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Efficient Estimation of Equilibria in Large Aggregative Games with Coupling Constraints

Paulin Jacquot, Cheng Wan, Olivier Beaudé, Nadia Oudjane

Abstract—Aggregative games have many industrial applications, and computing an equilibrium in those games is challenging when the number of players is large. In the framework of atomic aggregative games with coupling constraints, we show that variational Nash equilibria of a large aggregative game can be approximated by a Wardrop equilibrium of an auxiliary population game of smaller dimension. Each population of this auxiliary game corresponds to a group of atomic players of the initial large game. This approach enables an efficient computation of an approximated equilibrium, as the variational inequality characterizing the Wardrop equilibrium is of smaller dimension than the initial one. This is illustrated on an example in the smart grid context.

Index Terms—Aggregative Game - Variational Nash Equilibrium - Variational Inequalities - Population Game

I. INTRODUCTION

a) Motivation: Aggregative games [1] form a class of noncooperative games [2], finding practical applications—in particular for the subclass of splittable congestion games [3]—in various fields as traffic management [4], communications [5], [6] and electrical systems [7], [8].

The notion of Nash equilibrium (NE) [9] has emerged as the central solution concept in game theory. However, the computation of an NE is considered a challenging problem: for instance, [10] showed that, in finite two-players matricial games, computing a pure NE is PPAD-complete, while [11] showed that computing an NE in splittable congestion games with *player-specific affine* latency functions is also PPAD-complete. In continuous games, NEs can be characterized as solutions of Variational Inequalities (VI) [12], a characterization we adopt through this work. The efficiency of NEs computation depends on the dimension of these VIs, given by the number of players and of constraints. In some applications (transport, energy, etc), one may consider several thousands of heterogeneous agents. At this scale, computing an NE can be intractable. In the case where the game model involves coupling constraints between agents' actions, the concept generalizing the NE is simply referred to as generalized Nash equilibria (GNE) [13], and can be even harder to compute. The consideration of coupling constraints has emerged from applications: for instance to model the capacity of roads and communication channels or, in energy, to model the maximal variations of production plants [14]. Estimating an NE as the probable outcome of the model is of main importance for the operator of a system: for instance, in order to control some physical or managerial parameters and optimize the system (e.g. in bi-level programming [15]).

In this paper, we consider aggregative games with a finite but *large* number of players. We propose a method to compute an approximation of variational Nash equilibria (VNEs) [13], a subset of GNEs in

the presence of coupling constraints. The main idea of the proposed method is to reduce the dimension of the VIs characterizing VNEs: players are divided into groups with similar characteristics. Then, each group is replaced by a homogeneous population of nonatomic players. Last, we compute a variational Wardrop equilibrium (VWE) [16] in the obtained approximating nonatomic population game. The quality of the NE estimation depends on how well the characteristics (action sets and cost functions) of each homogeneous population approximate those of the atomic players it replaces.

b) Related works: The relation between NEs and WEs has been studied in different contexts: [17] consider network congestion game with each origin-destination pair having n players and show the convergence of NEs to a WE when n goes to infinity. [18] generalizes this result in the framework of composite games where nonatomic players and atomic splittable players coexist.

A work particularly in line with the present paper is [19] where the authors show, using a variational inequalities approach, that the distance between a VNE and a VWE in aggregative games converges to zero when the number of players tends to infinity. Their VWE corresponds to an equilibrium of the game where each atomic player is replaced by a population, as done in Sec. III-A of this paper. Their quantitative results show that the self impact of each player on the aggregate action is negligible when the number of players is large. Our paper shows a different result: VNEs are approximated by a VWE of a game with a *reduced dimension*. To this end, Sec. III-A extends the results of [19] to the subdifferentiable case. In [20], Jacquot and Wan show that, in congestion games with a continuum of heterogeneous players, the VWE can be approximated by a VNE of an approximating game with a finite number of players. In [21], those results are extended to aggregative games with nonsmooth cost functions. The approach developed in the present paper is actually the inverse of the one taken in [20] and [21]: here, the VWE in the auxiliary game serves as an approximation of a VNE of the original large game.

c) Main contributions:

- we define an approximating population game (Sec. III-B), with smaller dimension but close enough to the original large game—quantified through the Hausdorff distance between action sets and between subgradients of players' objective functions;
- we show that a particular VWE of the approximating reduced-dimension population game is close to any VNE of the original game with or without coupling constraints, and we provide an explicit expression of this approximation error (Thm. 2);
- we show that VNEs are close to VWEs in large aggregative games Thm. 1, extending [19, Thm. 1] to subdifferentiable cost functions and general *aggregatively monotone* games;
- we provide a numerical illustration of our results (Sec. IV)

based on a practical application: the decentralized charging of electric vehicles through a demand response mechanism [22]. This example illustrates the nondifferentiable case through piece-wise linear functions (“block rates tariffs”) and coupling constraints (capacities and limited variations). This example shows that the proposed method is implementable and reduces the time needed to compute an equilibrium by computing its approximation (six times faster for an approximation with a normalized distance to the actual equilibrium

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of less than 2%).

d) *Structure*: The remainder of this paper is organized as follows: Sec. II specifies the framework of aggregative games with coupling constraints, and recalls the notions of variational equilibria and monotonicity for VIs, as well as several results on the existence and uniqueness of equilibria. Sec. III formulates the main results: Sec. III-A shows that a VWE approximates VNEs in large games and then, Sec. III-B formulates the approximating population game with the approximation measures, and gives an upper bound on the distance between the VWE profile of the approximating game to an original VNE profile. Sec. IV presents a numerical illustration in the context of electricity demand response.

II. AGGREGATIVE GAMES WITH COUPLING CONSTRAINTS

A. Model and equilibria

We consider the model of *atomic aggregative* games: the cost function of each player depends only on her own action and on the *aggregate action*. The term *atomic* is opposed to the *nonatomic* model where players are infinitesimal and have a negligible weight [2]. As the aggregate action of players is considered in this model, all players have their feasible action sets lying in the same space \mathbb{R}^T with $T \in \mathbb{N}^*$ fixed. The formal definition is as follows.

Definition 1: An *atomic aggregative game with coupling constraints* $\mathcal{G}(\mathcal{A})$ is defined by a finite set of players $\mathcal{N} \stackrel{\text{def}}{=} \{1, \dots, n, \dots, N\}$ and for each player $n \in \mathcal{N}$:

- a set of feasible actions $\mathcal{X}_n \subset \mathbb{R}^T$, with a fixed $T \in \mathbb{N}^*$;
- an individual cost function $f_n : \mathcal{X}_n \times \mathcal{A} \rightarrow \mathbb{R}$;
- a set $\mathcal{A} \subset \mathbb{R}^T$ defining the *aggregative coupling constraint*: $\bar{\mathbf{X}} \in \mathcal{A}$ where $\bar{\mathbf{X}} \stackrel{\text{def}}{=} \frac{1}{N} \sum_n \mathbf{x}_n$ denotes the average action for individual actions of players $(\mathbf{x}_n)_{n \in \mathcal{N}}$.

The game is written as the tuple $\mathcal{G}(\mathcal{A}) \stackrel{\text{def}}{=} (\mathcal{N}, \mathcal{X}, (f_n)_{n \in \mathcal{N}}, \mathcal{A})$.

We denote by $\mathcal{X} \stackrel{\text{def}}{=} \mathcal{X}_1 \times \dots \times \mathcal{X}_N$ the product set of action profiles, and the set of feasible average profiles by:

$$\bar{\mathcal{X}} \stackrel{\text{def}}{=} \{ \bar{\mathbf{X}} \in \mathbb{R}^T : \forall n \in \mathcal{N}, \exists \mathbf{x}_n \in \mathcal{X}_n \text{ s.t. } \frac{1}{N} \sum_{n \in \mathcal{N}} \mathbf{x}_n = \bar{\mathbf{X}} \}.$$

Last, we denote the set of feasible action profiles satisfying the coupling constraint \mathcal{A} by:

$$\mathcal{X}(\mathcal{A}) \stackrel{\text{def}}{=} \{ \mathbf{x} \in \mathcal{X} : \frac{1}{N} \sum_{n \in \mathcal{N}} \mathbf{x}_n \in \bar{\mathcal{X}} \cap \mathcal{A} \}.$$

A particular class of aggregative games with many applications is given by *splittable congestion games* [3]:

