Multi-Scale Games: Representing and Solving Games on Networks with Group Structure

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Abstract

Network games provide a natural machinery to compactly represent strategic interactions among agents whose payoffs exhibit sparsity in their dependence on the actions of others. Besides encoding interaction sparsity, however, real networks often exhibit a multi-scale structure, in which agents can be grouped into communities, those communities further grouped, and so on, and where interactions among such groups may also exhibit sparsity. We present a general model of multi-scale network games that encodes such multi-level structure. We then develop several algorithmic approaches that leverage this multi-scale structure, and derive sufficient conditions for convergence of these to a Nash equilibrium. Our numerical experiments demonstrate that the proposed approaches enable orders of magnitude improvements in scalability when computing Nash equilibria in such games. For example, we can solve previously intractable instances involving up to 1 million agents in under 15 minutes.

Introduction

Strategic interactions among interconnected agents are commonly modeled using the network, or graphical, game formalism (Kearns, Littman, and Singh 2001; Jackson and Zenou 2015). In such games, the utility of an agent depends on his own actions as well as those by its network neighbors. Many variations of games on networks have been considered, with applications including the provision of public goods (Allouch 2015; Buckley and Croson 2006; Khalili, Zhang, and Liu 2019; Yu et al. 2020), security (Hota and Sundaram 2018; La 2016; Vorobeychik and Letchford 2015), and financial markets (Acemoglu et al. 2012).

While network games are a powerful modeling framework, they fail to capture a common feature of human organization: groups and communities. Indeed, investigation of communities, or close-knit groups, in social networks is a major research thread in network science. Moreover, such groups often have a hierarchical structure (Clauset, Moor, and Newman 2008; Girvan and Newman 2002). For example, strategic interactions among organizations in a marketplace often boil down to interactions among their constituent business units, which are, in turn, comprised of individual decision makers. In the end, it is those lowest-





Figure 1: An illustration of a multi-scale (3-level) network.

level agents who ultimately accrue the consequences of these interactions (for example, corporate profits would ultimately benefit individual shareholders). Moreover, while there are clear interdependencies among organizations, individual utilities are determined by a combination of individual actions of some agents, together with *aggregate* decisions by the groups (e.g., business units, organizations). For example, an employee's bonus is determined in part by their performance in relation to their co-workers, and in part by how well their employer (organization) performs against its competitors in the marketplace.

We propose a novel *multi-scale game model* that generalizes network games to capture such hierarchical organization of individuals into groups. Figure 1 offers a stylized example in which three groups (e.g., organizations) are comprised of 2-3 subgroups each (e.g., business units), which are in turn comprised of 2-5 individual agents. Specifically, our model includes an explicit hierarchical network structure that organizes agents into groups across a series of levels. Further, each group is associated with an action which deterministically aggregates the decisions by its constituent agents. The game is grounded at the lowest level, where the agents are associated with scalar actions and utility functions that have modular structure in the strategies taken at each level of the game. For example, in Figure 1, the utility function of an individual member a_j of level-3 group $a_3^{(3)}$ is a function of the strategies of (i) a_j 's immediate neighbors (represented by links between pairs of filled-in circles), (ii)

 a_j 's level-2 group and its network neighbor (the small hollow circles), and (iii) a_j 's level-3 group, $a_3^{(3)}$ (large hollow circle) and its network neighbors, $a_1^{(3)}$ and $a_2^{(3)}$.

Our next contribution is a series of iterative algorithms for computing pure strategy Nash equilibria that explicitly leverage the proposed multi-scale game representation. The first of these simply takes advantage of the compact game representation in computing equilibria. The second algorithm we propose offers a further innovation through an iterative procedure that alternates between game levels, treating groups themselves as pseudo-agents in the process. We present sufficient conditions for the convergence of this algorithm to a pure strategy Nash equilibrium through a connection to Structured Variational Inequalities (He, Yang, and Wang 2000), although the result is limited to games with two levels. To address the latter limitation, we design a third iterative algorithm that now converges even in games with an arbitrary number of levels.

Our final contribution is an experimental evaluation of the proposed algorithms compared to best response dynamics. In particular, we demonstrate orders of magnitude improvements in scalability, enabling us to solve games that cannot be solved using a conventional network game representation.

Related Work: Network games have been an active area of research; see e.g., surveys by Jackson and Zenou (2015) and Bramoullé and Kranton (2016). We now review the most relevant papers.

Conditions for the existence, uniqueness and stability of Nash equilibria in network games under general best responses are studied in (Parise and Ozdaglar 2019; Naghizadeh and Liu 2017; Scutari et al. 2014; Bramoullé, Kranton, and D'amours 2014). Variational inequalities (VI) are used in these works to analyze the fixed point and contraction properties of the best response mappings. It is identified in (Parise and Ozdaglar 2019; Naghizadeh and Liu 2017; Scutari et al. 2014) that when the Jacobian matrix of the best response mapping is a P-matrix or is positive definite, a feasible unique Nash equilibrium exists and can be obtained by best-response dynamics (Scutari et al. 2014; Parise and Ozdaglar 2019).

