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# 1 Theory and Methods

# 1.1 The Inverted Pendulum (IP)

# 1.1.1 Dynamics

The Inverted Pendulum is an inherently unstable system with highly nonlinear dynamics, and is under-actuated.

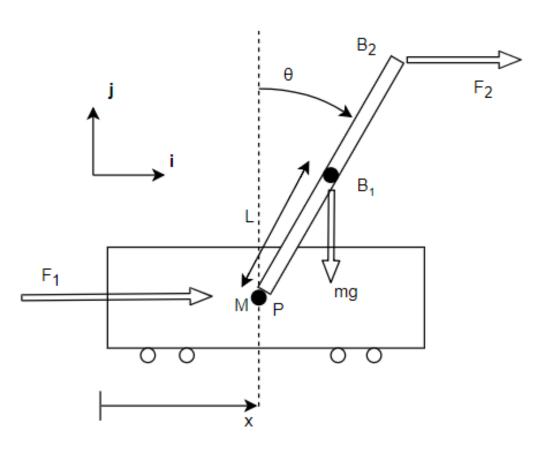


Figure 1.1: A free-body diagram of the inverted pendulum system. For the OpenAI IP the system is in discrete time and constrained by  $L=0.5\mathrm{m},\ m=0.1\mathrm{kg},\ M=1\mathrm{kg},$   $F=\pm 10\mathrm{N},\ x_{max}=\pm 2.4\mathrm{m},\ \theta_{max}=\pm 12^{\circ}.$ 

The full state space equations for the inverted pendulum as defined in fig. 1.1 are given

by:

$$\begin{bmatrix} \dot{x} \\ \ddot{x} \\ \dot{\theta} \\ \ddot{\theta} \end{bmatrix} = \begin{bmatrix} \dot{x} \\ \frac{(2M+m)F_2 - F_1(cos\theta + g(M+m)sin\theta - mL\dot{\theta}^2 sin\theta cos\theta)}{(M+msin^2\theta)} \\ \dot{\theta} \\ \frac{F_1 - F_2 cos(2\theta) + msin\theta(L\dot{\theta}^2 - gcos\theta)}{L(M+msin^2\theta)} \end{bmatrix}$$
(1.1)

Using Lyapunov's indirect method, we can write the linearised equations about the equilibrium,  $\mathbf{x}_e = [x_e, \dot{x}_e, \theta_e, \dot{\theta}_e]^T = [0, 0, 0, 0]^T$ , as:

$$\begin{bmatrix}
\delta \dot{x} \\
\delta \dot{x} \\
\delta \dot{\theta} \\
\delta \ddot{\theta}
\end{bmatrix} = \begin{bmatrix}
0 & 1 & 0 & 0 \\
0 & 0 & \frac{mg}{M} & 0 \\
0 & 0 & 0 & 1 \\
0 & 0 & \frac{(m+M)}{ML}g & 0
\end{bmatrix} \begin{bmatrix}
\delta x \\
\delta \dot{x} \\
\delta \theta \\
\delta \dot{\theta}
\end{bmatrix} + \begin{bmatrix}
0 & 0 \\
\frac{1}{M} & -\frac{1}{M} \\
0 & 0 \\
-\frac{1}{ML} & \frac{2M+m}{mML}
\end{bmatrix} \begin{bmatrix}
\delta F_1 \\
\delta F_2
\end{bmatrix}$$
(1.2)

The eigenvalues are given by  $det(\lambda I - A) = \lambda^2 (\lambda^2 - \frac{(m+M)}{ML}g) = 0$ . Therefore, the system is unstable about  $\boldsymbol{x}_e$  due to the right half plane pole,  $\lambda = \sqrt{\frac{(m+M)}{ML}g}$ . Additionally, the time constant of this unstable system is  $\tau = \sqrt{\frac{ML}{g(m+M)}} = 0.22s$ . Note, if  $M >> m, \tau \to \sqrt{\frac{L}{g}}$ , which is the time constant for a simple pendulum.

OpenAI's gym is a python package that supplies an inverted pendulum environment built-in (called CartPole). This environment was wrapped to use the dynamics above, whilst providing a rendering function shown in fig. 1.2.

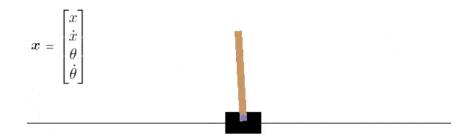


Figure 1.2: The OpenAI gym CartPole environment. The classical state representation is shown in the top left. Actions by the player and the adversary are taken as an impulse to the left or right as defined in fig. 1.1.

## 1.1.2 Cost and Value Function

For each step, a cost is calculated as in eq. (1.3), where  $\mathbf{w}^T = [w_1, w_2, w_3, w_4] = [0.4, 0.1, 0.7, 1]$  and  $0 \ge c(x_t, u_t) \ge -1$ . The weights,  $\mathbf{w}$ , were chosen through empirical measurement of the the importance of each state.

$$c(x_t, u_t) = -\frac{1}{\sum_i w_i} \cdot \hat{\boldsymbol{x}}^T \begin{bmatrix} w_1 & 0 & 0 & 0 \\ 0 & w_2 & 0 & 0 \\ 0 & 0 & w_3 & 0 \\ 0 & 0 & 0 & w_4 \end{bmatrix} \hat{\boldsymbol{x}}$$
(1.3)

where, 
$$\hat{\boldsymbol{x}}^T = \left[\frac{x_t}{x_{max}}, \frac{\dot{x}_t}{\dot{x}_{max}}, \frac{\theta_t}{\theta_{max}}, \frac{\dot{\theta}_t}{\dot{\theta}_{max}}\right]^T$$
 (1.4)

Weighting on the inputs was set to zero as there are only two inputs for this problem, thus the cost can be written as  $c(x_t)$ . The max values can be approximated experimentally (note,  $x_{max} = 2.4$  and  $\theta_{max} = 12^o$  are given constraints):

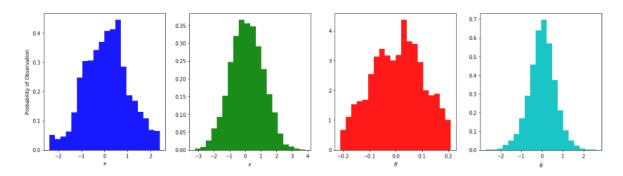


Figure 1.3: Histograms of typical state values. The frequencies greatly depend on the quality of the controller, with better controllers giving much narrower distributions. However, these are typical for a controller of medium efficacy over many episodes where the starting state is randomised (possibly to an uncontrollable state).

