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An energy-efficient process clustering assignment algorithm for distributed system



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ABSTRACT

This paper proposes a distributed assignment algorithm for selecting the optimal energy consumption during process execution, idling, and transmission in a distributed system. Selection criteria are based on identifying candidate processing units that are suitable for minimizing idle energy in task scheduling. The proposed algorithm tries to mimic as close to real situation as possible by assuming that each processing unit has multiple capabilities to execute different tasks with different characteristics. Task scheduling can be flexibly carried out to attain optimal energy consumption without any restrictions as those of comparative algorithms. Thus, the energy required by each processing unit varies considerably depending on the schedule. Experimental results show that the proposed algorithm yields the lowest idle, total energy consumption, and satisfactory execution energy. The extraneous transmission energy is a trade-off for scheduling flexibility.

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1. Introduction

Demands for energy conscious algorithms have increased in the recent years along with the fast-paced development of computational technology [1–4]. Various combined system architectures or operators were used to solve complicated problems in science and commercial including cloud computing and computer grid systems [5–9]. These systems were built with complex algorithms which required high computing power. Thus, one mandate that the supporting algorithms had to satisfy was minimizing the complexity of the computation process.

In general, when parallel code is pending for execution by a cluster of servers in a distributed system, the tasks, data, and functions are generated by the user on the client site. Only the required components are sent to the server cluster for execution. In most cases, each assigned server will only receive allotted jobs to perform. If the jobs coming from multiple sources must undergo additional security measures, a higher transferring rate to send data across networks will be required due to more processes to be finished. Therefore, ordinary server assignment algorithms are inadequate as the time complexity posts more power requirement to achieve the optimal energy consumption.

Concern of energy-aware issue calls for energy-efficient server assignment problem to be carefully studied [10–14]. Two minimum energy consumption techniques for real-time 2-level heterogeneous grid system were proposed by Terzopoulos and Karatza [15], i.e., Dynamic Voltage Scaling (DVS) and Dynamic Power Management (DPM). The predominant aspect is the reduction of energy consumption with minimum performance degradation. In a typical distributed system, some data or functions are kept at a local scheduler and some are kept in specific places due to required security. The task scheduler

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must determine the starting time of each task and the server to which the task should be sent. To attain the optimal energy consumption, scheduling scenarios could be quite complex due to the flexibility of assigning task to proper matching execution unit. Our proposed Energy-Efficient Process Clustering Assignment (EPC) algorithm aims to handle such complex scenarios to find the optimal energy efficient assignment in a reasonable computing time. The important aspects to the development of this algorithm are matching the servers to which the tasks should be sent, determining the execution time of each specific task to be run on the designated server, and clustering tasks for execution on the same server.

The paper is organized as follows. Related works and their background theory are briefly reviewed in Section 2. Details of the problem formulation and constraints are described in Section 3. The proposed algorithm is discussed in Section 4. Sections 5 presents the experimental details and results. Some notable points are discussed and concluded in Section 6.

2. Related works and background

Many parallel or distributed task scheduling algorithms have been proposed, such as list scheduling, clustering, and task duplication heuristics. A brief summary of each approach is described below.

2.1. List scheduling heuristics (LSH)

This is perhaps the most common scheduling algorithm owing to its practicality [16–18]. Most list scheduling algorithms have two parts: task prioritization and server selection. In the first part, the algorithm assigns a score to each task according to its average computation cost. The score will then be ranked in a descending order to determine task priority. In the second part, the algorithm estimates the execution time of each task used by individual server and assigns the task to the server which has the least execution time for that task. One of the interesting algorithms is the Min-Min algorithm proposed by Li et al. in 2011 [18]. For this algorithm, the authors modified the original Min-Min algorithm which treated each task graph as independent to treating each task graph as dependent. The algorithm finds the earliest finishing time for each task in each device and then schedules each task to be computed with the earliest finishing time.

2.2. Clustering heuristics (CH)

Heuristic clustering algorithms are algorithms aiming to group tasks into m groups according to their transmission edge, to reduce transmission cost, and then to map each group to n servers [19–22]. One of the examples of heuristic clustering algorithms is the algorithm introduced by Liou and Palis in 1997 [20]. The algorithm has four steps: clustering all tasks into groups, merging clusters to a number of servers, mapping individual cluster to a server, and determining the order of task execution. It balances both the load in each cluster or server and the communication traffic among the clusters. A cluster of tasks is assigned to a server starting from the cluster which has the highest communication cost until all clusters are assigned.