In this paper, we extended the analysis of equilibrium and best responses for a conventional network game to that in a multi-scale network game, where the utility functions are decomposed into separable utility components to which best responses are applied separately. This is similar to the generalization from a conventional VI problem to an SVI problem (He, Yang, and Wang 2000; He 2009; He and Yuan 2012; Bnouhachem, Benazza, and Khalfaoui 2013) problem.

Previous works on network games that involve group or community structure focus on finding such structures; e.g., community detection in networks using game theoretic methods have been studied in (Mcsweeney, Mehrotra, and Oh 2017; Newman 2004; Alvari, Hajibagheri, and Sukthankar 2014).

By contrast, our work focuses on analyzing a network game with a given group/community structure, and using the structure as an analytical tool for the analysis of equilibrium and best responses.

Preliminaries

A general *normal-form game* is defined by a set of agents (players) $I = \{1, ..., N\}$, with each agent a_i having an action/strategy space K_i and a utility function $u_i(x_i, \boldsymbol{x}_{-i})$ that i aims to maximize; $x_i \in K_i$ and x_{-i} denotes the actions by all agents other than i. We term the collection of strategies of all agents \boldsymbol{x} a *strategy profile*. We assume $K_i \subset \mathbb{R}$ is a compact set.

We focus on computing a *Nash equilibrium* (*NE*) of a normal-form game, which is a strategy profile with each agent maximizing their utility given the strategies of others. Formally, x^* is a *Nash equilibrium* if for each agent *i*,

$$x_i^* \in \underset{x_i \in K_i}{\operatorname{argmax}} \quad u_i(x_i, \boldsymbol{x}_{-i}^*). \tag{1}$$

A network game encodes structure in the utility functions such that they only depend on the actions by network neighbors. Formally, a network game is defined over a weighted graph (I, E), with each node an agent and E is the set of edges; the agent's utility $u_i(x_i, \boldsymbol{x}_{-i})$ reduces to $u_i(x_i, \boldsymbol{x}_{I_i})$, where I_i is the set of network neighbors of i, although we will frequently use the former for simplicity.

An agent's best response is its best strategy given the actions taken by all the other agents. Formally, the best response is a set defined by

$$BR_i(\boldsymbol{x}_{-i}, u_i) = \operatorname*{argmax}_{x_i} u_i(x_i, \boldsymbol{x}_{-i}).$$
(2)

Whenever we deal with games that have a unique best response, we will use the singleton best response set to also refer to the player's best response strategy (the unique member of this set).

Clearly, a NE of a game is a fixed point of this best response correspondence. Consequently, one way to compute a NE of a game is through *best response dynamics* (*BRD*), which is a process whereby agents iteratively and asynchronously (that is, one agent at a time) take the others' actions as fixed values and play a best response to them.

We are going to use this BRD algorithm as a major building block below. One important tool that is useful for analyzing BRD convergence is *Variational Inequalities (VI)*. To establish the connection between NE and VI we assume the utility functions $u_i, \forall i = 1, ..., N$, are continuously twice differentiable. Let $K = \prod_{i=1}^{N} K_i$ and define $F : \mathbb{R}^N \to \mathbb{R}^N$ as follows:

$$F(\boldsymbol{x}) := \left(-\nabla_{x_i} u_i(\boldsymbol{x}) \right)_{i=1}^N.$$
(3)

Then \boldsymbol{x}^* is said to be a solution to VI(K, F) if and only if

$$(\boldsymbol{x} - \boldsymbol{x}^*)^T F(\boldsymbol{x}^*) \ge 0, \ \forall \boldsymbol{x} \in K.$$
 (4)

In other words, the solution set to VI(K, F) is equivalent to the set of NE of the game. Now, we can define the condition that will guarantee the convergence of BRD.

Definition 1. The P_{Υ} condition: We denote $\alpha_i(F) = \inf_{\boldsymbol{x} \in K} ||\nabla_i F_i||_2$, $\beta_{i,j}(F) = \sup_{\boldsymbol{x} \in K} ||\nabla_j F_i||_2$, $i \neq j$. The

 Υ matrix generated from $F:\mathbb{R}^N\to\mathbb{R}^N$ is given as follows

$$\Upsilon(F) = \begin{bmatrix} \alpha_1(F) & -\beta_{1,2}(F) & \cdots & -\beta_{1,N}(F) \\ -\beta_{2,1}(F) & \alpha_2(F) & \cdots & -\beta_{2,N}(F) \\ \vdots & \vdots & \ddots & \vdots \\ -\beta_{N,1}(F) & -\beta_{N,2}(F) & \cdots & \alpha_N(F) \end{bmatrix},$$

If $\Upsilon(F)$ is a P-matrix, that is, if all of its principal minors have a positive determinant, then we say F satisfies the P_{Υ} condition.

Theorem 1. (Scutari et al. 2014) If F satisfies the P_{Υ} condition, then F is strongly monotone on K, and VI(K, F) has a unique solution. Moreover, BRD converges to the unique NE from an arbitrary initial state.

