

FN6815_Lect_08_Pricing_NI

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FN6815 Numerical Methods for Financial Instrument Pricing

Lecture 8: Basic Instrument Pricing and Pricing with Numerical Integration

- Dr. Yang Ye
- Email: yy@runchee.com
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1. Introduction

In this lecture, we'll explore the fundamentals of financial instrument pricing and how numerical integration can be applied in this context.

As we have learnt earlier, numerical integration is a technique used to approximate the definite integral of a function. It's extensively used in financial instrument pricing, particularly for exotic options. These options often have customized payoff structures with European-style, i.e. do not have pre-maturity optionality.

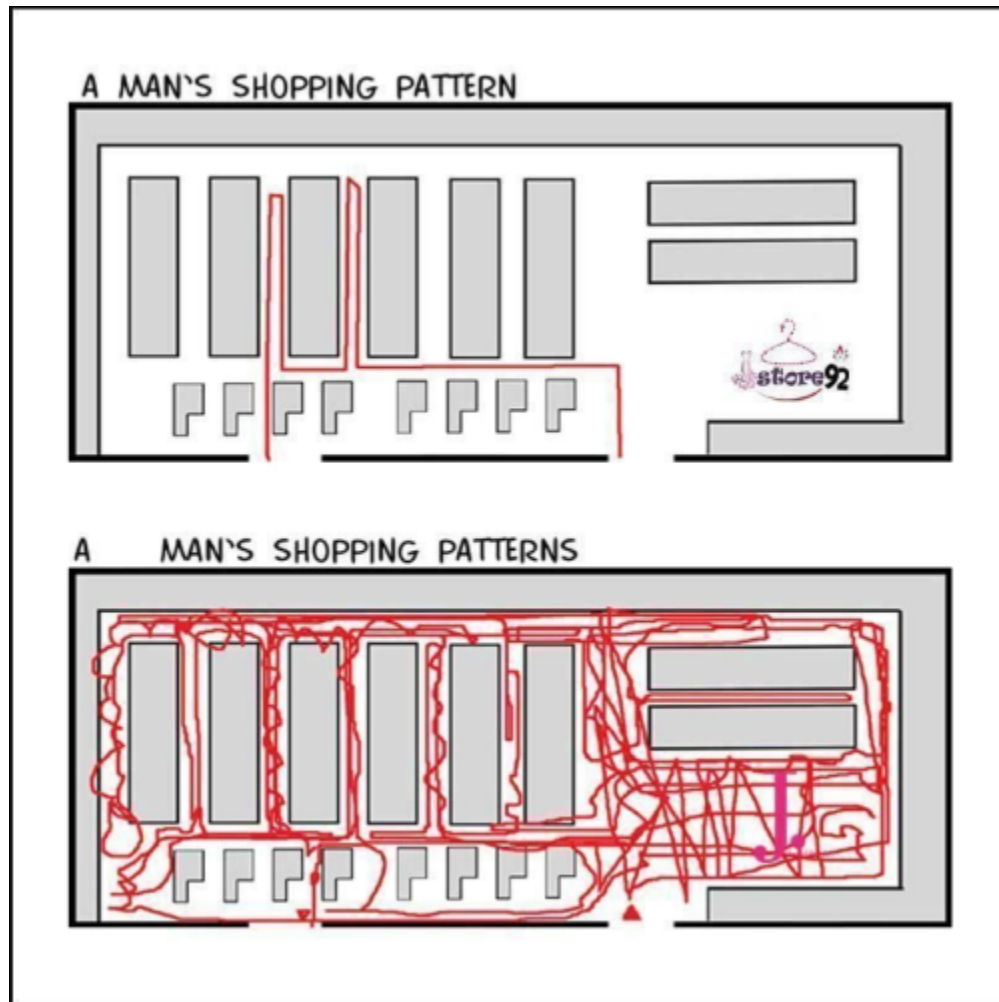
For instance, consider an exotic option with a payoff that depends on the square of the underlying asset's price.

1.2 Pricing Methods Overview

There are several methods available for pricing financial instruments:

- **Analytic/Semi-analytical methods (10%):** These include closed-form solutions, numerical integration, and tree-based methods. They are used when the pricing formula can be derived analytically or approximated semi-analytically.
- **Monte Carlo methods (60%):** These methods use random sampling to estimate the price of financial instruments, especially when the dimensionality of the problem is high.
- **Partial Differential Equations (PDEs) and Finite Difference methods (30%):** These methods solve the PDEs that arise in financial instrument pricing, typically by discretizing the problem and solving the resulting system of equations.

The percentages in parentheses indicate the approximate frequency of use of each method in practice.



1.3 Choosing the Right Pricing Method

Each pricing method has its advantages and disadvantages:

- **Analytic methods:** These have well-defined calculation steps but offer limited support for path-dependent options. Monte Carlo (MC) and Partial Differential Equations (PDE) methods can handle a wider variety of options.
- **Monte Carlo methods:** These are usually “more the better” algorithms but could have hit limit of computation resource easily, meaning their accuracy improves with more iterations. However, they require high-performance hardware to run efficiently.
- **PDE methods:** These run faster than MC methods but are more complex to set up. They can also suffer from the “curse of dimensionality”.

Curse of Dimensionality

- More complex models lead to a higher number of Stochastic Differential Equations (SDEs). If we have n SDEs, each one adds an extra dimension to the PDE.
- The computational cost of Monte Carlo methods is linear in the number of iterations N , but more assets require more iterations, leading to a complexity cost of $O(N \times n)$.

- The computational cost of PDE methods is exponential in $O(N * M^n)$, where N is the number of time intervals, M is the number of asset price intervals, and n is the number of assets.
- Typically, PDE methods are used for up to $n = 3$ dimensions. For example, an exchange rate product might have SDEs for the local interest rate, foreign interest rate, and exchange rate.

2. Closed-Form Solutions

Closed-form solutions for option pricing are typically the fastest. For instance, Vanilla European options can be priced using the Black-Scholes model.