Definition 2: A *splittable congestion game* is defined by a set of resources \mathcal{T} , each resource t having a cost function $c_t \in \mathbb{R}^{\mathbb{R}^+}$, and a set of player \mathcal{N} with actions set $(\mathcal{X}_n)_{n \in \mathcal{N}} \subset (\mathbb{R}_+^T)^{\mathcal{N}}$, and with individual utility function $u_n \in \mathbb{R}^{\mathcal{X}_n}$. The cost function $f_n : \mathcal{X}_n \times \bar{\mathcal{X}} \rightarrow \mathbb{R}$ of player $n \in \mathcal{N}$ is defined as:

$$\forall \mathbf{x}_n, \bar{\mathbf{Y}} \in \mathcal{X}_n \times \bar{\mathcal{X}}, f_n(\mathbf{x}_n, \bar{\mathbf{Y}}) \stackrel{\text{def}}{=} \sum_{t \in \mathcal{T}} x_{n,t} c_t(\bar{Y}_t) - u_n(\mathbf{x}_n). \quad (1)$$

As the action of n appears both in the first and second argument (through $\bar{\mathbf{X}}$) of her cost function f_n , we consider for each $n \in \mathcal{N}$, the *modified* cost function:

$$\hat{f}_n(\mathbf{x}_n, \bar{\mathbf{Y}}_{-n}) \stackrel{\text{def}}{=} f_n(\mathbf{x}_n, \bar{\mathbf{Y}}_{-n} + \frac{1}{N} \mathbf{x}_n)$$

for $\mathbf{x}_n \in \mathcal{X}_n$ and $\bar{\mathbf{Y}}_{-n} \in \bar{\mathcal{X}}_{-n} \stackrel{\text{def}}{=} \{ \frac{1}{N} \sum_{m \neq n} \mathbf{x}_m : \mathbf{x}_m \in \mathcal{X}_m, \forall m \}$.

In this paper, we adopt the following standard assumptions:

Assumption 1:

- for each player $n \in \mathcal{N}$, the set \mathcal{X}_n is a convex and compact subset of \mathbb{R}^T with nonempty relative interior;

- for each $n \in \mathcal{N}$, the cost function f_n and modified cost function \hat{f}_n are convex in the first variable;

- for each $n \in \mathcal{N}$, the cost function $\mathbf{x}_n, \bar{\mathbf{X}} \mapsto f_n(\mathbf{x}_n, \bar{\mathbf{X}})$ is Lipschitz continuous in $\bar{\mathbf{X}}$, with Lipschitz constant L_2 independent of n and \mathbf{x}_n ;

- \mathcal{A} is a convex closed set of \mathbb{R}^T , and $\bar{\mathcal{X}} \cap \mathcal{A}$ is not empty.

Let us consider the subgradient correspondences $\hat{H} : \mathcal{X} \rightrightarrows \mathbb{R}^{NT}$ and $H : \mathcal{X} \rightrightarrows \mathbb{R}^{NT}$, defined for any $\mathbf{x} \in \mathcal{X}$ as:

$$\begin{aligned} \hat{H}(\mathbf{x}) &\stackrel{\text{def}}{=} \{ (\mathbf{h}_n)_{n \in \mathcal{N}} \in \mathbb{R}^{NT} : \mathbf{h}_n \in \partial_1 \hat{f}_n(\mathbf{x}_n, \bar{\mathbf{X}}_{-n}), \forall n \in \mathcal{N} \} \\ &= \prod_{n \in \mathcal{N}} \partial_1 \hat{f}_n(\mathbf{x}_n, \bar{\mathbf{X}}_{-n}); \\ H(\mathbf{x}) &\stackrel{\text{def}}{=} \{ (\mathbf{h}_n)_{n \in \mathcal{N}} \in \mathbb{R}^{NT} : \mathbf{h}_n \in \partial_1 f_n(\mathbf{x}_n, \bar{\mathbf{X}}), \forall n \in \mathcal{N} \} \\ &= \prod_{n \in \mathcal{N}} \partial_1 f_n(\mathbf{x}_n, \bar{\mathbf{X}}), \end{aligned}$$

where ∂_1 denotes the partial differential w.r.t. the first variable of the function. The interpretation of $\hat{H}(\mathbf{x})$ is clear: \mathbf{h}_n is a subgradient of player n 's utility function \hat{f}_n w.r.t. her action \mathbf{x}_n . Let us leave the interpretation of $H(\mathbf{x})$ until Def. 4. For the moment, let us establish a relation between \hat{H} and H :

Lemma 1: For each $\mathbf{x} \in \mathcal{X}$, $\mathbf{h} \in \hat{H}(\mathbf{x})$ if and only if there exists $\mathbf{g}_1 \in \prod_{n \in \mathcal{N}} \partial_1 f_n(\mathbf{x}_n, \bar{\mathbf{X}}) = H(\mathbf{x})$ and $\mathbf{g}_2 \in \prod_{n \in \mathcal{N}} \partial_2 f_n(\mathbf{x}_n, \bar{\mathbf{X}})$ such that $\mathbf{h} = \mathbf{g}_1 + \frac{1}{N} \mathbf{g}_2$.

Proof: The proof can be obtained from [23, Prop. 16.6 and 16.7]. Details are omitted for brevity. ■

In the presence of coupling constraints, the notion of Nash equilibrium (NE) [9] is replaced by the one of Generalized Nash Equilibrium (GNE). A profile $\mathbf{x} \in \mathcal{X}(\mathcal{A})$ is a GNE if, for each player n :

$$\hat{f}_n(\mathbf{x}_n, \bar{\mathbf{X}}_{-n}) \leq \hat{f}_n(\mathbf{y}_n, \bar{\mathbf{X}}_{-n}), \forall \mathbf{y}_n \text{ s.t. } (\mathbf{y}_n/N + \bar{\mathbf{X}}_{-n}) \in \bar{\mathcal{X}} \cap \mathcal{A}.$$

For atomic games, Variational Nash Equilibria (VNE) [13], [24] form a special class of GNEs with symmetric properties: in some sense, the burden of constraint $\bar{\mathbf{X}} \in \mathcal{A}$ is shared symmetrically by players [13]. VNEs, in the subdifferentiable case, are characterized as the solution of a generalized VI (GVI) (2) stated below.

Definition 3 (Variational Nash Equilibrium (VNE), [13]): A VNE is a solution $\hat{\mathbf{x}} \in \mathcal{X}(\mathcal{A})$ to the following GVI problem:

$$\exists \mathbf{g} \in \hat{H}(\hat{\mathbf{x}}) \text{ s.t. } \langle \mathbf{g}, \mathbf{x} - \hat{\mathbf{x}} \rangle \geq 0, \forall \mathbf{x} \in \mathcal{X}(\mathcal{A}). \quad (2)$$

In particular, if $\bar{\mathcal{X}} \subset \mathcal{A}$, a VNE is an NE. We refer to [25] for further discussions on VNE and VI characterization. In this paper, we adopt the notion of VNE as the equilibrium concept in the presence of aggregate constraints.

As the first step of approximation, we define a *nonatomic* aggregative game $\mathcal{G}(\mathcal{A})'$ associated to $\mathcal{G}(\mathcal{A})$, where each player n is replaced by a continuum of identical nonatomic players. The set of nonatomic players is represented by the interval $[0, 1]$. Each player in population n has action set \mathcal{X}_n and cost function f_n . In nonatomic games, the concept of VNE is replaced by the concept of variational Wardrop equilibrium (VWE). We refer to [21] for the formal definition of a VWE and further discussions. In general, a VWE is characterized by an infinite dimensional VI. Here, we only consider *symmetric* VWE, characterized by a VI of finite dimension as defined below:

Definition 4: A *symmetric variational Wardrop equilibrium* (SVWE) of $\mathcal{G}(\mathcal{A})'$ is a solution $\mathbf{x}^* \in \mathcal{X}(\mathcal{A})$ to the GVI:

$$\exists \mathbf{g} \in H(\mathbf{x}^*) \text{ s.t. } \langle \mathbf{g}, \mathbf{x} - \mathbf{x}^* \rangle \geq 0, \forall \mathbf{x} \in \mathcal{X}(\mathcal{A}). \quad (3)$$

where, for each $n \in \mathcal{N}$, \mathbf{x}_n^* is the common action adopted by all nonatomic players in population n .

In addition to the population game interpretation, an SVWE can be interpreted as an equilibrium concept in the initial game, where the contribution of each atomic player on the aggregate action $\bar{\mathbf{X}}$ is negligible [19]. However, in this interpretation, an SVWE is not

stable, as the individual SVWE profile \mathbf{x}_n^* of a player n does not necessarily minimize her cost function $\hat{f}_n(\cdot, \bar{\mathbf{X}}_{-n})$.

The existence of an SVWE in the population game $\mathcal{G}(\mathcal{A})'$, with players identical in the same population, is given in Prop. 1.

Proposition 1 (Existence of equilibria): Under Asm. 1, $\mathcal{G}(\mathcal{A})$ (resp. $\mathcal{G}(\mathcal{A})'$) admits a VNE (resp. SVWE).

