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Génie Logiciel

RN quantifié pour un contexte deep learning embarqué

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Résumé

Les réseaux de neurones sont devenus très utilisés dans l'intelligence artificielle grâce à leurs performances de prédictions.
Ces réseaux de neurones sont parfois très complexes, et il n'est pas pratique de les utiliser dans des systèmes à ressources limités.

Pour cela, dans ce rapport nous allons introduire les réseaux de neurones binaires (BNNs), les formaliser, étudier quelques modèles binarisés et dériver leurs formules, implémenter les résultats trouvés dans une bibliothèque qu'on va nommer **binaryflow**.
Et finalement, nous allons implémenter et entraîner des modèles binarisés pour la classification et régression, et nous allons comparer les performances de prédiction, ainsi que les performances temps et mémoire de ces BNNs et leurs contrepart classique.

Remerciement

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Contenu

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Liste des Figures

Liste des Tableaux

List of Algorithms

Introduction

Avec l'explosion de l'intelligence artificielle, et surtout les modèles d'apprentissage profonds, la complexité des modèles a subi une croissance considérable, qui les rend inexploitable dans les systèmes à complexité limitée.

Dans ce rapport, nous allons étudier la quantification des paramètres et des entrées des couches du réseau de neurones sur un seul bit.

Ce rapport va détailler notre approche de la formalisation des BNNs, de l'analyse et généralisation des approches existantes, vers l'implémentation d'une bibliothèque unifiant les BNNs, et son utilisation. Il est composé de 6 chapitres.

Dans le premier chapitre, nous allons présenter la societé **dB Sense** et sa méthodologie.

Dans le deuxième chapitre, nous allons donner une petite histoire de l'apprentissage profond, et puis poser le problème de la grande complexitée de ces modèle, en posant la binarisation comme une solution.

Dans le troisième chapitre, nous allons formaliser notre approche, en définissant les BNNs. Après nous allons poser quelques problèmes dans l'entraînement. Après nous allons proposer les optimisations temps et mémoire qu'on peut exploiter avec les BNNs.
Finalement, nous allons proposer l'algorithme d'entraînement et d'interférence des BNNs.

Dans le quatrième chapitre, nous allons étudier et analyser quelques BNNs répandus dans la littérature, en conformant avec notre définition proposée.

Dans le cinquième chapitre, nous allons implémenter la bibliothèque **binaryflow**, en justifiant les paradigmes utilisés.

Dans le sixième chapitre, nous allons utiliser les différents BNNs étudiés sur 3 jeux de données.
Pour chacun de ces modèles nous allons analyser son performance de prédiction, sa taille mémoire au déploiement, et une estimation sur la complexité de son interférence en calculant le nombres d'instruction équivalents.

Chapter 1

Cadre du Stage

Introduction

Chapter 2

Analysis & Implementation

2.1 Introduction

2.2 Formalisation

To define a Mean Payoff Game, we will start by formalising a weighted di-graph[1].

2.2.1 Di-Graph

A Weighted Di-Graph $\mathcal{G}$ is a tuple $(\mathcal{V}, \mathcal{E}, \mathcal{W})$ where:

- $\mathcal{V}$ is the set of vertices.
- $\mathcal{E} \subseteq V \times V$ is the set of edges.
- $\mathcal{W} : \mathcal{E} \to \mathbb{G}$ is the weight function, assigning a weight for every edge, with $\mathbb{G}$ some ordered abelian group[2].

2.2.2 Mean Payoff Game

Formally, a **Mean Payoff Graph** is a tuple $(\mathcal{V}, \mathcal{E}, \mathcal{W}, \mathcal{P}, s, p)$ where:

- $\mathcal{G} = (\mathcal{V}, \mathcal{E}, \mathcal{W})$ is a di-graph.
- $s \in \mathcal{V}$ denotes the starting position.
- $\mathcal{P} = \{\text{Max}, \text{Min}\}$ is the set of players.
- $p \in \mathcal{P}$ the starting player

A **Mean Payoff Game** is a perfect information, zero-sum, turn based game played indefinitively on a Mean Payoff Graph as follow:

- The game starts at $u_0 = s$, with player $p_0 = p$ starting.
- For each $n \in \mathbb{N}$, Player p_n will choose a vertex $u_{n+1} \in \operatorname{Adj} u_n$, with a payoff $w_n \mathcal{W}(u_n, u_{n+1})$
- The winner of the game will be determined by the Mean Payoff. There are different winning conditions.

[1] Directed Graph

[2] This definition is too general. We will only consider $\mathbb{G} \in \{\mathbb{Z}, \mathbb{Q}, \mathbb{R}\}$. Also, $\mathbb{G}$ itself should be clear from the context.

Condition C1

Player Max wins **iff**:

$$\liminf_{n\in\mathbb{N}^*} \frac{1}{n}\sum_{k=0}^{n-1} w_k \geqslant 0$$

Otherwise, Player Min will win

Condition C2

Player Max wins **iff**:

$$\liminf_{n\in\mathbb{N}^*} \frac{1}{n}\sum_{k=0}^{n-1} w_k > 0$$

Otherwise, Player Min will win.

Condition C3

Player Max wins **if**:

$$\liminf_{n\in\mathbb{N}^*} \frac{1}{n}\sum_{k=0}^{n-1} w_k > 0$$

Player Min wins **if**:

$$\limsup_{n\in\mathbb{N}^*} \frac{1}{n}\sum_{k=0}^{n-1} w_k < 0$$

Otherwise, it is a draw.

2.2.3 Well Foundness

It is not very clear from the definition that the game is well founded.
In fact, there are choices for which the mean payoff does not converge. That is the sequence $\left(\frac{1}{n}\sum_{k=0}^{n-1} w_k\right)_{n\in\mathbb{N}^*}$ does not converge.
One such example is the sequence defined by:

$$w_n = (-1)^{\lfloor \log_2(n+1) \rfloor}$$

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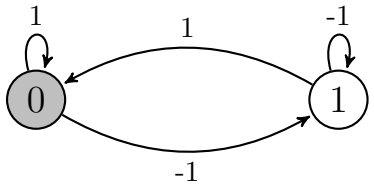
For that sequence, the $(2^r - 1)$-step mean payoff is equal to:

$$\begin{aligned}
\sum_{k=0}^{2^r-2} w_k &= \sum_{k=1}^{2^r-1} (-1)^{\lfloor \log_2(k) \rfloor} \\
&= \sum_{i=0}^{r-1} \sum_{j=2^i}^{2^{i+1}-1} (-1)^{\lfloor \log_2(j) \rfloor} \\
&= \sum_{i=0}^{r-1} \sum_{j=2^i}^{2^{i+1}-1} (-1)^{i} \\
&= \sum_{i=0}^{r-1} (2^{i+1} - 2^i)(-1^i) \\
&= \sum_{i=0}^{r-1} (-2)^i \\
&= \frac{1 - (-2)^r}{3} \\
\implies \frac{1}{2^r - 1} \sum_{k=0}^{2^r-2} w_k &= \frac{1}{3} \cdot \frac{1 - (-2)^r}{2^r - 1} \\
&= \frac{1}{3} \cdot \frac{2^{-r} - (-1)^r}{1 - 2^{-r}}
\end{aligned}$$

That sequence has two accumulation points $\pm\frac{1}{3}$, and thus, it does not converge.

On the other hand, the introduction of the supremum and infimum operators will solve the convergence problem, as the resulting sequences will become monotone.

An example of an execution that gives a rise to such payoffs is the following Meab Payoff Game instance[3]:



(a) Representation of the Mean Payoff Game

Pair of strategies defined as:

$$\begin{aligned}
\Phi : & V^+ \times P \to V \\
& (s_0 \dots s_r, p) \to B(r) \bmod 2
\end{aligned}$$

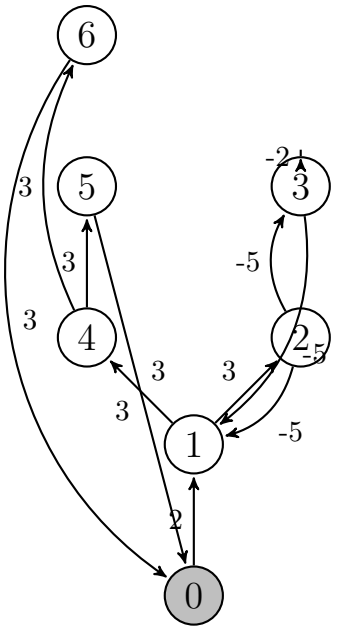
With $B(r)$ the position of the left-most bit in the binary representation of r

(b) Definition of both strategies

Figure 2.1: An example of an execution with non-convergent Mean Payoffs

2.2.4 Symmetries

[3]Note that the proposed pair of strategies is odd in the sense that it appears that both players cooperated on the construction of non-convergent mean payoffs instead of trying ot win the game.



2.2.5 Strategy

Deterministic Strategies

Let p be a player.
A (deterministic) strategy[4] is a function $\Pi^p : V^+ \to V$ such that:

$$\forall v_0 \dots v_r \in V^+, \quad \Pi_p(v_0 \dots v_r) \in \operatorname{Adj} v$$

If the strategy does only depend on the current vertex, we say it is a memoryless (deterministic) strategy $\Pi : V \to V$

Probabilistic Strategies

A probabilistic strategy is a random process that assigns for each sequence of vertices $v \in \mathcal{V}$ a probability distribution over $\operatorname{Adj} v$. This constitutes the most general strategy of a player:

$$\forall v_0 \dots v_r \in V^+, \quad \Pi_p(v_0 \dots v_r) \in \mathscr{D}(\operatorname{Adj} v)$$

Considered Strategies

Strategies that depends in complete past histories are in general intractable. For Mean Payoff Game, it is proven that the optimal strategy is a **deterministic** and **memoryless**.
For that we will only consider **memoryless** strategies. And for the scope of this report:

- A deterministic strategy should refer to memoryless deterministic strategy.
- A probabilistic strategy should refer to memoryless probabilistic strategy.
- A strategy should refer to memoryless deterministic strategy.

[4]By default, we refer to deterministic strategies. If the strategy is not deterministic, we will explicit it.

We will still consider (memoryless) probabilistic strategies as they reside in a smooth space, and thus they can be used for machine learning purposes.

Deterministic Optimal Strategy

There are three kinds of optimality:

Weak Optimality : In the deterministic case, a strategy Φ of player $p \in P$ is weakly optimal if one of the following is true:

- For each strategy Φ^p of player p, player $\bar{p}$ can win the game by finding a countering strategy $\Phi^{\bar{p}}$.
- Player p will not lose the game no matter his opponent's strategy

Strong Optimality : In the deterministic case, a strategy Φ of player $p \in P$ is strongly optimal if one of the following is true:

- For each strategy Φ^p of player p, player $\bar{p}$ can win or tie the game by finding a countering strategy $\Phi^{\bar{p}}$.
- Player p will win the game no matter his opponent's strategy

Payoff Optimality : In the deterministic case, a strategy Φ of player $p \in P$ is payoff optimal if independently of $\bar{p}$'s strategy it:

- Maximises the Mean Payoff if $p = \text{Min}$
- Minimises the Mean Payoff otherwise

Now we have the following hiearchy considering the set of optimal strategies:

$\forall$Mean Payoff Game $G, \forall p \in P, \quad \texttt{PayoffOptimal}(G,p) \subseteq \texttt{StrongOptimal}(G,p) \subseteq \texttt{WeakOptimal}(G,p)$

2.3 Evaluating Strategies

Suppose we have a pair of potentially probabilitic strategies $(\Phi^{\text{Max}}, \Phi^{\text{Min}})$. The problem is to evaluate the winner without doing an infinite simulation of the game.

2.3.1 Monte-Carlo Simulations

This is the most intuitive evaluation method

2.3.2 Memoryless Deterministic Strategies

If both strategies are deterministic and memoryless. Then the generated sequence of vertices $(s_n)_{n\in\mathbb{N}}$ will be completely determined by the recurrence relation:

$$s_n = \begin{cases} s & \text{if } n = 0 \\ \Phi^{\text{Max}}(s_{n-1}) & \text{if } n \text{ is odd} \\ \Phi^{\text{Min}}(s_{n-1}) & \text{otherwise} \end{cases}$$

This can be represented in the compact form:

$$\forall n \in \mathbb{N}^*, \quad (s_n, p_n) = (\Phi^{p_{n-1}}(s_{n-1}), \bar{p}_{n-1}) = F(s_{n-1}, p_{n-1})$$

Since $\mathcal{V} \times \mathcal{P}$ is a finite set and F is a function, such sequence will be eventually periodic, that is:

$$\exists N \in \mathbb{N}, \exists T \in \mathbb{N}^* / \quad \forall n \in \mathbb{N}_{\geqslant N}, \quad (s_n, p_n) = (s_{n+T}, p_{n+T})$$

We can calculate its eventual period using the turtle hare algorithm.