A Multi-Scale Game Model

Consider a conventional network (graphical) game with the set I of N agents situated on a network G = (I, E), each with a utility function $u_i(x_i, \boldsymbol{x}_{I_i})$, with I_i the set of i's neighbors, I the full set of agents/nodes and E the set of edges connecting them.¹ Suppose that this network G exhibits the following structure and feature of the strategic dependence among agents: agents can be partitioned into a collection of groups $\{S_k\}$, where k is a group index, and an agent a_i in the kth group (i.e., $a_i \in S_k$) has a utility function that depends (i) on the strategies of its network neighbors in S_k , and (ii) only on the aggregate strategies of groups other than k (see, e.g., Fig. 1). Further, these groups may go on to form larger groups, whose aggregate strategies impact each other's agents, giving rise to a *multi-scale* structure of the network. This kind of structure is very natural in a myriad of situations. For example, members of criminal organizations take stock of individual behavior by members of their own organization, but their interactions with other organizations (criminal or otherwise) are perceived in group terms (e.g., how much another group has harmed theirs). A similar multi-level interaction structure exists in national or ethnic conflicts, organizational competition in a market place, and politics. Indeed, a persistent finding in network science is that networks exhibit a multi-scale interaction structure (i.e., communities, and hierarchies of communities) (Girvan and Newman 2002; Clauset, Moor, and Newman 2008).

We present a general model to capture such multi-scale structure. Formally, an *L*-level structure is given by a hierarchical graph structure $\{G^{(l)}\}$ for each level $l, 1 \le l \le L$, where $G^{(l)} = (\{S_k^{(l)}\}_k, E^{(l)})$ represents the level-*l* structure. The first component, $\{S_k^{(l)}\}_k$ prescribes a partition, where agents in level l - 1 form disjoint groups given by this partition; each group is viewed as an agent in level l, denoted as $a_k^{(l)}$. Notationally, while both $a_k^{(l)}$ and $S_k^{(l)}$ bear the superscript (l), the former refers to a level-*l* agent, while the latter is the group (of level-(l - 1) agents) that the former represents. The set of level-*l* agents is denoted by $I^{(l)}$ and their total number $N^{(l)}$. The second component, $E^{(l)}$, is

a set of edges that connect level-l agents, encoding the dependence relationship among the groups they represent. This structure is anchored in level 1 (the lowest level), where sets $S_k^{(1)}$ are singletons, corresponding to agents a_k in the game, who constitute the set I.

To illustrate, the multi-scale structure shown in Fig. 1 is given by $G^{(1)} = G = (\{S_k^{(1)}\}_k = I, E^{(1)} = E)$, as well as how level-1 agents are grouped into level-2 agents, how level-2 agents are further grouped into level-3 agents, and the edges connecting these groups at each level.

It should be obvious that the above multi-scale representation of a graphical game is a generalization of a conventional graphical game, as any such game essentially corresponds to a L = 1 multi-scale representation. On the other hand, not all conventional graphical games have a meaningful L > 1 multi-scale representation (with non-singleton groups of level-1 agents); this is because our assumption that an agent's utility only depends on the *aggregate* decisions by groups other than the one they belong to implies certain properties of the dependence structure. For the remainder of this paper we will proceed with a given multi-scale structure defined above, while in the supplementary material we outline a set of conditions on a graphical game G that allows us to represent it in a (non-trivial) multi-scale fashion.

Since the resulting multi-scale network is strictly hierarchical, we can define a *direct supervisor* of agent $a_i^{(l)}$ in level-*l* to be the agent $a_k^{(l+1)}$ corresponding to the level-(l + 1) group *k* that the former belongs to. Similarly, two agents who belong in the same level-*l* group *k* are (level-*l*) group mates. Finally, note that any level-1 agent a_i belongs to exactly one group in each level *l*. We index a level-*l* group to which a_i belongs by k_{il} .

In order to capture the agent dependence on aggregate actions, we define an *aggregation function* $\sigma_k^{(l)}$ for each level-l group k that maps individual actions of group members to \mathbb{R} (a *group strategy*). Specifically, consider a level-l group $S_k^{(l)}$ with level-(l-1) agents in this group playing a strategy profile $\boldsymbol{x}_{S_k^{(l)}}$. The (scalar) group strategy, which is also the strategy for the corresponding level-(l+1) agent, is determined by the aggregation function,

$$x_k^{(l)} = \sigma_k^{(l)}(\boldsymbol{x}_{S_k^{(l)}}).$$
(5)

A natural example of this is linear (e.g., agents respond to total levels of violence by other criminal organizations): $\sigma_k^{(l)}(\boldsymbol{x}_{S_k^{(l)}}) = \sum_{i \in S_k^{(l)}} x_i^{(l)}.$ The *L*-level structure above is captured strategically by

The *L*-level structure above is captured strategically by introducing structure into the utility functions of agents. Let $I_{k_{il}}$ denote the set of neighbors of level-*l* group *k* to which level-1 agent a_i belongs; i.e., this is the set of level-*l* groups that interact with agent a_i 's group. This level-1 agent's utility function can be decomposed as follows:

$$u_i(x_i, \boldsymbol{x}_{-i}) = \sum_{l=1}^{L} u_{k_{il}}^{(l)} \left(x_{k_{il}}^{(l)}, \boldsymbol{x}_{I_{k_{il}}}^{(l)} \right).$$
(6)

In this definition, the level-*l* strategies $x_k^{(l)}$ are implicitly functions of the level-1 strategies of agents that comprise the

¹The edges are generally weighted, resulting in a weighted adjacency matrix on which the utility depends.

