Structured Network Games: Leveraging Relational Information in Equilibrium Analysis

Kun Jin, Parinaz Naghizadeh, Member, IEEE, and Mingyan Liu, Fellow, IEEE

Abstract—We study games with nonlinear best response functions on structured networks. These network structures can emerge from agents' communities or multi-relational interactions, where each relation may follow a different interaction graph. For these structured network games, we establish conditions for uniqueness and stability of pure strategy Nash equilibrium that are stronger yet more computationally efficient to verify than their counterparts in prior research on unstructured (mostly single-relational) network games. Specifically, the network structures enable us to determine Nash equilibrium uniqueness and stability conditions using low-dimensional matrices, often on the order of the number of partitions. This is in contrast to conventional analyses that rely on matrices with dimensions determined by the number of agents multiplied by the action space size. Additionally, we introduce a new degree centrality measure to assess partition influence and use it to establish new Nash equilibrium uniqueness and stability conditions. We compare our findings with prior unstructured network research both analytically and through numerical simulations.

ndex	Terms —network games,	multi-relational	networks,	Nash equilib	orium uniqueness	S.
						

1 Introduction

TETWORK games are commonly used to model the strategic interactions among interdependent agents, where the interdependencies are typically described by a weighted and directed (interaction) network, which also maps into an adjacency matrix. The utility of an agent in these games depends on its own actions as well as the actions of other agents in its local neighborhood as defined by the interaction network. This framework can be used to capture different forms of interdependencies between agents' decisions, such as allowing the action of an agent to be a strategic substitute or complement to that of its neighbors. While these problems have been extensively studied in prior works, see e.g., [1], [2], [3], [4], [5], [6], [7], [8], [9], [10], [11], the focus has been largely on network games with a single, generic underlying network governing the relationship among agents. These will be referred to as *unstructured* network games in this paper; accordingly, their corresponding adjacency matrices will be referred to as regular adjacency matrices. A main focus of this line of work is in identifying sufficient conditions under which the Nash equilibrium (NE) of such a network game is unique and stable, see e.g., [10], [11]. Such conditions are important from a number of perspectives. First, many equilibrium computation algorithms require assumptions that guarantee the uniqueness of the NE. To ensure the convergence or effectiveness of such algorithms, verification of such assumptions is important. Moreover, knowledge about properties like the uniqueness and stability of an NE can help a (mechanism) designer in devising more suitable mechanisms and interventions for the system.

In contrast to prior works, this paper focuses on network games with *structures* that arise naturally. Specifically, we consider two (non-mutually exclusive) families of structures in this study. (1) In the first, the underlying network enjoys certain special graphical properties, a prime example being agents forming *local*

communities, thereby creating sub-graphs more strongly dependent/connected within themselves. (2) In the second case, agents enjoy multi-relational interactions, whereby they are connected over multiple (parallel) networks, each of which governs the dependency relationship of a different action dimension in a high-dimensional action space. Our goal is to understand how the existence of these types of structures affects the resulting NE analysis, and how to exploit such structures, when they exist, to provide characterizations of conditions for the existence, uniqueness, and stability of NEs in such network games. Similar to [1], [7], [8], [10], [11], we consider utility functions with nonlinear best-response functions and use the equivalent variational inequality (VI) representation to study the properties of the NE.

It turns out that such network structures can be used to our advantage in analyzing the NEs of these games, as well as in substantially reducing the computational requirement for verifying the aforementioned conditions. This is done by adopting an *extended* or *generalized* form of adjacency matrix defined over the space of both agents and actions, and identifying *partitions* in this matrix as a result of the types of structures in the network and/or in the agents' dependent relationships described earlier.

Specifically, the computational effort in verifying the uniqueness and stability conditions derived using existing methods, also referred to as *unstructured conditions* in this paper, entails the study of the properties of a game Jacobian (this is the Jacobian matrix of the operator in the VI problem yielding the NE). This matrix is in general asymmetric, making the condition verification a co-NP-complete problem; this means the computational complexity grows faster than polynomial time in the size of the game (total number of agents for an unstructured network game, or total number of agent-action pairs for a structured network game). In the special case of a symmetric game Jacobian, the verification complexity of such conditions still grows in polynomial time in the size of the game.

In contrast, taking advantage of the partition structures in a structured network game, we derive a new set of conditions, also referred to as *structured conditions*, that depend only on the

Kun Jin and Mingyan Liu are with the EECS Department, University of Michigan, Ann Arbor, MI 48109. E-mails: {kunj, mingyan}@umich.edu.

Parinaz Naghizadeh is with the ECE Department, University of California, San Diego, CA 92093. E-mail: parinaz@ucsd.edu.

partitions in the game Jacobian and the size of the partitions. Both structured and unstructured conditions are based on matrices derived based on the game's utility functions and adjacency matrices. When the unstructured derived matrix is asymmetric (resp. symmetric), the structured condition reduces the verification complexity multiple degrees in exponential time (resp. polynomial time). Empirically, as shown in Section 7, the verification of the new condition takes only 2% of the CPU time needed to verify the unstructured conditions from existing literature and avoids memory overflow on large games.

Reducing the verification complexity is of great conceptual and practical interest: it allows us to obtain a high-level understanding of a game much faster in a more sustainable manner especially for large games, and enables early decisions. However, such computational efficiency gain does come at a cost, in that the structured conditions are stronger, i.e., they are sufficient conditions to the unstructured conditions. In other words, some games may satisfy the unstructured conditions (which guarantee the uniqueness and stability of the NE) but fail to satisfy the structured conditions. Our extensive numerical experiments show that this sufficiency gap is small in general, and the sample games that lead to such sufficiency gaps have concentrated features. We characterize such features and provide real-world interpretations.

In addition to the uniqueness and stability of the NE, we introduce a new notion of centrality for structured network games, which can be viewed as a generalization of degree centrality. This new centrality measures the influence or importance of a partition in the network game. We show that this notion can help identify additional conditions for the uniqueness and stability of the NE.

Our main contributions are summarized as follows.

- We provide sufficient conditions for the existence, uniqueness, and stability of the NE in network games, by taking advantage of the partition structures which may arise due to, e.g., the existence of communities or multi-relational interactions.
- We show that these conditions are sufficient conditions to those obtained in previous works, but that they are computationally much easier to verify than their counterparts obtained using conventional methods without utilizing the partition structures.
- The partition structure further allows us to define a new centrality measure, which can be used to verify the uniqueness and stability of the NE in these games.
- We conduct numerical experiments that compare our new conditions with that in previous works in terms of their verification complexity and strengths, which shed light on when using these new conditions is advantageous.

Related works. Conventional (unstructured) network games and their equilibrium outcomes have been studied in a variety of application areas, including the private provision of public goods [1], [2], [3], [12], security decision making in interconnected cyber-physical systems [13], [14], shock propagation in financial markets [15], and (anti)coordination of resources or opinions over networks [16], [17], [18].

A common line of research in this literature studies the effect of the network structure on the existence, uniqueness, and stability of the game equilibria (see [19] for a survey). In particular, equilibrium properties of unstructured network games with nonlinear best-response functions have been studied in [1], [7], [8], [9], [10], [11]. Allouch [1] introduces a sufficient condition for the uniqueness of NE called *network normality* which imposes lower and upper bounds on the derivative of Engel curves. Acemoglu *et al.* [7] consider a network game with idiosyncratic shocks and

show that if the best response mapping is either a contraction with a Lipschitz constant smaller than one or a bounded non-expansive mapping, then the game has a unique NE. Zhou *et al.* [8] establish a connection between nonlinear complementary problems (NCP) and network games, and use results from the NCP literature to find sufficient conditions for the uniqueness of NE.

Our work is closest to [9], [10], [11], which also use the framework of variational inequalities to study network games with nonlinear best-responses (for *unstructured* networks). Naghizadeh and Liu [11] show that a sufficient condition for the uniqueness and stability of the NE can be determined by the lowest eigenvalue of matrices constructed based on the slope of the agents' interaction functions and the intensity of their interactions. Parise and Ozdaglar [10] identify an operator in the variational inequality problem which involves the derivative of the agent's cost with respect to its action, and show that various properties of this operator and its Jacobian will determine conditions for the existence, uniqueness, and stability of a NE.

Our prior related work [20] shows that community structures can be utilized to derive sufficient conditions on NE uniqueness and stability. This paper extends [20] by introducing more general structured network games that further allow for multi-relational dependencies between agents, and obtains sufficient conditions for existence, uniqueness, and stability of NE in these network games. The conditions in [20] can be recovered as special cases of the structured conditions in Sections 4, 5, and 6, when we limit our attention to a single-relational, multi-partite network and use each partite as a partition.

The computational complexity of checking the uniqueness of NE has also been explored in a number of prior works; notably, [21], [22] show that verifying the uniqueness of NE is in general an NP-hard problem. More recently, the complexity of checking the existence of a NE in (unstructured) games on networks has been explored in [23], [24]. Our work contributes to this literature by providing such results for structured network games. A computationally efficient method for verifying the existence and uniqueness of NE can further serve as a precursor to algorithms for (efficient) computation of NE, or before implementing iterative or distributed algorithms which require uniqueness of the NE for convergence to it (e.g., [25], [26]).

2 MOTIVATING EXAMPLES

We elaborate on the idea of structured networks through a running example consisting of a number of vendors/store owners (strategic agents) selling similar merchandise. Please refer to Appendix B for values used in Figure 1-4.

Communities due to stronger connectivity. Community formation stems from situations where the strengths of agents' strategic interactions are (statistically) different within certain groups compared to those between these groups. One type of community structure, commonly studied and discovered using spectral analysis [27], [28], is characterized by groups with much stronger connectivity (higher density of connections or existence of edges, as well as higher edge weights on those edges) within themselves, and much weaker connectivity (lower density of edges and smaller edge weights) between them.

Consider the case of three store owners, with agents a_1 and a_2 in close proximity of each other and a_3 located far away. Further, consider the store owners' single action of selecting business hours. Since a_1 and a_2 offer similar goods, their individual

decisions on business hours (from complete overlap to mutually exclusive) will have direct consequences on the other's business volume and goods sold, resulting in a stronger dependence relationship between the two. Their dependence on (or influence of) a_3 may be far weaker. This is illustrated in the left of Figure 1, where the stronger relationship between a_1 and a_2 is indicated by a thicker edge; here a_1 and a_2 form a group or community.

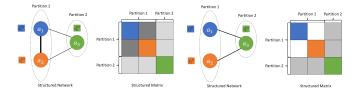


Fig. 1. Network with communities/groups: 3 agents and 1 action dimension; a_1 and a_2 form the first group, a_3 is a singleton group.

Fig. 2. Bipartite graph, where a_1 and a_2 are on one side, and exhibit the same type of dependence on a_3 on the other side.

The recognition of groups gives rise to a "block" view in the structured adjacency matrix shown in the right-hand side of Figure 1. Given there is only a single action, here every row and column is associated with an agent, resulting in a 3×3 interaction matrix. The colors of the diagonal entries match the identity of the agents depicted in the network as well as their indices. The diagonal elements in the regular interaction matrix represent an agent's self influence, whereas the off-diagonal elements represent the agents' mutual influence. Similarly, the diagonal blocks in the structured matrix represent a partition's self influence, whereas the off-diagonal blocks represent the partitions' mutual influence. These quantities will be precisely defined in the next section.

Communities due to similarity in function. Communities can also result from a logical relationship. As an example, consider again the three store owners, where a_1 and a_2 carry completely orthogonal merchandise (e.g., a bakery vs. a hardware store) but both rely on a_3 to provide store security and vehicle rental as needed, and consider the single action of staffing levels. In this case, the dependency only exists between a_1, a_3 , and between a_2, a_3 , but not between a_1, a_2 , resulting in a bipartite network shown in Figure 2. Yet in this case it is still appropriate to view a_1 and a_2 as belonging to the same group, because they each exhibit very similar dependence on another group.

Partition on action dimensions. Our next example is more complex and introduces a high-dimensional action space. The general idea is that when actions are high-dimensional, the interaction/dependency relationships among agents can be different for different action dimensions, effectively resulting in multiple parallel networks superimposed on each other; this will also be referred to as a *multi-relational* network game.

Consider again the three store owners, each of which now has both physical and online sales. It is reasonable to expect that the decisions a seller makes about what goods to display in the window or offer for sampling may have more impact on other similar stores in its physical proximity (one set of agents), whereas its decisions on webpage design, layout, picture quality, and payment options may only impact its online competitors (another set of agents). In such multi-relational games, the action dimensions naturally create action-based structures on the network.

Suppose we capture the above with two action dimensions (instore and online decisions), and further suppose in our case a_1 and a_2 are much more in direct competition in terms of physical-

store sales due to their proximity, but that a_2 and a_3 are much more in direct competition in terms of their online sales due to high similarity in the goods they carry. This means that agents form different groups in different actions, as illustrated in Figure 3, where agents are represented by different colors and action dimensions are represented by different shades. In the first (solid color) action, a_1 and a_2 form a group whereas in the second (faded color) action, a_2 and a_3 form a group.

To capture the multiple action dimensions, we will define an *extended* adjacency/interaction matrix where each row and each column represent an agent-action dimension pair; the regular adjacency matrix is then a special case of this extended definition when the action dimension is 1. Under this definition, we have a 6×6 interaction matrix for this example, shown in Figure 4. Given such a matrix, partitions can emerge either by agents (similar sets of agents form the same group regardless of the action dimension) or by actions (interactions along different dimensions tend to be orthogonal), as is the case in this example and shown in Figure 4.

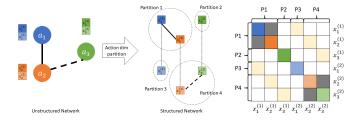


Fig. 3. Action dimension partitioned graph, with 3 agents and 2 actions. a_1 and a_2 are closely connected on dimension 1. a_2 and a_3 are closely connected on dimension 2.

Fig. 4. Action dimension partitioned matrix for the network in Figure 3.

