# Simulation Project

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#### Abstract

The project explores two distinct applications of Stochastic Approximation (SA) algorithms in optimization tasks. In **Project 1**, the focus was on optimizing investment strategies for an investor with capital distributed among three companies. The SA algorithm determined an allocation strategy yielding a maximum Sharpe ratio, suggesting approximately 29.17% of capital to company 1, 19.16% to company 2, and 51.67% to company 3. In **Project 2**, the objective was to minimize waiting times in a service station by optimizing the mean service time  $\theta$  using SA. Results demonstrated convergence to an optimal  $\theta$  value of approximately 3.4698, with a standard deviation of 0.0040. Additionally, an algorithm was developed to concurrently update  $\theta$  and estimate  $q_{\theta}$  within a fixed computational budget, showcasing efficient computational resource management.

Key words: Stochastic Approximation; Investment Optimization; Queuing System Optimization

### Project 1

#### 1 Introduction

The first project focuses on an investor with a total capital of 1, aiming to allocate this capital across three companies to maximize the risk-adjusted performance of the investment. The market value and returns of each company are influenced by independent uniformly distributed returns and various market factors, including common economic conditions and company-specific risks. The objective is to determine the optimal allocation vector that maximizes the Sharpe ratio, reflecting the trade-off between expected profit and risk. This is achieved by formulating and solving an optimization problem using an SA algorithm.

### 2 Description of Project 1

In this project, we aim to determine the optimal investment strategy for an investor with a total capital of 1 unit. The capital is to be distributed among n companies over a time period of t units. The market value of company i at time t is given by  $X_i$ . Let  $x_i \in \mathbb{R}$ , i = 1, ..., n be given thresholds. Company i will not be able to generate any profit if  $X_i < x_i$ . If  $X_i \ge x_i$ , the return on the investment is given by  $Y_i$ . Investing a fraction  $p_i$  of the capital yields an expected return of

$$p_i \mathbb{E}[Y_i \mathbf{1}_{X_i \ge x_i}] \tag{1}$$

for company i, where

$$\sum_{i=1}^{n} p_i = 1 \quad \text{and} \quad 0 \le p_i \le 1.$$
 (2)

Assume that n = 3 and  $Y_i$  are independent and uniformly distributed on  $[0, X_i]$ . We let

$$X_{i} = \frac{\rho V + \sqrt{1 - \rho^{2}} \eta_{i}}{\max(W, 1)}, \quad 1 \le i \le n,$$
(3)

with  $\eta_i$  normally distributed with mean 0 and variance i modeling the company's idiosyncratic risk, V standard normally distributed modeling the common factor that affects the economy, and W exponentially distributed with rate 1/0.3 modeling common market shocks. The variables V, W, and  $\eta_i$ 's are all independent. The weight factor is set to  $\rho = 0.6$ , and the thresholds are  $x_1 = 2$ ,  $x_2 = 3$ , and  $x_3 = 1$ .

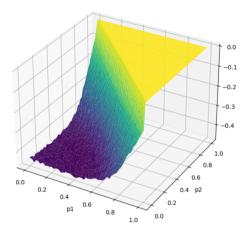
The investor seeks to find the optimal investment strategy by maximizing the risk-adjusted performance of the investment, also known as the Sharpe ratio, which leads to the objective function:

$$\max_{p_1, p_2, p_3 \ge 0, \sum p_i = 1} \mathbb{E}\left[\frac{\sum_{i=1}^3 p_i Y_i \mathbf{1}_{X_i \ge x_i}}{\operatorname{std}\left(\sum_{i=1}^3 p_i Y_i \mathbf{1}_{X_i \ge x_i}\right)}\right],\tag{4}$$

where std(X) denotes the standard deviation of the random variable X. To simplify the problem and solution process, we converted the objective function to:

$$\min_{p_1, p_2, p_3 \ge 0, \sum p_i = 1} -\mathbb{E}\left[\frac{\sum_{i=1}^3 p_i Y_i \mathbf{1}_{X_i \ge x_i}}{\operatorname{std}\left(\sum_{i=1}^3 p_i Y_i \mathbf{1}_{X_i \ge x_i}\right)}\right],$$
(5)

In order to visualize the objective function for the optimization, we drew the 3D plot of the function (Figure 1 and Figure 2).



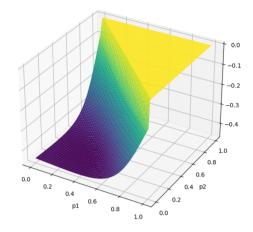


Figure 1: J function without random seed

Figure 2: J function with random seed = 20

### 3 Motivation for the Project

In real-life scenarios, financial data is often noisy and uncertain. Therefore, it is crucial to use optimization methods that can handle such noise effectively. Stochastic Approximation (SA) algorithms are particularly suited for this purpose as they are designed to deal with noisy observations, making them ideal for financial applications.

## 4 Optimization Approach

To address the optimization problem, we employed a Stochastic Approximation (SA) algorithm, specifically used the Simultaneous Perturbation Stochastic Approximation (SPSA) algorithm to calculate the gradient. Here's a detailed explanation of the approach.

First of all, this optimization problem is well-posed because it has a clear objective function to maximize (the Sharpe ratio), subject to the constraints that the sum of the investment fractions  $p_i$  equals 1 and each  $p_i$  is between 0 and 1. The constraints ensure that the solution space is bounded and feasible.

What's more, the vector field in this context is the gradient of the objective function with respect to the investment fractions  $p_i$ . The coercivity of this vector field is ensured by the constraints and the nature of the Sharpe ratio, which promotes exploration within the feasible region.