Suitable estimates for the the values of  $\dot{x}_{max}$  and  $\dot{\theta}_{max}$  are thus approximately 3 and 2 respectively. After the episode has completed, the "true value" of each state is computed as the discounted future costs over a finite horizon (eq. (1.5)). I.e. the value of a state,  $\mathbf{x}_t$ , is given by  $v(\mathbf{x}_t) \propto c(\mathbf{x}_t) + \gamma c(\mathbf{x}_{t+1}) + ... + \gamma^k c(\mathbf{x}_{t+k})$ , where k is the length of the horizon (chosen to be 14). A horizon constraint of  $\gamma^k < \frac{1}{20}$  was chosen as it is a standard factor for insignificance. Therefore  $\gamma$  is calculated as  $\gamma < \frac{1}{20}^{\frac{1}{k}} \Longrightarrow$  for k = 14,  $\gamma = 0.8$ .

$$v_0 = \frac{\sum_{\tau=0}^k \gamma^{\tau} c(x_{\tau})}{\sum_{\tau=0}^k \gamma_{\tau}}, \quad \text{where } \gamma^k < \frac{1}{20}$$
 (1.5)

Where for simplicity of notation,  $v_0 = v(t)$ , the state value at step t. Once a terminal state (either out of bounds, or the maximum number of steps) has been reached, the simulation continues for k steps in order for a value to be calculated for every step of the episode.

## 1.1.3 State Representations

The state can be represented in a number of ways. The simplest method would be  $\mathbf{x} = [x, \dot{x}, \theta, \dot{\theta}]$ . This has a number of advantages such as lower computational cost, greater numerical accuracy (if the process is fully observable) and simpler implementation. Conversely, a 2-dimensional (2D) representation may be used. There are several possibilities for this, all of which first require binning  $\mathbf{x}$ :

(1) A matrix stack of x vs  $\dot{x}$  and  $\theta$  vs  $\dot{\theta}$ , both of which would only have one non-zero entry. This scales as  $b^n$  where b = number of bins and n = number of states.

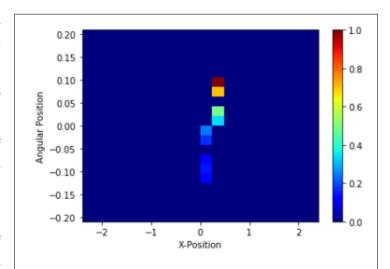


Figure 1.4: An example of a 2D state representation where there are 20 bins and 17 random actions have been taken.

- (2) A matrix stack of  $x_t$  vs  $x_{t-1}$  for all states. Similarly this scales as  $b^n$ , however the derivative states do not need to be plotted as these can be inferred. This has the advantage that, if the derivatives are not observable, they are built into the 2D representation, however, if they are observable then this is less accurate than (1).
- (3) A matrix of discounted previous states, forming a motion history image. An example of this is shown in section 1.1.3, and algorithm 1 shows the implementation details. This was the chosen representation.

A 2D representation such as this allows us to use a convolutional neural network (CNN), which has the benefit of various transformation invariances - these are particularly useful for the inverted pendulum since it is symmetric.

### Algorithm 1 Create 2D State

9: end function

```
1: function GETSTATE2D(\hat{x}_{t-1}^{(2D)}, binEdges, nBins)
          \hat{\boldsymbol{x}} \leftarrow \text{getNormedState}()
2:
          for all x_i \in \hat{\boldsymbol{x}}_t do
3:
                x_i \leftarrow \operatorname{argmin}|binEdges - x_i|
                                                                               ▷ Get the index of the nearest bin edge.
4:
          end for
5:
          HistEdges \leftarrow \texttt{linspace}(-0.5, nBins - 0.5, nBins + 1) \quad \triangleright \texttt{Centre by shifting -0.5}
6:
          \hat{\boldsymbol{x}}_{t}^{(2D)} \leftarrow \text{histogram2d}(x, \theta, \text{bins} = (histEdges, histEdges))
7:
                                                                                                                     ▶ Inbuilt function
          return \hat{\boldsymbol{x}}_{t}^{(2D)} + \lambda \hat{\boldsymbol{x}}_{t-1}^{(2D)}
8:
```

The linearised inverted pendulum (valid for small time steps) can be modelled as a state space model:

$$\boldsymbol{x}_{t}^{(2D)} = C\boldsymbol{x}_{t} + \boldsymbol{V}_{t}$$
  $\boldsymbol{V}_{t} \sim \mathcal{U}\left(\begin{bmatrix} \frac{1}{\delta x} \\ \frac{1}{\delta \theta} \end{bmatrix}\right)$  (1.6)

$$\boldsymbol{x}_t = A\boldsymbol{x}_{t-1} + B\boldsymbol{u}_t \tag{1.7}$$

Where A and B are the linearised system dynamics, and C is the linear transformation to a 2D state space, with quantisation noise  $\mathbf{V}$  modelled as a uniform random disturbance,  $\mathcal{U}$ , centred on (0, 0).

The initial mean squared error (MSE), and the propagation of the error (when estimated optimally) are given by eqs. (1.8) and (1.9) where  $\delta x = \delta \theta$  (derivation details can be found in appendix A.2).

$$\Sigma_0 = \sigma_v^2 I = \begin{bmatrix} \frac{\delta x^2}{12} & 0\\ 0 & \frac{\delta \theta^2}{12} \end{bmatrix} \tag{1.8}$$

$$\Sigma_{n+1} = (I - \bar{\Sigma}_{n+1}C^T(\sigma_v^2 I + C\bar{\Sigma}_{n+1}C^T)^{-1}C)\bar{\Sigma}_{n+1}$$
(1.9)

Where  $\bar{\Sigma}_{n+1} = A\Sigma_n A^T$ . It can be shown (see appendix A.2) that as  $t \to \infty$ , the covariance of quantisation decays to zero since  $\lim_{\delta x \to 0} \sigma_v^2 = 0$  and therefore this form for a 2D state becomes lossless (assuming that the neural network acts as an optimal Kalman filter). For a given bin size, the MSE decreases linearly with time, however, the MSE decreases quadratically with bin size. Therefore, a smaller bin size will quadratically improve the rate of decay but quadratically increase the computational load of when computing the state and through the neural network. Therefore, there is a tradeoff in choosing the number of bins between computational complexity and state error.

With a non-zero bin size the overall error can be reduced by binning more densely in regions in which the IP is expected to spend more time. Figure 1.3 shows that the state visit frequencies roughly correspond to normal distribution, therefore by transforming the bins with an inverse Gaussian c.f.d. a flattened distribution can be obtained with a greater density of bins in the centre (fig. 1.5). This has the additional benefit of allowing finer control where it is needed. For example, if the pole is far from the equilibrium the optimal action more easily determined, and subject to less change with small state variations, therefore coarser binning can be used.

## 1.1.4 Discretisation and Continuous Time

For a small enough discrete time step, the simulation will approximate the continuous process very well. Under a continuous action space, the sampling rate must be faster

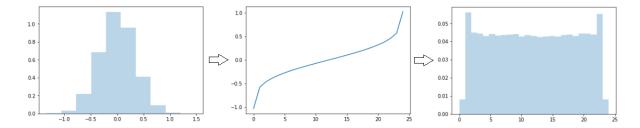


Figure 1.5: Binning of a Gaussian distribution with bin edges scaled with an inverse Gaussian c.f.d. For this example there are 25 bins.

than the system dynamics to ensure controllability. When the action space is restricted to  $u \in \{-1,1\}$ , there is a trade off between the sampling rate and the size of the action space. If the sampling rate is fast enough pseudo-continuous actions can be achieved via pulsewidth modulation. For the IP system described, the time constant,  $\tau$ , is approximately 0.022s. Moreover, the agents move alternately, therefore the sampling time step  $\delta t$ , must be less than 0.011s, however due to computational limitations, a time step of 0.01s was chosen.