Arbitrary partitions in the extended adjacency matrix. We note that the idea of a partitioned extended adjacency matrix is not restricted to the above example scenarios. In principle, this can be done across both agents and action dimensions in arbitrary ways. As such, our approach, which is based on a partitioned matrix, is applicable whether the structure arises due to special graphical properties or multi-relational interactions, or both.

We also note that in our analysis, we will be working with the Jacobian (∇F) of a best-response operator (F), defined precisely in subsequent sections, rather than the extended adjacency matrix (G) itself. This is because the former captures both first-order and second-order information about the utility functions, which is not captured in the latter. In particular, the extended adjacency matrix does not reveal cross-action dependencies induced by the utility functions. We will, however, also show that there is a direct correspondence between a partition on the extended adjacency matrix and that on the Jacobian. For this reason, the use of the adjacency matrix in this section is for illustration purposes only, as it is much more intuitive and straightforward to visualize the described partition structures in an adjacency matrix.

3 Model and Preliminaries

3.1 The Structured Network Game Model

We consider a structured network game among N agents $\mathcal{N} = \{a_1, \ldots, a_N\}$, each with K action dimensions.¹ The multi-

1. We set the same K for all agents for simplicity of presentation. Having heterogeneous action dimensions across agents will increase the number of agent-action components; we can apply the same analysis to this larger extended adjacency matrix, with the only difference being that this can increase the verification complexity of the uniqueness and stability conditions.

relational network is represented by a multi-relational graph with extended adjacency matrix G. The adjacency matrix on the k-th action dimension is denoted as $G^{(k)}$, and is a submatrix of G. The edge weight $G_{ij}^{(k)} \in \mathbb{R}$ is a real number representing the strength of influence agent a_j has on agent a_i (or a_i 's dependence on a_j) in the k-th action dimension.

We use $x_i^{(k)} \in \mathbb{R}$ to denote the k-th action of agent a_i , $\boldsymbol{x}_i = (x_i^{(k)})_{k=1}^K \in \mathbb{R}^K$ to denote the action vector of a_i , and $\boldsymbol{x}_i^{(-k)} = [x_i^{(1)}, \cdots, x_i^{(k-1)}, x_i^{(k+1)}, \cdots, x_i^{(K)}]^T$ to denote the action profile of a_i , excluding the k-th dimension. In addition, let $\boldsymbol{x}^{(k)} = (x_i^{(k)})_{i=1}^N \in \mathbb{R}^N$ be the action vector of all agents on the k-th action dimension, and \boldsymbol{x}_{-i} denotes the action profile of all agents other than a_i .

Each agent a_i has an action constraint $\boldsymbol{x}_i \in Q_i = \prod_{k=1}^K Q_i^{(k)}$, where $Q_i^{(k)} := [0, B_i^{(k)}]$ such that $B_i^{(k)}$ captures the physical or financial constraints (budgets) on the k-th action.

We consider games with utility functions consisting of an individual component and a network component:

$$u_i(\boldsymbol{x}_i, \boldsymbol{x}_{-i}, G) = d_i(\boldsymbol{x}_i) + f_i(\boldsymbol{x}_i, G^{(1)} \boldsymbol{x}^{(1)}, \dots, G^{(K)} \boldsymbol{x}^{(K)}).$$
 (1)

Here, $d_i(\cdot)$ is the individual component, which only depends on a_i 's own actions; conventionally, it contains a standalone benefit and cost of taking action \boldsymbol{x}_i . The network component $f_i(\cdot)$ depends on not only the agent's own but also others' actions. Here, the network influence on the k-th action dimension is captured by $G^{(k)}\boldsymbol{x}^{(k)}$. Throughout the paper, we make the following assumptions.

Assumption 1. The utility functions $u_i(\mathbf{x}_i, \mathbf{x}_{-i}, G)$ are concave and twice continuously differentiable, for all i.

An example of this type of utility function is given below.

Example 1. The multi-relational extension of linear-quadratic utility functions, studied in e.g., [29], has the following form,

$$u_i(\boldsymbol{x}_i, \boldsymbol{x}_{-i}, G) = \boldsymbol{x}_i^T \boldsymbol{b}_i + \sum_{k=1}^K \sum_{j \neq i} x_i^{(k)} g_{ij}^{(k)} x_j^{(k)} - \frac{1}{2} \boldsymbol{x}_i^T C_i \boldsymbol{x}_i,$$

where $x_i^T b_i - \frac{1}{2} x_i^T C_i x_i$ is the individual component containing the standalone benefit term and the cost term, and $\sum_{k=1}^K \sum_{j \neq i} x_i^{(k)} g_{ij}^{(k)} x_j^{(k)}$ is the (multi-relational) network influence component. (We can further expand this utility function with cross-relational influence terms $\sum_{k=1}^K \sum_{j \neq i} x_i^{(k)} g_{ij}^{(k,l)} x_j^{(l)}$, where $G^{(k,l)}$ is a submatrix in G modeling the network influence from the l-th action to the k-th action.)

A Nash equilibrium of the described structured network game can be found as a fixed point of the agents' best response mappings. The best response of an agent in the network game is defined as the action an agent takes to maximize its own utility, given other agents' actions and the network topology. For our model, we denote the best response of agent a_i as

$$BR_i(\boldsymbol{x}_{-i}, G) := \underset{\boldsymbol{x}_i \in Q_i}{\operatorname{max}} \ u_i(\boldsymbol{x}_i, \boldsymbol{x}_{-i}, G).$$

We also define an operator F_i as follows,

$$F_i(\boldsymbol{x}_i, \boldsymbol{x}_{-i}) = -\nabla_{\boldsymbol{x}_i} u_i(\boldsymbol{x}_i, \boldsymbol{x}_{-i}, G) \in \mathbb{R}^K. \tag{2}$$

Next, we introduce the Variational Inequality (VI) framework and its relation to Nash equilibria in network games.

3.2 The Variational Inequality (VI) Problem

Variational Inequalities (VIs) are a class of optimization problems with applications in game theory. In particular, the Nash equilibria of many games can be found as solutions to a corresponding VI problem [10], [30]. We state the VI problem formally below, using the set of notations introduced earlier so that the correspondence between the game model and the VI problem is clear.

Definition 1. A variational inequality VI(Q, F) consists of a set $Q \subseteq \mathbb{R}^N$ and a mapping $F: Q \to \mathbb{R}^N$, and is the problem of finding a vector $\mathbf{x}^* \in Q$ such that,

$$(\boldsymbol{x} - \boldsymbol{x}^*)^T F(\boldsymbol{x}^*) \ge 0, \ \forall \boldsymbol{x} \in Q. \tag{3}$$

Finding the Nash equilibrium of a structured network game is equivalent to solving a variational inequality problem $VI(Q,F(\boldsymbol{x}))$ with the appropriate choice of Q and F. An example choice of Q and F is: $Q=Q_1\times Q_2\times\ldots\times Q_N\subseteq\mathbb{R}^{NK}$, with Q_i being the action space of a_i , and $F(\boldsymbol{x})=(F_i(\boldsymbol{x}_i,\boldsymbol{x}_{-i}))_{i=1}^N\in\mathbb{R}^{NK}$, with $F_i(\cdot)$ given in (2). Then, since finding an NE is the problem of finding a fixed point of the best response mappings in (3.1), it is equivalent to solving the VI problem given in (3) with the choice of F and Q stated above (see, e.g., [31, Proposition 1.4.2]). Intuitively, if the equilibrium \boldsymbol{x}_i^* is an interior point of Q_i , then $F_i(\boldsymbol{x}_i^*, \boldsymbol{x}_{-i}^*) = \mathbf{0}$ if and only if \boldsymbol{x}_i^* is a best response to \boldsymbol{x}_{-i}^* ; if \boldsymbol{x}_i^* is on the boundary, we can still show $(\boldsymbol{x}_i - \boldsymbol{x}_i^*)^T F_i(\boldsymbol{x}^*) \geq 0$ holds. Note also that the above definition holds for Q and F that are permutations of the preceding example choice, as long as the permutations are consistent.

3.3 Partitions

We next formally define partitions in a structured network game. As shown in Section 2, a partition is essentially a set of indices, where each index corresponds to an agent-action pair (which also corresponds to a column or row in the extended adjacency/interaction matrix G). Figures 1, 2, and 4 show how, by grouping certain indices together (rearranging rows and columns), block structures may emerge in G.

It turns out a block structure in G translates to a similar block structure in the Jacobian of the operator F, when F is a suitable permutation of the operator $(F_i(\boldsymbol{x}_i, \boldsymbol{x}_{-i}))_{i=1}^N$ consistent with the partitions. In Eqn (2), we can think of F_i as the best response direction vector of a_i , where each element in F_i corresponds to an agent-action component. Therefore, operator $F(\boldsymbol{x})$ is the global best response direction vector containing all agent-action components; these components can be arranged in an arbitrary order, and $(F_i(\boldsymbol{x}_i, \boldsymbol{x}_{-i}))_{i=1}^N$ corresponds to a common order.

While F captures important first-order derivative information of the utility functions with respect to the agent-action components, the Jacobian of F, denoted as ∇F captures important second-order information of the utility functions and first-order information of the operator F. It is common, see e.g., [10], [11], [30], to use the properties of ∇F to derive unstructured conditions for equilibrium analysis.

When K=1, $\nabla F\in\mathbb{R}^{N\times N}$, where the off-diagonal elements in ∇F measure how an agent's action influences another agent's best response. For K>1, $\nabla F\in\mathbb{R}^{NK\times NK}$, and now the off-diagonal elements in ∇F measure how an agent-action component influences the best response direction of another agent-action component.

There is a one-to-one mapping between the dimensions in ∇F and G, since each dimension in ∇F and G corresponds to an agent-action component, and thus any partition on G can equally apply to ∇F , as partitions are nothing more than separating the agent-action components into disjoint sets. Using the same 3-agent, 2-action example in Figures 3 and 4, while $(F_i(\boldsymbol{x}_i, \boldsymbol{x}_{-i}))_{i=1}^N$ orders the agent-action components as $x_1^{(1)}, x_1^{(2)}, x_2^{(1)}, x_2^{(2)}, x_3^{(1)}, x_3^{(2)}, F$ orders them as $x_1^{(1)}, x_2^{(1)}, x_3^{(1)}, x_1^{(2)}, x_2^{(2)}, x_3^{(3)}$. Then, if we partition the network as shown in Figure 3, we can partition ∇F the same way, as illustrated in Figure 4.

We note that the partition based on (agent) group structure and the partition based on action dimensions are special cases. For the remainder of the paper, we will discuss network games with general, arbitrary partition structures. Specifically, for an arbitrary structure, we can partition all the NK agent-action components (and their corresponding indices) into an arbitrary number (M) of disjoint sets, denoted by $\mathcal{P}_1, \ldots, \mathcal{P}_M$. Accordingly, we will also denote by $N_i = |\mathcal{P}_i|$ the size of the partitions.

4 EXISTENCE AND UNIQUENESS

4.1 Existence of NE

We first state the conditions under which a NE exists. The following theorem from [30] guarantees the existence of NE:

Theorem 4.1. ([30, Theorem 3]) If F is continuous on Q, and Q is nonempty, compact and convex, then VI(Q, F) has a nonempty and compact solution set.

This is because in our problem, $Q=Q_1\times Q_2\times\ldots\times Q_N$, with $Q_i=[0,B_i^{(1)}]\times\ldots\times[0,B_i^{(K)}]$, so that together with Assumption 1, the conditions of Theorem 4.1 are satisfied for the network game defined in Section 3.1.

4.2 Uniqueness of NE

We next introduce sufficient conditions under which the network game defined in Section 3.1 has a unique NE. We begin by introducing the following definitions.

Definition 2. *P-Matrix:* A matrix $A \in \mathbb{R}^{N \times N}$ is a *P-matrix if* every principal minor of A has a positive determinant.

Definition 3. A mapping $F: Q \mapsto \mathbb{R}^{NK}$, where $Q \subseteq \mathbb{R}^{NK}$ is nonempty, compact and convex, and F is continuously differentiable on Q, is strongly monotone if there exists a c > 0 s.t.

$$(\boldsymbol{x} - \boldsymbol{y})^T (F(\boldsymbol{x}) - F(\boldsymbol{y})) \ge c \|\boldsymbol{x} - \boldsymbol{y}\|_2^2, \ \forall \boldsymbol{x}, \boldsymbol{y} \in Q.$$

Further, $F = (F_1, F_2, \dots, F_N)$ is a uniform block P-function w.r.t. the (agent-level) partition $Q = Q_1 \times Q_2 \times \dots \times Q_N$ if there exists a constant b > 0 such that

$$\max_{i \in \mathbb{N}[1,M]} (\boldsymbol{x}_i - \boldsymbol{y}_i)^T [F_i(\boldsymbol{x}) - F_i(\boldsymbol{y})] \ge b \|\boldsymbol{x} - \boldsymbol{y}\|_2^2, \ \forall \boldsymbol{x}, \boldsymbol{y} \in Q.$$

By setting b=c/N, it is easy to see that strong monotonicity is a sufficient condition for the uniform block P-condition. Parise and Ozdaglar [10] show that if $F(\boldsymbol{x})$ is a uniform block P-function, then VI(Q,F) has a unique solution, and the NE of the network game corresponding to VI(Q,F) is unique. Unfortunately, it is computationally costly to verify these conditions for the function $F(\boldsymbol{x})$, since checking whether a square matrix is a P-matrix is co-NP-complete [32], and typically, the complexity

grows (faster than polynomial time) in the size of the matrix. Below, we show that it is possible to take advantage of the block structure when it is present, and identify conditions for the uniqueness of NE that are of lower computational complexity to verify. To do so, we define a structured matrix Υ^S and its components, the *internal* and *external* impact levels of partitions.