Background

The Black-Scholes option pricing models for European-style options on a non-dividend paying stock are as follows:

For a call option:

$$c = S_0 N(d_1) - K e^{-rT} N(d_2)$$

For a put option:

$$p = K e^{-rT} N(-d_2) - S_0 N(-d_1)$$

$$d_1 = \frac{\log(S_0/K) + (r + \sigma^2/2)T}{\sigma\sqrt{T}}$$

$$d_2 = d_1 - \sigma\sqrt{T}$$

where:

- $N(x)$ is the cumulative density function of a standardized normal distribution.
- S_0 is the current price of the stock.
- K is the strike price of the option.
- T is the time to maturity of the option.
- r is the (continuously-compounded) risk-free rate of return.
- σ is the stock price volatility.
- $\log()$ is the natural logarithm.

```
[1]: import matplotlib.pyplot as plt
import numpy as np
import numpy.typing as npt
import scipy.stats as sps
```

```
[2]: S_0 = 100
strike = 100
ttm = 2
rf = 0.03
sigma = 0.3

def bs_call(S, strike, ttm, rf, sigma):
```

```

    d1 = (np.log(S / strike) + (rf + sigma**2 / 2.0) * ttm) / (sigma * np.
↪sqrt(ttm))
    d2 = d1 - sigma * np.sqrt(ttm)
    return S * sps.norm.cdf(d1) - strike * np.exp(-rf * ttm) * sps.norm.cdf(d2)

def bs_put(S, strike, ttm, rf, sigma):
    d1 = (np.log(S / strike) + (rf + sigma**2 / 2.0) * ttm) / (sigma * np.
↪sqrt(ttm))
    d2 = d1 - sigma * np.sqrt(ttm)
    return sps.norm.cdf(-d2) * strike * np.exp(-rf * ttm) - sps.norm.cdf(-d1) *
↪S

c = bs_call(S_0, strike, ttm, rf, sigma)
p = bs_put(S_0, strike, ttm, rf, sigma)
print("call:", c, "put:", p, c / p)

# Runs with different input don't affect performance
%timeit -n100 -r100 bs_call(100,90,2,.05,0.4)
%timeit -n100 -r100 bs_call(20,90,4.,.05,0.4)

```

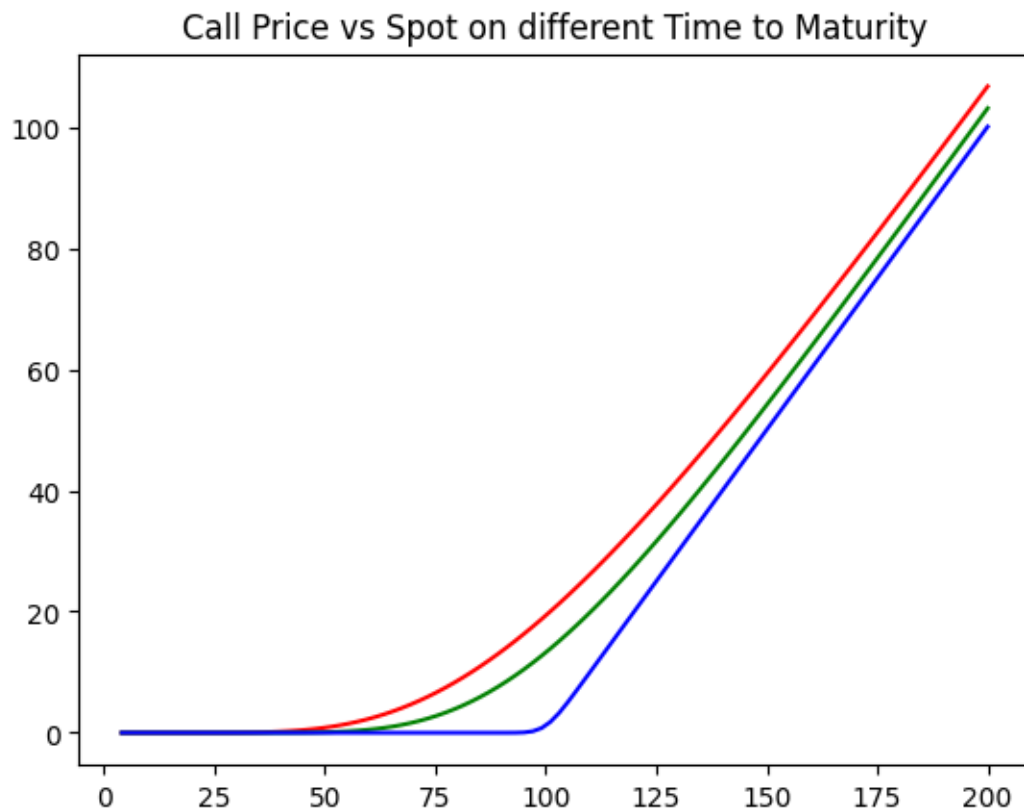
call: 19.38254929806086 put: 13.559002656485731 1.429496681217149
 59.2 μ s \pm 3.7 μ s per loop (mean \pm std. dev. of 100 runs, 100 loops each)
 55.8 μ s \pm 2.47 μ s per loop (mean \pm std. dev. of 100 runs, 100 loops each)

Rapid Option Pricing with Closed-Form Equations Closed-form equations enable rapid calculation of option prices for a variety of asset prices or strike prices. These equations provide a direct method to compute the price of an option given the necessary parameters, making them highly efficient for pricing a large number of options simultaneously.