The positional change per time step can be computed by considering the average maximum velocities: 3m/s and 2rad/s. Thus the maximum positional change between an agent's actions is  $\dot{x}_{i,max} \times \frac{\delta t}{x_{i,max}} = 2.5\%$  and 30% respectively. In practice a 30% change in  $\theta$  would only occur around  $\theta_{max}$  and would be uncontrollable. Therefore, the choice of  $\delta t = 0.01$  is not detrimental to the control, but will not give pseudo-continuous actions.

This maximal positional change per time step also informs the bin size needed. Ideally, each time step lands within a new bin, therefore for the x-position,  $\frac{1}{2.5\%} = 40$  is the minimum number of bins needed. Similarly,  $\frac{1}{30\%} = 3$  is the minimum number of bins needed for the angular position, however, empirically,  $\theta$  is a more "important" state variable due to its relative ease of exceeding  $\theta_{max}$ . Therefore, given that the bins are more finely space near (0, 0) and for simplicity, 40 bins for both x and  $\theta$  was chosen.

# 1.2 Self Play and Adversaries

### 1.2.1 Cost and Value Functions

In AlphaZero, the adversary is working to minimise the value function whereas the player is working to maximise it. For board games, where agents are on equal footing, a value function (-1 < v < 1) representing the expected outcome can be used. This has the advantage of symmetry - when the board is viewed from the other agent's perspective the value can be multiplied by -1, giving the expected outcome for that agent.

The adversary for the inverted pendulum has the additional advantage of gravity, making the play asymmetric. Given that both the state-value and cost are -1 < v, c < 0,

multiplying by -1 would mean the adversary is maximising a value function 0 < v < 1. State-values outside these ranges have no meaning. If a single neural network is used, values close to equilibrium may be predicted into the wrong range. Consequently, both the adversary and the player must predict true state-values. This also has the advantage of maintaining an intuitive meaning of the state-value.

#### Choice of Action 1.2.2

The inverted pendulum system is symmetric in both x and  $\theta$ . By taking advantage of this the number of training examples could be doubled by reflecting episodes along the axis of symmetry. However, as shown with AlphaZero [15], this provides minimal benefit and also hinders generalisation. The adversarial point of action was chosen to be at the top of the pole, acting horizontally (fig. 1.1), thus ensuring that two distinct policies must be learnt, rather than one being just inverse probabilities of the other. For simplicity  $u_{adv} \in h \cdot \{-1, 1\}$  was chosen, where h is a handicap applied to the adversary.

The handicap can be approximated using the linearised dynamics eq. (1.2) by investigating what force would be needed by each agent to achieve the same change in the state (eqs. (1.10a) and (1.10b)) (setting  $\delta x = 0$ ).

$$\delta \ddot{x} = \frac{1}{M} \delta F_1 - \frac{1}{M} \delta F_2 \implies \delta F_2 \approx \delta F_1$$
 (1.10a)

$$\delta \ddot{x} = \frac{1}{M} \delta F_1 - \frac{1}{M} \delta F_2 \quad \Longrightarrow \quad \delta F_2 \approx \delta F_1 \tag{1.10a}$$

$$\delta \ddot{\theta} = -\frac{1}{ML} \delta F_1 + \frac{2M+m}{mML} \delta F_2 \quad \Longrightarrow \quad \delta F_2 \approx \frac{m}{2M+m} \delta F_1 \tag{1.10b}$$

Empirically, the pendulum is more likely to fall over (exceed  $\theta_{max}$ ) rather than to exceed  $x_{max}$ . Therefore, a good estimate of h would be  $\frac{m}{2M+m} = \frac{0.1}{2+0.1} \approx 0.0476$ . When testing this it was found that if  $F_1$  was set near 0.04-0.05 initially, the adversary gets the upper-hand quickly and the pendulum falls very quickly, not allowing the player to gain any training experience. Therefore, for the first N iterations, the adversary force was set to 0, and increased as the number of iterations increased according to  $h = (1 - 0.7^{i-N+1})$ , where i is the iteration, and 0.7 was chosen such that the adversary could be implemented on the first iteration without overpowering the still-training player.

#### Agent Representation 1.2.3

There are two good ways of representing whether it is the adversary's turn for the neural network:

Multiply the board representation by -1 such that opponent pieces are negative. This has the disadvantages that it can only take two agents and, a network that outputs both state-values and action p.m.f.'s should predict the same state values for both player and adversary, but predict vastly different actions. Therefore, the values could be decoupled from the actions, which was one of the major benefits of using a single neural network. However, a single network is simpler, and negating the board more closely follows AlphaZero's methodology.

Record the agent with each example and use agent-labeled examples to train different neural networks. Using a neural network for each agent causes half of the examples to be lost as only the relevant agent's examples are used to train each network. However, this does not suffer from the problems above and can cope with agents with a different number of possible actions more easily.

Recording the agent with each example was therefore chosen. Note, in the case of the inverted pendulum, the optimal action is the inverse of the worst action. However, this is not a general result. For example, in a system with non-linear and asymmetric dynamics it is possible to have the target perpendicular to the closest edge of the invariant set, thus for the adversary it is better to push the system into instability rather than away from equilibrium.

## 1.2.4 Episode Execution and Policy Iteration

Pseudocode for episode execution following the sections above is shown in Algorithm 2.

```
Algorithm 2 Execute Episode
```

```
1: function EXECUTEEPISODE
 2:
           example \leftarrow []
           \boldsymbol{x}, \boldsymbol{x}^{(2D)}, c \leftarrow \text{resetEpisode}()
                                                                 \triangleright Set initial \boldsymbol{x} randomly and initialise the cost
 3:
           agent \leftarrow 0
 4:
           repeat
 5:
                \boldsymbol{\pi} \leftarrow \operatorname{getActionProb}(\boldsymbol{x}, \boldsymbol{x}^{(2D)}, agent)
                                                                                            ▶ Perform MCTS Simulations
 6:
                example.append((\boldsymbol{x}^{(2D)}, \boldsymbol{\pi}, c, agent))
 7:
                                                                       \triangleright Sample action. Note after S steps, \tau \to 0
                u \sim \pi
 8:
                \boldsymbol{x}, c \leftarrow \text{step}(\boldsymbol{x}, u)
 9:
                                                                                            \boldsymbol{x}^{(2D)} \leftarrow \text{getState2D}(\boldsymbol{x}, \boldsymbol{x}^{(2D)})
10:
                agent \leftarrow nextAgent(agent)
11:
12:
           until episodeEnded(\boldsymbol{x})
           example \leftarrow convertCostsToValues(example)
13:
           return example
14:
15: end function
```

The action, u, is sampled from  $\pi$ . However, after S steps, the temperature,  $\tau \to 0$ , which is equivalent to  $u = \operatorname{argmax} \pi$ .

The overall policy iteration algorithm is given by section 1.2.4.