Definition 4. We say a partition \mathcal{P}_i receives internal impact, α_i^S , and external impact, β_{ij}^S , defined as follows:

$$\alpha_i^S = \inf_{\boldsymbol{x} \in Q} \|\nabla_i F_i(\boldsymbol{x})\|_2, \ \forall i \in \mathbb{N}[1, M]$$

$$\beta_{ij}^S = \sup_{\boldsymbol{x} \in Q} \|\nabla_j F_i(\boldsymbol{x})\|_2, \ \forall i, j \in \mathbb{N}[1, M], \ i \neq j,$$

where $\nabla_j F_i(\boldsymbol{x}) \in \mathbb{R}^{N_i \times N_j}$ is a matrix with k, l-th entry $\frac{\partial F_i^{(k)}(\boldsymbol{x})}{\partial x_j^{(l)}}$. The structured matrix Υ^S is defined accordingly as:

$$\Upsilon^{S} = \begin{bmatrix}
\alpha_{1}^{S} & -\beta_{12}^{S} & \dots & -\beta_{1M}^{S} \\
-\beta_{21}^{S} & \alpha_{2}^{S} & \dots & -\beta_{2M}^{S} \\
\vdots & \vdots & \ddots & \vdots \\
-\beta_{M1}^{S} & -\beta_{M2}^{S} & \dots & \alpha_{M}^{S}
\end{bmatrix} .$$
(4)

We note that these impact measures are only defined between partitions and not agents. Accordingly, in this definition the subscripts are indices of partitions instead of agents.

To motivate the above definition, it helps to understand the matrix Υ^S in the context of what is typically used in prior works for checking the uniqueness of the NE. If we ignore the structure in the network and simply view each agent-action pair as a singleton partition, then using an existing methodology (such as in [10], [11], [30]) will give us the Jacobian Υ^U , which is a $NK \times NK$ matrix. Specifically, Υ^U is given by

$$\Upsilon^{U} = \begin{bmatrix}
\alpha_{1}^{U} & -\beta_{12}^{U} & \dots & -\beta_{1,NK}^{U} \\
-\beta_{21}^{U} & \alpha_{2}^{U} & \dots & -\beta_{2,NK}^{U} \\
\vdots & \vdots & \ddots & \vdots \\
-\beta_{NK,1}^{U} & -\beta_{NK,2}^{U} & \dots & \alpha_{NK,NK}^{U}
\end{bmatrix}, (5)$$

and contains the following elements:

$$\begin{array}{lcl} \alpha_k^U & = & \inf_{\boldsymbol{x} \in Q} |\nabla_k F_k(\boldsymbol{x})|, \; \forall k \in \mathbb{N}[1, NK] \\ \beta_{kl}^U & = & \sup_{\boldsymbol{x} \in Q} |\nabla_l F_k(\boldsymbol{x})|, \; \forall k, l \in \mathbb{N}[1, NK], \; k \neq l. \end{array}$$

To clarify, these will be referred to as the *component-level* internal and external impact, respectively. Suppose we rearrange the rows and columns of Υ^U in the following way: group together rows whose action dimensions are in \mathcal{P}_i , $i=1,\cdots,M$; group together columns whose agents are in \mathcal{P}_j , $j=1,\cdots,M$.

We can now view this rearranged Υ^U in blocks/submatrices denoted by $\Upsilon^U_{\mathcal{P}_i,\mathcal{P}_j}$, and there are $M\times M$ blocks. The matrix Υ^S is essentially a condensed version of this rearranged matrix, summarizing or abstracting each block into a single quantity as defined in Definition 4: α^S_i for the diagonal block $\Upsilon^U_{\mathcal{P}_i,\mathcal{P}_i}$ and β^S_{ij} for the off-diagonal block $\Upsilon^U_{\mathcal{P}_i,\mathcal{P}_i}$.

This abstraction aims at capturing the dependence relationship between partitions rather than between individual agent-action pairs. In particular, β^S_{ij} represents the largest influence level of partition \mathcal{P}_j on partition \mathcal{P}_i , and α^S_i represents the minimum influence level of \mathcal{P}_i on itself. Formally:

Lemma 4.2. We have
$$\|\Upsilon^U_{\mathcal{P}_i,\mathcal{P}_j}\|_2 \geq \beta^S_{ij}$$
 and $\|\Upsilon^U_{\mathcal{P}_i,\mathcal{P}_i}\|_2 \leq \alpha^S_i$.

In what follows, we show that the matrix Υ^S can be used to provide sufficient conditions on the uniqueness of NE in network games with partition structures. Such an abstraction takes advantage of the partition structure and reduces the dimension of the matrix used to check conditions for the uniqueness from $NK \times NK$ to a single $M \times M$ matrix and M matrices of size $|\mathcal{P}_k| \times |\mathcal{P}_k|, k \in [1, M]$. This greatly reduces the complexity of condition verification, since as mentioned earlier, P-matrix verification is co-NP-complete [32], and the complexity typically grows (faster than polynomial) in the size of the matrix. For example, in a special case where the matrix is symmetric, P-matrix verification is equivalent to examining its positive definiteness and the complexity becomes $O(N^3)$ for an $N \times N$ matrix. If there are N=20 agents on the network with K=5 action dimensions, then the verification complexity of unstructured conditions is $O(10^6)$; if we partition the game by action dimensions, the complexity of the structured conditions is $5 \cdot 8 \cdot O(10^3) + O(5^3)$, much lower than $O(10^6)$. This complexity gap will only increase if the matrix is asymmetric.

However, we also note that the strengths of the conditions obtained from the structured network and the unstructured network are not equivalent. Specifically, we will later show that the conditions obtained from the structured network are stronger (sufficient conditions to) their counterparts in the unstructured network. Numerical results presented in Section 7 also highlight the gap between these two sets of conditions.

The following result identifies a condition for the uniqueness of the NE in games with the structured network.

Theorem 4.3. If the following two conditions are satisfied,

- 1) Υ^S is a P-matrix, and
- 2) $\Upsilon^{U}_{\mathcal{P}_{i},\mathcal{P}_{i}}$, $\forall i$ (the diagonal blocks of Υ^{U}) are P-matrices,

then the network game has a unique NE.

Moreover, when both conditions are satisfied, then Υ^U is also a P-matrix; i.e., the uniqueness conditions (1 & 2) on the structured network are also a sufficient condition for the uniqueness of the NE in the underlying unstructured network (i.e., one that ignores the partitioned structures).

An interpretation of the above result is as follows. As noted earlier, β_{ij}^S represents the external impact of partition \mathcal{P}_j on partition \mathcal{P}_i , while α_i^S is the internal impact of \mathcal{P}_i . Typically, when β_{ij}^S has a relatively small value compared to α_i^S , then Υ^S is a P-matrix. Moreover, when Υ^S is (row or column) diagonally dominant, then Υ^S is a P-matrix [33]. In these types of networks, partitions' action profiles have a bounded influence on each other. Conversely, if at least one partition's action profile has an out-sized effect on other partitions, then its decision can shift the state of the network substantially and result in possibly multiple equilibria.

Remark 1. It is worth mentioning that the structured (resp. unstructured) network condition verification is of a much lower complexity if Υ^S (resp. Υ^U) is symmetric. By [34], a symmetric matrix is a P-matrix if and only if it is positive definite, which means that instead of checking the determinant for every principal minor, we only need to do an eigendecomposition. This reduces the complexity from solving a co-NP-complete problem down to polynomial time.

We next present two corollaries of Theorem 4.3 which provide alternative ways for verifying the uniqueness of NE in structured network games. We define

$$\Gamma^{S} = \begin{bmatrix} 0 & -\beta_{12}^{S}/\alpha_{1}^{S} & \dots & -\beta_{1M}^{S}/\alpha_{1}^{S} \\ -\beta_{21}^{S}/\alpha_{2}^{S} & 0 & \dots & -\beta_{2M}^{S}/\alpha_{2}^{S} \\ \vdots & \vdots & \ddots & \vdots \\ -\beta_{M1}^{S}/\alpha_{M}^{S} & -\beta_{M2}^{S}/\alpha_{M}^{S} & \dots & 0 \end{bmatrix}, (6)$$

where β_{ij}^S and α_i^S are as in Definition 4. Similar to before, by treating each agent as a singleton partition, we can obtain Γ^S 's unstructured counterpart, $\Gamma^U \in \mathbb{R}^{NK \times NK}$. We also denote the spectral radius of Γ^S as $\rho(\Gamma^S)$. By [35], if $\rho(\Gamma^S) < 1$, then Υ^S is a P-matrix. Therefore, we have the following Corollary of

Corollary 4.3.1. Assume both of the following conditions hold

- 1) $\rho(\Gamma^S) < 1$, 2) $\Upsilon^U_{\mathcal{P}_i,\mathcal{P}_i}, \forall i \ are \ P\text{-matrices}.$

Then, the network game has a unique NE and Υ^U is a P-matrix.

Intuitively, the conditions of Corollary 4.3.1 guarantees Υ^S is a P-matrix, and we can use Theorem 4.3 to show Υ^U is a P-matrix.

By [34], we know that a symmetric matrix is a P-matrix if and only if it is positive definite. Therefore, we have the following:

Corollary 4.3.2. Assume Γ^S is a symmetric matrix. Denote the eigenvalues of Γ^S by $\lambda_1(\Gamma^S) \leq \lambda_2(\Gamma^S) \leq \cdots \leq \lambda_M(\Gamma^S)$. Then, the network game has a unique NE if the following two conditions hold simultaneously,

- 1) $\lambda_1(\Gamma^S) > -1$, 2) $\Upsilon^U_{\mathcal{P}_i,\mathcal{P}_i}, \forall i \ are \ P\text{-matrices}.$

Moreover, under these conditions, Υ^U is a P-matrix.

These corollaries provide two alternative ways to check for the uniqueness of the NE. In terms of complexity, both finding the spectral radius and the eigendecomposition are of complexity $O(N^3)$ for an $N \times N$ matrix; these corollaries' conditions are therefore computationally easier to verify than the co-NPcomplete problem. However, the trade-off is that the conditions in Corollary 4.3.1 are stronger than those of Theorem 4.3, while Corollary 4.3.2 can only be used given a symmetric Γ^S (resp. Γ^{U}) on structured (resp. unstructured) networks. It is also worth mentioning that even when Υ^S (resp. Υ^U) are asymmetric, we can still have symmetric Γ^S (resp. Γ^U) matrices, and thus, Corollary 4.3.2 could still provide a computationally lighter alternative verification in these cases.

Lastly, we again note that $\Gamma^U \in \mathbb{R}^{NK \times NK}$ could be formed by treating each agent as a singleton partition. When we take this viewpoint, Corollary 4.3.2 reduces to Proposition 3 of [11]. On the other hand, by using the partition structure, Γ^S is an $M \times M$ matrix, and the conditions in Corollary 4.3.2 are computationally easier to verify as compared to those in [11]. Specifically, checking the eigenvalues of a matrix requires performing eigendecomposition over it. To elaborate on the comparison, suppose $N_i = \overline{N}$ for all \mathcal{P}_i (all partitions have the same size). Then, using the unstructured network, the complexity of the eigendecomposition on Υ^U or Γ^U is $O(K^3N^3)$ while using the partition structure, the complexity on Υ^S or Γ^S is reduced to $O(M^3)$. Of course, using the partition structure, we have to compute the α_i^S and β_{ij}^S values as well, which is of complexity $O(M^2\overline{N}^3)$. Altogether, the complexity of checking whether Υ^S is a P-matrix under the conditions in the above corollary is $O(M^3 + M^2 \overline{N}^3)$, which can be much lower than $O(K^3N^3)$ in the unstructured case.

4.3 Sufficiency Gaps on the Uniqueness Conditions

To close this section, we elaborate on the difference between using and not using information about partition structures in checking for the uniqueness of Nash equilibria.

The first thing to note is that, as mentioned earlier, the set of sufficient conditions derived when accounting for partition structures are generally stronger than their counterparts derived without using information about the structure: as Theorem 4.3, Corollary 4.3.1, and Corollary 4.3.2 indicate, if a structured network satisfies these uniqueness conditions, then it also satisfies the corresponding unstructured uniqueness conditions, but the opposite is in general not true. This means that there is a sufficiency gap between the conditions obtained from the structured and unstructured networks. The most important reason behind this sufficiency gap has to do with the way partition structures are abstracted. There are two forms of abstractions made during the creation of the Υ^S matrix; both are for partitions and each summarizes the component (agent-action) level internal and external impact, respectively:

- 1) The internal impact α_i^S of partition \mathcal{P}_i , is an abstraction of the component-level internal impact α_k^U for components (and corresponding indices) in \mathcal{P}_i , and component level external impact β_{kl}^U between different indices in \mathcal{P}_i ;
- 2) The external impact β_{ij}^S of \mathcal{P}_j on \mathcal{P}_i , is an abstraction of component-level external impact of indices in \mathcal{P}_j to indices in \mathcal{P}_i .

These types of abstractions inevitably introduce gaps in the sufficiency conditions. As mentioned in Lemma 4.2, the value of α_i^S is lower-bounded by $\|\Upsilon_{\mathcal{P}_i,\mathcal{P}_i}^U\|$ and highly depends on the agents with component-level internal impact in \mathcal{P}_i . Meanwhile, the β_{ij}^S value is upper bounded by $\|\Upsilon_{\mathcal{P}_i,\mathcal{P}_j}^U\|$, and highly depends on the strongest component-level external impact from one index in \mathcal{P}_j to another in \mathcal{P}_i . The significance of α_i^S in verifying conditions for the uniqueness of the Nash equilibria is akin to the observation "a chain is as strong as its weakest link"; we refer to this as the "weakest link effect". Similarly, the significance of the β_{ij}^S values is referred to as the "strongest link effect".

Recall an earlier observation in Theorem 4.3 that when each agent has stronger component-level internal impact than component-level external impact, the NE is unique. Similarly, in terms of structure, when each partition has stronger internal impact than external impact, the NE is unique and the conditions in Theorem 4.3 and Corollaries 4.3.1, 4.3.2 are sufficient to guarantee such impact differentials.