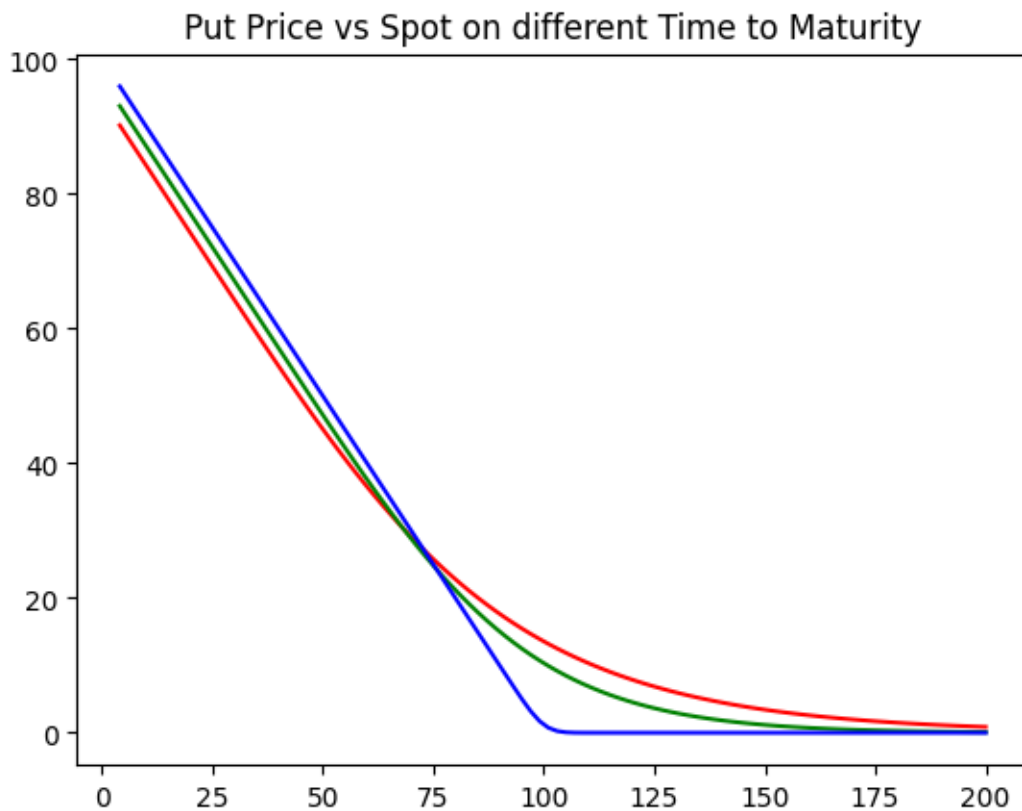
```

[3]: xs = np.linspace(4, 200, 100)
    cs = bs_call(xs, 100, 2, rf, sigma)
    plt.plot(xs, cs, "r-")
    cs = bs_call(xs, 100, 1, rf, sigma)
    plt.plot(xs, cs, "g-")
    cs = bs_call(xs, 100, 0.01, rf, sigma)
    plt.plot(xs, cs, "b-")
    plt.title("Call Price vs Spot on different Time to Maturity")
    plt.show()

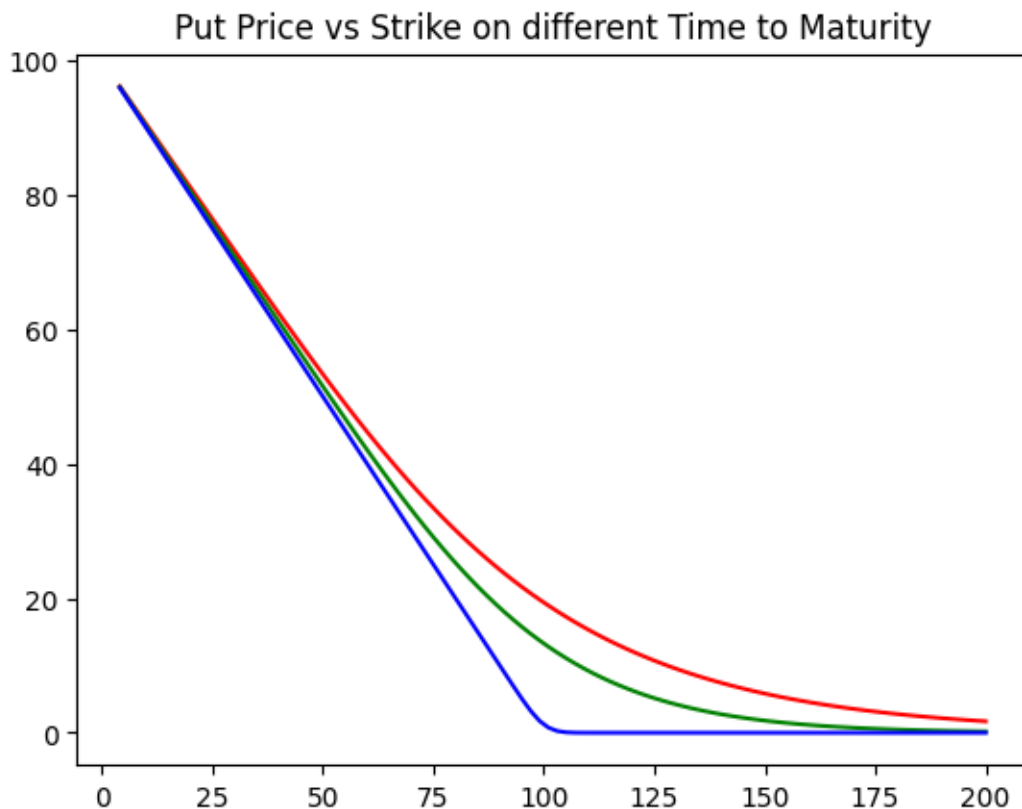
```



```
[4]: xs = np.linspace(4, 200, 100)
cs = bs_put(xs, 100, 2, rf, sigma)
plt.plot(xs, cs, "r-")
cs = bs_put(xs, 100, 1, rf, sigma)
plt.plot(xs, cs, "g-")
cs = bs_put(xs, 100, 0.01, rf, sigma)
plt.plot(xs, cs, "b-")
plt.title("Put Price vs Spot on different Time to Maturity")
plt.show()
```



```
[5]: xs = np.linspace(4, 200, 100)
cs = bs_call(100, xs, 2, rf, sigma)
plt.plot(xs, cs, "r-")
cs = bs_call(100, xs, 1, rf, sigma)
plt.plot(xs, cs, "g-")
cs = bs_call(100, xs, 0.01, rf, sigma)
plt.plot(xs, cs, "b-")
plt.title("Put Price vs Strike on different Time to Maturity")
plt.show()
```



3. Pricing by Integration

Flexibility of Integration in Option Pricing While closed-form solutions are faster, numerical integration offers more flexibility with respect to the payoff function.

For path-independent options, pricing by integration can accommodate any payoff function without the need for decomposition.

Consider, for example, a ‘wedding cake’ option, which has a payoff structure that resembles the tiers of a wedding cake.

This complex payoff structure can be handled directly by numerical integration methods, unlike closed-form solutions that require simpler payoff functions.

In the next section, we’ll discuss the payoff functions for digital call and put options.

```
[6]: S = np.array([1, 2, 3, 4])

# np.ones(shape=S.shape) if S >= 2 else np.zeros(shape=S.shape)

def digital_call_po(S, strike):
    po = np.ones(shape=S.shape)
```

```

    po[S < strike] = 0
    return po

def digital_put_po(S, strike):
    po = np.ones(shape=S.shape)
    po[S > strike] = 0
    return po

digital_put_po(S, 2), digital_call_po(S, 2)

```

[6]: (array([1., 1., 0., 0.]), array([0., 1., 1., 1.]))

Numerical integration can be performed using various methods, Two common methods are:

- **Linear/Quadratic Approximation:** This method approximates the function to be integrated with linear or quadratic functions, making the integral easier to compute.
- **Monte Carlo Integration:** This method uses random sampling to estimate the integral. It's particularly useful when dealing with high-dimensional integrals, but requires a large number of samples to achieve high accuracy

3.1 Numerical Integration

Geometric Brownian Motion and Option Pricing For Geometric Brownian Motion (GBM), we have the following stochastic differential equation, where W is the standard Wiener process:

$$\frac{dS}{S} = r dt + \sigma dW$$

The solution to the GBM is

$$S_t = S_0 e^{(\mu - \frac{\sigma^2}{2})dt + \sigma W(t)}$$

Here, the change in W over a time interval t is a normal random variable with $N(0, t)$, i.e., $W(t) = \sqrt{t}z$. So, we have:

$$S_t = S_0 e^{(\mu - \frac{\sigma^2}{2})t + \sigma \sqrt{t}z}$$

We define a payoff function PayOff, which is the reward at the outcome of the option. The payoff function can take any form and depend on any kind of input, with the asset price S_t being a primary dependent variable.

