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# A Path-following Polynomial Equations Systems Approach for Computing Nash Equilibria 

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#### Abstract

This paper presents a path-following combinatorial framework based on systems of polynomial equations to compute a mixed Nash equilibrium in $N$-person games. We provide the first detailed implementable description of Wilson's path-following method, extending Lemke-Howson's algorithm to N-person games and handling degenerate games. Our approach resembles, in some respects, support enumeration methods. We thus compare both approaches, theoretically and experimentally. Then, we show that the pathfollowing approach allows to deal with a large family of succinctly expressed games: hypergraphical games, graphical games and polymatrix games. The described algorithms have been implemented in Python, making use of Sagemath libraries to solve systems of polynomial equations, allowing an experimental comparison of the different combinatorial approaches on a large variety of games.


## KEYWORDS

Game Theory, Nash Equilibrium, Path Following Algorithm

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## 1 INTRODUCTION

The geometric path-following approach of [23] extends Lemke and Howson's 2-player games algorithm [17] and theoretically allows for the computation of exact mixed Nash equilibria of $N$-person games. However, while implementations of Lemke-Howson's type algorithms do exist for bimatrix and polymatrix games [14], Wilson's approach has never been implemented. Wilson only showed that computing a mixed Nash equilibrium amounts to finding a solution to a Polynomial Complementary Problem (PCP) and provided a high-level informal mathematical description of his approach. On different grounds, [21] provided a simple Nash equilibrium computation algorithm based on an exhaustive exploration of strategies' supports, exploiting polynomial systems and dominated strategies computation. This method is still state of the art for practical exact NE computation in $N$-person games. In this article, we compare both approaches theoretically. Then we extend the path-following approach in order to handle (i) degenerate games and (ii) polymatrix

[^0][24], graphical [16] and hypergraphical games [20]. Finally, we provide implementations of support enumeration and path-following approaches, allowing comparisons on different classes of games.

In Section 2, we present the PCP formulation of the Nash equilibrium search problem introduced in [23] and show how [21]'s support enumeration method can also be cast in the PCP framework. In Section 3 we present an original "combinatorial" view of [23]'s path-following algorithm and in Section 4 we extend it to handle degenerate PCP (approximately, in some cases). In Section 5, we show that Nash equilibrium computation in graphical and hypergraphical games [16,20] can be formulated as a PCP of size polynomial in the concise expression of the games. Finally, Section 6 provides an experimental comparison of both approaches ${ }^{1}$.

## 2 N-PLAYER GAMES AND POLYNOMIAL COMPLEMENTARITY PROBLEMS (PCP)

### 2.1 Nash equilibria as solutions of a PCP

Let us consider a $N$-player game $\Gamma^{N}=(P, \pi, a) . P=\{1, \ldots, N\}$ is the set of players. $\pi=S_{1} \times \ldots \times S_{N}$ is the set of joint pure strategies of the game ( $S_{n}$ is the finite set of pure strategies available to player $n \in P$ ). $a_{\omega}^{n}$ is the strictly positive disutility ${ }^{2}$ received by player $n$ when the joint pure strategy is $\omega \in \pi \cdot a=\left(a_{\omega}^{n}\right)_{\omega \in \pi, n \in P}$ can thus be represented as a matrix with $|\pi|=\prod_{n=1 . . N}\left|S_{n}\right|$ lines and $N$ columns. A mixed strategy $\xi^{n}=\left(\xi_{i}^{n}\right)_{i \in S_{n}}$ is a probability distribution over the pure strategies of player $n$. A joint mixed strategy is a tuple $\xi=\left(\xi^{n}\right)_{n \in P}$ of mixed strategies. $\xi^{-n}$ denotes the mixed strategies of all players except $n$. Hence, $\xi=\left(\xi^{n}, \xi^{-n}\right)$.

A mixed Nash equilibrium of a game $\Gamma^{N}$ is defined as:
Definition 2.1 (Mixed Nash equilibrium). Joint mixed strategy $\xi=$ $\left(\xi_{i}^{n}\right)_{n \in P, i \in S_{n}}$ is a mixed Nash equilibrium of game $\Gamma^{N}=(P, \pi, a)$ if and only if: $\forall n \in P, \operatorname{Dis}_{n}[\xi] \leq \operatorname{Dis}_{n}\left[\left(\bar{\xi}^{n}, \xi^{-n}\right)\right], \forall \bar{\xi}^{n} \neq \xi^{n}$, where

$$
\operatorname{Dis}_{n}[\xi]=\operatorname{def} \sum_{\omega=\left(\omega_{1}, \ldots, \omega_{n}\right) \in \pi} a_{\omega}^{n} \prod_{v=1}^{N} \xi_{\omega_{v}}^{v}
$$

and $\left(\bar{\xi}^{n}, \xi^{-n}\right)$ is the joint mixed strategy where the mixed strategy $\xi^{n}$ has been replaced with mixed strategy $\bar{\xi}^{n}$.
[23] showed that a mixed Nash equilibrium can be computed from a solution of a Polynomial Complementarity Problem (PCP). Let us define $E^{\pi}$, the Euclidian space of dimension $D=\sum_{n \in P}\left|S_{n}\right|$. Joint mixed strategies will be indirectly represented by lists of

[^1]vectors $x=\left(x^{n}\right)_{n \in P}$ of coordinates $\left(x_{i}^{n}\right)_{i \in S_{n}} \in E^{\pi}$. The $x^{n}$, s are non-negative but generally not normalized (we will build a probability distribution by normalization of $x^{n}$ ). Let us define multilinear polynomials $A_{i}^{n}$ :
\[

$$
\begin{align*}
A_{i}^{n}\left(x^{-n}\right) & =\operatorname{def} \quad \sum_{\substack{\omega \in \pi_{\omega} \\
\omega_{n}=i}} a_{\omega}^{n} \prod_{v \neq n} x_{\omega_{v}}^{v}, \forall n \in P, i \in S_{n}  \tag{1}\\
\text { and } A^{n}(x) & =\operatorname{def} \quad \sum_{i \in S_{n}} A_{i}^{n}\left(x^{-n}\right) x_{i}^{n}, \forall n \in P \tag{2}
\end{align*}
$$
\]

Where, by definition, $x^{-n}=\left(x_{i}^{v}\right)_{v \in P \backslash\{n\}, i \in S_{v}}$.
When every $x^{n}$ is normalized, i.e. $x^{n}=\xi^{n}$ are mixed strategies, $A^{n}(\xi)$ is the expected disutility of joint mixed strategy $\xi$ to player $n$ and $A_{i}^{n}\left(\xi^{-n}\right)$ is the expected disutility of player $n$ when $n$ plays pure strategy $i \in S_{n}$ instead of mixed strategy $\xi^{n}$.