Now, the mean payoff will be equal to the mean of weights that appears on the cycle. This can be proven as follow.:

$$\begin{aligned}
S_{aT+b+N} &= \sum_{k=0}^{aT+b+N-1} w_k \\
&= \sum_{k=0}^{N-1} w_k + \sum_{k=0}^{aT+b-1} w_{k+N} \\
&= \sum_{k=0}^{N-1} w_k + a \sum_{r=0}^{T-1} w_{r+N} + \sum_{r=0}^{b-1} w_{r+N} \\
\implies \left| S_{n+N} - \lfloor \frac{n}{T} \rfloor \sum_{r=0}^{T} w_{k+N} \right| &\leqslant (N+T-1) \max_{(u,v)\mathcal{E}} |\mathcal{W}(u,v)| \\
&\leqslant (N+T-1) \|\mathcal{W}\|_\infty \\
\implies \left| \frac{1}{n+N} S_{n+N} - \frac{1}{n+N} \cdot \lfloor \frac{n}{T} \rfloor \sum_{r=0}^{T-1} w_{k+N} \right| &\leqslant \frac{N+T-1}{n+N} \|\mathcal{W}\|_\infty
\end{aligned}$$

Now it can be proven that:

$$\lim_{n \to +\infty} \frac{1}{n+N} \cdot \lfloor \frac{n}{T} \rfloor \sum_{r=0}^{T} w_{k+N} = \frac{1}{T} \sum_{r=0}^{T-1} w_{r+N}$$

With that:

$$\lim_{n \to +\infty} \frac{1}{n} \sum_{k=0}^{n-1} w_k = \frac{1}{T} \sum_{r=0}^{T-1} w_{r+N} \quad \blacksquare$$

Now, our algorithm will be composed of 3 main parts:

- Calculating the transition function $F : V \times P \to V \times P$. This is straightforward from the construction.
- Calculating the period and the offset of the sequence. We will use Floyd's cycle finding algorithm for that.
- Calculating the Mean Payoff

This is an illustrative implementation of our algorithm.

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Algorithm 1 Deterministic strategies evaluation

Require: $G = (V, E, P, s, p)$ a mean payoff game
Require: $(\Phi^{\text{Max}}, \Phi^{\text{Min}})$ the edge probability
Ensure: R The mean payoff
  $F \leftarrow \text{Transition}(G, \Phi^{\text{Max}}, \Phi^{\text{Min}})$ ▷ Calculate the transition function
  $x_0 \leftarrow (s, p)$
  $(T, r) \leftarrow \text{FloydCycleFinding}(F, x_0)$ ▷ Find the period and the offset
  $S \leftarrow 0$ ▷ S represents the cumulative payoffs along a cycle
  $x \leftarrow x_0$
  for $k \in \{1, \ldots, r\}$ **do** ▷ Advance until arriving to the cycle
    $x \leftarrow F(x)$
  end for
  for $k \in \{1, \ldots, T\}$ **do**
    $y \leftarrow\!\leftarrow F(x)$
    $u \leftarrow \underset{V \times P \to V}{\text{projection}}(x)$ ▷ Extracts the current vertex
    $v \leftarrow \underset{V \times P \to V}{\text{projection}}(y)$ ▷ Extracts the next vertex
    $S \leftarrow S + W(u, v)$
    $x \leftarrow y$
  end for
  return $R \leftarrow \frac{S}{T}$

2.3.3 Probabilistic Strategies

Due to the undeterministic nature of probabilistic strategies, it does not make sense to evaluate the mean payoffs, as different executions may lead to different mean payoffs.
Instead, probabilistic strategies gives rise to a discrete distribution of mean payoffs.
Now two closely related, but different evaluations are possible

- Expected Mean Payoff
- Distribution of winners

Now, with both strategies fixed. A Mean Payoff Game can be considered as a Markov Chain.

2.4 Countering Strategies

By fixing the strategy of player pP to Φ^p, then the Markov Game Process reduces to a Markov Decision Process, which can be solve by Linear Programming tools
Moreover, in a deterministic Mean Payoffs, a counter strategy can be calculated efficiently by reducing the problem to finding a negative cycle in a di-graph.

2.4.1 Deterministic Counter Strategy

For simplicity, we assume that $p = \text{Max}$, the same results apply for Min player.
For a Mean Payoff Game (V, E, P, s, p'), we introduce the following graph:

$$G' = \left(V, E', W'\right)$$

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where:

- E' is defined as follow:
$$E' = \{(u, \Phi^p(v)), \quad (u, v) \in E\}$$
- Also, W' is defined adequately:
$$\forall (u, v) \in E : W'(u, \Phi^p(v)) = W(u, v) + W(v, \Phi^p(v))$$

The problem of finding a counter strategy will be reduced to finding a negative cycle $\mathcal{C}$ in G'
This can be done in $\mathcal{O}(|V|^3)$

2.4.2 Probabilistic Counter Strategy

On the other hand, if Φ^p is probabilistic, then the problem can be reduced to optimizing the mean payoff of an infinite-horizon Markov Decision Process.

2.5 Learning Strategies

2.5.1 Iterative Methods

Chapter 3

Dataset Generation

3.1 Introduction

3.2 Analysis

Generating a Mean Payoff Game can be decomposed into two subsequent objectives.

1. Generate the Graph itself.
2. Generate the Weights

3.3 Graph Distributions

There are many well studied graph distributions in the litterature.
One of the most explored ones are the $\mathcal{G}(n,p)$ and $\mathcal{G}(n,m)$ families.

3.3.1 $\mathcal{G}(n,p)$ Family

For $n \in \mathbb{N}, p \in [0,1]$, a graph G is said to follow a $\mathcal{G}(n,p)$ distribution if $|V| = n$ and:

$$\forall e \in \mathscr{E}, \quad \mathscr{P}(s \in \mathcal{E}) = p$$

Where $\mathscr{E}$ is a set of valid edges.

3.3.2 $\mathcal{G}(n,m)$ Family

For $n \in \mathbb{N}, m \in \mathbb{N}$, a graph G is said to follow a $\mathcal{G}(n,m)$ distribution if $|V| = n, |E| = m$ and the edges $e_1, \ldots, e_m$ were drawn from a set of valid edges $\mathscr{E}$.

3.3.3 Valid edges

The set of valid edges $\mathscr{E}$ is the set defining the potential edges of the graph. It is equal to:

1. $V \times V$ for directed graphs with loops
2. $(V \times V) \backslash V \odot V$ for directed graphs without loops
3. The set of subsets of size 2 of V denoted $\mathscr{P}_2(V)$ for undirected graphs with loops.
4. The set of subsets of size 2 of V denoted $\mathscr{P}_2(V)$ for undirected graphs with loops.

3.3.4 $\mathcal{D}(n,p)$ Graph Construction

Naive Method

The definition of $\mathcal{D}(n,p)$ gives a straightforward construction.
This is achieved by flipping a coin[1] for each pair of node $(u,v) \in V^2$, we add an edge if we get a Head.
This is implemented in the following algorithm:

Algorithm 2 $\mathcal{D}(n,p)$ Graph Generation

Require: $n \in \mathbb{N}^*$ the size of the graph
Require: $p \in \mathbb{N}^*$ the edge probability
Ensure: $G \sim \mathcal{D}(n,p)$
  $A : (u,v) \in V \times V \to 0$
  for $u \in V$ **do**
    for $v \in V$ **do**
      Generate $X \sim \mathcal{B}(p)$ ▷ $\mathcal{B}(p)$ is the bernoulli distribution
      $A(u,v) \leftarrow X$
    end for
  end for
  return $G \leftarrow$ `GraphFromAdjacencyMatrix`(A)

The complexity[2] of the following algorithm is $\mathcal{O}(n^2)$.

Optimized Method

Instead of iterating over all possible pair of nodes. For each vertex $v \in V$:

- We can sample a number d from the outgoing degree distribution[3]
- We then choose d numbers uniformly without replacement from an indexable representation of V

The following algorithm implements the optimized method:

Algorithm 3 $\mathcal{D}(n,p)$ Graph Generation Optimisation

Require: $n \in \mathbb{N}^*$ the size of the graph
Require: $p \in \mathbb{N}^*$ the edge probability
Ensure: $G \sim \mathcal{D}(n,p)$
  $A : u \in V \to \varnothing$
  for $u \in V$ **do**
    Generate $d \sim \mathcal{B}(n,p)$ ▷ d represents the degree, $\mathcal{B}(n,p)$ is the binomial distribution
    $A(u) \leftarrow \text{choice}(V,d)$
  end for
  return $G \leftarrow$ `GraphFromAdjacencyList`(A)

[1] The coin is potentially biased with a probability of obtaining head equal to $p \in [0,1]$
[2] We assume the cost of generating a Bernoulli random variable as $\mathcal{O}(1)$
[3] Or the ingoing degree distribution, they are in fact equal.

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Now, Let $C(n, m)$ be the cost of choice function. The expected complexity of this algorithm will be:

$$\tilde{\mathcal{O}}\left(n\mathbb{E}_d[C(n,d)]\right) \quad \text{where } d \sim \mathcal{B}(n,p)$$

We will show on the next section what choice function should we use.

3.3.5 Choice Function

First Proposition

We propose here a simple choice algorithm, but it is still efficient for our use case.
It works simply by drawing without replacement, but we ignore duplicate elements. This is implemented as follow

Algorithm 4 $\mathcal{D}(n,p)$ Choice without replacement

Require: S a list
Require: $m \in \{0, \dots |S|\}$ the number of chosen elements
Ensure: H a set of size m containing uniformly drawn elements without replacement.
  $H \leftarrow \varnothing$
  while $|H| < m$ **do**
    Generate $v \sim \mathcal{U}(S)$ ▷ Where $\mathcal{U}(S)$ is the uniform distribution over S
    $H \leftarrow H \cup \{v\}$
  end while
  return H

To estimate the cost of this algorithm, we will use probabilistic reasoning.
Let $X_{n,m} = C(n, m)$ the running time of an execution of algorithm 4 in a set S of size n, with m elements to be chosen. We have:

$$\begin{aligned}
&X_{n,0} \text{ is deterministic} \\
&X_{n,0} = \mathcal{O}(1) \\
\mathbb{E}[X_{n,m}] &= 1 + \frac{1}{n}\sum_{k=0}^{n-1} \mathbb{E}[X_{n,m} \mid \text{The last drawn number is } k] \\
&= 1 + \frac{1}{n}\sum_{k=0}^{m-2} \mathbb{E}[X_{n,m}] + \frac{1}{n}\sum_{k=m-1}^{n-1} \mathbb{E}[X_{n,m-1}] \\
&= 1 + \frac{m-1}{n}\mathbb{E}[X_{n,m}] + \frac{n-m+1}{n}\mathbb{E}[X_{n,m-1}]
\end{aligned}$$

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Now we arrived at a recurrent formula. We will simplify it as shown below:

$$\begin{aligned}
\frac{n-m+1}{n}\mathbb{E}[X_{n,m}] &= \frac{n-m+1}{n}\mathbb{E}[X_{n,m-1}] + 1 \\
\implies \mathbb{E}[X_{n,m}] &= \frac{n-m+1}{n-m+1}\mathbb{E}[X_{n,m-1}] + \frac{n}{n-m+1} \\
&= \mathbb{E}[X_{n,m-1}] + \frac{n}{n-m+1} \\
&= \sum_{k=1}^{m} \frac{n}{n-k+1} + \mathcal{O}(1) \\
&= \sum_{k=0}^{m-1} \frac{n}{n-k} + \mathcal{O}(1) \\
&= n \sum_{k=n-m+1}^{n} \frac{1}{k} + \mathcal{O}(1) \\
&= n(H_n - H_{n-m}) + \mathcal{O}(1)
\end{aligned}$$

Here $(H)_{n\in\mathbb{N}*}$ is the harmonic series, and we define $H_0 = 0$.

Complexity

The expected complexity of algorithm 4 depends on both n and m:

- If $m = o(n)$, then it is $\tilde{\mathcal{O}}(m)$.
- If $m = kn + o(n)$ with $k \in]0, 1[$, then it is $\tilde{\mathcal{O}}(m)$.
- If $m = n - o(n)$, It is[4] $\tilde{\mathcal{O}}(m \log m)$.

To prove this result, we use a well known asymptotic approximation of the Harmonic series[5]:

$$H_n = \ln n + \gamma - \frac{1}{2n} + \mathcal{O}\left(\frac{1}{n^2}\right)$$

We can prove this claim as follow:

$$\begin{aligned}
m = o(n), \quad \mathbb{E}[C(n,m)] &= -n \ln\left(1 - \frac{m}{n}\right) - \frac{1}{2}\left(1 - \frac{n}{n-m}\right) + \mathcal{O}\left(\frac{1}{n}\right) \\
&= m + o(m) \\
&= \mathcal{O}(m) \\
m = km + o(m), k \in]0,1[, \quad \mathbb{E}[C(n,m)] &= -n \ln\left(1 - \frac{m}{n}\right) - \frac{1}{2}\left(1 - \frac{n}{n-m}\right) + \mathcal{O}\left(\frac{1}{n}\right) \\
&= -n \ln(1 - k + o(1)) + \frac{1}{n}(1 - \tfrac{1}{1-k+o(1)}) + \mathcal{O}(\tfrac{1}{n}) \\
&= \mathcal{O}(m)
\end{aligned}$$

[4] Here we use the minus sign to emphasize that $m \leqslant n$

[5] This asymptotic approximation can be proven using the Euler–Maclaurin formula

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For $m = n - o(n)$, we prove it by noting that:

$$\begin{aligned}\mathbb{E}[C(n,m)] &\leqslant \mathbb{E}[C(n,n)] \\ &\leqslant nH_n \\ &\leqslant n\ln n + \gamma n + \frac{1}{2} + \mathcal{O}\left(\frac{1}{n}\right) \\ &= \mathcal{O}(m\log m)\end{aligned}$$

Refinement

If m tends to n, it is more hard to select m elements from a set of size n without replacement. This explains the extra logarithmic factor.
In that case, we can instead focus on the dual problem: "Find the $n-m$ elements that will not be selected". This can be calculated in $\mathcal{O}(n-m)$.
Once we find the elements that will not be selected, their set complement are exactly the m elements that will be selected. This new algorithm is guaranteed to be $\mathcal{O}(m)$ irrespective of n and m

Algorithm 5 Fine tuned $\mathcal{D}(n,p)$ Choice without replacement

Require: S a list
Require: $m \in \{0,\dots|S|\}$ the number of chosen elements
Require: choice The choice function defined on algorithm 4
Require: τ a fine tuned threshold. We will use $\tau = \frac{1}{2}$ for all practical purposes.
Ensure: H a set of size m containing uniformly drawn elements without replacement.
  if $\frac{m}{|S|} \leqslant \tau$ **then**
    $H \leftarrow \text{choice}(V,n)$
  else
    $H \leftarrow S \backslash \text{choice}(S, n-m)$
  end if
  return H

Also, an important point is that by combining the analysis of all possible cases, we can extract a constant factor that is independent of n. So that the Big-O notation is only a function of m.