Proof: From [26, Proposition 8.7], we obtain that \hat{H} and H are nonempty, convex, compact valued, upper hemicontinuous correspondences. Then [27, Corollary 3.1] shows that the GVI problems (2) and (3) both admit a solution on the finite dimensional convex compact $\mathcal{X}(\mathcal{A})$. ■

Before discussing the uniqueness of equilibria, let us recall some relevant definition of monotonicity for correspondences:

Definition 5: A correspondence $\Gamma : \mathcal{X} \rightrightarrows \mathbb{R}^{NT}$ is said:

i) *monotone* if for all $\mathbf{x}, \mathbf{y} \in \mathcal{X}, \mathbf{g} \in \Gamma(\mathbf{x}), \mathbf{h} \in \Gamma(\mathbf{y})$:

$$\sum_{n \in \mathcal{N}} \langle \mathbf{g}_n - \mathbf{h}_n, \mathbf{x}_n - \mathbf{y}_n \rangle \geq 0 ;$$

ii) *strictly monotone* if the equality holds iff $\mathbf{x} = \mathbf{y}$;

iii) *aggregatively strictly monotone* if equality holds iff $\sum_n \mathbf{x}_n = \sum_n \mathbf{y}_n$;

iv) α -*strongly monotone* if $\alpha > 0$ and, for all $\mathbf{x}, \mathbf{y} \in \mathcal{X}$:

$$\sum_n \langle \mathbf{g}_n - \mathbf{h}_n, \mathbf{x}_n - \mathbf{y}_n \rangle \geq \alpha \|\mathbf{x} - \mathbf{y}\|^2, \forall \mathbf{g} \in \Gamma(\mathbf{x}), \mathbf{h} \in \Gamma(\mathbf{y});$$

v) β -*aggregatively strongly monotone* on \mathcal{X} if $\beta > 0$ and, for all $\mathbf{x}, \mathbf{y} \in \mathcal{X}$ with $\bar{\mathbf{X}} = \frac{1}{N} \sum_n \mathbf{x}_n, \bar{\mathbf{Y}} = \frac{1}{N} \sum_n \mathbf{y}_n$:

$$\sum_n \langle \mathbf{g}_n - \mathbf{h}_n, \mathbf{x}_n - \mathbf{y}_n \rangle \geq N\beta \|\bar{\mathbf{X}} - \bar{\mathbf{Y}}\|^2, \forall \mathbf{g} \in \Gamma(\mathbf{x}), \mathbf{h} \in \Gamma(\mathbf{y}).$$

If $T = 1$, “monotone” corresponds to “increasing”. Besides, (aggregatively) strict monotonicity implies monotonicity, while strong (resp. aggregatively strong) monotonicity implies strict (resp. aggregatively strict) monotonicity.

Prop. 2 recalls some results on the uniqueness of VNE and SVWE, according to the monotonicity of \hat{H} and H :

Proposition 2 (Uniqueness of equilibria): Under Asm. 1:

i) if \hat{H} (resp. H) is strictly monotone, then $\mathcal{G}(\mathcal{A})$ (resp. $\mathcal{G}(\mathcal{A})'$) has a unique VNE (resp. SVWE);

ii) if \hat{H} (resp. H) is aggregatively strictly monotone, then all VNE (resp. SVWE) of $\mathcal{G}(\mathcal{A})$ (resp. $\mathcal{G}(\mathcal{A})'$) have the same aggregate profile.

Proof: The proof can be adapted from [12, Chapter 2].

As opposed to the monotonicity of H which happens for instance in congestion games with convex cost of resources [19, Lemma 3], \hat{H} is rarely monotone (except in some particular cases, e.g. with \mathbf{c} linear [3], [28]). In view of Prop. 2, the absence of monotonicity of \hat{H} can result in multiple VNEs [29]. However, the next section shows that, with a large number of players, those VNEs are close to each other, and are well approximated by an SVWE.

III. APPROXIMATING VNES OF A LARGE GAME

A. Considering SVWE instead of VNE

Let us denote by $\mathcal{X}_0 \subset \mathbb{R}^T$ the convex closed hull of $\bigcup_{n \in \mathcal{N}} \mathcal{X}_n$, and consider the radius:

$$R \stackrel{\text{def}}{=} \max_{\mathbf{x} \in \mathcal{X}_0} \|\mathbf{x}\| .$$

Thm. 1 gives the first step of approximation of VNEs, by bounding the distance between a VNE and an SVWE.

Theorem 1 (SVWE close to VNE): Under Asm. 1, let $\mathbf{x} \in \mathcal{X}(\mathcal{A})$ be a VNE of $\mathcal{G}(\mathcal{A})$ and $\mathbf{x}^* \in \mathcal{X}(\mathcal{A})$ an SVWE of $\mathcal{G}(\mathcal{A})'$:

i) if H is α -strongly monotone, then \mathbf{x}^* is unique and:

$$\|\mathbf{x} - \mathbf{x}^*\| \leq \frac{L_2}{\alpha} \frac{1}{\sqrt{N}}, \quad \text{with } \alpha \stackrel{\text{def}}{=} \min_{n \in \mathcal{N}} \alpha_n, \text{ besides} \quad (4)$$

$$\frac{1}{N} \sum_n \|\mathbf{x}_n - \mathbf{x}_n^*\| \leq \frac{L_2}{\alpha N} \quad \text{and} \quad \|\bar{\mathbf{X}} - \bar{\mathbf{X}}^*\| \leq \frac{L_2}{\alpha N}, \quad (5)$$

where L_2 is the Lipschitz constant defined in Asm. 1;

ii) if H is β -aggr. str. monotone, then \mathbf{X}^* is unique and:

$$\|\bar{\mathbf{X}} - \bar{\mathbf{X}}^*\| \leq \sqrt{\frac{2RL_2}{\beta N}}, \quad \text{with } \beta \stackrel{\text{def}}{=} \min_{t \in \mathcal{T}} \beta_t . \quad (6)$$

Proof: The proof for i) can be adapted from [19, Th. 1.2)] to the subdifferential case, by relying on Lemma 1. The proof for ii) generalizes [19, Th. 1.3] for *aggr. monotone* games. ■

While VNEs are not unique in general (see Prop. 2), by applying the triangle inequality to the results of Thm. 1 we observe that all VNEs are close to each other when N is large.

Thm. 1 shows that an (average) SVWE approximates an (average) VNE of $\mathcal{G}(\mathcal{A})$ when the number of players is large, but this does not reduce the dimension of the GVI to resolve, as the GVI characterizing the VNE and those characterizing SVWE have the same dimension. For this reason, the second step of approximation consists in regrouping the similar populations.

B. Classification of populations

In this subsection, we shall regroup the populations of the game $\mathcal{G}(\mathcal{A})'$ having similar strategy sets \mathcal{X}_n and utility subgradients $\partial(-u_n)$, into larger populations, endow them with a common strategy set and a common utility function, so that the SVWE of this new population game approximates the SVWE of $\mathcal{G}(\mathcal{A})'$. The similarity between two sets \mathcal{X} and \mathcal{Y} is measured through the Hausdorff distance:

$$d_H(\mathcal{X}, \mathcal{Y}) \stackrel{\text{def}}{=} \max \left(\max_{\mathbf{x} \in \mathcal{X}} d(\mathbf{x}, \mathcal{Y}), \max_{\mathbf{y} \in \mathcal{Y}} d(\mathbf{y}, \mathcal{X}) \right) .$$

Let us define the compact set $\mathcal{M} \stackrel{\text{def}}{=} \mathcal{X}_0 + B_0(\delta)$, where $B_0(\delta)$ is the ball centered at 0 and of radius δ , and $\delta \geq \max_{n, m \in \mathcal{N}} d_H(\mathcal{X}_n, \mathcal{X}_m)$. W.l.o.g, we assume that for each $n \in \mathcal{N}$, f_n can be extended to a neighborhood of \mathcal{M}^2 , and is bounded on \mathcal{M}^2 . We make the additional assumption:

Assumption 2: For each $n \in \mathcal{N}$ and each $\bar{\mathbf{X}} \in \bar{\mathcal{X}}$, $f_n(\cdot, \bar{\mathbf{X}})$ is Lipschitz continuous with constant L_{1n} , and $L_1 \stackrel{\text{def}}{=} \max_n L_{1n} < \infty$.

At the SVWE of the game with reduced dimension, all the nonatomic players in the same population play the same action, by definition of an SVWE. Therefore, in order for this new SVWE to well approximate the SVWE in $\mathcal{G}(\mathcal{A})'$, a condition is that similar populations in $\mathcal{G}(\mathcal{A})'$ do play similar actions at the SVWE of $\mathcal{G}(\mathcal{A})'$: this is obtained easily in the case without coupling constraint, as shown in Prop. 3.

