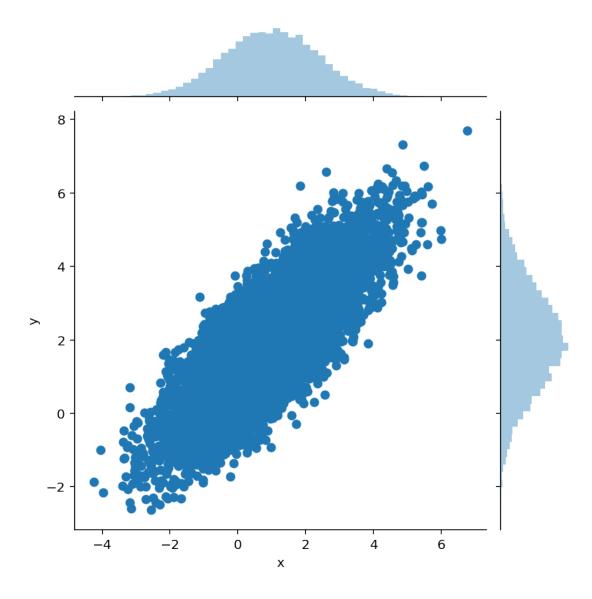
      9999
      2.232821
      2.760091
```

[10000 rows x 2 columns]

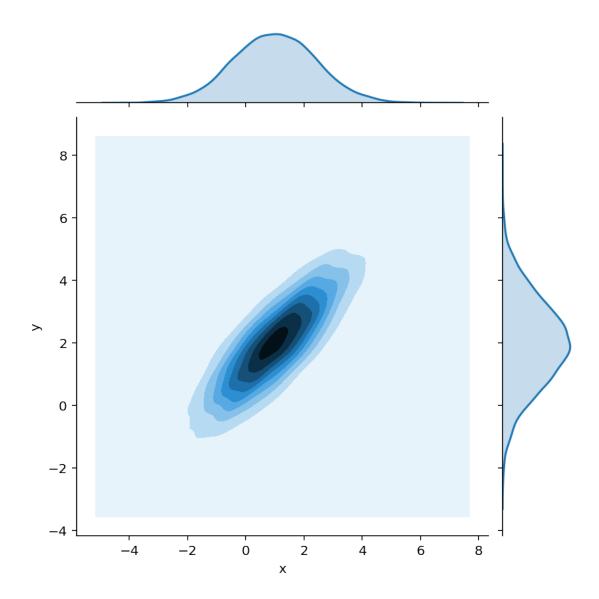
```
[36]: import seaborn as sns
sns.jointplot(x="x", y="y", data=df);
```



```
[37]: sns.jointplot(x="x", y="y", data=df);
```



```
[38]: sns.jointplot(x="x", y="y", data=df, kind="kde");
```



1.4 Crude Monte Carlo

Say that an insurer has n=10 claims each month, and each claim size is $X_i \overset{\text{i.i.d.}}{\sim} \mathsf{Pareto}(\alpha = \frac{3}{2})$. The reinsurer will cover the excess of $S_n = \sum_{i=1}^n X_i$ over the threshold $\gamma = 10^2$.

What is the probability of reinsurer having to payout?

[39]: n = 10 $\alpha = 3/2$ $\gamma = 10**2$ R = 10**6

```
[40]: %%time

rng = rnd.default_rng(1)
numPayouts = 0

for r in range(R):
    S_n = rng.pareto(α, size=n).sum()
    if S_n > γ:
        numPayouts += 1

print("Probability of reinsurer paying out:", numPayouts / R)
```

Probability of reinsurer paying out: 0.013021 Wall time: 7.31 s

This last cell is quite clunky and slow; please never write code like that. The preferred way is the vectorised code below. Firstly note that if we give a list to the **size=** parameter (actually we use an immutable list called a 'tuple') it returns a matrix of i.i.d. Paretos:

Now if we generate all the Pareto variable we need at once, everything is more efficient.

```
[42]: %%time

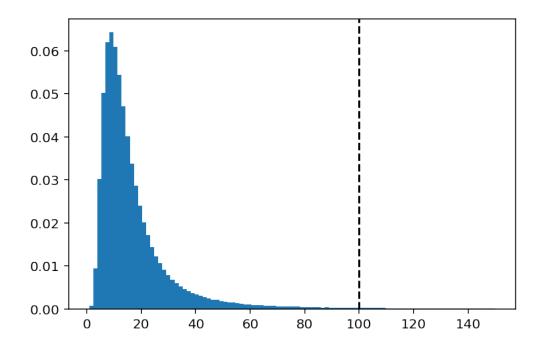
rng = rnd.default_rng(1)
losses = rng.pareto(α, size=(R,n)).sum(axis=1)

ests = losses > γ
ellHat = ests.mean()

print("Probability of reinsurer paying out:", ellHat)
```

Probability of reinsurer paying out: 0.013021 Wall time: 360 ms