Algorithm 1: BRD Algorithm

Initialize the game, $t = 0, x_i(0) = (\mathbf{x}_0)_i, i = 1, \dots, N;$ while not converged do for i = 1:N do $| x_i(t+1) = BR_i(\mathbf{x}_{-i}(t), u_i)$ end $t \leftarrow t+1$ end

group, per a recursive application of Eqn. (5). Consequently, the utility is an additive function of the hierarchy of grouplevel components for increasingly (with l) abstract group of agents. Note that conventional network games are a special case with only a single level (L = 1).

To illustrate, if we consider just two levels (a collection of individuals and groups to which they directly belong), the utility function of each agent a_i is a sum of two components:

$$u_i(x_i, \boldsymbol{x}_{-i}) = u_{k_{i1}}^{(1)} \left(x_{k_{i1}}^{(1)}, \boldsymbol{x}_{I_{k_{i1}}}^{(1)} \right) + u_{k_{i2}}^{(2)} \left(x_{k_{i2}}^{(2)}, \boldsymbol{x}_{I_{k_{i2}}}^{(2)} \right).$$

In the first component, $x_{k_{i1}}^{(1)} = x_i$, since level-1 groups correspond to individual agents, whereas $\boldsymbol{x}_{I_{k_{i1}}}^{(1)}$ is the strategy profile of *i*'s neighbors *belonging to the same group as i*, given by $E^{(1)}$. The second utility component now depends only on the aggregate strategy $x_{k_{i2}}^{(2)}$ of the group to which *i* belongs, as well as the aggregate strategies of the groups with which *i*'s group interacts, given by $E^{(2)}$.

Algorithms and Analysis

Consider the BRD algorithm (formalized in Algorithm 1) in which we iteratively select an agent who plays a best response to the strategy of the rest from the previous iteration.

The conventional BRD algorithm operates on the "flattened" utility function which evaluates utilities explicitly as functions of the strategies played by all agents $a_i \in I$. Our goal henceforth is to develop algorithms that take advantage of the special multi-scale structure and enable significantly better scalability than standard BRD, while preserving the convergence properties of BRD.

Taking Advantage of Multi-Scale Utility Representation

The simplest way to take advantage of the multi-scale representation is to directly leverage the structure of the utility function in computing best responses. Specifically, the multi-scale utility function is more compact than one that explicitly accounts for the strategies of all neighbors of i (which includes *all* of the players in groups other than the one i belongs to). This typically results in a direct computational benefit to computing a best response. For example, in a game with a linear best response, this can result in an exponential reduction in the number of linear operations.

Algorithm 2: Multi-Scale BRD (MS-BRD)

```
Initialize the game,
t = 0, x_i^{(1)}(0) = (\mathbf{x}_0)_i, i = 1, \dots, N
for l = 2:L do
      for k = 1:N^{(l)} do
          m{x}_{k}^{(l)}(0) = \sigma_{k}^{(l)}(m{x}_{S_{i}^{(l)}}(0));
      end
end
while not converged do
      for i = 1:N (Level-1) do
           x_i^{(1)}(t+1) = BR_i(\boldsymbol{x}_{-i}^{(1)}(t), u_i)
       end
      for l = 2:L do
            for k = I: N^{(l)} do

\mathbf{x}_{k}^{(l)}(t+1) = \sigma_{k}^{(l)}(\mathbf{x}_{S_{k}^{(l)}}(t+1));
            end
      end
      t \leftarrow t + 1;
end
```

The resulting algorithm, *Multi-Scale Best-Response Dynamics (MS-BRD)*, which takes advantage of our utility representation is formalized as Algorithm 2. The main difference from BRD is that it explicitly uses the multi-scale utility representation: in each iteration, it updates the aggregated strategies at all levels for the groups to which the most recent best-responding agent belongs. Since MS-BRD simply performs operations identical to BRD but efficiently, its convergence is guaranteed under the same conditions (see Theorem 1). Next, we present iterative algorithms for computing NE that take further advantage of the multi-scale structure, and study their convergence.

Taking Advantage of Multi-Scale Strategic Dependence Structure

In order to take full advantage of the multi-scale game structure, we now aim to develop algorithms that treat groups explicitly as agents, with the idea that iterative interactions among these can significantly speed up convergence. Of course, in our model groups are not actual agents in the game: utility functions are only defined for agents in level 1. However, note that we already have well-defined group strategies - these are just the aggregations of agent strategies at the level immediately below, per the aggregation function (5). Moreover, we have natural utilities for groups as well: we can use the corresponding group-level component of the utility of any agent in the group (note that these are identical for all group members in Eqn. (6)). However, using these as group utilities will in fact not work: since ultimately the game is only among the agents in level 1, equilibria of all of the games at more abstract levels must be consistent with equilibrium strategies in level 1. On the other hand, we need to enforce consistency only between neighboring levels, since that fully captures the across-level interdependence induced by the aggregation function. Therefore, we define the following *pseudo-utility functions* for agents at levels other than 1, with agent k in level l corresponding to a subset of agents from level l - 1:

$$\hat{u}_{k}^{(l)} = u_{k}^{(l)} \left(x_{k}^{(l)}, \boldsymbol{x}_{I_{k}}^{(l)} \right) - L_{k}^{(l,l-1)} \left(x_{k}^{(l)}, \sigma_{k}^{(l)}(\boldsymbol{x}_{S_{k}^{(l)}}) \right) - L_{k}^{(l,l+1)} \left(\sigma_{k}^{(l+1)}(\boldsymbol{x}_{S_{k}^{(l+1)}}), x_{k}^{(l+1)} \right).$$
(7)

