# Distributed Autonomous Systems M

Dante Piotto

spring semester 2024

# Contents

1	Intr	roduction and scenarios	5
	1.1	Distributed Autonomous System	5
	1.2	Scenarios and applications of distributed systems	5
	1.3	Measurement filtering in wireless sensor networks	5
	1.4	Parameter Estimation in Wireless Sensor Networks	6
		1.4.1 Opinion Dynamics in Social Influence Networks	6
	1.5	Main questions in averaging algorithms	6
	1.6	Distributed control in cooperative robotics	6
		1.6.1 Main questions in cooperative robotics	7
		1.6.2 Distributed optimal control	7
2	Pre	eliminaries on Algebraic Graph Theory	9
3	Ave	eraging Systems	11
	3.1	Distributed algorithm	11
	3.2	Discrete-time averaging systems	11
	3.3	Stochastic matrices	12
	3.4	Example: Metropolis-Hastings weights	12
	3.5	Time-varying digraphs	13
		3.5.1 Averaging distributed algorithms over time-varying graphs	13
		3.5.2 Discrete-time consensus over time-varying graphs	13
	3.6	Laplacian dynamics	14
		3.6.1 Properties of the Laplacian matrix	
		3.6.2 Consensus for Laplacian dynamics	15
4	Opt	timization basics	17
	4.1	Optimization algorithms	18
		4.1.1 A system theoretical perspective to the gradient method	
	4.2	steady-state analysis of the gradient method	
		4.2.1 Gradient method for quadratic programs	
		4.2.2 Gradient flow	
		4.2.3 Nesterov accelerated gradient method	19
5	Par	callel Optimization and Federated Learning	21
	5.1	Incremental gradient method	
	5.2	Stochastic Gradient Descent	
		5.2.1 Convergence	
	5.3	Beyond SGD: Adaptive Momentum Estimation (Adam)	
	5.4	Federated learning	
	5.5	Distributed learning	23
6	Lea	der Follower networks: Formation control	<b>2</b> 5
	6.1	analogy with mass-spring systems	25

4 CONTENTS

7	Distributed Aggregative Optimization 7.1 Aggregative optimization	30
8	Learning with Neural Networks 8.1 Basic element: the neuron model	<b>33</b>
9	Multi-Robot Safety Controllers         9.1 Control Barrier Functions	35

# Introduction and scenarios

## 1.1 Distributed Autonomous System

Each agent  $i \in \{1, ..., N\}$  has

- local physical and/or cyber state  $x_i$
- computational and sensing capabilities
- communication capability: exchange messages with "neighbours"

## 1.2 Scenarios and applications of distributed systems

- Averaging: distributed estimation, opinion dynamics
- Distributed control in cooperative robotics
- Distributed optimization
  - distributed machine learning
  - distributed decision-making in cooperative robotics
  - distributed optimal control in energy systems and cooperative robotics

# 1.3 Measurement filtering in wireless sensor networks

Consider a network of N sensors with local sensing, computation and communication. Agent  $i, i \in \{1, ..., N\}$ , takes a local measurement from the environment (temperature, pressure, etc.). Let  $x_{i0} \in \mathbb{R}$  be the scalar local measurement Agents are interested in agreeing on the average of the measurements,

$$x_{\text{avg}} = \frac{1}{N} \sum_{i=1}^{N} x_{i0}$$

to have a better estimate of the environment quantity

Consider the following "distributed algorithm" based on "local" linear averaging, for each  $i \in \{1, ..., N\}$ 

$$x_i^0 = x_{i0}$$
 
$$x_i^{k+1} = \text{average}(x_i^k, \{x_j^k, j \text{ "neighbour" of } i\}), \qquad k \in \mathbb{N}$$

generalizing coefficients of the update:

$$x_i^0 = x_{i0}$$

$$x_i^{k+1} = \sum_{j=1}^N a_{ij} x_j^k \qquad k \in \mathbb{N}$$

Remark.  $a_{ij} \geq 0$  and  $\sum_{j=1}^{N} a_{ij} = 1$ 

Remark.  $a_{ij} = 0$ , for some  $j \in \{1, ..., N\}$ , i.e.  $a_{ij} = 0$  if i does not have access to the estimate of j

### 1.4 Parameter Estimation in Wireless Sensor Networks

Consider a network of N sensors with local sensing, computation and communication aiming at estimating a common parameter  $\theta^* \in \mathbb{R}$  Each sensor i measures

$$y_i = B_i \theta^* + v_i$$

with  $y_i \in \mathbb{R}^{m_1}, B_i$  known matrix and  $v_i$  a random measurement noise. Assume  $v_1, \ldots, v_N$  independent and Gaussian, with zero mean and covariance  $E[v_i v_i^T] = \Sigma_i$ . Assume  $\sum_{i=1}^N m_i \ge m$  and  $\begin{bmatrix} B_1 \\ \vdots \\ B_N \end{bmatrix}$  full rank Compute a least-squares estimate

$$\hat{\theta} = \arg\min_{\theta} \sum_{i=1}^{N} (y_i - B_i \theta)^T \Sigma_i^{-1} (y_i - B_i \theta)$$

The optimal solution is

$$\hat{\theta} = \left(\sum_{i=1}^{N} B_i^T \Sigma_i^{-1} B_i\right)^{-1} \sum_{i=1}^{N} B_i^T \Sigma_i^{-1} y_i$$

$$= \left(\frac{1}{N} \sum_{i=1}^{N} B_i^T \Sigma_i^{-1} B_i\right)^{-1} \frac{1}{N} \sum_{i=1}^{N} B_i^T \Sigma_i^{-1} y_i$$

The optimal solution can be obtained by computing two averages  $\frac{1}{N}\sum_{i=1}^{N}\beta_i$  and  $\frac{1}{N}\sum_{i=1}^{N}\beta_i$ 

#### 1.4.1 Opinion Dynamics in Social Influence Networks

Group of N individuals, with  $x_i^k$  being the opinion of individual i at time k. Opinions are updated according to

$$x_i^{k+1} = \sum_{j=1}^{N} a_{ij} x_j^k$$

# 1.5 Main questions in averaging algorithms

- Do node estimates converge? Do they converge to a common value ("reach consensus")?
- Do they reach consensus to the average ("average consensus")?
- How can we model communication in general networks?
- Can we answer the above questions for general networks and communication protocols?
- What assumptions do we need on the communication network?

# 1.6 Distributed control in cooperative robotics

Team of N (mobile) robots aiming to execute complex tasks

#### Basic tasks

- rendevous, containment
- formation, flocking
- coverage

#### Complex tasks

- pickup and delivery
- $\bullet\,$  surveillance and patrolling
- exploration
- satellite constellation

### 1.6.1 Main questions in cooperative robotics

- Do robot states asymptotically converge?
- Do the asympototic staes satisfy the global, desired task?
- How can we model communication in (general) robotic networks?
- What assumptions do we need on the communication network?
- Can we answer the above questions for general networks and communication protocols?