3.3.6 Complexity of Optimised $\mathcal{D}(n,p)$ Graph Construction

We return to evaluate the asymptotic behaviour of $\mathbb{E}_d[C(n,d)]$.
Let $\delta \in \mathbb{R}_+^*$
The Chebychev Inequality implies that:

$$\mathscr{P}\left(\left|\frac{1}{n}\sum_{i=1}^{n} X_i - p\right| \geqslant \frac{\delta}{\sqrt{n}}\sqrt{p(1-p)}\right) \leqslant \frac{1}{\delta^2}$$

By setting: $\delta = \frac{1}{\sqrt{p}}$, we have:

$$\mathscr{P}\left(\left|\frac{1}{n}\sum_{i=1}^{n} X_i - p\right| \geqslant \frac{\sqrt{1-p}}{\sqrt{n}}\right) \leqslant p$$

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Let $I = [np - \sqrt{n(1-p)}, np + \sqrt{n(1-p)}]$.
We have:

$$\begin{aligned}\mathbb{E}\left[C(n,d)\right] &\leqslant \mathbb{E}\left[C(n,d) \mid d \in I\right] + p^2\mathbb{E}\left[C(n,d) \mid d \notin I\right] \\ &\leqslant \max_{m\in\mathbb{N}\cap I} C(n,m) + \max_{m\in\{0,\dots,n\}} p^2 C(n,m)\end{aligned}$$

To further simplify this, we need two key observations:

- The interval $m \in \mathbb{N} \cap [np - 1 + p, np + 1 - p]$ contains at most 3 integers, all of which are within a distance of 1 to np
- The complexity of $\mathbb{E}[C(n,m)]$ does only depend on m.

Thus, we have the following:

$$\begin{aligned}\max_{m\in\mathbb{N}\cap[np-1+p,np+1-p]} C(n,m) &= \mathcal{O}(np) \\ \max_{m\in\{0,\dots,n\}} p^2 C(n,m) &= \mathcal{O}(np^2)\end{aligned}$$

Now, by combining both estimations, we get:

$$\tilde{\mathcal{O}}(n\mathbb{E}_d[C(n,d)]) = \tilde{\mathcal{O}}(n^2 p) \quad \blacksquare$$

3.3.7 $\mathcal{D}(n,m)$ Construction

To construct a random $\mathcal{D}(n,m)$ graph, we only have to select m uniformly random elements from the set $V \times V$.
We will use algorithm 5 for this purpose[6]:

Algorithm 6 Fine tuned $\mathcal{D}(n,p)$ Choice without replacement

Require: $n \in \mathbb{N}^*$
Require: $m \in \{0,\dots,n^2\}$ the number of chosen elements
Ensure: $G \sim \mathcal{D}(n,m)$
  $E \leftarrow \text{choice}(\text{Lazy}(V) \times \text{Lazy}(V), m)$ ▷ We only need the m elements on-demand.
  return $G \leftarrow \text{GraphFromEdges}(E)$ ▷ This justifies using Lazy

Here $\text{Lazy}(V) \times \text{Lazy}(V)$ is a lazy implementation of cartesian product that supports bijective indexing[7] over $\{0,\dots,n^2-1\}$.
The complexity of this construction is: $\tilde{\mathcal{O}}(m)$

3.4 Sinkless Conditionning

Sampling from a graph distribution may lead to graphs that have at least one sink.
These graphs are problematic as Mean Payoff Graphs are exactly the sinkless graphs.
To migitate this, we will impose a conditionning on both distribution that will gives a guaranteed

[6] It is essential that the list $V \times V$ be lazy loaded. In particular, each element will only be loaded when it is indexed. This is essential to reduce the complexity. Otherwise, we will be stuck in an $\mathcal{O}(n^2)$ algorithm.
[7] Indexing is required for uniform sampling

Mean Payoff Graph.
We will explore such conditionning both distribution:

- $\mathcal{G}^S(n,p)$: This is the distribution of graphs following $\mathcal{G}(n,p)$ with the requirement that they do not have a sink.
- $\mathcal{G}^S(n,m)$: This is the distribution of graphs following $\mathcal{G}(n,m)$ with the requirement that they do not have a sink.

3.4.1 Repeating Construction

Algorithm

This method is very intuitive. It will repeat the sampling until getting the desired graph.
The following is an implemention of the repeating construction.

Algorithm 7 Fine tuned $\mathcal{D}(n,p)$ Choice without replacement

Require: $n \in \mathbb{N}^*$
Require: $m \in \{0, \dots |S|\}$ the number of chosen elements
Require: choice The choice function defined on algorithm 4
Require: Threshold τ
Ensure: H a set of size m containing uniformly drawn elements without replacement.
  if $\frac{m}{|S|} \leqslant \tau$ **then**
    $H \leftarrow \text{choice}(V, n)$
  else
    $H \leftarrow V \setminus \text{choice}(S, n-m)$
  end if
  return H

Analysis

We will analyse the runtime of generating a $\mathcal{G}^S(n,p)$.
We expect a similar runtime for $\mathcal{G}^S(n,m)$ due to the similarity between $\mathcal{G}(n,m)$ and $\mathcal{G}(n,p)$.
Let $F(n)$

3.5 Weights Distribution

3.5.1 Construction

Once the graph is constructed. We only have to generate the weights.
This will be done by creating a random weight function:

$$W(u,v) : (u,v) \to W_{u,v}$$

Here $W_{u,v}$ will be a sequence of real random variables.
In our case, we set $(W_{u,v})_{(u,v)\in E}$ to be independent and identically distributed over a real distribution $\mathcal{W}$.

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3.6 Proposed MPG Distribution

3.6.1 Desired Properties of Mean Payoff Game Distributions

Fairness in the Limit

This is essential, as we intend to generate a sequence of Mean Payoff Games that do not favour statistically a certain player, in the sense that, if we generate sufficient independent and identically distributed Mean Payoff Games $G_1, \dots, G_n$, we expect the following:

$$\lim_{n \to +\infty} |\mathtt{R}_{\text{Max}}(G_1, \dots, G_n) - \mathtt{R}_{\text{Min}}(G_1, \dots, G_n)| = 0$$

Where $\mathtt{R}$ is defined as follow:

$$\mathtt{R}_{\text{Op}}(G_1, \dots, G_n) = \frac{1}{n} \sum_{i=1}^{n} \mathscr{P}(\text{Op wins } G_i \text{ assuming optimal strategies})$$

Symmetric

A real distribution is said to be symmetric if:

$$\forall [a, b] \in \mathbb{R}, X \sim \mathcal{W}, \quad \mathscr{P}(X \in [a, b]) = \mathscr{P}(X \in [-b, -a])$$

We will define a symmetric Mean Payoff Game distribution as a distribution of Mean Payoff Game whose weights are independent and identically distributed on a symmetric real distribution. This property is stronger than Fairness in the Limit, as it implies that:

$$\mathscr{P}(\text{Max wins } G \text{ assuming optimal strategies}) = \mathscr{P}(\text{Min wins } G \text{ assuming optimal strategies})$$

We will require a Symmetric Mean Payoff Game as we do not want a player to have an inherit advantage other the other one[8]

3.6.2 Implemented Distributions

The following table resumes the implemented distributions:

[8] Other than the first move.

Distribution Family	Parameters	Type
$\mathcal{D}(n,p)$	• n : Graph size • p : Edge probability	Graph distribution
$\mathcal{D}(n,m)$	• n : Graph size • m : Number of edges	Graph distrbiution
$\mathcal{U}_{\text{discrete}}(-r,r)$	• r : The radius of the support	Weight distribution
$\mathcal{U}(-r,r)$	• r : The radius of the support	Weight distribution
$\mathcal{N}(0,\sigma)$	• σ : The standard deviation	Weight distribution

Table 3.1: Le tableau d'avancement des BNNs

3.7 MPG Generation

3.7.1 Distribution

- Each generated graph will follow a distribution $\mathcal{G}(n,p(n))$ for some $n \in \mathbb{N}^*$
- The weights will follow the discrete uniform distribution $\mathcal{D}(-1000,1000)$

We will generate two kinds of datasets, depending on the nature of the graph

3.7.2 Dense Graphs

- Let $\mathcal{P} = \{0.1, 0.2, 0.3, 0.5, 0.7, 0.8, 0.9, 1\}$
- $\mathcal{N} = \{10, 20, 30, 40, 50, 60, 70, 80, 90, 100, 110, 120, 130, 140, 150, 200, 250, 300, 400, 500\}$
- For each $(n,p) \in \mathcal{N} \times \mathcal{P}$, we will generate $K = 1000$ observations $G_1^{n,p}, \ldots, G_K^{n,p} \sim \mathcal{G}(n,p)$

The total number of examples is:

$$K \times |\mathcal{N}| \times |\mathcal{P}| = 160000$$

The generation was done on a 'haswell64' partition with 24 cores. and it took 02:12:38 hours.

3.8 Annotation

3.8.1 Approach

We used the CSP algorithm 12 to annotate the dataset, potentially augmented with some heuristics. We implemented a program that takes the path of the dataset, and solves the Mean Payoff Games one by one.
To maximize efficiency, the program launches many solver threads, with each one independently working on a single file, and the results are accumulated using a ConcurrentQueue.

3.8.2 Target Values

The solver will calculate the following targets:

- The optimal pair of strategies
- The mean payoffs for each starting position, turn.
- The winners for each starting position, turn.

Also, some additional metadata are generated for analysis:

- `dataset`: The name of the whole dataset
- `graph`: The name of the graph.
- `status`: The solver's status on the given graph. In particular, whether it succeeded to solve the instance or not[9]. Equal to "OK" if the execution is successful.
- `running_time`: The time needed to solve the instance.

3.8.3 Heuristics

To accelerate the annotation of the two datasets, we had to apply some heuristics to the algorithm. We made essentially two kinds of heuristics.

Linear Bound Heuristic

This is the heuristic based on the view that for almost all solutions of a Ternary Max Atom system extracted from our generated random games, either:

- All variables are infinite:
$$X(u) = -\infty \quad \forall u \in V$$
- The diameter of assignments is in the order of $\|W\|_\infty$
$$\Delta X = \sup_{u \in V} X(u) - \inf_{u \in V, X(u) > -\infty} X(u) = \mathcal{O}(\|W\|_\infty)$$

This heuristic suggests a much tighter search space to the worst case $\|W\|_1$ one. Going further, with uniform random weights:

$$\|W\|_1 = \mathcal{O}\left(|E| \times \|W\|_\infty\right)$$

We believe this heuristic arises due to the random property of graphs, because in general, one can build an infinite family of ternary max atom systems that violate this heuristic.
In fact, going further, one can build ternary max atom systems were the $\|W\|_1$ estimation is tight.

To generating the dataset, we applied this heuristic with $\Delta X = 4\|W\|_\infty$

$$D = \{-\infty, -2\|W\|_\infty, , -2\|W\|_\infty + 1, \ldots, 2\|W\|_\infty\}$$

[9] We expect that the solver may crash due to several reasons (corrupted file, out of memory, etc...). For that we made additional effort for exception handling, so that an error for a single instance does not propagate to the whole program.

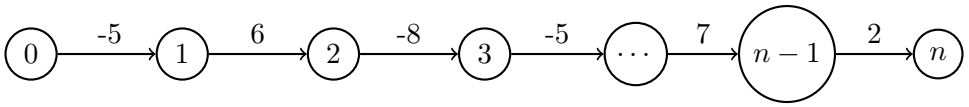

Figure 3.1: A counter example of the Linear Bound heuristic

Early Stopping

If after any iteration of arc consistency, $\max_{x \in V} \nu(x) < \sup D$. Then, $\nu(t)$ will converge to $-\infty$ for all t.
Thus, we stop the algorithm and sets $\nu(t) \leftarrow -\infty, \quad \forall t$

Proof suppose that in fact there is an assignment with:

$$-\infty < \max_{u \in V} \nu(u) < \sup D$$

We will take the u with the biggest such $\nu(u)$.
Now our system is a tropical max atom system, which means translations are also a polymorphism of this system, so for any assignment $\nu : V \to \mathbb{Z},\ X + t$ is also an assignment $\forall t \in \mathbb{Z}$. With that, $\nu + \sup D - \nu(u)$ is also an assignment.
This assignment has the property:

$$\forall s \in V, \quad \nu(s) + \sup D - \nu(u) \in D$$

Which is a contradiction, as it violates the consistency of arc consistency, and the maximality of the solution with respect to the domain D

The efficiency of the Early Stopping heuristic depends on the density of the graph. Empirically, for dense graphs. the analoguous ternary max atom system has two kind of assignments:

1. Either all variables are finite
2. Either all variables are $-\infty$

This translates back in a dense Mean Payoff setting, that the winner of the game does not depend in the starting position and the starting turn.
With that, the Early Stopping heuristic will quickly detect the second case, which is the usually hurdle of the algorithm.
On the other hand, for sparse graphs, we do not have this nice distinction between finite and infinite assignments, and they can overlap, and so will make this heuristic useless in practice.

