# A Fully Asynchronous Newton Tracking Method for Decentralized Optimization

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Abstract—We consider fully asynchronous decentralized optimization over a directed graph. While various algorithms have been proposed, real-world applications require relaxing the assumption and considering communication networks with asynchronous and heterogeneous nodes. To meet these challenges, we propose an efficient and robust Newton tracking mechanism for fully asynchronous optimization. Our proposed mechanism can be adapted to different asynchronous first-order methods as required by the practical context. Through the theoretical analysis we demonstrate the R-linear rate of our method and derive an explicit expression of decaying factor under local conditions. Furthermore, numerical comparison with existing algorithms support the efficiency and robustness of our method.

# I. INTRODUCTION

In this paper, we consider a fully asynchronous decentralized optimization problem, where n agents cooperate to minimize the average of local functions

$$x^* = \arg\min_{x \in \mathbb{R}^d} F(x) \coloneqq \frac{1}{n} \sum_{i=1}^n f_i(x), \tag{1}$$

where  $x \in \mathbb{R}^d$  is the decision variable and  $f_i : \mathbb{R}^d \to \mathbb{R}$  is the loss function privately owned by node *i*. The optimization problem (1) has applications in various fields, such as deep learning [1], sensor networking [2], statistical learning [3], and optimal transport [4]. To leverage the benefits of decentralized storage of data and parallel computing power of nodes, decentralized algorithms for solving problem (1) have been extensively studied, especially for the design of synchronous algorithms, such as DGD [5], EXTRA [6], gradient tracking [7] based on undirected graphs, and Push-Sum [8] and Push-Pull [9] based on directed graphs.

However, due to the heterogeneity of data and systems, and the impact of uncontrollable factors such as communication delays or errors, it is essential to design asynchronous algorithms. Asynchronous algorithms can reduce waiting time, alleviate communication costs, and make the algorithm more fault-tolerant. This paper considers the *fully* asynchronous scenario: agents can use local and neighbor information (possibly stale) for local computation at any time and send the results to neighbors without any synchronous coordination and scheduling. The difference between synchronous and our fully asynchronous settings is depicted in Fig. 1. Specifically, in the synchronous setting, all nodes perform computation and communication at the same time per iteration, while



Fig. 1. Comparison of Synchronous and Asynchronous Setting.

in the fully asynchronous setting, nodes can start at any time. We can observe that asynchronous algorithms are more efficient, as they can perform more iterations within the same time period.

Asynchronous Decentralized Methods. In recent years, there has been a noticeable shift in research focus towards asynchronous decentralized optimization methods. Some of these methods utilize random activation mechanisms that satisfy specific distributions, where a randomly selected group of agents are activated per iteration to perform computation and communicate with their neighbors [10]–[12]. Some works propose asynchronous algorithms that are robust to communication delays and investigate the effect of delay on convergence [13], [14]. However, these algorithms still require some degree of coordination among nodes and are not fully asynchronous. Several algorithms have been developed based on a fully asynchronous model [15]-[18], where agents can use local and neighbor information (possibly delayed) for local computation at any time and send the results to neighbors without any form of coordination or centralized scheduling. The asynchronous Gradient-Push (AGP) method proposed in [15] uses the buffer mechanism to store delayed information and is proven to converge to a neighborhood of the optimal solution. However, as shown in [18], AGP cannot converge precisely due to the uneven update frequency of each node. To tackle this issue, [18] introduces an auxiliary variable to adjust the update frequency of each node and proposes a subgradient-type method with exact convergence. However, the decaying stepsize rule makes convergence rate sub-linear. To improve the convergence rate, [16] proposes an

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asynchronous algorithm over a directed network, where each node estimates the global gradient using the idea of gradient tracking [19]. Inspired by Push-Sum [20], the method in [16] gets a linear convergence approaching to the optimal point. The APPG method proposed in [17] is based on the Push-Pull algorithm over a directed network. It uses an augmented graph to model the communication of the asynchronous algorithm and shows that APPG converges to the optimal solution at a linear rate. Apart from firstorder methods, some works study second-order methods that use curvature information to accelerate convergence. The work [21] adapts an approximate second-order information scheme under a partial asynchronous setting and can handle uncoordinated activation and packet loss. However, the work [21] requires that only one agent wakes up at each iteration, which requires a global clock. [12], [22] studied the asynchronous decentralized second-order algorithm, while they assume the activation of the node follows Poisson Process, this assumption make convergence analysis easier, but there is a strong dependence of the delays on the activation index which often contrasts the case in reality. To the best of our knowledge, there is currently no known decentralized second-order algorithm specifically designed for a fully asynchronous setting.