Proposition 3: Under Asm. 1, let $\mathbf{x}^* \in \mathcal{X}$ be an SVWE of $\mathcal{G}'(\mathcal{A})$ with $\mathcal{A} = \mathbb{R}^T$ (no coupling constraint). For populations $n, m \in \mathcal{N}$, if $f_n(\cdot, \bar{\mathbf{X}}^*)$ is α_n -strongly convex, $d_H(\mathcal{X}_n, \mathcal{X}_m) \leq \delta$, and if:

$$\sup_{\mathbf{g}_m \in \partial_1 f_m(\mathbf{x}_m^*, \bar{\mathbf{X}}^*)} d(\mathbf{g}_m, \partial_1 f_n(\mathbf{x}_m^*, \bar{\mathbf{X}}^*)) \leq d$$

then: $\|\mathbf{x}_n^* - \mathbf{x}_m^*\|^2 \leq \frac{1}{\alpha_n} ((L_{1n} + L_{1m})d + 2\delta R)$.

Proof: The proof is obtained using strong convexity, the definition of d , and Lipschitz convexity, following the same scheme as for Thm. 2. ■

In the case with coupling constraints, the proof for a similar result is more complicated, and we leave it to Cor. 1.

Let us now present the regrouping procedure. Denote by $\tilde{\mathcal{G}}^{\mathcal{I}}(\mathcal{A})$ an auxiliary game, with a set \mathcal{I} of p populations. Each population $i \in \mathcal{I}$ corresponds to a subset \mathcal{N}_i of populations in the game $\mathcal{G}(\mathcal{A})'$, such that $\bigcup_{i \in \mathcal{I}} \mathcal{N}_i = \mathcal{N}$ and for any $i, j \in \mathcal{I}, \mathcal{N}_i \cap \mathcal{N}_j = \emptyset$, i.e. $(\mathcal{N}_i)_{i \in \mathcal{I}}$ forms a partition of \mathcal{N} . Denote $p_i = |\mathcal{N}_i|$ the number of original populations now included in i . By abuse of notations, let i also denote the interval $[0, p_i]$, so that each nonatomic player in population i is represented by a point $\theta \in [0, p_i]$. The common action set of each nonatomic player in i is a compact convex subset

of \mathbb{R}^T , denoted by \mathcal{X}_i . Each player θ in each population i playing action \mathbf{x}_θ , let

$$\bar{\mathbf{X}} \stackrel{\text{def}}{=} \frac{1}{N} \sum_{i \in \mathcal{I}} \int_{\theta \in i} \mathbf{x}_\theta d\theta$$

denotes the aggregate action profile. The set of aggregate action-profile is a subset of \mathbb{R}^T given as:

$$\bar{\mathcal{X}}^{\mathcal{I}} \stackrel{\text{def}}{=} \left\{ \frac{1}{N} \sum_{i \in \mathcal{I}} \int_{\theta \in i} \mathbf{x}_\theta d\theta : \mathbf{x}_\theta \in \mathcal{X}_i, \forall \theta \in i, \forall i \in \mathcal{I} \right\}.$$

All players θ in population $i \in \mathcal{I}$ have a similar cost function denoted by $\mathbf{x}_\theta, \bar{\mathbf{X}} \mapsto f_i(\mathbf{x}_\theta, \bar{\mathbf{X}})$, convex w.r.t. \mathbf{x}_θ .

We are only interested in *symmetric* action profiles, where all the nonatomic players in each population i play the same action. Denote the set of symmetric action profiles by $\mathcal{X}^{\mathcal{I}} = \prod_{i \in \mathcal{I}} \mathcal{X}_i$. Considering the coupling constraint $\mathbf{X} \in \mathcal{A}$, let:

$$\mathcal{X}^{\mathcal{I}}(\mathcal{A}) \stackrel{\text{def}}{=} \{ \mathbf{x} \in \mathcal{X}^{\mathcal{I}} : \bar{\mathbf{X}} = \frac{1}{N} \sum_{i \in \mathcal{I}} p_i \mathbf{x}_i \in \mathcal{A} \}.$$

We introduce two indicators to quantify the clustering of $\bar{\mathcal{G}}^{\mathcal{I}}$:

- $\bar{\delta} = \max_{i \in \mathcal{I}} \delta_i$, where

$$\delta_i \stackrel{\text{def}}{=} \max_{n \in \mathcal{N}_i} d_H(\mathcal{X}_n, \mathcal{X}_i), \quad (7)$$

- $\bar{d} = \max_{i \in \mathcal{I}} d_i$, where

$$d_i \stackrel{\text{def}}{=} \max_{n \in \mathcal{N}_i} \sup_{\mathbf{x}_i \in \mathcal{X}_i, \bar{\mathbf{X}} \in \mathcal{M}} d_H(\partial_1 f_i(\mathbf{x}_i, \bar{\mathbf{X}}), \partial_1 f_n(\mathbf{x}_i, \bar{\mathbf{X}})). \quad (8)$$

The quantity δ_i measures the heterogeneity in strategy sets of populations within the group \mathcal{N}_i , while d_i measures the heterogeneity in the subgradients in the group \mathcal{N}_i .

Since the auxiliary game $\bar{\mathcal{G}}^{\mathcal{I}}$ is to be used to compute an approximation of a VNE of \mathcal{G} , the indicators δ_i and d_i should be minimized when defining $\bar{\mathcal{G}}^{\mathcal{I}}$. Thus, we assume that $(\mathcal{X}_i)_i$ and $(u_i)_i$ are chosen such that the following holds:

Assumption 3: For each $i \in \mathcal{I}$, we have:

- \mathcal{X}_i is in \mathcal{X}_0 , the convex hull of $\bigcup_{n \in \mathcal{N}_i} \mathcal{X}_n$, so that $\max_{\mathbf{x} \in \mathcal{X}_i} \|\mathbf{x}\| \leq R$. Moreover, for each $n \in \mathcal{N}_i$, $\text{aff } \mathcal{X}_n \subset \text{aff } \mathcal{X}_i$, where $\text{aff } S$ denotes the affine hull of S ;
- similarly, f_i is chosen such that $\partial_1 f_i(\mathbf{x}_i, \bar{\mathbf{X}})$ is contained in the convex hull of $\bigcup_{n \in \mathcal{N}_i} \partial_1 f_n(\mathbf{x}_i, \bar{\mathbf{X}})$ for all $\mathbf{x} \in \mathcal{X}_i$ and $\bar{\mathbf{X}} \in \mathcal{M}$, so that $\sup_{\mathbf{g}_i \in \partial_1 f_i(\mathbf{x}_i, \bar{\mathbf{X}})} \|\mathbf{g}_i\| \leq \max_{n \in \mathcal{N}_i} L_{1n} \leq L_1$.

An interesting case appears when \mathcal{N} can be divided into homogeneous populations, as in Ex. 1 below:

Example 1: The player set \mathcal{N} can be divided into a small number p of subsets $(\mathcal{N}_i)_{1 \leq i \leq p}$, with homogeneous players inside each subset \mathcal{N}_i (i.e., for each i and $n, m \in \mathcal{N}_i$, $\mathcal{X}_n = \mathcal{X}_m$ and $f_n = f_m$). In that case, consider an auxiliary game $\bar{\mathcal{G}}^{\mathcal{I}}$ with i populations and, for each $i \in \mathcal{I}$ and $n \in \mathcal{N}_i$, $\mathcal{X}_i \stackrel{\text{def}}{=} \mathcal{X}_n$ and $f_i \stackrel{\text{def}}{=} f_n$. Then, $\bar{\delta} = \bar{d} = 0$.

In order to approximate the SVWE of $\mathcal{G}(\mathcal{A})'$ by the SVWE of an auxiliary game $\bar{\mathcal{G}}^{\mathcal{I}}$, let us first state the following result on the geometry of the action sets for technical use.

Lemma 2: Under Asm. 1, there exists a strictly positive constant ρ and an action profile $\mathbf{z} \in \mathcal{X}(\mathcal{A})$ such that, $d(\mathbf{z}_n, \text{rbd } \mathcal{X}_n) \geq \rho$ for all $n \in \mathcal{N}$, where rbd stands for the relative boundary.