By definition of a mixed Nash equilibrium, $\xi$ is a mixed NE iff:

$$
\left\{\begin{align*}
A^{n}(\xi) & \leq A_{i}^{n}\left(\xi^{-n}\right), \forall n \in P, i \in S_{n},  \tag{3}\\
A^{n}(\xi) & =A_{i}^{n}\left(\xi^{-n}\right), \forall n \in P, i \in S_{n}, \text { s.t. } \xi_{i}^{n}>0
\end{align*}\right.
$$

By a change of variables, system (3) can be represented by the following Polynomial Complementarity Problem [23]:

Definition 2.2 (Polynomial Complementarity Problem). The PCP corresponding to $\Gamma^{N}=(P, \pi, a)$ is a system of equations/inequations in variables $\left(x_{i}^{n}\right)_{n \in P, i \in S_{n}} \in E^{\pi}$ :

$$
\forall(n, i) \in I_{N},\left\{\begin{aligned}
x_{i}^{n} & \geq 0 \\
A_{i}^{n}\left(x^{-n}\right) & \geq 1 \\
x_{i}^{n} \cdot\left(A_{i}^{n}\left(x^{-n}\right)-1\right) & =0
\end{aligned}\left(\mathcal{S}^{N}\right)\right.
$$

where $I_{N}={ }_{d e f}\left\{(n, i), n \in P, i \in S_{n}\right\}$.
$\mathcal{D}^{N}$ denotes the set of points $x \in E^{\pi}$ satisfying the set of inequations of the PCP: $\mathcal{D}^{N}=\left\{x \in E^{\pi}, x_{i}^{n} \geq 0, A_{i}^{n}\left(x^{-n}\right) \geq 1, \forall(n, i) \in I_{N}\right\}$. Problem $\mathcal{S}^{N}$ is called a polynomial complementarity problem since we look for a non-negative solution $x$ such that for any $(n, i) \in I_{N}$, either $x_{i}^{n}=0$ or $A_{i}^{n}\left(x^{-n}\right)=1$ holds (hence a complementary solution) and $A_{i}^{n}\left(x^{-n}\right)$ are multivariate polynomials in variables $\left(x_{i}^{n}\right)_{n \in P, i \in S_{i}} \cdot x^{n}$ representing an unnormalized probability distribution, the equalities in $\mathcal{S}^{N}$ correspond to usual best response arguments: if $i$ is not a best response to a NE for player $n$ (i.e. $A_{i}^{n}\left(x^{-n}\right)>1$ ), its probability should be 0 (thus, also $x_{i}^{n}$ ).

In the case where there are only two players ( $N=2$ ), one can check that, $\forall(n, i) \in I_{N}, A_{i}^{n}\left(x^{-n}\right)=1$ is a linear equation. Hence, in this case the obtained problem $\mathcal{S}^{2}$ is a Linear Complementarity Problem [17]. [23] shows the equivalence between the Nash equilibria of a game and the solutions of the corresponding PCP:

Proposition 2.3 (NE / PCP equivalence [23]). Let $\Gamma^{N}$ be a $N$-player game and $\mathcal{S}^{N}$ its PCP transformation. Then, the Nash equilibria of $\Gamma^{N}$ and the solutions of $\mathcal{S}^{N}$ are in one-to-one correspondence, i.e.:
(1) If $x$ is a solution of $\mathcal{S}^{N}$, then $\xi$ defined by $\xi_{i}^{n}=\frac{x_{i}^{n}}{\sum_{j \in S_{n}} x_{j}^{n}}$, $\forall(n, i) \in I_{N}$ is a Nash equilibrium of $\Gamma^{N}$.
(2) If $\xi$ is a mixed Nash equilibrium of $\Gamma^{N}, x$ defined as
$x_{i}^{n}=\left(\frac{\prod_{v \neq n} A^{v}(\xi)}{A^{n}(\xi)^{N-2}}\right)^{\frac{-1}{N-1}} \xi_{i}^{n}, \forall(n, i) \in I_{N}$ is a solution of $\mathcal{S}^{N}$.

A solution $x \in E^{\pi}$ of a PCP is called a complementary point. A point $x \in E^{\pi}$ is called a ( $n, i$ )-almost complementary point if it satisfies every constraints in $\mathcal{S}^{N}$, except possibly ${ }^{3}$ for the equation $x_{i}^{n} \cdot\left(A_{i}^{n}\left(x^{-n}\right)-1\right)=0$.

In the view of this result, finding a mixed Nash equilibrium of a game $\Gamma^{N}$ amounts to finding a solution to the $\operatorname{PCP} \mathcal{S}^{N}$.

### 2.2 Solving PCP through support enumeration

[21] suggested a simple support enumeration method to find a Nash equilibrium of a game, which can be described very simply in terms of PCP solution. This approach consists in enumerating every set of joint supports $W \subseteq I_{N}$ (such that $W$ contains at least one pair $(n, i), \forall n \in P$ ) and trying to solve a system of equations/inequations, which we write $\mathcal{S}^{\bar{W}, W}$ for a reason which will be made clear in the following section. $\mathcal{S}^{\bar{W}, W}$ is equal to the system $\mathcal{S}^{N}$ where the equalities are replaced with two distinct sets of equalities:

$$
\left\{\begin{aligned}
& x \in \mathcal{D}^{N}, \\
x_{i}^{n}=0, & \forall(n, i) \in \bar{W}, \\
A_{i}^{n}\left(x^{-n}\right)=1, & \forall(n, i) \in W
\end{aligned} \quad\left(\mathcal{S}^{\bar{W}, W}\right)\right.
$$

A solution $x$ of system $\mathcal{S}^{\bar{W}, W}$, if it exists, will correspond to an equilibrium $\xi$. However, there are exactly $\prod_{i=1}^{N}\left(2^{\left|S_{i}\right|}-1\right)$ possible supports $W \subseteq I_{N}$, so trying to solve a system of polynomial equations/inequations for every $W$ is not a good idea. [21] suggest to enumerate supports in increasing size of $W$, in an order compatible with inclusion (if $W \subset W^{\prime} W^{\prime}$ will be explored after $W$ ). As soon as a system $\mathcal{S}^{\bar{W}, W}$ with a solution is found, the algorithm exits and returns the corresponding Nash equilibrium. Iterated Removal of Dominated Alternatives [9] is applied before any new system $\mathcal{S}^{\bar{W}}, W$ is actually solved. If any alternative is removed, this means that we can reduce support $W$ to $W^{\prime \prime} \subset W$. However, given the enumeration order, the algorithm has already tried to solve $\mathcal{S}^{\overline{W^{\prime \prime}}, W^{\prime \prime}}$ and failed. So, $W$ is skipped before solving $\mathcal{S}^{\bar{W}}, W$. The algorithm is guaranteed to find a mixed Nash equilibrium of the game (or a continuum of solutions if a degenerate system $\mathcal{S}^{\bar{W}}, W$ is encountered).

Support enumeration is quite efficient in general, especially when an equilibrium with support of small size exists. However, for games with no NE of small support size, many joint supports may be explored and many systems of equations may have to be solved. Next, we describe an alternative approach to [21] to solve polynomial complementarity problems. This approach is inspired from [23]'s method, itself extending Lemke-Howson's algorithm [17]. The idea is, instead of enumerating blindly all possible pairs ( $\bar{W}, W$ ) until we find a solvable system, to solve an ordered sequence of systems $\mathcal{S}^{Z, W}$, where $Z$ may be different from $\bar{W}$ and the union $Z \cup W$ may be strictly included in $I_{N}$. This approach generally explores fewer pairs ( $Z, W$ ) than [21] but cannot exploit IRDA, so a larger proportion of systems may have to be solved.

[^2]
## 3 A COMBINATORIAL PATH-FOLLOWING ALGORITHM TO SOLVE PCP

Wilson [23], following the approach proposed by [17] to solve linear complementarity problems, proposed a mathematical path following approach to solve a non-degenerate PCP. Wilson's description leaves some steps of the algorithm undefined and does not deal with degenerate games. In this Section, we propose an original and operational rewriting of Wilson's approach.

The approach we propose is based on a definition of (almostcomplementary) nodes, arcs and paths, in terms of sets of multilinear equations - we detail these definitions in Section 3.1. Then, we show how paths can be extended through different levels of subPCP (Sections 3.2 and 3.3). Section 3.4 is devoted to the arc-traversal problem and Section 3.5 describes the full PCP solution algorithm.