- If the payoff is **path-independent**, it's a function of the ending price $S_t(z, T)$.
- If the payoff is **path-dependent**, it's a function of the price series $S_t(z, t)$ for $t \in (0, T)$.

When we have a **path-independent** payoff and the inputs T , S_0 , μ , and σ are constant (with μ and σ being the average value during the period), the only variable in the payoff function $\text{Payoff}(S_T(z))$ is the normal random variable z .

The expectation of the PayOff can be calculated as a direct integration of the product of the Payoff function and the probability density function at z , with respect to z (Radon–Nikodym theorem).

$$\mathbb{E}(\text{Payoff}(S_T(z))) = \int_{-\infty}^{+\infty} \text{Payoff}(S_T(z)) * \text{pdf}(z) dz$$

Applying no-arbitrage pricing, the option's present value (PV) is the expected value of PayOff discounted to the present.

$$PV = \text{discount} \times \mathbb{E}(\text{Payoff}(S_T(z)))$$

To combine the above, to price a path-independent option, we just need to perform an integration of the terminal payoff and the PDF of the normal distribution:

$$PV = \text{discount} \times \int_{-\infty}^{+\infty} \text{Payoff}(S_T(z)) * \text{pdf}(z) dz$$

In practice, we can only integrate over a smaller range $[-a, +b]$ instead of $[-\infty, +\infty]$. Thanks to the normal distribution being a **centralized distribution**, a range of $[-6, 6]$ can cover a large part of it. However, the range $[-a, +b]$ should be sufficiently large to cover the entire non-zero range of the payoff function with the option's parameters (`S_0`, `ttm`, `rf`, `sigma`). For example, for a call option with a strike of 110, the integration range should include GBM prices ≥ 110 . The difference between increases in the integration range should be lower than the tolerance.

$$PV = \text{discount} \times \int_{-a}^{+b} \text{Payoff}(S_T(z)) * \text{pdf}(z) dz$$

Let's get z for a range of prices to find the sufficient large range for integration.

```
[7]: def gbm(S_0, ttm, rf, sigma, z):
    S = S_0 * np.exp((rf - sigma**2 / 2) * ttm + sigma * np.sqrt(ttm) * z)
    return S

# z is dependent on all these variable
def z_value(S, S_0, ttm, rf, sigma):
    z = (np.log(S / S_0) - (rf - sigma**2 / 2.0) * ttm) / (sigma * np.sqrt(ttm))
    return z

z_1 = z_value(1, S_0, ttm, rf, sigma)
z_20 = z_value(20, S_0, ttm, rf, sigma)
z_110 = z_value(110, S_0, ttm, rf, sigma)
z_120 = z_value(120, S_0, ttm, rf, sigma)
z_200 = z_value(200, S_0, ttm, rf, sigma)

print((z_1, z_20, z_110, z_120, z_200))
```

```

print(gbm(S_0, ttm, rf, sigma, z=z_20))

# Changes to z when we increase the ttm

z_1 = z_value(1, S_0, ttm * 3, rf, sigma)
z_20 = z_value(20, S_0, ttm * 3, rf, sigma)
z_110 = z_value(110, S_0, ttm * 3, rf, sigma)
z_120 = z_value(120, S_0, ttm * 3, rf, sigma)
z_200 = z_value(200, S_0, ttm * 3, rf, sigma)
print((z_1, z_20, z_110, z_120, z_200))
print(gbm(S_0, ttm * 3, rf, sigma, z=z_20))

```

```

(-10.783779545315657, -3.7227708611509227, 0.2953589263044791, 0.5004467086703,
1.7044742505662334)
20.0
(-6.144368365276803, -2.0676931007240076, 0.25217521370222246,
0.37058270005863253, 1.0657283254839731)
20.0

```

Payoff Function Implementation Numerical Integration (NI) allows us to separate the Payoff and probability, providing an opportunity to implement a more structured approach in programming.

- We can define functions for different payoff types, such as `call_payoff()` and `put_payoff()`
- These can be integrated with the process (`gbm`) in `option_payoff()`.

Next, we'll implement the payoff function as described above. This function will calculate the value of the option at expiry based on the given inputs.

Note: We'll be using the `np.maximum` function from the numpy library, which returns the maximum of the input values.

```

[8]: # (Optional) Here I used npt to do Python typing for numpy

import numpy.typing as npt

def call_payoff(S: npt.ArrayLike, strike: npt.ArrayLike) -> npt.ArrayLike:
    return np.maximum(S - strike, 0)

def put_payoff(S, strike):
    return np.maximum(strike - S, 0)

def option_payoff(S_0, strike, ttm, rf, sigma, payoff_func, z):
    S = gbm(S_0, ttm, rf, sigma, z)
    return payoff_func(S, strike)

```

```

print("Call @ z = 0:", option_payoff(S_0, strike, ttm, rf, sigma, call_payoff, ↵
↵0))
print("Put @ z = 0:", option_payoff(S_0, strike, ttm, rf, sigma, put_payoff, ↵
↵0))

# Check the range of z
print((z_value(300, S_0, ttm, rf, sigma), z_value(20, S_0, ttm, rf, sigma)))

# Verification
print(option_payoff(S_0, 300, ttm, rf, sigma, lambda x, y: x, 2.6601))
print(option_payoff(S_0, 300, ttm, rf, sigma, lambda x, y: x, -3.72277))

```

```

Call @ z = 0: 0.0
Put @ z = 0: 2.9554466451491805
(2.660164675492299, -3.7227708611509227)
299.9917682666852
20.000007307109218

```

Visualizing the Payoff Function and Probability Next, we'll plot the payoff function alongside the probability derived from the PDF of the normal distribution. The value of the option is essentially the product of these two factors, discounted to its Net Present Value (NPV).