In some games, the indices k with weak component-level internal impact α_k^U also have weak component-level external impact β_{kl}^U while the indices l with strong component-level internal impact α_l^U have similarly strong component-level external impact β_{lk}^U . While we may be able to guarantee the uniqueness of an NE using the unstructured network, we may not be able to do so using structures when a partition contains both types of (strong and weak) indices. This is because the abstraction becomes inaccurate as the partition's internal impact α_i^S is weak and the external impact β_{ij}^S strong; thus the conditions obtained from the structured network may fail to guarantee the uniqueness of NE. The following example highlights these observations.

Example 2. Consider a 4-agent, single-action dimension, 2-partition game where agents a_1, a_2 form \mathcal{P}_1 and agents a_3, a_4 form \mathcal{P}_2 , with utility functions: $u_1(\mathbf{x}) = x_1(x_2 + 5x_3) - x_1(x_2 + 5x_3)$

$$5x_1^2, u_2(\mathbf{x}) = x_2(x_1 + x_4) - 2x_2^2, u_3(\mathbf{x}) = x_3(x_4 + 5x_1) - 5x_3^2, u_4(\mathbf{x}) = x_4(x_3 + x_2) - 2x_4^2$$
. Then, we have

$$\Upsilon^U = \begin{bmatrix} 10 & -1 & -5 & 0 \\ -1 & 4 & 0 & -1 \\ -5 & 0 & 10 & -1 \\ 0 & -1 & -1 & 4 \end{bmatrix} \succ \mathbf{0}, \ \Upsilon^S = \begin{bmatrix} 3.84 & -5 \\ -5 & 3.84 \end{bmatrix} \prec \mathbf{0},$$

which means that the unstructured condition guarantees the uniqueness of the NE in this game, yet the game fails to satisfy the sufficient structured condition. This is an example when the internal impact of a partition is weak but the external impact between partitions is strong.

We end this section by summarizing the types of games and partition structures that result in small sufficiency gaps between the two sets of conditions. The sufficiency gap is small if the partition members have similar component-level internal impacts; or if the members have similar connections to members in other partitions and have similar component-level external impacts; or if the member with the weakest component-level internal impact also has the strongest component-level external impact. We further elaborate on these comparisons through experiments in Section 7.

5 STABILITY

We next examine conditions for the stability of the NE in these games. When small changes occur to the underlying model parameters, a new NE may result. Intuitively, if the new NE is close enough to the original one, then we say the original NE is stable.

Formally, we generalize our utility functions in Eqn (1) to the family of parameterized functions $u_i(\boldsymbol{x}_i, \boldsymbol{x}_{-i}, \boldsymbol{p}_i)$, where $\boldsymbol{p}_i = [p_i^{(1)}, \cdots, p_i^{(K)}] \in \mathbb{R}^K$ is a vector valued perturbation parameter or shock on a_i , and $\boldsymbol{p} = [\boldsymbol{p}_1, \cdots \boldsymbol{p}_N] \in \mathbb{R}^{NK}$ denotes the vector of all perturbations/shocks. Moreover, let $\boldsymbol{x}^*(\boldsymbol{p})$ be the action profile at the NE of the game under perturbation vector \boldsymbol{p} and \boldsymbol{x}^* be the NE of the unperturbed game $(\boldsymbol{x}^* := \boldsymbol{x}^*(\boldsymbol{0}))$.

We denote a ball of radius r>0 centered at $\boldsymbol{x}\in\mathbb{R}^N$ by $B(\boldsymbol{x},r):=\{\boldsymbol{y}\in\mathbb{R}^N:\|\boldsymbol{x}-\boldsymbol{y}\|_2< r\}.$

Definition 5. ([36]) A NE x^* is stable if $\exists r > 0, d > 0$ such that $\forall p \in B(0, r)$, the NE $x^*(p)$ exists and satisfies

$$\|m{x}^*(m{p}) - m{x}^*\|_2 \leq d\|H(m{x}^*(m{p}), m{p}) - H(m{x}^*(m{p}), m{0})\|_2 \;,$$
 where $H(m{x}, m{p}) = (H_i(m{x}, m{p}))_{i=1}^M$ with

 $H_i(\boldsymbol{x},\boldsymbol{p}) = \boldsymbol{x}_i - BR_i(\boldsymbol{x}_{-i},G,\boldsymbol{p})$.

Definition 5 states that if an NE x^* is stable, the NE of the perturbed game $(x^*(p))$ remains close to the NE of the unperturbed game $(x^*(0))$.

5.1 Stability Condition Without Network Structure

To determine whether an NE x^* is stable, [11] proposed dividing the agents' action indices into three disjoint sets based on x^* :

$$\begin{split} &A(\boldsymbol{x}^*) := &\{j = \psi(i,k) \mid x_i^{(k)*} > 0, \ x_i^{(k)*} = \tilde{BR}_i^{(k)}(\boldsymbol{x}_{-i}^*, u_i)\}, \\ &I(\boldsymbol{x}^*) := &\{j = \psi(i,k) \mid x_i^{(k)*} = 0, \ x_i^{(k)*} > \tilde{BR}_i^{(k)}(\boldsymbol{x}_{-i}^*, u_i)\}, \\ &B(\boldsymbol{x}^*) := &\{j = \psi(i,k) \mid x_i^{(k)*} = 0, \ x_i^{(k)*} = \tilde{BR}_i^{(k)}(\boldsymbol{x}_{-i}^*, u_i)\}, \end{split}$$

where $\tilde{BR}_i(\boldsymbol{x}_{-i}^*, u_i)$ is the *unbounded* best response and can take negative values, $\tilde{BR}_i^{(k)}$ denotes the k-th action dimension

unbounded best response, and $\psi:\{1,\ldots,N\}\times\{1,\ldots,K\}\mapsto\{1,\ldots,KN\}$ maps the $N\times K$ agent-action indices pair to the KN indices in the unstructured operator F. $A(\boldsymbol{x}^*)$ is referred to as the set of active indices, $I(\boldsymbol{x}^*)$ the set of strictly inactive indices, and $B(\boldsymbol{x}^*)$ the set of borderline inactive indices. Intuitively, with a small parametric perturbation \boldsymbol{p} , agent action indices in $A(\boldsymbol{x}^*)$ remain active $(x_i^{(k)*}(\boldsymbol{p})>0)$ and agent action indices in $I(\boldsymbol{x}^*)$ remain inactive $(x_i^{(k)*}(\boldsymbol{p})=0)$, while agent action indices in $B(\boldsymbol{x}^*)$ can transform from inactive to active $(x_i^{(k)*}(\boldsymbol{p})>x_i^{(k)*}(\boldsymbol{0})=0)$. Under these definitions, [11] established the following sufficient condition for the solution to VI(Q,F) to be stable in the sense of Definition 5.

Theorem 5.1. ([11]) Consider the matrix

$$\nabla_{A,B}F_{A,B}(\boldsymbol{x}^*) = \begin{bmatrix} \nabla_A F_A(\boldsymbol{x}^*) & \nabla_B F_A(\boldsymbol{x}^*) \\ \nabla_A F_B(\boldsymbol{x}^*) & \nabla_B F_B(\boldsymbol{x}^*) \end{bmatrix}$$
(7)

where $\nabla_{S_1}F_{S_2}(\boldsymbol{x}^*)$ is a sub-matrix of $\nabla F(\boldsymbol{x}^*)$ with rows and columns corresponding to the agent action indices in sets S_1 and S_2 (not necessarily groups), respectively, and $\nabla_{A,B}F_{A,B}(\boldsymbol{x}^*)$ is generated by selecting rows and columns corresponding to $A \cup B$ from the game Jacobian $\nabla F(\boldsymbol{x}^*)$. If $\nabla_{A,B}F_{A,B}(\boldsymbol{x}^*)$ is positive definite on Q, then the solution \boldsymbol{x}^* to VI(Q,F) is stable.

Below we provide an easier to verify condition as compared to that in Theorem 5.1 by taking the partition structure into account.

5.2 Stability Condition with Partition Structure

Similar to [11], we divide *partitions* into active, strictly inactive, and borderline inactive sets. Specifically: (1) a partition is active if at least one agent action index in that partition is active at NE \boldsymbol{x}^* ; (2) if all agent action indices in a partition are strictly inactive, then the partition is strictly inactive; (3) if all agent action indices of a partition are inactive and at least one of them is borderline inactive, then the partition is considered as a borderline inactive partition. Formally, we have,

$$A_{S}(\boldsymbol{x}^{*}) := \{\mathcal{P}_{i} \mid \boldsymbol{x}_{\mathcal{P}_{i}}^{*} \neq \boldsymbol{0}\},$$

$$I_{S}(\boldsymbol{x}^{*}) := \{\mathcal{P}_{i} \mid \boldsymbol{x}_{\mathcal{P}_{i}}^{*} = \boldsymbol{0}, \boldsymbol{x}_{\mathcal{P}_{i}}^{*} > \tilde{BR}_{\mathcal{P}_{i}}(\boldsymbol{x}^{*})\},$$

$$B_{S}(\boldsymbol{x}^{*}) := \{\mathcal{P}_{i} \mid \boldsymbol{x}_{\mathcal{P}_{i}}^{*} = \boldsymbol{0}\} - I_{S}(\boldsymbol{x}^{*}),$$
(8)

where $A_S(\boldsymbol{x}^*)$, $I_S(\boldsymbol{x}^*)$, $B_S(\boldsymbol{x}^*)$ denote the set of active, strictly inactive and borderline inactive partitions, respectively. We use $\boldsymbol{x}_{\mathcal{P}_i}^*$, and $\tilde{BR}_{\mathcal{P}_i}(\boldsymbol{x}^*)$ to denote the vectors by choosing all indices in partition \mathcal{P}_i from \boldsymbol{x}^* , and $\tilde{BR}(\boldsymbol{x}^*)$.

Theorem 5.2. Consider an NE x^* of the network game. Re-index all partitions in $A_S(x^*) \cup B_S(x^*)$ with indices $1, 2, \dots, Z, Z = |A_S(x^*)| + |B_S(x^*)|$. Then, define

$$H^{S}(\boldsymbol{x}^{*}) = \begin{bmatrix} \theta_{1}^{S}(\boldsymbol{x}^{*}) & -\delta_{12}^{S}(\boldsymbol{x}^{*}) & \dots & -\delta_{1Z}^{S}(\boldsymbol{x}^{*}) \\ -\delta_{21}^{S}(\boldsymbol{x}^{*}) & \theta_{2}^{S}(\boldsymbol{x}^{*}) & \dots & -\delta_{2Z}^{S}(\boldsymbol{x}^{*}) \\ \vdots & \vdots & \ddots & \vdots \\ -\delta_{Z1}^{S}(\boldsymbol{x}^{*}) & -\delta_{Z2}^{S}(\boldsymbol{x}^{*}) & \dots & \theta_{Z}^{S}(\boldsymbol{x}^{*}) \end{bmatrix}$$

where $\delta_{ij}^S(\boldsymbol{x}^*) = \|\nabla_j F_i(\boldsymbol{x}^*)\|_2$, $\theta_i^S(\boldsymbol{x}^*) = \|\nabla_i F_i(\boldsymbol{x}^*)\|_2$. If the following conditions hold simultaneously:

1)
$$H^S(x^*) \succ 0$$
,

2)
$$\nabla_i F_i(\mathbf{x}^*) \succ \mathbf{0}, \forall \mathcal{P}_i \in A_S \cup B_S$$
,

then x^* is stable. Moreover, these conditions are sufficient for $\nabla_{A,B}F_{A,B}(x^*) \succ \mathbf{0}$, i.e., the condition for stability on an unstructured network (Theorem 5.1) holds under these conditions.

Intuitively, the matrix $H^S(\boldsymbol{x}^*)$ captures the mutual influence between active and borderline inactive partitions at the current NE profile. The borderline inactive partitions can turn into active partitions under parametric perturbations. When such flips are significant, large fluctuations can appear in the network, which can be further amplified through rebounds and reflections. In this case, new equilibria may not exist, and even if they do, they may be far away from the original equilibrium. However, when $H^S(\boldsymbol{x}^*) \succ \mathbf{0}$ holds, the maximum impacts of flipping partitions from (borderline) inactive to active are bounded, and therefore the current NE remains stable.

In terms of the complexity of verifying these conditions, note that H^S is a $Z \times Z$ matrix. Similar to the comparison shown in Section 4, if we denote $Y = |A(\boldsymbol{x}^*)| + |B(\boldsymbol{x}^*)|$, then the computational complexity of condition verification in Proposition 5.2 vs. Theorem 5.1 are $O(Z^3 + Z^2\overline{N}^3)$ vs. $O(K^3Y^3)$, where \overline{N} is the average group size. Therefore, since Z < KY ($Z\overline{N} \approx KY$), the computational complexity of condition verification in Proposition 5.2 is lower than that of Theorem 5.1.

We conclude this section with a stability condition on Υ^S .

Theorem 5.3. Assume Υ^S is symmetric. Then if both of the following conditions hold:

1) $\Upsilon^S \succ \mathbf{0}$,

2) $\Upsilon^{U}_{\mathcal{P}_i,\mathcal{P}_i}$, $\forall i$ are P-matrices,

the network game's NE is unique and stable.

6 CENTRALITY

In network games, notions of node centrality are used to measure the influence of individual nodes on network-level outcomes. Degree centrality is one of the centrality metrics which has gained attention in the literature [37], [38]. In a directed graph, two different measures of degree centrality are considered for each node: in-degree centrality, which is a count of edges directed to a given node, and out-degree centrality, which is the number of outward edges from the given node. In this section, we propose a generalization of the degree centrality for disjoint partitions.

Recall that we capture the influence of a partition using the Jacobian matrix $\nabla F(x)$. Matrix $\nabla_j F_i(x)$ measures the sensitivity of partition i to the action profile of partition j. Accordingly, we define our generalized centrality measure as follows.