The first term is the level-l component of the utility of any level-1 agent in group k. The second and third terms model the inter-level inconsistency loss that penalizes a level-l agent $a_k^{(l)}$, where $L_k^{(l,l+1)}$ and $L_i^{(l,l-1)}$ penalize its inconsistency with the level-(l + 1) and level-(l - 1) entities respectively. In general, $L_k^{(l,l+1)}$ is a different function from $L_k^{(l+1,l)}$; we elaborate on this further below. The central idea behind the second algorithm we propose

is simple: in addition to iterating best response steps at level 1, we now interleave them with best response steps taken by agents at higher levels, which we can since strategies and utilities of these pseudo-agents are well defined. This algorithm is similar to the augmented Lagrangian method in optimization theory, where penalty terms are added to relax an equality constraint and turn the problem into one with separable operators. We can decompose this type of problem into smaller subproblems and solve the subproblems sequentially using the alternating direction method (ADM) (Yuan and Li 2011; Bnouhachem, Benazza, and Khalfaoui 2013). The games at adjacent levels are coupled through the equality constraints on their action profiles given by Eqn (5), and the penalty functions are updated before starting a new iteration. The full algorithm, which we call Separated Hierarchical BRD (SH-BRD), is provided in Algorithm (3).

The penalty updating rule in iteration t of Algorithm (3) is:

1

$$\begin{aligned} \text{For } l &= 2, \dots, L, i = 1, \dots, N^{(l)} \\ & L_i^{(l,l-1)} \left(x_i^{(l)}, \sigma_i^{(l)}(\pmb{x}_{S_i^{(l)}}(t+1)) \right) \\ & = h_i^{(l)} \left[x_i^{(l)} - \sigma_i^{(l)}(\pmb{x}_{S_i^{(l)}}(t+1)) + \lambda_i^{(l)}(t) \right]^2. \end{aligned}$$

2. For $l = 1, ..., L - 1; i = 1, ..., N^{(l)}$, where $a_i^{(l)} \in S_k^{(l+1)}$

$$L_{k}^{(l,l+1)} \left(\sigma_{k}^{(l+1)}(\boldsymbol{x}_{S_{k}^{(l+1)}}), x_{k}^{(l+1)}(t) \right)$$

= $h_{k}^{(l+1)} \left[\sigma_{k}^{(l+1)}(\boldsymbol{x}_{S_{k}^{(l+1)}}) - x_{k}^{(l+1)}(t) - \lambda_{k}^{(l+1)}(t) \right]^{2}.$
(9)

3. For
$$l = 2, \dots, L, i = 1, \dots, N^{(l)}$$

 $\lambda_i^{(l)}(t+1)$
 $= \lambda_i^{(l)}(t) - h_i^{(l)} \bigg[\sigma_i^{(l)}(\boldsymbol{x}_{S_i^{(l)}}(t+1)) - x_i^{(l)}(t+1) \bigg].$
(10)

When updating, all other variables are treated as fixed, and $\lambda^{(l)}(0)$, $h_i^{(l)} > 0$ are chosen arbitrarily.

Algorithm 3: Separated Hierarchical BRD (SH-BRD)
Initialize the game,
$t = 0, x_i^{(1)}(0) = (\boldsymbol{x}_0)_i, i = 1, \dots, N^{(0)}$
for $l = 2:L$ do
for $k = 1: N^{(l)}$ do
$egin{array}{c c c c c c c c c c c c c c c c c c c $
end
end
while not converged do
for $l = 1:L$ do
for $i = I: N^{(t)}$ (l to $l - 1$ Penalty Update, if
l > 1) do
Update $L_i^{(i,i-1)}$
end
for $i = 1: N^{(l)}$ (l to $l + 1$ Penalty Update, if
l < L do $(l l+1)$ $(l) = c(l+1)$
Update $L_k^{(i,i+1)}$, where $a_i^{(i)} \in S_k^{(i+1)}$
end
for $i = 1: N^{(i)}$ (Best Response) do
$x_i^{(l)}(t+1) = BR_i \bigg(\sigma_i^{(l)}(\boldsymbol{x}_{S_i^{(l)}}(t+1)),$
$(l) (\cdot, (l+1) (\cdot, (l)))$
$\boldsymbol{x}_{I_i}^{(\prime)}(t), x_k^{(\prime+1)}(t), u_i^{(\prime)} \right)$
end
end
$t \leftarrow t + 1;$
end

Unlike MS-BRD, the convergence of the SH-BRD algorithm is non-trivial. To prove it, we exploit a connection between this algorithm and Structured Variational Inequalities (SVI) with separable operators (He 2009; He and Yuan 2012; Bnouhachem, Benazza, and Khalfaoui 2013). To formally state the convergence result, we need to make several explicit assumptions.