Since the problem is well-posed, and the gradient is coercive for the problem. Therefore, the update rule:  $\theta_{n+1} = \theta_n + \epsilon Y_n$  converges to a solution. This convergence is guaranteed by

the Stochastic Approximation Theory, which states that under conditions of unbiased gradient estimates and a properly decreasing gain sequence, the iterative updates will converge to the true solution of the optimization problem. Therefore, we employ the Stochastic Approximation (SA) algorithm to conduct the optimization.

The Simultaneous Perturbation Stochastic Approximation (SPSA) algorithm iteratively updates the investment fractions  $p_i$  using random perturbations and a decreasing gain size. The gain size decreases over iterations, which ensures convergence.

The whole algorithm can be described as follows:

- 1. Initialize  $p_1$  and  $p_2$  to 0, and compute  $p_3 = 1 p_1 p_2$ .
- 2. For each iteration i:
- Generate random perturbations  $r_1$  and  $r_2$ .  $P(r_1 = -1) = P(r_1 = 1) = \frac{1}{2}$ , same as  $r_2$ .
- Calculate the step size as  $\theta = \frac{1}{(i+10000)^{0.3}}$ .

Proving process: Because  $\sum_{i=n}^{\infty} \epsilon = \infty$ ,  $\sum_{i=n}^{\infty} \epsilon \theta^2 < \infty$  and  $\sum_{i=n}^{\infty} \left(\frac{\epsilon^2}{\theta^2}\right) < \infty$ , so  $\epsilon = \frac{1}{\alpha}$  which  $\alpha$  is linear polynomial about i and therefore  $\theta = \frac{1}{k^c}$  which k is linear polynomial about i. Also, we have to ensure that  $\frac{1}{i+100} < \frac{1}{\alpha} < \frac{1}{i+1}$  and  $\frac{1}{(i+10^5)^c} < \frac{1}{k^c} < \frac{1}{(i+10)^c}$ .

Since  $\frac{1}{i+100} < \frac{1}{\alpha} < \frac{1}{i+1}$ ,  $\sum_{i=n}^{\infty} \frac{1}{i+100} < \sum_{i=n}^{\infty} \frac{1}{\alpha} = \infty$  and  $\sum_{i=n}^{\infty} \epsilon \theta^2 < \sum_{i=n}^{\infty} \frac{1}{(i+1)^{2c+1}} < \infty$ , so 2c+1>1, here's the formula 1. Besides that, we also can find that  $\sum_{i=n}^{\infty} \frac{\epsilon^2}{\theta^2} < \sum_{i=n}^{\infty} \frac{(i+10^5)^{2c}}{(i+1)^2}$ , we want to make it convergence. In order to do that, we assume  $f(i) = \frac{(i+10^5)^{2c}}{(i+1)^2}$ , when  $i \to \infty$ , f(i) turns to  $i^{\frac{1}{2-2c}}$ . And because f(i) should be convergence, 2-2c>1, which is the formula 2. Combine both formulas together, and we will find a range of c that 0 < c < 0.5, and we choose c = 0.3 as our c in the  $\theta$  formula.

By referring to the "Taylor Expansion" and "Newton-Raphson Method" mentioned in the "Optimization and Learning via Stochastic Gradient Search", we determined the y1 and y2 with the formula list above.

• Estimate the gradients  $y_1$  and  $y_2$  using the objective function J:

$$y_1 = \frac{J(p_1 + \theta r_1, p_2 + \theta r_2) - J(p_1 - \theta r_1, p_2 - \theta r_2)}{2\theta r_1},$$
(6)

$$y_2 = \frac{J(p_1 + \theta r_1, p_2 + \theta r_2) - J(p_1 - \theta r_1, p_2 - \theta r_2)}{2\theta r_2}.$$
 (7)

• Update  $p_1$  and  $p_2$ :

$$p_1 \leftarrow \max(p_1 - \text{step\_size} \cdot y_1, 0),$$
 (8)

$$p_2 \leftarrow \max(p_2 - \text{step\_size} \cdot y_2, 0).$$
 (9)

According to the requirements of the problem,  $0 \le p_1 + p_2 \le 1$ , and both  $p_1$  and  $p_2$  are greater than or equal to 0. When the generated points exceed the range, we need to project them back into the constraints. If the newly generated  $p_1$  or  $p_2$  is greater than 1, then assign  $p_1$  and  $p_2$  the value of 1. If  $p_1$  or  $p_2$  is less than 0, then assign  $p_1$  or  $p_2$  the value of 0. If  $p_1 + p_2$  falls within the area where x + y > 1 and is bounded by x = 1 and y = 1 (denoted as area M), project the points onto the line y + x = 1.

For example, Figure 3 displays the projection process.

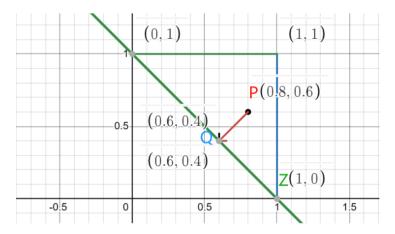


Figure 3: The Projection Process

If a point falls within area M, we will calculate its projection. Suppose the point on the line is Q, and arbitrarily take a point Z on the line y + x = 1 (here for the sake of calculation

convenience, we take (1,0) and perform the following operation:

$$\overrightarrow{QP} = (p_1 - x_0, p_2 - y_0), \overrightarrow{QZ} = (1 - x_0, -y_0)$$
 (10)

$$\overrightarrow{QP} \cdot \overrightarrow{QZ} = 0 \tag{11}$$

$$\Leftrightarrow \begin{cases} (p_1 - x_0)(1 - x_0) + (p_2 - y_0)(-y_0) = 0\\ x_0 + y_0 = 1 \end{cases}$$
 (12)

$$\Leftrightarrow (p_1 - 1 + y_0)y_0 = (p_2 - y_0)y_0 \tag{13}$$

$$\therefore y_0 \neq 0 \tag{14}$$

$$\therefore \begin{cases} x_0 = \frac{p_1 - p_2 + 1}{2} \\ y_0 = \frac{p_2 - p_1 + 1}{2} \end{cases}$$
 (15)

• Ensure that the sum of  $p_1$ ,  $p_2$ , and  $p_3$  equals 1 by adjusting  $p_1$  and  $p_2$  if necessary.