Definition 6. Generalized Degree Centrality (GDC): Following Definition 4, denote $\beta_{ij}^S = \sup_{\boldsymbol{x} \in Q} \|\nabla_j F_i(\boldsymbol{x})\|_2$, and $\alpha_i^S = \inf_{\boldsymbol{x} \in Q} \|\nabla_i F_i(\boldsymbol{x})\|_2$. The generalized degree centralities for partition \mathcal{P}_i are given by:

$$D_i^{in} = \sum_{j:j \neq i} \frac{\beta_{ij}^S}{\alpha_i^S}, \quad D_i^{out} = \sum_{j:j \neq i} \frac{\beta_{ji}^S}{\alpha_j^S}, \ \forall i, j \in \mathbb{N}[1, M] \ .$$

Moreover, the maximum GDCs are defined as follows:

$$D_{max}^{in} = \max_{i \in \mathbb{N}[1,M]} D_i^{in}, \quad D_{max}^{out} = \max_{i \in \mathbb{N}[1,M]} D_i^{out} \;.$$

The above definition can be interpreted as follows: out-degree centrality measures the influence of a given partition \mathcal{P}_i on the network based on three factors, (1) connectivity, or the number of links directed outward from \mathcal{P}_i , (2) the internal impact of the target partitions that receive impact from \mathcal{P}_i , and (3) the external impact \mathcal{P}_i has for every target group. In Definition 6, D_i^{out} and D_i^{in} capture these factors through the summation of $\frac{\beta_{ji}^s}{\alpha_i^s}$ and $\frac{\beta_{ji}^s}{\alpha_i^s}$.

In addition to capturing importance due to their roles in the network, each partition can be endowed with certain exogenous (non-network related) importance (not to be confused with external impact from another partition). This will result in an extended centrality measure. The following is a generalization of the extended centrality measure defined in [39].

Definition 7. Generalized Extended Degree Centrality (GEDC): Let $e \in \mathbb{R}^{M}_{>0}$ denote the vector of external importance, where $(e)_i = e_i > 0$ is \mathcal{P}_i 's external importance. The generalized extended degree centralities for \mathcal{P}_i are given by

$$D_i^{in}(\boldsymbol{e}) = \sum_{i:j \neq i} \frac{\beta_{ij}^S}{\alpha_i^S} \frac{e_j}{e_i}, \ D_i^{out}(\boldsymbol{e}) = \sum_{i:j \neq i} \frac{\beta_{ji}^S}{\alpha_j^S} \frac{e_i}{e_j}, \forall i,j \in \mathbb{N}[1,M]$$

and the maximum GEDCs are defined as

$$D_{max}^{in}(\boldsymbol{e}) = \max_{i \in \mathbb{N}[1,M]} D_i^{in}(\boldsymbol{e}), \ D_{max}^{out}(\boldsymbol{e}) = \max_{i \in \mathbb{N}[1,M]} D_i^{out}(\boldsymbol{e}) \ .$$

When $e = \alpha \mathbf{1}, \alpha > 0$, Definitions 6 and 7 are equivalent. We now connect our centrality measure and the uniqueness of the NE.

Theorem 6.1. *If both of the following conditions hold:*

1) $\Upsilon^{U}_{\mathcal{P}_{i},\mathcal{P}_{i}}$, $\forall i$ are P-matrices, 2) \exists $e \succ \mathbf{0}$ such that $D^{in}_{max}(e) < 1$, or $D^{out}_{max}(e) < 1$, then the NE is unique. If in addition Υ^{S} is symmetric, the NE is

Theorem 6.1 implies that if either the in-degree or out-degree GEDCs are bounded, then the NE is unique. On the other hand, if neither the indegree nor outdegree is bounded, then at least one partition has an outsized effect on the network. This partition's decision can change the state of the network significantly, resulting in possibly multiple equilibria.

Theorem 6.1 is similar to Proposition 7 in [30], but differs as follows. In our work, β_{ij}^S represent the influence of partitions on each other, while β_{ij} in [30] represent the component-level influence of agents on each other. Moreover, when both conditions in Theorem 6.1 hold, $\Upsilon^S \succ 0$; this then becomes a special case of the condition in Corollary 4.3.2 (where Υ^S is symmetric).

Moreover, if Υ^S is symmetric, Theorem 6.1 shows that the unique NE is also stable. Intuitively, this is because if a partition's degree centrality is not bounded, it has a considerable impact on the network, and the NE may not be stable as a small perturbation affecting this partition can influence the network dramatically.

7 NUMERICAL RESULTS

We now present numerical results to compare the verification complexity and sufficiency gaps on the two sets of structured and unstructured conditions. We first use games where both sets of conditions are satisfied to allow comparisons of verification complexity. We then generate games using different parameter settings, and measure how frequently the structured conditions fail while the unstructured conditions are satisfied, to assess the sufficiency gap. We detail the generation steps of the game instances on which these comparisons are done in Appendix J.2.

7.1 The Computational Complexity Gap

As discussed in Section 4, the verification of uniqueness conditions on structured networks is of lower complexity. There are several factors that determine how big this complexity gap is, which we examine in this section.

The first factor is the size of the network (total number of agents). Specifically, the complexity gap increases with the number of agents, and is at least quadratic in the number of groups. Table 1 lists the verification complexity (in floating point operations, FLOPs) of the conditions in Corollary 4.3.2 with a dense Γ^U matrix. We see that the verification complexity of structured conditions is orders of magnitude lower than that of the unstructured conditions, and the gap increases with the size of the network. Table 2 shows the difference in CPU times, where the results are averaged over 50 different instances. Moreover, the complexity reduction in verifying the conditions in Theorem 4.3 is much more significant than those in Corollary 4.3.2. For instance, the verification complexity of conditions in Theorem 4.3 on a game of size 10×10 (10 partitions of 10 agents each) on a dense Υ^U matrix is 1.08×10^{35} FLOPs while that on the corresponding Υ^S matrix is 1.26×10^6 FLOPs. We refer the interested reader to Appendix I for additional comparisons.

Size $(N_i \times M)$	Unstructured	Structured
10×10	6.67×10^5	5.83×10^4
20×20	4.27×10^7	1.99×10 ⁶
50×50	1.04×10^{10}	2.02×10 ⁸
100 × 100	6.67×10^{11}	6.57×10 ⁹
200 × 200	4.27×10^{13}	2.12×10 ¹¹

TABLE 1 Verification complexity in number of FLOPs of conditions in Corollary 4.3.2 over the number of agents.

Size $(N_i \times M)$	Unstructured (sec)	Structured (sec)
10×10	0.0031 ± 0.0011	$0.003 {\pm} 0.0002$
20×20	0.0575±0.0074	0.0213±0.0019
50×50	5.851 ± 0.197	$0.6768 {\pm} 0.0442$
100 × 100	332.8±1.6	7.052 ± 0.203
200 × 200	Memory Overflow	116.8±6.9

TABLE 2

Verification complexity in CPU times of conditions in Corollary 4.3.2 over the number of agents; all times are in seconds. All experiments were performed on a machine with A 6-core 2.60/4.50 GHz CPU with hyperthreaded cores, 12MB Cache, and 16GB RAM.

The second factor affecting the complexity gap is how (a given number of) agents are partitioned into groups. Figure 5 and 6 shows the complexity (in FLOPs) of verifying the structured condition (of Υ^S being a P-matrix) in two games of size 50 and 100 agents, respectively, both with a dense Υ^U matrix, as we vary the number of groups. Here, we set to groups to be of equal size, rounding off to the nearest integer if needed; e.g., with 10 agents and 3 groups, the partition sizes are 3, 3, 4.²

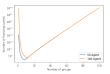
We see that in each game the complexity has a V shape, reaching a minimum when $M \approx \sqrt{N}$. To explain this, note that we can approximate the complexity of LU decomposition by $\frac{2}{3}p^3$ for a matrix of size $p \times p$; for a $p \times k$ matrix (p > k), the approximate number of FLOPs for singular value decomposition is $2pk^2 - \frac{2}{3}k^3$ ([40, p. 75], based on QR decomposition using

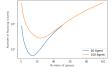
2. Here we are comparing the structured complexity between games of different sizes. For a given size N with M equal-sized partitions, the structured and unstructured conditions have the same complexity when M=1 or N.

Householder transformations). Then the complexity for checking the conditions in Theorem 4.3 is given by (S = N/M):

$$\binom{M}{2} \left(2S^3 - \frac{2}{3}S^3\right) + M \sum_{k=1}^{S} \binom{S}{k} \left(\frac{2}{3}k^3\right) + \sum_{k=1}^{M} \binom{M}{k} \left(\frac{2}{3}k^3\right).$$

As $\max\{M!, S!\}$ will be the dominant term in the (expanded) expressions, the minimum is achieved when $M = S = \sqrt{N}$.





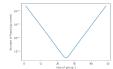


Fig. 5. Verification complexity (FLOPs) of conditions in Theorem 4.3 over the number of groups.

Fig. 6. Verification complexity (FLOPs) of conditions in Corollary 4.3.2 over the number of groups.

 $\begin{array}{lll} \text{Fig.} & 7. & \text{Complexity} \\ \text{(FLOPs)} & \text{over} \\ \text{different} & \text{partitions} \\ \text{in games} & \text{with} & 50 \\ \text{agents, 2 groups and} \\ \text{a dense } \Upsilon^U \text{ matrix.} \end{array}$

Note also that in Figure 5 the two curves overlap almost completely beyond the minimum. This is because checking if Υ^S is a P-matrix has two phases: generating Υ^S , then an eigendecomposition on it. When M is small the first phase dominates, whereas when M is larger than \sqrt{N} the second becomes dominant. Since Υ^S is $M \times M$, when M is large the complexity depends much more on M than on the network size N.

In another experiment, we verify the conditions in Corollary 4.3.2 on two games of sizes 50 and 100, respectively, both with a dense Υ^U matrix. The results are shown in Figure 6. This time the minimum occurs at some $M>\sqrt{N}$. This is because the approximated complexity is given by $\binom{M}{2}\left(2S^3-\frac{2}{3}S^2\right)+\frac{2}{3}M^3+\frac{2}{3}S^3$, which has a minimum at $M>\sqrt{N}$.

A third factor affecting the complexity gap is the size distribution of groups, given fixed N and M. Figure 7 shows the complexity of verifying whether Υ^S is a P-matrix in a 50-agent, 2-partition game with a dense Υ^U matrix, as we vary the size of the first group. We see that the complexity reaches its minimum when the two groups are equal sized.

More generally, when we have more action dimensions and create partitions based on them, we can expect that some of the off-diagonal elements in the structured matrix Υ^S will be computed with much lower complexity. To see why, consider two games, one with $N=100,\,K=1$ and 10 groups of size 10 each, the other with $N=50,\,K=2$ and 5 groups of size 10 each. If we create a partition for all agent-action components from the same group in the same action, we get 10 partitions for both games. While the first game is similar to the samples in this section, the second game has 20 off-diagonal elements in Υ^S computed from the coupled cost. Computing these 20 elements are easier than computing the 2-norm of a matrix since the corresponding block is a diagonal matrix. Moreover, if all agents have the same cost function, then these 20 elements are the same and based on one utility function, which makes the computation even easier.

7.2 The Sufficiency Gap

The structured conditions are stronger than their unstructured counterparts, thus may fail to discover the uniqueness and stability of an NE in a game that fails the former while satisfying the latter. We showed this using an example in Section 4. In what follows we

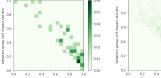
will use numerical results to measure this sufficiency gap. We will focus on the uniqueness conditions, and note that the comparison of stability conditions is very similar.

Each of the next set of figures is a heat map showing how often these two conditions in Corollary 4.3.2 (over Γ^S and Γ^U , respectively, with sample games that guarantee both Γ^S and Γ^U are symmetric) yield the same or different result. The former means either both are satisfied or not satisfied, while the latter necessarily means the structured condition fails and the unstructured condition holds. Each game is of size 400, with 20 groups of 20 members each. For each cell in the heat map, 50 sample games are generated using a set of parameters corresponding to the cell indicated on the figure; the cell color indicates the fraction of these games that resulted in a difference (sufficiency gap), the higher the fraction, the darker the color. In each of the heat maps, we can see regions of darker cells (clearly) separating the map into two lighter regions. In general, the bottom left represents parameter settings where both structured and unstructured conditions are satisfied and the top right represents settings where both conditions do not hold.

Specifically, in Figures 8, 9, and 10 we hold component-level external impact fixed (at weak, medium, and strong levels, respectively, corresponding to connection frequency at 0.2, 0.5, 0.8 respectively; in-partition connection strengths at $0.2/N_i, 0.5/N_i, 0.8/N_i$ respectively, for partition \mathcal{P}_i (normalizing the strengths to make external and internal impact comparable); between-partition connection strengths fixed at 0.2/N, 0.5/N, 0.8/N (normalize the strengths, similar as above), while changing the variances of component-level internal impact. We note that the normalized upper bound for the model parameters are chosen at 1 when we study the sufficiency gaps, because higher values cause both sets of conditions to fail thereby reducing the significance of such sample games. Please refer to Appendix J for more details on the settings.

Overall the sufficiency gap is quite low, i.e., the two types of conditions yield the same outcome in the vast majority of parameter settings (as evidenced by mostly 0 values/light-colored cells on these heat maps). The measured difference (obtained by adding the number of different results in every cell and dividing by the total number of sample games) in Figures 8, 9, and 10 are 0.26%, 5.58%, 0.05%, respectively.

Figures 8 and 9 show a similar pattern. In the top right corners where component-level internal impact has large variance both between and within groups, neither condition is true, while in the lower left corners both conditions are satisfied, giving rise to the broad agreement (light regions). In comparing the two, we see that the dark region expands and shifts leftward and downward in Figure 9, suggesting that the gap between the two conditions is bigger and triggered by lower variances in component-level internal impact when the component-level external impact increases from weak to medium. Furthermore, high within-partition variance combined with low between-partition variances in componentlevel internal impact results in the largest gap. The reason is the "weakest member effect" discussed in Section 4, where the α_i^S values depend highly on the minimum component-level internal impact $\alpha_k^{\bar{U}}$ of members in \mathcal{P}_i , which makes the abstraction Υ^S inaccurate, causing a larger gap. Interestingly, an increase in between-partition variances in component-level internal impact can mitigate the above effect and reduce the gap. On the other hand, when the component-level external impact is sufficiently high, Figure 10 shows that the sufficiency gap all but disappears as now most game instances do not satisfy either condition.