## Algorithm 3 Policy Iteration - Training the NeuralNetwork

```
1: function PolicyIteration
       nnet \leftarrow NeuralNetwork
                                                                   ▶ Initialise NeuralNetwork
 2:
       examples \leftarrow []
 3:
       for iter in policyIterations do
 4:
           incrementHandicap(iter)
                                                                      \triangleright F_2 = 0 for first n iters
 5:
           for ep in trainingEpisodes(iter) do
                                                             ▶ More episodes for first n iters
 6:
                                                                 ▷ Use the most current nnet
               resetMCTS(nnet)
 7:
               example \leftarrow executeEpisode()
 8:
               examples.append(example)
9:
           end for
10:
                                                                  ▶ Last N policys' examples
           recents \leftarrow getMostRecentExamples(examples)
11:
           nnet \leftarrow trainNeuralNetwork(recents, nnet)
12:
           saveNeuralNetwork(nnet)
13:
       end for
14:
15: end function
```

# 1.3 Neural Networks

### 1.3.1 Loss Functions and Pareto

The following loss function is minimised when training the neural networks:

$$\mathcal{L} = \underbrace{\sum_{t} (v_{\theta}(\boldsymbol{x}_{t}) - v_{t})^{2} + c_{pareto} \cdot \underbrace{\boldsymbol{\pi}_{t} \cdot log(\boldsymbol{p}_{\theta}(\boldsymbol{x}_{t}))}_{\mathcal{L}_{action}} + \underbrace{c||w_{nnet}||^{2}}_{regularisation}}$$
(1.11)

Assuming that policy action predictions and value predictions are of similar importance, similar magnitudes of action and value losses are desireable. The constant,  $c_{pareto}$ , was chosen such that  $(\mathcal{L}_{action} \approx \mathcal{L}_{value})$ . The minimum entropy for two possible actions is 1bit, and since the action loss is the negative log likelihood of a categorical distribution, the minimum loss is -1. The value losses are calculated using the MSE, which is positive, and has an average value of approximately 0.1 (figs. A.1 to A.4), therefore  $c_{pareto}$  was set to 0.1.

During training the agents were chosen to move probabilistically throughout for the first 14 moves, and then move greedily. 14 was chosen as this is the average episode length when the player moves randomly without an adversary.

## 1.3.2 Architectures

Two neural network architectures are used, one that takes the raw state,  $\mathbf{x} = [x \ \dot{x} \ \theta \ \dot{\theta}]^T$  as input (the "baseline" network), and another that takes the 2D state,  $\mathbf{x}^{(2D)}$  (the convolutional neural network, "CNN"). The adversary and player have separate networks, both predicting the state-value and action p.m.fs.

For the network with  $\mathbf{x}^{(2D)}$  as input, there are 2 fully connected convolutional layers with max-pooling, followed by a feedforward layer with ReLU activations. The two "heads" split here and each has fully-connected layer, with the value head outputting  $v_{\theta}$  and the action head outputting  $p_{\theta}$ . The network with  $\mathbf{x}$  as input is the same, except the convolutional layers are replaced with 2 feedforward layers.

Both architectures perform training with the *Adam Optimiser* using a learning rate of 0.0005, a mini-batch size of 64, a dropout of 0.3, batch normalisation and 1 epoch. The networks are implemented in pytorch<sup>1</sup>.

The neural network is evaluated once every step in the MCTS, therefore in an episode of length L and S MCTS simulations/step, the neural network is evaluated more than  $L \times S$  times. The GPU used to run experiments is a single NVIDEA GeForce GTX 1050. The convolutional layers of a neural network are the most computationally expensive part and, increasing with the image size, are more than 50% of the computational time in the architectures defined above. Therefore, there is a trade off between the number of convolutional layers used and computation time. The number of layers was chosen such that running the program takes less than 24hrs. Additionally, the memory of the GPU is 4GB, which limits the number of weights that can be used in the neural networks. It was found that adding an additional layer to the CNN exceeds this memory limit. The architecture of the baseline network was chosen to match that of the CNN.

## 1.4 MCTS

The general MCTS algorithm, outlined in ??, was implemented for the inverted pendulum as in algorithms 4 and 5.

## 1.4.1 Tree Nodes and Simulations

The state for the inverted pendulum is continuous ( $x \in \mathbb{R}^4$ ), as opposed to the discrete nature of board games. In order to be able to compare nodes, the states must be stored as unique integers. Each MCTS simulation ends with a leaf node being expanded, therefore after S simulations from a node, S states will be visited. Over the whole episode,  $S \times L$  distinct states will be visited in total. Thus, the probability of revisiting a state from a

<sup>&</sup>lt;sup>1</sup>www.pytorch.org

different trajectory is  $(S \times L)b^4$ , where b is the number of bins for each dimension. Given that L = 400 and, if S > 30, the maximum recursion depth of the MCTS will be reached - limiting the maximum value of S. Multiplying the states by  $10^{12}$  and converting to integers ensures a vanishingly small probability of encountering a state twice.

### 1.4.2 Action Selection

Action selection moving down the tree was modified to reflect the asymmetry of the state-value:

$$u_{player}^{*} = \underset{u \in \mathcal{U}_{player}}{\operatorname{argmax}} \left\{ Q(\boldsymbol{x}, u) + c_{explore} \cdot \boldsymbol{p}_{\theta}(\boldsymbol{x}, u) \frac{\sqrt{\sum_{b} N(\boldsymbol{x}, b)}}{1 + N(\boldsymbol{x}, b)} \right\} = \underset{u \in \mathcal{U}_{player}}{\operatorname{argmax}} U_{player}(\boldsymbol{x}, u)$$

$$(1.12)$$

$$u_{adv}^{*} = \underset{u \in \mathcal{U}_{adv}}{\operatorname{argmax}} \left\{ - Q(\boldsymbol{x}, u) + c_{explore} \cdot \boldsymbol{p}_{\theta}(\boldsymbol{x}, u) \frac{\sqrt{\sum_{b} N(\boldsymbol{x}, b)}}{1 + N(\boldsymbol{x}, b)} \right\} = \underset{u \in \mathcal{U}_{adv}}{\operatorname{argmax}} U_{adv}(\boldsymbol{x}, u)$$

$$(1.13)$$

Where c = 1 was used. Negating Q in  $U_{adv}$  causes the adversary to minimise the state-value whilst still maximising the upper confidence bound, thus ensuring exploration.

When the inverted pendulum exceeds a constraint the tree search should return -1. Note, the simulation terminates after 400 steps (the arbitrarily set "end" of the episode), but this should not return -1. Therefore, when 400 steps is reached, the neural network is evaluated and v is returned.

```
Algorithm 4 Get Action Probabilities (based on S.Nair's implementation [16])
```

```
1: function GETACTIONPROB(x_{root}, x^{(2D)}, agent)
         for i in nMCTSSims do
                                                                 \triangleright Simulate nMCTSSims episodes from x_{root}
2:
              resetSimulation(\boldsymbol{x}_{root})
3:
              MCTSSearch(\boldsymbol{x}^{(2D)}, agent)
4:
         end for
5:
         N(\boldsymbol{x}_{root}, u) \leftarrow \text{getCounts}()
                                                  \triangleright Count the times each edge, (\boldsymbol{x}_{root}, u), was visited.
6:
         N(\boldsymbol{x}_{root}, u) \leftarrow N(\boldsymbol{x}_{root}, u)^{\tau}
                                                                  \triangleright Control Sparsity with the temperature, \tau
7:
         return \pi \leftarrow \text{norm}(N(\boldsymbol{x}_{root}, u))
9: end function
```

# 1.5 Player and Adversary Evaluation

For problems in which one agent has a distinct advantage, determining how "good" their policy is compared to other policies becomes difficult. For example, if the pendulum stays

up for longer, is is difficult to determine whether the player has improved because it is controlling the system better, or the adversary is less effective, or both have improved but one improved more. The score used is the number of steps the agent stayed up for divided by the max number of steps. For the player this ranges between 0 and 1, and for the adversary score = (200 - steps)/200 is used.