The algorithm runs for a specified number of iterations, and the final values of  $p_1$ ,  $p_2$ , and  $p_3$  represent the optimal investment fractions.

### 5 Experimental Setting

Sample Size: we used a sample size of 10,000 for generating the random variables. This large sample size helps accurately estimate the expectations and standard deviations required for the Sharpe ratio calculation.

Confidence Bounds: to ensure the reliability of our results, we performed multiple experiments and recorded the means of the last 500 iterations for each experiment. This approach provides a measure of stability and confidence in the algorithm's convergence.

#### 5.1 Algorithm Implementation

We implemented the SPSA algorithm with the following settings:

- Initial Points:  $p_1$  and  $p_2$  start at 0, with  $p_3$  computed as  $1 p_1 p_2$ .
- Iteration Count: The algorithm runs for 3,000 iterations per experiment. The choice of 3,000 iterations is based on the total computational budget and time constraints for obtaining results. Also, 3,000 iterations are found to be sufficient for the algorithm to converge to a near-optimal solution (Figure 4), therefore it becomes a reasonable stopping criterion.

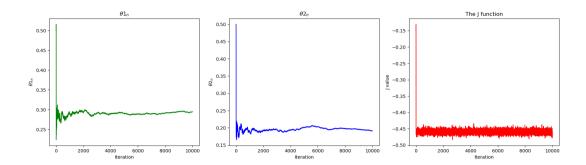


Figure 4: Result with 10, 000 Iterations

- Step Size: We tried fixed gain size 0.01 and decreasing gain size  $\frac{80}{(i+100)}$  and  $\frac{10}{(i+100)}$  to experiment with different strategies for adjusting the gain dynamically.
- Gradient Calculation: The gradient is estimated using random perturbations in both  $p_1$  and  $p_2$  directions.

## 6 Validation and Verification

Figure 4 presents the three-dimensional representation utilized in our prior analysis to elucidate the shape of the objective function. The annotated red circle in the figure indicates the position of the optimization result, thereby substantiating the validity of our experimental outcomes.

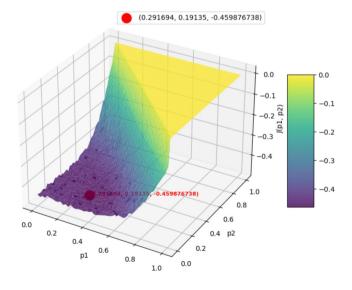


Figure 5: Result in 3D Plot of the Objective Function

## 7 Results and Analysis

The results are recorded in text files and visualized using plots. Each plot shows the evolution of  $p_1$ ,  $p_2$ , and the objective function J over iterations. The mean values of  $p_1$  and  $p_2$  over the last 500 iterations are used to determine the optimal allocation strategy.

#### 7.1 Gain Size Analysis

The figures show that optimization with a fixed gain size leads to weaker convergence, while a decreasing gain size behaves better and costs less. Specifically, we found that a gain size of  $\frac{80}{(i+100)}$  converges slower than  $\frac{10}{(i+100)}$ . Therefore, the gain size is updated as  $\frac{10}{(i+100)}$ , leveraging the strengths of the SA algorithm.

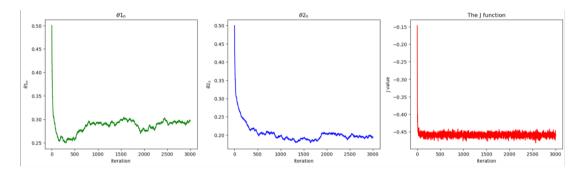


Figure 6: Fixed Gain Size 0.01

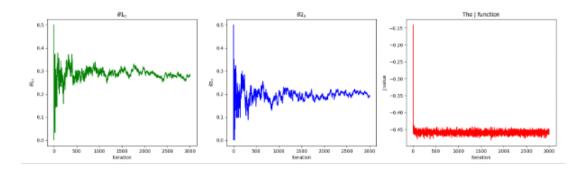


Figure 7: Decreased Gain Size  $\frac{80}{(i+100)}$ 

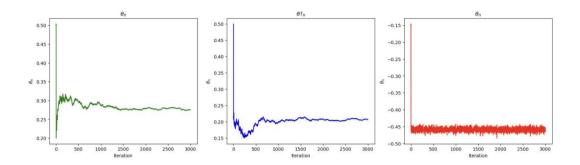


Figure 8: Decreased Gain Size  $\frac{10}{(i+100)}$ 

### 7.2 Initial Point Analysis

We tried to choose 2 different initial points for our task. However, both can converge to a minimum value, as seen in the figures. Hence, there is no significant difference between the 2 initial points. Therefore, we picked one of them as our initial point which is (0.5, 0.5).