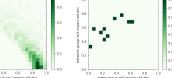
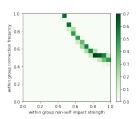


Fig. 8. Sufficiency gap over the variance of internal impact, weak ext. impact.

Fig. 9. Sufficiency gap over the variance of internal impact, medium ext. impact.

Fig. 10. Sufficiency gap over the variance of internal impact, strong ext. impact.

We next fix the agent's component-level internal impact at 1, the between-partition (resp. within-group) component-level external impact at strong, and vary the within-partition (resp. between-group) component-level external impact strength (x-axis) and connection frequency (y-axis), shown in Figure 11 (resp. Figure 12). We see that the gap is generally small (2% gap forboth). In both figures the disagreement is concentrated around a reciprocal curve, suggesting that when the product of the two parameters is around a critical level, the sufficiency gaps occur. They also suggest that the role of between-partition and withinpartition external impact is very similar under these two sets of conditions. The dark regions in these two figures suggest that when individuals have a homogeneous internal impact, games where the expected sum of component-level external impact is about 10% higher than the expected sum of component-level internal impact, are most likely to have sufficiency gaps. We provide more detailed discussions in Appendix J. Moreover, the dark cell curve in Figure 12 shows that when the between-partition connection frequency gets lower, the chance of having a sufficiency gap is higher. This is because a mismatch similar to Example 2 (an agent with weak component-level internal impact does not receive strong component-level external impact but a member in the same partition with strong component-level internal impact does) is more likely to happen in this case.



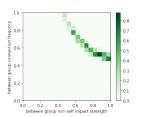


Fig. 11. Sufficiency comparison over the within partition connections, when every agent has the same internal impact and strong between partition influence.

Fig. 12. Sufficiency comparison over the between partition connections when every agent has the same internal impact, medium within partition influence.

These numerical results suggest that in general, when network communities are formed around homogeneous agents, i.e., those with similar component-level internal impact, component-level external impact and connectivity, then the two types of conditions yield identical verification outcomes.

More generally, when we have more action dimensions and create partitions based on them, we can expect some of the off-diagonal elements in the structured matrix Υ^S to be computed from the coupled cost function. Similar to Section 7.1, we compare the sufficiency gap in two games, one with $N=100,\,K=1$ and 10 groups of size 10 each, the other with $N=50,\,K=2$

and 5 groups of size 10 each, and create 10 partitions for both games. Again, the first game is similar to the samples in this section, while the second game has 20 off-diagonal elements in Υ^S computed from the coupled cost. When all agents have the same cost functions, these 20 elements do not suffer from the strongest link effect and thus further reduce the sufficiency gap.

We also perform similar experiments on the real-world email-Eu-core network [41]. This is a binary directed graph with edges indicating whether an email was ever sent from one node to another, where nodes represent people from a large European research institution. We compare three ways of partitioning the agent-action space in Fig 13, 14, 15: (1) original partition based on the nodes' department affiliation, (2) degree-based, which puts nodes with similar in- and out- degrees in the same partition, and (3) connection-type: which puts nodes with similar fraction of within-department connections in the same partition. We observe that the degree-based partition performs best, as it successfully identifies and retains the importance of strongly connected subgraphs in our reduced conditions. We provide further elaborate and additional experiments on this network, in Appendix J.3.

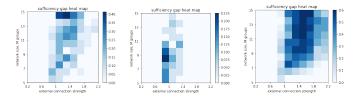


Fig. 13. Sufficiency Fig. 14. Sufficiency gap, original partition. gap, degree-based.

Fig. 15. Sufficiency gap, connection-type.

8 Conclusions and Discussions

We introduced and studied a family of structured network games with non-linear best response functions. Prior works on network games have found sufficient conditions for the uniqueness and stability of Nash equilibria which are mostly difficult to verify. In this work, we showed that the existence of structure in the network (e.g., in the form of communities, or when there are multirelational dependencies between agents), helps us find alternatives for such conditions, which we refer to as "structured conditions" as opposed to the "unstructured conditions" in previous works. In particular, we show that the structured conditions for the uniqueness and stability of NE are related to matrices which are possibly lower dimensional, with their dimensions depending on the number of partitions naturally arising in a network due to its structured nature. We also demonstrated both analytically and numerically that the structured conditions are sufficient conditions to the unstructured conditions, and that their verification is of much lower computational complexity. We used numerical experiment results to show that the sufficiency gap between the structured conditions and unstructured conditions is small in general and typically occurs in games with some specific characteristics. Moreover, we proposed a new notion of degree centrality to evaluate the influence of a partition in the network, and used it to identify additional conditions for uniqueness and stability.

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Kun Jin received his Ph.D. degree in electrical and computer engineering from the University of Michigan, Ann Arbor, in 2023. He is a Machine Learning Scientist at TikTok. His research interests are in algorithmic game theory, network economics, mechanism design, and data mining in cybersecurity.

Parinaz Naghizadeh received her Ph.D. degree in electrical engineering from the University of Michigan, Ann Arbor, in 2016. She is an Assistant Professor with the Department of Electrical and Computer Engineering at the University of California, San Diego. Her research interests include network economics, game theory, algorithmic economics, and reinforcement learning.

Mingyan Liu received her Ph.D in electrical engineering from the University of Maryland, College Park, in 2000. She is the Associate Dean of Academic Affairs and a professor of Electrical and Computer Engineering at the University of Michigan, Ann Arbor. Her interests are in sequential decision and learning theory, game theory and incentive mechanisms, with applications to large-scale networked systems. She is a Fellow of the IEEE and a member of the ACM.

APPENDIX A SUMMARY OF NOTATIONS

Notation	Meanings
a_i	i-th agent in the game
N	The set of agents
u_i	Utility function of a_i
d_i	Individual component of u_i
f_i	Network interaction component of u_i
x_i	Action profile of a_i
x_{-i}	Action profile of all agents other than a_i
x	Action profile of all agents
x^*	Equilibrium action profile
K	The action dimension number
G	Adjacency matrix
$G^{(k)}$	Adjacency matrix on the k-th action dimension
$G_{ij}^{(k)}$	Influence strength from a_j to a_i on the k -th action dimension
$\overline{\mathcal{P}_i}$	The <i>i</i> -th partition
\overline{F}	Best response operator
∇F	Game Jacobian
BR_i	Best response function of \mathcal{P}_i
$\overline{Q_i}$	Action space of \mathcal{P}_i
\overline{Q}	Action space
Υ^U	A matrix for unstructured network equilibrium uniqueness condition
Υ^S	A matrix for structured network equilibrium uniqueness condition
Γ^U	A matrix for an alternative unstructured network equilibrium uniqueness condition
Γ^S	A matrix for an alternative structured network equilibrium uniqueness condition
ρ	Spectral radius
λ_k	The k-th lowest eigenvalue
α_i^U	Internal impact term in Υ^U
β_{ij}^U	External impact term in Υ^U
α_i^S	Internal impact term in Υ^S
β_{ij}^{S}	External impact term in Υ^S
$A_S(\boldsymbol{x}^*)$	The set of active partitions at x^*
$A_S(\boldsymbol{x}^*)$	The set of borderline active partitions at $oldsymbol{x}^*$
$I_S(oldsymbol{x}^*)$	The set of inactive partitions at $oldsymbol{x}^*$
H^S	A matrix for structured network equilibrium stability condition
	C 1: 1: 1 CD
D_i^{in}	Generalized in-degree of \mathcal{P}_i
	Generalized in-degree of \mathcal{P}_i Generalized out-degree of \mathcal{P}_i
D_i^{in} D_i^{out} D_{max}^{in}	

TABLE 3 Summary of notation.

APPENDIX B VALUES USED IN FIGURE 1-4

We use Figures 1-4 to illustrate the high level idea of the partitioning on structured multi-relational networks. In Figure 1 and 2, using the linear quadratic utility functions $u_i = b_i x_i + \sum_{j \neq i} g_{ij} x_i x_j - \frac{1}{2} x_i^2$, the adjacency matrices are

$$G = \begin{bmatrix} 0 & 0.6 & 0.2 \\ 0.6 & 0 & 0.2 \\ 0.2 & 0.2 & 0 \end{bmatrix}, \ G = \begin{bmatrix} 0 & 0 & 0.2 \\ 0 & 0 & 0.2 \\ 0.2 & 0.2 & 0 \end{bmatrix}$$

respectively. In Figure 3 and 4, when using the utility function

$$u_{i} = b_{i}x_{i} + \sum_{j \neq i} g_{ij}^{(1)} x_{i}^{(1)} x_{j}^{(1)} + \sum_{j \neq i} g_{ij}^{(2)} x_{i}^{(2)} x_{j}^{(2)} - \frac{1}{2} ((x_{i}^{(1)})^{2} + (x_{i}^{(2)})^{2} + 0.2x_{i}^{(1)} x_{j}^{(1)}),$$

the adjacency matrices are

$$\begin{split} G^{(1)} &= \begin{bmatrix} 0 & 0.6 & 0 \\ 0.6 & 0 & 0 \\ 0 & 0 & 0 \end{bmatrix}, \ G^{(2)} &= \begin{bmatrix} 0 & 0 & 0 \\ 0 & 0 & 0.6 \\ 0 & 0.6 & 0 \end{bmatrix}, \\ G &= \begin{bmatrix} G^{(1)} & 0 \\ 0 & G^{(2)} \end{bmatrix}. \end{split}$$

APPENDIX C PROOF OF LEMMA 4.2

Proof. We will prove that $\|\Upsilon_{\mathcal{P}_i,\mathcal{P}_j}^U\|_2 \geq \beta_{ij}^S$. The proof establishing $\|\Upsilon_{\mathcal{P}_i,\mathcal{P}_i}^U\|_2 \leq \alpha_i^S$ is similar.

From the definition of β^S_{ij} in Eqn (4), we denote the action profile in Q that obtains the value of β^S_{ij} as $\hat{\boldsymbol{x}}(i,j)$, i.e.,

$$\beta_{ij}^S = \sup_{\boldsymbol{x} \in O} \|\nabla_j F_i(\boldsymbol{x})\|_2 = \|\nabla_j F_i(\hat{\boldsymbol{x}}(i,j))\|_2,$$

where $\nabla_j F_i(\hat{x}(i,j)) \in \mathbb{R}^{N_i \times N_j}$. We note here that since our action space Q is compact and ∇F is assumed to be continuous and differentiable on q, there exist maxima for $\|\nabla_j F_i(x)\|$ such that $\hat{x} \in Q$. Therefore, the supremum equals the maximum and can be achieved. Since we will only focus on a specific pair of (i,j) values where $i \neq j$ in this part, we will simply use \hat{x} to denote this action profile for convenience.

Also, based on the definition of β_{kl}^U , we have

$$eta_{kl}^U = \sup_{oldsymbol{x} \in Q} |\nabla_l F_k(oldsymbol{x})| \ge |\nabla_l F_k(\hat{oldsymbol{x}})|,$$

where $\nabla_l F_k(\hat{\boldsymbol{x}}) \in \mathbb{R}$.

Next, we perform Singular Value Decomposition on $\nabla_j F_i(\hat{\boldsymbol{x}})$ such that $\nabla_j F_i(\hat{\boldsymbol{x}}) = USV^T$. From this, we can get the left singular vector $\boldsymbol{u} \in \mathbb{R}^{N_i}$ and the right singular vector $\boldsymbol{v} \in \mathbb{R}^{N_j}$ that correspond to the largest singular value β_{ij}^S , where $\|\boldsymbol{u}\| = \|\boldsymbol{v}\| = 1$. Then

$$\boldsymbol{u}^T \nabla_j F_i(\hat{\boldsymbol{x}}) \boldsymbol{v} = \beta_{ij}^S \cdot \|\boldsymbol{u}\| \cdot \|\boldsymbol{v}\| = \beta_{ij}^S.$$

To proceed with the proof, we introduce the following matrix A obtained from matrix $\nabla_j F_i(\hat{\boldsymbol{x}}(i,j))$ by replacing every element with its absolute value, formally,

$$A = (|\nabla_l F_k(\hat{\boldsymbol{x}})|)_{k,l:a_k \in \mathcal{P}_i, a_l \in \mathcal{P}_i}.$$

We also introduce vectors u^+ and v^+ , where

$$u^+ = (|u_k|)_{k=1}^{N_i}, \ v^+ = (|v_l|)_{l=1}^{N_j}.$$

Then, $\| oldsymbol{u}^+ \| = \| oldsymbol{v}^+ \| = 1$, and it is easy to see that

$$(\boldsymbol{u}^+)^T A \boldsymbol{v}^+ \geq \boldsymbol{u}^T \nabla_j F_i(\hat{\boldsymbol{x}}) \boldsymbol{v},$$

Next, consider the product $u^T \Upsilon^U_{\mathcal{P}_i, \mathcal{P}_j} v$. We denote the index in $\Upsilon^U_{\mathcal{P}_i, \mathcal{P}_j}$ that corresponds to the element u_p and v_q as k_p and l_q respectively, so that,

$$oldsymbol{u}^T \Upsilon^U_{\mathcal{P}_i,\mathcal{P}_j} oldsymbol{v} = \sum_{p=1}^{N_i} \sum_{q=1}^{N_j} \Upsilon^U_{k_p l_q} \cdot u_p \cdot v_q \; .$$

then we have

$$\begin{split} \|\boldsymbol{\Upsilon}_{\mathcal{P}_{i},\mathcal{P}_{j}}^{U}\| &= \|-\boldsymbol{\Upsilon}_{\mathcal{P}_{i},\mathcal{P}_{j}}^{U}\| \cdot \|\boldsymbol{u}^{+}\| \cdot \|\boldsymbol{v}^{+}\| \\ &\geq (\boldsymbol{u}^{+})^{T}(-\boldsymbol{\Upsilon}_{\mathcal{P}_{i},\mathcal{P}_{j}}^{U})\boldsymbol{v}^{+} \\ &= -\sum_{p=1}^{N_{i}}\sum_{q=1}^{N_{j}}\boldsymbol{\Upsilon}_{k_{p}l_{q}}^{U} \cdot |u_{p}| \cdot |v_{q}| \\ &= \sum_{p=1}^{N_{i}}\sum_{q=1}^{N_{j}}\beta_{k_{p}l_{q}}^{U} \cdot |u_{p}| \cdot |v_{q}| \\ &\geq \sum_{p=1}^{N_{i}}\sum_{q=1}^{N_{j}}|\nabla_{l_{q}}F_{k_{p}}(\hat{\boldsymbol{x}})| \cdot |u_{p}| \cdot |v_{q}| \\ &= (\boldsymbol{u}^{+})^{T}A\boldsymbol{v}^{+} \\ &\geq \boldsymbol{u}^{T}\nabla_{j}F_{i}(\hat{\boldsymbol{x}})\boldsymbol{v} \\ &= \beta_{i:i}^{S_{i}}. \end{split}$$

This completes the proof.