### 3.1 Almost-complementary nodes, arcs and paths

Assume that $\operatorname{PCP} \mathcal{S}^{N}$ is given. For any $x \in \mathcal{D}^{N}$, let us write $Z(x)=$ $\left\{(n, i) \in I_{N}, x_{i}^{n}=0\right\}$ and $W(x)=\left\{(n, i) \in I_{N}, A_{i}^{n}\left(x^{-n}\right)=1\right\}$. Then, by definition, $x \in \mathcal{D}^{N}$ is a solution of $\mathcal{S}^{N}$ if and only if $Z(x) \cup W(x)=I_{N}$. A non-degenerate PCP at level $N$ is defined as:

Definition 3.1 (Non-degenerate PCP). PCP $\mathcal{S}^{N}$ is non-degenerate iff the following conditions hold:
(1) No point in $\mathcal{D}^{N}$ satisfies more than $\left|I_{N}\right|$ equations.
(2) No two distinct points satisfy the same set of $\left|I_{N}\right|$ equations.

In mathematical terms, condition 1 is equivalent to

$$
\forall x \in \mathcal{D}^{N},|Z(x)|+|W(x)| \leq D=\left|I_{N}\right|
$$

and condition 2 is equivalent to: $\forall x, y \in \mathcal{D}^{N}$,

$$
\left.\begin{array}{rll}
Z(x)=Z(y) & = & Z  \tag{4}\\
W(x)=W(y) & = & W \\
|Z|+|W| & = & \left|I_{N}\right|
\end{array}\right\} \Rightarrow x=y
$$

Intuitively, a PCP is non-degenerate if none of its polynomial constraints are redundant (it involves $\left|I_{N}\right|$ independent equations in a space of dimension $\left.\left|I_{N}\right|\right)$. For the moment, we assume nondegeneracy. Considering a non-degenerate PCP allows to distinguish particular points of interest, almost complementary nodes:

Definition 3.2 (Almost-complementary nodes). $x \in \mathcal{D}^{N}$ is an almost-complementary node of $\mathcal{S}^{N}$, if and only if:

$$
|Z(x)|+|W(x)|=\left|I_{N}\right| \text { and }|Z(x) \cap W(x)| \leq 1
$$

In particular, an almost-complementary node is said complementary if $Z(x) \cup W(x)=I_{N}$ and $(n, i)$-almost complementary if $(n, i) \notin Z(x) \cup W(x)$. Note that if $\mathcal{S}^{N}$ is non-degenerate, the almostcomplementary nodes of $\mathcal{D}^{N}$ are in one-to-one correspondence with the pairs $Z, W \subseteq I_{N}$ such that $|Z|+|W|=\left|I_{N}\right|,|Z \cap W| \leq 1$ and such that the system:

$$
\left\{\begin{array}{rc} 
& x \in \mathcal{D}^{N}, \\
x_{i}^{n}=0, & \forall(n, i) \in Z, \\
A_{i}^{n}\left(x^{-n}\right)=1, & \forall(n, i) \in W
\end{array} \quad\left(\mathcal{S}^{Z, W}\right)\right.
$$

has a solution. In the following we identify ( $n, i$ )-almost complementary node $x$ (denoted $\rho(Z, W)$ ) with the pair $(Z, W)$ which defines the system $\mathcal{S}^{Z, W}$ that $x$ solves.

We also define ( $n, i$ )-almost complementary arcs.
Definition 3.3 (Almost complementary arcs). The ( $n, i$ )-almost complementary arcs of non-degenerate PCP $\mathcal{S}^{N}$ are the subsets $\gamma(Z, W) \subseteq \mathcal{D}^{N}$, for all pairs $(Z, W) \subseteq I_{N}$ such that $Z \cap W=\emptyset$ and $Z \cup W=I_{N} \backslash\{(n, i)\}$, where $\gamma(Z, W)$ is formed by the points $x \in \mathcal{D}^{N}$ such that $Z(x)=Z$ and $W(x)=W$.

If almost complementary arc $\gamma(Z, W)$ is non-empty and the game is non degenerate, $\gamma(Z, W)$ is included in the set of solutions of a system of $D-1$ equations over $D$ variables. Non-degeneracy implies that this set has dimension 1 and can be parameterized by a single real-valued parameter. The extreme points of $\gamma(Z, W)$, belonging to the frontier of domain $\mathcal{D}^{N}$ are almost complementary nodes. An arc will typically have two extreme points (if bounded) or a single one if unbounded.
Wilson's approach consists in following a one-dimensional path in $\mathcal{D}^{N}$, by traversing arcs and nodes, until we eventually reach a complementary node. It relies on the following proposition ${ }^{4}$ :

Proposition 3.4 (Arcs neighbouring nodes). Let $\mathcal{S}^{N}$ be a non-degenerate PCP and $i \in S_{N}$. Two ( $N, i$ )-almost complementary arcs neighbour a $(N, i)$-almost-complementary node of $\mathcal{S}^{N}$ which is not complementary and a single ( $N, i$ )-almost complementary arc neighbours a complementary node.
The two arcs neighbouring almost complementary node $\rho(Z, W)$ are $\gamma(Z \backslash W, W)$ and $\gamma(Z, W \backslash Z)$, as shown in the proof of Proposition 3.4. Remark that Proposition 3.4 implies the following proposition:

Proposition 3.5 (Finite path). Let $\left(\mathcal{S}^{N}\right)$ be a non-degenerate $P C P$ and $i \in S_{N}$. There is a unique non-directed path, made of a finite sequence of ( $N, i$ )-almost complementary nodes and arcs, which reaches any complementary node of $\mathcal{S}^{N}$.

Proposition 3.5 does not tell us about the other end of the path. The path can end with another complementary node or with a ( $N, i$ )-almost complementary unbounded arc which neighbours a single ( $N, i$ )-almost complementary node. Let us see how this unique non-directed path can be extended through different "layers" of sub-PCP of the initial PCP, until it reaches an "easy-to-compute" initial node.

### 3.2 Defining a sequence of PCP

Let us consider a PCP $\mathcal{S}^{N}$, an arbitrary pure joint strategy $\omega^{0}=$ $\left(\omega_{1}^{0}, \ldots, \omega_{N}^{0}\right) \in \pi$, two integers $1 \leq n \leq k \leq N$ and a pure strategy $i \in S_{n}$. We write $A_{i}^{n, k}\left(x^{\{1, . ., k\} \backslash\{n\}}\right)$ the multivariate polynomial obtained from $A_{i}^{n}\left(x^{-n}\right)$ by fixing all values $x_{j}^{v}$ to 0 whenever $v>k$ and $j \neq \omega_{v}^{0}$ and to 1 whenever $v>k$ and $j=\omega_{v}^{0}$. Then,

$$
\begin{equation*}
A_{i}^{n, k}\left(x^{\{1, . ., k\} \backslash\{n\}}\right)=\sum_{\substack{\omega \in \pi, \omega_{n}=i \\ \omega_{m}=\omega_{m}^{0}, \forall m>k}} a_{\omega}^{n} \prod_{\substack{v \leq k, v \neq n}} x_{\omega_{v}}^{v} . \tag{5}
\end{equation*}
$$

$A_{i}^{n, k}$ is a multilinear polynomial of degree $k-1$. Note that if $\xi$ is a joint mixed strategy, $A_{i}^{n, k}(\xi\{1, \ldots, k\} \backslash\{n\})$ is the expected disutility