#### A. Major Contribution

In this paper, we propose an efficient and robust asynchronous newton tracking mechanism for fully asynchronous decentralized optimization. The main contributions are summarized as follows:

- We propose a novel newton-tracking mechanism for fully asynchronous optimization, which can be easily adapted in existing asynchronous first-order algorithms. Our mechanism enhances the performance and robustness of existing asynchronous first-order algorithms by incorporating global second-order information.
- Through our theoretical analysis, we have clearly shown that our algorithm enjoys a linear convergence rate and derived the explicit expression for decaying factor that surpasses that of existing asynchronous first-order algorithms under local conditions.
- Our numerical comparisons have further demonstrated the considerable improvements in efficiency and robustness of our mechanism when compared to both asynchronous first-order algorithms and synchronous second-order algorithms.

# B. Notation

We use k to denote the index of iteration and t(k) to denote the time of k-th iteration. We use  $\|\cdot\|$  to denote the Euclidean norm of a vector or the largest singular value of a matrix,  $\|\cdot\|_F$  to denote the Frobenius norm. We use [n] to denote the numbers from 1 to n. For the aggregated variable, we define  $\boldsymbol{x} = [x_1; \ldots; x_n] \in \mathbb{R}^{nd}$ ,  $\nabla f(\boldsymbol{x}^k) = [\nabla f_1(x_1^k); \ldots; \nabla f_n(x_n^k)], \boldsymbol{H}^k = [H_1^k; \ldots; H_n^k]$ other aggregated variables are defined similarly. For the average variable, the average decision variable over all nodes at the iteration k is defined as  $\bar{\boldsymbol{x}}^k = \frac{1}{n} \sum_{i=1}^n \boldsymbol{x}_i^k$ ,  $\overline{\nabla f}(\boldsymbol{x}^k) = \frac{1}{n} \sum_{i=1}^n \nabla f_i(\boldsymbol{x}_i^k) \in \mathbb{R}^d$ ,  $\nabla F(\bar{\boldsymbol{x}}^k) = \frac{1}{n} \nabla f_i(\bar{\boldsymbol{x}}^k) \in \mathbb{R}^d$ ,  $\bar{H}^k = \frac{1}{n} \sum_{i=1}^n H_i^k$ , other average variables are defined similarly. Given a buffer  $\mathcal{B}$ ,  $\operatorname{avg}(\mathcal{B})$  and  $\operatorname{sum}(\mathcal{B})$  return the average and return the sum of variables in the buffer, respectively.

## C. Organization

The remaining sections of this document are structured in the following manner. We provide the problem setting and proposed asynchronous newton tracking mechanism along with its implementations in Section II. We establish the convergence analysis in Section III and present the numeric examples in Section IV. Section V serves as the concluding part of the paper.

# **II. PROBLEM SETTING AND ALGORITHM DEVELOPMENT**

In this section, we give a detailed description of the problem setting, then we develop a newton tracking mechanism for asynchronous decentralized optimization and adapt the mechanism over existing asynchronous first-order methods.

# A. Problem Setting

We consider a directed graph  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$ , where  $\mathcal{V} = \{1, \ldots, n\}$  is the set of *n* nodes with  $|\mathcal{V}| = n$  and  $\mathcal{E}$  is the set of all edges. If  $(i, j) \in \mathcal{E}$ , it indicates that *i* can send message to *j*. For a node *i*, we defined in-neighbors which send information to node *i* as  $\mathcal{N}_{in}^i = \{j \mid (j, i) \in \mathcal{E}\}$ , and out-neighbors which receive information from node *i* as  $\mathcal{N}_{out}^i = \{j \mid (i, j) \in \mathcal{E}\}$ . We give the general assumptions for decentralized optimization [9].

Assumption 2.1: The communication network G is strongly connected.

Then we give some assumptions on objective functions.

Assumption 2.2: The entire objective function F is  $\mu$ -strongly convex for some constant  $\mu > 0$ , i.e.,

$$\nabla^2 F(x) \succeq \mu I_d, \forall x \in \mathbb{R}^d, \tag{2}$$

where  $\mu$  is the strong convexity parameter.

Assumption 2.3: Each  $f_i$  is twice-differentiable, and both the gradient and Hessian are Lipschitz continuous, i.e.,

$$\|\nabla f_i(x) - \nabla f_i(y)\| \le L_1 \|x - y\|, \forall x, y \in \mathbb{R}^d, \quad (3)$$

and

$$\|\nabla^2 f_i(x) - \nabla^2 f_i(y)\| \le L_2 \|x - y\|, \forall x, y \in \mathbb{R}^d,$$
 (4)

where  $L_1 \ge 0$  and  $L_2 \ge 0$  are the Lipschitz constants of the local gradient and local Hessian, respectively.

Finally we give a common assumption on transmission delays.

Assumption 2.4: For any  $(i, j) \in \mathcal{E}$ , the transmission delay from any two nodes is bounded by a constant D satisfy  $0 < D < \infty$ .