APPENDIX D PROOF OF THEOREM 4.3

Proof. Given the community level partition $Q = \prod_{i=1}^M Q_i$, we denote $\nabla F_i(\boldsymbol{z}) = ((\nabla_i F_j(\boldsymbol{z}))_{j=1}^M)^T \in \mathbb{R}^{N_i \times N}$.

We use the notation L(x, y) to denote the line segment between two points x and y in \mathbb{R}^N . Formally,

$$L(x, y) = {\alpha x + (1 - \alpha)y : 0 < \alpha < 1}.$$

Under Assumption 1 in Section II, $F_i: Q \to \mathbb{R}^{N_i}, Q \subseteq \mathbb{R}^N$ is continuously differentiable on Q, and for $\forall x, y$ in $Q_i, L(x, y) \subseteq Q_i$. According to [42, Theorem 12.9] we know that for every vector \mathbf{a} in \mathbb{R}^{N_i} , there is a point $\mathbf{z} \in L(x, y)$ such that:

$$\mathbf{a} \cdot (F_i(\mathbf{x}) - F_i(\mathbf{y})) = \mathbf{a} \cdot (\nabla F_i(\mathbf{z})(\mathbf{x} - \mathbf{y})).$$
 (9)

Let \boldsymbol{a} in equation (9) be $(\boldsymbol{x}_i - \boldsymbol{y}_i)^T$, and denote $\boldsymbol{l} = (l_j)_{j=1}^M$, where $l_j = \|\boldsymbol{x}_j - \boldsymbol{y}_j\|_2, \forall j \in \mathbb{N}[1, M]$, then,

$$(\boldsymbol{x}_{i} - \boldsymbol{y}_{i})^{T} (F_{i}(\boldsymbol{x}) - F_{i}(\boldsymbol{y}))$$

$$= (\boldsymbol{x}_{i} - \boldsymbol{y}_{i})^{T} (\nabla F_{i}(\boldsymbol{z})(\boldsymbol{x} - \boldsymbol{y}))$$

$$= (\boldsymbol{x}_{i} - \boldsymbol{y}_{i})^{T} [\sum_{j=1}^{M} \nabla_{i} F_{j}(\boldsymbol{z})(\boldsymbol{x}_{j} - \boldsymbol{y}_{j})]$$

$$\geq (\boldsymbol{x}_{i} - \boldsymbol{y}_{i})^{T} \nabla_{i} F_{i}(\boldsymbol{z})(\boldsymbol{x}_{i} - \boldsymbol{y}_{i})$$

$$-|\sum_{j \neq i} (\boldsymbol{x}_{i} - \boldsymbol{y}_{i})^{T} \nabla_{i} F_{j}(\boldsymbol{z})(\boldsymbol{x}_{j} - \boldsymbol{y}_{j})|$$

$$\geq \alpha_{i}^{S}(l_{i})^{2} - \sum_{j \neq i} \beta_{ij}^{S} \cdot l_{i} \cdot l_{j}$$

$$= l_{i} \cdot (\Upsilon^{S} \boldsymbol{l})_{i}$$

$$(10)$$

By [43, Theorem 3.3.4(b)], a real square matrix $M \in \mathbb{R}^{n \times n}$ is a P-matrix if it satisfies $\forall l \in \mathbb{R}^n$

$$\max_{i \in \mathbb{N}[1,M]} l_i(M\boldsymbol{l})_i > 0$$

Denote $b=\max_{i\in\mathbb{N}[1,M]} \frac{l_i\cdot[\Upsilon^{\mathcal{C}}l]_i}{\|l\|_2^2}>0.$ Then, we have

$$\max_{i \in \mathbb{N}[1,M]} (\boldsymbol{x}_i - \boldsymbol{y}_i)^T (F_i(\boldsymbol{x}) - F_i(\boldsymbol{y})) \ge \max_{i \in \mathbb{N}[1,M]} l_i \cdot [\Upsilon^C \boldsymbol{l}]_i$$

$$\ge b \cdot \|\boldsymbol{x} - \boldsymbol{y}\|_2^2$$

which, according to Definition 2.(b), shows that F satisfies uniform block P-condition. Therefore, by [10, Proposition 2 part (b)] and [31, Proposition 3.5.10 part (b)], the Nash equilibrium is unique.

Finally, we show that the two conditions

1) Υ^S is a P-matrix,

2) $\Upsilon^{U}_{\mathcal{P}_{i},\mathcal{P}_{i}}$ are P-matrices,

are sufficient for Υ^U to be a P-matrix. Using Lemma 4.2,

$$\begin{split} &\Upsilon^S \in \mathbb{R}^{M \times M} \text{ is a P-matrix} \\ \Leftrightarrow & \max_{j} \ l_j(\Upsilon^S \boldsymbol{l})_j > 0, \forall \boldsymbol{l} \neq \boldsymbol{0}, \boldsymbol{l} \in \mathbb{R}^M, \\ & i := \arg\max_{j} l_j(\Upsilon^C \boldsymbol{l})_j, \\ \Leftrightarrow & \boldsymbol{x}_i(\Upsilon^U_{\mathcal{P}_i,:} \cdot \boldsymbol{x})_{\mathcal{P}_i} \\ & \geq \ \|\Upsilon^U_{\mathcal{P}_i,\mathcal{P}_i}\| \cdot \|\boldsymbol{x}_i\|^2 - \sum_{j \neq i} \|\Upsilon^U_{\mathcal{P}_i,\mathcal{P}_j}\| \cdot \|\boldsymbol{x}_i\| \cdot \|\boldsymbol{x}_j\| \\ & \geq \ l_i \alpha^S_{ii} l_i - \sum_{j \neq i} l_i \beta^C_{ij} l_i = \ l_i (\Upsilon^C \boldsymbol{l})_i > 0, \\ \Rightarrow & \max_{k \in S_i} \ x_k (\Upsilon^U \boldsymbol{x})_k > 0, \\ \Rightarrow & \max_{k \in S_i} \ x_k (\Upsilon^U \boldsymbol{x})_k > 0, \\ \Leftrightarrow & \Upsilon^U \in \mathbb{R}^{N \times N} \text{ is a P-matrix} \end{split}$$

APPENDIX E PROOF OF COROLLARY 4.3.2

Proof. We can see that $I + \Gamma^S \succ \mathbf{0}$, iff $\lambda_1(\Gamma^S) > -1$, and from symmetry, $I + \Gamma^S$ is a P-matrix. The Υ^S matrix can be obtained by scaling the ith row of $I + \Gamma^S$ by α_i^S , which is a positive number. Therefore, the determinant of every principal minor of Υ^S has the same sign as the determinant of the corresponding principal minor of $I + \Gamma^S$, and thus Υ^S is also a P-matrix. By Theorem 4.3, the Nash equilibrium is unique and the sufficiency holds.

Appendix F Proof of Theorem 5.2

We prove for the K=1 case and the proof generates to an arbitrary K.

Proof. When we consider borderline inactive and active groups, the set of all their members are a superset of the union of the borderline inactive and active set of agents.

We denote

$$Y = \{a_k | a_k \in \mathcal{P}_i, s.t.\mathcal{P}_i \in A_S(\boldsymbol{x}^*) \cup B_S(\boldsymbol{x}^*)\}$$

as the set of all agents that belong to communities in $A_S(\mathbf{x}^*) \cup B_S(\mathbf{x}^*)$, then we have $A(\mathbf{x}^*) \cup B(\mathbf{x}^*) \subseteq \mathcal{S}$. Similar to Theorem

5.1, we denote $\nabla_{\mathcal{S}} F_{\mathcal{S}}(\boldsymbol{x}^*)$ as a sub-matrix of $\nabla F(\boldsymbol{x}^*)$ whose columns and rows correspond to the agents in set \mathcal{S} .

Our proof proceeds by the following logic:

$$H^{S}(\boldsymbol{x}^{*}) \succ \boldsymbol{0} \Rightarrow \nabla_{S}F_{S}(\boldsymbol{x}^{*}) \succ \boldsymbol{0} \Rightarrow \nabla_{A} {}_{B}F_{A} {}_{B}(\boldsymbol{x}^{*}) \succ \boldsymbol{0}.$$

For the first part, we adopt techniques similar to those we used in Appendix D. For $\forall \boldsymbol{v} \in \mathbb{R}^{|S|}_{\geq 0}$, we denote $\boldsymbol{l} = (l_i)_{i=1}^Z \in \mathbb{R}^Z_{\geq 0}$, where $l_i = \|\boldsymbol{v}_i\|_2, \forall i \in \mathbb{N}[1, Z]$. Then we have

$$\begin{split} & \boldsymbol{v}^T \nabla_Y F_{\mathcal{S}}(\boldsymbol{x}^*) \boldsymbol{v} \\ = & \sum_{i=1}^Z \sum_{j=1}^Z \boldsymbol{z}_j^T (\nabla_j F_i(\boldsymbol{x}^*)) \boldsymbol{v}_i \\ = & \sum_{i=1}^Z \boldsymbol{v}_i^T (\nabla_i F_i(\boldsymbol{x}^*)) \boldsymbol{v}_i + \sum_{i=1}^Z \sum_{j=1, j \neq i}^Z \boldsymbol{v}_j^T (\nabla_j F_i(\boldsymbol{x}^*)) \boldsymbol{v}_i \\ \geq & \sum_{i=1}^Z \theta_i^S \cdot \|\boldsymbol{v}_i\|_2^2 - \sum_{i=1}^Z \sum_{j=1, j \neq i}^Z \delta_{ij}^S \cdot \|\boldsymbol{v}_i\|_2 \cdot \|\boldsymbol{v}_j\|_2 \\ = & \sum_{i=1}^Z \theta_i^S \cdot l_i^2 - \sum_{i=1}^Z \sum_{j=1, j \neq i}^Z \delta_{ij}^S \cdot l_i \cdot l_j \\ = & \boldsymbol{l}^T H^S(\boldsymbol{x}^*) \boldsymbol{l} > 0 \end{split}$$

which shows $\nabla_{\mathcal{S}} F_{\mathcal{S}}(x^*) \succ \mathbf{0}$ and completes the proof of the first part.

For the second part, we know from the Sylvester's criterion that an $N\times N$ Hermitian matrix (for real valued matrices, it is symmetric) is positive definite if and only if every leading principal component of it (the top left $k\times k$ submatrices, for $k=1,\ldots,N$) has a positive determinant. Since $A(\boldsymbol{x}^*)\bigcup B(\boldsymbol{x}^*)\subseteq Y$, we know that $\nabla_{A,B}F_{A,B}(\boldsymbol{x}^*)$ is a leading principal minor of $\nabla_Y F_Y(\boldsymbol{x}^*)$ and thus $\nabla_{A,B}F_{A,B}(\boldsymbol{x}^*)\succ \mathbf{0}$, which completes the proof.

Appendix G Proof of Theorem 5.3

Proof. Since $\Upsilon^S \succ \mathbf{0}$, we know from Theorem 4.3 that the Nash equilibrium is unique. It remains to prove that this NE x^* is stable.

We denote $\Upsilon_{A,B}^S \in \mathbf{Z} \times \mathbf{Z}$ as the principal minor of Υ^S by picking out all the rows and columns corresponding to groups in $A_S(\mathbf{x}^*) \bigcup B_S(\mathbf{x}^*)$, and thus $\Upsilon_{A,B}^S \succ \mathbf{0}$.

Without loss of generality, suppose $\mathcal{P}_1,\mathcal{P}_2\in A_S(\boldsymbol{x}^*)\bigcup B_S(\boldsymbol{x}^*)$, then we know their new indices in $H^S(\boldsymbol{x}^*)$ remain unchanged. We know from the definition that $\alpha_1^S \leq \theta_1^S$ and $\beta_{12}^S \geq \delta_{12}^S$ (similar comparison can generalize to all other elements), and thus for $\forall \boldsymbol{v}\in\mathbf{R}^Z$, we have

$$v^{T}H^{S}(\boldsymbol{x}^{*})\boldsymbol{v}$$

$$= \sum_{i=1}^{Z} \sum_{j=1}^{Z} G_{ij}^{S}(\boldsymbol{x}^{*}) \cdot v_{i} \cdot v_{j}$$

$$= \sum_{i=1}^{Z} \theta_{i}^{S} \cdot v_{i}^{2} - \sum_{i=1}^{Z} \sum_{j=1, j \neq i}^{Z} \delta_{ij} \cdot v_{i} \cdot v_{j}$$

$$\geq \sum_{i=1}^{Z} \alpha_{i}^{S} \cdot v_{i}^{2} - \sum_{i=1}^{Z} \sum_{j=1, j \neq i}^{Z} \beta_{ij} \cdot v_{i} \cdot v_{j}$$

$$= \boldsymbol{v}^{T} \Upsilon_{A}^{S} \boldsymbol{N} > 0,$$

which shows $H^S(x^*) \succ \mathbf{0}$, and thus x^* is stable.

