# OREISTEIN\_Pierre\_TP3

December 15, 2018

# 1 TP 3: Reinforcement Learning with Function Approximation

In []:

# 2 0 - Packages

```
In [1]: # Mathematical packages
   import numpy as np

# Matplotlib packages
   %matplotlib inline
   import matplotlib.pyplot as plt
   from mpl_toolkits.mplot3d import Axes3D
   import seaborn as sns
   sns.set()

# Import tqdm as tqdm
   import tqdm as tqdm
   import tqdm as tqdm

# Local packages
   import lqg1d
   import utils
```

# 3 1 - On-Policy Reinforcement Learning with Parametric Policy

### 3.1 1.1 - Guassian policy model

```
In [2]: class gaussianPolicy(object):
    """Define a gaussian policy with parameter theta and sigma."""

def __init__(self, theta, sigma=0.4, sigma_constant=True):
    """Initialisation of the policy."""
    self.theta = theta
    self.sigma = sigma
    self.sigma_constant=sigma_constant
```

```
def mu_theta(self, s):
   """Definition of the mu fonction according to theta."""
   return self.theta * s
def d_mu_theta(self, s):
   """Definition of the differentiate of mu_theta."""
def sigma_w(self, s):
   """Definiton of the sigma for our model."""
   if self.sigma_constant:
       return self.sigma
   else:
       return np.sqrt((self.sigma * s) ** 2)
def d_sigma_w(self, s):
   """Definition of the differentiate of sigma_w for our model."""
   if self.sigma_constant:
       return 0
   else:
       return self.sigma
def draw_action(self, s):
   """Return a sampled action according to the gaussian policy at state s."""
   # Draw a sample of the gaussian policy
   result = np.random.normal(self.mu_theta(s), self.sigma_w(s))
   return result
def gradientGaussianPolicy(self, a, s):
   """Return the log-gradient of the gaussian policy taken in a and s."""
   # Initialisation of the gradient
   log_g = np.zeros((2,1))
   # Computation of the gradient in theta
   log_g[0] = (a - self.mu_theta(s)) / (self.sigma_w(s) ** 2) * self.d_mu_theta(s)
   # Computation of the gradient in w
   log_g[1] /= (self.sigma_w(s) ** 3)
   return log_g
```

#### 3.2 1.2 - Adam - Step Optimization

# Shape

n, n\_iter = np.shape(array)