2.3. Task duplication heuristics (TDH)

This algorithm pushes redundant tasks to different servers to reduce the communication cost between tasks [23,24]. Although it can reduce the computing time, it requires high energy consumption and execution cost as redundant tasks are executed on different servers. This approach is effective for systems with a high ratio of communication and execution cost. Unfortunately, its operation is very complex and the overhead to find an optimal solution is high.

2.4. Heterogeneous earliest finish time (HEFT)

The algorithm was proposed by Topcuoglu et al. in 2002 [25]. It finds the shortest time of overall finishing time of a system with insertion base. The algorithm consists of two steps, namely, task priority and server selection. The task priority calculates the priority of computation and communication for each task. The server selection determines idle time slots on each server to insert a task for execution at the earliest time.

Some predominant aspects of the algorithms can be summarized as follows. List scheduling heuristic yields the lowest computational cost, while clustering heuristic has a moderate communication cost. At any rate, HEFT gives a shorter finishing time at relatively low cost. Task duplication heuristics are strong in short finishing time but use high energy. By exploiting such strengths and weaknesses, we have derived an energy saving approach by grouping tasks in groups to be executed on the same server that requires the least energy. This in turn reduces the overall communication energy considerably. However, some tasks may carry their inherent restrictions which preclude them from being clustered or executed on the same server. Hence, the details of problem formulation and constraints of task characteristics and server competency are discussed below.

3. Problems formulation and constraints

Formulation of the proposed algorithm and its constraints are given in the following sections.

3.1. Problem formulation

Let G = (V, E) be a directed graph representing a process captured in the forms of tasks and their dependencies. Graph G is called dependent task graph. $V = \{v_1, \ldots, v_n\}$ is a set of tasks to be performed. $E = \{(v_i, v_k) | v_i, v_k \in V\}$ is a set dependency edges among the tasks. An edge (v_i, v_k) denotes that task v_i must be performed prior to task v_k . A set of tasks occurs on the client site, which may be a mobile phone or a PC. Each task can be performed either on a client site or a server site. Each server is capable of performing more than one task with different amount of energy consumption. A task is processed by a processing unit which can be either a client machine or a server machine. Thus, instead of distinguishing a client from a server, we will call them *processing unit*. Each processing unit is capable of processing some tasks. It is possible that there may be more than one processing unit that can process the same task. Let $e_a(v_i)$ be the amount of energy consumed to process task v_i by processing unit a. Assuming that the number of tasks is greater than the number of processing units, the problem studied in this paper is formulated as follows.

Given a dependent task graph G, find a method to partition tasks v_i 's $\in V$ into m groups of tasks $\{\mathbf{g}_1, ..., \mathbf{g}_m\}$ so that group \mathbf{g}_j can be performed on processing unit j in such a way that the total amount of energy consumed by all processing units is minimum. In other words, the objective is to minimize $\sum_{1 \le i \le m: v_i \in \mathbf{g}_i} e_i(v_j)$.

3.2. Constraints on energy consumption

The amount of energy $e_a(v_i)$ can be decomposed into three types, namely, execution energy, transmission energy, and idle energy. Execution energy depends primarily on server competency and task characteristics. Task v_i requires $t_a(v_i)$ amount of time to be executed on processing unit a. Let α_a be the energy per unit time consumed by processing unit a to execute a task. The execution energy to execute task v_i on processing unit a is computed by $\alpha_a t_a(v_i)$.

Transmission energy occurs when data are sent from one processing unit to other processing units according to their data dependencies. Let $d_a(v_i)$ be the amount of data of task v_i sent from current processing unit a to other processing units in the same dependency path of G. For each unit amount of data, the transmission energy consumed is set to λ_a . Therefore, the transmission energy is computed by $\lambda_a d_a(v_i)$.

The idle energy of processing unit a in this study is based on the assumption that if processing unit a does not process any task, then it is put into an idle state. Processing unit a is activated when there are some other processing units transferring the dependent data to be processed by it. Suppose the remaining jobs of task v_i are transferred to processing unit a. When processing unit a is activated, it must spends $w_a(v_i)$ unit time to receive all related data of v_i to be transferred before execution. Therefore, the idle energy must involve the energy consumed during the idle state and the energy wasted during the data transferring period. Let $\psi_a(v_i)$ be the time that processing unit a is in the idle state before being activated by task v_i . The total idle time is equal to $\psi_a(v_i) + w_a(v_i)$. We assume that each idle unit time of processing unit a consumes a0 unit of energy. Hence, the total energy is equal to a1 to a2 unit of energy.