[^3]of player $n$ playing action $i \in S_{n}$ when players $1, \ldots, k$ except $n$ play their mixed strategy in $\xi$, while players $k+1, \ldots, N$ play their pure strategy in $\omega^{0}$. For any $1 \leq k \leq N$, let us define $I_{k}=$ $\left\{(n, i), 1 \leq n \leq k, i \in S_{n}\right\}$. We also let $E_{k}^{\pi}$ denote the Euclidian space spanned by the variables $x_{i}^{n}$ with $(n, i) \in I_{k}$. The dimension of $E_{k}^{\pi}$ is $D_{k}=\left|I_{k}\right|$ (thus $D=D_{N}$ ). Then, from PCP $\mathcal{S}^{N}$, we can define the following sequence of sub-PCP $\mathcal{S}^{k}$, for $k=1, \ldots, N-1$ :

Definition 3.6 (Sub-PCP). Let $\mathcal{S}^{N}$ be a PCP and $\omega^{0}$ a fixed pure joint strategy. Then, for $2 \leq k<N$, we define sub-PCP $\mathcal{S}^{k}$ as the following system of polynomial equations/inequations over $E_{k}^{\pi}$ :

$$
\forall(n, i) \in I_{k},\left\{\begin{aligned}
x_{i}^{n} & \geq 0 \\
A_{i}^{n, k}\left(x^{\{1, \ldots, k\} \backslash\{n\}}\right) & \geq 1 \\
x_{i}^{n} \cdot\left(A_{i}^{n, k}\left(x^{\{1, . ., k\} \backslash\{n\}}\right)-1\right) & =0
\end{aligned}\left(\mathcal{S}^{k}\right)\right.
$$

For $n=1, \operatorname{PCP} \mathcal{S}^{1}$ is slightly different, although deduced from best responses principles as well:

$$
\left\{\begin{aligned}
x_{i}^{1} & \geq 0, \forall i \in S_{1} \\
x_{i}^{1} \cdot\left(\frac{a_{\left(i, \omega_{-1}^{0}\right)}^{1}}{\min _{j \in S_{1}} a_{\left(j, \omega_{-1}^{0}\right)}^{1}}-1\right) & =0, \forall i \in S_{1} \quad\left(\mathcal{S}^{1}\right) \\
\sum_{i \in S_{1}} x_{i}^{1} & =1
\end{aligned}\right.
$$

Sub-PCP $\mathcal{S}^{k}$ is constructed from $\Gamma^{k}\left(\omega^{0}\right)$, the game played by the first $k$ players of $\Gamma^{N}$ when the remaining players play following $\omega^{0}$, in the same way as $\mathcal{S}^{N}$ was constructed from $\Gamma^{N}$. As for the initial PCP, we assume that all sub-PCP $\mathcal{S}^{k}$ are non-degenerate. In particular, assuming that $\mathcal{S}^{1}$ is non-degenerate implies that the minimum $\min _{j \in S_{1}} a_{\left(j, \omega_{-1}^{0}\right)}^{1}$ is attained for a single index $j^{*}$. Thus, we easily get the complementary point at level 1 , characterized by $Z^{1}=S_{1} \backslash\left\{j^{*}\right\}$ and $W^{1}=\left\{j^{*}\right\}$. Then, we can define the polynomial systems $\mathcal{S}_{k}^{Z, W}$ for any pair $Z, W \subseteq I_{k}$ :

$$
\left\{\begin{aligned}
& x \in \mathcal{D}^{k}, \\
x_{i}^{n}=0, & \forall(n, i) \in Z, \\
A_{i}^{n, k}\left(x^{\{1 . . k\} \backslash\{n\}}\right)=1, & \forall(n, i) \in W,
\end{aligned} \quad\left(\mathcal{S}_{k}^{Z, W}\right)\right.
$$

where $\mathcal{D}^{k}$ is defined by the inequations of Definition 3.6.

### 3.3 Complementary nodes and initial nodes

We can now exploit the combinatorial point of view of the above sequence of systems of equations in order to design an algorithm computing a sequence of almost-complementary nodes until a complementary node at level $N$ is reached.

Proposition 3.7 (Complementary node lifting). Let $\mathcal{S}^{k}$ be a sub-PCP of $\mathcal{S}^{N}$ at level $1 \leq k<N$ and arbitrary pure joint strategy $\omega^{0}$. Assume that, for $Z, W \subseteq I_{k}, \mathcal{S}_{k}^{Z, W}$ defines a complementary node of $\mathcal{S}^{k}$ in $E_{k}^{\pi}$. Then, $\mathcal{S}_{k+1}^{Z^{\prime}, W}$ defines a $\left(k+1, \omega_{k+1}^{0}\right)$-almost complementary arc of sub-PCP $\mathcal{S}^{k+1}$, where $Z^{\prime}=Z \cup\left\{(k+1, j), j \neq \omega_{k+1}^{0}\right\}$. Furthermore, this arc neighbours a single $\left(k+1, \omega_{k+1}^{0}\right)$-almost complementary node at level $k+1$ (it is an unbounded arc).

The ( $k+1, \omega_{k+1}^{0}$ )-almost complementary node at level $k+1$ can be computed by trying to solve all systems $\mathcal{S}_{k+1}^{Z^{\prime} \cup\{(v, j)\}, W}$ with $(v, j) \in I_{k} \backslash Z^{\prime}$ and $\mathcal{S}_{k+1}^{Z^{\prime}, W \cup\{(v, j)\}}$ with $(v, j) \in I_{k+1} \backslash W$ until
we find a solution. Such a system with a solution exists (this is a consequence of Lemma 2 in [23]). Furthermore, it is unique, due to non-degeneracy. This solution is called initial node at level $k+1$.

Definition 3.8 (Initial node). An initial node at level $k+1$ is a $\left(k+1, \omega_{k+1}^{0}\right)$-almost complementary node, solution to $\mathcal{S}_{k+1}^{Z, W}$, with $Z, W \subseteq I_{k+1}$, such that only one of its neighbouring arcs is bounded. It also satisfies: $\left(k+1, \omega_{k+1}^{0}\right) \notin Z$ and $(k+1, j) \in Z, \forall j \neq \omega_{k+1}^{0}$.

With this characterization in mind, Proposition 3.5 can be reinterpreted. It states that from any complementary node at level $N$ there is a unique path, leading either to another complementary node, or to an initial node at level $N$. This is obviously true at any level $k \in\{2, \ldots, N\}$, since sub-PCP are derived from subgames. Complementary nodes are at one end of a path of $\left(k, \omega_{k}^{0}\right)$-almost complementary points which other end is either another complementary node or an initial node. The set of almost-complementary points in $E_{k}^{\pi}$ which satisfy $\mathcal{S}^{k}$ form disjoint paths (see Figure 1).


Figure 1: Example almost-complementary paths at level $k$.
Then, we can define a descent procedure from an initial node at level $k$ to a complementary node at level $k-1$, reciprocal to the previous lifting procedure.

Proposition 3.9 (Initial node descent). Let $\mathcal{S}_{k}^{Z, W}$ define an initial node at level $k>1$. Let $Z^{\prime}=Z \cap I_{k-1}$ and $W^{\prime}=W \cap I_{k-1}$. Then, either $Z^{\prime} \cap W^{\prime}=\{(v, j)\}$ or $Z^{\prime} \cap W^{\prime}=\emptyset$.