Assumption 1. The functions $u_i^{(l)}, \forall l = 1, ..., L, \forall i = 1, ..., N^{(l-1)}$ are twice continuously differentiable.

Assumption 2. $-\nabla_{x_i^{(l)}} u_i^{(l)}$ are monotone $\forall l = 1, \ldots, L, \forall i = 1, \ldots, N^{(l-1)}$. The solution set of $\nabla_{x_i^{(l)}} u_i^{(l)} = 0, \forall l = 1, \ldots, L, \forall i = 1, \ldots, N^{(l-1)}$ is nonempty, with solutions in the interior of the action spaces.

Let $F^{(l)}$ be defined as in Equation (3) for each level-l pseudo-utility.

Assumption 3. $F^{(l)}$ satisfy the P_{Υ} condition.

Note that these assumptions directly generalize the conditions required for the convergence of BRD to our multi-scale pseudo-utilities. The following theorem formally states that SH-BRD converges to a NE for *2-level games*.

Theorem 2. Suppose L = 2. If Assumptions 1 and 3 hold, SH-BRD converges to a NE, which is unique.

The full proof of this theorem, which makes use of the connection between SH-BRD and SVI, is provided in the Supplement due to space constraint. The central issue, however, is that there are no established convergence guarantees for ADM-based algorithms for SVI with 3 or more separable operators. Alternative algorithms for SVI can extend to the case of 3 operators using parallel operator updates with regularization terms, but no approaches exist that can handle more than 3 operators (He 2009). We thus propose an algorithm for iteratively solving multi-scale games that uses the general idea from SH-BRD, but packs all levels into two *meta-levels*. The two meta-levels each has to be comprised of consecutive levels. For example, if we have 5 levels, we can have $\{1, 2, 3\}$ and $\{4, 5\}$ combinations, but not $\{1, 2, 4\}$ and $\{3, 5\}$. Upon grouping levels together to obtain a metagame with only two meta-levels, we can apply what amounts to a 2-level version of the SH-BRD. This yields an algorithm, which we call *Hybrid Hierarchical BRD (HH-BRD)*, that now provably converges to a NE for an arbitrary number of levels L given assumptions 1-3.

As presenting the general version of HH-BRD involves cumbersome notation, we illustrate the idea by presenting it for a 4-level game (Algorithm 4). The fully general version is deferred to the Supplement. In this example, the objectives of the meta-levels are defined as

$$\begin{split} \hat{u}_{i}^{(sl_{1})} &= u_{i}^{(1)} + u_{k_{i2}}^{(2)} - L_{k_{i3}}^{(sl_{1},sl_{2})} \bigg(\sigma_{k_{i3}}^{(3)}(\pmb{x}_{S_{k_{i3}}^{(3)}}), x_{k_{i3}}^{(3)} \bigg), \\ \hat{u}_{k_{i3}}^{(sl_{2})} &= u_{k_{i3}}^{(3)} + u_{k_{i4}}^{(4)} - L_{k_{i3}}^{(sl_{2},sl_{1})} \bigg(x_{k_{i3}}^{(3)}, \sigma_{k_{i3}}^{(3)}(\pmb{x}_{S_{k_{i3}}^{(3)}}) \bigg) \,. \end{split}$$

Theorem 3. Suppose Assumptions 1-3 hold Then HH-BRD finds the unique NE.

Proof Sketch. We first "flatten" the game within each metalevel to obtain an effective 2-level game. We then use Theorem 2 to show this 2-level game converges to the unique NE of the game under SH-BRD. Finally, we prove that SH-BRD and HH-BRD have the same trajectory given the same initialization, thus establishing the convergence for HH-BRD. For full proof see Supplement.

HH-BRD combines the advantages of both MS-BRD and SH-BRD: not only does it exploit the sparsity embedded in the network topology, but it also avoids the convergence problem of SH-BRD when the number of levels is higher than three. Indeed, there is a known challenge in the related work on structured variational inequalities that convergence is difficult when we involve three or more operators (He 2009), which we leverage for our convergence results, with operators mapping to levels in our multi-scale game representation. One may be concerned that HH-BRD pseudocode appears to involve greater complexity (and more steps) than SH-BRD. However, this does not imply greater Algorithm 4: Hybrid Hierarchical BRD