3.8.4 Deployment

After some experiments, it was very clear that vertical scaling with the number of threads is not sufficient. By analysing the running time of some samples, we estimated the total running time solving both datasets to exceed 30 days.
As a result of this, we deployed a pipeline of 24 nodes, each with 24 threadss working simultaneously on a partition of the dataset.

Chapter 4

Model Design

4.1 Introduction

4.2 Objectives

The main objective

4.3 Properties

We expect the model $\mathcal{M}$ to verify the following properties:

4.3.1 Totality

Definition

A function is total if it is defined in the whole domain.
Also, we will say a model $\mathcal{M}$ is total if it works on all Mean Payoff Games..

Importance

This property implies two main characteristics:

1. The model works for any Mean Payoff Game whatever the size of $|V|$. This is difficult to not violate, as most machine learning models require static shapes.

2. The model works for any Mean Payoff Game whatever the representation is; This is implicitely verified by an encoding of the Mean Payoff Games.

4.3.2 Node Agnostic

This property tells that a model should not use any extra information about the nodes.

Formalisation

Let $G_1(V_1, E_1, W_1, s_1), G_2(V_2, E_2, W_2, s_2)$ two isomorphic Mean Payoff Games in the sense that there exists a bijection $\Phi : V_1 \to V_2$ such that:

$$\begin{aligned} E_2 &= \{(\Phi(u), \Phi(v)), \quad (u, v) \in E_1\} \\ \forall (u, v) \in E_1, \quad W_2(u, v) &= W_1(\Phi(u), \Phi(v)) \\ s_2 &= \Phi(s_1) \end{aligned}$$

Then:

$$\Phi\left(\mathcal{M}(G_1)\right) = \mathcal{M}(G_2)$$

Explanation

The property tells that if two graphs mean payoffs only differ by their node representation, then the results should also only differ by the representation of the nodes.

Importance

This property implies that we can simply encode a mean payoff as $G(V, E, W, s)$ as $G'(V', E', W', s')$ with $V' = \{0, ;|V| - 1\}$, and E', W', s' defined accordingly.
In fact, the encoding is done implicitely by our model[1].

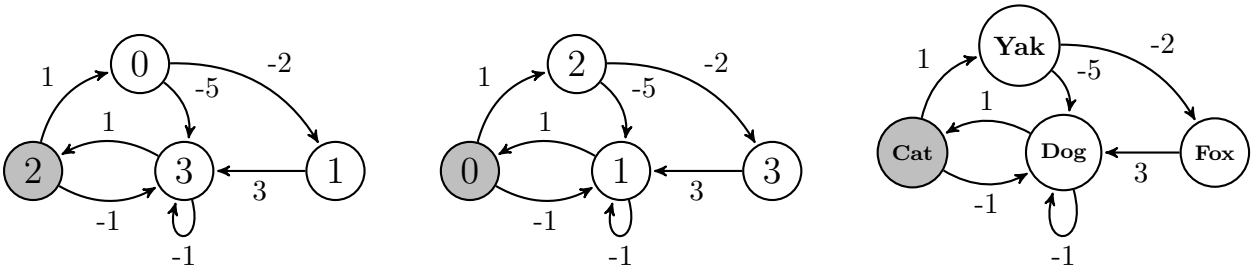

Figure 4.1: Three isomorphic Mean Payoff Games

In the figure 4.1, we have 3 equivalent Mean Payoff Games. To illustrate a node agnostic model:

- Let $\mathcal{M}$ be a model that predicts for a given Mean Payoff Game, the action to be played at the starting node, and the winner.
- Suppose also that $\mathcal{M}(G_1) = (3, \text{Max})$

If $\mathcal{M}$ is node agnostic, then:

$$\begin{aligned} \mathcal{M}(G_2) &= (1, \text{Max}) \\ \mathcal{M}(G_3) &= (\text{Dog}, \text{Max}) \end{aligned}$$

[1]While the encoding is done implicitely, the model itself can violate this property. An example of this is a Multi Layer Perceptron. Which lead to different results for different encodings.

4.3.3 Invariance under Positive Scaling

This property comes directly from the fact that both the winner and the set of optimal strategies[2] are invariant under a positive scaling of the weights.
Now, it is very easy to make augment a model $\mathcal{M}$ into such invariant model $\mathcal{M}'$. We only do the following:

$$\mathcal{M}'(E, W, s) = \mathcal{M}'(E, \text{Normalize}(W), s)$$

Where Normalize is any endomorphism of weight functions that verify the following constraint:

$$\forall_{\text{MPG}} G(E, W), \forall, \quad \text{Normalize}(W) = \frac{W}{H(W)}$$

With $H : \mathscr{F}(E, \mathbb{R}) \to \mathscr{F}(E, \mathbb{R})$ a function satisfying[3]:

$$\forall s \in \mathbb{R}_+^*, \quad H(sW) = sH(W)$$
$$H(W) > 0$$

Standard Scaling

This scaling treats the values of W as samples of random variables, and divides W by an estimate of their variance.

$$\text{Normalize}(W) = \frac{W}{\sqrt{\mathbb{V}[W]}}$$
$$\mathbb{V}[W] = \frac{1}{|E|} \sum_{(u,v)\in E} (W(u, v) - \mathbb{E}[W])^2$$
$$\mathbb{E}[W] = \frac{1}{|E|} \sum_{(u,v)\in E} W(u, v)$$

Maximum Scaling

This scaling reduces the interval of the weights to $[-1, 1]$ by dividing by the largest weight in terms of absolute value:

$$\text{Normalize}(W) = \frac{W}{\|W\|_\infty}$$

Implementation Notes: The weights function is implemented as a matrix W, which is equal to 0 for $(u, v) \notin E$.
It is important to igonre these zeros[4] in both normalisations, otherwise it may lead to biased scaling.

[2] Whatever the definition of optimality (Weak, Strong, Payoff).
[3] Special care must be when H is $\geqslant$ instead of $>$
[4] And only these zeros. If $W(u, v) = 0$ for $(u, v) \in E$, this term should be accounted.

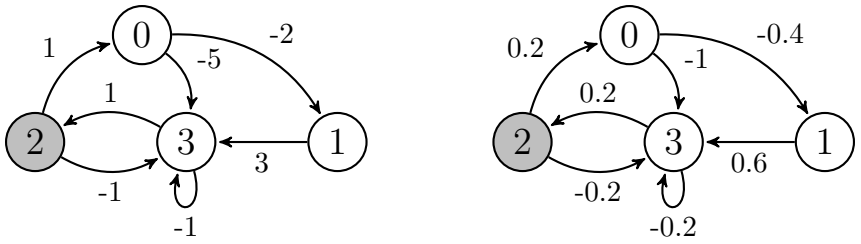

Figure 4.2: Mean Payoff Game, with a rescaled version using Maximum normalization

4.3.4 Permutation Equivariance

Definition

This property states that the model should ouptut the same results for permuted nodes, up to permutation of the results.

Importance

While it is a special case of the Node Agnostic property, but it is still important as the encoding of the graphs is done implicitely, and thus Permutation Equivariance is enough to get a Node Agnostic model.

4.3.5 Stability under Padding

Definition

Let $G_1(V_1, E_1, W_1, s_1), G_2(V_2, E_2, W_2, s_2)$ be two disjoint Mean Payoff Games, in the sense that $V_1 \cap V_2 = \varnothing$.
The padding of G_1 by G_2, denoted $G_1 \triangleright G_2$ defined as:

$$G_1 \triangleright G_2 = (V_1 \cup V_2, E_1 \cup E_2, W_1 \cup W_2, s_1)$$

A model is said to be stable under padding if:

$$\mathcal{M}(G_1 \triangleright G_2) = \mathcal{M}(G_1)$$

Importance

Deep Learning algorithms generally accept batches of data having a homogeneous shape.
In the other hand, graph input generally has different shapes, and thus are problematic to most learning algorithms.
While we succeeded in experimenting a learning algorithm with a ragged batch[5], it suffered the following limits:

- It greatly limits the choice of potential models.
- The training is not supported by GPU, as some core operations were only implemented in the CPU for ragged batches.

[5] A batch of inhomogeneous input

For this reasons, we opted to pad the graphs to get homogeneous batches, and this is why the stability under padding is important.

Removing Unreachable Nodes

Another major point for this property, is it gives the possibility to remove unreachable nodes[6] without affecting the model's results.

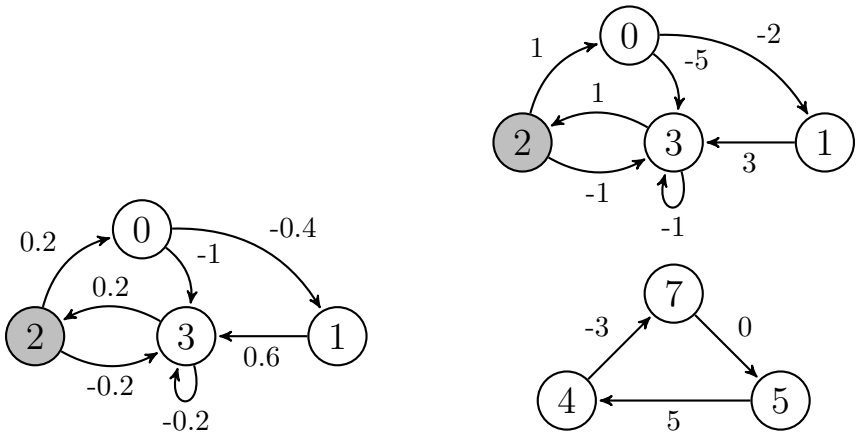

Figure 4.3: Original Mean Payoff Game with a padded version

Following the example under figure 4.3, if the model $\mathcal{M}$ is stable under padding, then the triangular subgraph should not be accounted in the output.

4.4 Considered Models

While there are many possible predictive models. We considered mainly two families of predictive models.

4.4.1 Value based Model

Such model predicts the evaluation of a certain position.
The evaluation is a function $\mathcal{M} : V \times P \to [-1, 1]$ with the following interpretation:

- $\mathcal{M}(s, p) = 1$ when the model predicts that player p will win the game given the position.
- $\mathcal{M}(s, p) = -1$ whens the model predicts that player p will lose the game given the position.
- $\mathcal{M}(s, p) = 0$ when the model predicts that the position will result in a draw.

Now, while such a model can be used to predict the winner. We claim that it is not efficient[7] to extract the strategy from such model, as the information of winning alone does not directly induce a strategy.
The following figure illustrates an example.

[6] If we ignore the orientation. A stronger property allowing unreachable nodes on the di-graph can also be defined, but this is beyond the scope.

[7] Efficient in the sense that we cannot extract the best strategy from the evaluation in linear time.

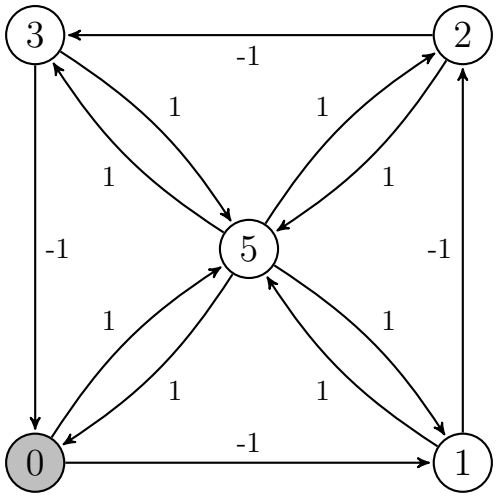

Figure 4.4: An MPG where the first player always wins.

4.4.2 Strategy based Model

Such model predicts the strategy for each player.
This is a function $\mathcal{M} : V \times P \to \mathscr{P}(V)$ such that:

$$\begin{aligned} \forall (u,p) \in V \times P, \quad & \mathcal{M}(u,p) \in \operatorname{Adj} u \\ \forall (u,p) \in V \times P, \forall v \in \operatorname{Adj} u, \quad & \mathcal{P}\left(\mathcal{M}(u,p) = v\right) \in [0,1] \\ \forall (u,p) \in V \times P, \quad & \sum_{v \in \operatorname{Adj} u} \mathcal{P}\left(\mathcal{M}(u,p) = v\right) = 1 \end{aligned}$$

4.5 Building the Model

Our model follows from well esta../Figures/PreprocessingBlock.pngblished graph neural network architectures, with minor modifications to suit our needs.
It is composed of blocks, each containing a **Weighted Graph Convolutional Network**.