# Compute the emprical mean

# Compute the empirical variances
emp\_variance = np.std(array, axis=0)

```
In [3]: class AdamStep(object):
            def __init__(self, d=2, alpha=0.001, beta_1=0.9, beta_2=0.999, epsilon=10e-8):
                self.alpha = alpha
                self.beta_1 = beta_1
                self.beta 2 = beta 2
                self.epsilon = epsilon
                self.m_t = 0
                self.v_t = 0
                self.t = 0
            def update(self, g_t):
                """Execute one step of the Adam algorithm.
                   cf https://arxiv.org/pdf/1412.6980.pdf"""
                self.t = self.t + 1
                self.m_t = self.beta_1 * self.m_t + (1 - self.beta_1) * g_t
                self.v_t = self.beta_2 * self.v_t + (1 - self.beta_2) * g_t * g_t
               hat_m_t = self.m_t / (1 - self.beta_1 ** self.t)
                hat_v_t = self.v_t / (1 - self.beta_2 ** self.t)
                step = self.alpha * hat_m_t / (np.sqrt(hat_v_t) + self.epsilon)
                return step
3.3 1.3 - Constant Step
In [4]: class ConstantStep(object):
            """Define a step optimisation class with a constant learning rate."""
            def __init__(self, learning_rate):
                self.learning_rate = learning_rate
            def update(self, gt):
                return self.learning_rate * gt
3.4 1.4 - REINFORCE Algorithm
In [5]: def computeCIBounds(array):
            """Compute the empirical 95% CI bounds."""
```

```
emp_mean = np.mean(array, axis=0)
            # Compute the bounds
            low_CI = emp_mean - 1.96 * emp_variance / np.sqrt(n)
            high_CI = emp_mean + 1.96 * emp_variance / np.sqrt(n)
            return low_CI, high_CI
In [6]: def REINFORCE(env, Policy, n_experiences=5, n_itr=100, N=100, T=100,
                      Step=ConstantStep(10e-5), sigma_w=0.4, addOptimism=None):
            """Executes the REINFORCE algorithm on env."""
            # Saving array for the aug of mean_parameters and aug_return
            mean_parameters = np.zeros((n_experiences, n_itr))
            avg_return = np.zeros((n_experiences, n_itr))
            for j in tqdm.tqdm(range(n_experiences)):
                # Intialisation of theta_k
                theta_k = np.zeros((2,1)) - 0.1
                # Saving array for the parameters theta and for the return
                parameters = []
                r = []
                \# Executes n_itr for optimising theta
                for i in range(n_itr):
                    # Update of the current policy
                    policy = Policy(theta_k[0, 0], sigma=sigma_w)
                    # Executes all the trajectories
                    paths = utils.collect_episodes(env, policy=policy, horizon=T, n_episodes=N
                    # Computation of the gradient step
                    g_t = np.zeros((2, 1))
                    # Initialisation of the avg_return
                    r.append(0)
                    for episode in paths:
                        # Extract the actions taken
                        actions = episode["actions"]
                        states = episode["states"]
                        rewards = episode["rewards"]
                        # Update the rewards
```

```
rewards = addOptimism(env, actions, states, rewards)
            # Array of the discounts
            discounts_1 = np.array([discount ** t for t in range(len(rewards))])
            for t in range(len(states)):
                # Extract current state
                a_t = actions[t, 0]
                s_t = states[t, 0]
                # Sum all the future rewards
                R_t = np.sum(rewards[t:] * discounts_l[t:] * discount ** (-t))
                # Computation of the gradient of the log
                g_log_pi_t = policy.gradientGaussianPolicy(a_t, s_t)
                # Add it to g_t
                g_t += g_log_pi_t * R_t
                # Update the avg_return
                if t == 0:
                    r[i] += R_t / N
        # Average of the gradient over the episodes
        g_t = g_t / N
        # Make an iteration of the gradient descent
        theta_k = theta_k + Step.update(g_t)
        # update of the mean parameter
        parameters.append(theta_k[0,0])
    # Update of mean parameters
    parameters = np.array(parameters)
    # Update of mean_parameters_l and avg_return_l
    mean_parameters[j,:] = parameters
    avg_return[j,:] = r
# Compute the CI bounds
mean_parameters_CI_low, mean_parameters_CI_high = computeCIBounds(mean_parameters)
avg_return_CI_low, avg_return_CI_high = computeCIBounds(avg_return)
# Normalisation of mean_parameters_l and avg_return_l
mean_parameters = mean_parameters.mean(axis=0)
avg_return = avg_return.mean(axis=0)
```

if addOptimism != None:

```
return [mean_parameters, avg_return, mean_parameters_CI_low, mean_parameters_CI_his avg_return_CI_low, avg_return_CI_high]
```

#### 3.5 1.5 - Graphic function

```
In [7]: def lineplotCI(ax, x_data, y_data, sorted_x, low_CI, upper_CI, x_label, y_label, title
            """Plot the points with the confidence interval"""
            # Plot the data, set the linewidth, color and transparency of the
            # line, provide a label for the legend
            ax.plot(x_data, y_data, lw=1, color='#539caf', alpha=1, label='Mean')
            # Shade the confidence interval
            ax.fill_between(sorted_x, low_CI, upper_CI, color='#539caf', alpha=0.4, label='95%
            # Label the axes and provide a title
            ax.set_title(title)
            ax.set_xlabel(x_label)
            ax.set_ylabel(y_label)
            # Display legend
            ax.legend(loc = 'best')
In [10]: def figures (mean_parameters, avg_return, mean_parameters_CI_low, mean_parameters_CI_h
                     avg_return_CI_low, avg_return_CI_high):
             """Display the two figures required."""
             fig, axs = plt.subplots(1, 2, figsize=(20, 8))
             plt.grid(True)
             # plot the average return obtained by simulating the policy
             # at each iteration of the algorithm (this is a rought estimate
             # of the performance
             lineplotCI(axs[0],
                        x_data=np.linspace(1, n_itr, n_itr),
                        y_data = avg_return,
                        sorted_x = np.linspace(1, n_itr, n_itr),
                        low_CI = avg_return_CI_low,
                        upper_CI = avg_return_CI_high,
                        x_label = 'Number of iteration of the gradient ascent',
                        y_label = 'Average reward',
                        title = 'Average reward over the gradient ascent')
             # plot the distance mean parameter
             # of iteration k
             # Call the function to create plot
             lineplotCI(axs[1],
```

```
x_label = 'Number of iteration of the gradient ascent',
                  y label = 'Theta',
                  title = 'Value of theta over the gradient ascent')
         plt.show()
3.6 1.6 - Results
# Define the environment and the policy
      env = lqg1d.LQG1D(initial_state_type='random')
      # Experiments parameters
      # We will collect N trajectories per iteration
      # Each trajectory will have at most T time steps
      T = 100
      # Number of policy parameters updates
      n_{itr} = 100
      # Set the discount factor for the problem
      discount = 0.9
      # Learning rate for the gradient update
      learning_rate = 10e-5
      # define the update rule (stepper)
      Step = ConstantStep(learning_rate) #AdamStep() # e.g., constant, adam or anything you
# REINFORCE
      # fill the following part of the code with
      # - REINFORCE estimate i.e. gradient estimate
      # - update of policy parameters using the steppers
      # - average performance per iteration
      \# - distance between optimal mean parameter and the one at it k
      # Number of experiences
      n_{experiences} = 5
      # REINFOIRCE algoritm
      [mean_parameters, avg_return,
```