Over one policy iteration, there are 4 different policies: the two current player and adversary policies,  $f_{\theta, curr}^{(player)}$  and  $f_{\theta, curr}^{(adv)}$ , and the challenger policies,  $f_{\theta, chal}^{(player)}$  and  $f_{\theta, chal}^{(adv)}$ . The improvement of the player can be determined by pitting the current player against the current adversary, and then the challenging player against the current adversary; and vice-versa for the adversary improvement<sup>2</sup>. This gives the four equations in ??.

does this make sense?

$$\mathbb{E}[s_{curr}^p] = \frac{1}{1 + 10^{\wedge} \left(\frac{\mu_{curr}^a - \mu_{curr}^p}{400}\right)} \qquad \qquad \mathbb{E}[s_{curr}^a] = \frac{1}{1 + 10^{\wedge} \left(\frac{\mu_{curr}^p - \mu_{curr}^a}{400}\right)} \tag{1.14}$$

$$\mathbb{E}[s_{curr}^{p}] = \frac{1}{1 + 10^{\wedge} \left(\frac{\mu_{curr}^{a} - \mu_{curr}^{p}}{400}\right)} \qquad \mathbb{E}[s_{curr}^{a}] = \frac{1}{1 + 10^{\wedge} \left(\frac{\mu_{curr}^{p} - \mu_{curr}^{a}}{400}\right)} \qquad (1.14)$$

$$\mathbb{E}[s_{chal}^{p}] = \frac{1}{1 + 10^{\wedge} \left(\frac{\mu_{curr}^{a} - \mu_{chal}^{p}}{400}\right)} \qquad \mathbb{E}[s_{chal}^{a}] = \frac{1}{1 + 10^{\wedge} \left(\frac{\mu_{curr}^{p} - \mu_{chal}^{a}}{400}\right)} \qquad (1.15)$$

where  $\mathbb{E}[s_{curr}^p] = \mathbb{E}[s(f_{\theta, \, curr}^{(player)})]$  denotes the score for the current player, and  $\mu_{curr}^p =$  $\mu(f_{\theta, curr}^{(player)})$  is the rating of the current player. The expected scores can be estimated via monte-carlo. Rearranging gives the relation between successive agent ratings as:

$$\mu_{chal}^{p} = \mu_{curr}^{p} + 400log_{10} \left( \frac{E[s_{chal}^{p}]}{1 - E[s_{chal}^{p}]} \cdot \frac{1 - E[s_{curr}^{p}]}{E[s_{curr}^{p}]} \right)$$
(1.17)

<sup>&</sup>lt;sup>2</sup>Note, if the ratings were correctly updated in the last policy iteration then  $\mathbb{E}[\mathbb{E}[s_{curr}^{agent}] - \mu_{curr}^{agent}] = 0$ 

## Algorithm 5 MCTS (based on S.Nair's implementation [16])

```
1: function MCTSSEARCH(\boldsymbol{x}^{(2D)}, agent)
 2:
            if x is terminal then
 3:
                  if fallen then
                         return -1
 4:
                  else
 5:
                         \boldsymbol{\pi}, v \leftarrow f_{\theta}(\boldsymbol{x}^{(2D)})
 6:
                         return v
                                                                                     \triangleright If the episode end is reached, return v_{\theta}
 7:
                  end if
 8:
 9:
            end if
10:
            if \boldsymbol{x} \notin Tree then
                                                                                                                         ▷ Expand Leaf Node
11:
                  \boldsymbol{\pi}, v \leftarrow f_{\theta}(\boldsymbol{x}^{(2D)})
12:
                  N(\boldsymbol{x},\cdot) \leftarrow 0
13:
                  P(\boldsymbol{x},\cdot) \leftarrow \boldsymbol{\pi}
14:
15:
                  return v
            end if
16:
17:
                                                                                                          ▶ Get best action using UCB
            if agent = player then
18:
                  u^* \leftarrow \operatorname{argmax} U_{player}(\boldsymbol{x}, u)
19:
                            u \in \mathcal{U}_{player}
20:
            else
                  u^* \leftarrow \operatorname{argmax} U_{adv}(\boldsymbol{x}, u)
                              u \in \mathcal{U}_{adv}
            end if
22:
            \boldsymbol{x}, c \leftarrow \text{step}(\boldsymbol{x}, u)

    ▶ Take next MCTS simulated step

23:
            \boldsymbol{x}^{(2D)} \leftarrow \text{getState2D}(\boldsymbol{x}, \boldsymbol{x}^{(2D)})
24:
            agent \leftarrow nextAgent(agent)
25:
            v \leftarrow \text{MCTSSearch}(\boldsymbol{x}^{(2D)}, agent)
                                                                                                                 ▶ Recursion to next node
26:
27:
            Q(\boldsymbol{x}, u^*) \leftarrow \frac{N(\boldsymbol{x}, u^*)Q(\boldsymbol{x}, u^*) + v}{N(\boldsymbol{x}, u^*) + 1}
28:
                                                                                                       ▶ Backup Q-values up the tree
            N(\boldsymbol{x}, u^*) \leftarrow N(\boldsymbol{x}, u^*) + 1
29:
30:
            return v
31: end function
```

# A Appendices

# A.1 Inverted Pendulum Dynamics Derivation

The state space equations for the Inverted Pendulum can be found using d'Alembert forces. Define the distance and velocity vectors to the important points:

$$egin{aligned} m{r}_P &= x m{i} \ m{r}_{B_1/P} &= L sin heta m{i} + L cos heta m{j} \ m{r}_{B_1} &= (x + L cos heta) m{i} + L \dot{ heta} sin heta m{j} \ \dot{m{r}}_{B_1} &= (\dot{x} + L \dot{ heta} cos heta) m{i} - L \dot{ heta} sin heta m{j} \end{aligned}$$

Linear Momentum,  $\rho = \sum_{i} m_i \dot{r}_{i/o} = m \dot{r}_{B_1} + M \dot{r}_P$ :

$$\boldsymbol{\rho} = \begin{bmatrix} (M+m)\dot{x} + mL\dot{\theta}cos\theta \\ -mL\dot{\theta}sin\theta \\ 0 \end{bmatrix}$$

Moment of momentum about P,  $\boldsymbol{h}_P = \boldsymbol{r}_{B_1/P} \times m \boldsymbol{\dot{r}}_{B_1}$ :

$$\mathbf{h}_{P} = -mL(L\dot{\theta} + \dot{x}cos\theta)\mathbf{k}$$
$$\therefore \dot{\mathbf{h}}_{P} = -mL(L\ddot{\theta} + \ddot{x}cos\theta - \dot{x}\dot{\theta}sin\theta)\mathbf{k}$$