4.5.1 Weighted Graph Convolutional Network

The weighted graph convolutional network **WGCN**, as its name suggests, is a convolutional operator acting on graphs, with many desirable properties such as "Permutation Equivariance" and "Stability under Padding".
It is based on the graph convolutional network as described on the following figure

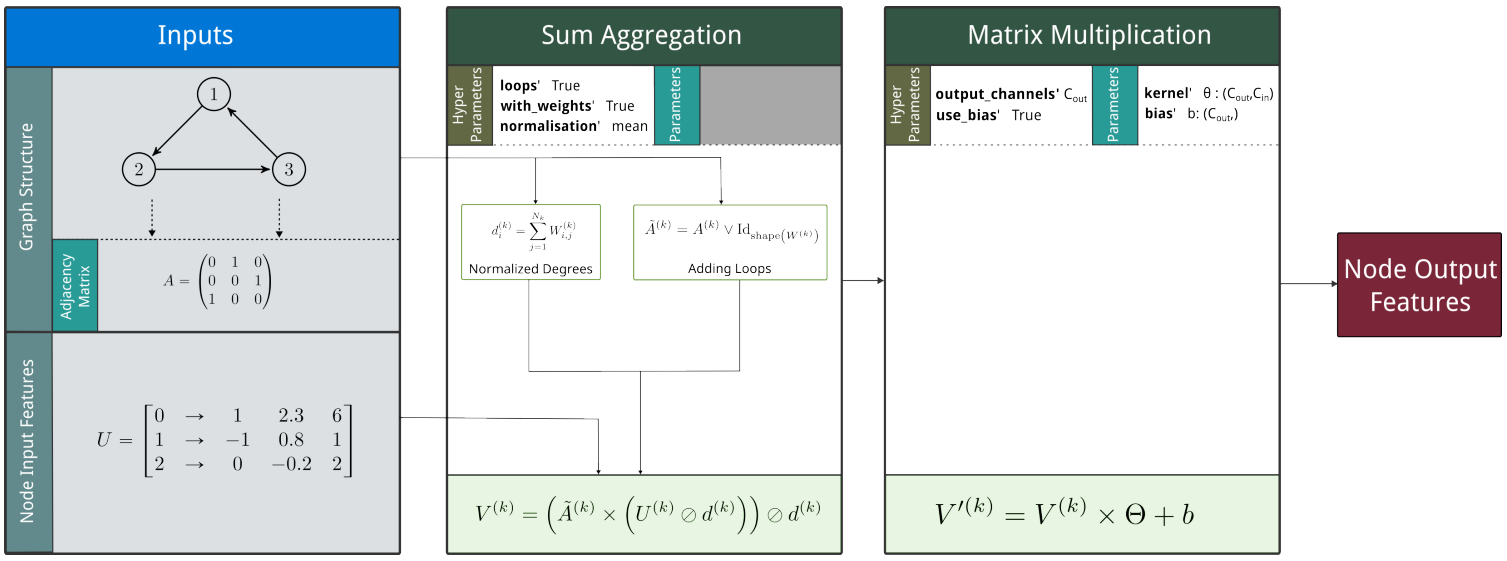

Figure 4.5: Graph Convolutional Network

The problem with **GCN** is that they ignore the weights information. For Mean Payoff Games, such information is crucial to determine the winner and the strategies.
For that, we introduced **WGCN** to capitalize on the weights information. The figure below shows how it is implemented.

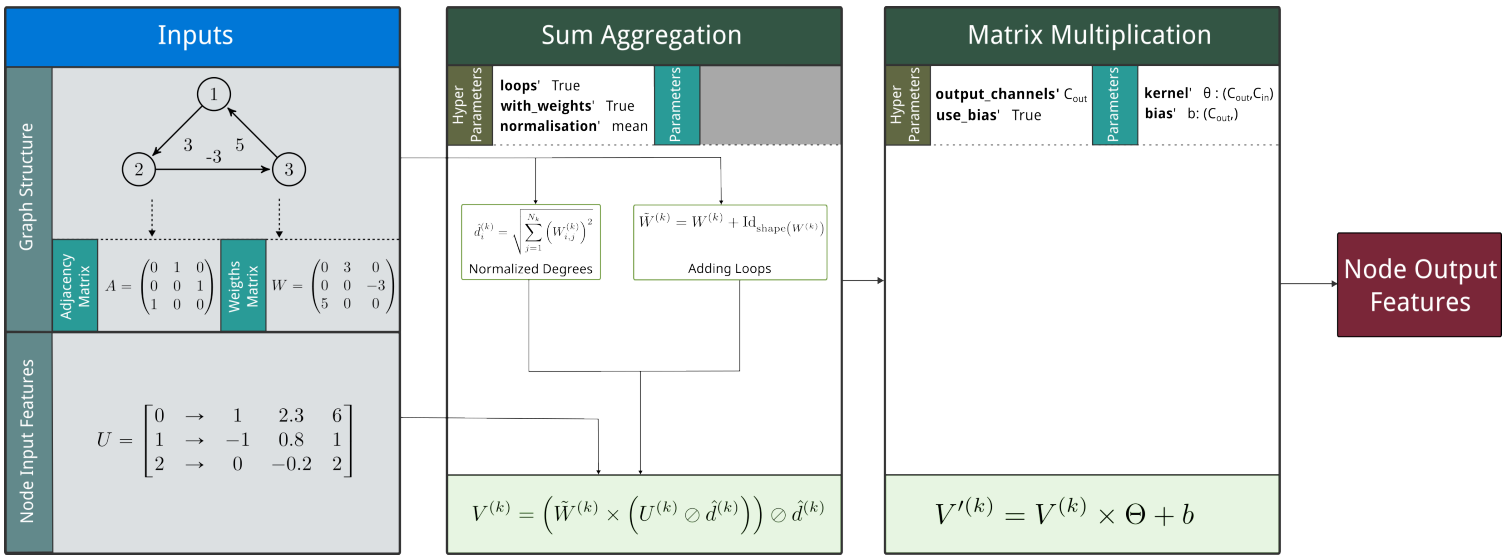

Figure 4.6: Weighted Graph Convolutional Network

The **WGCN** operator exhibits the desired properties described on 4.3, except property 4.3.3.

4.5.2 Preprocessing Block

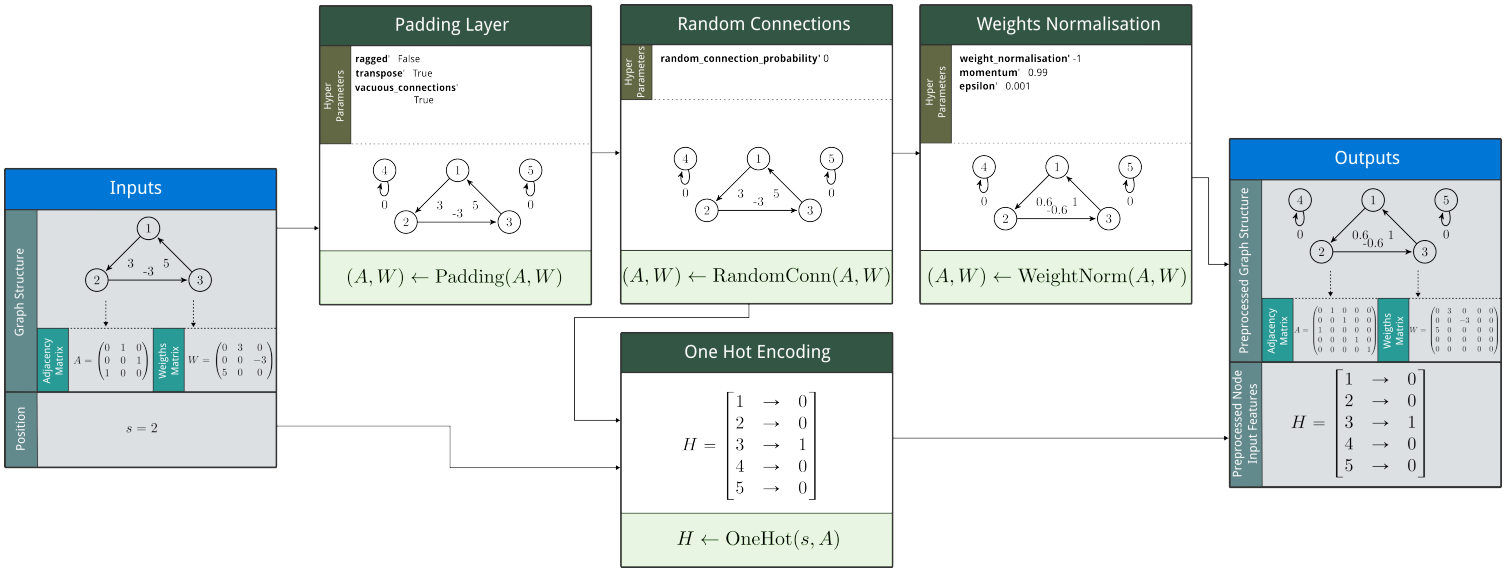

Figure 4.7: Preprocessing Block

4.5.3 Convolutional Block

Each intermediate block is composed of a graph convolution, and then a batch normalisation operator, as described by the figure below:

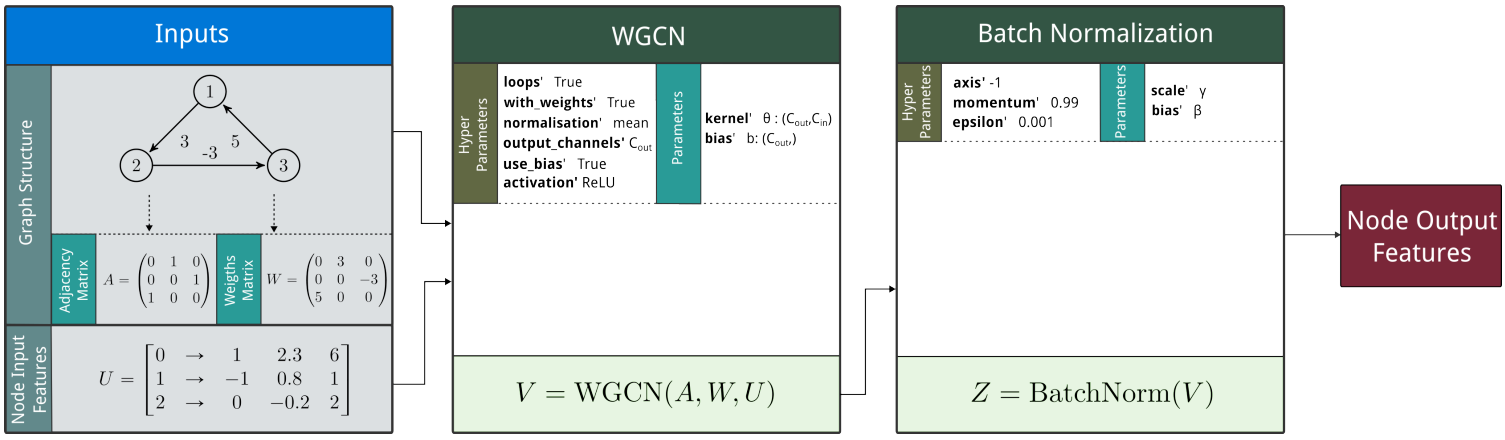

Figure 4.8: Graph Convolutional Block

4.5.4 Prediction Block

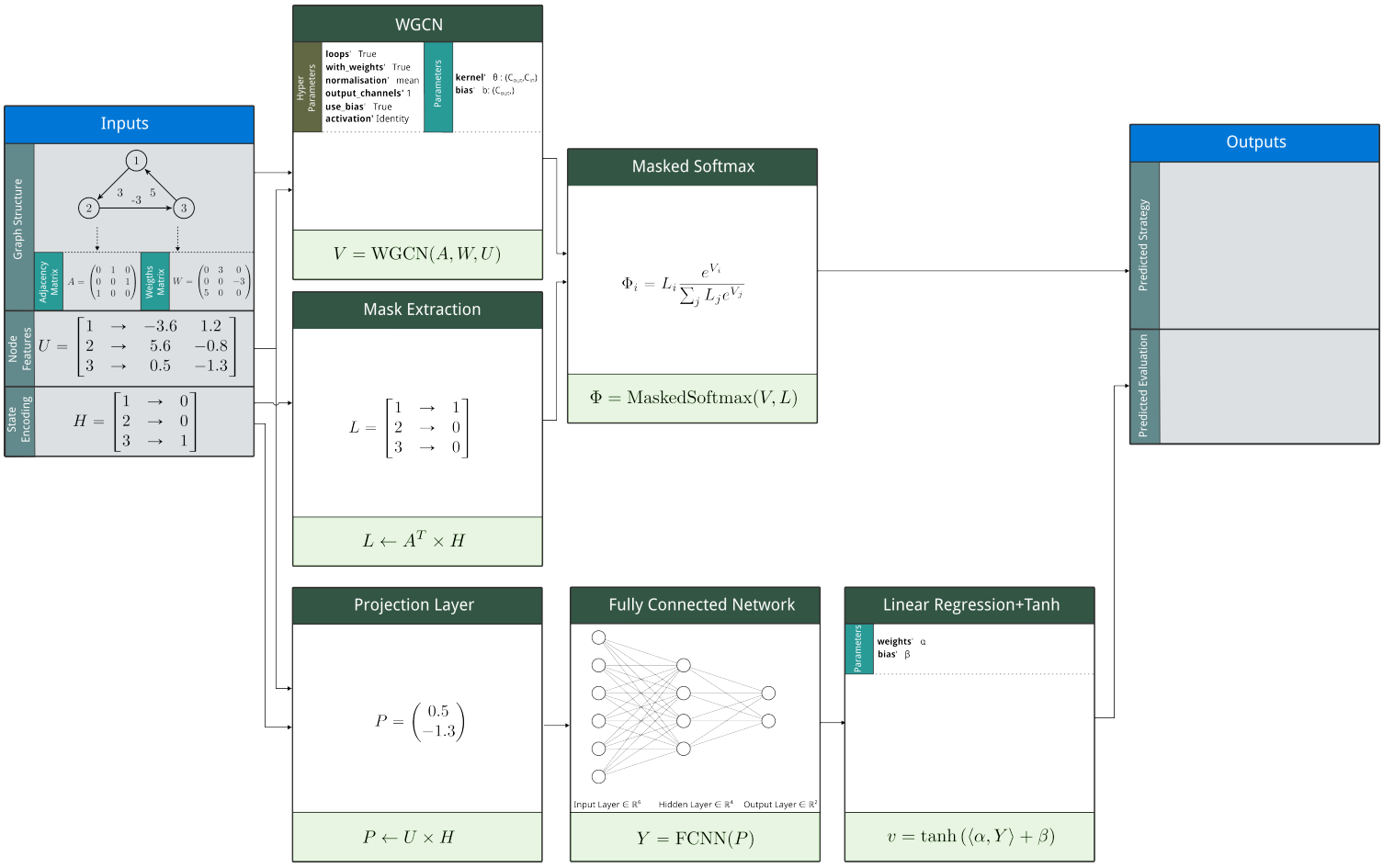

Figure 4.9: Graph Convolutional Block

4.5.5 Model Architecture

Putting all building block together, we designed a model that takes arbitrary graphs, encode them, predicts the strategy and the evaluation of the position.
This is detailed by the figure below:

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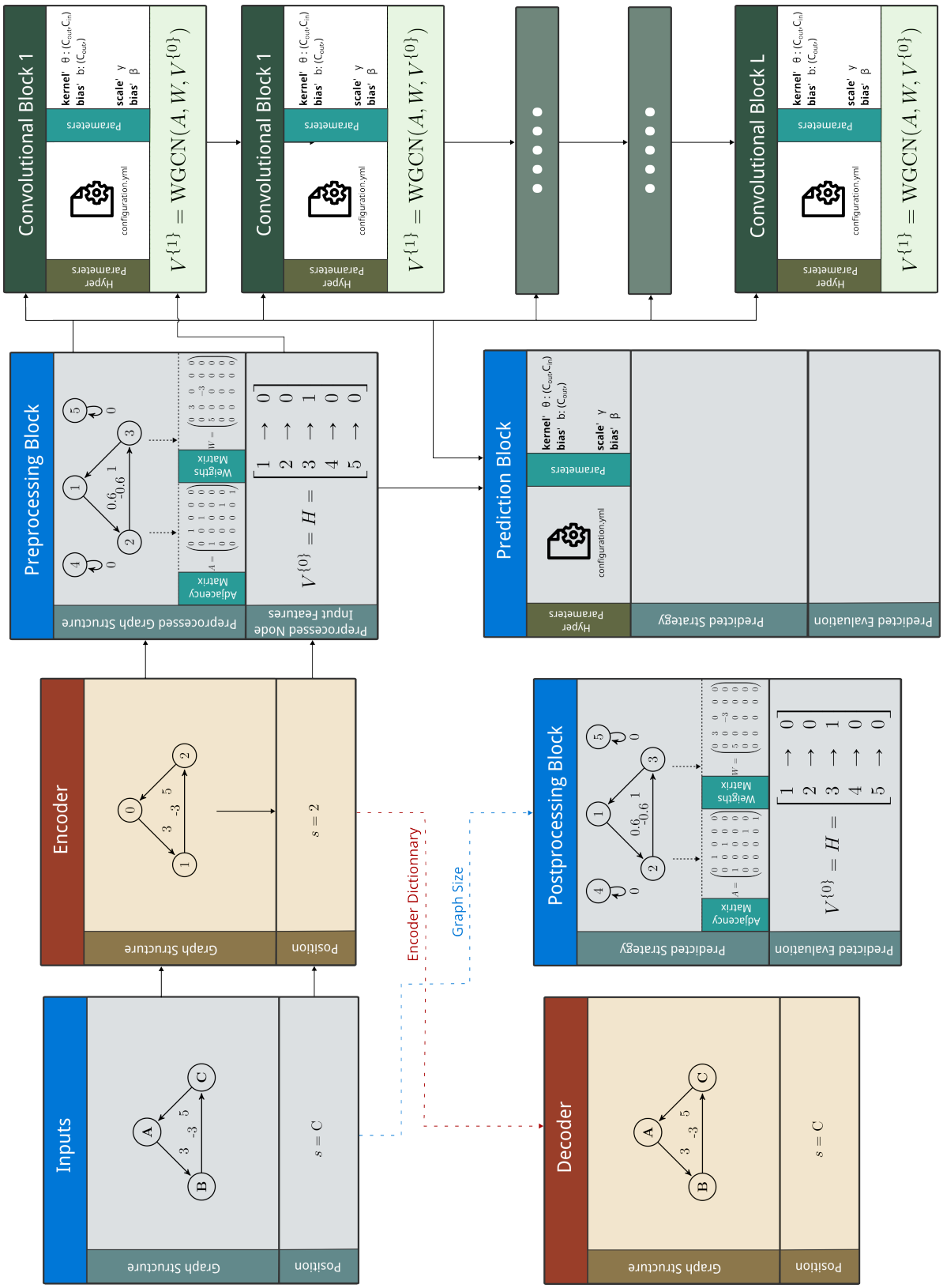

Figure 4.10: Model Architecture

Chapter 5

Reinforcement Learning Approach

5.1 Introduction

5.2 Monte Carlo Augmentation

5.2.1 Monte Carlo Tree Search

5.3 Pipeline

5.4 Services

5.5 Configuration

Chapter 6

Analyse

For $x \in \mathcal{X}, Y \subseteq \mathcal{X}^m, c \in I$, let $\mathrm{MA}(x, Y, c)$ be defined as follow:

$$\mathrm{MA}(x, Y, c) \iff x \leqslant \max Y + c$$

New York (NY) Hi

Conclusion

Durant ce stage, nous avons étudié les BNNs, implémenté une bibliothèque de BNNs basée sur tensorflow et larq qu'on a nommé binaryflow.

Dans le premier chapitre, nous avons présenté **dB Sense** et ses activités.

Dans le deuxième chapitre, nous avons présenté le problème de la grande complexité, introduit le concept des BNNs et proposé **binaryflow** comme notre solution pour les BNNs

Dans le troisième chapiter, nous avons formalisé les BNNs, et nous avons décris leurs optimisations possible, et les problèmes rencontrés dans leur implémentation.

Dans le quatrième chapitre, nous avons présenté des modèles BNNs, chacun en dérivant ses formules

Dans le cinquième chapitre, nous avons donné notre implémentation de **binaryflow** en jusitifiant les paradigmes utilisés

Dans le sixième chapitre, nous avons analysé 3 jeux de données en implémentant des modèles BNNs pour chacune de ces 3 jeux de données, et en comparant les performances de prédicition et la complexité temps et mémoire des modèles entraînés.

Notre travail n'est qu'une petite introduction des BNNs. En effet, la liste des BNNs proposés dans la littérature est très vaste, et il existe plusieurs autres approches pour faciliter l'entraînement et l'interférence des BNNs que nous n'avons pas considéré vu les contraintes de stage, y parmi:

1. Les méthodes d'optimisations discrètes
2. Les optimiseurs dédiés au BNNs
3. Les binarisations entraînables
4. Les méthodes ensemblistes pour régulaliser les BNNs

De plus, nous avons réussi à vérifier l'optimisation de la multiplication matricielle (L'éxecution est parfois 30 fois plus rapide), mais nous n'avons pas pu intégrer cette optimisation aux modèles tensorflow. Et malgré que larq supporte lui même un déploiement optimisé, nous n'avons pas aussi pu l'exploiter puisque ce déploiement ne supporte que les processeurs ARMv8, et nous utilisons une machine avec un processeur d'architecture x86-64.

Finalement, nous avons fait une petite intégration du code carbone comme une mésure de coût d'entraînement. Une fois le problème de déploiement est résoulu, nous recommenderions l'utilisation de ce même métrique pour estimer le coût d'interférence qui va justifier l'utilisation des BNNs.

Appendix A

On Constraint Satistfaction Problems

In the previous chapters, we described how the system works, without formalising the CSP approach.
On this chapter, we will describe the CSP systems that we have used, with an equivalence proof between them.

A.1 Constraint Satisfaction Problem

A.1.1 Definition

A constraint satisfaction problem

A.1.2 Assignment

An assignment of a $\texttt{CSP}(\mathcal{X}, D, \Gamma)$ is a function $X : \mathcal{X} \to D$ such that, by replacing each $x \in \mathcal{X}$, by $X(x)$, all the constraints will evaluate to **True**

A.1.3 Polymorphism

A function $F : \mathscr{F}(\mathcal{X}, D)^k \to \mathscr{F}(\mathcal{X}, D)$ is said to be a polymorphism if:

$\forall X_1, \ldots, X_k$ assignments of $\texttt{CSP}(\mathcal{X}, D, \Gamma)$, $\quad F(X_1, \ldots, X_k)$ is also an assignment of $\texttt{CSP}(\mathcal{X}, D, \Gamma)$

Now, in the next section, we will define an important class of CSPs that is used to solve Mean Payoff Games, with the polymorphisms required for the solution algorithm's correctness

A.2 Ternary Max Atom Systems

A.2.1 Definition

- Let $\mathcal{X}$ be a finite set of variables
- Let $D = I \cup \{-\infty\}$, with $I \subseteq \mathbb{R}$.
- For $x, y, z \in \mathcal{X}, c \in I$, let $\mathrm{MA}_3(x, y, z, c)$ be defined as follow:

$$\mathrm{MA}_3(x, y, z, c) \iff x \leqslant \max(y, z) + c$$

A ternary max atom system is $\texttt{CSP}(D, \Gamma)$ where:

$$\begin{aligned}&\Gamma = \{\mathrm{MA}_3(x, y, z, c), \quad (x, y, z, c) \in \mathscr{R}\}\\&\mathscr{R} \subseteq \mathcal{X}^3 \times I\\&\mathscr{R}\text{is finite}\end{aligned}$$

A.2.2 Example

An example of a ternary max atom system is the following $\texttt{CSP}(D, \Gamma)$ with $D = \mathbb{Z}$ and Γ represented as follow:

$$\begin{aligned}x &\leqslant \max(y, z) - 1\\y &\leqslant \max(z, x) - 1\\z &\leqslant \max(x, y) - 1\end{aligned}$$

A.3 Max Atom Systems

A.3.1 Definition

- Let $\mathcal{X}$ be a finite set of variables
- Let $D = I \cup \{-\infty\}$, with $I \subseteq \mathbb{R}$
- For $x \in \mathcal{X}, Y \subseteq \mathcal{X}^m, c \in I$, let $\mathrm{MA}(x, Y, c)$ be defined as follow:

$$\mathrm{MA}(x, Y, c) \iff x \leqslant \max Y + c$$

A max atom system is $\texttt{CSP}(D, \Gamma)$ where:

$$\begin{aligned}&\Gamma = \{\mathrm{MA}(x, Y, c), \quad (x, Y, c) \in \mathscr{R}\}\\&\mathscr{R} \subseteq \mathcal{X} \times (\mathscr{P}(\mathcal{X})\backslash\{\varnothing\}) \times I\\&\mathscr{R}\text{is finite}\end{aligned}$$

A.3.2 $\mathbf{MA} \leqslant \mathbf{MA}_3$

- Let $S = \texttt{CSP}(\mathcal{X}, D, \Gamma)$ a max atom system.
- Let $R \in \Gamma$
- Let $x \in \mathcal{X}, Y \in \mathscr{P}(\mathcal{X}), c \in I$ such that $R = \mathrm{MA}(x, Y, c)$ such that $|Y| > 2$

Recursive Reduction

We will reduce the arity of R as follow:

- Let $y, z \in Y$ such that $y \neq z$
- We introduce a variable $w \notin \mathcal{X}$
- Let $\mathcal{X}' = \mathcal{X} \cup \{w\}$

- Let $Y' = (Y \cup \{w\}) \backslash \{y, z\}$
- Let $R' = \mathrm{MA}(x, Y', c)$
- Let $R_w = \mathrm{MA}(w, \{y, z\}, 0)$
- Let $\Gamma' = (\Gamma \cup \{R', R_w\}) \backslash \{R\}$
- Let $S' = \mathtt{CSP}(\mathcal{X}', D, \Gamma)$

We will prove that S' is equivalent to S.

Implication Let $X : \mathcal{X}' \to D$ an assignment of S'. It is trivial that by removing $X(w)$, $X_{|\mathcal{X}}$ is an assignment of S

Equivalence

- Let $X : \mathcal{X}' \to D$ such that $X_{|\mathcal{X}}$ is an assignment of S.
- We will set $X(w) = \max(X(y), X(z))$

Then, X is an assignment of S'

Induction

Since the number of variables is finite, the arity of each constraint is finite. Also, as the the number of constraints is finite, Applying such reduction iteratively will eventually give a system S^* equivalent to S with:

- $\mathcal{X}^*$ the set of variables with $\mathcal{X} \subseteq \mathcal{X}^*$
- Γ^* is the set of constraints:
- Each constraint is of the form $\mathrm{MA}(x, Y, c)$ with $x \in \mathcal{X}^*, Y \subseteq \mathcal{X}^*, c \in I$ with $|Y| \leqslant 2$

Now such system can be transformed to a ternary system S_3 as follow:

- The set of variables is $\mathcal{X}^*$
- The domain is D
- For every relation $R = \mathrm{MA}(x, Y, c)$ we map it to the relation $R_3 = \mathrm{MA}(x, y, z, c)$ as follow:
  - If $|Y| = 2$, then y, z are the elements of Y.
  - Otherwise, $|Y| = 1$, and $y = z$ are the same element of the singleton[1] Y.