Proof: See Appendix I. ■

Given a *symmetric* action profile $\mathbf{x}^{\mathcal{I}} \in \mathcal{X}^{\mathcal{I}}(\mathcal{A})$ in the auxiliary game $\bar{\mathcal{G}}^{\mathcal{I}}(\mathcal{A})$, we can define a corresponding symmetric action profile of $\mathcal{G}(\mathcal{A})'$ such that all the nonatomic players in the populations regrouped in \mathcal{N}_i play the same action $\mathbf{x}_i^{\mathcal{I}}$ (it is allowed that $\mathbf{x}_i^{\mathcal{I}}$ be not in \mathcal{X}_n). Formally, we define the map $\psi : \mathbb{R}^{pT} \rightarrow \mathbb{R}^{NT}$:

$$\forall \mathbf{x}^{\mathcal{I}} \in \mathbb{R}^{pT}, \psi(\mathbf{x}^{\mathcal{I}}) = (\mathbf{x}_n)_{n \in \mathcal{N}} \text{ where } \mathbf{x}_n = \mathbf{x}_i^{\mathcal{I}}, \forall n \in \mathcal{N}_i.$$

Conversely, for a symmetric action profile \mathbf{x} in $\mathcal{G}(\mathcal{A})'$, we define a corresponding symmetric action profile in the auxiliary game $\bar{\mathcal{G}}^{\mathcal{I}}(\mathcal{A})$ by the following map $\bar{\psi} : \mathbb{R}^{NT} \rightarrow \mathbb{R}^{pT}$:

$$\forall \mathbf{x} \in \mathbb{R}^{NT}, \bar{\psi}(\mathbf{x}) = (\mathbf{x}_i^{\mathcal{I}})_{i \in \mathcal{I}} \text{ where } \mathbf{x}_i^{\mathcal{I}} = \frac{1}{p_i} \sum_{n \in \mathcal{N}_i} \mathbf{x}_n.$$

Thm. 2 below is the main result of this subsection. It gives an upper bound on the distance between the SVWE of the population game $\mathcal{G}(\mathcal{A})'$ —of same dimension as the original atomic game $\mathcal{G}(\mathcal{A})$ —and the SVWE of an auxiliary game $\bar{\mathcal{G}}^{\mathcal{I}}(\mathcal{A})$ with reduced dimension.

Theorem 2 (SVWE of $\bar{\mathcal{G}}^{\mathcal{I}}(\mathcal{A})$ is close to SVWE of $\mathcal{G}(\mathcal{A})'$): Under Asms. 1 and 3, in an auxiliary game $\bar{\mathcal{G}}^{\mathcal{I}}(\mathcal{A})$, $\bar{\delta}$ and \bar{d} are defined by (7) and (8), with $\bar{\delta} < \frac{\rho}{2}$. Let \mathbf{x} be an SVWE of $\bar{\mathcal{G}}^{\mathcal{I}}(\mathcal{A})$, and \mathbf{x}^* an SVWE of $\mathcal{G}(\mathcal{A})'$. Then:

- if H is α -strongly monotone, then \mathbf{x} and \mathbf{x}^* are unique and

$$\|\psi(\mathbf{x}) - \mathbf{x}^*\| \leq \sqrt{N \frac{K(\bar{\delta}, \bar{d})}{\alpha}}, \text{ besides} \quad (9)$$

$$\frac{1}{N} \sum_n \|\psi_n(\mathbf{x}) - \mathbf{x}_n^*\| \leq \sqrt{\frac{K(\bar{\delta}, \bar{d})}{\alpha}} \text{ and } \|\bar{\mathbf{X}} - \bar{\mathbf{X}}^*\| \leq \sqrt{\frac{K(\bar{\delta}, \bar{d})}{\alpha}};$$

- if H is β -aggr. str. monotone, then $\bar{\mathbf{X}}, \bar{\mathbf{X}}^*$ are unique and

$$\|\bar{\mathbf{X}} - \bar{\mathbf{X}}^*\| \leq \sqrt{\frac{K(\bar{\delta}, \bar{d})}{\beta}}, \quad (10)$$

where $K(\bar{\delta}, \bar{d}) \xrightarrow{\bar{\delta}, \bar{d} \rightarrow 0} 0$ is a quantity defined as:

$$K(\bar{\delta}, \bar{d}) \stackrel{\text{def}}{=} 2R(3 \frac{L_1}{\rho} \bar{\delta} + \bar{d}). \quad (11)$$

Proof: Let us start by two technical lemmas, for which the proofs are given in Appendix II.

Lemma 3:

- For each $i \in \mathcal{I}$ and $\mathbf{x} \in \mathcal{X}_i$, if $d(\mathbf{x}, \text{rbd } \mathcal{X}_i) > \delta_i$, then $\mathbf{x} \in \mathcal{X}_n$ for each $n \in \mathcal{N}_i$.
- For each $i \in \mathcal{I}$, $n \in \mathcal{N}_i$ and $\mathbf{x} \in \mathcal{X}_n$, if $d(\mathbf{x}, \text{rbd } \mathcal{X}_n) > \delta_i$, then $\mathbf{x} \in \mathcal{X}_i$.

Lemma 4: Under Asm. 1, if $\bar{\delta} < \frac{\rho}{2}$, then:

- for each $\mathbf{x} \in \mathcal{X}^{\mathcal{I}}(\mathcal{A})$, there is $\mathbf{w} \in \mathcal{X}(\mathcal{A})$ such that $\|\mathbf{w}_n - \psi_n(\mathbf{x})\| \leq 4R \frac{\bar{\delta}}{\rho}$ for each $n \in \mathcal{N}$;
- for each $\mathbf{x} \in \mathcal{X}(\mathcal{A})$, there is $\mathbf{w} \in \mathcal{X}^{\mathcal{I}}(\mathcal{A})$ such that $\|\mathbf{w}_i - \bar{\psi}_i(\mathbf{x})\| \leq 2Rp_i \frac{\bar{\delta}}{\rho}$ for each $n \in \mathcal{I}$.

Let $\mathbf{w} \in \mathcal{X}(\mathcal{A})$ be s.t. $\forall n \in \mathcal{N}$, $\|\mathbf{w}_n - \psi_n(\mathbf{x})\| \leq 4R\bar{\delta}/\rho$ (cf. Lemma 4). Since \mathbf{x}^* (resp \mathbf{x}) is an SVWE of $\mathcal{G}(\mathcal{A})'$ (resp. $\bar{\mathcal{G}}^{\mathcal{I}}(\mathcal{A})$), there exists $\mathbf{g}(\mathbf{x}^*) \in H(\mathbf{x}^*)$ (resp. $(\mathbf{h}_i(\mathbf{x}))_{i \in \mathcal{I}} \in \prod_i \partial_1 f_i(\mathbf{x}_i, \bar{\mathbf{X}})$).

$$\sum_n \langle \mathbf{g}(\mathbf{x}^*), \mathbf{x}_n^* - \mathbf{w}_n \rangle \leq 0 \text{ and } \sum_i p_i \langle \mathbf{h}_i(\mathbf{x}), \mathbf{x}_i - \mathbf{y}_i \rangle \leq 0.$$

for all $\mathbf{y} \in \mathcal{X}^{\mathcal{I}}(\mathcal{A})$. Moreover, for each i and $n \in \mathcal{N}_i$, by definition of d_n , let $\mathbf{r}_n(\mathbf{x}) \in \partial_1 f_n(\mathbf{x}_i, \bar{\mathbf{X}})$ s.t. $\|\mathbf{r}_n(\mathbf{x}) - \mathbf{h}_i(\mathbf{x})\| \leq d_n$. Then, $\langle \mathbf{g}(\mathbf{x}^*) - \mathbf{r}(\mathbf{x}), \mathbf{x}^* - \psi(\mathbf{x}) \rangle$ equals to:

$$\langle \mathbf{g}(\mathbf{x}^*), \mathbf{x}^* - \mathbf{w} \rangle + \langle \mathbf{g}(\mathbf{x}^*), \mathbf{w} - \psi(\mathbf{x}) \rangle + \langle \mathbf{r}(\mathbf{x}), \psi(\mathbf{x}) - \mathbf{x}^* \rangle.$$

The first term is nonpositive because \mathbf{x}^* is SVWE. By definition of \mathbf{w} , the second term is bounded by $L_1 N 4R\bar{\delta}/\rho$. Let us decompose the third term, using a $\mathbf{y} \in \mathcal{X}^{\mathcal{I}}(\mathcal{A})$ s.t. $\forall i$, $\|\mathbf{y}_i - \bar{\psi}_i(\mathbf{x}^*)\|_{\mathcal{I}} \leq 2Rp_i \bar{\delta}/\rho$ (cf. Lemma 4), into:

$$\langle \mathbf{r}(\mathbf{x}) - \mathbf{h}(\mathbf{x}), \psi(\mathbf{x}) - \mathbf{x}^* \rangle + \langle \mathbf{h}(\mathbf{x}), \psi(\mathbf{x}) - \mathbf{y} \rangle + \langle \mathbf{h}(\mathbf{x}), \mathbf{y} - \mathbf{x}^* \rangle.$$

The second term is nonpositive because \mathbf{x} is an SVWE of $\bar{\mathcal{G}}^{\mathcal{I}}(\mathcal{A})$. The first term is bounded by $2N R \bar{d}$. The third and last term is bounded by $L_1 2RN \bar{\delta}/\rho$.