Balance moments using  $\dot{\boldsymbol{h}}_P + \dot{\boldsymbol{r}}_P \times \boldsymbol{\rho} = \boldsymbol{Q}_e$  and  $\boldsymbol{Q}_e = \boldsymbol{r}_{B_1/P} \times -mg\boldsymbol{j} + \boldsymbol{r}_{B_2/P} \times F_2\boldsymbol{i}$ :

$$m{\dot{h}}_P + \dot{m{r}}_P imes m{
ho} = egin{bmatrix} 0 \\ 0 \\ -mL(\ddot{x}cos heta + L\ddot{ heta}) \end{bmatrix} = egin{bmatrix} 0 \\ 0 \\ -L(mgsin heta + 2F_2cos heta) \end{bmatrix} = m{Q}_e$$

And also balance linear momentum using  $F_e = \dot{\rho}$ :

$$\dot{\boldsymbol{\rho}} = \begin{bmatrix} (m+M)\ddot{x} + mL(\ddot{\theta}cos\theta - \dot{\theta}^2sin\theta) \\ -mL(\ddot{\theta}sin\theta + \dot{\theta}^2cos\theta) \\ 0 \end{bmatrix} = \begin{bmatrix} F_1 + F_2 \\ R - (M+m)g \\ 0 \end{bmatrix} = \boldsymbol{F}_e$$

Finally write the system dynamics in terms of  $\ddot{\theta}$  and  $\ddot{x}$ :

$$\ddot{\theta}(M+msin^2\theta)L = \left(\frac{2M+m}{m}F_2 - F_1\right)cos\theta + g(M+m)sin\theta - mL\dot{\theta}^2sin\theta cos\theta \quad (A.1)$$

$$\ddot{x}(M + m\sin^2\theta) = F_1 - F_2\cos(2\theta) + m\sin\theta(L\dot{\theta}^2 - g\cos\theta) \tag{A.2}$$

Simplifying by substituting in constants, the full state space equation are:

$$\begin{bmatrix} \dot{x} \\ \ddot{x} \\ \dot{\theta} \\ \ddot{\theta} \end{bmatrix} = \begin{bmatrix} \dot{x} \\ \frac{(2M+m)F_2-F_1)\cos\theta+g(M+m)\sin\theta-mL\dot{\theta}^2\sin\theta\cos\theta}{(M+m\sin^2\theta)} \\ \dot{\theta} \\ \frac{F_1-F_2\cos(2\theta)+m\sin\theta(L\dot{\theta}^2-g\cos\theta)}{L(M+m\sin^2\theta)} \end{bmatrix} = \begin{bmatrix} f_1(\boldsymbol{x},F_1,F_2) \\ f_2(\boldsymbol{x},F_1,F_2) \\ f_3(\boldsymbol{x},F_1,F_2) \\ f_4(\boldsymbol{x},F_1,F_2) \end{bmatrix}$$
(A.3)

Using Lyapunov's indirect method, the linearised equations about the equilibrium,  $\mathbf{x}_e = [x_e, \dot{x}_e, \theta_e, \dot{\theta}_e]^T = [0, 0, 0, 0]^T$ , are:

$$\begin{bmatrix} \delta \dot{x} \\ \delta \ddot{x} \\ \delta \dot{\theta} \\ \delta \ddot{\theta} \end{bmatrix} = \begin{bmatrix} \frac{\partial f_1}{\partial x} |_{\mathbf{x}_e} & \frac{\partial f_1}{\partial \dot{x}} |_{\mathbf{x}_e} & \frac{\partial f_1}{\partial \theta} |_{\mathbf{x}_e} & \frac{\partial f_1}{\partial \dot{\theta}} |_{\mathbf{x}_e} \\ \frac{\partial f_2}{\partial x} |_{\mathbf{x}_e} & \frac{\partial f_2}{\partial \dot{x}} |_{\mathbf{x}_e} & \frac{\partial f_2}{\partial \theta} |_{\mathbf{x}_e} & \frac{\partial f_2}{\partial \dot{\theta}} |_{\mathbf{x}_e} \\ \frac{\partial f_3}{\partial x} |_{\mathbf{x}_e} & \frac{\partial f_3}{\partial \dot{x}} |_{\mathbf{x}_e} & \frac{\partial f_3}{\partial \theta} |_{\mathbf{x}_e} & \frac{\partial f_3}{\partial \dot{\theta}} |_{\mathbf{x}_e} \\ \frac{\partial f_3}{\partial \dot{\theta}} |_{\mathbf{x}_e} & \frac{\partial f_4}{\partial \dot{\theta}} |_{\mathbf{x}_e} & \frac{\partial f_4}{\partial \dot{\theta}} |_{\mathbf{x}_e} & \frac{\partial f_4}{\partial \dot{\theta}} |_{\mathbf{x}_e} \end{bmatrix} \begin{bmatrix} \delta x \\ \delta \dot{x} \\ \delta \theta \\ \delta \dot{\theta} \end{bmatrix} + \begin{bmatrix} \frac{\partial f_1}{\partial F_1} |_{\mathbf{x}_e} & \frac{\partial f_1}{\partial F_2} |_{\mathbf{x}_e} \\ \frac{\partial f_2}{\partial F_1} |_{\mathbf{x}_e} & \frac{\partial f_2}{\partial F_2} |_{\mathbf{x}_e} \\ \frac{\partial f_3}{\partial F_1} |_{\mathbf{x}_e} & \frac{\partial f_3}{\partial F_2} |_{\mathbf{x}_e} \\ \frac{\partial f_4}{\partial F_1} |_{\mathbf{x}_e} & \frac{\partial f_4}{\partial F_2} |_{\mathbf{x}_e} \end{bmatrix} \begin{bmatrix} \delta F_1 \\ \delta F_2 \end{bmatrix}$$

$$(A.4)$$

$$\begin{bmatrix}
\delta \dot{x} \\
\delta \ddot{x} \\
\delta \dot{\theta} \\
\delta \ddot{\theta}
\end{bmatrix} = \begin{bmatrix}
0 & 1 & 0 & 0 \\
0 & 0 & \frac{mg}{M} & 0 \\
0 & 0 & 0 & 1 \\
0 & 0 & \frac{(m+M)}{ML}g & 0
\end{bmatrix} \begin{bmatrix}
\delta x \\
\delta \dot{x} \\
\delta \theta \\
\delta \dot{\theta}
\end{bmatrix} + \begin{bmatrix}
0 & 0 \\
\frac{1}{M} & -\frac{1}{M} \\
0 & 0 \\
-\frac{1}{ML} & \frac{2M+m}{mML}
\end{bmatrix} \begin{bmatrix}
\delta F_1 \\
\delta F_2
\end{bmatrix}$$
(A.5)

The eigenvalues are given by  $det(\lambda I - A) = \lambda^2 (\lambda^2 - \frac{(m+M)}{ML}g) = 0$ . Therefore, the system is unstable about  $\boldsymbol{x}_e$  due to the right half plane pole,  $\lambda = \sqrt{\frac{(m+M)}{ML}g}$ . Additionally, the time constant of this unstable system is  $\tau = \sqrt{\frac{ML}{g(m+M)}}$ . Note, if  $M >> m, \tau \to \sqrt{\frac{L}{g}}$ , which is the time constant for a simple pendulum.