x\_data = np.linspace(1, n\_itr, n\_itr),

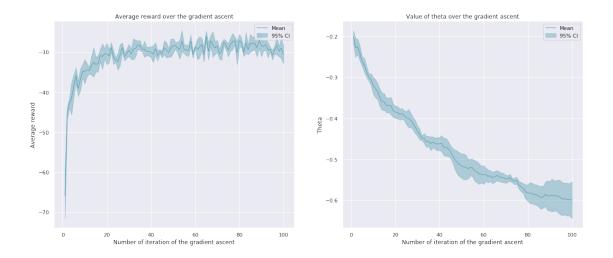
sorted\_x = np.linspace(1, n\_itr, n\_itr),

y\_data = mean\_parameters,

low\_CI = mean\_parameters\_CI\_low,
upper CI = mean parameters CI high,

# Figures

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**Commentaries**: \* As initialisation for  $\theta$  we choose  $\theta = -0.1$ . As it is not so far from the optimal  $\theta^*$  it allows to converge easier. \* As the gradient becomes really big at the first step we choose a small learning rate; we chose  $\alpha_t = 10^{-5}$ . The other parameters are left unchanged. \* For the confidence intervals, we use the empirical mean and empirical variance over the different experiences run with 100 optimisation iterations. Here, we use only 5 experiences which is really low but it was already computationally costly

#### 3.7 1.7 - Influence of the parameters

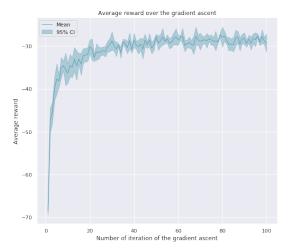
#### 3.7.1 1.7.1 Influence of N

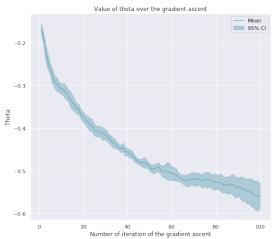
```
In [10]: ### High N
    N = 200

# REINFOIRCE algoritm
[mean_parameters, avg_return,
    mean_parameters_CI_low, mean_parameters_CI_high,
    avg_return_CI_low, avg_return_CI_high] = REINFORCE(env, gaussianPolicy, N=N)
```

#### # Figures

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#### In [11]: ### Low N N = 5

#### # REINFOIRCE algoritm

[mean\_parameters, avg\_return,
mean\_parameters\_CI\_low, mean\_parameters\_CI\_high,
avg\_return\_CI\_low, avg\_return\_CI\_high] = REINFORCE(env, gaussianPolicy, N=N)

#### # Figures

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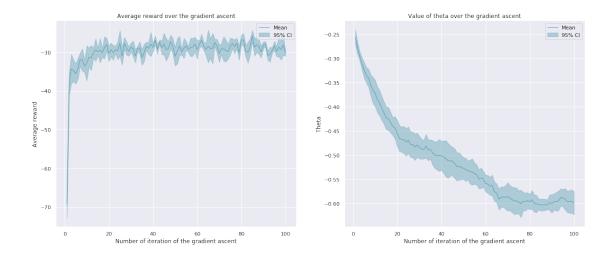
**Commentaries**: \* We can observe that N influences mainly the variance of the convergence of both the average reward and the mean theta. It is logic because more N is important, more trajectories are extracted. Therefore, the approximation of the expectation of the current policy is better when N is big and so better is the approximation of  $\nabla J_{theta}^{\pi_{\theta}}$ . Hence, it is easier to converge and there is less variance in the results.