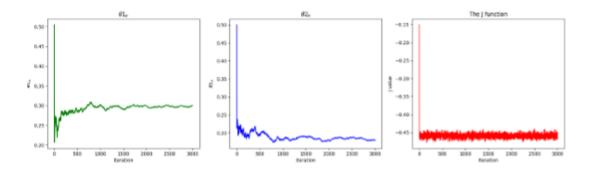


Figure 9: Initial Point (0.5, 0.5)

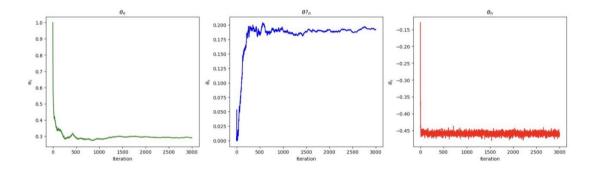
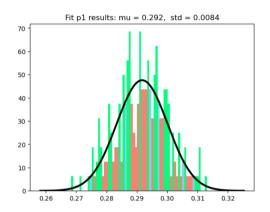


Figure 10: Initial Point (1, 0)

#### 7.3 Output Analysis

Upon initial analysis, it was determined that n=216 was insufficient to achieve normally distributed results for  $\theta_n(p_1, p_2)$ , as depicted in Figures 11 and 12. Subsequently, a larger sample size of 505 runs was utilized, resulting in approximately normally distributed outcomes, as illustrated in Figures 13 and 14. The estimated means  $\mu_{p_1} = 0.291681$  and  $\mu_{p_2} = 0.191590$  were derived from this sample size, indicating adequacy for assuming a normal distribution of  $\hat{p}$ .



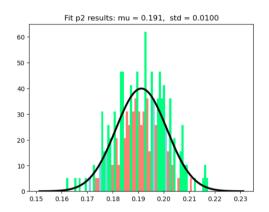
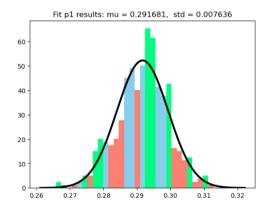


Figure 11: p1 Result Distribution of 216 Runs Figure 12: p2 Result Distribution of 216 Runs



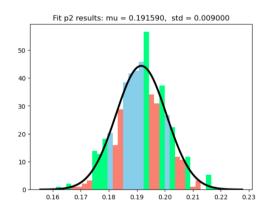


Figure 13: p1 Result Distribution of 505 Runs Figure 14: p2 Result Distribution of 505 Runs

For a sample size of 505, the 95% confidence interval for  $\hat{p}_1$  with distribution N(0.29168, 0.00763) is [0.27671, 0.29101]. The 95% confidence interval for  $\hat{p}_2$  with distribution N(0.19159, 0.00900) is [0.17395, 0.20923].

## 8 Conclusion

In conclusion, the optimal investment strategy for the investor with a total capital of 1 unit, distributed among three companies, has been determined using the Stochastic Approximation (SA) algorithm. The results suggest that allocating approximately 29.17% of the capital to company 1, 19.16% to company 2, and the remaining 51.67% to company 3 yields the highest Sharpe ratio, maximizing the risk-adjusted performance of the investment.

The experimental results confirm the effectiveness of the SA algorithm in handling noisy observations and optimizing financial portfolios. The analysis also demonstrates the importance of selecting an appropriate step size and gain size for convergence. The confidence intervals and distribution analysis further validate the robustness of the obtained solutions.

## Project 2 Question 1

### 1 Introduction

This project investigates optimising waiting times in a service station using a GI/GI/1 queueing model with a first-come, first-served (FCFS) discipline. Customers arrive according to a renewal point process with inter-arrival times  $\{A_n\}$  that are independent and identically distributed (iid) with finite mean  $E[A_n]$ . The service times  $\{S_n(\theta)\}$  are also iid and depend on a parameter  $\theta$ , representing the average service time, which is our key variable for optimization. The primary objective is to minimize the average waiting time for customers while maintaining a cost-effective operation of the service station.

The main goal is to determine the optimal mean service time  $\theta$  that minimizes the cost associated with the average waiting time  $W_n(\theta)$  for the first N customers. This involves applying a Stochastic Approximation (SA) algorithm to find the optimal  $\theta$  that satisfies the desired service level, specifically minimizing the 90% quantile of the average waiting time.

### 2 Methodology

#### 2.1 Model Formulation

Define the waiting time  $W_n(\theta)$  recursively:

$$W_{n+1}(\theta) = \max(0, W_n(\theta) + S_n(\theta) - A_{n+1}),$$

with initial condition  $W_0(\theta) = S_0(\theta) = 0$ .

Calculate the average waiting time over the first N customers:

$$W^{N}(\theta) = \frac{1}{N} \sum_{n=1}^{N} W_{n}(\theta).$$

#### 2.2 Optimization Problem

Aim to minimize the expected squared deviation of the 90% quantile  $q_{\theta}$  of  $W^{N}(\theta)$  from a target value z:

$$\min_{0<\theta} E[(q_{\theta}-z)^2].$$

#### 2.3 Stochastic Approximation Algorithm

Implement the SA algorithm to estimate the quantile and optimize  $\theta$ . Use the estimator  $F_{\theta}(w) = P(W^{N}(\theta) \leq w)$ , which is monotone decreasing in  $\theta$  for fixed w.

### 3 Experiment Setting

#### 3.1 Parameter Selection

Inter-arrival times are exponentially distributed with mean value 5. Service times are exponentially distributed with mean value  $\theta$ . Set N=10 and z=8.

#### 3.2 Algorithm Implementation

We initially tested multiple different functions, including Simultaneous Perturbation Stochastic Approximation (SPSA) and Stochastic Approximation (SA), each with multiple different data. After evaluating their performance, we decided to use the SA algorithm. This decision was based on its superior performance in terms of stability and convergence towards the optimal mean service time  $\theta$ .