## 1.6.2 Distributed optimal control

$$\begin{aligned} \min_{\substack{x_1,\dots,x_N\\u_1,\dots,u_N\\i=1}} \sum_{i=1}^N (\sum_{\tau=0}^{T-1} \ell_i(z_{i,\tau},u_{i,\tau}) + m_i(z_{i,T})) \\ \text{subj to } \sum_{i=1}^N H_i z_{i,\tau} \leq h, & \tau \in [0,T] \\ z_{i,\tau+1} = A_i z_{i,\tau} + B_i u_{i,\tau} & \forall i,\tau \in [0,T] \\ z_{i,\tau} \in Z_i, \quad u_{i,\tau} \in U_i, & \forall i,\tau \in [0,T] \end{aligned}$$

# Preliminaries on Algebraic Graph Theory

```
Definition 2.1 (Digraph)
A digraph is a pair G = (I, E) where I = 1, ..., N is a set of elements called nodes and E \subset I \times I is a set of
ordered node pairs called edges
Edge: the pair (i, j) denotes an edge from i to j
Self-loop: edge from a node to itself, i.e. (i,i)
Definition 2.2 (Undirected (di)graph)
if for any (i, j) \in E then (j, i) \in E
Definition 2.3 (Subgraph)
(I',E') subgraph of (I,E) if I'\subset I and E'\subset E. Spanning subgraph if I'=I
Definition 2.4 (In-neighbours of i)
j \in I is an in-neighbour of i \in I if (j,i) \in E
Definition 2.5 (Set of in-neighbours of i)
\mathcal{N}_{i}^{\text{IN}} = \{ j \in \{1, \dots, N\} | (j, i) \in E \}
Definition 2.6 (Out-neighbours of i)
j \in I is an out-neighbour of i \in I if (i, j) \in E
Definition 2.7 (Set of out-neighbours of i)
\mathcal{N}_{i}^{\text{IN}} = \{ j \in \{1, \dots, N\} | (i, j) \in E \}
Definition 2.8 (In-degree \deg_i^{IN})
number of in-neighbours, i.e. car
inality of \mathcal{N}_i^{\text{IN}}(\deg_i^{\text{IN}} = |\mathcal{N}_i^{\text{IN}}|)
     Out-degree analogous
 \begin{array}{l} \textbf{Definition 2.9} \text{ (Balanced digraph)} \\ \textbf{A digraph } G \text{ is balanced if } \deg_i^{\text{IN}} = \deg_i^{\text{OUT}} \text{ for all } i \in \{1, \dots, N\} \end{array} 
Definition 2.10 (Complete graph)
Unweighted graph such that \forall i, j \exists (i, j), (j, i) \in E
```

# **Averaging Systems**

# 3.1 Distributed algorithm

Given a network of N agents communicating according to a fixed digraph G, i.e. each agent i can receive messages only from in-neighbours in the graph, i.e. from  $j \in \mathcal{N}_i^{\text{IN}}$ . We start by considering a fixed graph, thus, each agent communicates with the same neighbours at each iteration  $k \in \mathbb{N}$ 

$$x_i^{k+1} = \text{stf}_i(x_i^k, \{x_i^k\}_{i \in \mathcal{N}^{\text{IN}}}), \qquad i \in \{1, \dots, N\}$$

where  $\mathrm{stf}_i$  is a function depending only on state  $x_i$  and states  $x_j, j \in \mathcal{N}_i^{\mathrm{IN}}$ .

Alternative version with out-neighbours:

$$x_i^{k+1} = \operatorname{stf}_i(\{x_j\}_{j \in \mathcal{N}_i^{\text{OUT}}})$$

# 3.2 Discrete-time averaging systems

Let  $G^{\text{comm}} = (I, E)$  be a fixed (communication) digraph (self loops included). A linear averaging distributed algorithm can be written as:

$$x_i^{k+1} = \sum_{J \in \mathcal{N}_i^{\text{IN}}} a_{ij} x_j^k \qquad i \in \{1 \dots, N\}$$

where  $x_i^k \in \mathbb{R}$  is the state of agent i at k and  $a_{ij} > 0$  are positive weights.

Remark. The weights  $a_{ij}$  are defined only for  $(i,j) \in E$ 

Wach i uses only the states of neighbours  $j \in \mathcal{N}_i^{\text{IN}}$ , thus distributed algorithm.

For analysis purposes, let us define weights  $a_{ij} = 0$  for  $(j,i) \notin E$ . Thus we can rewrite the distributed algorithm as

$$x_i^{k+1} = \sum_{i=1}^{N} a_{ij} x_j^k \qquad i \in \{1, \dots, N\}$$

This is a LTI autonomous system

$$\begin{bmatrix} x_1^{k+1} \\ \vdots \\ x_N^{k+1} \end{bmatrix} = \begin{bmatrix} a_{11} & \cdots & a_{1N} \\ \vdots & \ddots & \vdots \\ a_{N1} & \cdots & a_{NN} \end{bmatrix} \begin{bmatrix} x_1^k \\ \vdots \\ x_N^k \end{bmatrix}$$

Which can be compactly written as

$$x^{k+1}0Ax^k$$

Remark. The matrix A can be seen as the weighted adjacency matrix of the reverse digraph  $G^{\text{comm,rev}}$  of the digraph  $G^{\text{comm}}$ 

If instead of in-neighbours we use out-neighbours, we call the digraph a sensing digraph  $G^{\text{sens}}$ . In this case the notation becomes consistent with graph theory, so we get

$$x^{k+1} = Ax^k$$

where A can be seen as the weighted adjacency matrix of the sensing digraph  $G^{\text{sens}}$ 

## 3.3 Stochastic matrices

The non-negative square matrix  $A \in \mathbb{R}^{N \times N}$  is

- row stochastic if  $A\mathbf{1} = \mathbf{1}$  (each row sums to 1)
- column stochastic if  $A^{\top} \mathbf{1} = \mathbf{1}$  (each column sums to 1)
- doubly stochastic if both row and column stochastic.

**Lemma.** Let A be a row-stochastic matrix and G the associate digraph. If G is strongly connected and aperiodic, then

- 1. the eigenvalue  $\lambda = 1$  is simple;
- 2. all the other eigenvalues  $\mu$  satisfy  $|\mu| < 1$

Remark. The condition "G contains a globally reachable node and the subgraph of globally reachable noes is aperiodic" is necessary and sufficient

#### **Theorem 3.1** (Consensus)

Consider a (discrete-time) averaging system with associated digraph G and wieghted adjacency matrix A. Assume G is strongly connected and aperiodic, and A is row stochastic. Then

1. there exists a left eigenvector  $w \in \mathbb{R}^N$ , w > 0 (i.e. with positive components  $w_i > 0$  for all i = 1, ..., N) such that

$$\lim_{k \to \infty} x^k = \mathbf{1} \frac{w^\top x^0}{w^\top \mathbf{1}} = \begin{bmatrix} 1 \\ \vdots \\ 1 \end{bmatrix} \frac{\sum_{i=1}^N w_i x_i^0}{\sum_{i=1}^N w_i}$$

i.e., consensus is reached to  $\frac{\sum_{i=1}^{N}w_{i}x_{i}^{0}}{\sum_{i=1}^{N}w_{i}}$ 

2. if additionally A is doubly stochastic, then

$$\lim_{k \to \infty} x^k = \begin{bmatrix} 1 \\ \vdots \\ 1 \end{bmatrix} \frac{\sum_{i=1}^N x_i^0}{N}$$

i.e., average consensus is reached

# 3.4 Example: Metropolis-Hastings weights

Given an undirected unweighted graph G with edge set E and degrees  $d_1, \ldots, d_n$ 

$$a_{ij} = \begin{cases} \frac{1}{1 + \max\{d_i, d_j\}} & \text{if } (i, j) \in E \text{ and } i \neq j \\ 1 - \sum_{h \in \mathcal{N}_i \setminus \{i\}} a_{ih} & \text{if } i = j \\ 0 & \text{otherwise} \end{cases}$$

Result: the matrix A is symmetric and doubly-stochastic.