#### 3.7.2 1.7.2 - Influence of T

```
In [51]: ### High T
       T = 300
       # REINFOIRCE algoritm
       [mean_parameters, avg_return,
        mean_parameters_CI_low, mean_parameters_CI_high,
        avg_return_CI_low, avg_return_CI_high] = REINFORCE(env, gaussianPolicy, T=T)
       # Figures
       figures (mean_parameters, avg_return, mean_parameters_CI_low, mean_parameters_CI_high,
              avg_return_CI_low, avg_return_CI_high)
 0%1
            | 0/5 [00:00<?, ?it/s]
20%1
           | 1/5 [02:47<11:10, 167.70s/it]
40%1
         | 2/5 [05:34<08:22, 167.36s/it]
60%1
       | 3/5 [08:20<05:34, 167.07s/it]
```

```
80%| | 4/5 [11:08<02:47, 167.22s/it]
```

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```
In [52]: ### High T
T = 25
```

#### # REINFOIRCE algoritm

[mean\_parameters, avg\_return,

 ${\tt mean\_parameters\_CI\_low, mean\_parameters\_CI\_high,}$ 

avg\_return\_CI\_low, avg\_return\_CI\_high] = REINFORCE(env, gaussianPolicy, T=T)

#### # Figures

#### 

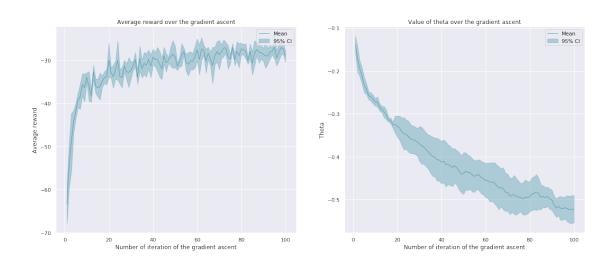
0%| | 0/5 [00:00<?, ?it/s]

20%| | 1/5 [00:14<00:56, 14.16s/it]

40%| | 2/5 [00:28<00:42, 14.32s/it]

60%| | 3/5 [00:42<00:28, 14.23s/it]

80%| | 4/5 [00:57<00:14, 14.23s/it]



**Commentaries**: \* With these two graphics on the influence of T, we can remark that when T is bigger, the convergence is quicker. It seems quite logic because when T is bigger, better is our approximation of the expectation of the reward of our current policy. However, when T is bigger, we have also more chance to have bias because we use the sub-trajectories of the trajectories but it seems that it is not the case here.

#### 3.7.3 1.7.3 - Influence of sigma\_w

figures(mean\_parameters, avg\_return, mean\_parameters\_CI\_low, mean\_parameters\_CI\_high, avg\_return\_CI\_low, avg\_return\_CI\_high)

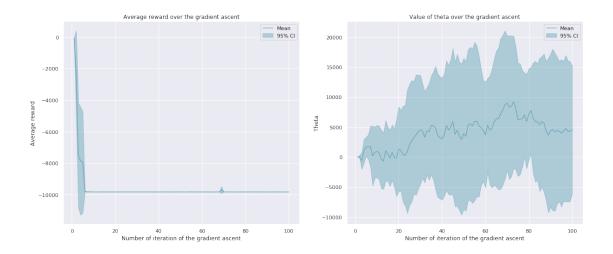
```
0%| | 0/5 [00:00<?, ?it/s]
20%| | 1/5 [00:55<03:41, 55.47s/it]
```

```
40%| | 2/5 [01:50<02:46, 55.45s/it]

60%| | 3/5 [02:46<01:50, 55.38s/it]

80%| | 4/5 [03:40<00:55, 55.19s/it]

100%|| 5/5 [04:36<00:00, 55.39s/it]
```



#### # REINFOIRCE algoritm

[mean\_parameters, avg\_return,
mean\_parameters\_CI\_low, mean\_parameters\_CI\_high,
avg\_return\_CI\_low, avg\_return\_CI\_high] = REINFORCE(env, gaussianPolicy, sigma\_w=si

#### # Figures

#### 

0%| | 0/5 [00:00<?, ?it/s]

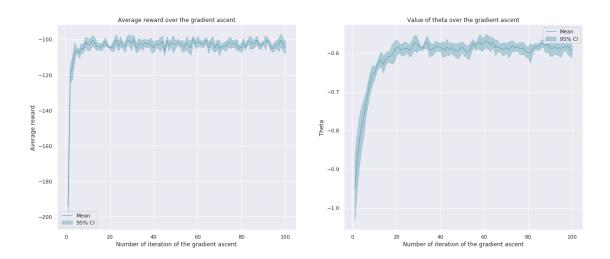
20%| | 1/5 [00:55<03:40, 55.24s/it]

40%| | 2/5 [01:52<02:47, 55.91s/it]