```
[43]: plt.hist(losses[losses < 1.5*γ], bins=100, density=True)
plt.axvline(γ, color="black", linestyle="dashed");</pre>
```



```
[44]: sigmaHat = ests.std()
widthCI = 1.96 * sigmaHat / np.sqrt(R)
CIs = (ellHat - widthCI, ellHat + widthCI)
print(f"Probability of reinsurer paying out: {ellHat} (+- {widthCI:f}))")
```

Probability of reinsurer paying out: 0.013021 (+- 0.000222))

Bonus question: Can compare to series expansion by ISFA's Quang Huy NGUYEN & Christian ROBERT.

How much is the reinsurer paying out on average?

```
[45]: rng = rnd.default_rng(1)
losses = rng.pareto(α, size=(R,n)).sum(axis=1)
payouts = np.maximum(losses - γ, Θ)
np.mean(payouts)
```

[45]: 2.5494052977235837

Note: We can't calculate confidence intervals here using the normal approach. We're in the unlucky case that our variables have infinite variance and the CLT doesn't apply.

What is the expected payout for the of reinsurer conditioned on the event of a payout?

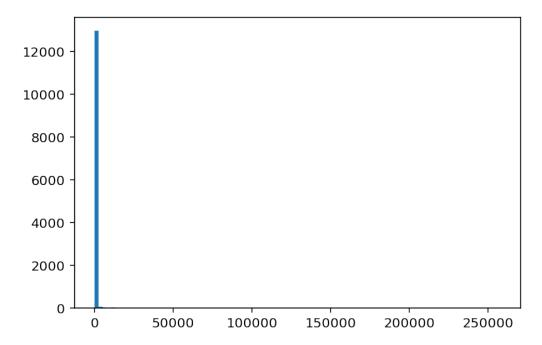
```
[46]: rng = rnd.default_rng(1)
  losses = rng.pareto(α, size=(R,n)).sum(axis=1)
  bigLosses = losses[losses > γ]
```

```
payouts = bigLosses - γ

np.mean(payouts)
```

[46]: 195.79182072986583

We had 13021 reinsurer payouts out of 10⁶ simulations.



What about the 99.9% Value-at-Risk for the reinsurer?

```
[48]: rng = rnd.default_rng(1)

losses = rng.pareto(α, size=(R,n)).sum(axis=1)
payouts = np.maximum(losses - γ, 0)

np.quantile(payouts, 0.999)
```

[48]: 377.0792235808656

Let's consider a financial example. Say that X_i is the future stock price for company i at expiry time T. We assume the Black-Scholes model, so $X_i \sim \mathsf{Lognormal}(\mu_i, \sigma^2)$, and assume a constant correlation ρ between each pair of stocks.

Let's imagine we have a simple index which tracks n of these stocks, so at time T is will have the value

$$S_T = \sum_{i=1}^n X_i.$$

What would be the value of a call option on this index, i.e., what is

$$\mathbb{E}[e^{-rT}(S_T - K)_+]?$$

(Let's ignore the \mathbb{Q} measure here.)

Set n = 2, r = 0.05, T = 1, $\mu_i = \frac{i}{10}$, $\sigma^2 = \frac{1}{10}$, $\rho = 0.25$, K = 2.

```
[49]: # Problem constants
      n = 2
      r = 0.05
      T = 1
      K = 3
      \rho = -0.5
      \sigma 2 = 1/10
      R = 10**6
      # Mean vector and covariance matrix
      \mu = np.arange(1, n+1) / 10
      \Sigma = \sigma 2 * ((1-\rho) * np.eye(n) + \rho * np.ones(n))
      # Simulating the index value at T
      rng = rnd.default_rng()
      normals = rng.multivariate_normal(\mu, \Sigma, size=R)
      Xs = np.exp(normals)
      Ss = Xs.sum(axis=1)
      # Calculating the MC estimate and CIs
      ests = np.exp(-r*T) * np.maximum(Ss - K, 0)
      ellHat = ests.mean()
      sigmaHat = ests.std()
      widthCI = 1.96 * sigmaHat / np.sqrt(R)
      print(f"Option value: {ellHat} (+- {widthCI:f}))")
      ests = (Ss > K)
      ellHat = ests.mean()
      sigmaHat = ests.std()
      widthCI = 1.96 * sigmaHat / np.sqrt(R)
      print(f"Probability of payout: {ellHat} (+- {widthCI:f}))")
```

Option value: 0.024134250001376582 (+- 0.000208))
Probability of payout: 0.094762 (+- 0.000574))

[50]: Ss.mean()

[50]: 2.4458187502697877