# 3.5 Time-varying digraphs

A time-varying digraph is a sequence of digraphs  $\{G(k)\}_{k\geq 0}$ .

Remark. The main definitions of in/out neighbours, in/out degree, adjacency matrix can be generalized by considering time-varying versions, i.e.  $\mathcal{N}_i^{\text{IN}}(k)$ ,  $\mathcal{N}_i^{\text{OUT}}(k)$ ,  $\deg_i^{\text{IN}}(k)$ ,  $\deg_i^{\text{OUT}}(k)$ , A(k) associated to each graph G(k). Connectivity requires new definitions as assuming each G(k) to be connected is too conservative.

**Definition 3.1** (Jointly strongly connected digraph) if  $\bigcup_{\tau=k}^{+\infty} G(\tau)$  is strongly connected  $\forall k \geq 0$ 

**Definition 3.2** (Uniformly jointly strongly connected (or *B*-strongly connected) digraph) if there exists  $B \in \mathbb{N}$  such that  $\bigcup_{\tau=k}^{k+B} G(\tau)$  is strongly connected  $\forall k \geq 0$ 

*Remark.* The graph can be disconnected at some time k.

### 3.5.1 Averaging distributed algorithms over time-varying graphs

Let  $\{G(k)\}_{k\geq 0}$  be a time-varying digraph (with self loops for each G(k)). Consider the distributed algorithm

$$x_i^{k+1} = \sum_{j \in \mathcal{N}_i^{\text{IN}}(k)} a_{ij}(k) x_j^k \qquad \forall i \in \{1, \dots, N\}$$

or the out-neighbours version

$$x_i^{k+1} = \sum_{j \in \mathcal{N}_i^{\text{OUT}}(k)} a_{ij}(k) x_j^k \qquad \forall i \in \{1, \dots, N\}$$

where  $x_i^k \in \mathbb{R}$  is the state of agent i at k and  $a_{ij}(k) > 0$ .

For analysis purposes, let us define weights  $a_{ij}(k) = 0$  for  $(i, j) \notin E(k)$ . Thus we can rewrite the distributed algorithm as

$$x_i^{k+1} = \sum_{j=1}^{N} a_{ij}(k) x_j^k \qquad i \in \{1, \dots, N\}$$

This is a Linear Time-Varying system

$$x^{k+1} = A(k)x^k$$

with state  $x := [x_1, \dots, x_N]^{\top}$  and state matrix

$$A(k) := \begin{bmatrix} a_1 1k & \cdots & a_{1N}(k) \\ \vdots & \ddots & \vdots \\ a_N 1k & \cdots & a_{NN}(k) \end{bmatrix}$$

being a weighted adjacency matrix associated to the digraph G(k).

#### 3.5.2 Discrete-time consensus over time-varying graphs

#### Theorem 3.2

Let  $\{A(k)\}_{k\geq 0}$  be a sequence of row-stochastic matrices with associated digraphs  $\{G(k)\}_{k\geq 0}$ . Assume

- 1. each digraph G(k) has a self-loop at each node;
- 2. each non-zero edge weight  $a_{ij}(k)$ , including the self-loop wights  $a_{ii}(k)$ , is larger than a constant  $\epsilon > 0$ ;
- 3. there exists  $B \in \mathbb{N}$  such that, for all times  $k \geq 0$ , the union digraph  $G(k) \cup \cdots \cup G(k+B)$  is strongly connected.

Then

1. there exists a non-negative vector  $w \in \mathbb{R}^N$  such that the solution to  $x^{k+1} = A(k)x^k$  converges (exponentially) to  $\mathbf{1} \frac{w^\top x^0}{w^\top 1}$ , i.e.

$$\lim_{k \to \infty} x^k = \mathbf{1} \left( \frac{w^\top x^0}{w^\top \mathbf{1}} \right)$$

2. if additionally each matrix in the sequence is doubly-stochastic, then

$$\lim_{k \to \infty} x^k = \mathbf{1} \frac{1}{N} \sum_{i=1}^N x_i^0$$

i.e., average consensus is achieved

# 3.6 Laplacian dynamics

Consider a network of dynamical systems with dynamics

$$\dot{x}(t) = u_i(t) \qquad i \in \{1, \dots, N\}$$

with states  $x_i \in \mathbb{R}$  and inputs  $u_i \in \mathbb{R}$ , communicating (or interacting) according to a digraph G. Consider a (distributed) "proportional" feedback control

$$u_i(t) = -\sum_{j \in \mathcal{N}_i^{\text{IN}}} a_{ij} (x_i(t) - x_j(t))$$

or the out-neighbour version

$$u_i(t) = -\sum_{j \in \mathcal{N}_i^{\text{OUT}}} a_{ij}(x_i(t) - x_j(t))$$

For analysis purposes, let us define weights  $a_{ij}(k) = 0$  for  $(i,j) \notin E(k)$ . Thus we can rewrite the distributed control systems as

$$\dot{x}_i(t) = -\sum_{j=1}^N a_{ij}(x_i(t) - x_j(t)) \qquad \forall i \in \{1, \dots, N\}$$

Defining  $x := [x_1 \cdots x_N]^{\top}$ , it can be shown that it can be rewritten as the following Linear Time Invariant continuous-time system

$$\dot{x}(t) = -Lx(t)$$

where L is the (weighted) Laplacian associated to the digraph G with (weighted) adjacency matrix A Let

$$\dot{x}_i(t) = -\sum_{j=1}^{N} a_{ij}(x_i(t) - x_j(t)) \quad \forall i \in \{1, \dots, N\}$$

rearranging terms

$$\dot{x}_i(t) = -\left(\sum_{j=1}^N a_{ij}\right) x_i(t) + \sum_{j=1}^N a_{ij} x_j(t) = -\deg_i^{\text{OUT}} x_i(t) + (Ax(t))_i$$

where  $(Ax(t))_i$  is the *i*-th element of Ax(t). Writing the previous dynamics in a compact form

$$\dot{x}(t) = -(D^{OUT} - A)x(t)$$

where we recall that  $D^{\text{OUT}}$  is the (weighted) out-degree matrix. Recalling that  $L = D^{\text{OUT}} - A$ , it holds that

$$\dot{x}(t) = -Lx(t)$$

Remark. if the in-neighbours version is considered, then  $\dot{x}(t) = -L^{\text{IN}}x(t)$ , where  $L^{\text{IN}} = D^{\text{IN}} - A^T$  is the in-degree Laplacian (i.e. the Laplacian of the reverse graph of G)

### 3.6.1 Properties of the Laplacian matrix

It can be easily verified that

$$L\mathbf{1} = D^{\text{OUT}}\mathbf{1} - A\mathbf{1} = \begin{bmatrix} \deg_1^{\text{OUT}} \\ \vdots \\ \deg_i^{\text{OUT}} \end{bmatrix} - \begin{bmatrix} \deg_1^{\text{OUT}} \\ \vdots \\ \deg_i^{\text{OUT}} \end{bmatrix} = 0$$

i.e.,  $\lambda = 0$  is an eigenvalue of L and 1 is an associated eigenvector.