In the first case, either $\mathcal{S}_{k-1}^{Z^{\prime} \backslash\{(v, j)\}, W^{\prime}}$ or $\mathcal{S}_{k-1}^{Z^{\prime}, W^{\prime} \backslash\{(v, j)\}}$ defines a complementary node at level $k-1$. In the second case, $\mathcal{S}_{k-1}^{Z^{\prime}, W^{\prime}}$ defines a complementary node.

### 3.4 Algebraic arc traversal

Our path-following algorithm heavily relies on the problem of traversing a $\left(k, \omega_{k}^{0}\right)$-almost complementary $\operatorname{arc}^{5} \gamma^{k}(Z, W)$, at level $k$ $\left(Z, W \subseteq I_{k}\right)$, that leaves an almost-complementary node $\rho^{k}\left(Z^{\prime}, W^{\prime}\right)$ with either (i) $Z=Z^{\prime} \backslash W^{\prime}$ and $W=W^{\prime}$ or (ii) $Z=Z^{\prime}$ and $W=W^{\prime} \backslash Z^{\prime}$. This arc traversal problem can be stated in algebraic terms. First, remark that, by definition:

[^4]$\gamma^{k}(Z, W)=\mathcal{V}\left(\mathcal{S}_{k}^{Z, W}\right) \cap \mathcal{D}^{k}$ and $\rho^{k}\left(Z^{\prime}, W^{\prime}\right)=\mathcal{V}\left(\mathcal{S}_{k}^{Z^{\prime}, W^{\prime}}\right) \cap \mathcal{D}^{k}$, where $\mathcal{V}(\mathcal{S})$ is the set of solutions of $(\mathcal{S})$, ignoring the domain constraint. In algebraic terms, $\mathcal{V}(\mathcal{S})$ is called an affine variety [5]. When the system $(\mathcal{S})$ is non-degenerate, the variety is a single point for a node and has dimension 1 for an arc. The systems can be solved approximately numerically using off-the-shelf solvers. Instead, we use an "exact" algebraic solver, based on Groebner basis computation. This kind of solver may be slower in practice and may limit the size of games which can be solved. However, it provides guaranteed "exact" finite representations of node coordinates and is especially useful in the case of degenerate games. We will discuss in the conclusion how to adapt our approach in order to use faster, approximate solvers. In our implementation, we use the functions implemented in the Singular toolbox, accessible from the Sagemath environment ${ }^{6}$, to compute Groebner bases and varieties.

The arc traversal problem at level $k$ consists, given the two pairs $\left(Z^{\prime}, W^{\prime}\right)$ and $(Z, W)$, in computing $\left(Z^{\prime \prime}, W^{\prime \prime}\right) \in I_{k}$ corresponding to the following $(n, i)$-almost complementary node $\rho^{k}\left(Z^{\prime \prime}, W^{\prime \prime}\right)$. Algorithm 1 computes this almost-complementary node by trying every possible constraint additions ${ }^{7}$.

```
Algorithm 1: TraverseArc \(\left((Z, W),\left(Z^{\prime}, W^{\prime}\right), I_{k}\right)\).
    /* Computes the end node of \(\operatorname{arc} \gamma^{k}(Z, W)\),
        given starting almost-complementary node
        \(\rho^{k}\left(Z^{\prime}, W^{\prime}\right)\). */
    /* Initialization */
    Sol \(\leftarrow \emptyset\);
    for \((v, j) \in I_{k}\) do
        if \((v, j) \in I_{k} \backslash Z^{\prime}\) then
            \(Z_{l o c} \leftarrow Z \cup\{(v, j)\}, W_{l o c} \leftarrow W\);
            if \(\operatorname{dim}\left(\mathcal{V}\left(\mathcal{S}_{k}^{Z_{\text {loc }}, W_{\text {loc }}}\right)\right)=0\) then
                \(\rho_{l o c} \leftarrow \mathcal{V}\left(\mathcal{S}_{k}^{Z_{l o c}, W_{l o c}}\right) \cap \mathcal{D}^{k} ;\)
                if \(\rho_{\text {loc }} \neq \emptyset\) then
                    Sol \(\leftarrow \operatorname{Sol} \cup\left\{\left(Z_{l o c}, W_{l o c}, \rho_{l o c}\right)\right\} ;\)
        if \((v, j) \in I_{k} \backslash W^{\prime}\) then
            \(Z_{l o c} \leftarrow Z, W_{l o c} \leftarrow W \cup\{(v, j)\} ;\)
            if \(\operatorname{dim}\left(\mathcal{V}\left(\mathcal{S}_{k}^{Z_{l o c}, W_{l o c}}\right)\right)=0\) then
                \(\rho_{l o c} \leftarrow \mathcal{V}\left(\mathcal{S}_{k}^{Z_{\text {loc }}, W_{\text {loc }}}\right) \cap \mathcal{D}^{k} ;\)
                if \(\rho_{\text {loc }} \neq \emptyset\) then
                    Sol \(\leftarrow \operatorname{Sol} \cup\left\{\left(Z_{l o c}, W_{l o c}, \rho_{l o c}\right)\right\} ;\)
    return Sol
```

When the game is non-degenerate, a unique constraint addition will lead to an almost-complementary node:

Proposition 3.10 (Arc traversal correctness). When the PCP is non-degenerate, Algorithm 1 returns a single triple $\left(Z^{\prime \prime}, W^{\prime \prime}, \mathcal{V}\left(\mathcal{S}_{k}^{Z^{\prime \prime}, W^{\prime \prime}}\right)\right.$ and $\mathcal{V}\left(\mathcal{S}_{k}^{Z^{\prime \prime}, W^{\prime \prime}}\right)$ is a single point.

[^5]
### 3.5 The path-following procedure

Now, we have nearly all the elements necessary to build a pathfollowing procedure to reach a complementary node of a PCP. We only lack an initialization procedure. This procedure uses the arbitrary pure joint strategy $\omega^{0}$ and consists in solving $\left(\mathcal{S}^{1}\right)$, decribed in Definition 3.6. Once we have this complementary node at level 1 , we compute the corresponding initial node at level 2 , using Proposition 3.7 and follow a path at level 2 , starting from this node. If we reach a complementary node at level 2 , we climb to level 3 , etc. If, at some level $k$, we reach an initial node, then we compute a new complementary node at level $k-1$, using node descent (Proposition 3.9), from which we go on. In the case where the game is non-degenerate, since we start from a complementary node at level 1 and since the property that every node except the initial node at level 1 and complementary node at level $N$ has exactly two neighbours, the path followed is unique, for a given $\omega^{0}$. Furthermore, it can only end in a complementary node at level $N$ : A solution of the PCP/game. Figure 2 illustrates a few steps of the algorithm.


Figure 2: Portion of a path followed by the algorithm.

## 4 HANDLING DEGENERATE PCP

Recall (Definition 3.1) that there are two ways in which a PCP at level $k$ can be degenerate: (i) if there are points satisfying more than $\left|I_{k}\right|$ equations and (ii) if some systems admit more than one solution. In the first case, which is far more frequent than the second, things may go wrong with the arc traversal algorithm (Algorithm 1). When we follow, during arc traversal or node lifting, an arc $\gamma^{k}(Z, W)$, we normally encounter a unique feasible almost complementary node, either $\rho^{k}(Z \cup\{z\}, W)$ or $\rho^{k}(Z, W \cup\{w\})$. However, for some PCP, it may happen that an arc-traversal makes more than one new constraint binding. For example, there may exist $w_{1}$ and $w_{2}$, such that $\rho^{k}\left(Z, W \cup\left\{w_{1}\right\}\right)$ and $\rho^{k}\left(Z, W \cup\left\{w_{2}\right\}\right)$ define the same feasible solution (identical coordinates). Such a node is degenerate. While traversal of non-degenerate nodes is unambiguous, it is ambiguous for degenerate nodes, since they neighbour three or more almostkomplementary arcs. In this case, an arbitrary choice of the next arc may lead the algorithm to cycle. Figure 3 shows an example game with some degenerate nodes, and paths linking them.