Some processing unit may not consume any idle energy and transmission energy if it executes the last task of G which is a leaf vertex. To handle this situation, constants $\kappa_a \in \{0, 1\}$ and $\mu_a \in \{0, 1\}$ are introduced to indicate whether a task v_i executed by processing unit a is a leaf vertex or an internal vertex in G. $\kappa_a = 0$ if the number of out-degrees of v_i is equal to zero and $\mu_a = 0$ if the number of in-degrees of v_i is equal to zero. Otherwise, it is equal to 1. Hence, the total energy of G, denoted by Γ_G , is computed by this equation.

$$\Gamma_{G} = \sum_{1 \leq a \leq m; v_{j} \in \mathbf{g}_{a}} e_{a}(v_{j}) + \sum_{1 \leq a \leq m; v_{i} \in \mathbf{g}_{a}} \beta_{a} \psi_{a}(v_{i})$$

$$= \sum_{1 \leq a \leq m; v_{j} \in \mathbf{g}_{a}} \alpha_{a} t_{a}(v_{j}) + \kappa_{a} \lambda_{a} d_{a}(v_{j}) + \mu_{a} \beta_{a} w_{a}(v_{j})$$

$$+ \sum_{1 \leq a \leq m; v_{i} \in \mathbf{g}_{a}} \beta_{a} \psi_{a}(v_{i})$$
(2)

4. Energy-Efficient Process Clustering Assignment (EPC) algorithm

We assume that, for any processing unit a, the values of $t_a(v_i)$, α_a , $d_a(v_i)$, λ_a , $w_a(v_i)$, and β_a can be estimated prior to the assignment of tasks to processing units. This assumption is feasible and practical enough for this study. Furthermore, each processing unit in this study is assumed to have the capability to process different kinds of task unless a specific capability is defined to the processing unit. This means that, in general, more than one task can be assigned to any processing unit. Let P_{v_i} be a set of processing units capable of processing task v_i . The task assignment consists of the following steps.

1. Identify a set of candidate processing units for each task v_i . A candidate processing unit for v_i is a processing unit whose consumed energy is less than or equal to the minimum energy estimated from every processing unit in P_{v_i} . This candidate processing unit is called a *primary candidate* processing unit. Any processing units that are not satisfied with the above

Table 1Power consumption of benchmark desktop CPUs at peak and idle states. The values were taken from [26,27] and were used to estimate energy consumption of each processing unit in the experiment.

Processing unit	CPU	Clock (GHz)	Power consumption (W)	
			Peak	Idle
а	Intel Core i7-975 XE	3.33	240	105
b	Intel Core 2 Extreme QX6850	3.00	233	110
С	Intel Core 2 Extreme QX6700	2.66	229	106
d	Intel Core i7-920	2.66	224	105

- condition will be classified as secondary or tertiary candidate processing units, depending on the following conditions. Processing unit a is called a secondary candidate processing unit if $\alpha_a t_a(v_i) \leqslant \tau$, otherwise it is called tertiary candidate processing unit. Details are given in Algorithm 1.
- 2. Identify a set of actual candidate processing units. For any task v_i and its candidate processing unit a, other dependent tasks of v_i executable by a are virtually assigned to a. The energy consumption of primary candidate and secondary candidate processing units are re-computed and compared. Any candidate processing unit a is replaced by a secondary candidate processing unit b if the energy consumption of a is higher than that of b. Details are given in Algorithm 2.
- 3. Schedule the set of tasks assigned to each processing unit. The scheduling process consists of two phases. In the first phase, a set of dependent tasks assigned to a processing unit is orderly arranged according to their dependent paths in dependent task graph *G* in the form of processing time duration. For those independent tasks, they can be arbitrarily scheduled. Therefore, the scheduled sequence in any processing unit will consist of alternate slots of task duration and slots of idle duration. The second phase is to minimize the idle energy of the whole system. If there are some tasks possibly executed by more than one processing unit, then these tasks are re-assigned to other processing units under the condition that the new total energy must not be increased. The detail of phase 1 is given in Algorithm 3 and that of phase 2 is given in Algorithm 4.