**Lemma.** Given a weighted digraph with Laplacian L, then all eigenvalues of L different from zero have strictly positive real part

**Lemma.** Given a weighted digraph with Laplacian L, the following statements are equivalent:

- 1. G is weight-balanced, i.e.  $D^{\text{IN}} = D^{\text{OUT}}$
- 2. 1L = 0

#### Theorem 3.3

A weighted digraph with Laplacian L contains a globally reachable node if and only if  $\lambda = 0$  is simple.

Corollary. If a weighted digraph is strongly connected, then  $\lambda = 0$  is simple

#### 3.6.2 Consensus for Laplacian dynamics

#### Theorem 3.4

let L be a (weighted) Laplacian matrix with associated strongly connected (weighted) digraph G. Consider the Laplacian dynamics  $\dot{x}(t) = -Lx(t), \ t \ge 0$ , then

1.

$$\lim_{t \to \infty} x(t) = \mathbf{1} \left( \frac{w^{\top} x(0)}{w^{\top} \mathbf{1}} \right)$$

with  $w^{\top}L=0$ , i.e. w is a left eigenvector for the eigenvalue  $\lambda=0$ ;

2. if additionally G is weight-balanced then

$$\lim_{t \to \infty} x(t) = \mathbf{1} \frac{\sum_{i=1}^{N} x_i(0)}{N}$$

# Optimization basics

#### Convexity and gradient monotonicity

If a convex function  $\ell$  is also differentiable, then its gradient  $\nabla \ell : \mathbb{R}^d \to \mathbb{R}^d$  satisfies

$$\left(\nabla \ell(z_A) - \nabla \ell(z_B)\right)^{\top} (z_A - a_B) \ge 0$$

for all  $z_A, z_B$ . That is, the gradient  $\nabla \ell$  is a monotone operator

#### Strict convexity and gradient monotonicity

A function  $\ell$  is strictly convex if for  $z_A \neq z_B$  and  $\theta \in (0,1)$ 

$$\ell(\theta z_A + (1 - \theta)z_B) < \theta \ell(z_A) + (1 - \theta)\ell(z_B)$$

If the strictly convex function  $\ell$  is also differentiable, then its gradient satisfies

$$\left(\nabla \ell(z_A) - \nabla \ell(z_B)\right)^{\top} (z_A - z_B) > 0$$

for all  $z_A, z_B$ . That is, the gradient  $\nabla \ell$  is a strictly monotone operator

### Strong convexity and gradient monotinicity

A function  $\ell$  is strongly convex with parameter  $\mu > 0$  if for  $z_A \neq z_B$  and  $\theta \in (0,1)$ 

$$\ell(\theta z_A + (1 - \theta)z_B) < \theta \ell(z_A) + (1 - \theta)\ell(z_B) - \mu \theta(1 - \theta) \|z_A - z_B\|^2$$

The gradient of a differentiable strongly convex function satisfies

$$\left(\nabla \ell(z_A) - \nabla \ell(z_B)\right)^{\top} (z_A - z_B) \ge \mu \|z_A - z_B\|^2$$

for all  $z_A, z_B$ . That is, the gradient  $\nabla \ell$  is a strongly monotone operator

#### Convexity and Lipschitz continuity of the gradient

Consider a differentiable convex function  $\ell$  with a Lipschitz continuous gradient with parameter L > 0, i.e.

$$\|\nabla \ell(z_A) - \nabla \ell(z_B)\| \le L\|z_A - z_B\|$$

for all  $z_A, z_B$ . Then, the following characterization holds

$$\left(\nabla \ell(z_A) - \nabla \ell(z_B)\right)^{\top} (z_A - z_B) \ge \frac{1}{L} \|\nabla \ell(z_A) - \nabla \ell(z_B)\|^2$$

for all  $z_A, z_B$ . That is, the gradient  $\nabla \ell$  is a co-coercive operator

#### Strong convexity and Lipschitz continuity of the gradient

Consider a strongly convex (with parameter  $\mu > 0$ ) function  $\ell$  with Lipschitz continuous gradient (with parameter L > 0). The the followin characterization holds

$$(\nabla \ell(z_A) - \nabla \ell(z_B))^{\top} (z_A - z_B) \ge \frac{\mu L}{\mu + L} \|z_A - z_B\|^2 + \frac{1}{\mu + L} \|\nabla \ell(z_A) - \nabla \ell(z_B)\|^2$$

for all  $z_A, z_B$ .

## 4.1 Optimization algorithms

We consider optimization algorithms based on iterative descent.

Notation. notatioWe denote by  $z^k \in \mathbb{R}^d$  the estimate at iteration  $k \in \mathbb{N}$  of a local minimum.

The algorithm starts at a given initial guess ecc ecc we know iterative descent

### 4.1.1 A system theoretical perspective to the gradient method

Let  $\ell$  be  $\mu$ -strongly convex and with L-Lipschitz continuous gradient. The gradient method is a discrete-time integrator in feedback interconnection with a static map

$$z^{k+1} = z^k - \alpha u^k,$$
  $z^0$  given  $u^k = \nabla \ell(z^k)$ 

where  $u^k$  is a "control input" obtained through a static (nonlinear) state feedback. The block diagram is This is known as the Lur'e problem

## 4.2 steady-state analysis of the gradient method

Assume that the state  $z^k$  converges to some value  $z_{eq}$ . Then, such an equilibrium must satisfy:

$$z_{eq} = z_{eq} - \alpha \nabla \ell(z_{eq}) \implies z_{eq} : \nabla \ell(z_{eq})$$
  
 $\implies z_{eq} = z^*$ 

Consider the change of coordinates  $z^k \to \tilde{z}^k - z_{eq} = z^k - z^*$ . Then, the error dynamics is

$$\tilde{z}^{k+1} = \tilde{z}^k - \alpha u^k$$
$$u^k = \nabla \ell (\tilde{z}^k + z^*) - \nabla \ell (z^*)$$

where  $u^k$  and  $\tilde{z}^k$  satisfy, in light of the assumption on  $\ell$ , the following inequality<sup>1</sup>

$$-(u^k)^T \tilde{z}^k \le -\gamma_1 \|\tilde{z}^k\|^2 - \gamma_2 \|u^k\|^2$$

Consider a Lyapunov function  $V: \mathbb{R}^d \to \mathbb{R}_{\geq 0}$  given by  $V(\tilde{z}) = \|\tilde{z}\|^2$ . Then

$$\begin{split} V(\tilde{z}^{k+1}) - V(\tilde{z}^k) &= \|\tilde{z}^{k+1}\|^2 - \|\tilde{z}^k\|^2 \\ &= \|\tilde{z}^k\|^2 - 2\alpha(u^k)^\top \tilde{z}^k + \alpha^2 \|u^k\|^2 - \|\tilde{z}^k\|^2 \\ &\leq -2\alpha\gamma_1 \|\tilde{z}^k\|^2 + \alpha(\alpha - 2\gamma_2) \|u^k\|^2 \end{split}$$