We conclude to obtain the bounds stated above from the definition of H being α -strongly monotone (resp. β -aggr. strongly monotone). ■

Remark 1: The bound (9) given on the *individual profiles* diverges with the number of players N . This is a consequence of the fact that individual errors, $\|\mathbf{x}_n - \mathbf{x}_i^*\|$ for each n within a population \mathcal{N}_i , may accumulate, and this accumulation is captured by the Euclidean norm, as $\|\mathbb{1}_N\| = \sqrt{N}$.

We have pointed out that the approximation error depends on how the populations are clustered according to \mathcal{I} , and is related to the heterogeneity of players in \mathcal{N} rather than their number. In particular, in the case of Ex. 1, Thm. 2 states that the (aggregate) SVWE of the auxiliary game $\tilde{\mathcal{G}}^{\mathcal{I}}(\mathcal{A})$ is exactly equal to the (aggregate) SVWE of the large game $\mathcal{G}(\mathcal{A})'$.

A corollary of Thm. 2-i) is that two populations in $\mathcal{G}(\mathcal{A})'$ with similar characteristics have similar behavior at an SVWE, which extends Prop. 3 in the presence of coupling constraints.

Corollary 1: Let $\mathbf{x}^* \in \mathcal{X}$ be an SVWE of game $\mathcal{G}(\mathcal{A})'$. Under Asm. 1, for two populations n and m in \mathcal{N} , if $d_H(\mathcal{X}_n, \mathcal{X}_m) \leq \delta$, $\sup_{\mathbf{x} \in \mathcal{M}/p} d_H(\partial(-u_m)(\mathbf{x}), \partial(-u_n)(\mathbf{x})) \leq d$, and u_n (resp. u_m) is α_n - (resp. α_m -) strongly concave, then

$$\|\mathbf{x}_n^* - \mathbf{x}_m^*\| \leq \left(\frac{1}{\sqrt{\alpha_n}} + \frac{1}{\sqrt{\alpha_m}} \right) K(\delta, d)^{1/2}.$$

C. Combining the two steps to approximate a VNE of $\mathcal{G}(\mathcal{A})$

The following theorem is the main result of the paper, which is immediately obtained as the combination of the two steps of approximation given in Thm. 1 and in Thm. 2 in the computation of a VNE of the original game $\mathcal{G}(\mathcal{A})$.

Corollary 2 (SVWE of $\tilde{\mathcal{G}}^{\mathcal{I}}(\mathcal{A})$ is close to VNEs of $\mathcal{G}(\mathcal{A})$): Under Asms. 1 and 3, in an auxiliary game $\tilde{\mathcal{G}}^{\mathcal{I}}(\mathcal{A})$, $\bar{\delta}$ and \bar{d} are defined by Eqs. (7) and (8), with $\bar{\delta} < \frac{\rho}{2}$. Let \mathbf{x}^* be an SVWE of $\tilde{\mathcal{G}}^{\mathcal{I}}(\mathcal{A})$, $\mathbf{x} \in \mathcal{X}(\mathcal{A})$ be a VNE of $\mathcal{G}(\mathcal{A})$, $\bar{\mathbf{X}}^* = \frac{1}{N} \sum_{i \in \mathcal{I}} p_i \mathbf{x}_i^*$, $\bar{\mathbf{X}} = \frac{1}{N} \sum_{n \in \mathcal{N}} \mathbf{x}_n$, and $K(\bar{\delta}, \bar{d})$ the constant given by (11).

i) if H is α -strongly monotone, then \mathbf{x}^* is unique and

$$\begin{aligned} \|\psi(\mathbf{x}^*) - \mathbf{x}\| &\leq \frac{L_2}{\alpha} \frac{1}{\sqrt{N}} + \sqrt{N \frac{K(\bar{\delta}, \bar{d})}{\alpha}}, \\ \frac{1}{N} \sum_n \|\psi_n(\mathbf{x}^*) - \mathbf{x}_n\| &\leq \frac{L_2}{\alpha N} + \sqrt{\frac{K(\bar{\delta}, \bar{d})}{\alpha}} \\ \text{and } \|\bar{\mathbf{X}} - \bar{\mathbf{X}}^*\| &\leq \frac{L_2}{\alpha N} + \sqrt{\frac{K(\bar{\delta}, \bar{d})}{\alpha}}; \end{aligned}$$

ii) if H is β -aggr. str. monotone, then $\bar{\mathbf{X}}^*$ is unique, and

$$\|\bar{\mathbf{X}} - \bar{\mathbf{X}}^*\| \leq R \sqrt{\frac{2TC}{N\beta}} + \sqrt{\frac{K(\bar{\delta}, \bar{d})}{\beta}}.$$

Given the large game $\mathcal{G}(\mathcal{A})$ and a certain $p \in \mathbb{N}^*$, Cor. 2 suggests that we should find the auxiliary game $\tilde{\mathcal{G}}^{\mathcal{I}}$ with $\mathcal{I} = \{1, \dots, p\}$ that minimizes $K(\bar{\delta}, \bar{d})$ in order to have the best possible approximation of the equilibria. This would correspond to a ‘‘clustering problem’’ given as follows:

$$\min_{(\mathcal{N}_i)_i \in \mathcal{P}^p(\mathcal{N})} \min_{(\mathcal{X}_i)_i} \min_{(u_i)_i} K(\bar{\delta}, \bar{d}), \quad (12)$$

where $\mathcal{P}^p(\mathcal{N})$ denotes the set of all partitions of \mathcal{N} of cardinal p , while $(\mathcal{X}_i)_i$ and $(u_i)_i$ are chosen according to Asm. 3.

The value of the optimal solutions of problem (12), and thus of the quality of the approximation in Cor. 2, depends on the homogeneity of the N players in \mathcal{N} in terms of action sets and utility functions. The ‘‘ideal’’ case is given in Ex. 1 where \mathcal{N} is composed of a small number p of homogeneous populations and thus $K(\bar{\delta}, \bar{d}) = 0$.

Solving (12) is a hard problem in itself: it generalizes the k -means clustering problem [30] (with $k = p$ and considering a function of Hausdorff distances), which is NP-hard [31]. In Sec. IV, we illustrate how we use directly the k -means algorithm to compute efficiently an approximate solution $(\mathcal{N}_i, \mathcal{X}_i, u_i)_{i \in \mathcal{I}}$ of (12) in the parametric case.

Finally, the number p in the definition of the auxiliary game should be chosen a priori as a trade-off between the minimization of $K(\bar{\delta}, \bar{d})$ and a sufficient minimization of the dimension. Indeed, with $\mathcal{I} = \mathcal{N}$, $\mathcal{X}_n = \mathcal{X}_i$ and $u_i = u_n$, we get $\bar{d} = \bar{\delta} = 0$. However, the aim of Cor. 2 is to find an auxiliary game $\tilde{\mathcal{G}}^{\mathcal{I}}$ with $p \ll N$ so that the dimension of the GVIs characterizing the equilibria (and thus the time needed to compute their solutions) is significantly reduced, while ensuring a relatively small error, measured by \bar{d} and $\bar{\delta}$.