This visualization will help us understand how the final option price is influenced by both the payoff at expiry and the probability of that payoff.

```

[9]: import matplotlib.pyplot as plt
import numpy as np
import scipy.stats as sps

S = np.arange(0, 200, 10)
payoff = call_payoff(S, strike)
plt.plot(S, payoff)
plt.title(f"Payoff with {strike=}")
plt.show()

def normal_pdf(z, mean=0, sd=1):
    # return 1/np.sqrt(2*np.pi) when mean = 0, sd = 1
    return 1 / sd / np.sqrt(2 * np.pi) * np.exp(-((z - mean) ** 2) / sd**2 / 2)

def verify_normal_pdf(z, mean=0, sd=1):
    a = normal_pdf(z, mean, sd)
    b = sps.norm.pdf(z, mean, sd)
    print("verify:", a, b, np.isclose(a, b))

```

```

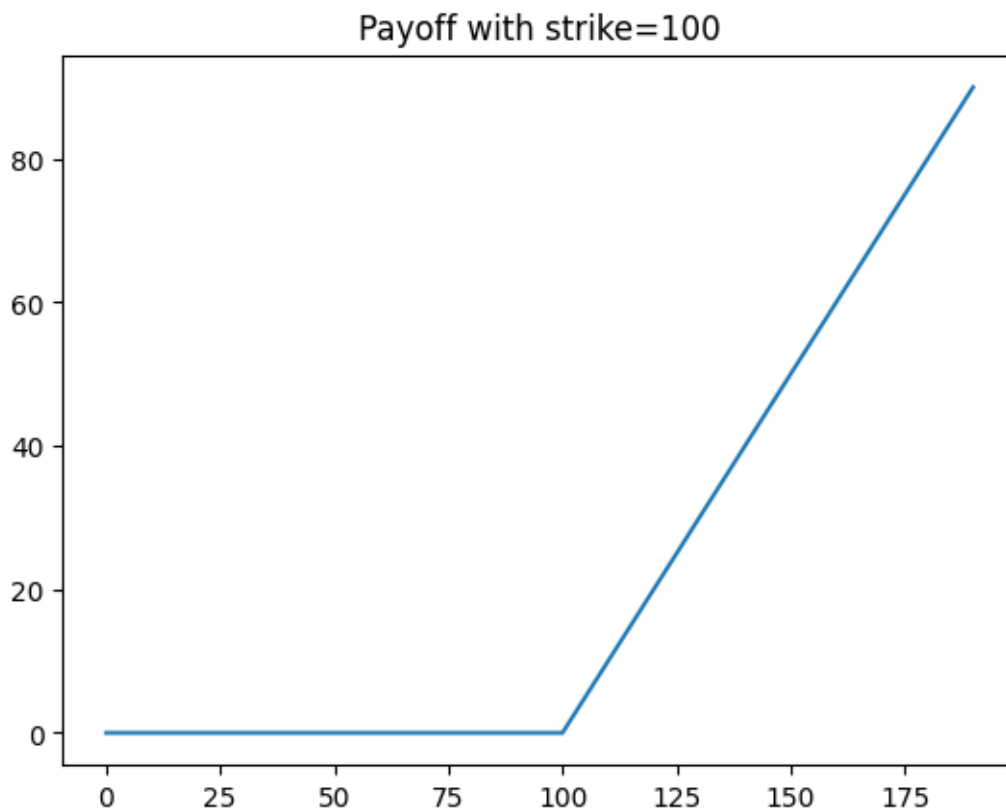
verify_normal_pdf(normal_pdf(0))
verify_normal_pdf(normal_pdf(0, 0.1, 0.05))
verify_normal_pdf(normal_pdf(-0.312, 0.3, 0.05))

def plot_po(z):
    prob = sps.norm.pdf(z)
    S_t = S_0 * np.exp((rf - sigma**2) * ttm + sigma * np.sqrt(ttm) * z)

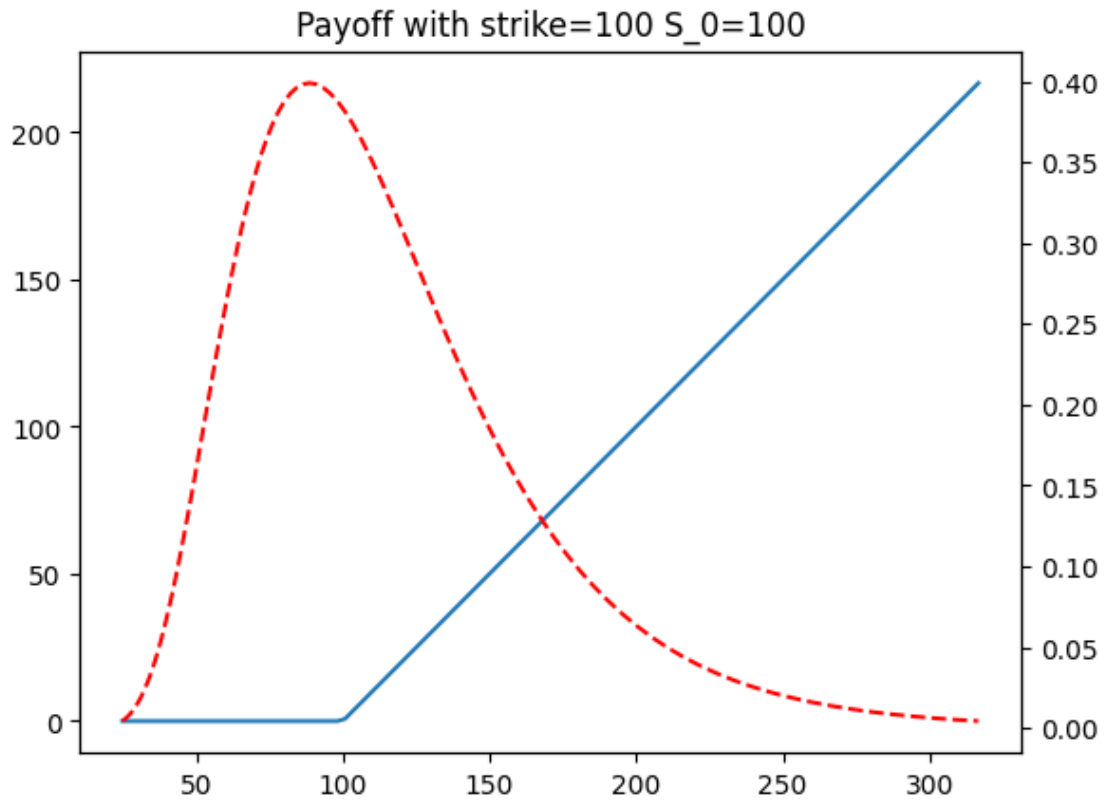
    payoff = call_payoff(S_t, strike)
    plt.plot(S_t, payoff)
    plt.gca().twinx().plot(S_t, prob, "r--")
    plt.title(f"Payoff with {strike=} {S_0=}")
    plt.show()
    print(
        f"S_t/prob range: {np.min(S_t):.3f}/{prob[0]:.3f} {np.max(S_t):.3f}/
        ↪{prob[-1]:.3f} for z in {np.min(z)} to {np.max(z)}"
    )

plot_po(np.linspace(-3, 3, 101))
plot_po(np.linspace(-6, 6, 101))
plot_po(np.linspace(-10, 10, 101))

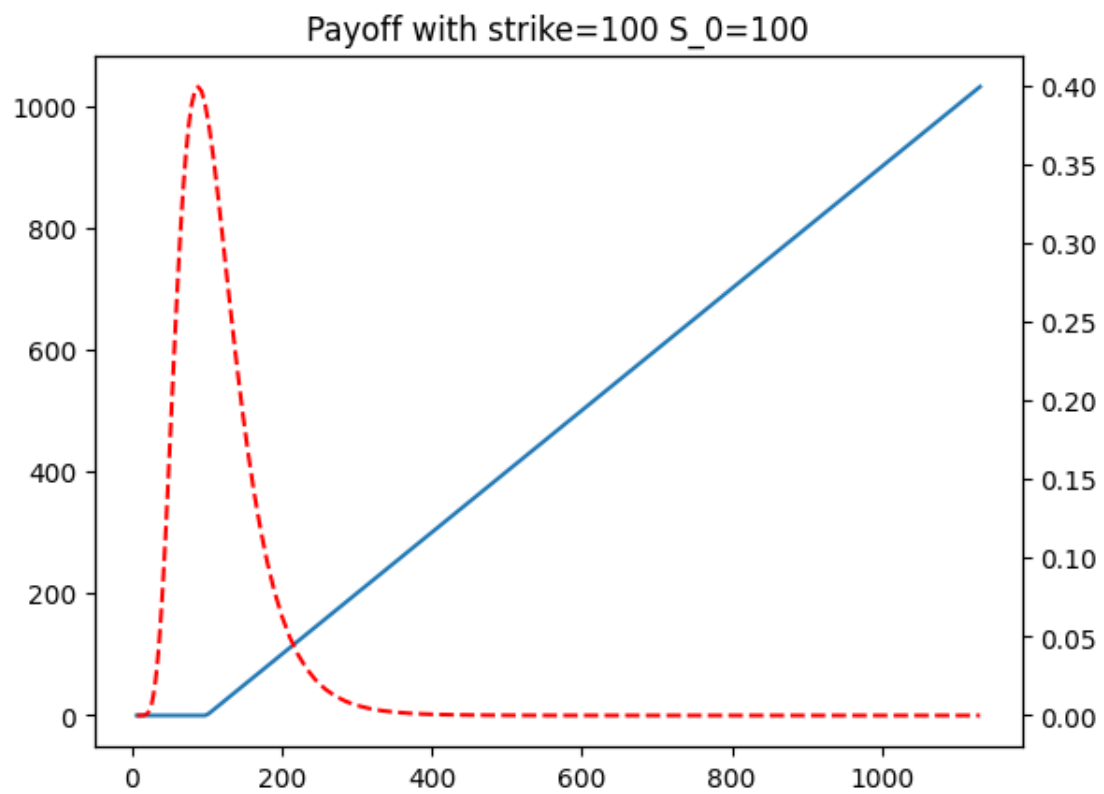
```