### B. Proposed Asynchronous Newton Tracking Mechanism

To model the process of sending messages between nodes in the network, we introduce an induced matrix  $W \in \mathbb{R}^{n \times n}$ with entries  $w_{ij}$ . The matrix  $W \in \mathbb{R}^{n \times n}$  is non-negative where  $w_{ji} > 0$  if and only if node *i* can send information to node *j*, and the value of  $w_{ji}$  represents the weight assigned by node *i* to the information sent to node *j*. To estimate the global Hessian matrix, one common method is to use the dynamic average consensus method (DAC) [19], [23]. We first present DAC method under synchronization, showing clearly that this synchronous method cannot be directly applied to the asynchronous case. Then we propose our fully asynchronous Newton tracking mechanism. To this end, the synchronous DAC method is

$$H_{i}^{k+1} = H_{i}^{k} - \gamma \sum_{j \in \mathcal{N}_{\text{in}}^{i}} w_{ij} \left( H_{i}^{k} - H_{j}^{k} \right) + \nabla^{2} f_{i} \left( x_{i}^{k+1} \right) - \nabla^{2} f_{i} \left( x_{i}^{k} \right),$$

$$\tag{5}$$

where  $H_i^k$  hold on node *i* is the estimation of the global Hessian matrix at time step *k*. Under synchronization, to estimate the global Hessian, we require that *W* is column stochastic, i.e.,  $\sum_{i=1}^{n} w_{ij} = 1, \forall j \in [n]$  [9], [24]. With this condition, by taking the average of both sides of (5) over all the nodes, we have

$$\frac{1}{n}\sum_{i=1}^{n}H_{i}^{k} = \frac{1}{n}\sum_{i=1}^{n}\nabla^{2}f_{i}(x_{i}^{k}), \quad \forall k,$$
(6)

where we use the fact that W is column stochastic and the initialization  $H_i^0 = \nabla^2 f_i(x_i^0)$ ,  $\forall i \in [n]$ . The property (6) is very important for each node to successfully track the global Hessian. However, in the asynchronous setting, node i may fail to receive  $H_j^k$  from node j at time step k due to the uncoordinated activations and staled information. This causes the property (6) to no longer hold, thus affecting the performance of the algorithm.

To tackle this issue, we use an augmented graph to store the delayed information from the neighbors before it is used [25], [26]. An example of an augmented graph is given in Fig.2, the left of the Fig.2 is the original graph including two nodes i and j. On the right is the augmented graph, where node j sends  $\tilde{H}_j^k$  to node i at some time t(k) and node i uses it at some time t(k + 3). Two matrices are corresponding induced matrices W and  $\tilde{W}$  of the original graph and the augmented graph, respectively. Let  $\hat{n} = n(D + 1)$  be the number of nodes in the augmented graph,  $\tilde{H}_i^k$  be the Hessian information on node  $i \in [\hat{n}]$  at step k, and  $\tilde{W} \in \mathbb{R}^{\hat{n} \times \hat{n}}$  be the induced matrix of the augmented graph. It has been shown that  $\tilde{W} \in \mathbb{R}^{\hat{n} \times \hat{n}}$  is column stochastic [17]. Thus, for the augmented graph we have

$$\frac{1}{\hat{n}}\sum_{i=1}^{\hat{n}}\tilde{H}_{i}^{k} = \frac{1}{n}\sum_{i=1}^{n}\nabla^{2}f(x_{i}^{k}).$$
(7)

This means that we still have the key property (7) in the asynchronous setting similar to (6) in the synchronous setting through the augmented graph. Note that the mixing of the Hessian information on node i, weighted by the i-th row



Fig. 2. Augmented graph and induced matrix

of the induced matrix in the augmented graph, is equivalent to introducing a buffer to store  $\frac{1}{|\mathcal{N}_{out}^j|}H_j$  and summing over the Hessian information in the buffer. For simplicity, we use a buffer to implement our asynchronous Newton tracking mechanism, as shown in Algorithm 1.

Algorithm 1 Proposed Asynchronous Newton Tracking Mechanism on node *i* 

**Initialize**:  $\gamma > 0$ .  $H_i^0 = \nabla^2 f_i(x_i^0)$ . Create local buffers  $\mathcal{H}_i$ . Broadcast  $\tilde{H}_i^0 \coloneqq \frac{H_i^0}{|\mathcal{N}_{\text{out}}^i|}$  to out-neighbors.

Repeat

Keep receiving  $\tilde{H}_j$  from  $j \in \mathcal{N}_{in}^i$  and add it to the local buffer  $\mathcal{H}_i$ , until node *i* is activated.

if node i is activated then

 $\begin{array}{l} H_{i}^{k+1} = (1-\gamma)H_{i}^{k} + \gamma \mathbf{sum}(\mathcal{H}_{i}) + \nabla^{2}f_{i}(x_{i}^{k+1}) - \nabla^{2}f_{i}(x_{i}^{k}) \\ \text{Empty local buffer } \mathcal{H}_{i} \text{ and broadcast } \tilde{H}_{i}^{k+1} = \frac{H_{i}^{k+1}}{|\mathcal{N}_{\text{out}}|} \\ \text{to its out-neighbors.} \end{array}$ 

end if Output  $H_i^{k+1}$ 

# C. Implementations of Proposed Asynchronous Newton Tracking Mechanism

In this section, we combine the proposed mechanism with two existing asynchronous first-order algorithms, APPG [17] and ASY-SONATA [16], resulting in two asynchronous Newton algorithms which converge faster than the original APPG and SONATA. After that, we provide the communicationefficient implementation of the proposed Mechanism, which uses compression to avoid sending full matrices. Note that the existence of Hessian approximation in the algorithms makes convergence analysis challenging, and the analysis of APPG and SONATA is no longer applicable.