In this example, nodes $1,2,3$ and 4 are degenerate. The corresponding graph contains cycles ( $\{1,2,3\},\{2,3,4\},\{1,2,4,3\}$ ). However, this graph always contains at least one path from the initial


Figure 3: A PCP where four degenerate almostcomplementary nodes are encountered. Node 5 is a complementary node.
node at level 0 toward a complementary node at level $N^{8}$. Taking this fact into account, we may alter the path-following algorithm by applying depth-first search over arcs issued from degenerate nodes, as shown in Algorithm 2.

```
Algorithm 2: Depth-First Graph Traversal
    Data: A PCP and a first degenerate almost-complementary
                node, firstdegennode
    Result: A complementary node
    listofarcs \(\leftarrow \emptyset\);
    listofarcs.append(firstdegennode.outgoingarcs);
    while listofarcs \(=\emptyset\) do
        currentarc \(\leftarrow\) listofarcs.pop ();
        nextnode \(\leftarrow\) traversepath(currentarc);
        if nextnode.is_complementary then break;
        if nextnode \(=\emptyset\) then currentarc \(\leftarrow\) listofarcs.pop();
        else listofarcs.append(nextnode.outgoingarcs);
    return nextnode
```

The procedure traversepath() performs node traversal along a path of non-degenerate nodes, until either (i) the complementary node is reached, (ii) a new degenerate node is reached or (iii) a degenerate node which has already been explored is reached. In cases (i) and (ii) the variable nextnode is set to the corresponding node, while in case (iii) it takes value $\emptyset$.

Applying Algorithm 2 to the example of Figure 3, assuming that we append degenerate nodes' outgoing arcs from top to bottom in listofarcs (different heuristics may be used to decide in which order arcs should be appended), we get the following sequence of lists of arcs: $\{(1, u p),(1$, down $)\} \rightarrow\{(1, u p),(2, u p),(2$, middle $),(2$, bottom $)\}$ $\rightarrow\{(1, u p),(2, u p),(2$, middle $)\} \rightarrow\{(1, u p),(2, u p),(4$, up $),(4$, bottom $)\}$ $\rightarrow$ Node 5 is returned.

Note that in the case of LCP (i.e. for bimatrix games or polymatrix games), a heuristic to order the arcs to append to the list is known that allows Algorithm 2 to never backtrack [22]. This heuristic, based on a lexicographic perturbation of linear systems, is useful since it avoids unnecessary arc traversals, which are costly. The same kind of approach is likely to be applicable to PCP as well. However, computing the heuristic for PCP would require to solve polynomial systems, so the overall benefit is unclear. We leave

[^6]the design of a lexicographic approach for PCP and its evaluation for further research. We should also mention the second case of degeneracy, corresponding to the case where some $\mathcal{V}\left(\mathcal{S}_{k}^{Z, W}\right)$ encountered in the course of the algorithm is a variety of strictly positive dimension (an arc, an hypersurface...) because its defining equations are redundant. While the use of a Groebner basis approach allows to exactly represent such positive-dimensional varieties in parameterized form, it is not clear how such varieties should be traversed in the course of our procedure ${ }^{9}$. We defer the design of a non-zero dimensional varieties traversal method to further work. Instead, what we suggest to do in this case of degeneracy is to apply a small (non-infinitesimal) random perturbation to the coefficients of the initial PCP and solve this new perturbed PCP. The perturbed PCP will generally not be degenerate. So doing, we obtain an approximate Nash equilibrium when this form of degeneracy occurs.

## 5 GRAPHICAL/HYPERGRAPHICAL GAMES

The PCP approach extends naturally to polymatrix games [24], graphical games [16] and hypergraphical games [20]. These are succinct representations of $N$-player games where the utilities of the players are local, i.e. depend on the strategies of subsets of $P$ only. Hypergraphical games are defined as follows:

Definition 5.1 (Hypergraphical game). A $N$-player hypergraphical game, $\Gamma^{N}$, is defined as:
$\Gamma^{N}=\left(\left(P_{g}\right)_{g=1, \ldots, G},\left(S_{n}\right)_{n \in P},\left(a^{g}=\left(a_{\omega_{P_{g}}}^{g, n}\right)_{n \in P_{g}, \omega_{P_{g}} \in \pi_{P_{g}}}\right)_{g=1, \ldots, G}\right)$.

- $P_{g} \subseteq P=\{1, \ldots, N\}, \forall g \in 1, \ldots, G . P_{g}$ is the set of players of the $g^{\text {th }}$ local game and $\cup_{g=1, \ldots, G} P_{g}=P$.
- $\left(S_{n}\right)_{n \in P}$ is the list of players pure strategies sets.
- $a_{\omega_{P_{g}}}^{g, n}$ is the (positive) disutility that player $n$ gets in local game number $g$ (provided that $n$ belongs to $P_{g}$ ), when the joint strategy of all players in game $g$ is $\omega_{P_{g}}$.
The disutility of player $n \in P$ for joint strategy $\omega$ is $a_{\omega}^{n}=\sum_{g, n \in P_{g}} a_{\omega_{P_{g}}}^{g, n}$.
Polymatrix games and graphical games are specific cases of hypergraphical games. Polymatrix games are characterized by the fact that any local game $g$ involves exactly two players: $\left|P_{g}\right|=$ $2, \forall g=1, \ldots, G$. Graphical games are hypergraphical games where the utility of any player $n \in P$ only depends on the strategies of a subset $P_{n} \in P$ of players. They are characterized by the fact that $G=N$ (there is one local game attached to each player) and that $a_{\omega_{P_{g}}}^{g, n}=0, \forall \omega \in \pi, \forall n \neq g$.

Since a hypergraphical game can be represented as a normal form game (potentially exponentially larger to express), NE computation in hypergraphical games admits an exponential size PCP formulation. Fortunately, we can exploit the factorisation of the disutility functions of a hypergraphical game in order to compute a corresponding PCP of "reasonable" size. First, we can show that:

Proposition 5.2 (PCP factorization).

$$
A_{i}^{n}\left(x^{-n}\right)=\sum_{g, n \in g} Q_{g}\left(x^{P \backslash P_{g}}\right) R_{n, i, g}\left(x^{P_{g} \backslash\{n\}}\right) \text {, where }
$$

[^7]\[

$$
\begin{aligned}
Q_{g}\left(x^{P \backslash P_{g}}\right) & =\prod_{v \in P \backslash P_{g}}\left(\sum_{\omega_{v} \in S_{v}} x_{\omega_{v}}^{v}\right) \text { and } \\
R_{n, i, g}\left(x^{P_{g} \backslash\{n\}}\right) & =\sum_{\substack{\omega_{P_{g}} \\
\omega_{n}=i}} a_{\omega_{P_{g}}}^{g, n} \prod_{v \in P_{g} \backslash\{n\}} x_{\omega_{v}}^{v} .
\end{aligned}
$$
\]