IV. APPLICATION TO DEMAND RESPONSE AND ELECTRICITY FLEXIBILITIES

Demand response (DR) [32] refers to the techniques to optimize the electric consumption of distributed consumers [8]. The increasing number of electric vehicles (EV) offers a new source of flexibility in the optimization of the production and demand. To deal with privacy and distributed decisions and information issues, many works consider game theoretical approaches (considering consumers as players) for the management of consumption flexibilities [33]. Let us consider the charging of EVs on a set of 24-hour time-periods $\mathcal{T} \stackrel{\text{def}}{=} \{1, \dots, T\}$, with $T = 24$, indexing the hours from 10 PM to 9PM the day after (to include night time periods).

a) *Price functions: block rates energy prices:* The operator imposes electricity prices on each time-period, given as *inclining block-rates tariffs* [34], i.e. a piece-wise affine function $c(\cdot)$ which depends on the aggregate demand $X_t = \sum_n x_{n,t}$, where $x_{n,t}$ is the power used by consumer n at time period t (one can easily transform it to a function of the average action \bar{X}_t by scaling coefficients):

$$\begin{aligned} c(X) &= 1 + 0.1X \text{ if } X \leq 500 \\ c(X) &= -49 + 0.2X \text{ if } 500 \leq X \leq 1000 \\ c(X) &= -349 + 0.5X \text{ if } 1000 \leq X. \end{aligned} \quad (13)$$

This function c is continuous and convex. Prices are transmitted by the operator to each consumer (EV owner) n who minimizes, with respect to \mathbf{x}_n , a cost function of the *congestion* form (1) with an energy cost determined by (13) and a utility function u_n defined below. An equilibrium gives a stable situation where each consumer minimizes her cost and has no interest to deviate from her current profile.

b) *Consumers' constraints and parameters:* We consider a set of $N = 2000$ consumers who have demand constraints of the form: $\mathcal{X}_n \stackrel{\text{def}}{=} \{\mathbf{x}_n \in \mathbb{R}_+^T : \sum_t x_{n,t} = E_n \text{ and } \underline{x}_{n,t} \leq x_{n,t} \leq \bar{x}_{n,t}\}$ (14)

where E_n is the total energy needed by n , and $\underline{x}_{n,t}, \bar{x}_{n,t}$ (physical) bounds on the power allowed at time t . The utility functions have the form $u_n(\mathbf{x}_n) \stackrel{\text{def}}{=} -\omega_n \|\mathbf{x}_n - \mathbf{y}_n\|^2$, with \mathbf{y}_n a preferred charging profile. Simulation parameters are chosen as follows:

- E_n is drawn uniformly between 1 and 30 kWh, which corresponds to a typical charge of a residential EV.
- $\underline{x}_n, \bar{x}_n$: First, we generate, in two steps, a continual set of charging time-periods $\mathcal{T}_n = \{h_n - \frac{\tau_n}{2}, \dots, h_n + \frac{\tau_n}{2}\}$;
 - the duration τ_n is uniformly drawn from $\{4, \dots, T\}$;
 - h_n is then uniformly drawn from $\{1 + \frac{\tau_n}{2}, \dots, T - \frac{\tau_n}{2}\}$.
- Next, for $t \notin \mathcal{T}_n$, let $\underline{x}_{n,t} = \bar{x}_{n,t} = 0$ and, for $t \in \mathcal{T}_n$, $\underline{x}_{n,t}$ (resp. $\bar{x}_{n,t}$) is drawn uniformly from $[0, \frac{E_n}{\tau_n}]$ (resp. $[\frac{E_n}{\tau_n}, E_n]$).
- ω_n is drawn uniformly from $[1, 10]$.
- \mathbf{y}_n is taken equal to $\bar{x}_{n,t}$ on the first time periods of \mathcal{T}_n (first available time periods) until reaching E_n (consumption profile ‘‘plug and charge’’).

We consider the following *coupling constraints*:

$$-50 \leq X_T - X_1 \leq 50 \quad (15)$$

$$X_t \leq 1400, \quad \forall t \in \mathcal{T} \quad (16)$$

Constraint (15) imposes that the demand X_T at the end of the time horizon is close to the first aggregate X_1 , so that the profiles computed for the finite time set \mathcal{T} can be applied on a day-to-day, periodical basis. Constraint (16) is a capacity constraint, (e.g. induced by production capacities).

c) *Computing populations with k -means:* To apply the clustering procedure of Sec. III-B, we use the k -means algorithm [30], where “ k ” = p is the number of populations (groups) to replace the large set of N players. For each player n , we define her parametric description vector:

$$\mathbf{v}_n \stackrel{\text{def}}{=} [\omega_n, \mathbf{y}_n, E_n, \underline{\mathbf{x}}_n, \bar{\mathbf{x}}_n] \in \mathbb{R}^{3T+2}. \quad (17)$$

The k -means algorithm finds a partition $(S_i)_{1 \leq i \leq p}$ of \mathcal{N} into p clusters, approximately solving the combinatorial problem:

$$\min_{S_1, \dots, S_p} \sum_{1 \leq i \leq p} \sum_{\mathbf{v} \in S_i} \|\mathbb{E}_{S_i}(\mathbf{v}) - \mathbf{v}\|^2 = \min_{S_1, \dots, S_p} \sum_{1 \leq i \leq p} |S_i| \text{Var}(S_i),$$

where $\mathbb{E}_{S_i}(\mathbf{v}) = \frac{1}{|S_i|} \sum_{\mathbf{v} \in S_i} \mathbf{v}_n$ denotes the average value of \mathbf{v} over the set S_i , taken to define w_i , \mathbf{y}_i , E_i , $\underline{\mathbf{x}}_i$ and $\bar{\mathbf{x}}_i$.

The simulations are run with different population numbers, with $p \ll N$ chosen among $\{5, 10, 20, 50, 100\}$.

The k -means algorithm provides an efficient method to obtain an approximate solution of (12): the solution obtained can be sub-optimal in terms of $K(\delta, d)$, as the algorithm minimizes the squared distance of the average vector of parameters in S_i to the vectors of parameters of the points in S_i . Finding efficient methods to address the problem (12) is an avenue for further research.

d) *Computation methods:* We compute a VNE (Def. 3) with the original set of N players and the approximating SVWE (Def. 4) as solutions of the associated GVI (2).

We employ the iterative projection method exposed in [19, Algo. 2], adapted to our subdifferentiable case by replacing the fixed step τ used in with a variable step $\tau^{(k)} = 1/k$, in the spirit of subgradient algorithms [35], [36]. The algorithm stops when the Euclidean distance between two iterates is smaller than 10^{-3} . We refer to [19] for more details on the algorithm.

e) *A trade-off between precision and computation time:* Simulations were run using Python on a single core Intel Xeon @3.4Ghz and 16GB of RAM. In this simple example, the computation of the actual VNE $\hat{\mathbf{x}}$ (and aggregate profile $\hat{\mathbf{X}}$) of the original game with $N = 2000$ players is still possible, and is used to measure the precision of our approximations.

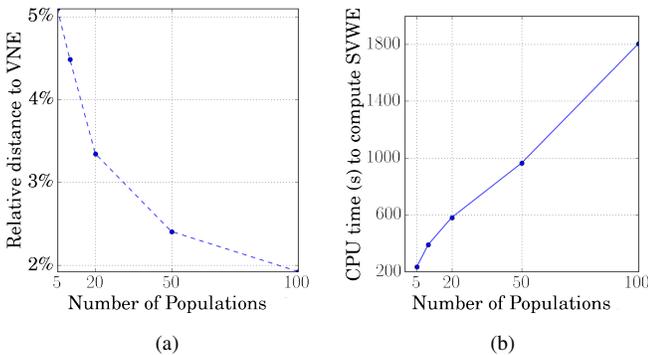


Fig. 1: (a) Relative error to VNE ; (b) SVWE CPU Time.

Fig. 1a and Fig. 1b show the two main metrics to consider to choose a relevant number of populations p : the precision of the SVWE approximating the equilibrium (measured by the relative distance of the aggregate SVWE profile \mathbf{X}^* to the aggregate VNE

profile $\hat{\mathbf{X}}$ computed along, $\|\mathbf{X}^* - \hat{\mathbf{X}}\| / \|\mathbf{X}^*\|$), and the CPU time needed to compute the SVWE. The distance between the aggregate equilibrium profile and its estimation decreases with N at a sublinear rate Fig. 1a. This is explained in light of Thm. 2 and in addition with the following remarks:

- the Hausdorff distance between parameterized polyhedra is Lipschitz continuous w.r.t their parameter vectors (see [37]);
- similarly, as subgradients of utility functions are reduced to a point, one has, for all n , $d_i = \max_{n \in \mathcal{N}_i} \max_{\mathbf{x} \in \mathcal{M}} 2 \|\omega_i(\mathbf{x} - \mathbf{y}_i) - \omega_n(\mathbf{x} - \mathbf{y}_n)\| = \mathcal{O}(\max_{n \in \mathcal{N}_i} |\omega_i - \omega_n| + \|\mathbf{y}_i - \mathbf{y}_n\|)$.

Fig. 1b shows the CPU time needed to compute the SVWE with our stopping criterion. The CPU time evolves linearly with the number of populations p , which is explained by the *distributed* structure of the algorithm, requiring a number of iterations proportional to p . The computation of a solution of the clustering problem with the k -means algorithm is negligible (less than ten seconds for each value of p).

Computing the VNE of the original game with the same configuration and stopping criterion took 3 hours and 26 minutes were needed, more than six times longer than to compute the SVWE with one hundred populations.

Last, the error between the aggregate demand profile at equilibrium and its approximation (Fig. 1a) is between 2% and 5%, which remains significant. However, as pointed out in Sec. III, the quality of the approximation depends on the heterogeneity of the set of players \mathcal{N} : as the parameters are drawn uniformly in this experiment, the set of players \mathcal{N} presents a large variance, and provides us with a “worst” case (as opposed to Ex. 1 which offers an optimal situation).