Initialize the game, $t = 0, x_i^{(1)}(0) = (\mathbf{x}_0)_i, i = 1, \dots, N^{(0)}$ for l = 2:4 do for $k = 1:N^{(l)}$ do $\boldsymbol{x}_{k}^{(l)}(0) = \sigma_{k}^{(l)}(\boldsymbol{x}_{S_{k}^{(l)}}(0));$ end end while not converged do for $k = 1:N^{(3)}$ (Meta-Level-1 Penalty Update) do Update $L_k^{(sl_1,sl_2)}$ for $i = 1 : N^{(1)}$ (Level-1) do $\begin{aligned} & x_i^{(1)}(t+1) = \\ & BR_i \left(\boldsymbol{x}_{I_i}^{(1)}(t), \boldsymbol{x}_{I_{k_{i2}}}^{(2)}(t), x_{k_{i3}}^{(3)}(t), \hat{u}_i^{(sl_1)} \right) \end{aligned}$ end for $j = 1:N^{(2)}$ (Level-2) do $\boldsymbol{x}_{j}^{(2)}(t+1) = \sigma_{j}^{(2)}(\boldsymbol{x}_{S_{s}^{(2)}}(t+1))$ end for $k = 1:N^{(3)}$ (Meta-Level-2 Penalty Update) do Update $L_k^{(sl_2,sl_1)}$ end for $k = 1 : N^{(3)}$ (Level-3) do $x_k^{(3)}(t+1) = BR_i \bigg(\sigma_k^{(3)}(\boldsymbol{x}_{S_k^{(3)}}(t+1)), \boldsymbol{x}_{I_k}^{(3)}(t),$ $\boldsymbol{x}_{-p}^{(4)}(t), \hat{u}_{k}^{(sl_{2})}$, $(a_{k}^{(3)} \in S_{p}^{(4)})$ end for $p = 1:N^{(4)}$ (Level-4) do $\boldsymbol{x}_{p}^{(4)}(t+1) = \sigma_{p}^{(4)}(\boldsymbol{x}_{S_{r}^{(4)}}(t+1))$ end $t \leftarrow t + 1;$ end

algorithmic complexity, but is rather due to our greater elaboration of the steps within each super level. Indeed, as our experiments below demonstrate, the superior theoretical convergence of HH-BRD also translates into a concrete computational advantage of this algorithm.

Numerical Results and Analysis

In this section, we numerically compare the three algorithms introduced in Section, as well as the conventional BRD. We only consider settings which satisfy Assumptions 1-3; consequently, we focus comparison on computational costs. We use two measures of computational cost: floating-point operations (FLOPs) in the case of games with a linear best response (a typical measure for such settings), and CPU time for the rest. All experiments were performed on a machine with A 6-core 2.60/4.50 GHz CPU with hyperthreaded

Size	BRD	MS-BRD	SH-BRD
30^{2}	$(2.51\pm0.18)\times10^{6}$	$(1.03\pm0.07)\times10^5$	(9.81±0.81)×10 ⁴
50^{2}	$(2.53\pm0.18)\times10^7$	$(5.33 \pm 0.04) \times 10^5$	(4.35±0.07)×10 [‡]
100^{2}	$(4.46\pm0.32)\times10^8$	$(4.36 \pm 0.31) \times 10^6$	(3.56±0.29)×10 ⁶
200^{2}	$(6.73 \pm 0.58) \times 10^9$	$(3.48 \pm 0.29) \times 10^7$	(2.79±0.21)×10 ⁷
500^{2}	$(2.84 \pm 0.21) \times 10^{1}$	$^{1}(5.69\pm0.41)\times10^{8}$	(4.04±0.29)×10 ⁸

Table 1: Convergence and complexity (flops) comparison with linear best response under multiple initialization.

cores, 12MB Cache, and 16GB RAM.

Games with a Linear Best Response (GLBRs) GLBRs (Bramoullé, Kranton, and D'amours 2014; Candogan, Bimpikis, and Ozdaglar 2012; Miura-Ko et al. 2008) feature utility functions such that an agent's best response is a linear function of its neighbors' actions. This includes quadratic utilities of the form

$$u_i(x_i, x_{I_i}) = a_i + b_i x_i + \left(\sum_{j \in I_i} g_{ij} x_j\right) x_i - c_i x_i^2, \quad (11)$$

since an agent's best response is:

$$BR_i(x_{I_i}, u_i) = \frac{\sum_{j \in I_i} g_{ij} x_j}{2c_i} - b_i$$

We consider a 2-level GLBR and compare three algorithms: BRD (baseline), MS-BRD, and HS-BRD (note that in 2-level games, HH-BRD is identical to HS-BRD, and we thus don't include it here). We construct random 2-level games with utility functions based on Equation (11). Specifically, we generalize this utility so that Equation (11) represents only the level-1 portion, $u_i^{(1)}$, and let the level-2 utilities be

$$u_k^{(2)}(x_k, \pmb{x}_{I_k}) = x_k^{(2)} \sum_{p \neq k} v_{kp} x_p^{(2)}$$

for each group k. At every level, the existence of a link between two agents follows the Bernoulli distribution where $P_{exist} = 0.1$. If a link exists, we then generate a parameter for it. The parameters of the utility functions are sampled uniformly in [0, 1] without requiring symmetry. Please refer to Supplement for further details. Results comparing BRD, MS-BRD, and SH-BRD are shown in Table 1. We observe dramatic improvement in the scalability of using MS-BRD compared to conventional BRD. This improvement stems from the representational advantage provided by multi-scale games compared to conventional graphical games (since without the multi-scale representation, we have to use the standard version of BRD for equilibrium computation). We see further improvement going from MS-BRD to SH-BRD which makes algorithmic use of the multi-scale representation.