For the question, it's an ill-posed problem, because  $\theta > 0$ . In order to make it a well-posed problem, we have to make the  $\theta$  range bigger or equal to 0. Set

$$\min_{0 \le \theta} J(\theta) = \min_{0 \le \theta} E\left[ (q_{\theta} - z)^2 \right]$$
 (16)

$$\nabla \hat{J}(\theta) = 2E[(q_{\theta} - z)] \tag{17}$$

 $\nabla J(\theta) = 2E[(q_{\theta} - z)] \times \nabla(q_{\theta} - z)$ , Because  $q_{\theta}$  approximates the average service time for a G/G/1 queue with a mean service time of  $\theta$ , the average waiting time for the top 90% of people. As  $\theta$  increases,  $q_{\theta}$  is a strictly monotonically increasing function, because as the average service time lengthens, the waiting time will definitely become longer. Therefore, the derivative of  $(q_{\theta} - z)$  is positive. In this project we dropped it, which means setting this value to 1 to approximate the gradient.

### 4 Validation and Verification

The result we find from the normal distribution is exactly the same as the red spot we allocated in the figure below, which proves the validity of our experiment.

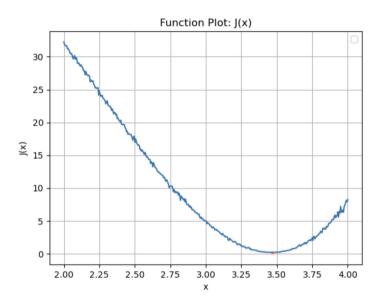


Figure 1: Result in the Objective Function

## 5 Results and Analysis

In this part, we will present the results of the SA algorithm, showing the convergence of  $\theta$  towards the optimal value. The figure represented a normal curve, which is symmetric and centered about the mean value of 3.469775865. With a standard deviation of 0.0039682, the data points are tightly clustered around the mean, suggesting that the values are relatively close to each other.

The 95% confidence interval for  $\hat{\theta}$  with distribution N(3.469775865, 0.0039682) is [3.4619980479, 3.47755368204], which means we can be 95% confident that the true population mean falls within the range of values provided by this confidence interval. The small breadth of the confidence interval indicates that the population mean was estimated with a high degree of precision. This suggests there is little variability between the observed data points and the underlying mean value.

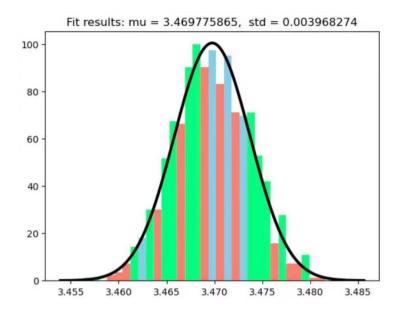


Figure 2: Result Distribution

## 5.1 Convergence of $\theta_n$ and $J(\theta)$

- 1. Left Diagrams:
- $\theta_n$ : Initial  $\theta = 5$ . Sharp decline initially, stabilizing at a lower value.
- $J(\theta)$ : High initial  $\theta$  value ( $J(\theta) \approx 1200$ ). Steep initial decline, stabilizing at a lower value.
  - 2. Middle Diagrams:
- $\theta_n$ : Initial  $\theta \approx 3$ . Significant fluctuations, indicating instability.
- $J(\theta)$ : Mid-range initial  $\theta$  value. Significant oscillations, indicating poor convergence.
  - 3. Right Diagrams:
- $\theta_n$ : Initial  $\theta \approx 3$ . Rapid initial decline, then stabilization.
- $J(\theta)$ : Lower initial  $\theta$  value  $(J(\theta) \approx 60)$ . Rapid initial decline, then stabilization.

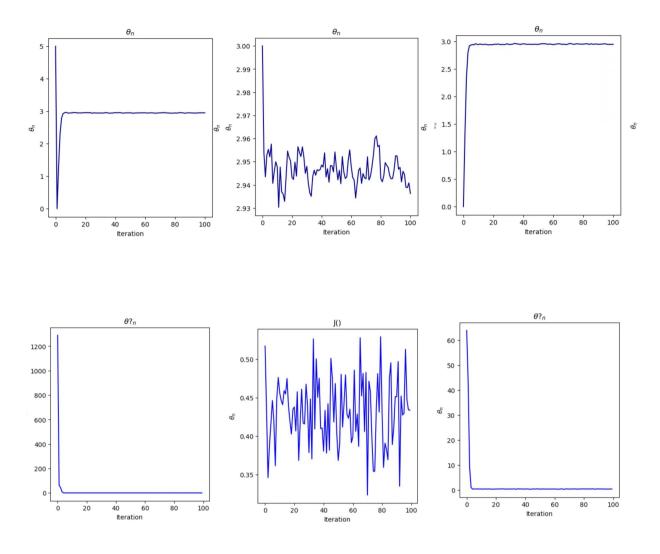


Figure 3: convergence diagrams

### Project 2 Question 2

### 1 Introduction

In the given task, we have to perform the mk = C, optimizing (4) and make  $Wn(\theta)$  approximate to w within a limited computational budget. To achieve an optimal result that converges well and correctly estimates the 90th percentile, we need to perform m iterations for the update of theta and k iterations to calculate  $Wn(\theta)$ . By optimizing the ratio of m to k, hence we can attain the desired outcome.

### 2 Implementation

The second task requires us to design an algorithm to simultaneously perform two types of updates: one for the standard Stochastic Approximation (SA) algorithm to find the optimal average service time  $\theta$ , and another to estimate the quantile  $q_{\theta}$  of the waiting time  $W^{N}(\theta)$  under  $\theta$ . We need to determine a good number of updates m and k, within a fixed budget mk, to improve computational efficiency.

We have to solve optimization problem below:

$$\min_{\theta>0} \mathbb{E}[(q_{\theta}-z)^2],$$

where  $F_{\theta}(w) := \mathbb{P}(W^N(\theta) \leq w)$  is the distribution function of  $W^N(\theta)$ , and  $F_{\theta}$  is a monotonically decreasing function of  $\theta$ .