V. CONCLUSION

This paper shows that equilibria in large splittable aggregative games can be approximated with a Wardrop equilibrium of an auxiliary population game of smaller dimension. Our results give explicit bounds on the distance of this approximating equilibrium to the equilibria of the original large game. These theoretical results can be used in practice to solve, by an iterative method, complex nonconvex bilevel programs where the lower level is the equilibrium of a large aggregative game, for instance, to optimize tariffs or tolls for the operator of a network. The analysis of such a procedure will constitute an interesting extension of this work. Besides, this study highlighted the need for efficient algorithms for solving clustering problems of type (12), as well as distributed algorithms for solving GVI problems without a strong monotonicity assumption [19], [35]. Those algorithmic considerations are also interesting avenues for further research.

APPENDIX I

PROOF OF LEMMA 2: EXISTENCE OF INTERIOR PROFILE

Let $\tilde{\mathbf{x}} \in \mathcal{X}$ be such that for all n :

$$d(\tilde{\mathbf{x}}_n, \text{rbd } \mathcal{X}_n) = \max_{\mathbf{x} \in \mathcal{X}_n} d(\mathbf{x}, \text{rbd } \mathcal{X}_n) \stackrel{\text{def}}{=} \eta_n.$$

Denote $\bar{\mathbf{X}} = \frac{1}{N} \sum_n \tilde{\mathbf{x}}_n$ and $\eta = \min_n \eta_n > 0$. Let $\mathbf{y} \in \mathcal{X}(\mathcal{A})$ and $\bar{\mathbf{Y}} = \frac{1}{N} \sum_n \mathbf{y}_n$ be s.t.:

$$d(\bar{\mathbf{Y}}, \text{rbd } \mathcal{A}) = \max_{\mathbf{X} \in \bar{\mathcal{X}} \cap \mathcal{A}} d(\bar{\mathbf{X}}, \text{rbd } \mathcal{A}).$$

Let us denote $t = d(\bar{\mathbf{Y}}, \text{rbd } \mathcal{A})/3R$ and let us define $\mathbf{z} = \mathbf{y} - t(\mathbf{y} - \tilde{\mathbf{x}}) \in \mathcal{X}$ and $\bar{\mathbf{Z}} = \frac{1}{N} \sum_n \mathbf{z}_n$. We obtain:

$$\|\bar{\mathbf{Y}} - \bar{\mathbf{Z}}\| = t \|\bar{\mathbf{Y}} - \bar{\mathbf{X}}\| \leq t2R \leq \frac{2}{3} d(\bar{\mathbf{Y}}, \text{rbd } \mathcal{A}),$$

hence $\bar{\mathbf{Z}} \in \bar{\mathcal{X}} \cap \text{ri } \mathcal{A}$, where ri means the relative interior. Besides, for any n , $\mathbf{z}_n = \mathbf{y}_n - t(\mathbf{y}_n - \tilde{\mathbf{x}}_n)$. Since $d(\tilde{\mathbf{x}}_n, \text{rbd } \mathcal{X}_n) \geq \eta$,

$\mathbf{y}_n \in \mathcal{X}_n$, and \mathcal{X}_n is convex, we have:

$$d(\mathbf{z}_n, \text{rbd } \mathcal{X}_n) \geq \eta t = \frac{\eta}{3R} d(\bar{\mathbf{Y}}, \text{rbd } \mathcal{A}).$$

We can conclude by defining $\rho \stackrel{\text{def}}{=} \frac{\eta}{3R} d(\bar{\mathbf{Y}}, \text{rbd } \mathcal{A})$.

APPENDIX II

PROOF OF THE LEMMAS OF THM. 2

Proof: Lemma 3 i) Suppose $\mathbf{x} \notin \mathcal{X}_n$. Let $\mathbf{y} \stackrel{\text{def}}{=} P_{\mathcal{X}_n}(\mathbf{x}) \neq \mathbf{x}$.

As $\mathbf{y} \in \text{aff } \mathcal{X}_n \subset \text{aff } \mathcal{X}_i$, then $\mathbf{x} - \mathbf{y} \in \text{aff } \mathcal{X}_i$. Let $\mathbf{z} \stackrel{\text{def}}{=} \mathbf{x} + \delta_i \frac{\mathbf{x} - \mathbf{y}}{\|\mathbf{x} - \mathbf{y}\|}$. Then, $\mathbf{z} \in \mathcal{X}_i$ because $\|\mathbf{z} - \mathbf{x}\| \leq \delta_i$. By the convexity of \mathcal{X}_n and the definition of \mathbf{y} , we have $d(\mathbf{z}, \mathcal{X}_n) = d(\mathbf{x}, \mathcal{X}_n) + \delta_i > \delta_i$, contradicting the fact that $\delta_i \geq d_H(\mathcal{X}_n, \mathcal{X}_i)$. The proof of ii) is symmetric. ■

Proof: [Proof of Lemma 4]

i) For $\mathbf{x} \in \mathcal{X}^{\mathcal{I}}(\mathcal{A})$, define $\mathbf{w} \in \mathcal{X}$ as follows: $\forall i \in \mathcal{I}, \forall n \in \mathcal{N}_i$, let $\mathbf{w}_n \stackrel{\text{def}}{=} \mathbf{x}_i + t(\mathbf{z}_n - \mathbf{x}_i)$ where \mathbf{z} is defined in Lemma 2, with $t \stackrel{\text{def}}{=} 2\bar{\delta}/\rho < 1$. On the one hand:

$$\forall i \in \mathcal{I}, \forall n \in \mathcal{N}_i, d(\mathbf{z}_n, \text{rbd } \mathcal{X}_i) \geq \rho - \bar{\delta}$$

$$\text{implies } d(\mathbf{w}_n, \text{rbd } \mathcal{X}_i) \geq t(\rho - \bar{\delta}) > t\rho/2 = \bar{\delta},$$

because each point in the ball with radius $t(\rho - \bar{\delta})$ centered at \mathbf{w}_n is on the segment linking \mathbf{x}_i and some point in the ball with radius $\rho - \bar{\delta}$ centered at \mathbf{z}_n which is contained in \mathcal{X}_n .

Thus, $\mathbf{w}_n \in \mathcal{X}_n \forall n \in \mathcal{N}_i$ according to Lemma 3.i). On the other hand, the linear mapping $S : \mathbb{R}^{IT} \ni \mathbf{v} \mapsto \frac{1}{N} \sum_{n \in \mathcal{N}} \mathbf{v}_n$ maps the segment linking $\psi(\mathbf{x})$ and \mathbf{z} in $\mathcal{X}(\mathcal{A})$ to a segment linking $\bar{\mathbf{X}} = \frac{1}{N} \sum_i p_i \mathbf{x}_i$ and $\bar{\mathbf{Z}}$ in the convex \mathcal{A} . We get:

$$\frac{1}{N} \sum_{n \in \mathcal{N}} \mathbf{w}_n = t\bar{\mathbf{Z}} + (1-t)\bar{\mathbf{X}} \in \mathcal{A}$$

and thus $\mathbf{w} \in \mathcal{X}(\mathcal{A})$. Finally, $\|\mathbf{w}_n - \psi_n(\mathbf{x})\| = t\|\mathbf{z}_n - \psi_n(\mathbf{x})\| \leq t2R = 4R\frac{\bar{\delta}}{\rho}$.

ii) For $\mathbf{x} \in \mathcal{X}(\mathcal{A})$, let $\mathbf{y} \stackrel{\text{def}}{=} \mathbf{x} + t(\mathbf{z} - \mathbf{x})$ with $t \stackrel{\text{def}}{=} \frac{\bar{\delta}}{\rho}$. Then, by similar arguments as above, $d(\mathbf{y}_n, \text{rbd } \mathcal{X}_n) \geq \bar{\delta}$ hence $\mathbf{y}_n \in \mathcal{X}_i$ and $\bar{\psi}(\mathbf{y}) \in \mathcal{X}^{\mathcal{I}}$. From the convexity of \mathcal{A} :

$$\frac{1}{N} \sum_n \mathbf{y}_n = t\bar{\mathbf{Z}} + (1-t)\left(\frac{1}{N} \sum_n \mathbf{x}_n\right) \in \mathcal{A}.$$

Hence $\mathbf{w} \stackrel{\text{def}}{=} \bar{\psi}(\mathbf{y}) \in \mathcal{X}^{\mathcal{I}}(\mathcal{A})$. Finally, $\|\mathbf{w}_i - \psi_i(\mathbf{x})\| = t\|\sum_{n \in \mathcal{P}_i} (\mathbf{z}_n - \mathbf{x}_n)\| \leq 2Rp_i\frac{\bar{\delta}}{\rho}$. ■

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