Games with a Non-Linear Best Response Next, we study the performance of the proposed algorithms in 2- and 3-level games, with the same number of groups in each level (we

Size	BRD	MS-BRD	SH-BRD
30^{2}	$1.50{\pm}0.05$	$1.02{\pm}0.02$	$0.54{\pm}0.01$
50^{2}	26.70±0.36	3.70±0.14	1.81±0.04
100^{2}	1512±9	23.81±0.69	12.10±0.13
200^{2}	> 18000	287.2±5.4	133.6±2.5
500^{2}	nan	5485±13	2524±10

Table 2: CPU times on a single machine on 2-Level games with general best response functions; all times are in seconds.

Size	BRD	MS-BRD	SH-BRD
30^{2}	1.21 ± 0.04	0.63 ± 0.01	0.037±0.003
50^{2}	23.88±0.16	$1.99{\pm}0.04$	0.079±0.004
100^{2}	1461±14	$15.49 {\pm} 0.24$	0.304±0.006
200^{2}	> 18000	192.0±1.2	1.87±0.05
500^{2}	nan	4258±56 s	28.79±0.37

Table 3: CPU times on a single machine for 2-Level, linear/nonlinear best-response games; all times are in seconds.

systematically vary the number of groups). Since SH-BRD and HH-BRD are identical in 2-level games, the latter is only used in 3-level games. All results are averaged over 30 generated sample games. The non-linear best response fits a much broader class of utility functions than the linear best response. The best responses generally don't have closed-form representations. In this case, we can't use linear equations to find the best response and instead have to apply gradient-based methods. In our instances, the utility with non-linear best responses is generated by adding an exponential cost term to the utility function used in GLBRs. Please refer to Supplement for further details.

Table 2 shows the CPU time comparison between all algorithms. The scalability improvements from our proposed algorithms are substantial, with orders of magnitude speedup in some cases (e.g., from ~ 25 minutes for the BRD baseline, down to ~ 12 seconds for SH-BRD for games with 10K agents). Furthermore, BRD fails to solve instances with 250K agents, which can be solved by SH-BRD in ~ 42 min. Again, we separate here the representational advantage of multi-scale games, illustrated by MS-BRD, and algorithmic advantage that comes from SH-BRD. Note that SH-BRD, which takes full advantage of the multi-scale structure, also exhibits significant improvement over MS-BRD, yielding a factor of 2-3 reduction in runtime.

Our next set of experiments involves games in which level-1 utility has a linear best response, but level-2 utility has a non-linear best response. The results are shown in Table 3. We see an even bigger advantage of SH-BRD over the others: it is now typically orders of magnitude faster

Size BRD	MS-BRD	SH-BRD	HH-BRD
10^3 1.23 ± 0.03	$0.59{\pm}0.01$	$0.76{\pm}0.03$	0.43±0.02
20^3 696.0±8.7	3.78±0.09	$6.05{\pm}0.08$	3.35±0.09
$30^3 > 18000$	$15.70{\pm}0.11$	25.13±0.14	13.39±0.11
50^3 nan	$68.59{\pm}0.75$	138.8±1.1	57.98±0.69
100^3 nan	1126±6	2343±21	877.1±11.5

Table 4: CPU times in seconds on a single machine on 3-Level, general best response games; all times are in seconds.

than even MS-BRD, which is itself an order of magnitude faster than BRD. For example, in games with 250K agents, in which BRD fails to return a solution, MS-BRD takes more than 1 hour to find a solution, whereas SH-BRD finds a solution in under 30 seconds.

Finally, Table 4 presents the results of HH-BRD in games with > 2 levels compared to SH-BRD, which does not provably converge in such games. In this case, HH-BRD outperforms the other alternatives, with up to 22% improvement over MS-BRD; indeed, we find that SH-BRD is considerably worse even than MS-BRD.

Conclusions and Future Directions

We proposed a novel representation of games that have a multi-scale network structure. These generalize network games, but with special structures that agent utilities are additive across the levels of hierarchy, with utility at each level depending only on the aggregate strategies of other groups. We present several iterative algorithms that make use of the multi-scale game structure, and show that they converge to a pure strategy Nash equilibrium under similar conditions as for best response dynamics in network games. Our experiments demonstrate that the proposed algorithms can yield orders of magnitude scalability improvement over conventional best response dynamics. Our multi-scale algorithms can reveal to what extent one's group affiliation impacts one's strategic decision making, and how strategic interactions among groups impact strategic interactions among individuals.

While the issue of multi-scale networks abounds in the network science literature (e.g., hierarchical clustering, etc.), the "multi-scale" part is primarily concerned with community structure in networks, rather than modeling how *how communities interact*, which is critical for us in describing a formal multi-scale structure for games. Thus a very important future direction is to identify and obtain relevant field data for experiments, and create realistic benchmarks for multi-scale games. This would involve identifying ways to obtain data about how communities (and not just individuals) interact. Once we have the ability to collect data about interactions at multiple scales (e.g., among members and among groups), we can apply our algorithms to such multi-scale networks. To use criminal networks (criminal organizations and their members) as an example, given game

models constructed with the help of domain expertise, we can:

- 1. compute equilibria predicting, say, criminal activity as a function of structural changes to organizations;
- 2. infer utility models from observational data at multiple scales;
- study policies (including strengthening or weakening connections between agents or groups, endowing agents/groups with more resources (lower costs of effort), etc.) that would induce more desirable equilibrium outcomes.

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