It is trivial that S^* is equivalent to S_3. With that, S is equivalent to S_3.

[1] A set with only one element

Algorithm 8 Converting a Max Atom System to Ternary Max Atom System

Require: S an N-ary Max Atom system
Ensure: S' a ternary Max Atom system
  $S' \leftarrow \varnothing$
  $H \leftarrow \varnothing$ ▷ H is a symmetric map between variable,variable to variables
  $V \leftarrow \text{Variables}(S)$ ▷ V is a set containing all variables
  for $\mathcal{C} \in S$ **do** ▷ Iterate over constraints
    c is the constant in the right hand side of $\mathcal{C}$
    Y is the variables in the right hand side of $\mathcal{C}$
    x is the variable in the left hand side of $\mathcal{C}$
    while $|Y| > 2$ **do**
      $y \leftarrow \text{pop}(Y)$
      $z \leftarrow \text{pop}(Y)$
      if $(y, z) \notin \text{domain}\, H$ **then**
        $w \leftarrow \text{newVariable}(V)$ ▷ Generate a new formal variable not included in V
        $V \leftarrow V \cup \{w\}$
        $H(y, z) \leftarrow w$
        $H(z, y) \leftarrow w$
      end if
      $w \leftarrow H(y, z)$
      $S' \leftarrow S' \cup \{\text{MA}(w, y, z, c)\}$
      $Y \leftarrow Y \cup \{w\}$
    end while
  end for
  return S'

A.3.3 Polymorphisms

Two main family of polymorphisms are defined for Max Atom systems:

- The max polymorphisms M^k defined by:

$$M^k(X_1, \ldots, X_k)(x) = \max_{k \in \{1,\ldots,k\}} X_k(x)$$

- The translation polymorphisms T_c defined by:

$$T_c(X)(x) = X(x) + c$$

A.4 Min-Max System

- Let $\mathcal{X}$ be a finite set of variables
- Let I be the domain of the variables.
- Let $D = I \cup \{-\infty\}$, with $I \subseteq \mathbb{R}$
- For $x \in \mathcal{X}, Y \subseteq \mathcal{X}^m, C \in I^m$, let $\mathrm{MA}(x, Y, C)$ be defined as follow:

$$\mathrm{MA}(x, Y, c) \iff x \leqslant \max(Y + C)$$

- For $x \in \mathcal{X}, Y \subseteq \mathcal{X}^m, C \in I^m$, let $\mathrm{MI}(x, Y, C)$ be defined as follow:

$$\mathrm{MI}(x, Y, C) \iff x \leqslant \min(Y + C)$$

A min-max system is $\texttt{CSP}(D, \Gamma)$ where:

$$\begin{aligned} \Gamma &= \{O(x, Y, C), \quad (O, x, Y, C) \in \mathscr{R}\} \\ \mathscr{R} &\subseteq \{\mathrm{MA}, \mathrm{MI}\} \times \mathcal{X} \times (\mathcal{X} \times I)^+ \\ \mathscr{R} &\text{ is finite} \end{aligned}$$

A.4.1 Transforming to Max Atom Systems

A Max Atom system is trivially a Min Max system. So we will only prove the latter implication.
Let $S' = \texttt{CSP}(D, \Gamma)$ be a Min Max system, and let:

- Γ_{MI} be the constraints that has MI
- Γ_{MA} be the constraints that has MA

Transforming MI constraints

For each $\mathrm{MI}(x, Y, c) \in \Gamma_{\mathrm{MI}}$. we replace it with the following constraints:

$$\Gamma_{\mathrm{MI}}^{x,Y,C} = \{\mathrm{MA}(x, \{y\}, c), \quad y, c \in \mathrm{zip}(Y, C)\}$$

Tranforming MA constraints

For each $(y, c) \in \mathcal{X} \times I$ present in a max constraint of the system:

- We add a formal variable $z^{y,c}$ if $c \neq 0$.
- Else, we will simply represent by $z^{y,c}$ the variable y.

By denoting $Z^{Y,C}$ the following set:

$$Z^{Y,C} = \{z^{y,c}, \quad (y, c) \in \text{zip}(Y, C)\}$$

We build the following constraints:

$$\Gamma_{\text{MA}}^{x,Y,C} = \left\{\text{MA}(x, Z^{Y,C}, 0)\right\} \cup \{\text{MA}(z^{y,c}, \{y\}, c), \quad (y, c) \in \text{zip}(Y, C)\}$$

Building the Max Atom System

Now, let:

$$\Gamma' = \bigcup_{\text{OP} \in \{\text{MI},\text{MA}\}} \bigcup_{\text{OP}(x,Y,C) \in \Gamma_{\text{OP}}} \Gamma_{\text{OP}}^{x,Y,C}$$

The system $\texttt{CSP}(D, \Gamma')$ is an equivalent max system.

Equivalence

Let:

- $\mathcal{X}' = \mathcal{X} \cup \mathcal{X}_{\text{Generated}}$ the augmented set of variables.
- $\mathcal{C} = \text{OP}(x, Y, C) \in \texttt{CSP}(D, \Gamma)$ be a constraint.
- $X : \mathcal{X}' \to D$ an assignment of $\texttt{CSP}(D, \Gamma')$.

If OP = MI, it is trivial that $\text{OP}(x, Y, C)$ is equivalent to $\Gamma_{\text{OP}}^{x,Y,C}$.
Otherwise, for each $(y, c) \in \text{zip}(Y, C)$ we have:

$$X(z^{y,c}) \leqslant \max\{X(y)\} + c = X(y) + c$$

Now, we also have:

$$X(x) \leqslant \max_{(y,c) \in \text{zip}(Y,C)} X(z^{y,c}) + 0 \leqslant \max_{(y,c) \in \text{zip}(Y,C)} (X(y) + c)$$

With that, $X_{|\mathcal{X}}$ is an assignment of $\texttt{CSP}(D, \Gamma)$

Algorithm 9 Converting a Min-Max System to Max Atom

Require: S a Min-Max system
Ensure: S' an N-ary Max Atom system
  $S' \leftarrow \varnothing$
  $H \leftarrow \varnothing$ ▷ H is a map between variable,offsets to variables
  $V \leftarrow \text{Variables}(S)$ ▷ V is a set containing all variables
  for $\mathcal{C} \in S$ **do** ▷ Iterate over constraints
    C is the constants in the right hand side of $\mathcal{C}$
    Y is the variables in the right hand side of $\mathcal{C}$
    x is the variable in the left hand side of $\mathcal{C}$
    if $\mathcal{C}$ is a min constraint **then**
      $S' \leftarrow S' \cup \{\text{MA}(x, \{y, y\}, c), \quad (y, c) \in \text{zip}(Y, C)\}$
    else
      $Z \leftarrow \varnothing$
      for $(y, c) \in \text{zip}(Y, C)$ **do**
        if $(y, c) \notin \text{domain}\, H$ **then**
          $z \leftarrow \text{newVariable}(V)$ ▷ Generate a new formal variable not included in V
          $V \leftarrow V \cup \{z\}$
          $H(y, c) \leftarrow z$
        end if
        $S' \leftarrow S' \cup \{\text{MA}(H(y, c), \{y, y\}, c)\}$
        $Z \leftarrow Z \cup \{H(y, c)\}$
      end for
      $S' \leftarrow S' \cup \{\text{MA}(x, Z, 0)\}$
    end if
  end for

Algorithm 10 Converting a Mean Payoff Game to a Min Max system

Require: G a Mean Payoff Game
Ensure: S an Min-Max system
  $E \leftarrow E(G)$ ▷ The edges of G
  $V \leftarrow V(G)$ ▷ The variables of G
  $W \leftarrow W(G)$ ▷ The weight function of G
  $P \leftarrow P(G)$ ▷ The set of player of G
  for $(u, p) \in V \times P$ **do**
    $x \leftarrow (u, p)$
    $A \leftarrow \text{Adj}(x)$
    $Y = \{(a, \bar{p}), \quad a \in A\}$
    $C \leftarrow W(A)$ ▷ Calculating the weights element wise.
    if p is Max **then**:
      OP $\leftarrow$ MA
    else
      OP $\leftarrow$ MI
    end if
    $S \leftarrow S \cup \{\text{OP}(x, Y, C)\}$
  end for

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Algorithm 11 AC3 for Ternary Max Atom systems

Require: $\mathcal{C}$ a ternary Max Atom constraint
Require: $\nu : V \to \mathscr{P}(D)$ the admissible values function
Require: Q a Queue of pending updates
Ensure: ν update the admissible values function.

```
x ← the left-hand side of C
(y, z) ← the right-hand side variables of C
c ← the right-hand side constant C
Z ← [x, y, z]
for o ∈ {0, 1, 2} do                                   ▷ Iterate over all rotations
    admissible ← False
    x ← Z[o]
    y ← Z[(o + 1) mod 3]
    z ← Z[(o + 2) mod 3]
    μ ← ∅                                              ▷ Set of values to delete
    for a ∈ ν(x) do
        for (b, c) ∈ ν(y) × ν(z) do
            T ← [a, b, c]
            p ← T[(−o) mod 3]
            q ← T[(1 − o) mod 3]
            q ← T[(2 − o) mod 3]
            if p ⩽ max(q, r) + c then
                admissible ← True
            end if
        end for
        if not admissible then
            μ ← μ ∪ {a}
        end if
    end for
    if |μ| > 0 then
        ν(x) ← ν(x)\μ
        append(Q, Adj x)
    end if
end for
```


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Algorithm 12 AC3 Optimized for Ternary Max Atom systems

Require: $\mathcal{C}$ a ternary Max Atom constraint
Require: $\nu : V \to D$ the maximum admissible value for each variable.
Require: Q a Queue of pending updates
Ensure: ν update the maximum admissible values function.

```
x ← the left-hand side of C
(y, z) ← the right-hand side variables of C
c ← the right-hand
r ← max(ν(y), ν(z)) + c
if r < L then
    r ← −∞
end if
if r < ν(x) then
    ν(x) ← r
    append(Q, x)
end if
```

Algorithm 13 Solving a Mean Payoff Graph for all states

Require: G a Mean Payoff Graph
Ensure: $\Phi : V \times P \to V$ The optimal strategy of each player

```
Φ ← ∅
for p ∈ P do
    if p is Max then
        G' ← G
    else
        G' ← Ḡ
    end if
    S ← transform_{MPG→Min-Max}(G')
    S' ← transform_{Min-Max→Max}(S)
    S'' ← transform_{Max→Max3}(S')
    X ← arcconsistency(S'')
    for C ∈ S'' do                                   ▷ Iterate over constraints of S''
        OP ← Operator(C)                             ▷ Get the operator of C. Either Max or Min
        Y the right-hand side variables of C
        C the right-hand side constants of C
        x the left-hand side variable of C
        u ← projection_{V×P→V}(x)                    ▷ Extract the vertex
        if OP is Max then
            y*, c* ← argmax_{(y,c)∈zip(Y,C)} {X(y) + c}   ▷ Extracts the maximum assignment
            Φ(u, p) ← projection_{V×P→V}(y*)          ▷ The Strategy is the vertex of the maximum assignment
        end if
    end for
end for
return Φ
```


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Appendix B

On Random Graphs

In the previous chapters, we gave a rough analysis of graph generation.
In this chapter, we will dive into a more detailed analysis.

B.1 Introduction

B.2 Sinkless $\mathcal{D}(n,p)$ Graph

B.2.1 Property

Let $\mathtt{P}$ be the property[1] "Graph has not sink".
This property is increasing in the sense that:

$$\forall H \text{ spanning subgraph of } G, \quad H \in \mathtt{P} \implies G \in \mathtt{P}$$

As a consequence:

$$\forall n \in \mathbb{N}, p, p' \in [0,1]/ \quad p \leqslant p', \quad \mathscr{P}(\mathcal{D}(n,p) \in \mathtt{P}) \leqslant \mathscr{P}(\mathcal{D}(n,p') \in \mathtt{P})$$

We will be interested in two properties:

- The property "Vertex v has no sinks". We denote it by $\text{NoSink}(v)$.
- The property "Graph G has no sinks at all". We denote it by $\text{Sinkless}(G)$.

B.2.2 Basic Comparison with Normal $\mathcal{D}(n,p)$

We will calculate the expected value of $\deg v$. By applying the law of total expectancy:

$$\begin{aligned}\mathbb{E}[\deg v] &= \mathbb{E}[\deg v \mid \deg v > 0] \times \mathscr{P}(\deg v > 0) + \mathbb{E}[\deg v \mid \deg v = 0] \times \mathscr{P}(\deg v = 0) \\ &= \mathbb{E}[\deg v \mid \text{Sinkless}(G)] \times \mathscr{P}(\text{NoSink}(v))\end{aligned}$$

[1] Formally, a property is a just a set of graphs. In practice, it is a set that has desirable "properties".