#### 2.1 Algorithm Design

We applied the following approach to design this algorithm:

- 1. Outer Loop: Update  $\theta$ .
- 2. Inner Loop: For each  $\theta$ , estimate the quantile  $q_{\theta}$  of the waiting time  $W^{N}(\theta)$ . Steps are as below.
- 1. Initialize  $\theta_0$  and  $q_{0,0}$  (initial quantile estimate).
- 2. For each outer loop iteration i: a. Calculate the sequence of waiting times  $\{W_n(\theta_i)\}$  under the current  $\theta_i$ . b. Update  $q_{i,j}$  k times using the inner loop to estimate the new quantile. c. Update  $\theta$  based on the quantile estimate.
  - 3. Return the final estimates of  $\theta$  and  $q_{\theta}$ .

The number of updates for the outer loop is m, and the number of updates for the inner loop is k, with a total budget of mk.

#### 2.2 Choosing m and k

Within the fixed budget mk, we need to allocate m and k efficiently. Typically, the choice of m and k depends on the following factors:

- 1. Convergence Speed: The inner loop k should be large enough to ensure the convergence of the quantile estimate  $q_{\theta}$ .
- 2. Update Frequency: The outer loop m should be large enough to ensure sufficient updates of the  $\theta$  parameter.

We can choose m and k based on experience or experimentation. For example:

- Choose m and k such that each outer update has enough inner iterations to ensure the accuracy of  $q_{\theta}$ . - Ensure that mk equals the fixed budget.

Assuming the total budget is B, we can set m and k as:

$$m = \sqrt{B}, \quad k = \sqrt{B}.$$

This ensures that  $m \times k = B$  and that each update of  $\theta$  has sufficient inner iterations to estimate the quantile.

#### 3 Proof

To prove that the quantile estimate  $q_{\theta}$  updated via the Stochastic Approximation (SA) algorithm will converge to the target quantile p after infinite iterations, we can leverage the convergence theory of the SA algorithm, specifically the Robbins-Monro algorithm's convergence conditions. The detailed information is in 3 between pages 137-139.

## 4 Result and Analysis

Here's the table of experiment data and final figure we choose.

| Items | m    | k     | theta        | q(theta)     | q-8           |
|-------|------|-------|--------------|--------------|---------------|
| 1     | 1000 | 500   | 3. 474143025 | 8.040816327  | 0.040816327   |
| 2     | 707  | 707   | 3. 469593627 | 8.055918367  | 0. 055918367  |
| 3     | 600  | 833   | 3. 47511119  | 8.02755102   | 0. 02755102   |
| 4     | 500  | 1000  | 3. 475664233 | 7. 790380762 | -0. 209619238 |
| 5     | 400  | 1250  | 3. 466433973 | 7. 72375     | -0. 27625     |
| 6     | 300  | 1667  | 3. 469458013 | 7.621333333  | -0. 378666667 |
| 7     | 200  | 2500  | 3. 472145236 | 7. 4085      | -0. 5915      |
| 8     | 150  | 3333  | 3. 486775509 | 8. 043877551 | 0.043877551   |
| 9     | 100  | 5000  | 3. 468914156 | 8.002564103  | 0.002564103   |
| 10    | 75   | 6667  | 3. 467788228 | 7. 998489796 | -0.001510204  |
| 11    | 50   | 10000 | 3. 460175492 | 7.603448276  | -0. 396551724 |
| 12    | 30   | 16667 | 3. 425933218 | 7.5          | -0. 5         |

Figure 4: m=500, k=100000

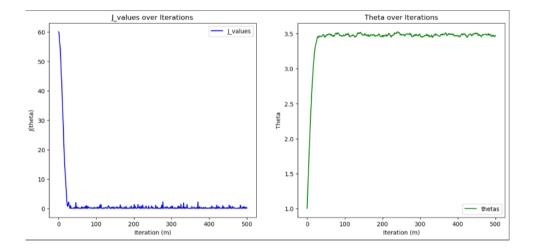


Figure 5: m=500, k=100000

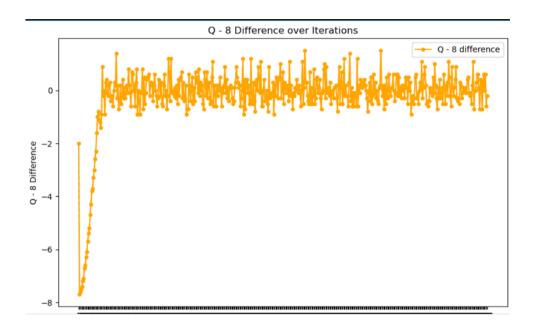


Figure 6: Q-8 Difference Over Iterations

And our final result is m=500, k=10000.

#### 5 Conclusion

For the first task, our objective was to minimize the average waiting time for customers by finding the optimal mean service time  $\theta$  through the application of a Stochastic Approximation (SA) algorithm. The results showed that the optimized  $\theta$  converged to a mean value of approximately 3.4698 with a standard deviation of 0.0040.

For the second task, we used an algorithm to perform simultaneous updates for the average service time  $\theta$  and the quantile estimate  $q_{\theta}$  within a fixed computational budget. By optimizing the ratio of outer (m) and inner (k) loop iterations, we ensured efficient and accurate computation. Our experiments demonstrated that the chosen values of m = 500 and k = 100000 provided optimal results, with  $\theta$  and  $J(\theta)$  converging effectively.

Overall, this project successfully applied stochastic approximation techniques to optimize queueing performance, achieving precise and reliable results with efficient computational resource usage.