For a small enough stepsize  $\alpha$  (i.e.,  $\alpha \leq 2\gamma_2$ ), we can write

$$V(\tilde{z}^{k+1}) - V(\tilde{z}^k) < -2\alpha\gamma_1 \|\tilde{z}^k\|^2 \implies \|\tilde{z}^{k+1}\|^2 \le (1 - 2\alpha\gamma_1) \|\tilde{z}^k\|^2 \le (1 - 2\alpha\gamma_1)^k \|\tilde{z}^0\|^2$$

Therefore  $\{\tilde{z}^k\}_{k\in\mathbb{N}}$  goes exponentially/geometrically fast to zero

$$(\nabla \ell(z_A) - \nabla \ell(z_B))^{\top} (z_A - z_B) \ge \frac{\mu L}{\mu + L} \|z_A - z_B\|^2 + \frac{1}{\mu + L} \|\nabla \ell(z_A) - \nabla \ell(z_B)\|^2$$

<sup>&</sup>lt;sup>1</sup>For all  $z_A, z_B$  it holds that

### 4.2.1 Gradient method for quadratic programs

Consider a quadratic program

$$\min_{z} \frac{1}{2} z^{\top} Q z + r^{\top} z$$

With  $Q = Q^{\top} > 0$  The gradient method is an affine linear system

$$z^{k+1} = z^k - \alpha(Qz^k + r)$$
  $z^k$  given  
=  $(I - \alpha Q)z^k - \alpha r$ 

For a sufficiently small  $\alpha$ , the state matrix  $(I - \alpha Q)$  is Schur. Hence, the state trajectory is <sup>2</sup>

$$z^{k} = (I - \alpha Q)^{k} z^{0} - \alpha \sum_{i=0}^{k-1} (I - \alpha Q)^{i} r \to^{k \to \infty} -\alpha \left( \sum_{i=0}^{\infty} (I - \alpha Q)^{i} \right) r = -Q^{-1} r$$

### 4.2.2 Gradient flow

Let us swap the roles of the plant (the static nonlinearity) and the controller (the integrator) We obtain the so-called gradient flow (continuous-time dynamics)

$$\dot{z}(t) = -\nabla \ell(z(t)) \qquad z(0) = z_0$$

Remark. A solution to the ODE exists if the vector field is Lipschitz continuous

#### 4.2.3 Nesterov accelerated gradient method

Consider the following two-step algorithm: for all  $k \in \mathbb{N}$ 

$$\zeta^{k+1} = \zeta^k + \alpha_1(\zeta^k - \zeta^{k+1}) - \alpha_2 \nabla \ell \left( \zeta^k + \alpha_1(\zeta^k - \zeta^{k+1}) \right), \quad \zeta^0, \zeta^{-1} \text{ given}$$

for some  $\alpha_1, \alpha_2 > 0$ . It admits the state-space representation More general updates can also be considered: for all  $k \in \mathbb{N}$ 

$$\zeta^{k+1} = \zeta^k + \alpha_1(\zeta^k - \zeta^{k+1}) - \alpha_2 \nabla \ell \left( \zeta^k + \alpha_3(\zeta^k - \zeta^{k-1}) \right), \quad \zeta^0, \zeta^{-1} \text{ given}$$

for some  $\alpha_1, \alpha_2, \alpha_3 > 0$ 

<sup>&</sup>lt;sup>2</sup>The geometric series  $\sum_{i=0}^{\infty} \rho^i$  is equal to  $(1-\rho)^{-1}$  for all  $\rho < 1$ 

# Parallel Optimization and Federated Learning

#### Cost-coupled optimization for learning

In learning applications, we usually consider optimization problems in the form

$$\min_{z \in \mathbb{R}^d} \sum_{i=1}^N \ell_i(z)$$

where, for all  $i=1,\ldots,N$ , the cost function  $\ell_i:\mathbb{R}^d\to\mathbb{R}$  is local and private

#### (Batch) gradient method for learning

Consider the optimization problem

$$\min_{z} \sum_{i=1}^{N} \ell_i(z)$$

The (batch) gradient method is: for each iteration  $k \in \mathbb{N}$ 

$$z^{k+1} = z^k - \alpha \sum_{i=1}^{N} \nabla \ell_i(z^k)$$

Remark. computation can be expensive

# 5.1 Incremental gradient method

Consider the optimizatio problem

$$\min_{z} \sum_{i=1}^{N} \ell_i(z)$$

Idea: rather than using the whole batch gradient at each  $k \in \mathbb{N}$ , just select one single "sample" per iteration. The incremental gradient method is: for each iteration  $k \in \mathbb{N}$ 

$$z^{k+1} = z^k - \alpha \nabla \ell_{i^k}(z^k)$$

where  $i^k \in \{1, ..., N\}$  Two rules for choosing index  $i^k$  at iteration k:

- Cyclic rule
- Randomized rule

## 5.2 Stochastic Gradient Descent

consider the stochastic optimization problem

$$\min_{z} \mathbb{E}_{\mathcal{W}}[\ell(z, \mathcal{W})]$$

where  $\mathbb{E}_{\mathcal{W}}[\cdot]$  denotes the expected value with respect to the random variable  $\mathcal{W}$  (possibly having an uknown probability distribution  $p_{\mathcal{W}}(w)$ )

Remark. for all z, also  $\ell(z, \mathcal{W})$  is a random variable, whose probability distribution depends on  $p_{\mathcal{W}}$  and  $\ell$ . Moreover, the gradient  $\nabla \ell(z, \mathcal{W})$  at each z is a random quantity

Assumption: There exsists an oracle that, given a realization  $\bar{w}$  of  $\mathcal{W}$ , returns the corresponding realization of the gradient  $\nabla \ell(\bar{z}, \bar{w})$  at any query point  $\bar{z}$ 

The stochastic gradient descent is: for each iteration  $k \in \mathbb{N}$  draw a realization  $w^k$  of  $\mathcal{W}$  and update

$$z^{k+1} = z^k - \alpha \nabla \ell(z^k, w^k)$$

#### 5.2.1 Convergence

### Proposition 5.1

Assume:

- $\ell$  is a  $\mu$ -strongly convex function with L-Lipschitz continuous gradient (uniformly in its second argument)
- $\nabla \ell(z, \mathcal{W})$  is an unbiased estimate of  $\nabla_z \mathbb{E}_{\mathcal{W}}[\ell(z, \mathcal{W})]$
- $\|\nabla \ell(z, \mathcal{W})\| \leq M$  almost surely for some M > 0

# 5.3 Beyond SGD: Adaptive Momentum Estimation (Adam)

The ADAM algorithm reads as follows

• Mean and Variance (first momentum and second momentum)

$$m^{k+1} = \beta_1 m^k + (1 - \beta_1) \nabla \ell(z^k, w^k), \qquad \text{for some } \beta \in (0, 1)$$
$$v^{k+1} = \beta_2 v^k + (1 - \beta_2) [\nabla \ell(z^k, w^k)]^2, \qquad \text{for some } \beta_2 \in (0, 1)$$

where the square operation  $[\cdot]^2$  is meant component-wise

• Construct the descent direction

$$\begin{split} \hat{m} &= \frac{1}{1 - \beta_1^{k+1}} m^{k+1} \\ \hat{v} &= \frac{1}{1 - \beta_2^{k+1}} v^{k+1} \\ d^k &= -\frac{\hat{m}}{\sqrt{\hat{v} + \epsilon}}, \quad \text{ for some } \epsilon > 0 \end{split}$$

where the division in the last equation is meant element-wise

• Update of the solution estimate

$$z^{k+1} = z^k + \alpha d^k$$

<sup>&</sup>lt;sup>1</sup>A sequence of random variables  $\{\mathcal{X}_k\}_{k\in\mathbb{N}}$  converges almost surely to the rv  $\mathcal{X}$  if  $\mathbb{P}(\lim_{k\to\infty}\mathcal{X}_k=\mathcal{X})=1$ 