```
60%| | 3/5 [02:49<01:52, 56.16s/it]

80%| | 4/5 [03:46<00:56, 56.36s/it]

100%|| 5/5 [04:41<00:00, 56.15s/it]
```



**Commentaries**: \* With these two graphics we understand better the influence of sigma\_w when it is constant. For the both cases, we have considered sigma\_w constant but with different value. In the first case, sigma\_w is equal to 0.01. In the second case it is equal to 2.5. When sigma\_w is low it means that we explore less the space of action and hence of state. So, it seems that by itself our policy does not explore the space of state-action. Therefore it is impossible to converge. When sigma\_w is higher, we explore more and we converges much faster. \* Also, even if we converge much faster towards the optimal  $\theta^*$ , we can observe that the reward is lower because of this high exploration.

#### 3.7.4 1.7.5 - Influence of the learning rate

```
In [65]: # Learning rate for the gradient update
    learning_rate = 0.1

# define the update rule (stepper)
Step = ConstantStep(learning_rate)

# REINFOIRCE algoritm
[mean_parameters, avg_return,
    mean_parameters_CI_low, mean_parameters_CI_high,
    avg_return_CI_low, avg_return_CI_high] = REINFORCE(env, gaussianPolicy, Step=Step)
```

#### # Figures

#### 

```
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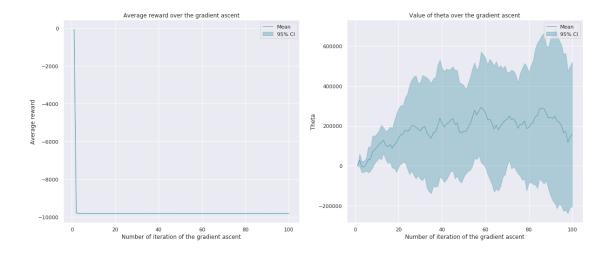
20%| | 1/5 [00:58<03:53, 58.29s/it]

40%| | 2/5 [01:55<02:53, 57.90s/it]

60%| | 3/5 [02:51<01:54, 57.35s/it]

80%| | 4/5 [03:47<00:56, 56.97s/it]

100%|| 5/5 [04:43<00:00, 56.85s/it]
```



```
In [58]: # Learning rate for the gradient update
    learning_rate = 10e-10

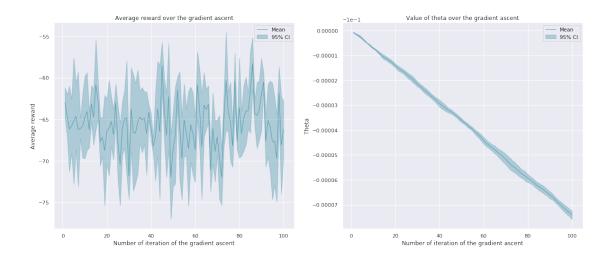
# define the update rule (stepper)
Step = ConstantStep(learning_rate)

# REINFOIRCE algoritm
[mean_parameters, avg_return,
    mean_parameters_CI_low, mean_parameters_CI_high,
    avg_return_CI_low, avg_return_CI_high] = REINFORCE(env, gaussianPolicy, Step=Step)
```

#### 

#### # Figures

#### 



#### # REINFOIRCE algoritm

[mean\_parameters, avg\_return,
 mean\_parameters\_CI\_low, mean\_parameters\_CI\_high,
 avg\_return\_CI\_low, avg\_return\_CI\_high] = REINFORCE(env, gaussianPolicy, Step=Step)

#### # Figures

#### 

figures (mean\_parameters, avg\_return, mean\_parameters\_CI\_low, mean\_parameters\_CI\_high, avg\_return\_CI\_low, avg\_return\_CI\_high)

```
0%| | 0/5 [00:00<?, ?it/s]

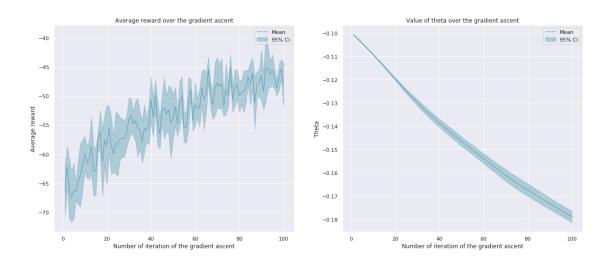
20%| | 1/5 [00:57<03:51, 57.83s/it]