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#### Appendix A Code from Project 1

```
import numpy as np
  import matplotlib.pyplot as plt
  def J(p1,p2):# In order to make things easier we decrease the
      amount of the variable: 3->2
       p3=round(1-p1-p2,7) #Get the value of p3
       #Generating the random variables
       eta1 = np.random.normal(0, 1, 10000)
       eta2 = np.random.normal(0, np.sqrt(2), 10000)
       eta3 = np.random.normal(0, np.sqrt(3), 10000)
       w = np.random.exponential(0.3, 10000)
10
       V = np.random.normal(0, 1, 10000)
       #Calculate the Xi
12
       X1 = (0.6 * V + 0.8 * eta1) / np.maximum(w, 1)
13
       X2 = (0.6 * V + 0.8 * eta2) / np.maximum(w, 1)
14
       X3 = (0.6 * V + 0.8 * eta3) / np.maximum(w, 1)
15
       scale = [] #To calculate the sum of three companies
17
       for i in range (10000):
18
           scale1, scale2, scale3 = 0.00
19
           if X1[i]>=2: #To get Xi>=xi
20
               scale1 = round(np.random.uniform(0,X1[i])*p1,7)
21
           if X2[i] > = 3:
               scale2 = round(np.random.uniform(0, X2[i])*p2,7)
23
           if X3[i] > =1:
24
               scale3 = round(np.random.uniform(0, X3[i])*p3,7)
25
           scale.append(round(scale1+scale2+scale3,7))
26
       summary=sum(scale)/10000 #calculate the Sharpe ratio
       std=np.std(scale)
```

```
expect=round (summary/std,9)
29
       return -expect #To find the maxima of the J(theta) could
           be convert to find the minima of -J(tehta)
   def J_graint(J,p1,p2,i): #Use the SPSA algorithm to calculate
      the gradient
       y1, y2=0,0 #gradient from the two directions
32
       r1=np.random.choice([1,-1])
33
       r2=np.random.choice([1,-1])
       theta=1/(i+10000)**0.3 #Using the decrasing theta
       y1 = (J(p1+theta*r1, p2+theta*r2)-J(p1-theta*r1, p2-r2*theta))
36
           /(2*theta*r1)
       y2 = (J(p1+theta*r1, p2+theta*r2)-J(p1-theta*r1, p2-r2*theta))
37
           /(2*theta*r2)
       return y1, y2
   def SA(J_graint, J, theta_list1, theta_list2, number): # The
      stochastic approximation algorithm
       for i in range (number):
40
            step\_size = 10/(i+100) #Uing the decreasing learning
41
               rate
            Jvalue.append(J(theta_list1[i], theta_list2[i]))
42
           y1, y2=J_graint(J, theta_list1[i], theta_list2[i], i)
43
            suml=theta_list2 [i]+theta_list1 [i]
44
            if suml>1: #In this case we do the projection
45
                theta_list2[i] = theta_list2[i] - (suml-1)/2
46
                theta_list1 [i]=theta_list1 [i]-(suml-1)/2
47
            theta_list1.append(\max(\text{theta_list1}[i]-\text{step_size}*y1,0))
48
            theta_list2.append(\max(\text{theta_list2}[i]-\text{step_size}*y2,0))
49
       return theta_list1, theta_list2
50
51
   p1_list = [] #It records the last 500 experiments' mean of p1
   p2_list = [] #It records the last 500 experiments' mean of p2
```

```
Experiment_number=4#The total experiment times
   for num in range (Experiment_number):
       Jvalue = [] #To record the J(theta) values
56
       number=3000 #iteration times
57
       theta_list1 = [0.5] #The starting point of p1
58
       theta_list2 = [0.5] #The starting point of p2
59
       pls, p2s=SA(J_graint, J, theta_list1, theta_list2, number)
60
       p1s=np.array(p1s[-501:-1]) #Get the final 500 points
       p2s=np. array (p2s[-501:-1])
       mean_p1 = np.mean(p1s) #Get the mean of the converge data
63
       mean_p2 = np.mean(p2s)
64
       pl_list.append(np.mean(pls))
65
       p2_list.append(np.mean(p2s))
       file_path1 = r"point.txt"
       with open(file_path1, 'a') as f:
68
            f.write(f'Experiment \{num\}: \text{-Mean \p1} = \{mean_p1\}, \text{-Mean \cdot}
69
               p2 = {\text{mean}_p2} \ n' #Write the p1 and p2 to a txt
               file
       fig, axs = plt.subplots (1,3, figsize = (20,5))
       axs[0].plot(theta_list1, color="green")
71
       axs[0].set_xlabel("Iteration")
72
       axs[0].set_ylabel(r"\$\theta1_n\$")
73
       axs[0].set_title(r"\$\theta1_n\$")
74
       axs[1].plot(theta_list2, color="blue")
75
       axs[1].set_xlabel("Iteration")
       axs[1].set_ylabel(r"\$\theta2_n\$")
77
       axs[1].set_title(r"\$\theta2_n\$")
78
       axs[2].plot(Jvalue, color="red")
79
       axs[2].set_xlabel("Iteration")
80
       axs[2].set_ylabel("J-value")
81
       axs [2]. set_title ("The-J-function")
```

```
save_path = f"{num}.png"

fig.savefig(save_path) #Get the diagram

plt.close(fig)
```