# 5.4 Federated learning

Consider the optimization problem

$$\min_{z} \sum_{i=1}^{N} \ell(z; \mathcal{D}^{i}, p^{i})$$

Paradigm:

- local private data  $\mathcal{D}^i = ([\mathcal{D}^{\flat}]_1, [\mathcal{D}^{\flat}]_2, \dots, [\mathcal{D}^{\flat}]_d)$  and  $p^i$
- learn common parameters  $z^* \in \mathbb{R}^d$  (common neural network)
- communication with a parameter server

# 5.5 Distributed learning

Consider the optimization problem

$$\min_{z} \sum_{i=1}^{N} \ell(z; \mathcal{D}^i, p^i)$$

Paradigm:

- local private data  $\mathcal{D}^i = ([\mathcal{D}^{\flat}]_1, [\mathcal{D}^{\flat}]_2, \dots, [\mathcal{D}^{\flat}]_d)$  and  $p^i$
- learn common parameters  $z^* \in \mathbb{R}^d$  (common neural network)
- communication with neighbours only

# Leader Follower networks: Formation control

#### 6.1 analogy with mass-spring systems

Consider a platoon of N masses such that each mass i is connected with mass i-1 and i+1 through a spring with elastic constants respectively  $a_{i-1,i} = a_{i,i-1} > 0$  and  $a_{i+1,i} = a_{i,i+1} > 0$ . Let  $x_i \in \mathbb{R}$  be the position of

The elastic force at mass i,  $F_{e,i}(x)$  is given by

$$F_{e,i}(x) = -a_{i,i-1}(x_i - x_{i-1}) - a_{i,i+1}(x_i - x_{i+1})$$

For each spring, we can write the associated elastic force as the negative gradient of the elastic (potential) energy, so that

$$F_{e,i}(x) = -\frac{\partial}{\partial x_i} \left( \frac{1}{2} a_{i,i-1} \|x_i - x_{i-1}\|^2 + \frac{1}{2} a_{i,i+1} \|x_i - x_{i+1}\|^2 \right)$$

Let us suppose that each spring can be written as the parallel of two springs with elastic constants respectively  $\frac{1}{2}a_{i,i-1} > 0$  and  $\frac{1}{2}a_{i,i+1} > 0$ .

Let  $x_i \in \mathbb{R}$  be the positino of mass i. The elastic force at mass i,  $F_{e,i}(x)$  is given by

$$F_{e,i} = -\left(\frac{1}{2}a_{i,i-1} + \frac{1}{2}a_{i,i-1}\right)(x_i - x_{i-1}) - \left(\frac{1}{2}a_{i,i+1} + \frac{1}{2}a_{i,i+1}\right)(x_i - x_{i+1})$$

As before, we can write the elastic force as the negative gradient of the elastic (potential) energy, i.e.

$$F_{e,i}(x) = -\frac{\partial}{\partial x_i} \left( \frac{1}{2} \frac{a_{i,i-1}}{2} \|x_i - x_{i-1}\|^2 + \frac{1}{2} \frac{a_{i,i-1}}{2} \|x_i - x_{i-1}\|^2 + \frac{1}{2} \frac{a_{i,i+1}}{2} \|x_i - x_{i+1}\|^2 + \frac{1}{2} \frac{a_{i,i+1}}{2} \|x_i - x_{i+1}\|^2 \right)$$

The total elastic (potential) energy of the mass-spring system can be written as

$$V(x) = \sum_{i=1}^{N} \sum_{j \in \mathcal{N}_i} \frac{1}{2} \frac{a_{i,j}}{2} ||x_i - x_j||^2$$
$$= \sum_{i=1}^{N} \sum_{j \in \mathcal{N}_i} V_{ij}(x_i, x_j)$$

where we have defined  $\mathcal{N}_i := \{i-1, i+1\}$  and  $V_{ij}(x_i, x_j) := \frac{1}{2} \frac{a_{i,j}}{2} \|x_i - x_j\|^2$ . Thus, the elastic force at mass i can be seen as the negative gradient of the energy, i.e.

$$F_{e,i}(x) = -\frac{\partial}{\partial x_i} \sum_{j \in \mathcal{N}_i} (V_{ij}(x_i, x_j) + V_{ji}(x_j, x_i))$$
$$= -\frac{\partial}{\partial x_i} V(x)$$

This formulation can be extended to more general systems in which masses are interconnected according to a topology described by an undirected graph  $G = (\{1, ..., N\}, E)$ 

By adding a damping term on each mass, the system dynamics can be written as

$$\dot{x}_i = v_i$$

$$m_i \dot{v}_i = -v_i - \frac{\partial}{\partial x_i} V(x) \quad \forall i \in \{1, \dots, N\}$$

where we have considered the damping coefficient equal to one.

If we assume that masses are small, we may written

$$v_i \approx -\frac{\partial}{\partial x_i} V(x)$$

so that the dynamics may be approximated by the following firt order dynamics

$$\dot{x}_i = -\frac{\partial}{\partial x_i} V(x) \quad \forall i \in \{1, \dots, N\}$$

Consider a network of N agents communicating/interacting according to a fixed, undirected graph G. Let  $x_i(t) \in \mathbb{R}^d$  be the state of agent i. Let agents run a Laplacian dynamics

$$\dot{x}_i = -\sum_{j \in \mathcal{N}_i} a_{ij} (x_i - x_j) \quad \forall i \in \{1, \dots, N\}$$

We can rewrite it as

$$\dot{x}_i(t) = -\sum_{j \in \mathcal{N}_i} \frac{\partial}{\partial x_i} \left( V_{ij}(x_i, x_j) + V_{ji}(x_j, x_i) \right) \quad \forall i \in \{1, \dots, N\}$$

with  $V_{ij}(x) = \frac{1}{2} \frac{a_{i,j}}{2} ||x_i - x_j||^2$ . By recalling the definition of the total energy

$$V(x) = \sum_{i=1}^{N} \sum_{j \in \mathcal{N}_i} V_{ij}(x_i, x_j)$$

the Laplacian dynamics

$$\dot{x} = -Lx$$

can be seen as a "gradient flow", i.e.

$$\dot{x} = -\nabla V(x)$$

Thus, the consensus configuration can be seen as a stationary point of V. This idea can be extended to general potential functions and applied to distributed control systems.