It can be proved that the inverted pendulum system is controllable by showing:

$$rank[\mathbf{B} \mathbf{A} \mathbf{B} \mathbf{A}^2 \mathbf{B} \mathbf{A}^3 \mathbf{B}] = 4 \tag{A.6}$$

Therefore for any initial condition we can reach  $x_e$  in finite time under these linear assumptions.

# A.2 Propagation of Quantisation Error

The state space model for the quantisation of the linearised inverted pendulum can be written as:

$$\boldsymbol{x}_{t}^{(2D)} = C\boldsymbol{x}_{t} + \boldsymbol{V}_{t}$$
  $\boldsymbol{V}_{t} \sim \mathcal{U}\left(\begin{bmatrix} \frac{1}{\delta x} \\ \frac{1}{\delta \theta} \end{bmatrix}\right)$  (A.7)

$$\boldsymbol{x}_t = A\boldsymbol{x}_{t-1} + B\boldsymbol{u}_t \tag{A.8}$$

Where A and B are the linearised system dynamics (valid for small time steps), and C is the linear transformation to a 2D state space, with quantisation noise V.

Assuming the quantisation bin sizes,  $\delta x$  and  $\delta \theta$ , are small and that x and  $\theta$  are independent within the bin, the quantisation noise can be modelled as uniform random variables with covariance,  $cov(\mathbf{V}, \mathbf{V}) = \mathbb{E}[\mathbf{V}\mathbf{V}^T]$ :

$$= \mathbb{E} \begin{bmatrix} x^2 & x\theta \\ \theta x & \theta^2 \end{bmatrix} = \begin{bmatrix} \int_{-\delta x/2}^{\delta x/2} x^2 \cdot \frac{1}{\delta x} dx & 0 \\ 0 & \int_{-\delta \theta/2}^{\delta \theta/2} \theta^2 \cdot \frac{1}{\delta \theta} d\theta \end{bmatrix} = \begin{bmatrix} \frac{\delta x^2}{12} & 0 \\ 0 & \frac{\delta \theta^2}{12} \end{bmatrix}$$
(A.9)

For simplicity, let  $\delta x = \delta \theta$ , and therefore,  $cov(\mathbf{V}, \mathbf{V}) = \sigma_v^2 I$ .

Kalman filtering can be used to find an optimal estimate  $\hat{\boldsymbol{x}}_n = K[\boldsymbol{x}_n | \boldsymbol{y}_{n-k:n}]$  using the algorithm (derived in [3]).

## Algorithm 6 Multivariate Kalman Filtering

1: Given: 
$$\hat{\boldsymbol{x}}_n = K[\boldsymbol{x}_n | \boldsymbol{y}_{1:n}]$$
 and  $\Sigma_n = \mathbb{E}[(\boldsymbol{x}_n - \hat{\boldsymbol{x}}_n)(\boldsymbol{x}_n - \hat{\boldsymbol{x}}_n)^T]$ 

**Prediction:** 

2: 
$$\bar{\boldsymbol{x}}_{n+1} = K[\boldsymbol{x}_n | \boldsymbol{y}_{1:n}] = A\hat{\boldsymbol{x}}_n + B\boldsymbol{u}_{n+1}$$
  
3:  $\bar{\Sigma}_{n+1} = \mathbb{E}[(\boldsymbol{x}_{n+1} - \bar{\boldsymbol{x}}_{n+1})(\boldsymbol{x}_{n+1} - \bar{\boldsymbol{x}}_{n+1})^T] = A\Sigma_n A^T + \Sigma_w$   $\triangleright \Sigma_w = \mathbf{0}$ 

Update:

4: 
$$\tilde{\boldsymbol{y}}_{n+1} = \boldsymbol{y}_{n+1} - C\bar{\boldsymbol{x}}_{n+1}$$
  $\Rightarrow$  Calculate Innovation residual  
5:  $S_{n+1} = \Sigma_v + C\bar{\Sigma}_{n+1}C^T$   $\Rightarrow$  Calculate Innovation Covariance  
6:  $\hat{\boldsymbol{x}}_{n+1} = \bar{\boldsymbol{x}}_{n+1} + \bar{\Sigma}_{n+1}C^TS_{n+1}^{-1}\tilde{\boldsymbol{y}}_{n+1}$   $\Rightarrow \Sigma_v = \sigma_v^2I$   
7:  $\Sigma_{n+1} = (I - \bar{\Sigma}_{n+1}C^TS_{n+1}^{-1}C)\bar{\Sigma}_{n+1}$ 

Thus we can find the optimal linear estimate of the covariance of Equation (A.7) from algorithm 6 line 7:

$$\Sigma_{n+1} = (I - \bar{\Sigma}_{n+1} C^T S_{n+1}^{-1} C) \bar{\Sigma}_{n+1}$$

$$= (I - \bar{\Sigma}_{n+1} C^T (\sigma_v^2 I + C \bar{\Sigma}_{n+1} C^T)^{-1} C) \bar{\Sigma}_{n+1}, \qquad where \ \bar{\Sigma}_{n+1} = A \Sigma_n A^T \quad (A.11)$$

For the univariate case this reduces to:

$$\sigma_{n+1}^2 = \bar{\sigma}_{n+1}^2 \left(1 - \frac{C^2 \bar{\sigma}_{n+1}^2}{\sigma_v^2 + C^2 \bar{\sigma}_{n+1}^2}\right) \tag{A.12}$$

$$= A^{2}\sigma_{n}^{2}\left(1 - \frac{C^{2}A^{2}\sigma_{n}^{2}}{\sigma_{v}^{2} + C^{2}A^{2}\sigma_{n}^{2}}\right)$$
(A.13)

It is obvious from this that as  $\sigma_v^2 \to 0$ ,  $\sigma_{n+1}^2 \to 0$ . Furthermore, if the spectral radius,  $\rho\left(A^2(1-\frac{C^2A^2\sigma_n^2}{\sigma_v^2+C^2A^2\sigma_n^2})\right) < 1$ , then the mean squared error from quantisation will reduce to zero. Since  $0 < ||\frac{C^2A^2\sigma_n^2}{\sigma_v^2+C^2A^2\sigma_n^2}||_2^2 < 1$  if  $\sigma_v^2 > 0$ , a sufficient condition is that  $A^2 \le 1 \implies \rho(A) \le 1$ .