Proof of Proposition 5.2: The disutility of player $n \in N$ for joint strategy $\omega$ is:

$$
a_{\omega}^{n}=\sum_{g, n \in P_{g}} a_{\omega_{P_{g}}}^{g, n}
$$

Using this expression of $a_{\omega}^{n}$, we can rewrite $A_{i}^{n}\left(x^{-n}\right)$ :

$$
\begin{aligned}
& A_{i}^{n}\left(x^{-n}\right)=\sum_{\substack{\omega \in \pi \\
\omega_{n}=i}}\left(\sum_{g, n \in P_{g}} a_{\omega_{P_{g}}}^{g, n}\right) \prod_{v \neq n} x_{\omega_{v}}^{v}, \forall n \in P, i \in S_{n}, \\
&=\sum_{g, n \in P_{g}} \sum_{\substack{\omega \in \pi \\
\omega_{n}=i}} a_{\omega_{P_{g}}}^{g, n} \prod_{v \neq n} x_{\omega_{v}}^{v}, \\
&\left.=\sum_{g, n \in P_{g}} \sum_{\substack{\omega_{P_{g}} \\
\omega_{n}=i}}^{a_{\omega_{P_{g}}}^{g, n}} \prod_{v, n \in P_{g}} \prod_{\omega_{\omega_{2}}=i} x_{\omega_{g} \backslash\{n\}}^{v} a_{\omega_{v}}^{g, n}\right) \sum_{\omega_{P \backslash \backslash P_{g}}} \prod_{v \in P \backslash P_{g}} x_{\omega_{v}}^{v}, \\
&\left.\prod_{v \in P_{g} \backslash\{n\}} x_{\omega_{v}}^{v}\right)\left(\prod_{v \in P \backslash P_{g}}\left(\sum_{\omega_{v} \in S_{v}} x_{\omega_{v}}^{v}\right)\right) . \\
& \text { Remark that equality } \sum_{\omega_{P \backslash P_{g}}} \prod_{v \in P \backslash P_{g}} x_{\omega_{v}}^{v}=\prod_{v \in P \backslash P_{g}}\left(\sum_{\omega_{v} \in S_{v}} x_{\omega_{v}}^{v}\right)
\end{aligned}
$$

results from the repeated application of the distributivity of the product over the addition. The proposition follows.

Note that $A_{i}^{n}\left(x^{-n}\right)$ is still a polynomial in variables $x_{j}^{m},(m, j) \neq$ ( $n, i$ ) with a number of terms of degree $N-1$ equal to $\left|\pi_{-n}\right|$. But it is known, for example, that polymatrix games admit a linear complementarity problem formulation [14]. We show that it is possible to get similar savings in representation size and solution complexity for hypergraphical games, through the addition of auxiliary variables.

Definition 5.3 (PCP auxiliary variables). In addition to the variables $\left\{x_{i}^{n}\right\}$, let us consider an additional set of variables, $\left\{y_{g}^{n}\right\}_{g=1 . . G, n \in P}$.

We define these variables as:
$y_{g}^{n}=\prod_{v=1}^{n} \alpha_{g}^{v}$, where $\alpha_{g}^{v}=1$ if $v \in P_{g}$ and $\alpha_{v}^{g}=\sum_{\omega_{v} \in S_{v}} x_{\omega_{v}}^{v}$ if $v \in P \backslash P_{g}$.
With this definition, we have $Q_{g}\left(x^{P \backslash P_{g}}\right)=y_{g}^{N}$ and

$$
\begin{equation*}
y_{g}^{1}=\alpha_{g}^{1} \text { and } y_{g}^{n}=y_{g}^{n-1} \times \alpha_{g}^{n}, \forall n=2, \ldots, N \tag{6}
\end{equation*}
$$

Note that there are exactly $G \times N$ additional variables $y_{g}^{n}$ and their values are defined by equations of degree 1 when $n \in P_{g}$ and degree 2 when $n \in P \backslash P_{g}$. This gives the following expression for $A_{i}^{n}$ :

$$
A_{i}^{n}\left(x^{-n},\left\{y_{g}^{N}\right\}_{g=1, \ldots, G}\right)=\sum_{g, n \in P_{g}} y_{g}^{N} \sum_{\substack{\omega_{P_{g}} \\ \omega_{n}=i}} a_{\omega P_{g}}^{g, n} \prod_{v \in P_{g} \backslash\{n\}} x_{\omega_{v}}^{v}
$$

Remark that the polynomials $A_{i}^{n}\left(x^{-n},\left\{y_{g}^{N}\right\}_{g=1, . ., G}\right)$ now have degree at most $\max _{g=1, \ldots, G}\left|P_{g}\right|$ and the size of the expression of these polynomials is comparable to that of the hypergraphical game.

Our algorithm can now be extended to deal with this new expression with additional variables. The only change in the algorithm is the definition of the polynomial subsystems which now have to consider the $y_{g}^{n}$ variables. We extend the definition of the polynomial systems $\mathcal{S}_{k}^{Z, W}$ to $\overline{\mathcal{S}}_{k}^{Z, W}$ for pairs $Z, W \subseteq I_{k}$ :

$$
\left\{\begin{aligned}
& x \in \mathcal{D}, \\
x_{\omega_{n}^{0}}^{n}=1 \text { and } x_{i}^{n}=0, & \forall n>k, \forall i \neq \omega_{n}^{0} \\
x_{i}^{n}=0, & \forall(n, i) \in Z, \\
y_{g}^{1}=\alpha_{g}^{1}, & \forall g=1, \ldots, G \\
y_{g}^{n}=y_{g}^{n-1} \times \alpha_{g}^{n}, & \forall g=1 . . G, \forall n=2 . . k \\
A_{i}^{n, k}\left(x^{\{1 . . k\} \backslash\{n\}},\left\{y_{g}^{k}\right\}_{g=1, \ldots, G}\right)=1, & \forall(n, i) \in W
\end{aligned}\right.
$$

With this generalization to hypergraphical games, it results that the number of terms of the polynomials of the systems is reduced and their degrees are bounded by the number of players of the largest subgame. This has an importance when it comes to computing Groebner bases. The Ideal Membership Problem, which is the core problem of the Groebner basis computation problem is known to be ExpSpace-complete. However, when the ideal has dimension 0 (which is the case when $\left(\overline{\mathcal{S}}_{k}^{Z, W}\right)$ has a finite number of solutions), it is solvable in single exponential time [7]. Furthermore, being able to upper bound the degrees of the polynomials involved in the PCP has a positive impact on the complexity of the algorithm:

Proposition 5.4 (Complexity of the path-following algoRITHM). The time complexity of the path-following algorithm for a non-degenerate graphical/hypergraphical game is simply exponential in $\left|I_{N}\right|$ and doubly exponential in the maximal number of players of any local game.

Proof of Proposition 5.4: Indeed, The worst-case time complexity of the best known Groebner basis computation algorithm is doubly exponential in the maximal degree of the involved polynomials (see e.g. [19]). The number of arc traversal steps is bounded by the number of almost complementary nodes at any level (with respect to a fixed $\omega_{0}$. Thus, at level $k$ it is bounded by \#Nodes $(k)=$ $\left|\left\{(Z, W),|Z|+|W|=\left|I_{k}\right|,|Z \cap W| \leq 1\right\}\right|$. Then, remark that, for any hypergraphical game, including normal form games: \#Nodes $(k) \leq$ $\left|I_{k}\right| 2^{\left|I_{k}\right|}$. Indeed, potential nodes can be built by choosing an arbitrary $Z \subseteq I_{k}$ and then there exist at most $\left|I_{k}\right| W$ potentially leading to a $(k, j)$-almost complementary node. So, the total number of almost-complementary nodes for all levels is $O\left(N\left|I_{N}\right| 2^{\left|I_{N}\right|}\right)$. Since each arc traversal requires at most $O\left(\left|I_{N}\right|\right)$ Groebner basis computations, we get the result ${ }^{10}$.