Consider a network of N autonomous agents communicating/interacting according to a fixed, undirected graph. Let  $x_i(t) \in \mathbb{R}^d$  be the state of agent i. Consider a global potential function defined as

$$V(x) = \sum_{i=1}^{N} \sum_{j \in \mathcal{N}_i} V_{ij}(x_i, x_j)$$

such that (local) minima of the potential correspond to desired configurations of the team. The gradient flow dynamics  $\dot{x} = -\nabla V(x)$  turns out to be distributed. That is,

$$\dot{x}_i(t) = \sum_{j \in \mathcal{N}} \frac{\partial}{\partial x_i} \left( V_{ij}(x_i, x_j) + V_{ji}(x_j, x_i) \right) \qquad \forall i \in \{1, \dots, N\}$$

We can define a desired formation by assigning a set of distances,  $d_{ij}$ , between neighbouring agents i and J in a suitable graph

The main idea for formation control is to define a potential function matching the sparsity of  $G, V^{\text{form}}(x) =$  $\sum_{i=1}^{N} \sum_{j \in \mathcal{N}_i} V_{ij}^{\text{form}}(i,j)$ , such that a configuration  $x^{textform}$  satisfying

$$||x_i^{\text{form}} - x_j^{\text{form}}|| = d_i j \quad \forall (i, j) \in E$$

is a minimum of V.

In order to reach a formation with assigned distances  $d_{ij}$ , let us define

$$V_{ij}^{\text{form}}(x) = \frac{1}{8} (\|x_i - x_j\|^2 - d_{ij}^2)^2$$

with corresponding (global) potential function  $V^{\text{form}}(x) = \sum_{i=1}^{N} \sum_{j \in \mathcal{N}_i} V_{ij}^{\text{form}}(x_i, x_j)$ .

The gradient flow dynamics of each agent i is given by

$$\dot{x}_i = -\sum_{j \in \mathcal{N}_i} \frac{\partial}{\partial x_i} \left( V_{ij}^{\text{form}}(x_i, x_j) + V_{ji}(x_j, x_i) \right) \qquad \forall i \in \{1, \dots, N\}$$

which reads as

$$\dot{x}_i = -\sum_{j \in \mathcal{N}_i} (\|x_i - x_j\|^2 - d_{ij}^2) (x_i - x_j) \qquad \forall i \in \{1, \dots, N\}$$

This dynamics has multiple equilibrium points, including the desired formation in which the agents are at the assigned distances. In particular, the consensual solution  $x_1 = x_2 = \cdots = x_N$  is an (undesired) equilibrium.

Such a "degenerate" equilibrium can be avoided by means of additional "collision avoidance" potential functions  $V_{ij}^{\text{ca}}(x_i, x_j)$  such that

$$\lim_{\|x_i - x_j\| \to 0} V_{ij}^{\text{ca}}(x_i, x_j) = +\infty$$

A possible solution is a barrier function given

$$V_{ij}^{\text{ca}} = -\log(x_i - x_j)$$

Similarly, barrier potential functions  $V^{\text{obs}}(x_i)$ , depending only on the state of agent  $x_i$ , can be used to avoid obstacles.

The formation control dynamics becomes

$$\dot{x}_i = -\frac{\partial V^{\text{form}}}{\partial x_i} - \frac{\partial V^{\text{ca}}(x)}{\delta x_i} - \nabla V^{\text{obs}}(x) \qquad \forall i \in \{1, \dots, N\}$$

# Distributed Aggregative Optimization

Consider N robots in the plane that want to optimize their positions  $z_i \in \mathbb{R}^2$ , for all i = 1, ..., N to perform multi-robot surveillance. Let:

- $r_0 \in \mathbb{R}^2$  be a target to protect
- $r_i \in \mathbb{R}^2$  be the intruder associated to robot i
- $\sigma(z) = \frac{1}{N} \sum_{i=1}^{N} z_i$  is the barycenter of the robots
- ullet Local cost function of robot i

$$\ell_i(z_i, \sigma(z)) = \gamma_i ||z_i - r_i||^2 + ||\sigma(z) - r_0||^2$$

with  $z \in \mathbb{R}^{2N}$  the stack of  $z_1, \dots, z_N$  and  $\gamma_i > 0$  being a tradeoff parameter.

# 7.1 Aggregative optimization

Let us consider aggregative optimization problems in the form

$$\min_{z_1, \dots, z_N} \sum_{i=1}^N \ell_i(z_i, \sigma(z))$$

where the aggregative variable  $\sigma(z)$  is defined as

$$\sigma(z) = \frac{1}{N} \sum_{i=1}^{N} \phi_i(z_i)$$

where

- $z = (z_1, \ldots, z_N)$ , with each  $z_i \in \mathbb{R}^{n_i}$ , for all  $i = 1, \ldots, N$
- $\ell_i: \mathbb{R}^{n_i} \times \mathbb{R}^d \to \mathbb{R}$  and  $\phi_i \mathbb{R}^{n_i} \to \mathbb{R}^d$ , for all  $i = 1, \dots, N$

For scalar states,  $z_i \in \mathbb{R}$ , the *centralized gradient* method at iteration k reads as

$$z_i^{k+1} = z_i^k - \alpha \frac{\partial}{\partial z_i} \left( \sum_{j=1}^N \ell_j(z_j, \sigma(z_1, \dots, z_N)) \right) \bigg|_{z_1 = z_1^k, \dots, z_N = z_N^k}$$

for all i = 1, ..., N where  $\alpha > 0$  is the stepsize

#### Gradient computation (scalar case)

Since the cost function is a composite function, we need the chain rule to compute its derivative with respect to  $z_i$ 

$$\begin{split} & \frac{\partial}{\partial z_i} \left( \sum_{j=1}^N \ell_j(z_j, \sigma(z_1, \dots, z_N)) \right) \bigg|_{z_1 = z_1^k, \dots, z_N = z_N^k} \\ &= \frac{\partial}{\partial z_i} \ell_i(z_i, \sigma) \bigg|_{z_i = z_i^k, \sigma = \frac{1}{N} \sum_{j=1}^N \phi_j(z_j^k)} \\ &+ \sum_{j=1}^N \frac{\partial}{\partial \sigma} l_j(z_j, \sigma) \bigg|_{z_j = z_j^k, \sigma = \frac{1}{N} \sum_{j=1}^N \phi_j(z_j^k)} \cdot \frac{\partial \sigma(z_1, \dots, z_N)}{\partial z_i} \bigg|_{z_1 = z_1^k, \dots, z_N = z_N^k} \end{split}$$

Notice that  $\frac{\partial \sigma(z_1,\ldots,z_N)}{\partial z_i} = \frac{1}{N} \frac{d}{dz_i} \phi_i(z_i)$  can be computed locally As in the scalar case, we use the chain rule to compute the gradient of the composite function. The *i*-th

block of the gradient, denoted as 
$$\left[\nabla\left(\sum_{j=1}^N \ell_j(z_j, \sigma(z_1, \dots, z_N))\right)\right]_i \in \mathbb{R}^{n_i}$$
, is given by

$$\left[ \nabla \left( \sum_{j=1}^{N} \ell_j(z_j, \sigma(z_1, \dots, z_N)) \right) \right]_i 
= \nabla_1 \ell_i(z_i, \sigma)|_{z_j = z_j^k, \sigma = \frac{1}{N} \sum_{j=1}^{N} \phi_j(z_j^k)} 
+ \left. \frac{1}{N} \nabla \phi_i(z_i) \right|_{z_i = z_i^k} \cdot \sum_{j=1}^{N} \nabla_2 \ell_j(z_j, \sigma) \right|_{z_i = z_i^k, \sigma = \frac{1}{N} \sum_{i=1}^{N} \phi_i(z_i^k)}$$