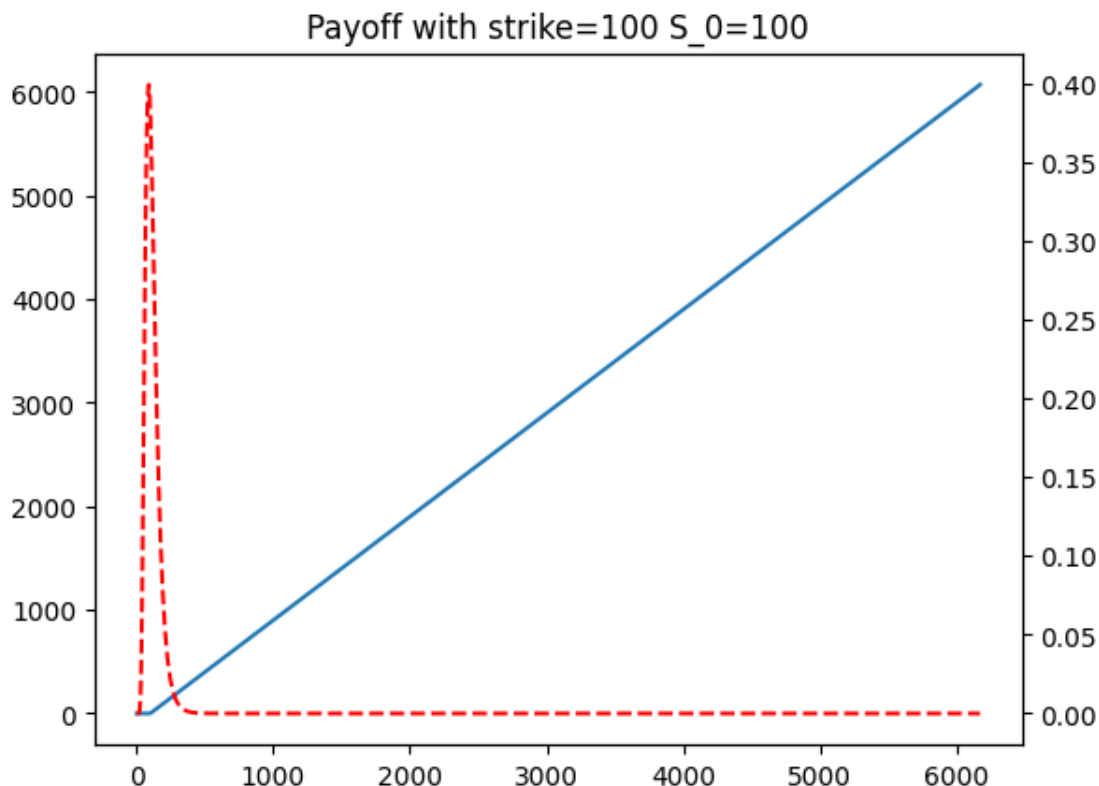
verify: 0.36842577779469815 0.36842577779469815 True
verify: 0.22269694424480851 0.22269694424480854 True
verify: 0.3989422804014327 0.3989422804014327 True