Using Gelfand's Theorem,  $\rho(M_1...M_n) \leq \rho(M_1)...\rho(M_n)$ , and the eigenvalue identity,  $(M+cI)v = (c+\lambda)v$ . The MSE of the multivariate case can be reduced to:

$$1 > \rho \left( (I - \bar{\Sigma}_{n+1} C^T (\sigma_v^2 I + C \bar{\Sigma}_{n+1} C^T)^{-1} C) \bar{\Sigma}_{n+1} \right)$$
(A.14)

$$> \rho(\bar{\Sigma}_{n+1}) \left[ 1 - \rho(\bar{\Sigma}_{n+1}C^T(\sigma_v^2 I + C\bar{\Sigma}_{n+1}C^T)^{-1}C) \right]$$
 (A.15)

Therefore, if  $\rho\left(\bar{\Sigma}_{n+1}C^T(\sigma_v^2I + C\bar{\Sigma}_{n+1}C^T)^{-1}C\right)$  is less than one, and  $\rho(\bar{\Sigma}_{n+1}) < 1$ , then this is true. This is equivalent to showing that if  $0 < \kappa(\Sigma_n)\kappa(A)^2\kappa(CC^T) \le 1$  and  $\rho(AA^T) < 1$ , where  $\kappa(\cdot)$  denotes the condition number, then the MSE will decay:

$$\rho\left(\bar{\Sigma}_{n+1}C^{T}(\sigma_{v}^{2}I + C\bar{\Sigma}_{n+1}C^{T})^{-1}C\right)\right) \leq \rho(\bar{\Sigma}_{n+1})\rho(CC^{T})\rho((\sigma_{v}^{2}I + C\bar{\Sigma}_{n+1}C^{T})^{-1})$$

$$\leq \rho(\bar{\Sigma}_{n+1})\rho(CC^{T})\frac{1}{|\lambda_{min}(\sigma_{v}^{2}I + C\bar{\Sigma}_{n+1}C^{T})|}$$

$$= \rho(\bar{\Sigma}_{n+1})\rho(CC^{T})\frac{1}{|\lambda_{min}(C\bar{\Sigma}_{n+1}C^{T})| + \sigma_{v}^{2}}$$

$$\leq \frac{|\lambda_{max}(\bar{\Sigma}_{n+1})\lambda_{max}(CC^{T})|}{|\lambda_{min}(C\bar{\Sigma}_{n+1}C^{T})|}$$

$$\leq \frac{|\lambda_{max}(\bar{\Sigma}_{n+1})\lambda_{max}(CC^{T})|}{|\lambda_{min}(\bar{\Sigma}_{n+1}C^{T})|}$$

$$\leq \frac{|\lambda_{max}(\bar{\Sigma}_{n+1})\lambda_{max}(CC^{T})|}{|\lambda_{min}(\bar{\Sigma}_{n+1})\lambda_{min}(CC^{T})|}$$

$$= |\kappa(\Sigma_{n})|\kappa(A)^{2}\kappa(CC^{T})$$

where the above uses the results  $\lambda_i(X\Sigma X^*) = \lambda_i(\Sigma X^*X)$ , where  $\Sigma$  is symmetric (this equality also holds for spectral radii) and  $\lambda_{min}(\bar{\Sigma}_{n+1}CC^T) > \lambda_{min}(\bar{\Sigma}_{n+1})\lambda_{min}(CC^T)$ . Note that  $\kappa(\Sigma_n) = 1$ . Therefore sufficient conditions for the decay of the covariance are:

(1) 
$$\rho(AA^T) \le 1 \implies |\lambda_{max}(AA^T)| \le 1 \tag{A.16}$$

(2) 
$$\kappa(A)^2 \le 1 \implies |\lambda_{min}(A)| = |\lambda_{max}(A)|$$
 (A.17)

$$(3) \kappa(CC^T) \le 1 (A.18)$$

For the linearised dynamics derived in eq. (A.5) the eigenvalues of  $AA^T$ ,  $\lambda$ , are given by  $\lambda^2(1-\lambda)^2=0$ , and the eigenvalues of A are  $\lambda=(0,0,\pm\sqrt{\frac{(m+M)}{ML}g})$ . Therefore (1) and (2) are true.

Therefore, the covariance decays to zero if  $\rho(CC^T) \leq 1$  in the linearised multivariate case and the 2D state becomes a lossless estimate of the state. Given that C is a 2x4 projection matrix of the form

$$C = \begin{bmatrix} c_{11} & 0 & 0 & 0 \\ 0 & 0 & c_{23} & 0 \end{bmatrix} \implies CC^{T} = \begin{bmatrix} c_{11}^{2} & 0 \\ 0 & c_{23}^{2} \end{bmatrix}$$
 (A.19)

By the choice of  $\boldsymbol{x}^{(2D)}$  as square,  $c_{11} = c_{23}$ . By the same logic as above,  $\kappa(CC^T) = 1$ . The rate of decay for the univariate case can be determined with a first order approximation about  $\sigma_v^2 = 0$ :

$$\sigma_{n+1}^2(\sigma_v^2) = A^2 \sigma_n^2 \left(1 - \frac{C^2 A^2 \sigma_n^2}{\sigma_v^2 + C^2 A^2 \sigma_n^2}\right)$$
(A.20)

$$\approx \sigma_{n+1}^2(0) + \delta \sigma_v^2 \frac{\partial \sigma_{n+1}^2(\sigma_v^2)}{\partial \sigma_v^2} |_{\sigma_v^2 = 0}$$
(A.21)

$$= \frac{\delta \sigma_v^2}{(AC\sigma_n)^2} \tag{A.22}$$

# A.3 Neural Network Losses

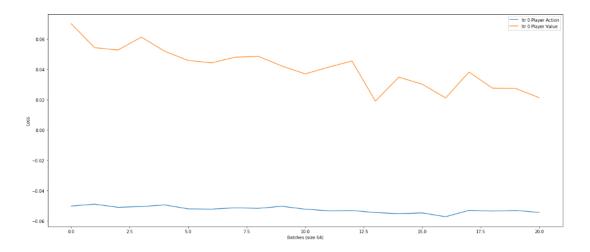


Figure A.1: The losses for both the player and adversary neural networks after one set of training examples.

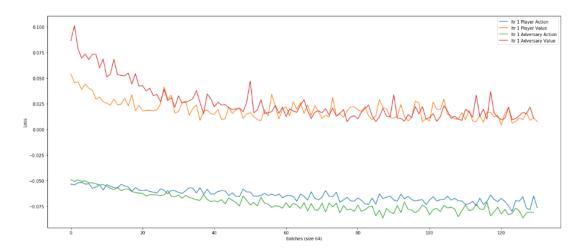


Figure A.2: The losses for both the player and adversary neural networks after two sets of training examples.

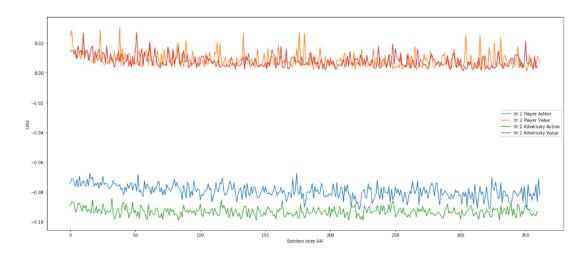


Figure A.3: The losses for both the player and adversary neural networks after three sets of training examples.

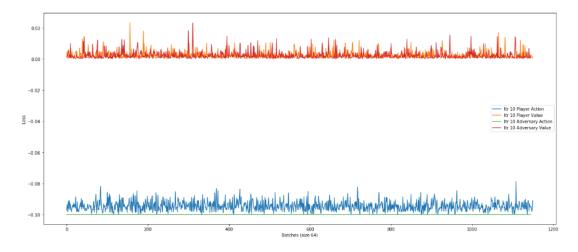


Figure A.4: The losses for both the player and adversary neural networks after ten sets of training examples.

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