Controlling Large-Scale Self-Organized Networks with Lightweight Cost for Fast Adaptation to Changing Environments

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Self-organization has potential for high scalability, adaptability, flexibility, and robustness, which are vital features for realizing future networks. Convergence of self-organizing control, however, is slow in some practical applications compared to control with conventional deterministic systems using global information. It is therefore important to facilitate convergence of self-organizing controls. In controlled self-organization, which introduces an external controller into self-organizing systems, the network is controlled to guide systems to a desired state. Although existing controlled self-organization schemes could achieve this feature, convergence speed for reaching an optimal or semioptimal solution is still a challenging task. We perform potential-based self-organizing routing and propose an optimal feedback method using a reduced-order model for faster convergence at low cost. Simulation results show that the proposed mechanism improves the convergence speed of potential-field construction (i.e., route construction) by at most 22.6 times with low computational and communication cost.

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Additional Key Words and Phrases: Controlled self-organization, robust control, potential-based routing, fast convergence

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1. INTRODUCTION

Self-organization, where components behave individually and autonomously, is a natural phenomenon in natural distributed systems [Pintea 2014; Yang et al. 2013]. In a self-organizing system, each component follows simple rules using locally available information. Through direct or indirect interactions among components, a global behavior or pattern emerges on a macroscopic level without a central control entity. In a self-organizing system, up-to-date information regarding the entire system or many other components is unnecessary, which considerably reduces computational cost

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and communication overhead for collecting global information. This localized control leads to a capability of handling local failures and small environmental changes by interaction of local components. Thus, self-organizing systems are expected to automatically recover from failures and adapt to environmental changes, without involving centralized control. These are the reasons a variety of self-organization-based models have been applied to information networking, such as routing, synchronization, and task assignment [Zhang et al. 2013; Zheng and Sicker 2013]. In future large-scale, complex networks, we can expect features such as scalability, adaptability, and robustness to be improved to an extent not possible by conventional network control methods [Balasubramaniam et al. 2011].

Although self-organization control without global knowledge of the current network state has various benefits, such control has critical disadvantages that complicate practical use in industrial and business systems [Dressler 2008]. It may take a long time for global patterns to emerge in large-scale systems, because they appear as a consequence of interactions between autonomous components. This property also leads to slow adaptation to large environmental changes, which is difficult to solve solely by local interaction in self-organizing systems. In addition, self-organizing systems that use only local information sometimes fall into local optima. On the other hand, in conventional centralized systems, global information can reach an optimal solution, although the required computational cost to do so often leads to unrealistic solutions.

Such disadvantages raised from real applications brought about the idea of controlled (guided, managed) self-organization, where the self-organizing system is controlled through some constraints Branke et al. 2006; Schmeck et al. 2010; Prokopenko 2014]. For example, Arakawa et al. and Kominami et al. use the concept of controlled self-organization, where an external observer/controller guides self-organizing optical network [Arakawa et al. 2011] and sensor network [Kominami et al. 2013] systems through a feedback mechanism that leads them to a desired state. The self-organizing system can then be controlled through fully observed information. However, enhancing the convergence speed to reach an optimal or semioptimal solution remains as an outstanding task.

We previously introduced an external controller with an optimal feedback mechanism to self-organizing systems in Kuze et al. [2014]. The external controller collects information regarding the network, such as node states and network topology, via a partial set of nodes directly monitored by the controller, and estimates system dynamics using a mathematical model that describes the network dynamics, then determines optimal control inputs based on robust control theory [Zhou et al. 1995] for facilitating the convergence. Optimal feedback mechanisms for controlling dynamical systems have been researched for many years in the field of control theory [Kirk 2012]. A controller monitors a system and provides an optimal control feedback that minimizes the cost function [Kirk 2012] based on a mathematical model of the system. Note that there are various errors (e.g., modeling errors) and unexpected disturbances, and thus the optimal control feedback is not necessarily optimal at any time in real systems. Simulation results in Kuze et al. [2014] showed that the mechanism improves the convergence speed of self-organization; however, in large-scale networks especially, it is generally difficult for the controller to collect detailed network information and to estimate the network dynamics [Mulvey et al. 1995], because doing so requires considerable costs and, even worse, loses the advantage of scalability of self-organizing systems.

The contribution of this article is to converge self-organizing network systems with lightweight cost while retaining a high performance (e.g., high convergence and adaptability in our proposal) of robust control. To that end, we first regulate the area in which the external controller collects node states to reduce the cost for collecting information (Figure 1). If the controller collects all node states, the communication overhead is extremely large and traffic congestion would occur. Therefore, we limit the area from

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 $\label{eq:Fig.1.} Fig. \ 1. \ \ Advantages \ of optimal \ feedback \ with \ model \ reduction \ (\mathbb{N}^+ \ is \ the \ set \ of \ all \ nonzero \ natural \ numbers).$

which the controller can collect information to nodes that can be reached within several hops from the monitored nodes. The controller can estimate the states of all nodes from only a part of them given the network topology, because the state of each node is determined in accordance with local interactions based on prescribed rules [Zhou et al. 1995]. In other words, the controller acquires network information with lower communication overhead and then calculates optimal feedback inputs. In our scheme, a controller has an internal model estimating the actual network system. The order (N_{dim}) of that model is proportional to that of the actual network model, namely the number of nodes. Note that we need $O(N_{dim}^2)$ for controller state updates at each time instance, as well as $O(N_{dim}^{3})$ computation for controller redesign when the network topology changes. The computational cost is thus extremely large in large-scale networks when we use the original model, whose order is proportional to the number of nodes. We therefore reduce the order of the network model (consequently that of N_{dim}) to decrease the computational cost via a model reduction technique [Antoulas et al. 2006]. As shown in Figure 1, model reduction refers to reducing the number of state variables in dynamical systems while retaining suitable input/output characteristics. Note that there is a trade-off between the optimization of control input and the number of state variables (i.e., between the convergence speed and the computational cost). Intuitively, a reduced-order model leads to the loss of the original model while decreasing the computational cost. This trade-off relation requires us to further examine the number h of state variables of the reduced-order model.

The effectiveness of our proposal is evaluated through computer simulation studies where we consider potential-based routing—a self-organizing routing mechanism for

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wireless sensor networks—with optimal feedback, and we evaluate the convergence speed after environmental changes. Wireless sensor networks represent one suitable application of our proposal. This is because our proposal requires the expensive computation of control input for an external powerful controller and not for each component of a self-organizing system. Therefore, the performance of the whole system can be improved even though each component in the systems does not need much computational power and energy consumption. Optimality of our feedback mechanism is analytically guaranteed in synchronous systems but not in asynchronous systems. Then we first assume a wireless sensor network where nodes behave asynchronously and the controller can observe the network state via the monitored nodes experiencing communication delay. To evaluate the robustness of our proposal against information loss in the controller, we next simulate an environment where several nodes fail and the controller cannot immediately detect node failures. Through these evaluations, we will show that optimal feedback using a reduced-order model can enhance the convergence speed of self-organizing systems at fairly low cost, even if the network is large. Moreover, optimal feedback does not have a negative impact on the robustness, which is a notable property of self-organizing systems. Note that potential-based routing for wireless sensor networks is just an example of application for our proposal.

The remainder of this article is organized as follows. First, we briefly explain potential-based routing in Section 2. We propose and explain potential-based routing with optimal feedback in Section 3 and then explain a reduced-order model with which the controller estimates the network dynamics in Section 4. We then show fast adaptation to an environmental change of the proposed method through simulation and give a discussion of our proposal in Section 5. Finally, in Section 6, we present our conclusions and suggest areas for future work.

2. POTENTIAL-BASED ROUTING

Potential-based routing is a self-organizing routing mechanism that is active in the fields of wireless sensor networks, mobile ad hoc networks, and information-centric networks [Kominami et al. 2013; Basu et al. 2003; Jung et al. 2009; Wu et al. 2008; Sheikhattar and Kalantari 2013; Eum et al. 2014; Lee et al. 2014]. Here we assume that potential-based routing is used in wireless sensor networks. In potential-based routing, each node has a scalar value called its *potential*, and data packets are forwarded to a neighbor whose potential is smaller than that of the forwarder. In wireless sensor networks, data packets are generally sent to a sink node, and the fewer the hops from the sink node a node is, the lower the potential value assigned to the node. The simple forwarding rule "forward data to a neighboring node with lower potential" can therefore result in data packet collection toward sink nodes, as illustrated in Figure 2. Potential-based routing has high scalability because each node uses only local information for calculating potentials and a local rule for forwarding data. Furthermore, it can achieve load balancing and consequently network lifetime improvement by calculating potentials using information such as flow rates, queue lengths, or remaining energy [Wu et al. 2008]. In Sections 2.1 and 2.2, we describe a potential field construction method and show how to select a next hop node using the gradient of the field.

2.1. Potential Field Construction

Sheikhattar and Kalantari [2013] focused on the convergence of potential-based routing and achieved enhancement of the potential convergence speed. They proposed a potential calculation method based not only on current potentials but also on last potentials to accelerate potential convergence. Node *n*'s potential at time t, $\theta_n(t)$, is



Fig. 2. Potential-based routing.

calculated by Equation (1):

$$\begin{aligned} \theta_n(t+1) &= (\alpha+1)\theta_n(t) - \alpha\theta_n(t-1) \\ &+ \beta\sigma_n\left(\sum_{k \in \mathcal{N}_b(n)} \{\theta_k(t) - \theta_n(t)\} + f_n(t)\right), \end{aligned} \tag{1}$$

where $\mathcal{N}_b(n)$ is the set of neighbors of node n and α is a parameter that determines the weights of the current and the last potential values when calculating the next potential. Larger α means that the weight of the last potential value is larger, and therefore the system becomes less subject to current noise, although the convergence speed is slower. Parameter β determines the influence amount of neighbor node potentials. σ_n is defined as $\sigma_0/|\mathcal{N}_b(n)|$ (σ_0 is a parameter), and $f_n(t)$ corresponds to the flow rate of node n at time t. If $f_n(t)$ is a positive value, it means the data generation rate of node n, whereas if $f_n(t)$ is negative, it means the rate of data packets delivered to node n. For sink node n, $f_n(t)$ corresponds to targeted flow rates that are given by the network manager. If the flow conservation constraint is upheld—that is, $\sum_{n \in \{1, ..., N\}} f_n(t) = 0$ —a potential field is constructed so that actual rates of data packets delivered to nodes satisfy given flow rates (i.e., all gradients, which are potential differences between next hop nodes).

The convergence speed based on Equation (1) is faster than simple Jacobi iterations (e.g., our previous work [Kominami et al. 2013]) but still takes a long time to converge due to its calculation being based only on local information (see Section 5 for an example convergence). In both cases, potentials of nodes converge as a result of the iterative behavior (i.e., potential updates of nodes), and therefore the convergence speed is faster with a shorter interval of potential updates. Instead of relying on only local interactions among nodes, we introduce into potential-based routing an external controller that observes network states (potential values), estimates its future state, and regulates potentials of a partial set of nodes for faster convergence.

2.2. Routing

If a node has a data packet, it forwards the data packet according to the potential values of itself and its neighbors. In our potential-based routing, when a sensor node generates or receives a data packet, it probabilistically selects a next node that is assigned a lower potential value than itself, and the packet eventually arrives at a sink node. Specifically, a next-hop node is selected proportionally with potential values—that is, the probability $p_{i \to j}(t)$ that sensor node *i* selects a neighbor node *j* as the

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Fig. 3. Potential-based routing with a controller's feedback, where an external controller collects potential values from observable nodes and introduces control inputs to controllable nodes periodically.

next-hop node for a data packet at time *t* is given by

$$p_{i \to j}(t) = \begin{cases} \frac{\theta_i(t) - \theta_j(t)}{\sum_{k \in \mathcal{N}_l(i)} \{\theta_i(t) - \theta_k(t)\}}, & \text{if } j \in \mathcal{N}_l(i) \\ 0, & \text{otherwise} \end{cases}$$

where $\mathcal{N}_l(i)$ is the set of node *i*'s neighbor nodes that are assigned lower potential values than node *i*. If node *i* has no neighbor node with lower potential (i.e., $|\mathcal{N}_l(i)| = 0$), the data packet is not sent to any node and is dropped, but such cases generally happen only in transient cases, such as node failures or changes of potential values at the sink node.

3. POTENTIAL-BASED ROUTING WITH OPTIMAL FEEDBACK

3.1. Overview

In this section, we describe a model of network dynamics and explain our optimal control scheme. A controller monitors network information, in particular the potential values of a partial set of nodes, which we call *observable nodes*. The controller then returns suitable control inputs to a partial set of nodes, which we call *controllable nodes*, for accelerating convergence of the potential distribution toward the target potential distribution. In this article, target potentials are estimates of converged potential values derived from current information, specified in Section 3.2. We assume that the controller and sink nodes are power supplied so that these sink nodes can have direct reliable connections to the controller at all times. Therefore, in our proposal, the controller monitors network information and provides control inputs via them, as illustrated in Figure 3.

In our proposal, the area over which the controller monitors potential values is limited to several hops from sink nodes for reducing communication overhead. Of course, the controller cannot directly get node potentials outside the area, but it can estimate them by utilizing the potential dynamics model, which describes potential changes based on local node interactions. When receiving control inputs from the controller, sink nodes diffuse the information, changing potential amounts over the entire network through local interactions of sensor nodes (Equation (2) in Section 3.2). This estimation of the controller requires high computational cost, and the reduced-order model can reduce this computational cost for estimating potential values of nonobservable nodes, as will be explained in Section 4. Note that information of the network topology is needed for designing a controller, as is flow rates of nodes for calculating

target potential values. Such information is difficult to estimate but is reported to the controller only when it changes, as we assume that changing intervals of the network topology and flow rates are lower than the convergence time of potentials. This assumption is feasible, because the potential convergence is generally achieved as a result of the iterative behavior (nodes' potential updates and the controller's feedback) in potential-based routing with an optimal feedback, so frequencies of potential updates and controls need to be much higher than those of changes in the network topology and flow rates.

3.2. Network Dynamics

Let the dynamics of potentials be given by a deterministic discrete-time model. In our proposal, we represent $f_n(t) = \bar{f}_n + d_n(t)$, where \bar{f}_n is the stationary flow rate and d_n is nonstationary fluctuation (white Gaussian noise) at node *n*. Then, as in Sheikhattar and Kalantari [2013], each potential is updated by

$$\theta_n(t+1) = (\alpha+1)\theta_n(t) - \alpha\theta_n(t-1) + \beta\sigma_n\left(\sum_{k\in\mathcal{N}_b(n)} \{\theta_k(t) - \theta_n(t)\} + \bar{f}_n + d_n(t)\right) + \eta_n(t),$$
(2)

where η_n represents feedback input received from the controller. If node *n* does not receive any feedback input directly from the controller, then $\eta_n(t) = 0$. In our proposal, the controller collects and estimates the node potentials and then provides to the network feedback inputs $\mathbf{u}(t) = [\eta_1(t) \ \eta_2(t) \ \cdots \ \eta_{N_{ctrl}}(t)]^T$ (where N_{ctrl} denotes the number of nodes that receive feedback from the controller), as described later. Because of feedback inputs, we consider that node potentials can converge faster than in the non-control scheme (Equation (1)) where each node updates its potential based only on local interactions with neighbors. In Sheikhattar and Kalantari [2013], σ_n is set to $\sigma_0/|\mathcal{N}_b(n)|$, but this value may lead to oscillation of potentials in some situations since Equation (9) (which will be shown later) has no solution. In this article, we therefore set σ_n to the constant value σ for all n ($n \in \{1, 2, \ldots, N\}$). Now we define $\theta_i(t)$ as $[\theta_i(t) \ \theta_i(t+1)]$, and Equation (2) is rewritten by

$$\boldsymbol{\theta}_{n}(t+1) = \boldsymbol{A}_{1}\boldsymbol{\theta}_{n}(t) + \boldsymbol{A}_{0}\left(\sum_{k\in\mathcal{N}_{b}(n)} \{\boldsymbol{\theta}_{k}(t) - \boldsymbol{\theta}_{n}(t)\} + \boldsymbol{M}\left(\bar{f}_{n} + d_{n}(t)\right)\right) + \boldsymbol{M}\eta_{n}(t), \quad (3)$$

where

$$\boldsymbol{A}_{1} = \begin{bmatrix} 0 & 1 \\ -\alpha & \alpha + 1 \end{bmatrix}, \, \boldsymbol{A}_{0} = \begin{bmatrix} 0 & 0 \\ 0 & \beta \sigma \end{bmatrix}, \, \boldsymbol{M} = \begin{bmatrix} 0 \\ 1 \end{bmatrix}.$$

In our proposal, the controller estimates potentials of unobservable nodes and future potentials and then determines control inputs with H^{∞} optimization (described later), allowing us to describe the system dynamics in the form of the state space model. We first formulate the potential dynamics of all nodes. For example, we consider the case of a network of four nodes, as shown in Figure 4. From Equation (3), the potential dynamics of node 1 is given by

$$\boldsymbol{\theta}_{1}(t+1) = \boldsymbol{A}_{1}\boldsymbol{\theta}_{1}(t) + \boldsymbol{A}_{0}\left(\sum_{k \in \{2,3\}} \{\boldsymbol{\theta}_{k}(t) - \boldsymbol{\theta}_{1}(t)\} + \boldsymbol{M}\left(\bar{f}_{1} + d_{1}(t)\right)\right).$$
(4)



Fig. 4. Example of a network with three sensor nodes and a sink node. f_1 , f_2 , f_3 , and f_4 are stationary flow rates for each node. η_3 is the control input to node 3 from the controller (the other nodes are not directly controlled by the controller, and η_1 , η_2 , and η_4 is 0 at all times). d_1 , d_2 , d_3 and d_4 are nonstationary fluctuations at each node, although they are not described in this figure.

Similarly to Equation (4), the potential dynamics of nodes 2, 3, and 4 are given, respectively, by

$$\boldsymbol{\theta}_{2}(t+1) = \boldsymbol{A}_{1}\boldsymbol{\theta}_{2}(t) + \boldsymbol{A}_{0}\left(\sum_{k \in \{1,3\}} \{\boldsymbol{\theta}_{k}(t) - \boldsymbol{\theta}_{2}(t)\} + \boldsymbol{M}\left(\bar{f}_{2} + d_{2}(t)\right)\right),$$
(5)

$$\boldsymbol{\theta}_{3}(t+1) = \boldsymbol{A}_{1}\boldsymbol{\theta}_{3}(t) + \boldsymbol{A}_{0}\left(\sum_{k \in \{1,2,4\}} \{\boldsymbol{\theta}_{k}(t) - \boldsymbol{\theta}_{3}(t)\} + \boldsymbol{M}\left(\bar{f}_{3} + d_{3}(t)\right)\right) + \boldsymbol{M}\eta_{3}(t), \quad (6)$$

$$\boldsymbol{\theta}_{4}(t+1) = \boldsymbol{A}_{1}\boldsymbol{\theta}_{4}(t) + \boldsymbol{A}_{0}\left(\sum_{k \in \{3\}} \{\boldsymbol{\theta}_{k}(t) - \boldsymbol{\theta}_{4}(t)\} + \boldsymbol{M}\left(\bar{f}_{4} + d_{4}(t)\right)\right).$$
(7)

From Equations (4) through (7), the potential dynamics of the network shown in Figure 4 is given by

$$\begin{bmatrix} \theta_{1}(t+1)\\ \theta_{2}(t+1)\\ \theta_{3}(t+1)\\ \theta_{4}(t+1) \end{bmatrix} = \begin{pmatrix} I_{4\times4} \otimes A_{1} - \begin{bmatrix} 2 & -1 & -1 & 0\\ -1 & 2 & -1 & 0\\ -1 & -1 & 3 & -1\\ 0 & 0 & -1 & 1 \end{bmatrix} \otimes A_{0} \end{pmatrix} \begin{bmatrix} \theta_{1}(t)\\ \theta_{2}(t)\\ \theta_{3}(t)\\ \theta_{4}(t) \end{bmatrix} + \begin{pmatrix} \beta\sigma \left(\begin{bmatrix} f_{1}\\ f_{2}\\ f_{3}\\ f_{4} \end{bmatrix} + \begin{bmatrix} d_{1}\\ d_{2}\\ d_{3}\\ d_{4} \end{bmatrix} \right) + \begin{bmatrix} 0\\ 0\\ \eta_{3}\\ 0 \end{bmatrix} \end{pmatrix} \otimes M,$$
(8)

where $I_{N \times N}$ is the identity matrix of $N \times N$. The matrix $\begin{bmatrix} 2 & -1 & -1 & 0 \\ -1 & 2 & -1 & 0 \\ -1 & -1 & 3 & -1 \\ 0 & 0 & -1 & 1 \end{bmatrix}$ corresponds

to the graph Laplacian, which indicates the network topology. We refer to the graph Laplacian as Γ in this article. The element l_{ij} of graph Laplacian Γ is given by

$$l_{ij} = \left\{ egin{array}{ll} \deg(i), & ext{if } i = j \ -1, & ext{if } i
eq j \land ext{node } i ext{ is adjacent to node } j \ 0, & ext{otherwise} \end{array}
ight.$$

where deg(*i*) corresponds to the degree of node *i*. The operation \otimes that is called the *Kronecker product* is an operation on two matrices. Given an $m \times n$ matrix J and a

 $p \times q$ matrix **K**, the Kronecker product $\boldsymbol{J} \otimes \boldsymbol{K}$ is given by

$$\boldsymbol{J}\otimes\boldsymbol{K}=\left[\begin{array}{ccc} j_{11}\boldsymbol{K}&\cdots&j_{1n}\boldsymbol{K}\\ \vdots&&\vdots\\ j_{m1}\boldsymbol{K}&\cdots&j_{mn}\boldsymbol{K}\end{array}\right].$$

Then we define a vector $\Theta(t)$ that shows potential values for all nodes as

$$\boldsymbol{\Theta}(t) := \begin{bmatrix} \boldsymbol{\theta}_1(t) \ \boldsymbol{\theta}_2(t) \ \cdots \ \boldsymbol{\theta}_N(t) \end{bmatrix}^T.$$

As in Equation (8), the potential dynamics of all nodes is described with $\Theta(t)$ as

$$\Theta(t+1) = \boldsymbol{A}\Theta(t) + (\beta\sigma(\boldsymbol{F} + \boldsymbol{D}(t)) + \boldsymbol{E}\boldsymbol{u}(t)) \otimes \boldsymbol{M},$$

where

$$\boldsymbol{A} = \boldsymbol{I}_{N \times N} \otimes \boldsymbol{A}_1 - \boldsymbol{\Gamma} \otimes \boldsymbol{A}_0.$$

The flow rate vector \mathbf{F} is defined as $[\bar{f}_1 \cdots \bar{f}_N]^T$, and the fluctuation rate vector $\mathbf{D}(t)$ is defined as $[d_1(t) \cdots d_N(t)]^T$. Note that the $(N \times N_{ctrl})$ -matrix \mathbf{E} specifies the controllable node—that is, the element $e_{ij} \in \{0, 1\}$ of \mathbf{E} is 1 if and only if node *i* receives the *j*-th element of $\mathbf{u}(t)$ as control input $\eta_i(t)$ (in the case with the network shown in Figure 4, \mathbf{E} is set to $[0 \ 0 \ 1 \ 0]^T$). With the larger number of controllable nodes (i.e., with denser \mathbf{E}), the influence of the optimal control is larger, so the convergence speed is faster. On the contrary, significant properties originating from self-organization such as scalability and adaptability could be lost with too many controllable nodes.

Under these dynamics, the target potential distribution is given by a solution of

$$(\boldsymbol{I}_{2N\times 2N} - \boldsymbol{A})\boldsymbol{\bar{\Theta}} = \beta \sigma \boldsymbol{F} \otimes \boldsymbol{M}.$$
(9)

Next we define X(t) as the regulation error $\overline{\Theta} - \Theta(t)$, which shows the differences between the convergence values and the current values of potentials for each node. The potential dynamics can be rewritten using X(t) in the form of the state space representation by Equation (10), with which the controller can estimate potentials of unobservable nodes and future potentials:

$$X(t+1) = AX(t) + B_1 d(t) + B_2 u(t).$$
(10)

Here the N_{ctrl} -dim vector \boldsymbol{u} (respectively, N-dim vector \boldsymbol{d}) concatenates $\eta_n(t)$ for controllable nodes (respectively, d_n). \boldsymbol{B}_1 and \boldsymbol{B}_2 characterize the effect of noises and control inputs, respectively, given by

$$oldsymbol{B}_1 = oldsymbol{B}_1^\prime \otimes \left[egin{array}{c} 0 \ eta\sigma \end{array}
ight], oldsymbol{B}_2 = oldsymbol{E} \otimes \left[egin{array}{c} 0 & 0 \ 0 & 1 \end{array}
ight].$$

 B'_1 characterizes the variance/co-variance of the noise distribution injected at each node. For simplicity, we choose $B'_1 = I_{N \times N}$ in this work, which implies that each node has independent noise with unit variance. Note that the dynamics (Equation (10)) depend on the graph Laplacian Γ but not on the flow rate \overline{f} .

3.3. Optimal Controller Design

Next we explain the controller dynamics. We consider the case in which the controller can observe

$$\begin{aligned} \mathbf{Y}(t) &= (\mathbf{H}^T \otimes \mathbf{I}_{2 \times 2}) \mathbf{X}(t) \\ &= (\mathbf{H}^T \otimes \mathbf{I}_{2 \times 2}) (\bar{\mathbf{\Theta}} - \mathbf{\Theta}(t)), \end{aligned}$$
(11)

where $\mathbf{Y}(t)$ is a $2N_{obs}$ -dim vector, with N_{obs} being the number of observable nodes, and an $(N \times N_{obs})$ matrix \mathbf{H} determines observable nodes in the same manner as \mathbf{E}

(if the controller can only observe node 3 in the network shown in Figure 4, \boldsymbol{H} is set to $[0 \ 0 \ 1 \ 0]^T$). Note that the controller can estimate the potential values of all nodes based on the information of only a partial set of nodes, and therefore not all elements of \boldsymbol{H} need to be set to 1. If more nodes are monitored by the controller, it can estimate potential values of all nodes more precisely while the communication overhead for collecting information of observable nodes becomes much larger. The future node potentials can be estimated by Equation (10) using observable information expressed with Equation (11). Considering the future potentials, the controller calculates a $\boldsymbol{u}(t)$, which accelerates the convergence speed of potentials.

Then the control input is calculated according to

$$\dot{\mathbf{X}}(t+1) = \mathbf{A}_c \, \mathbf{X}(t) + \mathbf{B}_c \, \mathbf{Y}(t), \tag{12}$$

$$\boldsymbol{u}(t) = \boldsymbol{C}_c \boldsymbol{X}(t) + \boldsymbol{D}_c \boldsymbol{Y}(t), \tag{13}$$

where 2N-dim vector \mathbf{X} is an internal model for the controller and the controller estimates potentials of all nodes with the model. A_c , B_c , C_c , and D_c are design parameters.

Concerning the performance criteria, let us define

$$\phi(k) = \boldsymbol{X}(k)^T \boldsymbol{X}(k) + r \boldsymbol{u}(k)^T \boldsymbol{u}(k)$$

as the stage cost, where r specifies the trade-off between convergence speed and input energy. With a larger r, control inputs become smaller and the stability of the system is enhanced. In other words, potentials change more gently, whereas the convergence speed of potentials becomes slower. Our design objective is then to minimize the worstcase error

$$\sup_{\boldsymbol{d}} \frac{\sum_{k=0}^{\infty} \phi(k)}{\sum_{k=0}^{\infty} \boldsymbol{d}(k)^T \boldsymbol{d}(k)}$$

This min-max type problem is called H^{∞} optimization [Zhou et al. 1995]. Given A, B_1 , B_2 , and H, it is known that the optimal A_c , B_c , C_c , D_c can be obtained based on semidefinite programming (see also Section 5). In general, the degree of the optimal controller (i.e., the size of \tilde{X}) is the same as that of the system to be controlled, which is 2N in this case. Note that we need to solve this optimization problem only when the connection topology changes. In other words, we only have to update \tilde{X} in Equation (12) and compute Equation (13) so long as the connection topology does not change.

In Kuze et al. [2014], we previously proposed potential-based routing with an optimal control that is explained in this section, but we only evaluate the situation where the controller observes all nodes (i.e., $N_{obs} = N$ and $\mathbf{H} = \mathbf{I}_{N \times N}$). The upper bound of the order of the communication cost for observation is $O(NP_{avg})$, where P_{avg} corresponds to the average path length of the network. Note that the average path length of the network generally depends on the number N of nodes. For example, the average path length of the Erdös-Rényi network (a random network) [Erdős and Rényi 1961] and the Barabási-Albert model (a scale-free network) [Albert and Barabási 2002] is $O(\log N)$, and that of the Watts-Strogatz network (a small-world network) [Watts and Strogatz 1998] is O(N) if $N < N^*$ and $O(\log N)$ if $N > N^*$.¹ Therefore, in this article, we show that the controller can estimate potentials of nonobservable nodes based on only a part of nodes' potentials with simulation evaluations in Section 5. Note that if the average path length is quite small against the network size, as in the case with WWW, the external controller can observe a large number of nodes with small communication

 $^{^{1}}N^{*}$ is a length that depends on the randomness p of a small-world network. See Albert and Barabási [2002] for details.

cost, which allows the external controller to estimate the potential dynamics more correctly.

Moreover, as the number of nodes increases (i.e., the size of vector $\mathbf{X}(t)$ becomes large), it becomes unfeasible to control the system using the model expressed in Equations (10) and (11), owing to the considerable computational cost. The design parameters (\mathbf{A}_c , \mathbf{B}_c , \mathbf{C}_c , and \mathbf{D}_c) are calculated with $O(N^3)$ because the computational cost is proportional to the cube of the size of \mathbf{X} [Gahinet and Apkarian 1994]. Control inputs ($\mathbf{u}(t)$) are calculated with $O(N^2)$ because they are calculated by Equations (12) and (13). We therefore approximate the model with a reduced-order model for reducing the computational cost.

4. MODEL REDUCTION

In our proposal, the controller uses a reduced-order model for estimating network dynamics to reduce the computational cost. In this section, we explain the model reduction of the network dynamics model described in Section 3.

In the reduced-order model, the dynamics of the system is expressed as an $(h \times 1)$ -vector $\mathbf{X}_r(t)$ whose elements are linear transformations of the original model $\mathbf{X}(t)$, and the reduced-order model is given by Equations (14) and (15):

$$\boldsymbol{X}_{r}(t+1) = \boldsymbol{A}_{r}\boldsymbol{X}_{r}(t) + \boldsymbol{B}_{r1}\boldsymbol{d}(t) + \boldsymbol{B}_{r2}\boldsymbol{u}_{r}(t), \qquad (14)$$

$$\boldsymbol{Y}_{r}(t) = \boldsymbol{C}_{r}\boldsymbol{X}_{r}(t) + \boldsymbol{D}_{r}\boldsymbol{d}(t).$$
(15)

Here, X_r is an *h*-dim vector with h (< 2N) being the order of the approximate model, and $u_r(t)$ corresponds to control inputs provided by the controller with the reduced-order model. We need to choose matrices A_r , B_{r1} , B_{r2} , C_r , and D_r of compatible dimensions such that $Y_r(t) \approx Y(t)$ for all input $u_r(t)$, $d_r(t)$. Many researchers have studied a variety of methods to approximate a model with a reduced-order model to control large, complex systems [Antoulas et al. 2006]. In our proposal, we approximate the original model based on a "balanced realization" that is highly compatible with the model expressed in the state space representation [Zhou et al. 1995; Antoulas et al. 2006]. In model reduction, a reduced-order model needs to have the same response characteristics as the original model for the accurate estimation of potentials. Here, response characteristics mean the effectiveness of inputs to the system.

The control input is then calculated in the same manner as $u_r(t)$ according to

$$\boldsymbol{X}_{r}(t+1) = \boldsymbol{A}_{cr}\boldsymbol{X}_{r}(t) + \boldsymbol{B}_{cr}\boldsymbol{Y}_{r}(t), \qquad (16)$$

$$\boldsymbol{u}_{r}(t) = \boldsymbol{C}_{cr} \tilde{\boldsymbol{X}}_{r}(t) + \boldsymbol{D}_{cr} \boldsymbol{Y}_{r}(t), \qquad (17)$$

where A_{cr} , B_{cr} , C_{cr} , and D_{cr} are design parameters. Now we can use Y(t) for $Y_r(t)$ in Equation (17) under the well-designed C_r , and D_r because of the relation of $Y_r(t) \approx Y(t)$. We similarly obtain the H^{∞} optimal controller of order h for the reduced system. These parameters determine how potential values observed by the controller affect feedback inputs.

A reduced-order model can describe the system dynamics with a constant number h of state variables, and the computational cost can be reduced using it. With a reducedorder model, the design parameters $(\mathbf{A}_{cr}, \mathbf{B}_{cr}, \mathbf{C}_{cr}, \text{ and } \mathbf{D}_{cr})$ are calculated with $O(h^3)$, and control inputs $(\mathbf{u}_r(t))$ are calculated with $O(h^2)$. In general, a model that has more state variables lets the controller estimate more correctly, but the computational cost is larger. In contrast, the computational cost is smaller, but the estimation error can be larger in a model that has fewer state variables. Therefore, h needs to be properly determined in accordance with the requirements or system properties (see also Section 5). In Section 5, we conduct a network simulation to reveal the packet-level dynamics and show that Equations (2) through (17), assuming synchrony of nodes, can be adopted to asynchronous network systems.

5. PERFORMANCE EVALUATION

In this section, we evaluate our proposal to show that the convergence speed is enhanced by introducing optimal feedback using the reduced-order model. We conduct a computer simulation and evaluate the convergence speed comparing our proposal (potential-based routing with optimal feedback (PBR-opt) and one using the model reduction (PBR-opt-mr)) with the non-control scheme proposed in Sheikhattar and Kalantari [2013]. First, in Section 5.2, we evaluate the potential convergence speed after traffic changes to show that our proposed methods (PBR-opt and PBR-opt-mr) enhance the convergence speed of potentials. Moreover, we show that PBR-opt-mr can be adapted to large-scale networks in Section 5.3. Next, in Section 5.4 for demonstrating the robustness of our proposal, we consider a changing environment where several nodes fail but the controller does not collect network topology information and thus cannot detect these failures.

In conducting simulation experiments, for the network simulator we use an eventdriven packet-level simulator written in Visual C++ that calls MATLAB functions $dhinflmi^2$ to design the optimal controller and $balred^3$ to obtain a reduced-order model on a 64-bit operating system with Intel Xeon CPU with 2.7GHz and 64GB memory. The simulator is of our own making, and in the MAC layer, each node sends information about its own potential to its neighbors for their potential updates using intermittent receiver-driven data transmission (IRDT) [Kominami et al. 2013], an asynchronous receive-drive data transmission protocol, in the MAC layer. Note that this setting does not mean that our proposal depends on specific MAC layer protocols such as IRDT. In the physical layer, we use a simple disk model for wireless communication, where the packet reception between two nodes is successful if two nodes are within the predefined communication range. Additionally, we use a simple packet collision model where a receiver node always drops a packet if other packets arrive during the reception of the packet. In the simulator, the asynchrony of components (i.e., the controller and all nodes) is implemented as follows. The controller and all nodes do not match their timing to provide feedback and update their potentials. The asynchronous behavior of all nodes is implemented on a single thread with an event queue in our simulator. Note that we set the interval of the control feedback by the controller and the intervals of potential updates in each node to be equal in accordance with Equations (10) and (14) so that the controller can estimate potentials of nodes with small errors.

5.1. Simulation Settings

We evaluate changes of potentials in potential-based routing and the number of data packets delivered to each sink node after traffic changes or node failures. Network models used in evaluations consist of an external controller and sensor/sink nodes. Only sink nodes are directly connected to the controller, so the controller monitors network states via sink nodes and sends suitable control inputs to sink nodes. The

²*dhinflmi* is a function for designing the H^{∞} optimal controller [Gahinet and Apkarian 1994] of a discrete system. Given parameter matrices A, B_1 , B_2 , C, D_1 , and D_2 for the system dynamics $X(t + 1) = AX(t) + B_1D(t) + B_2u(t)$ and the information $Y(t) = CX(t) + D_1D(t) + D_2u(t)$ that is observed by the controller, *dhinflmi* computes design parameters A_c , B_c , C_c , and D_c , which are referred to in Equations (12) and (13). The matrix D(t) shows nonstability fluctuations, u(t) are control inputs.

³*balred* is a function for reducing the model order [Varga 1991]. Given design parameter matrices A, B_1, B_2 , C, and D_1 for the original model and the degree h of a reduced-order model, *balred* computes a h-th order approximation of the original model.

controller provides feedback to each sink node at interval T_f , whereas at interval T_p , each node updates its next potential value. Typically, $T_f = T_p$ for matching with the potential dynamics described by Equation (2). In real situations, it is difficult to monitor up-to-date potential values of all nodes in the network because of the overhead for collecting potential values, especially when the number of nodes increases. Therefore, here we consider that the controller can directly monitor only nodes within p hops from the sink nodes. In other words, nodes within p hops from the sink nodes send a control message to the sink nodes at interval T_i to notify the controller of their potential values. Otherwise, the controller does not monitor node potentials outside p hops from sink nodes and only estimates them using the model (Equation (14)), which describes potential dynamics through local interactions among nodes. The interval of control message emission T_i is set to 50s, p is set to 2, and the *i*-th element h_i of **H** is set to

$$h_i = \begin{cases} 1, & \text{if node } i \text{ is within } p \text{ hops from sink nodes} \\ 0, & \text{otherwise} \end{cases}$$

At the beginning of the simulation, potential values of all nodes, including sensor nodes and sink nodes, are initialized to 0. During the first 1,000s, each node exchanges its potential values with neighbor nodes and updates its potential value at interval T_p according to Equation (2) so that the potential values are stabilized. For this, we do not generate data packets during that time duration. At 1,000s, data packets begin to be generated at sensor nodes according to the Poisson process with their flow rate.

We evaluate the convergence speed of potentials and data packets delivered to each sink node after traffic changes. To measure the convergence speed of potentials, we define the degree of the potential convergence $\epsilon_n(t)$ for each node, which is given by

$$\epsilon_n(t) = \frac{|\theta_n - \theta_n(t)|}{|\bar{\theta}_n|},$$

where $\bar{\theta}_n$ corresponds to the target potential value of node *n*. We consider convergence to be achieved when $\epsilon_n(t)$ for all nodes becomes sufficiently small. Convergence time is defined as the minimum time taken by all sensor and sink nodes to satisfy the condition

$$\epsilon_n(t) < c, \tag{18}$$

where $c (\geq 0)$ is a constant value. In an ideal situation where all nodes are completely synchronized and there is no noise, all sensor and sink nodes satisfy Equation (18) in the end even if c = 0, but in an actual situation, not all nodes can satisfy Equation (18). The larger the value of c, the shorter the time needed for the achievement of the potential convergence. Therefore, c needs to be set carefully according to the purpose of the evaluation. In this article, we set c to 0.2 to evaluate the convergence speed strictly.

Energy efficiency is a significant challenging task in wireless sensor networks. With respect to our proposal, the computational cost of each sensor node is fairly low because each sensor node behaves based only on local information and simple rules. This is an inherent characteristic of self-organizing systems. Moreover, our proposal can achieve load balancing by setting flow matrix F so that each sink node receives data packets equally, which we explain in Section 5.2. Finally, our proposal does not depend on specific MAC layer protocols. Therefore, we can reduce energy for data transmission between adjacent nodes by introducing energy-efficient MAC layer protocols, such as IRDT, that we use in this evaluation.

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Fig. 5. Network topology (N = 104).

Parameters are summarized in Table I. Optimal parameters (α , $\beta\sigma$) for PBR-no-ctrl are given in Sheikhattar and Kalantari [2013],⁴ but potentials of nodes diverge to infinity and do not converge. Therefore, we use original settings shown in Table I even for PBR-no-ctrl. All results presented in the following are averaged over 40 simulation runs for each parameter setting.

5.2. Performance Evaluation of the Reduced-Order Controller

Here we evaluate the convergence speed considering constraints in wireless sensor networks.

Figure 5 shows the network model with 104 nodes (including 4 sink nodes) used for this evaluation. Sink nodes are illustrated with squares and sensor nodes with dots in the figure. In this network model described in Figure 5, every sensor node is within three hops from a sink node. The number of sensor nodes within two hops from a sink node is 86—that is, the external controller observes potentials of 86 sensor nodes in this network. In this evaluation, at 10,000s from the beginning of the simulation, data

⁴Optimal α and $\beta\sigma$ are respectively given by $\frac{1-\xi}{1+\xi}$ and $\frac{\alpha+1}{2}$ for the fastest convergence rate of potentials,

where $\xi = \sqrt{1 - \nu_1^2}$. ν_1 is the spectrum radius of the matrix $\mathbf{S} = \mathbf{I}_{N \times N} - \mathbf{G}^{-1} \Gamma$, where \mathbf{G} is the degree matrix and Γ is the graph Laplacian of the network. The spectrum radius of the matrix \mathbf{S} is defined as the max value among absolute values of eigenvalues of \mathbf{S} . See Sheikhattar and Kalantari [2013] for more details.

packet generation rates of sensor nodes are changed to examine the convergence speed of our method. Data packet generation rates are initially set to be 0.0005 packets/s for sensor nodes in the left half of Figure 5 (illustrated with black dots) and 0.0015 packets/s for sensor nodes in the right half (illustrated with gray dots). After traffic changes at 10,000s, data packet generation rates are increased to 0.0015 packets/s for the left-half sensor nodes and decreased to 0.0005 packets/s for the right-half nodes. The average data generation rate of a node of 0.001 packets/s corresponds to $\bar{f_n} = 1$; therefore, before traffic changes, the flow rate vector $\mathbf{F} = [\bar{f_1} \cdots \bar{f_N}]^T$ is given by

$$\bar{f}_{i} = \begin{cases} 0.5, & \text{if } i \in \mathcal{N}_{sen_{l}} \\ 1.5, & \text{if } i \in \mathcal{N}_{sen_{r}} \\ -\frac{|\mathcal{N}_{sen}|}{|\mathcal{N}_{sin}|}, & \text{if } i \in \mathcal{N}_{sin} \end{cases}$$
(19)

where \mathcal{N}_{sen} corresponds to the set of sensor nodes and \mathcal{N}_{sin} corresponds to that of sink nodes. \mathcal{N}_{senl} is the set of sensor nodes in the left half of Figure 5, and \mathcal{N}_{senr} is that of sensor nodes in the right half. Note that we construct the potential fields such that all sink nodes can receive data packets equally because load balancing is a challenging task in wireless sensor networks. Thus, the flow rate at each sink node is ideally given by $|\mathcal{N}_{sen}|/|\mathcal{N}_{sin}|$ (=25). Similarly, after the traffic changes, the *i*-th element of **F** is given by

$$\bar{f}_{i} = \begin{cases} 1.5, & \text{if } i \in \mathcal{N}_{sen_{l}} \\ 0.5, & \text{if } i \in \mathcal{N}_{sen_{r}} \\ -\frac{|\mathcal{N}_{sin}|}{|\mathcal{N}_{sin}|}, & \text{if } i \in \mathcal{N}_{sin} \end{cases}$$
(20)

In this evaluation, we conduct simulation in the case in which flows of data packets massively change the network as mentioned earlier to show that our proposal can adapt to massive environmental changes. A system that can adapt to massive environmental changes is considered to be able to adapt to minor fluctuations. Moreover, we assume that the data generation rate does not change until 10,000s from the beginning of the simulation to compare our proposal and the non-control scheme under the same conditions.

In the evaluation of PBR-opt-mr, degree h of the reduced-order model is set to 3 and 100. As explained in Section 4, with lower h, the computational cost is lower but the estimation error is larger, which results in the slow convergence speed of potentials. Therefore, we evaluate not only in the case with low h (=3) but also in the case with high h (=100) for clarifying how the degree of the reduced-order model effects potential convergence.

In the evaluation of PBR-opt, the design parameters $(\mathbf{A}_c, \mathbf{B}_c, \mathbf{C}_c, \mathbf{D}_c)$ cannot be calculated with the *dhinflmi* function when N = 104 (i.e., the size of the model is 208) due to the considerable computational cost. This motivates us to reduce the computational burden upon controller parameter design and controller state updates. For comparison, we utilize an optimal static controller designed by the *dlqr* function,⁵ where the controller is assumed able to monitor the latest conditions of all nodes with no communication delay or control overhead (i.e., $h_i = 1$ for all nodes). It is known that static controllers can achieve satisfactory performance with such state feedback settings [Zhou et al. 1995].

⁵The function dlqr is a linear-quadratic state-feedback regulator for a discrete-time state space system. Given parameter matrices A, B, q, and r for the system dynamics X(t + 1) = X(t) + Bu(t), dlqr computes K, S, and e. Then control inputs u(t) = -KX(t) using the optimal gain matrix K minimize the quadratic cost function $J(u) = \sum_{n=1}^{\infty} (q||X(n)|| + r||u(n)||)$ for the system. Note that the controller needs to observe all nodes for calculating control inputs by u(t) = -KX(t), although the computational cost for calculating K with the dlqr function is much lower than that for calculating A_c , B_c , C_c , and D_c with the dhinflmi function.



Fig. 6. Potential convergence in case of traffic changes.

Scheme	Convergence Time [s]
PBR-no-ctrl	70,417
PBR-opt	9,174
PBR-opt-mr $(h = 3)$	11,765
PBR-opt-mr ($h = 100$)	9,313

Table II. Potential Convergence Time

Figure 6 shows changes of potential values of the non-control scheme, PBR-opt, and PBR-opt-mr with h = 3 and 100. More exactly, Figure 6 plots $\mathbf{X}(t) = \mathbf{\Theta} - \mathbf{\Theta}(t)$ against time t, and the potential convergence is achieved when each element of $\mathbf{X}(t)$ is sufficiently close to 0—that is, Equation (18) is satisfied. In the figure, thick lines correspond to potential changes of the two sink nodes, and thin lines correspond to those of other sensor nodes. Sink node potentials change more than those of sensor nodes, because sink nodes receive feedback inputs \mathbf{u} directly from the controller. Feedback inputs provided to each sink node are different, so the potential change of each sink node varies. On the other hand, sensor nodes are indirectly affected by feedback inputs via sink nodes through potential updates by Equation (2). As shown in the figure, if the potential value of a sink node increases (respectively, decreases) due to feedback inputs from the controller, potential values of sensor nodes near the sink node also increase (respectively, decrease).

As shown in Figure 6, our proposal can enhance the convergence speed of potentials, as compared to the case of no control scheme, and Table II shows the time needed from traffic changes until the potential convergence is achieved when c = 0.2. Figure 6 shows potential changes only within 10,000 to 30,000s, but it takes 70,417s for the potential

convergence (i.e., satisfying Equation (18)) with PBR-no-ctrl as shown in Table II. As a result, potential convergence is accelerated by about 5.99 times and 7.56 times due to PBR-opt-mr with h = 3 and 100, respectively. Compared to PBR-opt, potential convergence speed is 0.780 times and 0.985 times using PBR-opt-mr with h = 3 and 100, respectively. These results indicate that if h is properly chosen, the convergence speed of our proposal can be approximately equal to that of PBR-opt (about 7.68 times as fast as the original). The convergence speed of PBR-opt-mr with h = 3 is a little bit slower than that of PBR-opt-mr with h = 100, because the smaller the h, the larger the approximation error. Specifically, if h is smaller than 3, potentials diverge and cannot converge to targeted potential values, although we do not show the result in this article. However, from the point of view of computational cost, when h is 3, PBR-opt-mr can enhance the convergence speed with much lower computational cost than PBR-opt, as the cost of the design parameter calculation is reduced from $O(N^3)$ to $O(h^3)$ and the cost of the control input calculation is reduced from $O(N^2)$ to $O(h^2)$ by the model reduction. Note that the computational cost for designing the controller

by the model reduction. Note that the computational cost for designing the controller of PBR-opt using *dhinflmi* is too large to compute in a practical amount of time, as explained earlier. Therefore, PBR-opt cannot adapt to large-scale networks. It is worth mentioning that the controller of PBR-opt-mr observes only node potentials within two hops of sink nodes (i.e., p = 2). The controller estimates potentials of nonobservable and future potentials using Equation (14) and then determines the optimal feedback inputs. This indicates that the controller does not need to collect information of the entire network to enhance the convergence speed of potentials. In general, it becomes more difficult for the controller to estimate potentials correctly as the number of nodes observed by the controller becomes smaller. We will investigate the trade-off between the number of observable nodes and the accuracy of the potential estimation in future work.

At about 14,000s in Figure 6(c), potentials of nodes rapidly fluctuate. This is because the controller cannot receive control messages, which collect potential information and inform the controller about them, due to packet drops, so the controller temporarily provides irrelevant control inputs. However, such fluctuations are temporary, and potentials immediately converge again. Therefore, these fluctuations have little effect on the transmission of data packets as shown in Figure 7(c), which is mentioned later. Moreover, in the case in which the controller collects potential information via sensor nodes (i.e., not sink nodes), traffic congestion around sink nodes is reduced, and therefore these fluctuations caused by packet drops will be reduced.

Estimation errors of potentials cannot be completely avoided in real networks. This is because nodes are not synchronized unlike the potential dynamics described by Equation (2) and the potential values the controller collects are not always correct due to communication delays, dropped data, or interference. In this evaluation, traffic congestion occurs around sink nodes because the controller collects the network information via sink nodes, and 1.32% to 1.33% of the control messages for collecting potential values are dropped when using PBR-opt-mr with h = 3 and 100, respectively, leading to estimation errors of potentials. Moreover, the asynchrony of the controller and nodes is also a cause of estimation errors because PBR-opt-mr (and also PBR-opt) inherently assumes synchronous systems. Nevertheless, our proposal can achieve fast convergence of potentials despite such errors, which clearly shows that our proposal works well in an asynchronous environment where noise or disturbance exists.

Figure 7 shows the average number of data packets delivered to each sink node every 1,000s. In each case, the number of data packets delivered to each sink node becomes disproportionate after the traffic changes at 10,000s. Then sink nodes gradually become able to receive data packets equally, because potentials are updated to adapt to the current packet rate. We can observe that the traffic convergence is also accelerated by



Fig. 7. Data packets delivered to each sink node in case of traffic changes.

optimal feedback. This is because the potential convergence speed is enhanced by the optimal feedback mechanism.

One problem we can find is that our proposal reduces the average number of data packets delivered to each sink node immediately after traffic changes. This is because some sink nodes temporarily have the largest potential values within their communication ranges according to the control inputs, so data packets cannot arrive at sink nodes. Therefore, a partial set of data packets would drop when the controller makes large changes to the potentials, which contributes to the faster convergence speed of potentials. However, the data packet drops are immediately reduced and the traffic finally converges faster than the non-control scheme because of the faster potential convergence. Note that in an actual situation, data packets may be retransmitted instantly. Here we evaluate only the case in which data packets are never retransmitted, because the main purpose of this article is to reveal the upper limit of convergence speed of self-organizing systems. Moreover, Figure 7(c) and (d) show worst-case scenarios for temporary packet drops, because the controller changes sink node potentials (in other words, data packet destinations), so many data packets are dropped when a sink node temporarily gets the highest potential within its communication range due to control inputs. If the controller provides an optimal feedback to several sensor nodes where only some data packets arrive, the number of data packet drops will be smaller. With a lower r, sink nodes are more likely to be assigned higher potential values since the controller can make large changes to the potentials, whereas the recovery speed of data packets delivered to each sink node becomes faster. This indicates that there is a trade-off between the convergence speed of potentials and potential fluctuations.



Network condition (Y(t)) — Control input (u(t))

Fig. 8. Network topology (N = 309). Sink nodes are illustrated with squares and sensor nodes with dots. Only one sink node seems to be connected to the external controller, but all sink nodes actually are wired with the external controller.

It is also worth mentioning that the traffic convergence when using PBR-opt-mr with h = 3 is as fast as that with h = 100, as shown in Figure 7(c) and (d), although potentials oscillate more frequently until the potential convergence is achieved with h = 3 compared to the case of h = 100 (Figure 6(c) and (d)). This indicates that temporary oscillations of potential values because of approximation errors have little effect on data packet transmissions.

In this section, we have shown that an optimal control by the external controller is effective in wireless sensor networks where the capacity and energy of each node are limited. Moreover, PBR-opt-mr enhanced the convergence speed of potentials with much lower computational cost than PBR-opt. That indicates that a reduced-order model reflects the dominant characteristics of the original model. It is also worth mentioning that even when some amount of approximation error exists, fast convergence of potentials can be achieved. In this evaluation, however, we have assumed that the controller has up-to-date topology information. Therefore, in Section 5.4, we conduct a simulation experiment where several nodes fail and the controller cannot detect the failures (i.e., the controller provides feedback inputs that lead potentials to converge to values that are not adequate to the network).

5.3. Scalability of the Reduced-Order Controller

Next we conduct a simulation on a network with 309 nodes including 9 sink nodes to reveal the scalability of our proposal in large-scale networks. Specifically, we prove that our proposal can facilitate convergence of a self-organizing system with lightweight cost even if the network size is large. Figure 8 shows the network model that is used in this evaluation. In this network model described in Figure 8, every sensor node is within



Fig. 9. Potential convergence in the large-scale network.

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Scheme	Convergence Time [s]
PBR-no-ctrl	363,675
PBR-opt	51,712
PBR-opt-mr $(h = 20)$	16,234
PBR-opt-mr ($h = 100$)	16,064

three hops from a sink node. The number of sensor nodes within two hops from a sink node is 271—that is, the external controller observes potentials of 271 sensor nodes in this network. In this evaluation, at 10,000s from the beginning of the simulation, data packet generation rates of sensor nodes are changed to examine the convergence speed of our method similarly to the scenario of the evaluation in Section 5.2. Data packet generation rates are initially set to be 0.0005 packets/s for sensor nodes in the left half of Figure 8 (illustrated with black dots) and 0.0015 packets/s for sensor nodes in the right half (illustrated with gray dots). After traffic changes at 10,000s, data packet generation rates are increased to 0.0015 packets/s for the left-half sensor nodes and decreased to 0.0005 packets/s for the right-half nodes. The average data generation rate of a node of 0.001 packets/s corresponds to $\bar{f}_n = 1$, so before traffic changes, the flow rate vector $\boldsymbol{F} = [\bar{f}_1 \cdots \bar{f}_N]^T$ is given by Equation (19), and after the traffic changes, \boldsymbol{F} is given by Equation (20).

Figure 9 shows changes of potential values of the non-control scheme, PBR-opt, and PBR-opt-mr with h = 20 and 100, and Table III shows the time needed from traffic changes until the potential convergence is achieved when c = 0.2. Figure 6 shows

potential changes only within 10,000 to 30,000s, but it takes more than 20,000s for the potential convergence with PBR-no-ctrl and PBR-opt as shown in Table III. As a result, the potential convergence speed is accelerated by about 7.03 times, 22.4 times, and 22.6 times due to PBR-opt and PBR-opt-mr with h = 20 and 100, respectively, compared to the case with PBR-no-ctrl. Table III indicates that potentials converge faster with the larger number of state variables. This characteristic is valid because the controller is generally able to estimate potentials of nodes more correctly with the larger number of state variables. Table III also shows that the convergence speed of potentials with PBR-opt-mr is faster than that with PBR-opt. Feedback inputs provided by the external controller with PBR-opt are optimal in an ideal situation where the external controller and all nodes are synchronized and the external controller can obtain potential information without any communication delays. However, in a real situation, the optimality is not guaranteed because there are estimation errors due to the asynchrony of the system, data packet drops, and communication delays. Therefore, the convergence speed with PBR-opt is not always faster than that with PBR-opt-mr. On the contrary, the convergence speed of potentials with PBR-opt seems not to be so different from that with PBR-opt-mr in Figure 9. This indicates that PBR-opt works as well as PBR-opt-mr in the real setting of wireless sensor networks.

Compared to the evaluation of the smaller-scale network in Section 5.2, degree hof a reduced-order model needs to be larger. This is because a reduced-order model needs a larger number of state variables to keep the dominant characteristics of the original model with a larger number of nodes. When h is set to 3 as similar to the evaluation in Section 5.2, the controller cannot estimate correctly potentials of nodes and the potential convergence speed is quite slow, although we do not show the result in Figure 9. In this article, we only show the result of evaluations using PBR-opt-mr with h = 20 and 100, but our proposal can achieve the potential convergence even if h = 1 that is, the lower bound of degree h for the network shown in Figure 8 is 1, which is smaller than the lower bound of *h* for the network of 104 nodes shown in Figure 5 (=3). The lower bound of h deeply depends on the network topology, so we will investigate the scalability of the reduced-order model in the future. In PBR-opt-mr, the computational cost for designing the controller is $O(h^3)$ and that for calculating control inputs is $O(h^2)$. Therefore, the computational cost of PBR-opt-mr with h = 100 in Section 5.2 and that in Section 5.2 is approximately the same, although the number of nodes is 104 and 309, respectively. Moreover, the number of controllable and observable nodes increases compared to the evaluation of the smaller network. That is because the communication overhead for collecting the information of nodes far from observable nodes is large, and these nodes also take a long time to receive the effect of the optimal feedback that is provided to sink nodes.

Figure 10 shows the number of data packets delivered to each sink node every 1,000s. As shown in this figure, the traffic convergence speed is accelerated because the potential convergence speed is improved by an optimal feedback. As shown in Figure 9(c) and (d), many sink nodes have potential values higher than their target potential values according to control inputs provided by the controller. Therefore, the decreased rate of data packet arrivals in sink nodes is higher than that in the evaluation of 104 nodes. However, this is not always true because control inputs deeply depend on the network topology and flow rates. From Figure 10, the average number of data packets delivered to each sink node with PBR-opt-mr is a bit smaller than that with PBR-no-ctrl and PBR-opt. This is because a partial set of data packets and control messages are delivered to sink nodes. The number of data packet and control messages dropped will be reduced if the controller collects potential information via several sensor nodes.



Fig. 10. Data packets delivered to each sink node in the large-scale network.

In conclusion, our proposal can accelerate the convergence speed even in large-scale networks with the lower computational cost. However, the lower-bound degree of a reduced-order model deeply depends on the network topology.

5.4. Robustness Against Node Failures

We finally investigate the robustness of PBR-opt-mr. The network model of 104 nodes is used in this evaluation. Here, several nodes fail at 10,000s from the start of the simulation, and we examine changes of potentials and data packets delivered to each sink node after node failures. At 10,000s from the beginning of the simulation, q nodes fail. Failing nodes are randomly selected from nodes more than p hops from the sink nodes, so the controller cannot detect them. Instead, each sensor node constantly detect its current neighbors in a self-organizing manner. To avoid the influence of traffic changes, data packet generation rates of all sensor nodes are fixed at 0.001 packets/s, so the *i*-th element f_i of **F** is given by

$$ar{f_i} = \left\{ egin{array}{c} 1, & ext{if } i \in \mathcal{N}_{sen} \ -rac{|\mathcal{N}_{sen}|}{|\mathcal{N}_{sin}|}, & ext{if } i \in \mathcal{N}_{sin} \end{array}
ight.$$

The controller monitors only nodes within p (=2) hops from the sink nodes, as in Section 5.2. In other words, nodes within two hops from the sink nodes send a control message to the sink nodes at intervals of T_i (50s) and transfer their potentials to the controller. The degree h of a reduced-order model is set to 3. Other simulation settings are the same as in Section 5.2.



Fig. 11. Potential convergence in case of node failures (PBR-opt-mr with h = 3).

Figure 11 shows the potential changes with q = 0, 3, and 10. With q = 0, no nodes fail. The figure plots $\mathbf{X}(t) = \bar{\mathbf{\Theta}} - \mathbf{\Theta}(t)$ against time t. Note that $\bar{\mathbf{\Theta}}$ of each case is the same as $\bar{\mathbf{\Theta}}$ in the case in which no node fails, because the controller cannot detect the node failures. The reason the convergence speed is faster than Figure 6 in Section 5.2 is that data generation rates of all sensor nodes are fixed in this evaluation. As the figure shows, potentials converge immediately after the node failures, but potential values at the time when they converge are different from $\bar{\mathbf{\Theta}}$. This is because $\bar{\mathbf{\Theta}}$ changes due to the topology change, but the controller cannot detect the change. However, the potential convergence is achieved soon after node failures, and these failures only slightly affect the data packets delivered to each sink node, as explained in the following.

Figure 12 shows the number of data packets delivered to each sink node every 1,000s and the average number of packets. As the figure shows, each sink node receives data packets, equally even in the case in which node failures occur and the controller cannot detect them. With q = 10, the number of data packets delivered to sink nodes varies, but these differences are not very large. Compared to the case in which node failures do not occur, the average number of data packets delivered to each node is smaller when q = 3 or 10. This is because the number of data packets generated in the entire network is decreased due to the node failures. The data arrival rate is shown in Table IV. The data arrival rate corresponds to the ratio of data packets that arrive at any sink nodes to all generated data packets. This result indicates that the data arrival rate in the case in which node failures occur is approximately equal to that in the case in which no node fails. In conclusion, optimal feedback by the external controller can enhance convergence while maintaining the high robustness against inaccurate information due to node failures, which is an essential characteristic of self-organizing systems.



Fig. 12. Data packets delivered to each sink node in case of node failures (PBR-opt-mr with h = 3). Table IV. Data Arrival Bate

q	Data Arrival Rate [%]
0	99.12
3	99.03
10	99.18

6. CONCLUSION AND FUTURE WORK

In self-organizing systems, each component behaves according to only local information, which leads to slow convergence. We propose and evaluate potential-based routing with optimal feedback using a reduced-order model, where a controller monitors and estimates system states and provides optimal feedback for the fastest convergence. Simulation results have shown that optimal feedback using a reduced-order model can facilitate the convergence of potentials while reducing costs for collecting system information and estimating system dynamics. Moreover, our proposal has high scalability and robustness against information loss from node failures.

On the contrary, our proposal continues to have some challenging tasks. First, the optimal feedback mechanism assumes that the controller has the information of the network topology and the flow rates, which reduces the practicality of our proposal. Second, the potential convergence is achieved as a result of the iterative behavior—that is, the controller's optimal feedback and nodes' potential updates—so that potential cannot converge if environmental changes occur more frequently than the iterative behavior. Third, the controllability and the stability of networks depends on the network topology. Finally, the optimal control improves the convergence speed of potentials but also causes potential fluctuations, as shown by simulation results. These fluctuations

lead to data packet drops because sink nodes temporarily have the highest potentials among their neighbors. There is a trade-off between the improvement of the potential convergence speed and potential fluctuations.

We are now studying a distributed version where several controllers independently monitor, estimate, and control the system dynamics. For this purpose, it would be a promising direction to design hierarchical controllers based on the clustered model reduction [Ishizaki et al. 2014]. With optimal control by a controller, considerable control overhead is needed for collecting network states and estimating network dynamics information when the network is large. By introducing a distributed control mechanism, control overhead can be expected to be reduced because it is not necessary to collect network-wide information.

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