40%| | 2/5 [01:55<02:52, 57.65s/it]

60%| | 3/5 [02:53<01:55, 57.94s/it]

80%| | 4/5 [03:50<00:57, 57.61s/it]

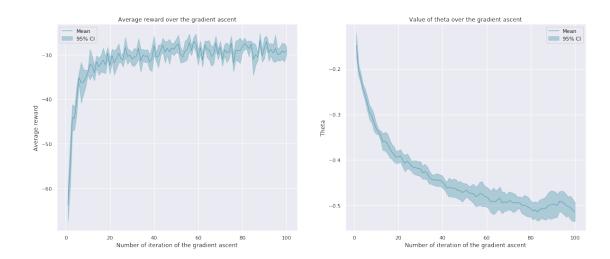
100%|| 5/5 [04:47<00:00, 57.48s/it]
```



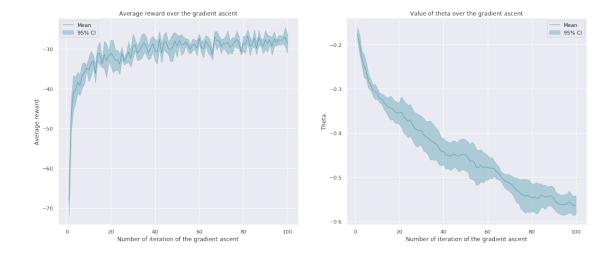
**Commentaries**: \* With these three past graphics we can understand better the influence of the learning rate  $\alpha_t$ . When  $\alpha_t$  is big as in the figure 1, we are going to do big steps but the convergence is not stable (we can observe a big variance) and so we do not converge in practice. When  $\alpha_t$  is low as in Figure 2, we can observe that the convergence is really stable (a very small variance). However as we do very small steps, the convergence is very slow. Finally in the last graphic we implement the Adam Step. Here, we can observe we have a balance between the two past solutions. However, among all our simulations, it appears that  $\alpha_t = 10^{-5}$  is the best trade-off between speed of convergence and stability.

#### 3.8 1.8 - Exploration in Policy Gradient

```
In [12]: def addOptimism(env, actions, states, rewards, beta=0, n_bins_A=20, n_bins_S=10):
            """This function adds optimism to rewards."""
            # Initialisation of N(\phi)
            N_phi_t = np.zeros((n_bins_S + 1, n_bins_A + 1), dtype=int)
            for i in range(len(actions)):
                # Extract action and state
                a_t = np.clip(actions[i, 0], -env.max_action, env.max_action)
                s_t = np.clip(states[i, 0], -env.max_pos, env.max_pos)
                # Extract the revelant bin of (s_t, a_t)
               bin_id_a_t = int((a_t + env.max_action) // (2 * env.max_action) * n_bins_A)
               bin_id_s_t = int((s_t + env.max_pos) // (2 * env.max_pos) * n_bins_S)
                # Update the counter
               N_phi_t[bin_id_s_t, bin_id_a_t] += 1
                # Update reward
               rewards[i] += beta * 1 / (np.sqrt(N_phi_t[bin_id_s_t, bin_id_a_t]))
            return rewards
In [13]: # n_exploration
        n_{experiences} = 5
        # Optimism
        def optimism(env, actions, states, rewards, beta=20):
            return addOptimism(env, actions, states, rewards)
        # REINFOIRCE algoritm
        [mean_parameters, avg_return,
         mean_parameters_CI_low, mean_parameters_CI_high,
         avg_return_CI_low, avg_return_CI_high] = REINFORCE(env, gaussianPolicy,
                                                         n_experiences=n_experiences,
                                                         addOptimism=optimism)
        # Figures
        figures(mean_parameters, avg_return, mean_parameters_CI_low, mean_parameters_CI_high,
                avg_return_CI_low, avg_return_CI_high)
100%|| 5/5 [05:21<00:00, 64.22s/it]
```