1) APPG with Newton Tracking: The first algorithm is based on APPG [17], The first algorithm is based on APPG, which is an asynchronous robust algorithm developed from the push-pull algorithm [9]. APPG employs the DAC method to track the global gradient and use a buffer to store unused information that has been sent to reduce the impact of asynchrony on convergence. After each node obtains an approximation of the global gradient, we leverage our

# Algorithm 2 APPG with Newton Tracking

**Initialize:**  $\alpha > 0, M > 0, x_i^0, d_i^0 = y_i^0 = \nabla f_i(x_i^0), H_i^0 = \nabla^2 f_i(x_i^0)$ . Create local buffers  $\mathcal{X}_i, \mathcal{Y}_i, \mathcal{H}_i$ . Broadcast  $\tilde{x}_i^0 \coloneqq x_i^0, \tilde{y}_i^0 \coloneqq \frac{y_i^0}{|\mathcal{N}_{out}^i|}, \tilde{H}_i^0 \coloneqq \frac{H_i^0}{|\mathcal{N}_{out}^i|}$  to out-neighbors **Repeat** 

Keep receiving  $\tilde{x}_j, \tilde{y}_j, \tilde{H}_j$  from  $j \in \mathcal{N}_{in}^i$  and add them to the local buffer  $\mathcal{X}_i, \mathcal{Y}_i, \mathcal{H}_i$ , respectively. if node *i* is activated **then** 

In node *i* is activated then  

$$\begin{aligned} x_i^{k+1} &= \mathbf{avg}(\mathcal{X}_i) - \alpha d_i^k \\ y_i^{k+1} &= \mathbf{sum}(\mathcal{Y}_i) + \nabla f_i(x_i^{k+1}) - \nabla f_i(x_i^k) \\ \text{Apply algorithm 1 to get } H_i^{k+1} \\ d_i^{k+1} &= (H_i^{k+1} + MI_d)^{-1} y_i^{k+1} \\ \text{Empty buffer and broadcast } \tilde{x}_i^{k+1} &= x_i^{k+1}, \, \tilde{y}_i^{k+1} = \frac{y_i^{k+1}}{|\mathcal{N}_{\text{out}}|}, \, \tilde{H}_i^{k+1} &= \frac{H_i^{k+1}}{|\mathcal{N}_{\text{out}}|} \text{ to out-neighbors} \\ \text{end if} \\ \text{until stopping criterion} \end{aligned}$$

proposed mechanism to generate an approximation of the global Hessian, allowing each node to approximate the global Newton direction. This direction, with its inclusion of curvature information, is capable of accelerating convergence. The resulting algorithm is summarized in Algorithm 2.

2) ASY-SONATA with Newton Tracking: We next combine the proposed asynchronous mechanism with ASY-SONATA [16] which presents a method for robustly estimating the global gradient asynchronously. In contrast to APPG that is developed from the push-pull algorithm, ASY-SANATA is based on the push-sum algorithm [6]. By integrating our Newton tracking mechanism, we enable each node to generate an approximation of the global Newton direction and get a second-order version of ASY-SONATA, as outlined in Algorithm 3.

Algorithm 3 Asy-SONATA with Newton Tracking

**Initialize:**  $\alpha > 0$ , M > 0,  $x_i^0$ ,  $d_i^0 = y_i^0 = \nabla f_i(x_i^0)$ ,  $\rho_i^0 = 0$ ,  $H_i^0 = \nabla^2 f_i(x_i^0)$ . Create local buffers  $\mathcal{X}_i$ ,  $\mathcal{H}_i$ ,  $\tilde{\rho}_i$ . Broadcast  $\tilde{x}_i^0 \coloneqq x_i^0$ ,  $\rho_i^0$ ,  $\tilde{H}_i^0 \coloneqq \frac{H_i^0}{|\mathcal{N}_{out}^i|}$  to out-neighbors **Repeat** 

Keep receiving  $\tilde{x}_j$ ,  $\rho_j$ ,  $\tilde{H}_j$  from  $j \in \mathcal{N}_{in}^i$ , replace buffer variables  $\tilde{x}_{ij}$  with  $\tilde{x}_j$ , add  $\tilde{H}_j$  to  $\mathcal{H}_i$ .

if node i is activated then

$$\begin{aligned} x_i^{k+1} &= x_i^k - \alpha d_i^k \\ x_i^{k+1} &= \frac{1}{|\mathcal{N}_{\text{in}}^i|} x_i^{k+1} + \sum_{j \in \mathcal{N}_{\text{in}}^i} \frac{1}{|\mathcal{N}_{\text{in}}^i|} \tilde{x}_{ij} \\ \epsilon^{k+1} &= \nabla f(x_i^{k+1})) - \nabla f(x_i^k)) \\ y_i^{k+1} &= \frac{1}{|\mathcal{N}_{\text{out}}^i|} (y_i^k + \sum_{j \in \mathcal{N}_{\text{in}}^i} (\rho_{ij} - \tilde{\rho}_{ij}) + \epsilon^{k+1} \\ \rho_i^{k+1} &= \rho_i^k + y_i^{k+1} \\ \text{replace buffer variables } \tilde{\rho}_{ij} \text{ with } \rho_j \\ \text{Apply algorithm 1 to get } H_i^{k+1} \\ d_i^{k+1} &= (H_i^{k+1} + MI_d)^{-1} y_i^{k+1} \\ \text{Empty } \mathcal{H}_i \text{ and broadcast } x_i^{k+1}, \ \rho_i^{k+1}, \ \tilde{H}_i^{k+1} := \frac{H_i^{k+1}}{|\mathcal{N}_{\text{out}}^i|} \text{ to out-neighbors} \\ \text{end if } \\ \text{until stopping criterion} \end{aligned}$$

#### **III. CONVERGENCE ANALYSIS**

In this section, we provide a unified analysis for Algorithm 2 and Algorithm 3. For the sake of clarity, we provide the compact form over the augmented graph for both algorithms.

$$\widetilde{\boldsymbol{x}}^{k+1} = \widetilde{A}^{k} (\widetilde{\boldsymbol{x}}^{k} - \alpha (\widetilde{\boldsymbol{H}}^{k} + MI)^{-1} \widetilde{\boldsymbol{y}}^{k}), \\ \widetilde{\boldsymbol{y}}^{k+1} = \widetilde{B}^{k} \widetilde{\boldsymbol{y}}^{k} + \nabla (k+1) - \nabla (k),$$
(8)

where

7

$$\begin{split} \widetilde{\boldsymbol{x}}^{k} &= [x^{k}; x^{k}(1); \cdots; x^{k}(D)] \in \mathbb{R}^{\hat{n}d} \\ \widetilde{\boldsymbol{y}}^{k} &= [y^{k}; y^{k}(1); \cdots; y^{k}(D)] \in \mathbb{R}^{\hat{n}d} \\ \widetilde{\boldsymbol{H}}^{k} &= [H^{k}; H^{k}(1); \cdots; H^{k}(D)] \in \mathbb{R}^{\hat{n}d \times d} \\ \nabla(k) &= [\nabla f(\widetilde{\boldsymbol{x}}^{k}); \boldsymbol{0}_{(\hat{n}-n)d}], \end{split}$$

and  $\widetilde{A}^k$  is a row-stochastic matrix and  $\widetilde{B}^k$  is a columnstochastic matrix which are already defined in [17]. It is worth mentioning that  $\widetilde{A}^k$  and  $\widetilde{B}^k$  are time-varying matrices depending on the way of delay in each timestamp. Through the proof, we assume d = 1 (scale variables), which is quite standard in asynchronous optimization, see [16], [17].

Inspired by work [16], we provided the convergence analysis for (8) with Theorem 3.2 below under local condition, which establishes linear convergence for our mechanism. We leave global convergence analysis for further discussions.

Assumption 3.1: For local convergence analysis, we assume each node is close to optimal point  $\mathbf{x}^*$ . Specifically, we have (9) for  $k \ge N_1$ .

$$\|H_{i}^{k} - H^{*}\|_{F} \le r, \forall i \in [n],$$
(9)

where  $H^*$  is the hessian matrix of the optimal point  $\mathbf{x}^*$ .

Theorem 3.2: Let  $\mathbf{x}^*$  denote the unique optimal solution of (1) and  $\{(\mathbf{x}_i^k)_{i=1}^n\}_k$  be the sequence generated by (8). With prior assumptions hold, there exists a constant  $\bar{\alpha} < 1$ such that when  $\alpha \leq \bar{\alpha}$ , the asynchronous algorithms holds

$$\left\|\mathbf{x}^{k} - \mathbf{1}_{n} \otimes \mathbf{x}^{\star}\right\| = \mathcal{O}\left(\lambda^{k}\right) \tag{10}$$

where  $\lambda \in (0, 1)$  is given by (20).

*Remark 3.3:* By properly choosing the stepsize, we can show that our decay factor  $\lambda$  is better than the factor of first-order methods derived in [16]. It indicates that our algorithm enjoys a sharper local linear convergence rate.

*Proof:* Build upon on prior idea in [16], we take the advantage of the Hessian approximation's curvature information to further bound the error quantities. Specifically, we utilize the locally quadratic convergence rate of the centralized Newton's method to bound the optimization error.

#### A. Step I

We adapt a unified model to study the dynamics of the consensus and optimization errors of the asynchronous framework, which consists in pulling out the tracking update and treating the variables term  $-\alpha (\tilde{H}^k + MI)^{-1}\tilde{y}^k$  as an exogenous perturbation  $\delta^k$ . Our framework has the following scheme in compact form as

$$\widetilde{\boldsymbol{x}}^{k+1} = \widetilde{A}^k (\widetilde{\boldsymbol{x}}^k + \delta^k), \tag{11}$$

Applying (11) recursively, we have

$$\widetilde{\boldsymbol{x}}^{k+1} = \widetilde{A}^{k:0}\widetilde{\boldsymbol{x}}^0 + \sum_{l=0}^k \widetilde{A}^{k:l}\delta^l.$$
(12)

where  $\widetilde{A}^{k:0} \triangleq \widetilde{A}^k \widetilde{A}^{k-1} \dots \widetilde{A}^0$ . Based on Proposition 18 in [16], we have

$$\|\widetilde{\boldsymbol{x}}^{k+1} - \mathbf{1}x_{\psi}^{k+1}\| \le C_2 \rho^k \|\widetilde{\boldsymbol{x}}^0 - \mathbf{1}x_{\psi}^0\| + C_2 \sum_{l=0}^k \rho^{k-l} |\delta^l|,$$
(13)

where  $x_{\psi}^{k+1}$  satisfies  $x_{\psi}^{k+1} = x_{\psi}^{k} + \psi_{i}^{k} \delta^{k}$  for a sequence of stochastic vectors  $\{\psi^{k}\}$  and  $\rho \in (0, 1), C_{2} > 0$  are given in [16, Lemma 17].

## B. Step II

In this subsection, we focus on consensus agreement, newton direction tracking error and optimization error for our framework, the consensus error and optimization error at iteration k are defined as

$$E_c^k \triangleq \|\widetilde{\boldsymbol{x}}^k - \boldsymbol{1} x_\psi^k|, \quad E_o^k \triangleq |x_\psi^k - x^*|.$$

Similar to the definition of Gradient Tracking Error in [16], we give the gradient tracking error and newton direction tracking error at iteration k along with the magnitude of the tracking variables are defined as

$$E_{\mathbf{t}}^{k} \triangleq \left| y_{i}^{k} - \xi_{i}^{k-1} \bar{g}^{k} \right|, E_{\mathbf{y}}^{k} \triangleq \left| y_{i}^{k} \right|, \bar{g}^{k} \triangleq \frac{1}{n} \sum_{i=1}^{n} \nabla f_{i} \left( x_{i}^{k} \right)$$
$$E_{d}^{k} \triangleq \left| d_{i}^{k} \right|, E_{dt}^{k} \triangleq \left| d_{i}^{k} - \xi_{i}^{k-1} \bar{d}^{k} \right|,$$

where  $\bar{d}^k = \frac{1}{n} \sum_{i=1}^n (\nabla^2 f_i(x_i))^{-1} \nabla f_i(x_i)$ ,  $\{\xi_k\}$  is the sequence of stochastic vector introduced in [16, Lemma 15]. Next, we build the connection between these error quantities.

Proposition 3.4: The error quantities satisfy: for all  $k \in \mathcal{N}_0$ , there exists some positive constants  $C_0, C_2$ , such that

$$E_{c}^{k+1} \le C_{2}\rho E_{c}^{0} + C_{2}\sum_{l=0}^{k}\rho^{k-l}\alpha^{l}E_{d}^{l}$$
(14)

$$E_t^{k+1} \le C_0 \rho^k \|\mathbf{g}^0\| + 3C_0 L_1 \sum_{l=0}^k \rho^{k-l} (E_c^l + \alpha^l E_d^l) \quad (15)$$

$$E_{d}^{k} \leq \frac{1}{M_{1}} E_{t}^{k} + \frac{L_{1}}{\mu \sqrt{n}} E_{c}^{k} + E_{o}^{k}$$
(16)

$$E_{dt}^k \le \frac{1}{M_1} E_t^k \tag{17}$$

$$E_{\mathbf{o}}^{k+1} \leq \sum_{l=0}^{k} \left( \prod_{t=l+1}^{k} \left( 1 - \eta^{2} \alpha + \eta^{2} \alpha \delta^{t} \right) \right) \left( \frac{L_{1}}{\mu \sqrt{n}} E_{\mathbf{c}}^{l} + E_{\mathbf{dt}}^{l} \right) \alpha^{l} + \prod_{t=0}^{k} \left( 1 - \eta^{2} \alpha + \eta^{2} \alpha \delta^{t} \right) E_{\mathbf{o}}^{0}$$

$$(18)$$

#### C. Step III

Based on [16, Theorem 23] and Proposition 3.4, we can build the linear dynamics system as (21), where  $|E_{\mathbf{d}}|^{\lambda,N}$ ,  $|E_{\mathbf{c}}|^{\lambda,N}$ ,  $|E_{\mathbf{t}}|^{\lambda,N}$ ,  $|E_{\mathbf{o}}|^{\lambda,N}$  are defined based on [16, Lemma 21]. By [16, Theorem 23], if  $\rho(K) < 1$ , the algorithm vanish with linear convergence rate. According to Lemma 24 in [16],  $\rho(K) < 1$  if and only if  $p_K(1) > 0$ . The characteristic polynomial of **K** is

$$x^{4} - \frac{b_{1}C_{2}x^{2}\alpha + b_{1}C_{2}x\alpha^{2}(\lambda - \mathcal{L}(\alpha))}{(\lambda - \rho)(\lambda - \mathcal{L}(\alpha))} - \frac{b_{2}C_{2}x\alpha + b_{2}x^{2}\alpha(\lambda - \rho)}{M_{1}(\lambda - \rho)^{2}} - \frac{b_{2}C_{2}\alpha^{2}}{M_{1}(\lambda - \rho)^{2}(\lambda - \mathcal{L}(\alpha))} - \frac{b_{2}x\alpha^{2}}{M_{1}(\lambda - \rho)(\lambda - \mathcal{L}(\alpha))}$$
(19)

Consider the the continuity of equation (19) and  $p_K(1) > 0$ , let x = 1,  $\lambda = 1$ , we obtain  $\bar{\alpha}_1 = \frac{1}{J_1}$ , where

$$\begin{split} J_1 &= \frac{b_1 C_2 M_1 (1-\rho) + b_2 C_2 + b_2 (1-\rho)}{M_1 (1-\rho)^2} + \frac{b_1 C_2}{(1-\rho)(\eta^2 - \eta^2 \delta)} \\ &+ \frac{b_2 C_2}{M_1 (1-\rho)^2 (\eta^2 - \eta^2 \delta)} + \frac{b_2}{M_1 (1-\rho)(\eta^2 - \eta^2 \delta)} \end{split}$$

For term  $\frac{b_1C_2}{(1-\rho)\eta^2(1-\delta)} \geq b_1C_2 \geq \frac{L_1}{\mu\sqrt{n}}2\sqrt{(D+2)n} = \frac{L_1}{\mu}2\sqrt{D+2} > 1$ , it's easy to check other terms are all positive, so we have  $\overline{\alpha}_1 < 1$ . Therefore, when  $0 < \alpha < \overline{\alpha}_1$ , we have

$$E_c^k = \mathcal{O}(\lambda^k), E_d^k = \mathcal{O}(\lambda^k), E_t^k = \mathcal{O}(\lambda^k), E_o^k = \mathcal{O}(\lambda^k)$$

with  $\lambda \in (0,1)$  given by

$$\lambda = \max(\rho + \sqrt{(b_1 C_2 + \frac{b_2 C_2}{M_1} + \frac{b_2}{M_1})(1 + \frac{1}{\epsilon})\alpha}, 1 - \alpha(\eta^2 - \eta^2 \delta - \epsilon)),$$
(20)

where  $\delta > 0$  is arbitrary small,  $\epsilon > 0$  is properly chosen.

# **IV. NUMERICAL EXPERIMENTS**

In this section, we numerically test our theoretical findings and compare with existing asynchronous methods and synchronous newton tracking method in the decentralized case over two classes of problems. The network is generated randomly, firstly generate a cycle graph which guarantee strong connected of the graph, then given specific number of out-neighbors  $\mathcal{N}_{\text{out}} \leq n$ , we add edges randomly. The optimal point of the problem  $x^*$  will be precomputed by centralized newton method.

The asynchronous model we adapt to test iteration performance is common in asynchronous optimization, see [16], [27]. In fully asynchronous setting, a virtual global clock that different from [20] which need real global clock for coordination is usually assumed to record the whole iteration of the network, no matter which node complete a update, k would increase one. The agents follow the generated activation list and transmit information to their out-neighbors immediately after complete their own local update. The information from node i will not be available to j after the virtual global iteration  $k + T_{ij}^k$ . Transmitted delay  $T_{ij}^k$ is modeled by (integer) traveling time which is sampled

$$\begin{bmatrix} |E_{\mathbf{d}}|^{\lambda,N} \\ |E_{\mathbf{c}}|^{\lambda,N} \\ |E_{\mathbf{t}}|^{\lambda,N} \\ |E_{\mathbf{c}}|^{\lambda,N} \end{bmatrix} \preccurlyeq \underbrace{\begin{bmatrix} 0 & b_{1} & \frac{1}{M_{1}} & 1 \\ \frac{C_{2\alpha}}{\lambda-\rho} & 0 & 0 & 0 \\ \frac{b_{2\alpha}}{\lambda-\rho} & \frac{b_{2}}{\lambda-\rho} & 0 & 0 \\ 0 & \frac{b_{1\alpha}}{\lambda-\mathcal{L}(\alpha)} & \frac{\alpha}{M_{1}(\lambda-\mathcal{L}(\alpha))} & 0 \end{bmatrix}}_{\triangleq \mathbf{K}} \begin{bmatrix} |E_{\mathbf{d}}|^{\lambda,N} \\ |E_{\mathbf{c}}|^{\lambda,N} \\ |E_{\mathbf{c}}|^{\lambda,N} \\ |E_{\mathbf{c}}|^{\lambda,N} \end{bmatrix} + \mathbf{c}, \qquad (21)$$

where  $b_1 \triangleq \frac{L_1}{\mu\sqrt{n}}, b_2 \triangleq 3C_0L_1, \mathcal{L}(\alpha) = 1 - \eta^2 \alpha + \eta^2 \alpha \delta.$ 



Fig. 3. Iteration Performance for Least Square



#### A. Least Square

Least Square is an typical instance of (1), where  $f_i = ||A_ix_i - b_i||^2$ ,  $A_i$  is the feature matrix and  $b_i$  is the label vector privately known by the agent *i*. The whole data *A* and *b* are randomly generated from standard Guassian distribution and normalize it by its spectral norm, then we partition it into n parts privately stored in local agents.

1) Comparison with Asynchronous First Order Methods: We compare the asynchronous newton tracking mechanism Algorithm 2 and Algorithm 3 with all existing asynchronous methods for iteration performance. Figure 3(a) and 3(b) show the iteration performance for 30 agents with 8 out\_neighbors in the asynchronous model, where *MaxTravelTime* is 10 and 40 respectively. We manually adjust the step-size in order to get best performance of each algorithm. From figure we observe that methods belong to the same kind mechanism share the similar feature, and our mechanism efficiently accelerate the asynchronous first-order algorithms. Additionally, in more asynchronous scenarios, we observe that these first order algorithms especially APPG suffers a lot if the network become more asynchronous, while our mechanism are more robust.

2) Comparison with Synchronous Newton Tracking: We compare the robustness of our mechanism with synchronous newton tracking from two perspectives, link failure and transmission delay. Synchronous newton tracking adapts synchronous peer to peer communication in mpi4py, while



Fig. 4. Wall-Clock Time Performance for Least Square

asynchronous method uses asynchronous peer to peer communication which needs  $Attach\_buffer()$  initially. Fig 4(a) shows the running time performance for 10 agents with different link failure probability  $0.01(1 + \frac{i}{4})$ . Fig 4(b) shows the running time performance for 10 agents with different delays delay<sub>i</sub> =  $0.05 * (1 + \frac{i}{5})$ , while each node keeps 100 observations, the dimension of features is 20. In addition, the iterations of asynchronous algorithms for agents range from 177 to 472 and the iterations of synchronous ones runs about 221 iterations. Results show our asynchronous newton tracking methods is more robust to the link failure, delay, and heterogeneous nodes, especially the slow nodes.

#### B. Logistic Regression

To further confirm the efficiency of the proposed mechanism, we solved the Two-Class logistic regression problem as follows.

$$x^* = \underset{x \in \mathbb{R}^d}{\operatorname{argmin}} \frac{\rho}{2} \|x\|^2 + \sum_{i=1}^n \sum_{j=1}^{m_i} \ln\left(1 + \exp\left(-\left(\mathbf{o}_{ij}^T x\right) \mathbf{p}_{ij}\right)\right),$$

where each node *i* privately owns  $m_i$  training samples  $(\mathbf{o}_{ij}, \mathbf{p}_{ij}) \in \mathbb{R}^d \times \{-1, +1\}, j = 1, \ldots, m$ . The elements of  $\mathbf{o}_{ij}$  are randomly generated following the standard Gaussian distribution and those of  $\mathbf{p}_{ij}$  are generated following the uniform distribution on  $\{-1, 1\}, \frac{\rho}{2} ||x||^2$  is a regularization term to avoid over-fitting. The network setting is the same as ridge regression, figure 5(a), and figure 5(b) collaborate the results in the least square. In asynchronous iteration performance over synthetic data, and for wall-clock time comparison with synchronous newton tracking, we tested it over real dataset, Covtype. Each agent keeps 1000 observations with different delays delay<sub>i</sub> =  $0.1 * (1 + \frac{i}{5})$ , the dimension of Covtype is 54.



Fig. 5. Numeric Experiments For Logistic Regression

#### V. CONCLUSIONS

In this paper, we have proposed a novel asynchronous newton tracking mechanism, which explicit second-order information under the fully asynchronous setting. Theoretical and empirical results demonstrate the efficiency and robustness of our mechanism compared with existing asynchronous methods. We leave several interesting directions for future work. In theoretical part, local condition is assumed, while it is possible to relax this condition and get a global convergence analysis. The other interesting direction is to introduce multi-consensus mechanism for our asynchronous newton tracking, there lies interesting balance between computation and communication for the fully asynchronous decentralized optimization.

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