S_t/prob range: 24.838/0.004 316.702/0.004 for z in -3.0 to 3.0



S_t/prob range: 6.956/0.000 1130.884/0.000 for z in -6.0 to 6.0



S_t/prob range: 1.274/0.000 6172.202/0.000 for z in -10.0 to 10.0

3.2.1 Quadrature

```
[10]: import scipy.integrate as spi
import scipy.stats as sps

# integrate a normal distribution's CDF
spi.quad(lambda x: sps.norm.pdf(x), -10, 10)
```

```
[10]: (1.0000000000000002, 8.671029888166837e-10)
```

```
[11]: import scipy.integrate as spi
import scipy.stats as sps

# Below uses a very concise form of integration with lambda functions.

ni_call = (
    np.exp(-rf * ttm)
    * spi.quad(
        lambda x: np.maximum(
            S_0 * np.exp((rf - 0.5 * sigma**2) * ttm + sigma * np.sqrt(ttm) * x)
            - strike,

```

```

        0,
    )
    * sps.norm.pdf(x),
    -10,
    10,
)[0]
)
ni_put = (
    np.exp(-rf * ttm)
    * spi.quad(
        lambda x: np.maximum(
            strike
            - S_0 * np.exp((rf - 0.5 * sigma**2) * ttm + sigma * np.sqrt(ttm) *
↪x),
            0,
        )
        * sps.norm.pdf(x),
        -10,
        10,
    )[0]
)
print(
    "NI vs. Analytical:",
    ni_call,
    ni_put,
    bs_call(S_0, strike, ttm, rf, sigma),
    bs_put(S_0, strike, ttm, rf, sigma),
)

# To reduce code density, let's re-write

def integrand(S_0, strike, ttm, rf, sigma, payoff_func):
    def _inner(z, payoff_func=payoff_func):
        return option_payoff(
            S_0, strike, ttm, rf, sigma, payoff_func, z
        ) * sps.norm.pdf(z)

    return _inner

ni_call = (
    np.exp(-rf * ttm)
    * spi.quad(integrand(S_0, strike, ttm, rf, sigma, call_payoff), -10, 10)[0]
)
ni_put = (
    np.exp(-rf * ttm)

```



```

    * spi.quad(integrand(S_0, strike, ttm, rf, sigma, put_payoff), -10, 10)[0]
)
print(
    "NI (quad) vs. Analytical:",
    ni_call,
    ni_put,
    bs_call(S_0, strike, ttm, rf, sigma),
    bs_put(S_0, strike, ttm, rf, sigma),
)

```

```

NI vs. Analytical: 19.382549294727323 13.559002656588023 19.38254929806086
13.559002656485731
NI (quad) vs. Analytical: 19.382549294727323 13.559002656588023
19.38254929806086 13.559002656485731

```

3.2.2 Trapezoidal Integration We can utilize the `spi.trapezoid` function from the SciPy library for numerical integration. This function applies the trapezoidal rule to compute the integral of a function. We can *reuse* our previously defined `integrand` function with this method to calculate the option price.

```

[12]: ll = np.linspace(-10, 10, 1001)
ni_call = np.exp(-rf * ttm) * spi.trapezoid(
    integrand(S_0, strike, ttm, rf, sigma, call_payoff)(ll), ll
)
ni_put = np.exp(-rf * ttm) * spi.trapezoid(
    integrand(S_0, strike, ttm, rf, sigma, put_payoff)(ll), ll
)
print(f"NI (trapezoidal): {ni_call, ni_put}")

```

```

NI (trapezoidal): (19.382810242040463, 13.55926360046533)

```

3.2.3 By-hand We write a customized `hand_int` function but still with the previously defined `integrand` function.

```

[13]: def hand_int(int_func, upper_b=10, lower_b=-10, points=1000):
    h = (upper_b - lower_b) / points
    ps = np.linspace(-10, 10, points + 1)
    int_ps = int_func(ps)
    return h * np.sum(int_ps)

ni_call = np.exp(-rf * ttm) * hand_int(
    integrand(S_0, strike, ttm, rf, sigma, call_payoff)
)
ni_put = np.exp(-rf * ttm) * hand_int(
    integrand(S_0, strike, ttm, rf, sigma, put_payoff)
)

```

```
print(f"by-hand: {ni_call, ni_put}")
```

by-hand: (19.38281024204046, 13.55926360046533)

3.2.5 Monte Carlo Integration Monte Carlo integration is a technique that uses random sampling to approximate definite integrals. For a function $q(x)$ over the interval $[a, b]$, the integral can be estimated as:

$$\int_a^b q(x)dx = (b-a) \int_a^b q(x) \frac{1}{b-a} dx = (b-a) \int_a^b q(x) f_U(x) dx = (b-a) \mathbb{E}(q(U))$$

Here, U is a continuous uniform random variable with density function $f_U(x) = \frac{1}{b-a}$.

Moving to high-dimension

- One dimension

$$\int_c^d q(x)dx = (b-a) \mathbb{E}(q(U_x))$$

- Two and higher Dimension

$$\int_c^d \int_a^b q(x, y) dx dy = (d-c)(b-a) \mathbb{E}(q(U_x, U_y))$$

Next, we'll implement a Monte Carlo integration function. We'll find our previously defined `integrand()` function useful here. Run the following code a few times to observe the results.

```
[14]: import numpy.random as npr

def mc_int(func, rg, N=int(1e6)):
    nr = npr.uniform(rg[0], rg[1], size=N)
    sums = func(nr)
    return np.mean(sums) * (rg[1] - rg[0])

ni_call = np.exp(-rf * ttm) * mc_int(
    integrand(S_0, strike, ttm, rf, sigma, call_payoff), (-6, 6)
)
ni_put = np.exp(-rf * ttm) * mc_int(
    integrand(S_0, strike, ttm, rf, sigma, put_payoff), (-6, 6)
)
print(
    "mc:",
    ni_call,
    ni_put,
    bs_call(S_0, strike, ttm, rf, sigma),
```

```
bs_put(S_0, strike, ttm, rf, sigma),
)
```

mc: 19.373052349299517 13.554526838807101 19.38254929806086 13.559002656485731

[15]: # 2-d MC integration

```
def mc_int_2d(func, rg1, rg2, N=int(1e6)):
    nr1 = npr.uniform(rg1[0], rg1[1], size=N)
    nr2 = npr.uniform(rg2[0], rg2[1], size=N)
    sums = func(nr1, nr2)
    print(sums.shape)
    res = np.mean(sums) * (rg1[1] - rg1[0]) * (rg2[1] - rg2[0])
    std = np.std(sums) * (rg1[1] - rg1[0]) * (rg2[1] - rg2[0]) / np.sqrt(N)
    return res, std

print(
    mc_int_2d(lambda x, y: np.sqrt(x**2 + y**2) < 1, (0, 1), (0, 1), int(1000 *
↪100)),
    np.pi / 4,
)

def rec_int_2d(func, rg1, rg2, N=int(1e6)):
    nr1, nr2 = np.meshgrid(
        np.linspace(rg1[0], rg1[1], N), np.linspace(rg2[0], rg2[1], N)
    )
    sums = func(nr1, nr2)
    print(sums.shape)
    return sums.sum() / N**2 * (rg1[1] - rg1[0]) * (rg2[1] - rg2[0])

print(
    rec_int_2d(lambda x, y: np.sqrt(x**2 + y**2) < 1, (0, 1), (0, 1),
↪int(1000)),
    np.pi / 4,
)
```

```
(100000,)
(0.78421, 0.0013008638510620548) 0.7853981633974483
(1000, 1000)
0.784785 0.7853981633974483
```

3.3 Designing Interfaces for Functions

In the following section, we'll present a complete version of the code, broken down into step-by-step segments.