## 6 EXPERIMENTS

The support enumeration and path-following algorithms were implemented in Python 3 using the Sagemath software for the computation of varieties. Each game was solved on a single node of a

[^8]HPC platform. Nodes were identical Intel Xeon E5-2680 v4 2,4 Ghz bi-processors with 128 Gb of RAM, under Linux OS.

The benchmark was composed of normal-form and hypergraphical games with 3 to 7 players, 2 to 4 actions and hyperedges of size 3 (for HGG). They were generated using the GAMUT ${ }^{11}$ suite. We generated Covariant and Random games with integer valued utilities in range $0 . .100$. In total, 100 games were solved for each configuration, using support enumeration (SE) and path-following (PF). Hypergraphical games were converted to normal-form prior to being solved using support enumeration. Before applying SE or PF , pure equilibrium search was performed. After this step there remained from 10 to 40 unsolved games in each configuration. For these, we computed the average NE support size per player found using SE or PF , as well as the time required to solve them. A timeout of 5 minutes per run was used for SE and PF. Figure 4 shows the results obtained for random games with 5 players and 3 actions ( 3 local games of size 3 for hypergraphical games). All other tested configurations are shown in the supplementary material.


Figure 4: Experiment on 100 Random game. 66 games with PNE in NFG, 60 in HGG.

Figure 4, top, shows some features generally observed for all configurations. First, as can be expected, PF solutions will present a larger variety of support sizes than SE and larger in general. Still, it happens, rarely, that PF finds solutions of smaller support size than SE. This can (and does) happen in the case of degenerate games where the equilibrium with smallest support is at one end of an arc of solutions of dimension 1 . SE will not find this solution while PF may encounter this complementary node. This is why PF finds equilibria of support size 6 when SE does not, in Figure 4.

Globally, we found that SE was more efficient than PF for games with no more than 6 players and 2 actions. PF is more efficient for games with at least 3 actions per player. With 7 players, 2 actions,

[^9]both approaches have similar performances. It confirms that when the total number of potential supports increases, Wilson's approach becomes more efficient than blind enumeration and removal of dominated alternatives. Similar conclusions hold for random and covariant games. As far as HGG are concerned, we found that the direct PF approach did not outperform a prior translation to NFG. In the study cases, it looks like the cost of adding variables $y$ is not compensated by the decreased polynomial degrees and number of terms. Studying whether it is possible to improve the PF approach's performance in case of HGG is left for further research.

## 7 CONCLUDING REMARKS

There is an abundant literature on algorithms for approximate equilibrium search in $N$-player games, including homotopy methods [ $3,10,11,13$ ]. These are based on the definition of a parametrized continuum of games, joining an arbitrary game with known equilibrium, to the game of interest. An "arc" of equilibria of the parametrized games is followed and [6] have used Groebner bases to solve the specifically designed (easy) initial game. Homotopy approaches are prone to numerical errors and potential non-convergence. Uniform strategy enumeration methods [1, 2, 18], on the other hand, suggest to enumerate a space of discretized mixed strategies in order to find a $\varepsilon$-approximate strategy. Polynomial systems-based approaches, such as [21,23] also rely on a form of enumeration. [21] enumerates and solves a sequence of systems formed from mixed strategies supports of increasing size, when [23] explores a deterministic sequence of systems. We provided the first implementation of [23], extended to degenerate and succinct games. Our approach explores a sequence of neighbour almost-complementary points defined as solutions of systems of equations at different levels. It differs from [21] which blindly explores every possible systems in increasing order of supports' sizes. When there exists a NE with small support, the method [21] is faster, while in other cases [23] is more efficient.

The building block of PCP methods, the polynomial system solver, can be easily adapted to follow the progresses in solvers' development. We used an exact algebraic solver. So doing, node's coordinates are algebraic numbers [4] that is, roots of polynomials which coefficients are either rational numbers, or algebraic numbers themselves. Thus they are finitely exactly represented. While exactness is desirable in Wilson's algorithm in order to distinguish degenerate from non-degenerate nodes, it comes at a computational cost. Wilson's arc traversal procedure can be adapted to use approximate solvers, by using depth-first search in the way we use it for dealing with degeneracy. It is enough to relax the condition that a node is degenerate whenever it corresponds to points $\rho^{k}(Z, W)$ and $\rho^{k}\left(Z^{\prime}, W^{\prime}\right)$ of identical coordinates: We may consider that $(Z, W)$ and ( $Z^{\prime}, W^{\prime}$ ) correspond to a single degenerate node whenever the coordinates of $\rho^{k}(Z, W)$ and $\rho^{k}\left(Z^{\prime}, W^{\prime}\right)$ are close enough. So doing, system solving will be faster, but the number of explored almost complementary nodes may increase, due to "false" degeneracy.

Finally, our approach may be naturally extended to other kinds of games, including Bayesian games [12] and stochastic games [8]. Indeed, it is known that two-player bayesian games can be represented as polymatrix games [15]. The same relation is likely to hold between $N$-player bayesian games and hypergraphical games, making a concise PCP formulation possible.

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[^0]:    Proc. of the 21st International Conference on Autonomous Agents and Multiagent Systems (AAMAS 2022), P. Faliszewski, V. Mascardi, C. Pelachaud, M.E. Taylor (eds.), May 9-13, 2022, Online. © 2022 International Foundation for Autonomous Agents and Multiagent Systems (www.ifaamas.org). All rights reserved.

[^1]:    ${ }^{1}$ Python/Sagemath codes as well as tutorial Jupyter notebooks are available online: https://forgemia.inra.fr/game-theory-tools-group/gtnash-git/.
    ${ }^{2}$ The PCP formulation considers that players are disutility minimizers. This is without loss of generality since any utility maximization game can be equivalently expressed as a disutility minimization game.

[^2]:    ${ }^{3}$ Thus, a complementary point is ( $n, i$ )-almost complementary for every pair ( $n, i$ ).

[^3]:    ${ }^{4}$ The proofs of this Section are modern rewritings of the proofs of Wilson's original paper. The reader will find them, as well as a tutorial example and additional experiments in https://doi.org/10.5281/zenodo. 5850463.

[^4]:    ${ }^{5}$ The exponent $k$ in $\gamma^{k}$ or $\rho^{k}$ indicates that we are at level $k$.

[^5]:    ${ }^{6}$ https://www.sagemath.org/index.html.
    ${ }^{7}$ In practice, as soon as a solution $\rho_{\text {loc }}$ is found in line 9 of Algorithm 1, the main loop is exited and it is tested whether the current node is degenerate, by checking whether additional constraints indexed in $Z$ or $W$ are satisfied by the coordinates of $\rho_{\text {loc }}$.

[^6]:    ${ }^{8}$ This is a consequence of the fact that a random infinitesimal perturbation of the coefficients of the PCP will resolve degeneracy by making some of the binding constraints non-binding. In the process, the new non-degenerate nodes will be defined by pairs ( $Z^{\prime}, W^{\prime}$ ) where $Z^{\prime} \subseteq Z$ and $W^{\prime} \subseteq W$.

[^7]:    ${ }^{9}$ When such varieties correspond to sets of complementary points at level $N$, we can directly provide the continuous set of equilibrium strategies in parameterized form.

[^8]:    ${ }^{10}$ This complexity bound is theoretical. In practice the length of the paths is often small.

[^9]:    ${ }^{11} \mathrm{http}: / /$ gamut.stanford.edu/