```
In [12]: \# n_{exploration}
       n_{experiences} = 5
       # Optimism
       def optimism(env, actions, states, rewards, beta=0):
           return addOptimism(env, actions, states, rewards)
       # REINFOIRCE algoritm
       [mean_parameters, avg_return,
        mean_parameters_CI_low, mean_parameters_CI_high,
        avg_return_CI_low, avg_return_CI_high] = REINFORCE(env, gaussianPolicy,
                                                    n_experiences=n_experiences,
                                                    addOptimism=optimism)
       # Figures
       figures (mean_parameters, avg_return, mean_parameters_CI_low, mean_parameters_CI_high,
              avg_return_CI_low, avg_return_CI_high)
100%|| 5/5 [05:23<00:00, 64.68s/it]
```



**Commentaries:** \* The first graphic is obtained with  $\beta = 20$  and the second one with  $\beta = 0$ . Thanks to these two graphic we understand its influence on the exploration. In fact, when  $\beta$  is big as in the first graphic, we explore more and so the variance is reduced between the experiences because all of them have more chance to explore the same domain of the space state-action. However, we contast also that the convergences towards the optimal  $\theta$  is slower precisely because of this exploration.

# 4 2 - Off-Policy Reinforcement Learning with Value Function Approximation

#### 4.1 2.0 - Definition of the environment

#### 4.2 2.1 - Definition of the behavorial policy

```
In [17]: class behavorialPolicy(object):
    """Define a behavorial policy which explores the space of actions and states
    uniformly."""

def __init__(self, actions):
    """Initialisation of the policy."""
    self.actions = actions

def draw_action(self, s):
    """Return a sampled action according to the uniform policy at state s."""

# Draw a sample of the gaussian policy
    result = np.random.choice(self.actions)

return result
```

#### 4.3 2.2 - Generation of the dataset

#### 4.4 2.3 - Definition of FQI

#### 4.4.1 2.3.1 - Auxiliary functions

```
In [19]: def index(value, Array):
          """Return the index of value in Array."""
```

```
In [20]: def phi(S=states, A=actions):
             """Definition of the feature fonction."""
             # Parameter
             n S = len(S)
             n_A = len(A)
             # Initialisation of the resulting array
             result = np.zeros((n_S, n_A, 3))
             # Compute the result of phi for each pair of (state, action)
             for i, s in enumerate(S):
                 for j, a in enumerate(A):
                     # Update result
                     result[i, j, :] = np.array([a, s * a, s ** 2 + a ** 2])
             return result
In [21]: def phiCouple(D):
             """Compute the feauture value over the couples in D."""
             # Parameter
             n = len(D)
             # Initialisation of the resulting array
             result = np.zeros((n, 3))
             # Compute the result of phi for each pair of (state, action)
             for i, couple in enumerate(D):
                 \# Extract s_i and a_i
                 s = couple[0]
                 a = couple[1]
                 # Update result
                 result[i, :] = np.array([a, s * a, s ** 2 + a ** 2])
             return result
In [22]: def Q_max(s, theta_k, A=actions, r_max=1000, discount=0.9):
             """Return the maximum value of Q among A for the given state and theta."""
```

# Index of value

return i

i, = np.where(Array == value)

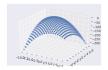
```
# Computation of Q
             phi_matrix = phi([s], A).reshape((len(A), 3))
             Q = np.dot(phi_matrix, theta_k).reshape(-1)
             # Return the maximum
             value = np.clip(np.max(Q), -r_max / (1 - discount), r_max / (1 - discount))
             return value
4.4.2 2.3.2 - Definition of the FOI function
```

```
In [23]: def fqi(env, D n, Phi_Mat, S=S, discount=0.9, n=50, lmbda=0.01):
             """Executes the fitted Q-iteration algorithm."""
             # Computation of number of iterations K
             K = len(D_n) // n
             # Initialisation of theta and phi
             theta_k = np.zeros((3, 1))
             # Saving array of all the theta
             theta_l = [theta_k]
             # Iteration of optimisation
             for k in range(K):
                 # Extract n samples
                 \# D_batch = D_n[n * k: n*(k + 1),:]
                 idx = np.random.randint(len(D_n), size=n)
                 D_batch = D_n[idx, :]
                 # Build testing set
                 y_t = np.zeros((len(D_batch), 1))
                 for i in range(len(D_batch)):
                     # Extract a_i, s_i
                     s_i = D_batch[i, 0]
                     a_i = D_batch[i, 1]
                     \# Sample s_i' and r_i
                     r_i = -np.dot(s_i, np.dot(env.Q, s_i)) -
                           np.dot(a_i, np.dot(env.R, a_i))
                     s_i_prime = np.clip(np.dot(env.A, s_i) + np.dot(env.B, a_i),
                                         -env.max_pos, env.max_pos)[0, 0]
                     # Compute y_i
                     y_i = r_i + discount * Q_max(s_i_prime, theta_k)
```