#### Appendix B Code from Project 2.1

```
import numpy as np
   def J(theta,z):
       #np.random.seed(10)
       s=0
       d=0
       for k in range (100):
            W_{mean} = []
            for j in range (700):
                A=np.random.exponential(5, 12)
                 S=list (np.random.exponential(theta, 12))
                 S. insert (0,0)
11
                W = [0]
12
                 for i in range (11):
13
                     W. append (\max(0, W[i] + S[i] - A[i+1]))
14
                 W_{mean.append(np.mean(W[1:11]))
15
            W_mean.sort()
16
            l=W_mean[int(0.9*len(W_mean))]
17
            s + = (1 - z) * * 2
18
            d+=l-z
19
       return s/100,d/100
20
   theta = []
   def SA(J, theta_list, z, number):
       for i in range(number):
24
            theta = 10/(i+100)
25
```

```
#theta_list.append(max(0, theta_list[i]-SPSA(J,
26
               theta_list[i],z)*theta))
            s, d=J(theta_list[i], z)
27
            theta_list.append(\max(0, \text{theta_list}[i]-2*(d)*\text{theta}))
28
            Jvalue.append(s)
20
       return theta_list [-1]
30
31
   for i in range (500):
       Jvalue=[]
33
       number=20
34
       z=8
35
       theta_list = [1]
36
       SA(J, theta_list, z, number)
37
       theta.append(np.mean(theta_list[15:-1]))
39
   plt.figure(1)
  mu = np.mean(theta)
   std = np.std(theta)
   number_of_bins=40
   bins = np.linspace(mu-4*std, mu+4*std, number_of_bins)
   n, bins, patches = plt.hist(theta, bins, density=True)
46
   def true_value(i):
       return norm.pdf(((mu-4*std+i*2*4*std/(number_of_bins-1))
48
          +2*4*std/(2*(number_of_bins_{-1})), mu, std)
49
   for i in range (number_of_bins -1):
       if n[i] < true_value(i):</pre>
                                           # Give bar red color
51
           if lower than PDF
            patches [i]. set_fc('salmon')
52
```

```
if n[i] > true_value(i):
                                        # Give bar green color
           if higher than PDF
               patches[i].set_fc('springgreen')
54
       if n[i] \le 0.05*n[i] + true\_value(i) and n[i] >= -0.05*n[i]
          | + true_value(i):
           patches[i].set_fc('skyblue') # Give bar blue color
              if approx. equal to PDF
57
  # Plotting the PDF of the normal distribution
  x = np. linspace (mu-4*std, mu+4*std, 1000)
  y = norm.pdf(x, mu, std)
                                         # PDF: 1/((2)) * exp
     (-(x-)^2/2^2)
  plt.plot(x, y, lw=3, color='black', label = 'PDF')
  # print(" mean theta_n = ", mu, "standard deviation theta_n"
     , std)
  title = "Fit-results: mu-=-%.9f, --std-=-%.9f" % (mu, std)
  plt.title(title)
  #plt.title("SA experiment")
  plt.show()
68
  import matplotlib.pyplot as plt
  fig, axs = plt.subplots (1,2, \text{ figsize} = (10,5))
  # Plot the iterate
  axs[0].plot(theta_list, color="darkblue")
  axs[0].set_xlabel("Iteration")
  axs[0].set_ylabel(r"\$\theta_n\$")
  axs[0].set_title(r"\$\theta_n\$")
  axs[1].plot(Jvalue, color="blue")
  axs[1].set_xlabel("Iteration")
  axs[1].set_ylabel(r"\$\theta_n\$")
```

```
_{79} axs[1].set_title('J()')
```

#### Appendix C Code from Project 2.2

```
def J_function(theta0,q,k):
       #np.random.seed(20)
        s=0
        d=0
        z=8
        alpha=0.1
       w = 0.9
       A=np.random.exponential(5,20002)
        S=np.random.exponential(theta0,20002)
        S=list(S)
        S.insert(0,0)
11
       W = [0]
12
       W_{mean} = []
13
        for o in range (20001):
14
            W. append (\max(0, W[o] + S[o] - A[o+1]))
15
        W_{\text{-}}mean.append(np.mean(W[1:20001]))
16
        W_mean.sort()
        l=np.mean(W_mean[int(0.9*len(W_mean))-10:int(0.9*len(W_mean))]
18
           W_{mean}) + 10
        s+=(1-z)**2
19
        d+=l-z
20
        for j in range(k): #k times inner loop
21
             for ww in W_mean:
                 if(ww \leq q):
23
                      q=alpha*(1-w)
24
                 else:
25
                      q=alpha*(0-w)
26
```

```
return s,d,q
28
29
   import matplotlib.pyplot as plt
30
   import numpy as np
31
   theta0=1
   q=6
   z=8
  w = 0.9
  m = 1000
   k = 1000
   alpha=0.1
   learning_rate=0.01
   thetas=[theta0]
   J_{\text{values}} = [J_{\text{function}}(\text{theta0}, z, k) [0]]
   q = list = [q]
   l_list = []
   for i in range (m):
       s,d,q=J_function(theta0,q,k)
       q_list.append(q)
       \#theta0=theta0 -2*(10/(i+100))*d
47
       theta0 = theta0 - 2*0.01*d
48
       thetas.append(theta0)
       J_values.append(s)
51
   plt. figure (figsize = (14, 6))
   plt.subplot(1, 2, 1)
   plt.plot(J_values, label='J_values', color='blue')
   plt.xlabel('Iteration (m)')
  plt.ylabel('J(theta)')
   plt.title('J_values over Iterations')
```

```
plt.legend()
plt.subplot(1, 2, 2)
plt.plot(thetas, label='thetas', color='green')
plt.xlabel('Iteration - (m)')
plt.ylabel('Theta')
plt.title('Theta-over-Iterations')
plt.legend()
differences = [q - 8 \text{ for } q \text{ in } q \text{-list}]
plt. figure (figsize = (10, 6))
plt.plot(range(1, len(q_list) + 1), differences, label='Q--8-
   difference', color='orange', marker='o', linestyle='-',
   markersize=4)
for i in range(1, len(q_list)):
    plt.plot([i, i+1], [differences[i-1], differences[i]],
        color='orange', linestyle='--')
plt.legend()
plt.title('Q--8-Difference-over-Iterations')
plt.xticks(range(1, len(q_list) + 1))
plt.ylabel('Q--8-Difference')
plt.show()
```