Design Guidelines

1. **Separate Data from Logic:** This makes the code easier to understand and maintain. It also allows for independent testing and reuse of logic.
2. **Make Logic Step-able:** Breaking down the logic into smaller, manageable steps makes the code easier to debug and understand.
3. **Make Code Reusable:** Design your code such that minimal changes are required to add new functionality. This increases the longevity and utility of your code.

```
[16]: import numpy as np
import scipy.integrate as spi
import scipy.stats as sps

def call_payoff(S, strike):
    return np.maximum(S - strike, 0)

def put_payoff(S, strike):
    return np.maximum(strike - S, 0)

def gbm(S_0, strike, ttm, rf, sigma, z):
    return S_0 * np.exp((rf - 0.5 * sigma**2) * ttm + sigma * np.sqrt(ttm) * z)

def option_payoff(S_0, strike, ttm, rf, sigma, payoff_func, z):
    S = gbm(S_0, strike, ttm, rf, sigma, z)
    return payoff_func(S, strike)

def integrand(S_0, strike, ttm, rf, sigma, payoff_func):
    def _inner(z, payoff_func=payoff_func):
        return option_payoff(
            S_0, strike, ttm, rf, sigma, payoff_func, z
        ) * sps.norm.pdf(z)

    return _inner

S_0 = 100
strike = 100
```

```

ttm = 2
rf = 0.03
sigma = 0.3

ni_call = (
    np.exp(-rf * ttm)
    * spi.quad(integrand(S_0, strike, ttm, rf, sigma, call_payoff), -10, 10)[0]
)
ni_put = (
    np.exp(-rf * ttm)
    * spi.quad(integrand(S_0, strike, ttm, rf, sigma, put_payoff), -10, 10)[0]
)
print("spi.quad:", ni_call, ni_put)

```

```
spi.quad: 19.382549294727323 13.559002656588023
```

Utilizing the functools Module In our code, we use an `_inner` function to return a function from another function. Alternatively, we could use `functools.partial` to return a partially initialized function with a subset of the required inputs.

This technique is known as “currying”. The returned function can be executed later by calling it with the remaining parameters. This approach can enhance code readability and reusability.

```

[17]: import functools

def f(a, b):
    return a - b

# f(3, 4)

print(
    (
        # add the name of the argument to be assigned
        (
            functools.partial(f, b=3)(4),
            functools.partial(f, 3)(4),
        ),
        (
            (lambda a, b=3: f(a, b))(
                4
            ), # The function definition limits the the default value to be
            ↪ set for later not earlier arguments.
            (lambda b, a=3: f(a, b))(4),
            ), # Need to put the earlier argument to the backseat to allow for
            ↪ default value.
        # Continue to use partial to get a function with ()
    )
)

```

```

        functools.partial(functools.partial(f, 3), 4)(),
    )
)

```

```
((1, -1), (1, -1), -1)
```

Refactoring with `functools.partial` Next, we'll revisit our original code and improve it by using `functools.partial`. This will allow us to create partially initialized functions, enhancing the modularity and reusability of our code.

```
[18]: import functools

# An improved version to integrand()
def integrand2(func, z):
    return func(z) * sps.norm.pdf(z)

def priceEuropeanOption(S, strike, ttm, rf, sigma, payoff_func):
    op_x = functools.partial(option_payoff, S, strike, ttm, rf, sigma,
    ↪payoff_func)
    int_x = functools.partial(integrand2, op_x)
    discountFactor = np.exp(-rf * ttm)
    R = 10
    return discountFactor * spi.quad(int_x, -R, R)[0]

print(priceEuropeanOption(S_0, strike, ttm, rf, sigma, call_payoff))
print(priceEuropeanOption(S_0, strike, ttm, rf, sigma, put_payoff))

```

```
19.382549294727323
```

```
13.559002656588023
```

Recap and Discussion

- **Recap:** Always remember to include the discount factor in your calculations. However, be careful not to apply the discounting twice.
- **Discussion:** The choice between Trapezoidal Integration, Gaussian Quadrature Integration, and Monte Carlo Integration depends on the specific requirements of your problem:
 - **Trapezoidal Integration:** This method is simple and easy to implement. It's best suited for problems where the function is relatively smooth and the integral is over a low-dimensional space.
 - **GQ Integration:** This method is more accurate than trapezoidal integration for functions that are not smooth. It's also efficient for low-dimensional integrals.
 - **Monte Carlo Integration:** This method is best for high-dimensional integrals, where traditional numerical integration methods become inefficient or impractical. However,

it requires a *very* large number of samples to achieve high accuracy.

4. Assignment

1. **Binary Options Pricing:** Apply Numerical Integration to price binary call/put options. The payoff of a binary call is 1 if $S > K$ and 0 otherwise, while for a binary put it's 1 if $S < K$ and 0 otherwise. Compare the results between the three numerical integration methods.
2. **Butterfly Options Pricing:** Calculate the price of a series of adjacent butterfly options ranging from \$0 to \$200 using numerical integration. Each butterfly option has a difference of strike prices of \$10. Repeat the calculation with a difference of strike prices of \$1.

The option parameters used in tasks 1 and 2 are:

```
S_0 = 100
strike = 100
ttm = 2
rf = 0.03
sigma=0.3
```

Note: A butterfly call option's payoff is composed of four call options with strikes $K_1 < K_2 < K_3$ and is calculated as $Call(K_1) - 2 * Call(K_2) + Call(K_3)$. In this question, $K_2 - K_1 = K_3 - K_2 = 10$ or \$1.

3. **Volume Calculation:** Use Monte Carlo integration to calculate the volume of a cube with the left-bottom point at non-zero coordinates (a, b, c) . Calculate for 1) a cube where each side has a length of 1 unit, and 2) a cuboid where each side has lengths of 1, 2, and 3 units respectively.

Appendix: timestamp

```
[19]: from datetime import datetime

print(f"Generated on {datetime.now()}")
```

Generated on 2024-03-07 12:20:41.343401