```
# Append the training set
                   y_t[i, :] = y_i
               # Resoltuion of the least square problem
               Z_t = phiCouple(D_batch)
               inv = np.linalg.inv(np.dot(Z_t.T, Z_t) + lmbda * np.eye(3))
               theta_k = np.dot(inv, np.dot(Z_t.T, y_t))
               theta l.append(theta k)
           return theta_1
In [24]: class greedyPolicy(object):
            """Define the greedy policy according to the given theta."""
           def __init__(self, theta, A=actions):
               """Initialisation of the policy."""
               self.actions = A
               self.theta = theta
           def draw_action(self, s):
               """Return a sampled action according to the greedy policy at state s."""
               # Draw of the greedy policy
               phi_matrix = phi([s], self.actions).reshape((len(self.actions), 3))
               Q = np.dot(phi_matrix, theta_k).reshape(-1)
               sample = self.actions[np.argmax(Q)]
               return sample
4.5 2.4 - Results
# Show the optimal Q-function
         K, cov = env.computeOptimalK(discount), 0.001
         print('Optimal K: {} Covariance S: {}'.format(K, cov))
         Q_fun_ = np.vectorize(lambda s, a: env.computeQFunction(s, a, K, cov, discount, 1))
         Q_{fun} = lambda X: Q_{fun}(X[:, 0], X[:, 1])
         Q_opt = Q_fun(SA)
         # Figure
         fig = plt.figure()
         ax = fig.add_subplot(111, projection='3d')
         ax.scatter(S, A, Q_opt)
```

```
ax.mouse_init()
plt.show()
```

Optimal K: [[-0.58840335]] Covariance S: 0.001

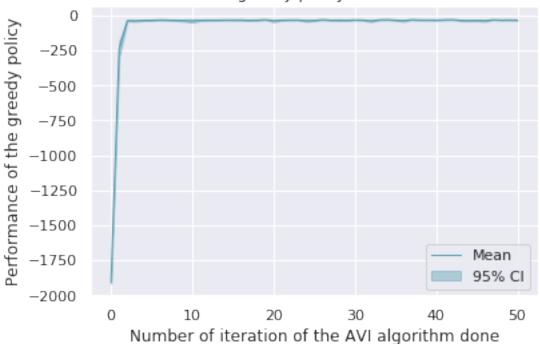


```
# Collect the samples using the behavioural policy
        # You should use discrete actions
       beh_policy = behavorialPolicy(actions)
       D_n = generateDataset(beh_policy, horizon=horizon, n_episodes=n_episodes)
        # Generation of the feature matrix
       Phi_Matrix = phi(states, actions)
        # Number of experiences
       n_{experiences} = 20
        # define FQI
       performances = [[] for exp in range(n_experiences)]
       for exp in tqdm.tqdm(range(n_experiences)):
           theta_l = fqi(env, D_n, Phi_Matrix)
           for theta_k in theta_l:
              # Computation of the greedy policy
              policy_k = greedyPolicy(theta_k)
              # plot obtained Q-function against the true one
              J = utils.estimate_performance(env, policy=policy_k, horizon=horizon,
                                       n_episodes=n_episodes, gamma=discount)
              performances[exp].append(J)
100%|| 20/20 [06:18<00:00, 18.43s/it]
# Display the performance
```

```
%matplotlib inline
# Converting resulting list
performances = np.array(performances)
# Initialisation of the figure
fig, axs = plt.subplots()
# X data
x = np.linspace(0, n_episodes, n_episodes+1)
y = np.mean(performances, axis=0)
# Compute the CI bounds
low_CI, upper_CI = computeCIBounds(performances)
# Plot the data, set the linewidth, color and transparency of the
# line, provide a label for the legend
axs.plot(x, y, lw=1, color='#539caf', alpha=1, label='Mean')
# Shade the confidence interval
axs.fill_between(x, low_CI, upper_CI, color = '#539caf', alpha=0.4, label='95% CI')
# Label the axes and provide a title
axs.set_title("Peformance of the greedy policy over the iterations of AVI")
axs.set_xlabel("Number of iterations of the AVI algorithm done")
axs.set_ylabel("Performance of the greedy policy")
# Display legend
axs.legend(loc='best')
```

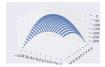
plt.show()







```
ax = fig.add_subplot(111, projection='3d')
ax.scatter(S, A, Q_obtained)
ax.mouse_init()
plt.show()
```



## In []: