# Cooperative Computing with Fragmentable and Mergeable Groups \*

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#### Abstract

This work considers the problem of performing a set of N tasks on a set of P cooperating message-passing processors  $(P \leq N)$ . The processors use a group communication service (GCS) to coordinate their activity in the setting where dynamic changes in the underlying network topology cause the processor groups to change over time. GCSs have been recognized as effective building blocks for fault-tolerant applications in such settings. Our results explore the efficiency of fault-tolerant cooperative computation using GCSs. Prior investigation of this area by Dolev et al. [8] focused on competitive lower bounds, non-redundant task allocation schemes and workefficient algorithms in the presence of fragmentation regroupings. In this work we investigate work-efficient and message-efficient algorithms for fragmentation and merge regroupings. We present an algorithm that uses GCSs and implements a coordinator-based strategy. This algorithm is motivated by the results in [8]. It achieves similar work complexity of  $O(N \cdot f + N)$ for fragmentations, where f is the number of new groups created by dynamic fragmentations. Additionally, our algorithm achieves substantially better message complexity of  $O(N \cdot f + N)$ , and it is able to deal with more general types of group changes. For the analysis of our algorithm we introduce the notion of *view-graphs* that represent the partially-ordered view evolution history witnessed by the processors. For fragmentations and merges, the work W of the algorithm (defined as the worst case total number of task executions counting multiplicities) is not more than  $\min\{N \cdot f + N, N \cdot P\}$ , and the message complexity M is no worse than  $4(N \cdot f + N + P \cdot m)$ , where f and m denote the number of new groups created by fragmentations and merges respectively. Note that the constants are very small and that, interestingly, while the work efficiency depends on the number of groups f created as the result of fragmentations, work does *not* depend on the number of groups m created as the result of merges.

Keywords: Groups, distributed algorithms, work, communication, complexity.

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## 1 Introduction

The problem of cooperatively performing a set of tasks in a decentralized setting where the computing medium is subject to failures is one of the fundamental problems in distributed computing. Variations on this problem have been studied in a variety of settings, e.g., in message-passing models [9, 6] and in shared-memory models [14]. This problem was also studied in the setting of processor groups in partitionable networks [8]. In this setting, the computation can take advantage of group communication services [4], and the processors must perform the tasks and must learn the results of the tasks efficiently, despite the dynamically changing group memberships.

Group communication services (GCS) can be used as effective building blocks for constructing fault-tolerant distributed applications. These services enable the application components at different processors to operate collectively as a group, using the service to multicast messages. The basis of a group communication service is a *group membership service*. Each processor, at each time, has a unique *view* of the membership of the group. The view includes a list of the processors that are members of the group. Views can change and may become different at different processors. There is a substantial amount of research dealing with specification and implementation of GCSs and group-oriented applications, e.g., [1, 2, 13, 7, 10, 16, 19, 21], and verification of GCSs and group-oriented systems, e.g., [5, 15, 11].

When developing group-oriented, and especially partition-aware applications, it is also important to understand the effectiveness of group communication services [23] and the efficiency benefits that can be expected when using group communication services [8]. One of the features of GCSs is their group management facilities that map a variety of failures in the underlying computing medium to changes in group memberships. Faulty communication links can partition the system into several connected components. Failures and recoveries trigger group membership activity that aims to establish a group for every connected component. An adversary that causes frequent and arbitrary failures may prevent applications from making steady computational progress. Thus, it is interesting to study restricted, yet realistic, models of adversaries for which efficient specific algorithms can be developed with the help of common group communication services. Studying the problem of performing a set of tasks on a set of processors in the group-oriented setting provides a convenient and powerful abstraction for understanding the efficiency of cooperative computation. A work-efficient algorithm is presented for this problem by Dolev et al. in [8], along with a lower bound and a scheduling strategy that minimizes redundant work. That algorithm is tolerant of arbitrary sequences of group fragmentations. In this work we continue the study of algorithms that are work-efficient and message-efficient, and that are able to deal with more general changes in group memberships.

Following [8], we investigate an approach whose goal is to utilize the resources of every component of the system during the entire computation. Thus the problem [8] has the following setting: a set of N independent and idempotent tasks must be performed by P processors in a distributed system, where each processor must learn all results. Group communication is used to coordinate the execution of the tasks. Our distributed system model, in addition to the processors and the network, includes a set of input/output ports accessible to the processors. In this model we enable any client of the required computation to query any processor for the results. This makes it mandatory, even for isolated processors, to be able to provide the results of the computation regardless of whether any other processors may already have the results. Thus, it is not sufficient to know that each of the tasks have been performed somewhere. It is also necessary for each processor to learn the results. We refer to this problem as the OMNI-DO problem.

Note that any algorithm that solves the problem in a way where in any group the processors perform no more than  $\Theta(N)$  tasks (counting multiplicities), will have work complexity of  $O(r \cdot N)$ , where r is the total number of new views installed. This makes it not very interesting to study the problem for adversaries that impose arbitrary view changes. Our major goal is develop *precise* upper bounds that describe the efficiency, work and messaging of solving OMNI-DO as functions of the number of tasks N, the number of processors P, and the numbers of distinct group views of *specific* types (fragmentations and merges in this work) installed by the group membership services.

We present an algorithm for the OMNI-DO problem for N tasks and P message-passing processors  $(P \leq N)$  that are interconnected by a network, which is subject to dynamic group fragmentations and merges. We assume a group communication service that provides group management and vieworiented messaging service (Section 2.1). The main complexity result is for the adversary that is restricted to causing fragmentations of groups and merges of groups. This extends the results in [8], which consider only the fragmentation adversary. Our analysis for the fragmentation-and-merge adversary yields analysis for the fragmentations-only adversary as a corollary.

For the fragmentation-and-merge adversary, we distinguish between the views that are installed as the result of fragmentations and the views installed as the result of merges. If r is the total number of views installed, then for the fragmentation-and-merge adversary we have that r = f + m, where f is the number of views due to fragmentations and m is the number of views due to merges. It is also not difficult to see that m < f when all processors initially start in a single group.

We now summarize our results.

- We present a new algorithm, called algorithm AX, that solves the OMNI-DO problem and we analyze it for the fragmentation-and-merge adversary. The algorithm employs a coordinatorbased approach and relies on the underlying group communication service. The algorithm is specified in Section 4 and it extends the approaches in [8] and [3].
- We introduce the notion of *view-graphs* that represent the partially-ordered view evolution history collectively witnessed by the processors (Section 3). We show that these digraphs are acyclic for the fragmentation-and-merge adversary and we use these view-graphs in the complexity analysis of the algorithm. We believe that view-graphs have the potential of serving as a general tool for studying cooperative computing with group communication services.
- For any pattern of fragmentations and merges, the work W of the algorithm is no more than min{N · f + N, N · P}, and the message complexity M is no worse than 4(N · f + N + P · m). Note that f ≤ r and here it is significant that we are expressing the upper bounds using explicit constants instead of the big-oh notation. Both complexity results depend on f, but only the message complexity depends on m. The fact that the work complexity does not depend on m, substantiates the intuition that merges lead to a more efficient computation. The upper bounds are tight for a broad range of view changes. This analysis is presented in Sections 5.1 and 5.2.
- For any pattern of fragmentations (i.e., when m = 0) our algorithm achieves work complexity of O(min{N · f + N, N · P}). This result is essentially the same as the result in [8]. However, our algorithm achieves substantially better message complexity O(N · f + N) as compared to the at least quadratic message complexity of the algorithm in [8]. Message optimization was outside of the scope of [8], yet this improvement was one of our goals. The improvement is largely due to our use of the coordinator-based strategy. These results are in Section 5.3.

Note that it is not difficult to show that when  $f \ge N$ ,  $W = \Omega(N \cdot P)$  and when f < N,  $W = \Omega(N \cdot f)$ . Thus,  $W = \Omega(\min\{N \cdot f, N \cdot P\})$  is a lower bound for OMNI-DO. So our algorithm is optimal with respect to work for the adversaries we consider. Considering optimality for the message complexity is less interesting, since the problem can be solved without any communication.

**Related work.** The problem of efficiently performing a set of tasks using a network of processors in the setting where the network is subject to dynamic changes was considered by Dolev, Segala and Shvartsman [8]. For the N-processor, N-task problem defined in that work, it was shown that for dynamic changes the termination time of any on-line task algorithm can be greater than the termination time of an off-line algorithm by a factor linear in N. An algorithm was also presented in [8] that for arbitrary fragmentations has work  $O(N \cdot f' + N)$ , where f' is the increase in the number of groups due to fragmentations. In comparing our result with the result in [8], we note that our definition of f is slightly different from the definition of fragmentation failures f' in [8]. In order to compare our complexity results with those in [8], we show in this paper that for any pattern of fragmentations allowed by [8] we have f' < f < 2f'. In [8] the work is counted in terms of the rounds executed by the processors. In our analysis we count only the number of task executions (including redundancies). However in our algorithm, for as long as any tasks remain undone in a given group, the processors perform the tasks in rounds, except for the last round. Therefore the difference in work complexity for these two algorithms is at most  $f \cdot N$ . Thus the different definitions of f and f' and of work can be subsumed in the big-oh analysis without substantial variation in the constants.

Group communication services (GCS) have become important as building blocks for faulttolerant distributed systems. Such services enable processors located in a fault-prone network to operate collectively as a group, using the services to multicast messages to group members. Examples of GCS include Isis [2], Transis [7], Totem [19], Newtop [10], Relacs [1], Horus [21] and Ensemble [13]. Examples of recent work dealing with primary groups are [5, 16]. An example of an application using a GCS for load balancing is by Fekete, Khazan and Lynch [15]. To evaluate the effectiveness of partitionable GCSs, Sussman and Marzulo [23] proposed the measure (*cushion*) precipitated by a simple partition-aware application.

Our definition of work follows that of Dwork, Halpern and Waarts [9]. Our fragmentation model creates a setting, within each fragment, that is similar to the setting in which the network does not fragment but the processors are subject to crash failures. Performing a set of tasks in such settings is the subject of several works [3, 6, 9, 12], however the analysis is quite different when work in all fragments has to be considered.

Our distributed problem has an analogous counterpart in the shared-memory model of computation, called the *collect* problem. The collect problem was originally abstracted by Saks, Shavit and Woll [22] (it also appears in Shavit's Ph.D. thesis). Although the algorithmic techniques are different, the goal of having all processors to learn a set of values is similar.

## 2 Definition and Models

A distributed system consists of P processors connected by communication links. Each processor has a unique identifier from the set  $\mathcal{P} = \{1, 2, \dots, P\}$ .

We define a *task* to be any computation that can be performed by a single processor in constant time. We assume that the tasks are independent and idempotent. Our distributed system is charged with the responsibility of performing a set of N tasks that are initially known to all processors. Each task has a unique identifier from the set  $\mathcal{T}$ .

To require that all processors acquire the results of all tasks, our system also includes a set of input/output ports. These ports are only used by the clients of the system to query individual processors for computation results. We do not make any failure assumptions about the input/output ports, in particular, our algorithm does not depend on the failure status of these ports, or the requests from them.

**Definition 2.1** The problem of performing a set of N independent tasks on a set of P message passing processors, where each processor must learn the results of all N tasks, is called the OMNI-DO problem.

The algorithm specification in this paper is done in terms of I/O automata of Lynch and Tuttle [17, 18]. Each automaton models a state machine with states and transitions between states, where actions are associated with sets of state transitions. There are input, output and internal actions. A particular action is enabled if the preconditions of that action are satisfied. The statements given as effects are executed as a program started in the existing state and atomically producing the next state as the result of the transition.

An execution  $\alpha$  of an I/O automaton Aut is a finite or infinite sequence of alternating states and actions (events) of Aut starting with the initial state, i.e.,  $\alpha = s_0, e_1, s_1, e_2, \ldots$ , where  $s_i$ 's are states ( $s_0$  is the initial state) and  $e_i$ 's are actions (events). We denote by execs(Aut) the set of all executions in Aut.

We next state our assumptions about the group communication services and define the work and message complexity measures.

## 2.1 Group Communication Service

We assume a group communication service (GCS) with certain properties. The assumptions are basic, and they are provided by several group communication systems and specifications [24]. The

service maintains group membership information and it is used to communicate information concerning the executed tasks within each group. The GCS provides the following primitives:

NEWVIEW $(v)_p$  informs processor p of a new view  $v = \langle id, set \rangle$ , where id is the identifier of the view and set is the set of processor identifiers in the group. When a NEWVIEW $(v)_p$  primitive is invoked, we say that processor p installs view v.

 $GPMSND(message)_p$  allows processor p to multicast a message to the current group members.

- $GPMRCV(message)_p$  enables processor p to receive multicasts from other processors.
- GP1SND(message, destination)<sub>p</sub> allows processor p to unicast a message to another member of the current group.

 $GP1RCV(message)_p$  enables processor p to receive unicasts from another processor.

To distinguish between the messages sent in different send events, we assume that each message sent by the application is tagged with a unique message identifier.

We assume the following safety properties on any execution  $\alpha$  of an algorithm that uses GCSs:

- 1. A processor is always a member of its view ([24] Prop. 3.1). If NEWVIEW $(v)_p$  occurs in  $\alpha$  then  $p \in v.set$ .
- 2. The view identifiers of the views that each processor installs are monotonically increasing ([24] Prop. 3.2). If event  $NEWVIEW(v_1)_p$  occurs in  $\alpha$  before event  $NEWVIEW(v_2)_p$ , then  $v_1.id < v_2.id$ . This property implies that:
  - (a) A processor does not install the same view twice.
  - (b) If two processors install the same two views, they install these views in the same order.
- 3. For every receive event, there exists a preceding send event of the same message ([24] Prop. 4.1). If  $\text{GPMRCV}(m)_p$  ( $\text{GP1RCV}(m)_p$ ) occurs in  $\alpha$ , then there exists  $\text{GPMSND}(m)_q$ ( $\text{GP1SND}(m, p)_q$ ) earlier in execution  $\alpha$ .
- 4. Messages are not duplicated ([24] Prop. 4.2). If  $\text{GPMRCV}(m_1)_p$  (GP1RCV $(m_1)_p$ ) and GPMRCV $(m_2)_p$  (GP1RCV $(m_2)_p$ ) occur in  $\alpha$ , then  $m_1 \neq m_2$ .
- 5. A message is delivered in the same view it was sent in ([24] Prop. 4.3). If processor p receives message m in view  $v_1$  and processor q (it is possible that p = q) sends m in view  $v_2$ , then  $v_1 = v_2$ .
- 6. In the initial state  $s_0$ , all processors are in the initial view  $v_0$ , such that  $v_0.set = \mathcal{P}$  ([24] Prop. 3.3 with [11, 20]).

We assume the following additional liveness properties on any execution  $\alpha$  of an algorithm that uses GCSs (cf. [24] Section 10):

- 7. If a processor p sends a message m in a view  $v_1$ , then for each processor q in  $v_1$ .set, either q delivers m in  $v_1$ , or p installs a next view  $v_2$ .
- 8. If a new view event occurs at any processor p in view v, then a view change will eventually occur at all processors in  $v.set \{p\}$ .

## 2.2 Regrouping-Numbers and Measures of Efficiency

In this section we define regrouping-numbers and complexity measures. We define the regroupingnumber r of an execution to be the number of NEWVIEW events with distinct view identifiers. (Note that if the same view is installed at multiple processors, this counts for a single regrouping.)

**Definition 2.2** Given an execution  $\alpha$ , we define the regrouping-number  $r_{\alpha}$  as:  $r_{\alpha} = |\{v : \text{NEWVIEW}(v)_p \text{ occurs in } \alpha\}|.$ 

When it is clear from the context, we use r instead of  $r_{\alpha}$  to denote the regrouping-number of execution  $\alpha$ .

We define *adversary models*, in the context of a specific algorithm, in terms of the collections of executions in the presence of an adversary. In the following definitions we assume that the algorithms use a group communication service as presented in Section 2.1.

**Definition 2.3** For an algorithm A, let  $\mathcal{F}_R(A)$  be the adversary model that includes all the possible executions of A, i.e.,  $\mathcal{F}_R(A) = execs(A)$ .

**Definition 2.4** For an algorithm A, let  $\mathcal{F}_{\emptyset}(A)$  be the adversary model that does not cause any NEWVIEW events, i.e.,  $\mathcal{F}_{\emptyset}(A) = \{\alpha : \alpha \in execs(A) \land r_{\alpha} = 0\}.$ 

When it is clear from the context, we use  $\mathcal{F}_{\emptyset}$  instead of  $\mathcal{F}_{\emptyset}(A)$  and  $\mathcal{F}_R$  instead of  $\mathcal{F}_R(A)$ .

It is easy to see that  $\mathcal{F}_{\emptyset} \subseteq \mathcal{F}_R$ . Let  $\mathcal{F}$  be some adversary model such that  $\mathcal{F}_{\emptyset} \subseteq \mathcal{F} \subseteq \mathcal{F}_R$ . In the following definitions we formalize the measures of work and message complexity for the specific  $\mathcal{F}$ . Our definition of work follows that of Dwork, Halpern and Waarts [9].

**Definition 2.5** The work  $W_{\alpha}(N, P)$  of an execution  $\alpha$  of algorithm A in the adversary model  $\mathcal{F}$ , is defined to be  $\sum_{i \in \mathcal{P}} W^{i}_{\alpha}$ , where  $W^{i}_{\alpha}$  is the number of tasks performed by processor i.

**Definition 2.6** For algorithm A in the adversary model  $\mathcal{F}$ , the work complexity  $W_{\mathcal{F}}(N, P, r)$  is defined as:  $W_{\mathcal{F}}(N, P, r) = \max_{\alpha \in \mathcal{F}, r_{\alpha} < r} \{W_{\alpha}(N, P)\}.$ 

**Definition 2.7** The message cost  $M_{\alpha}(N, P)$  of an execution  $\alpha$  of algorithm A in the adversary model  $\mathcal{F}$ , is defined to be  $\sum_{i \in \mathcal{P}} M_{\alpha}^i$ , where  $M_{\alpha}^i$  is the number of messages sent by processor i.

**Definition 2.8** For algorithm A in the adversary model  $\mathcal{F}$ , the message complexity  $M_{\mathcal{F}}(N, P, r)$  is defined as:  $M_{\mathcal{F}}(N, P, r) = \max_{\alpha \in \mathcal{F}, r_{\alpha} \leq r} \{M_{\alpha}(N, P)\}.$ 

## **3** View-Graphs and Specific Adversary Models

This section introduces view-graphs that are used to represent and analyze changes of processors' views in executions. View-graphs are directed graphs (digraphs) that are defined by the states and the NEWVIEW events of executions of algorithms that use group communication services. Representing view changes as digraphs enables us to use common graph analysis techniques to formally reason about the properties of executions. In this paper we deal with adversary models that cause group fragmentations and merges. Our view-graph approach to the analysis of executions is general, and we believe it can be used to study the complexity of computation and other properties of group communication services and algorithms for different adversary models.

#### 3.1 Executions and View-Graphs

Consider an algorithm A that uses a group communication service (GCS). We modify algorithm A by introducing, for each processor *i*, the history variable  $cv_i$  that keeps track of the current view at *i* as follows: In the initial state, we set  $cv_i$  to be  $v_0$ , the distinguished initial view for all processors  $i \in \mathcal{P}$ . In the effects of the NEWVIEW $(v)_i$  action for processor *i*, we include the assignment  $cv_i := v$ . In the rest of the paper, we assume that algorithms are modified to include such history variables. We now define *view-graphs* by specifying how a view-graph is induced by an execution of an algorithm.

**Definition 3.1** Given an execution  $\alpha$  of algorithm A, the view-graph  $\Gamma_{\alpha} = \langle V, E, L \rangle$  is defined to be the labeled directed graph as follows:

- 1. Let  $V_{\alpha}$  be the set of all views v that occur in NEWVIEW $(v)_i$  events in  $\alpha$ . The set V of nodes of  $\Gamma_{\alpha}$  is the set  $V_{\alpha} \cup \{v_0\}$ . We call  $v_0$  the initial node of  $\Gamma_{\alpha}$ .
- 2. The set of edges E of  $\Gamma_{\alpha}$  is a subset of  $V \times V$  determined as follows. For each NEWVIEW $(v)_i$ event in  $\alpha$  that occurs in state s, the edge  $(s.cv_i, v)$  is in E.

3. The edges in E are labeled by  $L: E \to 2^{\mathcal{P}}$ , such that  $L(u, v) = \{i : \text{NEWVIEW}(v)_i \text{ occurs in state s in } \alpha \text{ such that } s.cv_i = u\}.$ 

Observe that the definition ensures that all edges in E of  $\Gamma_{\alpha}$  are labeled.

**Example 1:** Consider the following execution  $\alpha$  (we omit all events other than NEWVIEW and any states that do not precede NEWVIEW events):

 $\alpha = s_0, \text{NEWVIEW}(v_1)_{p_1}, \dots, s_1, \text{NEWVIEW}(v_2)_{p_2}, \dots, s_2, \text{NEWVIEW}(v_3)_{p_4}, \dots, s_3, \text{NEWVIEW}(v_4)_{p_1}, \dots, s_4, \text{NEWVIEW}(v_1)_{p_3}, \dots, s_5, \text{NEWVIEW}(v_4)_{p_2}, \dots, s_6, \text{NEWVIEW}(v_4)_{p_3}, \dots, \text{ where } v_1.set = \{p_1, p_3\}, v_2.set = \{p_2\}, v_3.set = \{p_4\} \text{ and } v_4.set = \{p_1, p_2, p_3\}. \text{ Additionally, } v_0.set = \mathcal{P} = \{p_1, p_2, p_3, p_4\}.$ 

The view-graph  $\Gamma_{\alpha} = \langle V, E, L \rangle$  is given in Figure 1. The initial node of  $\Gamma_{\alpha}$  is  $v_0$ . The set of nodes of V of  $\Gamma_{\alpha}$  is  $V = V_{\alpha} \cup \{v_0\} = \{v_0, v_1, v_2, v_3, v_4\}$ . The set of edges E of  $\Gamma_{\alpha}$  is E = $\{(v_0, v_1), (v_0, v_2), (v_0, v_3), (v_1, v_4), (v_2, v_4)\}$ , since for each of these  $(v_j, v_k)$  the event NEWVIEW $(v_k)_p$ occurs in state  $s_\ell$  where  $s_\ell.cv_p = v_j$  for some certain p (by the definition of the history variable). The labels of the edges are  $L(v_0, v_1) = \{p_1, p_3\}, L(v_0, v_2) = \{p_2\}, L(v_0, v_3) = \{p_4\}, L(v_1, v_4) = \{p_1, p_3\}$ and  $L(v_2, v_4) = \{p_2\}$ , since for each  $p_i \in L(v_j, v_k)$  the event NEWVIEW $(v_k)_i$  occurs in state  $s_\ell$  where  $s_\ell.cv_{p_i} = v_j$ .



Figure 1: Example of a view-graph

Given a graph S and a node v of S, we define indegree(v, S) (outdegree(v, S)) to be the indegree (outdegree) of v in S.

**Lemma 3.2** For any execution  $\alpha$ , indegree $(v_0, \Gamma_\alpha) = 0$ .

**Proof:** In the initial state  $s_0$ ,  $s_0.cv$  is defined to be  $v_0$  for all processors in  $\mathcal{P}$  and  $v_0.set = \mathcal{P}$ . Assume that  $indegree(v_0, \Gamma_\alpha) > 0$ . By the construction of view-graphs, this implies that some processor  $i \in \mathcal{P}$  installs  $v_0$  a second time. But this contradicts property 2(a) of GCS.  $\Box$  **Lemma 3.3** Let  $\alpha$  be an execution and  $\Gamma_{\alpha}|_i$  be the projection of  $\Gamma_{\alpha}$  on the edges whose label includes *i*, for some  $i \in \mathcal{P}$ .  $\Gamma_{\alpha}|_i$  is an elementary path and  $v_0$  is the path's source node.

**Proof:** Let execution  $\alpha$  be  $s_0, e_1, s_1, e_2, \ldots$ . Let  $\alpha^{(k)}$  be the prefix of  $\alpha$  up to the  $k^{th}$  state. i.e.,  $\alpha^{(k)} = s_0, e_1, s_1, e_2, \ldots, s_k$ . Let  $\Gamma^k_{\alpha}$  be the view-graph that is induced by  $\alpha^{(k)}$ . Then define  $\Gamma^k_{\alpha}|_i$  to be the projection of  $\Gamma^k_{\alpha}$  on the edges whose label includes i, for some  $i \in \mathcal{P}$ .

For an elementary path  $\pi$ , we define  $\pi$ .sink to be its sink node.

We prove by induction on k that  $\Gamma_{\alpha}^{k}|_{i}$  is an elementary path, that  $\Gamma_{\alpha}^{k}|_{i}.sink = s_{k}.cv_{i}$  and that  $v_{0}$  is the path's source node.

Basis: k = 0.  $\Gamma_{\alpha}^{0}|_{i}$  has only one vertex,  $v_{0}$ , and no edges  $(\alpha^{(0)} = s_{0})$ . Thus,  $\Gamma_{\alpha}^{0}|_{i}$ . sink =  $s_{0}$ .  $cv_{i} = v_{0}$  and  $v_{0}$  is the source node of this path.

Inductive Hypothesis: Assume that  $\forall n \leq k$ ,  $\Gamma_{\alpha}^{n}|_{i}$  is an elementary path, that  $\Gamma_{\alpha}^{n}|_{i}$ .  $sink = s_{n}.cv_{i}$  and that  $v_{0}$  is the path's source node.

Inductive Step: n = k + 1. For state  $s_{k+1}$  we consider two cases:

Case 1: If event  $e_{k+1}$  is not a NEWVIEW event involving processor i, then  $\Gamma_{\alpha}^{k+1}|_i = \Gamma_{\alpha}^k|_i$ . Thus, by inductive hypothesis,  $\Gamma_{\alpha}^{k+1}|_i$  is an elementary path and  $v_0$  is its source node. From state  $s_k$  to state  $s_{k+1}$ , processor i did not witness any new view. By the definition of the history variable,  $s_{k+1}.cv_i = s_k.cv_i$ . Thus,  $\Gamma_{\alpha}^{k+1}|_i.sink = s_k.cv_i = s_{k+1}.cv_i$ .

Case 2: If event  $e_{k+1}$  is a NEWVIEW $(v)_i$  event that involves processor *i*, then by the construction of the view-graph,  $(s_k.cv_i, v)$  is a new edge from node  $s_k.cv_i$  to node *v*. By inductive hypothesis,  $\Gamma_{\alpha}^k|_{i.sink} = s_k.cv_i$ . Since our GCS does not allow the same view to be installed twice (property 2(a)),  $v \neq u$  for all  $u \in \Gamma_{\alpha}^k|_i$ . Thus,  $\Gamma_{\alpha}^{k+1}|_i$  is also an elementary path, with  $v_0$  its source node and  $\Gamma_{\alpha}^{k+1}|_{i.sink} = v$ . From state  $s_k$  to state  $s_{k+1}$ , processor *i* installs the new view *v*. By the definition of the history variable,  $s_{k+1}.cv_i = v$ . Thus,  $\Gamma_{\alpha}^{k+1}|_{i.sink} = s_{k+1}.cv_i$ . This completes the proof.  $\Box$ 

**Corollary 3.4** Any view-graph  $\Gamma_{\alpha}$ , induced by any execution  $\alpha$  of algorithm A is a connected graph. **Proof:** The result follows from Definition 3.1(2), from the observation that all edges of the viewgraph are labeled and from Lemma 3.3

**Definition 3.5** For a view-graph  $\Gamma_{\alpha} = \langle V, E, L \rangle$ , a fragmentation subgraph is a connected labeled subgraph  $S = \langle V_S, E_S, L_S \rangle$  of  $\Gamma_{\alpha}$  such that:

1. S contains a unique node v such that indegree(v, S) = 0; v is called the fragmentation node of S.

- 2.  $V_S = \{v\} \cup V'_S$ , where  $V'_S$  is defined to be  $\{w : (v, w) \in E\}$
- 3.  $E_S = \{(v, w) : w \in V'_S\}$
- 4.  $L_S$  is the restriction of L on  $E_S$ .
- 5.  $\bigcup_{w \in V'_{c}} (w.set) = v.set$
- 6.  $\forall u, w \in V'_S$  such that  $u \neq w$ ,  $u.set \cap w.set = \emptyset$
- 7.  $\forall w \in V'_S, L_S(v, w) = w.set$

In the analysis of algorithms, we are going to be referring to all NEWVIEW events that collectively induce a fragmentation subgraph for a fragmentation node v as a *fragmentation*.

**Example 2:** The shaded area A in Figure 1 shows the fragmentation subgraph  $S = \langle V_S, E_S, L_S \rangle$  of  $\Gamma_{\alpha}$  from Example 1. Here  $V_S = \{v_0, v_1, v_2, v_3\}, E_S = \{(v_0, v_1), (v_0, v_2), (v_0, v_3)\}$  and the labels are the labels of  $\Gamma_{\alpha}$  restricted on  $E_S$ . We can confirm that S is a fragmentation subgraph by examining the individual items of Definition 3.5.

**Definition 3.6** For a view-graph  $\Gamma_{\alpha} = \langle V, E, L \rangle$ , a merge subgraph is a connected labeled subgraph  $S = \langle V_S, E_S, L_S \rangle$  of  $\Gamma_{\alpha}$  such that:

- 1. S contains a unique node v such that outdegree(v, S) = 0 and indegree(v, S) > 1; v is called the merge node of S.
- 2.  $V_S = \{v\} \cup V'_S$ , where  $V'_S$  is defined to be  $\{w : (w, v) \in E\}$

3. 
$$E_S = \{(w, v) : w \in V'_S\}$$

- 4.  $L_S$  is the restriction of L on  $E_S$ .
- 5.  $\bigcup_{w \in V'_{\varsigma}} (w.set) = v.set$
- 6.  $\forall u, w \in V'_S$  such that  $u \neq w$ ,  $u.set \cap w.set = \emptyset$
- 7.  $\bigcup_{w \in V'_S} L_S(w, v) = v.set$

A regrouping of a group  $g_1$  to a group  $g_2$  such that  $g_1.set = g_2.set$  can be represented either as a fragmentation subgraph or as a merge subgraph. In this paper we choose to represent it as a fragmentation subgraph by requiring that indegree(v, S) > 1 for any merge node v.

In the analysis of algorithms, we are going to be referring to all NEWVIEW events that collectively induce a merge subgraph for a merge node v as a *merge*.

**Example 3:** The area B in Figure 1 of Example 1 shows the merge subgraph  $S = \langle V_S, E_S, L_S \rangle$  of  $\Gamma_{\alpha}$ , where  $V_S = \{v_1, v_2, v_3, v_4\}$ ,  $E_S = \{(v_1, v_4), (v_2, v_4)\}$  and the labels are the labels of  $\Gamma_{\alpha}$  restricted on  $E_S$ . We can verify this by examining all conditions of Definition 3.6.  $\Box$ .

**Definition 3.7** Given an execution  $\alpha$  of algorithm A, let  $frag(\Gamma_{\alpha})$  be the set of all the distinct fragmentation nodes in the induced view-graph  $\Gamma_{\alpha}$ .

**Definition 3.8** Given an execution  $\alpha$  of algorithm A, let  $merg(\Gamma_{\alpha})$  be the set of all the distinct merge nodes in the induced view-graph  $\Gamma_{\alpha}$ .

**Definition 3.9** A view-graph  $\Gamma_{\alpha}$  such that all of its non-terminal nodes are in frag( $\Gamma_{\alpha}$ ), is called a fragmentation view-graph.

**Definition 3.10** A view-graph  $\Gamma_{\alpha}$  such that each of its non-terminal nodes is either in frag( $\Gamma_{\alpha}$ ) or it is an immediate ancestor of a node that is in  $merg(\Gamma_{\alpha})$  is called an fm view-graph.

For  $\Gamma_{\alpha}$  in the example in Figure 1 we have  $v_0 \in frag(\Gamma_{\alpha})$  by Definition 3.7. Also,  $v_4 \in merg(\Gamma_{\alpha})$ by Definition 3.8; additionally, the nodes  $v_1$  and  $v_2$  are immediate ancestors of  $v_4 \in merg(\Gamma_{\alpha})$ . By Definition 3.10,  $\Gamma_{\alpha}$  is an fm view-graph. Observe that  $\Gamma_{\alpha}$  is a DAG. This is true for all fm view-graphs:

**Theorem 3.11** Any fm view-graph  $\Gamma_{\alpha} = \langle V, E, L \rangle$  is a Directed Acyclic Graph (DAG).

**Proof:** Assume that  $\Gamma_{\alpha}$  is not a DAG. Thus, it contains at least one cycle. Let  $((v_1, v_2)(v_2, v_3) \dots (v_k, v_1))$  be an elementary cycle of  $\Gamma_{\alpha}$ . By the construction of view-graphs (Definition 3.1(3)) and by the monotonicity property (property 2) of GCS,  $v_i.id < v_{i+1}.id$  for  $1 \le i \le k$  and  $v_k.id < v_1.id$ . But, by the transitivity of "<",  $v_1.id < v_k.id$ , a contradiction.

However, not all view-graphs are DAGs. It is also not difficult to see that any fragmentation view-graph is a rooted tree.

In the complexity analysis we use the following fact.

**Fact 3.12** In any (non-empty) DAG, there is at least one vertex, such that all of its descendants have outdegree 0.

#### 3.2 Adversary Models

Let A be an algorithm that uses GCS, as presented in Section 2.1. We now define two adversary models that are more restrictive than  $\mathcal{F}_R(A)$ , but less restrictive than  $\mathcal{F}_{\emptyset}(A)$ .

**Definition 3.13** For any algorithm A the fragmentation adversary  $\mathcal{F}_F(A)$  is the set of all executions of A, such that each execution induces a fragmentation view-graph.

**Definition 3.14** For any algorithm A the fragmentation-and-merge adversary  $\mathcal{F}_{FM}(A)$  is the set of all executions of A, such that each execution induces an fm view-graph.

It is easy to see that  $\mathcal{F}_{\emptyset}(A) \subseteq \mathcal{F}_{F}(A) \subseteq \mathcal{F}_{FM}(A) \subseteq \mathcal{F}_{R}(A)$ .

**Definition 3.15** Given an execution  $\alpha$  of algorithm A, and  $\Gamma_{\alpha} = \langle V, E, L \rangle$ , we define:

- 1. the fragmentation-number  $f_{\alpha} = |\{w : \text{NEWVIEW}(w)_p \text{ occurs in } \alpha \land (v, w) \in E \land v \in frag(\Gamma_{\alpha})\}|$
- 2. the merge-number  $m_{\alpha} = |\{v : \text{NEWVIEW}(v)_p \text{ occurs in } \alpha \land v \in merg(\Gamma_{\alpha})\}|$

Note that for an algorithm A and for an execution  $\alpha \in \mathcal{F}_{FM}(A)$ , by Definitions 2.2 and 3.15,  $r_{\alpha} = f_{\alpha} + m_{\alpha}$ . Also, by Definitions 3.7, 3.8 and 3.15,  $f_{\alpha} > m_{\alpha}$ . Observe that in the adversary model  $\mathcal{F}_{F}$ ,  $r_{\alpha} = f_{\alpha}$  and  $m_{\alpha} = 0$ .

## 4 Algorithm AX

We now present the algorithm, called algorithm AX, that deals with *regroupings* and that relies on a GCS as specified in Section 2.1. The analysis of the algorithm is in Section 5.

Algorithm AX uses a coordinator approach within each group view. The high level idea of the algorithm is that each processor performs (remaining) tasks according to a load balancing rule, and a processor completes its computation when it learns the results of all the tasks.

**Task allocation.** The set T of the initial tasks is known to all processors. During the execution each processor i maintains a local set D of tasks already done, a local set R of the corresponding results, and the set G of processors in the current group. (The set D may be an underestimate of the set of tasks done globally.) The processors allocate tasks based on the shared knowledge of the processors in G about the tasks done. For a processor i, let rank(i, G) be the rank of i in G when processor identifiers are sorted in ascending order. Let U be the tasks in T - D. For a task u in U, let rank(u, U) be the rank of u in U when task identifiers are sorted in ascending order. Our load balancing rule for each processor i in G is that:

- if  $rank(i, G) \leq |U|$ , then processor *i* performs task *u* such that rank(u, U) = rank(i, G);
- if rank(i, G) > |U|, then processor *i* does nothing.

Algorithm structure. The algorithm code is given in Figure 2 using I/O automata notation [18]. The algorithm uses the group communication service to structure its computation in terms of *rounds* numbered sequentially within each group view.

Initially all processors are members of the distinguished initial view  $v_0$ , such that  $v_0.set = \mathcal{P}$ . Rounds numbered 1 correspond to the initial round either in the original group or in a new group upon a regrouping as notified via the NEWVIEW event. If a regrouping occurs, the processor receives the new set of members from the group membership service and starts the first round of this view (NEWVIEW action). At the beginning of each round, denoted by a round number Rnd, processor iknows G, the local set D of tasks already done, and the set R of the results. Since all processors know G, they "elect" the group coordinator to be the processor which has the highest processor id (no communication is required since the coordinator is uniquely identified). In each round each processor reports D and R to the coordinator of G (GP1SND action). The coordinator receives and collates these reports (GP1RCV action) and sends the result to the group members (GPMSND action). Upon the receipt of the message from the coordinator, processors update their D and R, and perform work according to the load balancing rule (GPMRCV action).

For generality, we assume that the messages may be delivered by the GCS out of order. The set of messages within the current view is saved in the local variable A. The saved messages are also used to determine when all messages for a given round have been received. Processing continues until each member of G knows all results (the processors enter the *sleep* stage). When requests for computation results arrive from a port q (REQUEST action), each processor keeps track of this in a local variable *requests*, and, when all results are known, sends the results to the port (REPORT action).

**Correctness:** We now show the safety of algorithm AX.

We first show that no processor stops working as long as it knows of any undone tasks.

## **Theorem 4.1 (Safety 1)** For all states of any execution of Algorithm AX it holds that $\forall i \in \mathcal{P} : D_i \neq T \Rightarrow Phase \neq sleep$

**Proof:** The proof follows by examination of the code of the algorithm, and more specifically from the code of the input action  $\text{GPMRCV}(\langle j, H, Q, round \rangle)_i$ .

Note that the implication in Theorem 4.1 cannot be replaced by iff ( $\Leftrightarrow$ ). This is because if  $D_i = T$ , we may still have  $Phase \neq sleep$ . This is the case where processor *i* becomes a member of a group in which the processors do not know all the results of all the tasks.

## Data types:

 $\begin{array}{l} \mathcal{T}: \text{ tasks} \\ \mathcal{T}: \text{ tasks} \\ \mathcal{R}: \text{ results} \\ \textit{Result}: \mathcal{T} \to \mathcal{R} \\ \mathcal{M}es: \text{ messages} \\ \mathcal{P}: \text{ processor ids} \\ \mathcal{G}: \text{ group ids} \\ \textit{views} = \mathcal{G} \times 2^{\mathcal{P}}: \text{ views, selectors } id \text{ and } set \\ \mathcal{IO}: \text{ input/output ports} \\ \textbf{States:} \\ \mathcal{T} \in 2^{\mathcal{T}}, \text{ the set of } N = |\mathcal{T}| \text{ tasks} \\ D \in 2^{\mathcal{T}}, \text{ the set of done tasks, initially } \emptyset \\ \textit{R} \in 2^{\mathcal{R}}, \text{ the set of known results, initially } \emptyset \\ \textit{G} \in 2^{\mathcal{P}}, \text{ current members, init. } v_{0}.set = \mathcal{P} \\ \textit{A} \in 2^{\mathcal{M}es}, \text{ messages since last NEWVIEW, init. } \emptyset \\ \textit{Rnd} \in \mathbf{N}, \text{ round number, initially } 1 \\ \textit{requests} \in 2^{\mathcal{IO}}, \text{ set of output ports, initially } \emptyset \\ \textit{Phase} \in \{send, receive, sleep, mcast, mrecv\}, \\ \text{ initially send} \\ \end{array}$ 

#### Transitions at i:

input REQUEST<sub>a,i</sub> Effect: requests  $\leftarrow$  requests  $\cup \{q\}$ **input** NEWVIEW $(v)_i$ Effect:  $G \leftarrow v.set$  $A \leftarrow \emptyset$  $Rnd \leftarrow 1$  $Phase \leftarrow send$ cv := v**output** GP1SND $(m, j)_i$ Precondition: Coordinator(j)Phase = send $m = \langle i, D, R, Rnd \rangle$ Effect:  $MSG := MSG \cup \{m\}$  $Phase \leftarrow receive$ **input** GP1RCV( $\langle j, H, Q, round \rangle$ )<sub>i</sub> Effect:  $A \leftarrow A \cup \{\langle j, H, Q, \textit{round} \rangle\}$  $R \leftarrow R \cup Q$  $D \leftarrow D \cup H$ if  $G = \{j : \exists_{H',Q'} \langle j, H', Q', Rnd \rangle \in A\}$ then  $Phase \leftarrow mcast$ **output** GPMSND $(m)_i$ Precondition: Coordinator(i) $m = \langle i, D, R, Rnd \rangle$ Phase = mcastEffect:  $MSG := MSG \cup \{m\}$  $Phase \leftarrow mrecv$ 

 $m \in \mathcal{M}es$   $i, j \in \mathcal{P}$   $v \in views$   $H \in 2^{\mathcal{T}}$   $Q \in 2^{\mathcal{R}}$   $round \in \mathbf{N}$   $results \in 2^{\mathcal{R}}$  $q \in \mathcal{IO}$ 

#### Derived variables:

 $\begin{array}{l} U: \ T-D, \ \text{the set of remaining tasks}\\ Coordinator(i): \ \text{Boolean; if } i=\max(j:j\in G),\\ \text{then } True \ \text{else } False\\ \text{History variable } cv_i\in views \ (i\in \mathcal{P}),\\ \text{initially } \forall i, \ cv_i:=v_0.\\ \text{History variable } \operatorname{MSG}_i\in 2^{\mathcal{M}es} \ (i\in \mathcal{P}),\\ \text{initially } \forall i, \ \operatorname{MSG}_i:=\emptyset. \end{array}$ 

**input** GPMRCV $(\langle j, H, Q, round \rangle)_i$ Effect:  $D \leftarrow D \cup H$  $R \leftarrow R \cup Q$ if D = T then  $Phase \leftarrow sleep$ else if rank(i, G) < |U| then let u be such that rank(u, U) = rank(i, G) $R \leftarrow R \cup \{Result(u)\}$  $D \leftarrow D \cup \{u\}$ fi  $Rnd \leftarrow Rnd + 1$  $Phase \leftarrow send$ fi **output** REPORT(*results*)<sub>*a*,*i*</sub> Precondition:

T = D  $q \in requests$  results = REffect:  $requests \leftarrow requests - \{q\}$ 

Figure 2: Algorithm AX.

Next we show that if some processor does not know the result of some task, this is because it does not know that this task has been performed (Theorem 4.5 below). We show this using the history variables  $MSG_i$   $(i \in \mathcal{P})$ , which we define now.

**Definition 4.2** We define  $MSG_i$  to be a history variable that keeps on track all the messages sent by some processor  $i \in \mathcal{P}$  in all GP1SND(message, destination)<sub>i</sub> and GPMSND(message)<sub>i</sub> events of an execution of algorithm AX. Formally: In the effects of the GP1SND(m, j)<sub>i</sub> and GPMSND(m)<sub>i</sub> actions we include the assignment  $MSG_i := MSG_i \cup \{m\}$ . Initially, in state  $s_0$ ,  $\forall i$ ,  $MSG_i = \emptyset$ .

**Definition 4.3** Define  $MSG = \bigcup_{i \in P} MSG_i$ .

**Lemma 4.4** If m is a message received by processor  $i \in \mathcal{P}$  in a  $\text{GP1RCV}(m)_i$  or  $\text{GPMRCV}(m)_i$  event of an execution of algorithm AX, then  $m \in \mathcal{MSG}$ .

**Proof:** Property 3 of the GCS (Section 2.1) requires that for every receive event there exists a preceding send event of the same message (the GCS does not generate messages). Hence, m must have been sent by some processor  $q \in \mathcal{P}$  (possible q = i) in some earlier event of the execution. Messages can be sent only in GP1SND $(m,i)_q$  or GPMSND $(m)_q$  events. By definition,  $m \in MSG_q$ .

**Theorem 4.5 (Safety 2)** For all states of any execution of Algorithm AX:

- (a)  $\forall t \in T, \forall i \in \mathcal{P} : result(t) \notin R_i \Rightarrow t \notin D_i, and$
- (b)  $\forall t \in T, \forall \langle i, D', R', Rnd \rangle \in \mathcal{MSG} : result(t) \notin R' \Rightarrow t \notin D'.$

**Proof:** Let  $\alpha$  be an execution of AX and  $\alpha^k$  be the prefix of  $\alpha$  up to the  $k^{th}$  state, i.e.,  $\alpha^k = s_0, e_1, s_1, e_2, \ldots, s_k$ . The proof is done by induction on k.

Basis: k = 0. In  $s_0$ ,  $\forall i \in \mathcal{P}, D_i = \emptyset, R_i = \emptyset$  and  $\mathcal{MSG} = \emptyset$ . Thus, the basis case holds. Inductive hypothesis: For a state  $s_n$  such that  $n \leq k$ ,  $\forall t \in T$ ,  $\forall i \in \mathcal{P} : result(t) \notin R_i \Rightarrow t \notin D_i$ , and  $\forall t \in T, \forall \langle i, D', R', Rnd \rangle \in \mathcal{MSG} : result(t) \notin R' \Rightarrow t \notin D'$ .

Inductive step: n = k + 1. Consider the following seven types of actions leading to the state  $s_{k+1}$ :

- 1.  $e_{k+1} = \text{NEWVIEW}(v')_i$ : The effect of this action does not affect the invariant. By the inductive hypothesis, in state  $s_{k+1}$ , the invariant holds.
- e<sub>k+1</sub> = GP1SND(m, j)<sub>i</sub>: Clearly, the effect of this action does not affect part (a) of the invariant but it affects part (b). Since m = ⟨i, D<sub>i</sub>, R<sub>i</sub>, Rnd⟩, by the inductive hypothesis part (a), the assignment m ∈ MSG reestablishes part (b) of the invariant. Thus, in state s<sub>k+1</sub>, the invariant is reestablished.

- 3. e<sub>k+1</sub> = GP1RCV(⟨j, H, Q, round⟩)<sub>i</sub>: Processor i updates R<sub>i</sub> and D<sub>i</sub> according to Q and H respectively. The action is atomic, i.e., if R<sub>i</sub> is updated, then D<sub>i</sub> must be also updated. By Lemma 4.4, ⟨j, H, Q, round⟩ ∈ MSG. Thus, by the inductive hypothesis part (b), ∀t ∈ T : result(t) ∉ H ⇒ t ∉ Q. From the fact that D<sub>i</sub> and R<sub>i</sub> are updated according to H and Q respectively and by the inductive hypothesis part (a), in state s<sub>k+1</sub>, the invariant is reestablished.
- 4. e<sub>k+1</sub> = GPMSND(m)<sub>i</sub>: Clearly, the effect of this action does not affect part (a) of the invariant but it affects part (b). Since m = ⟨i, D<sub>i</sub>, R<sub>i</sub>, Rnd⟩, by the inductive hypothesis part (a), the assignment m ∈ MSG reestablishes part (b) of the invariant. Thus, in state s<sub>k+1</sub>, the invariant is reestablished.
- 5. e<sub>k+1</sub> = GPMRCV(⟨j, H, Q, round⟩)<sub>i</sub>: By Lemma 4.4, ⟨j, H, Q, round⟩ ∈ MSG. By the inductive hypothesis part (b), ∀t ∈ T : result(t) ∉ H ⇒ t ∉ Q. Processor i updates R<sub>i</sub> and D<sub>i</sub> according to Q and H respectively. Since H and Q have the required property, by the inductive hypothesis part (a), the assignments to D<sub>i</sub> and R<sub>i</sub> reestablish the invariant.

In the case where  $D_i \neq T$ , processor *i* performs a task according to the load balancing rule. Let  $u \in T$  be this task. Because of the action atomicity, when processor *i* updates  $R_i$  with result(u), it must also update  $D_i$  with *u*. Hence, in state  $s_{k+1}$ , the invariant is reestablished.

- 6.  $e_{k+1} = \text{REQUEST}_{q,i}$ : The effect of this action does not affect the invariant.
- 7.  $e_{k+1} = \text{REPORT}(results)_{q,i}$ : The effect of this action does not affect the invariant.

This completes the proof

## 5 Analysis of Algorithm AX

We express the work complexity of algorithm AX in the model  $\mathcal{F}_{FM}$  as  $W_{\mathcal{F}_{FM}}(N, P, r) = W_{\mathcal{F}_{FM}}(N, P, f + m)$ . The message complexity is expressed as  $M_{\mathcal{F}_{FM}}(N, P, r) = M_{\mathcal{F}_{FM}}(N, P, f + m)$ . Our analysis focuses on assessing the impact of the fragmentation number f and the merge number m on the work and message complexity, and in the rest of this section for clarity we let  $\mathcal{W}_{f,m}$  stand for  $W_{\mathcal{F}_{FM}}(N, P, f + m)$ .

### 5.1 Work Complexity

In this section we show the following result:

**Theorem 5.1**  $\mathcal{W}_{f,m} \leq \min\{ N \cdot f + N, N \cdot P \}$ 

Observe that  $\mathcal{W}_{f,m}$  does not depend on m (this of course does not imply that for any given execution, the work does not depend on merges). This observation substantiates the intuition that merges lead to a more efficient computation. We begin by providing definitions and proving several lemmas that lead to the above result.

**Definition 5.2** Let  $\alpha^{\mu}$  be any execution of algorithm AX in which all the processors learn the results of all tasks and that includes a merge of groups  $g_1, \ldots, g_k$  into the group  $\mu$ , where the processors in  $\mu$  undergo no further view changes.

We define  $\bar{\alpha}^{\mu}$  to be the execution we derive by removing the merge from  $\alpha^{\mu}$  as follows:

(1) We remove all states and events that correspond to the merge of groups  $g_1, \ldots, g_k$  into the group  $\mu$  and all states and events for processors within  $\mu$ .

(2) We add the appropriate states and events such that the processors in groups  $g_1, \ldots, g_k$  undergo no further view changes and perform any remaining tasks.

**Definition 5.3** Let  $\alpha^{\varphi}$  be any execution of algorithm AX in which all the processors learn the results of all tasks and that includes a fragmentation of the group  $\varphi$  to the groups  $g_1, \ldots, g_k$  where the processors in these groups undergo no further view changes.

We define  $\bar{\alpha}^{\varphi}$  to be the execution we derive by removing the fragmentation from  $\alpha^{\varphi}$  as follows:

(1) We remove all states and events that correspond to the fragmentation of the group  $\varphi$  to the groups  $g_1, \ldots, g_k$  and all states and events of the processors within the groups  $g_1, \ldots, g_k$ .

(2) We add the appropriate states and events such that the processors in the group  $\varphi$  undergo no further view changes and perform any remaining tasks.

**Note:** In Definitions 5.2 and 5.3, we claim that we can remove states and events from an execution and add some other states and events to it. This is possible because if the processors in a single view installed that view and there are no further view changes, then the algorithm will continue making computation progress. So, if we remove all states and events corresponding to a view change, then the algorithm can always proceed as if this view change never occurred.

**Lemma 5.4** In algorithm AX, for any view v, including the initial view, if the group is not subject to any regroupings, then the work required to complete all tasks in the view is no more than  $N - \max_{i \in v.set} \{|D_i|\}$ , where  $D_i$  is the value of the state variable D of processor i at the start of its local round 1 in view v. **Proof:** In the first round, all the processors send messages to the coordinator containing  $D_i$ . The coordinator computes  $\bigcup_{i \in v.set} \{D_i\}$  and broadcasts this result to the group members. Since the group is not subject to any regroupings, the number of tasks, t, that the processors need to perform is:  $t = N - |\bigcup_{i \in v.set} \{D_i\}|$ . In each round of the computation, by the load balancing rule, the members of the group perform distinct tasks and no task is performed more than once. Therefore, t is the work performed in this group. On the other hand,  $\max_{i \in v.set} \{|D_i|\} \leq |\bigcup_{i \in v.set} \{D_i\}|$ , thus,  $t \leq N - \max_{i \in v.set} \{|D_i|\}$ .

In the following lemma, groups  $\mu, g_1, \ldots, g_k$  are defined as in Definition 5.2.

**Lemma 5.5** Let  $\alpha^{\mu}$  be an execution of Algorithm AX as in Definition 5.2. Let  $W_1$  be the work performed by the algorithm in the execution  $\alpha^{\mu}$ . Let  $W_2$  be the worked performed by Algorithm AX in the execution  $\bar{\alpha}^{\mu}$ . Then  $W_1 \leq W_2$ .

**Proof:** Let W' be the work performed by all processors in  $\mathcal{P} - \bigcup_{1 \leq i \leq k} (g_i.set) - \mu.set$  in the execution  $\alpha^{\mu}$ . Observe that the work performed by all processors in  $\mathcal{P} - \bigcup_{1 \leq i \leq k} (g_i.set)$  in the execution  $\bar{\alpha}^{\mu}$  is equal to W'.

The work that is performed by processor j in  $g_{i.set}$  prior to the NEWVIEW $(\mu)_{j}$  event in  $\alpha^{\mu}$ , is the same in both executions. Call this work  $W_{i,j}$ . Define  $W'' = \sum_{i=1}^{k} \sum_{j \in g_{i.set}} W_{i,j}$ .

Define W = W' + W''. Thus, W is the same in both executions,  $\alpha^{\mu}$  and  $\bar{\alpha}^{\mu}$ .

Define  $W_{\mu}$  to be the work performed by all processors in  $\mu$ .set in execution  $\alpha^{\mu}$ .

For each processor j in  $g_{i.set}$ , let  $D_{j}$  be the value of the state variable D just prior to the NEWVIEW $(\mu)_{j}$  event in  $\alpha^{\mu}$ . For each  $g_{i}$ , define:  $d_{i} = |\bigcup_{j \in g_{i.set}} D_{j}|$ . Thus there are at least  $N - d_{i}$  tasks that remain to be done in each  $g_{i}$ .

In execution  $\bar{\alpha}^{\mu}$ , the processors in each group  $g_i$  proceed and complete these remaining tasks. This requires work at least  $N - d_i$ . Define this work as  $W_{g_i}$ . Thus,  $W_{g_i} \ge (N - d_i)$ .

In execution  $\alpha^{\mu}$ , groups  $g_1, \ldots, g_k$  merge into group  $\mu$ . The number of tasks that need to be performed by the members of  $\mu$  is at most  $N - d_j$ , where  $d_j = \max_i \{d_i\}$  for some j. By Lemma 5.4,  $W_{\mu} \leq N - d_j$ . Observe that:

$$W_1 = W + W_\mu \le W + N - d_j \le W + \sum_{i=1}^k (N - d_i) \le W + \sum_{i=1}^k W_{g_i} = W_2.$$

In the following lemma, groups  $\varphi, g_1, \ldots, g_k$  are defined as in Definition 5.3.

**Lemma 5.6** Let  $\alpha^{\varphi}$  be an execution of Algorithm AX as in Definition 5.3. Let  $W_1$  be the work performed by the algorithm in the execution  $\alpha^{\varphi}$ . Let  $W_2$  be the worked performed by Algorithm AX in the execution  $\bar{\alpha}^{\varphi}$ . Then  $W_1 \leq W_2 + W_3$ , where  $W_3$  is the work performed by all processors in  $\bigcup_{1 \leq i \leq k} (g_i.set)$  in the execution  $\alpha^{\varphi}$ .

**Proof:** Let W' be the work performed by all processors in  $\mathcal{P} - \bigcup_{1 \leq i \leq k} (g_i.set) - \varphi.set$  in the execution  $\alpha^{\varphi}$ . Observe that the work performed by all processors in  $\mathcal{P} - \varphi.set$  in the execution  $\bar{\alpha}^{\varphi}$  is equal to W'.

The work that is performed by processor j in  $\varphi$ .set prior to the NEWVIEW $(g_i)_j$  event in  $\alpha^{\varphi}$ , is the same in both executions. Call this work  $W_{\varphi,j}$ . Define  $W'' = \sum_{j \in \varphi.set} W_{\varphi,j}$ .

Define W = W' + W''. Thus, W is the same in both executions,  $\alpha^{\varphi}$  and  $\bar{\alpha}^{\varphi}$ .

Define  $W_{\varphi}$  to be the work performed by all processors in  $\varphi$ . set in execution  $\bar{\alpha}^{\varphi}$ . Let  $W''' = W_{\varphi} - W''$ . Observe that:  $W_1 = W + W_3 \leq W + W_3 + W''' = W_2 + W_3$ 

Lemma 5.7  $W_{f,m} \leq N \cdot P$ 

**Proof:** By the construction of algorithm AX, when processors are not able to exchange information about task execution due to regroupings, in the worst case, each processor has to perform all Ntasks by itself. Since  $P \leq N$ , the work is:  $\mathcal{W}_{f,m} \leq N \cdot P$ .

Lemma 5.8  $\mathcal{W}_{f,m} \leq N \cdot f + N$ .

**Proof:** By induction on the number of views, denoted by r, occurring in an execution. For a specific execution  $\alpha_r$  with r views, let  $f_r$  be the fragmentation-number and  $m_r$  the merge-number.

Basis: r = 0. Since  $f_r$  and  $m_r$  must also be 0, the basis follows from Lemma 5.4.

Inductive hypothesis: Assume that for all  $r \leq k$ ,  $\mathcal{W}_{f_r,m_r} \leq N \cdot f_r + N$ .

Inductive step: Need to show that for r = k + 1,  $\mathcal{W}_{f_{k+1}, m_{k+1}} \leq N \cdot f_{k+1} + N$ .

Consider a specific execution  $\alpha_{k+1}$  with r = k + 1. Let  $\Gamma_{\alpha_{k+1}}$  be the view-graph induced by this execution. The view-graph has at least one vertex such that all of its descendants are sinks (Fact 3.12). Let  $\nu$  be such a vertex. We consider two cases.

(1)  $\nu$  has a descendant  $\mu$  that corresponds to a merge in the execution. Therefore all ancestors of  $\mu$  in  $\Gamma_{\alpha_{k+1}}$  have outdegree 1. Since  $\mu$  is a sink vertex, the group that corresponds to  $\mu$  performs all the remaining (if any) tasks and does not perform any additional work.

Let  $\alpha_k = \bar{\alpha}_{k+1}^{\mu}$  (per Definition 5.2) be an execution in which this merge does not occur. In execution  $\alpha_k$ , the number of views is k. Also,  $f_{k+1} = f_k$  and  $m_{k+1} = m_k + 1$ . By inductive hypothesis,  $\mathcal{W}_{f_k,m_k} \leq N \cdot f_k + N$ . By Lemma 5.5, the work performed in execution  $\alpha_{k+1}$ , is no worse than the work performed in execution  $\alpha_k$ . The total work complexity is:

 $\mathcal{W}_{f_{k+1},m_{k+1}} \leq \mathcal{W}_{f_k,m_k} \leq N \cdot f_k + N = N \cdot f_{k+1} + N.$ 

(2)  $\nu$  has no descendants that correspond to a merge in the execution. Therefore, the group that corresponds to  $\nu$  must fragment, say into q groups. These groups correspond to sink vertices in  $\Gamma_{\alpha_{k+1}}$ , thus they perform all the remaining (if any) tasks and do not perform any additional work.

Let  $\alpha_{k+1-q} = \bar{\alpha}_{k+1}^{\nu}$  (per Definition 5.3) be an execution in which the fragmentation does not occur. In execution  $\alpha_{k+1-q}$ , the number of views is  $k+1-q \leq k$ . Also,  $f_{k+1-q} = f_{k+1} - q$  and  $m_{k+1-q} = m_{k+1}$ . By inductive hypothesis,  $\mathcal{W}_{f_{k+1-q},m_{k+1-q}} \leq N \cdot f_{k+1-q} + N$ . From Lemma 5.4, the work performed in each new group caused by the fragmentation is no more than N. Let  $W_{\sigma}$  be the total work performed in all q groups. Thus,  $W_{\sigma} \leq qN$ . By Lemma 5.6, the work performed in execution  $\alpha_{k+1}$ , is no worse than the work performed in execution  $\alpha_{k+1-q}$  and the work performed in all q groups. The total work complexity is:

$$\mathcal{W}_{f_{k+1},m_{k+1}} \leq \mathcal{W}_{f_{k+1-q},m_{k+1-q}} + W_{\sigma} \leq N \cdot f_{k+1-q} + N + W_{\sigma} = N \cdot (f_{k+1} - q) + N + W_{\sigma}$$
$$\leq N \cdot (f_{k+1} - q) + N + qN = Nf_{k+1} - qN + N + qN = N \cdot f_{k+1} + N.$$

The main result in Theorem 5.1 follows directly from Lemma 5.7 and Lemma 5.8.

The work complexity result is tight for a broad range of regroupings.

**Theorem 5.9** Given any execution  $\alpha$  in the  $\mathcal{F}_{FM}$  model, there is a pattern of fragmentations that causes algorithm AX to perform  $\Omega(N \cdot f_{\alpha})$  work.

**Proof:** Consider the following pattern. At the beginning of the execution, all P processors belong in one group. Before the processors communicate, a fragmentation occurs that splits the group into some number of distinct groups (more than one). Before the processors communicate or perform any tasks, arbitrary numbers of fragmentations occurs. This continues until we reach a point in the execution, where no more failures occur. Define the number of the existing groups after this point to be g. The processors in these g groups must perform all N tasks. Let W be the total work of the execution. Thus,  $W \ge N \cdot g$ . From Lemma A.1 (page 28) we know that  $g > \frac{f_{\alpha}}{2}$ . Thus,  $W > N \cdot \frac{f_{\alpha}}{2} = \Omega(N \cdot f_{\alpha})$ .

#### 5.2 Message Complexity

In this section we show the following result:

**Theorem 5.10**  $\mathcal{M}_{f,m} < 4(N \cdot f + N + P \cdot m)$ 

We start by showing several lemmas that lead to the message complexity result.

**Lemma 5.11** For algorithm AX, in any view v, including the initial view, if the group is not subject to any regroupings, and for each processor  $i \in v.set$ ,  $D_i$  is the value of the state variable D at the start of its local round 1 in view v, then the number of messages M that are sent until all tasks are completed is  $2(N-d) \leq M < 2(p+N-d)$  where p = |v.set|, and  $d = |\bigcup_{i \in v.set} D_i|$ .

**Proof:** By the load balancing rule, the algorithm needs  $\lceil \frac{N-d}{p} \rceil$  rounds to complete all tasks. In each round each processor sends one message to the coordinator and the coordinator responds with a single message to each processor. Thus  $M = 2p \cdot (\lceil \frac{N-d}{p} \rceil)$ . Using the properties of the *ceiling*, we get:  $2(N-d) \leq M < 2(p+N-d)$ .

In the following lemma, groups  $\mu, g_1, \ldots, g_k$  are defined as in Definition 5.2.

**Lemma 5.12** Let  $\alpha^{\mu}$  be an execution of Algorithm AX as in Definition 5.2. Let  $M_1$  be the message cost of the algorithm in the execution  $\alpha^{\mu}$ . Let  $M_2$  be the message cost of Algorithm AX in the execution  $\bar{\alpha}^{\mu}$ . Then  $M_1 < M_2 + 2P$ .

**Proof:** Let M' be the number of messages sent by all processors in  $\mathcal{P} - \bigcup_{1 \leq i \leq k} (g_i.set) - \mu.set$  in the execution  $\alpha^{\mu}$ . Observe that the number of messages sent by all processors in  $\mathcal{P} - \bigcup_{1 \leq i \leq k} (g_i.set)$  in the execution  $\bar{\alpha}^{\mu}$  is equal to M'.

The number of messages sent by any processor j in  $g_i.set$  prior to the NEWVIEW $(\mu)_j$  event in  $\alpha^{\mu}$ , is the same in both executions. Call this message cost  $M_{i,j}$ . Define  $M'' = \sum_{i=1}^k \sum_{j \in g_i.set} M_{i,j}$ . Define M = M' + M''. Thus, M is the same in both executions,  $\alpha^{\mu}$  and  $\bar{\alpha}^{\mu}$ .

Define  $M_{\mu}$  to be the number of messages sent by all processors in  $\mu$ .set in execution  $\alpha^{\mu}$ .

For each processor j in  $g_{i.set}$ , let  $D_{j}$  be the value of the state variable D just prior to the NEWVIEW $(\mu)_{j}$  event in  $\alpha^{\mu}$ . For each  $g_{i}$ , define:  $d_{i} = |\bigcup_{j \in g_{i.set}} D_{j}|$ . Thus there are at least  $N - d_{i}$  tasks that remain to be done in each  $g_{i}$ .

In execution  $\bar{\alpha}^{\mu}$ , the processors in each group  $g_i$  proceed and complete these remaining tasks. Let  $M_{g_i}$  be the number of messages sent by all processors in  $g_i$ .set in order to complete the remaining tasks. By Lemma 5.11,  $M_{g_i} \geq 2(N - d_i)$ .

In execution  $\alpha^{\mu}$ , groups  $g_1, \ldots, g_k$  merge into group  $\mu$ . The number of tasks that need to be performed by the members of  $\mu$  is at most  $N - d_j$ , where  $d_j = \max_i \{d_i\}$  for some j. By Lemma 5.11,  $M_{\mu} < 2(p + N - d_j)$ , where  $p = |\mu.set|$ . Observe that:

$$M_1 = M + M_\mu < M + 2(p + N - d_j) \le M + 2p + 2\sum_{i=1}^k (N - d_i) \le M + 2p + \sum_{i=1}^k M_{g_i}$$
  
=  $M_2 + 2p \le M_2 + 2P$ 

In the following lemma, groups  $\varphi, g_1, \ldots, g_k$  are defined as in Definition 5.3.

**Lemma 5.13** Let  $\alpha^{\varphi}$  be an execution of Algorithm AX as in Definition 5.3. Let  $M_1$  be the message cost of the algorithm in the execution  $\alpha^{\varphi}$ . Let  $M_2$  be the message cost of Algorithm AX in the

execution  $\bar{\alpha}^{\varphi}$ . Then  $M_1 \leq M_2 + M_3$ , where  $M_3$  is the number of messages sent by all processors in  $\bigcup_{1 \leq i \leq k} (g_i.set)$  in the execution  $\alpha^{\varphi}$ .

**Proof:** Let M' be the number of messages sent by all processors in  $\mathcal{P} - \bigcup_{1 \leq i \leq k} (g_i.set) - \varphi.set$  in the execution  $\alpha^{\varphi}$ . Observe that the number of messages sent by all processors in  $\mathcal{P} - \varphi.set$  in the execution  $\bar{\alpha}^{\varphi}$  is equal to M'.

The number of messages sent by processor j in  $\varphi$ .set prior to the NEWVIEW $(g_i)_j$  event in  $\alpha^{\varphi}$ , is the same in both executions. Call this message cost  $M_{\varphi,j}$ . Define  $M'' = \sum_{j \in \varphi.set} M_{\varphi,j}$ .

Define M = M' + M''. Thus, M is the same in both executions,  $\alpha^{\varphi}$  and  $\bar{\alpha}^{\varphi}$ .

Define  $M_{\varphi}$  to be the number of messages sent by all processors in  $\varphi$ .set in execution  $\bar{\alpha}^{\varphi}$ . Let  $M''' = M_{\varphi} - M''$ . Observe that:  $M_1 = M + M_3 \leq M + M_3 + M''' = M_2 + M_3$ 

The proof of Theorem 5.10 is done by induction similarily to the proof of Lemma 5.8:

**Proof:** By induction on the number of views, denoted by r, occurring in any execution. For a specific execution  $\alpha_r$  with r views, let  $f_r$  be the fragmentation number and  $m_r$  be the merge-number.

Basis: r = 0. Since  $f_r$  and  $m_r$  must also be 0, the basis follows from Lemma 5.11.

Inductive hypothesis: Assume that for all  $r \leq k$ ,  $\mathcal{M}_{f_r,m_r} < 4(N \cdot f_r + N + P \cdot m_r)$ .

Inductive step: Need to show that for r = k + 1,  $\mathcal{M}_{f_{k+1}, m_{k+1}} < 4(N \cdot f_{k+1} + N + P \cdot m_{k+1})$ .

Consider a specific execution  $\alpha_{k+1}$  with r = k + 1. Let  $\Gamma_{\alpha_{k+1}}$  be the view-graph induced by this execution. The view-graph has at least one vertex such that all of its descendants are sinks (Fact 3.12). Let  $\nu$  be such a vertex. We consider two cases.

(1)  $\nu$  has a descendant  $\mu$  that corresponds to a merge in the execution. Therefore all ancestors of  $\mu$  in  $\Gamma_{\alpha_{k+1}}$  have outdegree 1. Since  $\mu$  is a sink vertex, the group that corresponds to  $\mu$  performs all the remaining (if any) tasks and no further messages are sent.

Let  $\alpha_k = \bar{\alpha}_{k+1}^{\mu}$  (per Definition 5.2) be an execution in which this merge does not occur. In execution  $\alpha_k$ , the number of new views is k. Also,  $f_{k+1} = f_k$  and  $m_{k+1} = m_k + 1$ . By inductive hypothesis,  $\mathcal{M}_{f_k,m_k} < 4(N \cdot f_k + N + P \cdot m_k)$ . The total message complexity, using Lemma 5.12 is:  $\mathcal{M}_{f_{k+1},m_{k+1}} < \mathcal{M}_{f_k,m_k} + 2P < 4(N \cdot f_k + N + P \cdot m_k) + 2P = 4(N \cdot f_{k+1} + N + P \cdot m_{k+1} - P) + 2P$  $= 4Nf_{k+1} + 4N + 4Pm_{k+1} - 4P + 2P \leq 4(N \cdot f_{k+1} + N + P \cdot m_{k+1}).$ 

(2)  $\nu$  has no descendants that correspond to a merge in the execution. Therefore, the group that corresponds to  $\nu$  must fragment, say into q groups. These groups correspond to sink vertices in  $\Gamma \alpha_{k+1}$ , thus they perform all of the remaining (if any) tasks and do not send any additional messages.

Let  $\alpha_{k+1-q} = \bar{\alpha}_{k+1}^{\nu}$  (per Definition 5.3) be an execution in which the fragmentation does not

occur. In the execution  $\alpha_{k+1-q}$ , the number of new views is  $k+1-q \leq k$ . Also,  $f_{k+1-q} = f_{k+1}-q$ and  $m_{k+1-q} = m_{k+1}$ . By inductive hypothesis,  $\mathcal{M}_{f_{k+1-q},m_{k+1-q}} < 4(N \cdot f_{k+1-q} + N + P \cdot m_{k+1-q})$ . From Lemma 5.11, the message cost in each new group caused by a fragmentation is no more than 4N. Let  $M_{\sigma}$  be the total number of messages sent in all q groups. Thus,  $M_{\sigma} \leq 4qN$ . By Lemma 5.13, the number of messages sent in execution  $\alpha_{k+1}$ , is less than the number of messages sent in execution  $\alpha_{k+1-q}$  and the number of messages sent in all q groups. The total message complexity is:  $\mathcal{M}_{f_{k+1},m_{k+1}} \leq \mathcal{M}_{f_{k+1-q},m_{k+1-q}} + M_{\sigma} < 4(N \cdot f_{k+1-q} + N + P \cdot m_{k+1-q}) + M_{\sigma}$  $= 4(N \cdot f_{k+1} - qN + N + P \cdot m_{k+1}) + M_{\sigma} < 4N f_{k+1} - 4qN + 4N + 4Pm_{k+1} + 4qN$ 

$$= 4(N \cdot f_{k+1} + N + P \cdot m_{k+1}).$$

The message complexity result is tight for a broad range of regroupings.

**Theorem 5.14** Given any execution  $\alpha$  in the  $\mathcal{F}_{FM}$  model, there is a pattern of fragmentations that causes algorithm AX to send  $\Omega(N \cdot f_{\alpha})$  messages.

**Proof:** Consider the following pattern. At the beginning of the execution, all P processors belong in one group. Before the processors communicate, a fragmentation occurs that splits the group into some number of distinct groups (more than one). Before the processors communicate or perform any tasks, arbitrary numbers of fragmentations occurs. This continues until we reach a point in the execution, where no more failures occur. Define the number of the existing groups after this point to be g. The processors in these g groups must perform all N tasks. By Lemma 5.11, the message cost of each of the g groups is at least 2N. Let M be the total work of the execution. Thus,  $M \ge 2N \cdot g$ . From Lemma A.1 (page 28) we know that  $g > \frac{f_{\alpha}}{2}$ . Thus,  $M > 2N \cdot \frac{f_{\alpha}}{2} = \Omega(N \cdot f_{\alpha})$ .

Observe that the same scenario provides a lower bound both for work and message cost.

#### 5.3 Analysis for the Fragmentation Adversary

We express the work complexity of algorithm AX in the model  $\mathcal{F}_F$  as  $W_{\mathcal{F}_F}(N, P, r) = \mathcal{W}_f$  and the message complexity as  $M_{\mathcal{F}_F}(N, P, r) = \mathcal{M}_f$  (note that r = f for  $\mathcal{F}_F$ ). The following corollary is derived from Theorems 5.1 and 5.10.

## **Corollary 5.15** $\mathcal{W}_f \leq \min\{N \cdot f + N, N \cdot P\}$ and $\mathcal{M}_f < 4(N \cdot f + N)$

With Theorems 5.9 and 5.14 the results of Corollary 5.15 are tight.

In the failure model of [8] a group is not allowed to "fragment" into a single group with the same membership. Such fragmentation is allowed by our definition of  $\mathcal{F}_F$ . In order to compare our results with the results of [8], we define a more restricted adversary  $\mathcal{F}'_F$  that requires that any group may only fragment into 2 or more other groups. Clearly  $\mathcal{F}'_F \subseteq \mathcal{F}_F$ , and from Corollary 5.15 we have the following.

# $\textbf{Corollary 5.16} \hspace{0.2cm} W_{\mathcal{F}'_{F}}(N,P,f) \hspace{-0.1cm}=\hspace{-0.1cm} O(N \cdot f + N) \hspace{0.2cm} and \hspace{0.2cm} M_{\mathcal{F}'_{F}}(N,P,f) \hspace{-0.1cm}=\hspace{-0.1cm} O(N \cdot f + N)$

In the rest of this section we deal with the model  $\mathcal{F}'_F$ . Our definition of the fragmentation-number f is slightly different from the definition of fragmentation failures f' in [8]. When a group fragments into k groups, f is defined to be equal to k, but f' is defined to be equal to k-1. The next Lemma relates f and f'.

**Lemma 5.17** If f is the fragmentation-number and f' the number of fragmentation failures as defined in [8], then f' < f < 2f'.

**Proof:** Assume that k fragmentations occur. Let the number of the newviews in the  $i^{th}$  fragmentation be  $f_i$ . By the definition of  $f'_i$ ,  $f'_i = f_i - 1$ . Thus,  $f'_i + 1 = f_i$  which implies that  $f_i < f'_i + f'_i = 2f'_i$ . But  $f' = \sum_{i=1}^k f'_i$  and  $f = \sum_{i=1}^k f_i$ . Thus, f < 2f'. Now observe that,  $f' = \sum_{i=1}^k f'_i = \sum_{i=1}^k (f_i - 1) = \sum_{i=1}^k f_i - k = f - k$ . Thus f > f'

In [8] the work is counted in terms of the rounds executed by the processors. In our analysis we count only the number of task executions (including redundancies). However in our algorithm, for as long as any tasks remain undone in a given group, the processors perform the tasks in rounds, except for the last round. Therefore the difference in work complexity for these two algorithms is at most  $f \cdot N$ . Thus the different definitions of f, f' and work are subsumed in the big-oh analysis, and without substantial variation in the constants. On the other hand, the message complexity of our algorithm, as shown in Corollary 5.16, is substantially better than the at least quadratic message complexity of the algorithm from [8].

### 6 Discussion

We have considered the problem of performing a set of N tasks on a set of P cooperating messagepassing processors, where the processors must perform and learn the results of the tasks efficiently, subject to dynamically changing group memberships. To analyze our algorithm we introduce viewgraphs – digraphs that we use to represent and analyze changes of processors' views in executions. Our complexity analysis shows that group communication services can be used efficiently to solve distributed problems that are similar to our OMNI-DO problem. We believe that our view-graph approach can be generalized and used to study other dynamic group reconfiguration patterns. Ongoing work is pursuing this research direction.

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## A Appendix

We prove a Lemma that we use in the proofs of Theorems 5.9 and 5.14.

**Lemma A.1** In an execution  $\alpha$  of algorithm AX in the fragmentation adversary model where groups fragment to two or more groups, for N tasks and P processors it holds that  $f_{\alpha} < 2g$ , where g is the number of groups at a point in the execution where no more regroupings occur.

**Proof:** By induction on the number of new views, denoted r, occurring in any execution. For a specific execution  $\alpha_r$  with r views, let  $f_r$  be the fragmentation-number.

Basis: For r = 0,  $f_0 = 0 < 2$ .

Inductive hypothesis: Assume that for all  $r \leq k$ ,  $f_r < 2g$ .

Inductive step: Need to show that for r = k + 1,  $f_{k+1} < 2g$ .

Consider a specific execution  $\alpha_{k+1}$  with r = k + 1. Let  $\Gamma_{\alpha_{k+1}}$  be the view-graph induced by this execution. The view-graph has at least one vertex such that all of its descendants are sinks (Fact 3.12). Let  $\nu$  be such vertex. Since there are no merges, the group that corresponds to  $\nu$ , fragments into q groups. These groups correspond to sink vertices in  $\Gamma_{\alpha_{k+1}}$ , thus they perform all of the remaining (if any) tasks.

Let  $\alpha_{k+1-q} = \bar{\alpha}_{k+1}^{\nu}$  (per Definition 5.3) be an execution in which the fragmentation does not occur. In the execution  $\alpha_{k+1-q}$ , the number of new views is  $k+1-q \leq k$  and  $f_{k+1-q} = f_{k+1}-q$ . By inductive hypothesis,  $f_{k+1-q} < 2(g-q)$ .

Thus,  $f_{k+1} = f_{k+1-q} + q < 2(g-q) + q = 2g - q < 2g.$