APPENDIX H PROOF OF THEOREM 6.1

Proof. First of all, when $e = 1, \forall t > 0, D_{max}^{out}(1) < 1$ implies

$$\alpha_i^S > \sum_{j \neq i} \beta_{ij}^S, \forall i = 1, \dots, M,$$

and this shows that the Υ^S matrix is diagonally row dominant, and thus is a P-matrix [33] (Generating method 4.1, this matrix is a *positively stable* P-matrix).

For an arbitrary $e \succ \mathbf{0}$, we define the following matrix E^S , where the ith column of E^S is equal to e_i times the ith column of Υ^S . Then since $e_i > 0$, every principal minor's determinant of E^S is equivalent to the corresponding principal minor's determinant of Υ^S and thus E^S is a P-matrix iff Υ^S is a P-matrix. Then, it's easy to see that $D_{max}^{out}(e) < 1$ implies the diagonal row dominance of E^S , which is sufficient to show E^S is a P-matrix. This completes the proof of the out-degree part and the in degree part is similar. Moreover, the stability result follows from Theorem 5.3 and the in and out degree parts of this theorem.

APPENDIX I COMPLEXITY OF CONDITION VERIFICATION IN THEOREM 4.3

Size	Unstructured	Structured	
10 × 10	1.08×10^{35}	1.26 ×10 ⁶	
20×20	1.39×10^{127}	1.69 ×10 ¹⁰	
50×50	4.90×10^{761}	6.34 ×10 ²⁰	

TABLE 4
Verification complexity in FLOPs of conditions in Theorem 4.3 over number of agents.

APPENDIX J SUPPLEMENTARY REASONING IN SECTION 7

We begin with the claim that the upper bounds are large enough for the study of sufficiency gaps.

When the parameters that control the external impact are all chosen at their normalized upper bound, the expected sum of the off-diagonal elements will be twice as high as the mean value of the internal impact. For our choices of $S_{low}^{in}, S_{high}^{in}, S_{low}^{out}, S_{high}^{out}$, the Υ^U will not be a P-matrix with probability 1 and thus Υ^S will not be a P-matrix. For example, we assume that Υ^U is symmetric, and for a matrix with positive diagonal elements and non-positive off-diagonal elements, as long as the sum of the absolute values of the off-diagonal elements is greater than the sum of the diagonal elements (the sum of all its elements is negative), then we know $\mathbf{1}^T\Upsilon^U\mathbf{1} < 0$ and it is not positive definite and thus not a P-matrix. Increasing such upper bounds will only increase the upper right area above the curve formed by dark cells.

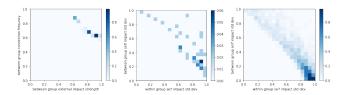
For the parameters that control the internal impact variances, the argument is similar, when the values of $c_{\mathcal{P}_i}^{high}-c_{\mathcal{P}_i}^{low}$ are $c^{high}-c^{low}$ close to 1, then the agent with the lowest internal impact will have an internal impact close to 0 and Υ^U will not be a P-matrix. Figures 18 and 19 shows the comparison after we

double such normalized upper bounds, and support our argument above.

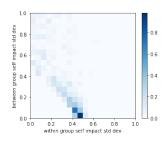
Next, we elaborate more on the "critical values" in Figures 11 and 12.

If agents have homogeneous internal impact, i.e., α_k^U are constant, then when $\mathbf{E}[\sum_{k \neq l} \beta_{kl}^U]$ is about 10% higher than $\sum_{k=1}^N \alpha_k^U$, the games are most likely to have sufficiency gaps. We discussed above that when Υ^U is symmetric, and $\sum_{k \neq l} \beta_{kl}^U > \sum_{k=1}^N \alpha_k^U$, then the unstructured condition is not satisfied, which is sufficient to conclude that the structured condition does not hold either. But when it's $\mathbf{E}[\sum_{k \neq l} \beta_{kl}^U] > \sum_{k=1}^N \alpha_k^U$, there are realizations (sample games under our choices of $S_{low}^{in}, S_{high}^{in}, S_{low}^{out}, S_{high}^{out})$ that have $\sum_{k \neq l} \beta_{kl}^U < \sum_{k=1}^N \alpha_k^U$ and thus are possible candidates for causing sufficiency gaps between the two sets of conditions. However, if $\mathbf{E}[\sum_{k \neq l} \beta_{kl}^U]$ is sufficiently larger (30% or more) $\sum_{k=1}^N \alpha_k^U$, all sample games will have neither set of conditions satisfied, which corresponds to the top right corner of the figures.

J.1 Sufficiency gap heatmaps for multipartite graphs



16. Sufficiency Fig. 17. Sufficiency Fig. 18. Sufficiency frequency over gap frequency over gap frequency over impact the between group the internal impact the internal external impact, fre-variances, with weak variances, with quency 0.83%. external impact, medium external frequency 0.18%. impact. frequency 6.31%.



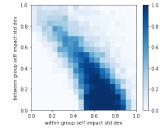


Fig. 19. Sufficiency gap frequency over the internal impact variances, with medium external impact, frequency 2.09%, doubled upper bounds.

Fig. 20. Sufficiency gap frequency over the internal impact variances, with strong external impact, frequency 26.93%.

J.2 Procedure for Synthetic Game Instance Generation

A game is generated for a specified size (number of agents N, number of action dimensions K=1, number of partitions M, and size of each partition N_i) and using utility functions $u_i(\boldsymbol{x},G)=x_i(b_i+\sum_j g_{ij}x_j)-\frac{c_i}{2}x_i^2$. The rest of the game is given by the interaction matrix G, and the vectors $\mathbf{b}=(b_i)_{i=1}^N$ and $\mathbf{c}=(c_i)_{i=1}^N$. In generating a random G, the diagonal elements are set to 0 without loss of generality. Each partition is associated with consecutive agent indices and thus the diagonal blocks

represent each partition and the off-diagonal blocks represent cross-partition interdependencies. The off-diagonal elements in a diagonal block are generated using a Bernoulli distribution with parameter P_{exist}^{in} , the probability for an connection (non-zero element) to exist between any pair of agent-action indices. If such a connection exists, its value (strength of the connection), g_{ij} , is drawn from a uniform distribution on the interval $[S_{low}^{in}, S_{high}^{in}]$. We generate g_{ij} and g_{ji} independently. These elements also determine the off-diagonal elements in the diagonal blocks of Υ^U and the diagonal elements α_i^S in the Υ^S matrix.

The off-diagonal blocks of G are generated using the same approach, with parameters P_{exist}^{out} and $[S_{low}^{out}, S_{high}^{out}]$. These determine the connection frequencies and strengths between groups. These elements also determine the off-diagonal blocks of the Υ^U matrix as well as the off-diagonal elements $\beta_{i,j}^S$ in the Υ^S matrix.

matrix as well as the off-diagonal elements β^S_{ij} in the Υ^S matrix. In subsequent numerical results, a *dense* matrix G refers to both $P^{in}_{exist}=1$ and $P^{out}_{exist}=1$, i.e., the entries in Υ^U and Υ^S are non-zero with probability 1.

The vector of individual cost c is generated by first choosing a fixed mean value $\overline{c}=\frac{1}{2}$ (so that $\mathbb{E}[\alpha_k^U]=1$) for all the cost terms. We then generate a partition mean $\overline{c}_{\mathcal{P}_i}$ by sampling uniformly at random from an interval $[c^{low},c^{high}]$, where $c^{low}+c^{high}=1$. Next, within each group, we choose an interval $[c^{low}_{\mathcal{P}_i},c^{high}_{\mathcal{P}_i}]$, where $(c^{low}_{\mathcal{P}_i}+c^{high}_{\mathcal{P}_i})/2=\overline{c}_{\mathcal{P}_i}$ and then sample uniformly at random the individual cost terms from this. All individual cost terms are set to strictly positive values, otherwise neither structured nor unstructured conditions hold, making the game instances trivial. The individual benefit is set to b=1 as it does not affect either Υ^U or Υ^S . In generating these values we fix the global mean but change the variance.

J.3 Sufficiency gap on email-Eu-core network data

We next perform a similar set of experiments using a real-world directed graph. This dataset is from the email-Eu-core network [41] and provides a binary directed graph with edges indicating whether an email was ever sent from one node to another, where nodes represent people from a large European research institution and each node has a department affiliation. We will use this graph to run game simulations and test the identified conditions on this network's structured and unstructured Υ matrices. In doing so, we add an edge weight ϕ to all the edges; this is shown as the x-axis "external connection strength" in the figures below. We also need to equip the game with a utility function (detailed shortly).

This graph is highly asymmetric. As discussed in Section 4, the conditions on asymmetric Υ matrices is much more computationally costly to verify than the conditions on symmetric Γ matrices of the same size. As a result, verifying the conditions on the entire network with 1005 nodes leads to memory overflow. Consequently, our experiments are on down-sampled versions of this graph. We employ the following three types of network sampling methods to generate sub-graphs for our experiments:

- Sample M departments (groups) uniformly at random with all nodes in each;
- Sample N_{sample} nodes uniformly at random;
- Sample M departments (groups) with at least M members uniformly at random first, and then sample M nodes from each department.

We will use the following utility function of a node a_i (an agent) given the sampled network:

$$u_i(x_i, \boldsymbol{x}_{-i}) = \frac{\phi}{\overline{D}} \sum_{j \neq i} (G_{sample})_{ij} x_i x_j - \frac{1}{2} x_i^2,$$

where $\overline{D}=\frac{\sum_{i=1}^{N_{sample}}D_{i}^{in}+D_{i}^{out}}{2N_{sample}}$, N_{sample} is the number of sampled nodes, and G_{sample} is the corresponding sampled submatrix.

We want to compare different ways of partitioning the agentaction space. Since each node is associated with a department, we can naturally partition the network in terms of departments; we will call this the *original partition*. We can also create partitions on the nodes using the following criteria:

- Degree-based: rank the nodes by the average of the in-degree and out-degree of a node, and then put nodes with similar degrees in the same partition;
- Connection-type-based: rank the nodes by the fraction of their within-department connections (number of connections in the department over the node's total number of connections), and create partitions that include nodes with similar fraction in the same partition.

As results in Figures 21-29 show, degree-based partition results in significantly lower sufficiency gap and clearer patterns on when the structured conditions fail compared to the original partition. In contrast, the connection-type-based partition results in higher sufficiency gaps and thus is not ideal. The average sufficiency gap is given for each figure in the caption. We can see from the numerical results that when we re-group the nodes based on their degree information, not only the overall sufficiency gap is lower, but the sufficiency gaps are also less sensitive to the network size and only sensitive to the relevant internal/external impact strengths.

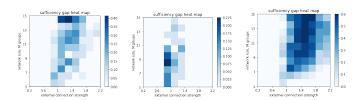


Fig. 21. Sufficiency gap when sampling M departments, original partition, 6.22%.

22. Sufficiency Fig. gap when sampling departments, degree-based partition, 1.84%.

Fig. 23. Sufficiency gap when sampling Mdepartments, connection-typebased partition, 20.40%.

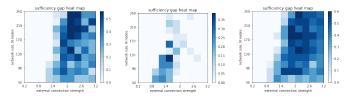
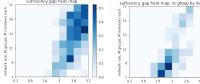


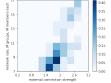
Fig. 24. Sufficiency gap when sampling Nnodes, original partition, 21.39%.

Fig. 25. Sufficiency gap when sampling Nnodes, degree-based partition, 3.73%.

26. Sufficiency gap when sampling N nodes, connectiontype-based partition, 28.63%.

Intuitively, these observations can be explained as follows. The high degree nodes in our sample network are frequently connected





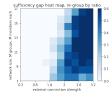


Fig. 27. Sufficiency gap when sampling M departments, each with ${\cal M}$ nodes, original partition, 9.94%.

28. Sufficiency Fig. gap when sampling Mdepartments, each with M nodes. degree-based partition, 3.49%.

Fig. 29. Sufficiency when sampling gap Mdepartments, each with M nodes. connection-typebased partition, 23.06%

with each other. Therefore, the degree-based partition perform best, as they successfully identify and retain the importance of such strongly connected sub-graphs in our reduced conditions. Moreover, we note that the singleton nodes who have no connections to others are put in the same partition. From the previous analysis, this means that the structured and unstructured conditions hold if and only if the corresponding conditions hold on the nonsingleton nodes' sub-graph.

J.4 Sufficiency gap heatmaps for stability conditions

We show the sufficiency gap experiment results for stability condition verification in Theorem 4.3, which show very similar trends as the uniqueness condition verification in Theorem 5.2. We note that using the GEDC for stability condition verification is equivalent in terms of computation complexity and sufficiency gaps.

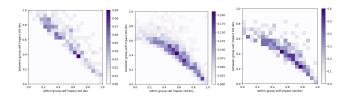
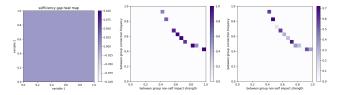


Fig. 30. Stability Fig. 31. Stability con- Fig. Stability 32. condition sufficiency dition sufficiency gap condition sufficiency frequency over frequency over the gap frequency over internal impact internal impact vari- the internal impact variances, with weak ances, with medium variances, with strong external impact, external impact, fre-external impact. frequency 0.55%. quency 1.97%. frequency 5.50%.



Stability Fig. 34. Stability con- Fig. Fig. 35. condition sufficiency dition sufficiency gap condition sufficiency gap frequency over frequency over the gap frequency over the internal impact internal impact vari- the internal impact variances, with overly ances, with medium variances, with strong strong external external impact, fre-external impact. and game quency 1.90%. impact frequency 1.19%. known to be unstable. frequency 0.0%.