With that:

$$\mathbb{E}[\deg v \mid \text{Sinkless}(G)] = \frac{\mathbb{E}[\deg v]}{\mathscr{P}(\text{NoSink}(v))} = \frac{np}{1-(1-p)^n} \leqslant \frac{np}{1-e^{-1}}$$
$$\mathbb{E}[|\mathcal{E}|] = \sum_{v \in V} \mathbb{E}[\deg v \mid \text{Sinkless}(G)] = \frac{n^2 p}{1-(1-p)^n} \leqslant \frac{n^2 p}{1-e^{-1}}$$

This shows that the conditional distribution does inflict a small multiplicative bias on the expected number of edges and expected degree.
This serves as an evidence that $\mathcal{D}^S(n,p)$ is similar enough to $\mathcal{D}(n,p)$

B.2.3 Property Probability

- Let $G \sim \mathcal{D}(n,p)$
- Let v a vertex of G

The probability that $\text{NoSink}(v)$ occurs is:

$$\begin{aligned}\mathcal{P}(\text{NoSink}(v)) &= 1 - \mathcal{P}(\text{Adj}\, v = \varnothing) \\ &= 1 - \mathcal{P}(\deg v = 0) \\ &= 1 - (1-p)^n\end{aligned}$$

Now, it is clear that the sequence of events $(\text{NoSink}(v))_{v \in V}$ is independent.
With that, the probability that the whole graph is sinkless is:

$$\begin{aligned}\mathcal{P}(\text{Sinkless}(G)) &= \mathcal{P}(\text{Adj}\, v \neq \varnothing \quad \forall v \in V) \\ &= \mathcal{P}\left(\bigwedge_{v \in V} \text{NoSink}(v)\right) \\ &= \prod_{v \in V} \mathcal{P}(\text{NoSink}(v)) \\ &= (1-(1-p)^n)^n\end{aligned}$$

B.2.4 Asymptotic Analysis For Dense $\mathcal{D}(n,p)$

Let $c > 0$. We have for large enough n:

$$(1-p)^n \leqslant \frac{c}{n}$$

Which implies:

$$(1-\tfrac{c}{n})^n \leqslant (1-(1-p)^n)^n \leqslant 1$$

If we take the limit, we have:

$$e^{-c} \leqslant \lim_{n \to +\infty} (1-(1-p)^n)^n \leqslant 1 \quad \forall c > 0$$

By tending c to 0, we get:

$$\lim_{n \to +\infty} (1-(1-p)^n)^n = 1$$

B.2.5 Asymptotic Analysis For Sparse $\mathcal{D}(n,p)$

Let:

$$\begin{aligned} f : \mathbb{R}_+^* \times \mathbb{R}_+ \times \mathbb{R} &\to \mathbb{R}_+ \\ x, k, c &\to (1 - g(x,k,c))^x \\ g : \mathbb{R}_+^* \times \mathbb{R}_+ \times \mathbb{R} &\to \mathbb{R}_+ \\ x, k, c &\to \left(1 - \frac{k \ln x + c}{x}\right)^x \end{aligned}$$

By construction, $f(n,k,c)$ is the probability of a graph following $\mathcal{G}(n, \frac{k \ln n + c}{n})$ to contain no sink.

We have:

$$\begin{aligned} \ln g(k,x,c) &= x \ln \left(1 - \frac{k \ln x + c}{x}\right) \\ &= -k \ln x - c - \frac{(k(\ln x) + c)^2}{2x} + o\left(\frac{(\ln x)^3}{x^2}\right) \\ \implies g(x,k,c) &= \exp\left(-k \ln x - c - \frac{(k \ln x + c)^2}{2x} + o\left(\frac{(\ln x)^3}{x^2}\right)\right) \\ &= \frac{e^{-c}}{x^k} \times e^{\frac{-(k \ln x + c)^2}{2x} + o\left(\frac{(\ln x)^3}{x^2}\right)} \\ &= \frac{e^{-c}}{x^k}\left(1 - \frac{(k \ln x + c)^2}{2x} + o\left(\frac{(\ln x)^3}{x^2}\right)\right) \\ &= \frac{e^{-c}}{x^k} - e^{-c}\frac{k^2(\ln x)^2}{2x^{k+1}} + o\left(\frac{(\ln x)^3}{x^{k+2}}\right) \\ &= \frac{e^{-c}}{x^k} + o\left(\frac{1}{x^k}\right) \\ \implies 1 - g(x,k,c) &= 1 - \frac{e^{-c}}{x^k} + o\left(\frac{1}{x^k}\right) \\ \implies x \ln(1 - g(x,k,c)) &= -\frac{e^{-c}}{x^{k-1}} + o\left(\frac{1}{x^{k-1}}\right) \\ &\sim -\frac{e^{-c}}{x^{k-1}} \\ \implies f(x,k,x) &= e^{-\frac{e^{-c}}{x^{k-1}} + o(\frac{1}{x^{k-1}})} \end{aligned}$$

Now with that:

$$\lim_{x \to +\infty} x \ln(1 - g(x,k,c)) = \begin{cases} -\infty & \text{if} k \in [0,1[\\ -e^{-c} & \text{if} k = 1 \\ 0 & \text{otherwise if} k \in]1, +\infty[\end{cases}$$

Finally, we can conclude that:

$$\lim_{x\to+\infty} f(x,k) \begin{cases} 0 & \text{if} k \in [0,1[\\ e^{-e^{-c}} & \text{if} k = 1 \\ 1 & \text{otherwise if} k \in]1,+\infty[\end{cases}$$

B.3 Repeating Construction

B.3.1 Estimating Complexity

Now the method of rejecting graphs that have sinks and retrying give us a natural question about how many times will the algorithm reject graph until finding a desirable one.
The number of such rejections will follow a geometric law $\mathcal{G}(h(n,p))$ where:

$$h(n,p) = \mathscr{P}\left(\text{Sinkless}(\mathcal{D}(n,p))\right) = (1-(1-p)^n)^n$$

With that, the expected complexity of the algorithm will be:

$$\tilde{\mathcal{O}}\left(\frac{C(n,p)}{h(n,p)}\right) = \tilde{\mathcal{O}}\left(\frac{C(n,p)}{(1-(1-p)^n)^n}\right)$$

With $C(n,p)$ the cost of building the graph, depending on the algorithm[2].

B.3.2 Dense Graph case

Now it is clear for dense enough graphs, in particular with $p(n) \geqslant \frac{k\ln(n)}{n}$ for large enough n, the expected complexity will reduce to $\mathcal{O}(C(n,p))$. And thus, we consider the rejection method to be efficient.

B.3.3 Sparse Graph case

If $p(n) = \frac{k\ln n}{n} + c$ with $k < 1$. We have:

$$(1-(1-p)^n)^n = e^{-e^{-c}x^{1-k}+o(x^{1-k})}$$

With that, the expected complexity of the rejection method will be:

$$\tilde{\mathcal{O}}\left(C(n,p) \times \exp\left(e^{-c}x^{1-k} + o(x^{1-k})\right)\right)$$

which is an exponential algorithm, and thus inefficient for large graphs. Since property `P` is increasing, this argument generalises to $p(n) \leqslant \frac{k\ln n}{n} + c$ for large enough n

[2]The two algorithms that we have discussed are the naive $\mathcal{O}(n^2)$ algorithm and the more optimized $\mathcal{O}(pn^2+n)$ algorithm.

B.4 Binomial Rejection Construction

Instead of throwing the whole graph at once. For every vertex $u \in V$, we try to construct the adjacenty vertices of u, and repeat if the procedure gives $\operatorname{Adj} u = \varnothing$
With this trick, the expected complexity will reduce for both algorithms to:

$$\tilde{\mathcal{O}}\left(\frac{C(n,p)}{1-(1-p)^n}\right)$$

Now for our case, it is natural to assume that $p(n) \geqslant \frac{1}{n}$, as a Mean Payoff Graph does not have a sink. With that:

$$1-(1-p)^n \geqslant 1-(1-\tfrac{1}{n})^n \geqslant 1-e^{-1}$$

Therefore, the expected complexity will simplify to:

$$\tilde{\mathcal{O}}\left(C(n,p)\right)$$

Moreover, the cost of the conditionning makes only a constant $\frac{1}{1-e^{-1}} \approx 1.582$ factor slowdown, which is effectively neligible.

B.5 Expected Mean Payoff

B.5.1 Definition

- Let $\mathcal{G} = (\mathcal{V}, \mathcal{E})$ be a mean-payoff game
- For $u \in \mathcal{V}$, Let $\mathscr{P}(u)$ be the set of probability distributions over the set $\operatorname{Adj}(u)$
- We define a fractional strategy as a function $\Phi \in \mathscr{P}$

B.5.2 Matrix Form

- Let $n = |\mathcal{V}|$
- Let $u_1, \ldots, u_n$ an enumeration of elements of $\mathcal{V}$ A fractional strategy can be represented as a matrix A such that:

$$\mathcal{P}(\Phi(u_i) = u_j) = A_{i,j}$$

B.5.3 Fractional Strategies

Notations

- Let A, B be a pair of fractional strategies
- Let P_m, Q_m two random variables defining the mean-payoffs for the respective players after turn m

Expected Cost of $\frac{1}{2}$-turn

Let $\Pi \in \{A, B\}$
We have:

Rami ZOUARI
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$$\mathbb{E}\left[w(u,\Pi(u))\right] = \sum_{v\in \mathrm{Adj}\, u} w(u,v)\cdot \mathcal{P}(\Pi(u)=v)$$

Expected Cost of full turn

Let h be the cost of a turn

We have:

$$\mathbb{E}\left[h(u,A(u),B\circ A(u))\right] = \mathbb{E}[w(u,A(u))] + \sum_{v\in \mathrm{Adj}\, u} \mathbb{E}[w(v,B(v))]\cdot \mathcal{P}(A(u)=v)$$

Expected Total Payoff

- Let $\Pi = B\circ A$
- Let $(X_m)_{m\in\mathbb{N}}$ defined as follow:

$$\begin{cases} X_0 & = s \\ \forall m\in\mathbb{N}^*, \quad X_m & = \Pi(X_{m-1}) \end{cases}$$

- Let $(R_m)_{m\in\mathbb{N}}$ defined as follow:

$$\begin{cases} R_0 & = 0 \\ \forall m\in\mathbb{N}^*, \quad R_m & = R_{m-1} + \sum_{u\in V} h(u,A(u),\Pi(u))\cdot \mathcal{P}(X_{m-1}=u) \end{cases}$$

We have:

$$\begin{aligned} \mathbb{E}[R_m] &= \mathbb{E}[R_{m-1}] + \sum_{u\in V} \mathbb{E}[h(u,A(u),\Pi(u))]\cdot \mathcal{P}(X_{m-1}=u) \\ &= \mathbb{E}[R_{m-1}] + \sum_{u\in V} P^{m-1}(s,u)\times q(u) \\ &= \mathbb{E}[R_{m-1}] + (P^{m-1}\cdot q)(s) \quad \text{(Matrix Multiplication)} \\ &= \sum_{k=1}^{m} (P^{k-1}\cdot q)(s) + \mathbb{E}[R_0] \\ &= \left(\sum_{k=0}^{m-1} P^k\cdot q\right)(s) + \mathbb{E}[R_0] \\ &= \left(\sum_{k=0}^{m-1} P^k\cdot q\right)(s) \end{aligned}$$

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Now, we may see that the formula is easy generalisable to any starting vertex:

$$\mathbb{E}[R_m] = \sum_{k=0}^{m-1} P^k \cdot q$$

Expected Mean Payoffs

The mean-payoff is defined as:

$$K_m = \frac{R_m}{m}$$

We define K_∞ as:

$$K_\infty = K_{+\infty} = \lim_{m \to +\infty} \frac{R_m}{m}$$

Let $m \in \mathbb{N} \cup \{+\infty\}$ Now, the expected mean-payoff can act as a the judge for who is winning:

1. Player 0 wins if $\mathbb{E}[K_m] > 0$
2. Player 1 wins if $\mathbb{E}[K_m] < 0$
3. Else, it is a tie

Now, if m is finite, we can calculate $\mathbb{E}[K_m]$ directly. Otherwise, we have:

$$\mathbb{E}[K_\infty] = \lim_{m \to +\infty} \frac{1}{m} \sum_{k=0}^{m} P^k \cdot q$$

Now, P can be seen as a stochastic matrix. Thus it has a simple eigenvalue of value 1, and all its other eigenvalue λ satisfies:

$$\lambda \neq 1 \wedge |\lambda| \leqslant 1$$

Also, we have (Proof?):

$$|\lambda| = 1 \implies \lambda \text{is a simple eigenvalue}$$

With that, it can be proven that the $\lim_{m \to +\infty} \frac{1}{m} \sum_{k=0}^{m} P^k$ converges so some matrix T
This matrix can be constructed as follow. Let $P = VJV^{-1}$ the jordan normal form of P
Without a loss of generality, we will suppose that the first eigenvalue of this decomposition is 1 We have then:

$$T = V \begin{pmatrix} 1 & 0 \\ 0 & \mathbf{0}_{n-1} \end{pmatrix} V^{-1}$$

Discounted Payoffs

Using the same approach as the mean payoffs. Let R_m be the discounted payoff. It can be shown that:

$$\mathbb{E}[R_m] = \sum_{n \in \mathbb{N}} \gamma^n P^n \cdot q = (\mathrm{Id} - \gamma P)^{-1} q$$

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