#### 7.2 Distributed aggregative optimization

In a distributed contex, each agent i

- knows only  $\ell_i$  and  $\phi_i$
- maintains an estimate  $z_i^k$  of  $z_i^*$
- maintains an estimate  $s_i^k$  of  $\phi(z^k) = \frac{1}{N} \sum_{i=1}^N \phi_j(z_j^k)$
- maintains an estimate  $v_i^k$  of  $\sum_{i=1}^N \nabla_2 \ell_j(z_j^k, \sigma(z^k))$

The "tracking" idea of gradient tracking algorithm is applied to aggregative optimization

$$z_{i}^{k+1} = z_{i}^{k} - \alpha \left( \nabla_{1} \ell_{i}(z_{i}^{k}, s_{i}^{k}) + \nabla \phi_{i}(z_{i}^{k}) v_{i}^{k} \right) \qquad z_{i}^{0} \in \mathbb{R}^{n_{i}}$$

$$s_{i}^{k+1} = \sum_{j \in \mathcal{N}_{i}} a_{ij} s_{j}^{k} + \phi_{i}(z_{i}^{k+1}) - \phi_{i}(z_{i}^{k}) \qquad s_{i}^{0} = \phi_{i}(z_{i}^{0})$$

$$v_{i}^{k+1} = \sum_{j \in \mathcal{N}_{i}} a_{ij} v_{j}^{k} + \nabla_{2} \ell_{i}(z_{i}^{k+1}, s_{i}^{k+1}) - \nabla_{2} \ell_{i}(z_{i}^{k}, s_{i}^{k}) \qquad v_{i}^{0} = \nabla_{2} \ell_{i}(z_{i}^{0}, s_{i}^{0})$$

**Theorem 7.1** (aggregative tracking distributed optimization algorithm: convergence)

Assume G is a strongly connected and aperiodic digraph, and A is doubly stochastic. Assume that each function  $\ell_i$  is strongly convex, the gradients  $\nabla_1 \ell_i$  and  $\nabla_2 \ell_i$  are Lipschitz continuous, and  $\phi_i$  is differentiable and Lipschitz continuous.

Then, there exists  $\alpha^*$  such that for all  $\alpha \in (0, \alpha^*)$  the sequences of local solution estimates  $\{z_1^k, \dots, z_N^k\}_{k \in \mathbb{N}}$ generated by the aggregative tracking distributed opitmization algorithm satisfy

$$\lim_{k \to \infty} \|z_i^k - z_i^\star\| = 0$$

at a linear rate, for all i = 1, ..., N

#### 7.2.1Extension to online aggregative optimization

Consider a time-varying instance of the problem

$$\min_z \sum_{i=1}^N \ell_i^k(z_i,\sigma^k(z))$$
 subj. to  $z_i \in Z_i^k, \qquad \forall i=1,\dots,N$ 

where  $Z_i^k \subset \mathbb{R}^{n_i}$  are local constraint sets. The goal is to design an algorithm generating a sequence  $\{z_i^k\}_{k\in\mathbb{N}}$  that "tracks" the solution  $z^{k,\star}=(z_1^{k,\star},\ldots,z_N^{k,\star})$  of the k-th problem instance.

Remark. A regret  $R_K$  can be introduced for the analysis. Under suitable assumptions, it can be proven that

# Learning with Neural Networks

In supervised learning we have labeled data whose generic element(sample) is composed by

- an input point  $\mathcal{D} \in \mathbb{R}^d$ , i.e. a vector  $\mathcal{D} = ([\mathcal{D}]_1, \dots, [\mathcal{D}]_d)$
- a lebel  $p \in \mathbb{R}$  associated to the input point

A dataset is usually made by  $\mathcal{M}$  samples, i.e.  $((\mathcal{D}^1, p^1), \dots, (\mathcal{D}^{\mathcal{M}}, p^{\mathcal{M}}))$ 

Goal: reconstruct the uknown input-output map between inputs and labels

Strategy: approximate the map as a nonlinear function  $\phi(\underline{\approx};\cdot):\mathbb{R}^d\to\mathbb{R}$  parametrized by  $\underline{\approx}$ 

Problem: find the best parameters  $\cong^*$  based on data

## 8.1 Basic element: the neuron model

A generic neuron h is a computational unit that

- has a set of weights  $u_h \in \mathbb{R}^d$
- elaborates a vector  $x \in \mathbb{R}^d$
- computes a scalar quantity

$$x_h^+ = \phi(x^\top u_h)$$

with  $\sigma: \mathbb{R} \to \mathbb{R}$  being the activation function

# Multi-Robot Safety Controllers

Consider a nonlinear system

$$\dot{x}(t) = f(x(t), u(t)), \qquad x \in \mathbb{R}^n \quad u \in U \subset \mathbb{R}^m$$

Let us define a safe (state) set

$$X^s := \{ x \in \mathbb{R}^n | V^s(x) \ge 0 \}$$

for some sufficiently regular function  $V^s: \mathbb{R}^n \to \mathbb{R}$ 

Goal: design a feedback control law  $\kappa(x), \kappa: \mathbb{R}^n \to \mathbb{R}^m$  such that the set  $X^s$  is forward invariant

### 9.1 Control Barrier Functions

Consider the time derivative of  $V^s(x(t))$  along the system trajectories

$$\frac{d}{dt}V^{s}(x(t)) = \nabla V^{s}(x(t))^{\top} f(x(t)) + \sum_{h=1}^{m} \nabla V^{s}(x(t))^{\top} g_{h}(x(t)) u_{h}(x(t))$$
$$= L_{f}V^{s}(x(t)) + L_{g}V^{s}(x(t)) u(t)$$

where  $L_f$  and  $L_g$  represent Lie derivatives. We say that  $V^s$  is a Control Barrier Function (CBF) if there exists a continuous, strictly increasing function  $\gamma : \mathbb{R} \to \mathbb{R}$ , with  $\gamma(0) = 0$ , such that

$$\sup_{u \in U} \{ L_f V^s(x) + L_g V^s(x) u + \gamma(V^s(x)) \ge 0 \}$$

for all  $x \in \mathbb{R}^n$ .

The above inequality induced by a CBF is called a control barrier certificate.

#### 9.1.1 Admissible Control Space

We can define the set of admissible (safe) controllers for a given state x as

$$U^{s}(x) = \{ u \in \mathbb{R}^{m} | L_{f}V^{s}(x) + L_{q}V^{s}(x)u + \gamma(V^{s}(x)) \ge 0 \}$$

i.e., the set of inputs satisfying the control barrier certificates.

Remark. It can be shown that (sufficiently regular) feedback controllers satisfying the above condition render  $X^s$  forward invariant and asymptotically stable.

## 9.1.2 Safety Filters via Control Barrier Certificates

Let  $u^{\text{ref}} \in \mathbb{R}^m$  be a (possibly unsafe) reference input, e.g., obtained by a higher-level controller. The safety controller is designed to be "minimally invasive", i.e. it aims at altering the reference controller as little as possible. A safe policy  $u = \kappa(x)$  can be designed as

$$\kappa(x) = \arg\min_{u \in \mathbb{R}^m} \|u - u^{\text{ref}}(x)\|^2$$
 subj. to  $-L_f V^s(x) - L_g V^s(x) u - \gamma(V^s(x)) \le 0$