Algorithm 1. Identifying Preliminary Candidate Processing Units

```
Require: v_i, P_{v_i}, and G.
 1. for v_i \in G do
       Let o be the number of out-degrees of v_i.
       \tau = \min_{a \in P_{v_i}}(\alpha_a t_a(v_i)) + \max_{a \in P_{v_i}}(o\kappa_a \lambda_a d_a(v_i) + \mu_a \beta_a w_a(v_i))
 4.
        for each processing unit a \in P_{v_i} do
 5.
          if \alpha_a t_a(v_i) \leq \tau then
 6.
             mark a as a secondary candidate processing unit.
 7.
 8.
             mark a as a tertiary candidate processing unit.
 9.
          end if
10.
       end for
11. end for
```

Algorithm 2. Identifying Actual Candidate Processing Units

```
Require: v_i, P_{v_i}, \mathbf{g}_a, and G.
 1. for v_i \in G do
 2.
        for each secondary candidate processing unit a \in P_{v_i} do
 3.
           Let \mathbf{g}_a = \phi for all a.
           for all ancestors v_i's of v_i do
 4.
 5.
              if v_i is executable by a then
                 Let \mathbf{g}_a = \mathbf{g}_a \cup \{v_j\}.
 6.
 7.
              end if
 8.
           end for
           for all descendants v_k's of v_i do
 9.
10.
              if v_k is executable by a then
11.
                 Let \mathbf{g}_a = \mathbf{g}_a \cup \{v_k\}.
```

```
end if
12.
13
        end for
14.
      end for
      Descendingly sort |\mathbf{g}_a|.
15.
      Mark any processing unit a \in P_{\nu_i} having maximum |\mathbf{g}_a| and lowest execution energy as a primary candidate
    processing unit.
17.
      Mark the rest of processing units in P_{\nu_i} as a secondary candidate processing units.
18. end for
19. for v_i \in G do
      for each secondary candidate processing unit b \in P_{v_i}do
20.
21.
        Compute energy e_b(v_i).
22.
        if \exists candidate processing unit a \in P_{v_i} such that e_b(v_i) < e_a(v_i) then
23.
           Mark processing unit a as a tertiary candidate processing unit.
           Mark processing unit b as a primary candidate processing unit.
24.
25.
         end if
26.
      end for
27. end for
```

Algorithm 3. Phase-1 Preliminary Task Scheduling in Processing Units

```
Require: v_i, P_{v_i}, and G.
 1. Let v_f be the latest task in its primary processing unit in P_{v_f}.
 2. Let \zeta_{v_f} be the finishing time of task v_f.
 3. Descendingly sorted task v_i at first level of G by execution time on its primary processing unit.
 4. for each sorted task v_i at first level of G do
       if primary processing unit P_{v_i} is an empty slot then
 6.
         Assign v_i to its primary processing unit in P_{v_i}.
 7.
       else
 8.
         Assign v_i to its primary processing unit in P_{v_i} at time \zeta_{v_t} + 1.
 9.
       end if
10. end for
11. while \exists unassigned v_i do
      Let \delta be a set of v_k having all ancestor tasks already assigned to their primary processing units.
       for each task v_k \in \delta \mathbf{do}
         Let A_{v_k} be a set of ancestor tasks of v_k already assigned to their primary processing units.
14.
15.
         Find a task v_j \in A_{v_k} having the latest finishing time at time \zeta_{v_i}.
16.
       end for
17.
       Find a task v_k \in \delta having earlier finishing time \zeta_{v_i} and the shortest execution time on its primary processing
18.
       if time slot \zeta_{\nu_i} + w_a(\nu_k) of primary processing unit in P_{\nu_k} is empty then
19.
         Assign v_k to its primary processing unit in P_{v_k} at time slot \zeta_{v_i} + w_a(v_k).
20. else
21.
         Assign v_k to its primary processing unit in P_{v_k} at time slot \zeta_{v_f} + 1.
22.
       end if
23. end while
```

Algorithm 4. Phase-2 Minimizing Idle Energy in Task Scheduling

```
Require: a list of scheduled tasks in each \mathbf{g}_a and G.
 1. Let S be a set of |\mathbf{g}_a| sorted in descending order.
```

- 2. **for** each corresponding $\mathbf{g}_a \in S$ **do**
- Let v_k be the last task in \mathbf{g}_a . 3.
- **while** task v_k is not the first task in \mathbf{g}_a **do**

(continued on next page)

```
Traverse the task list upward starting at v_k until an empty slot E is found and let v_i be the task next to this idle
 5.
    slot.
 6.
        Let A_{v_i} be a set of ancestors of v_i.
 7.
         for all tasks v_i \in A_{v_i} do
 8.
           if there exists a processing unit b with an empty slot and the data dependency of v_i and its ancestors in any
    processing units is not violated then
 9.
             Temporarily remove v_i from the present slot.
10.
             Insert v_i to the beginning of this new empty slot.
11.
             Reschedule all tasks whose starting times are after the ending time of v_i by using Algorithm 3.
             Let G' be the new dependent task graph after the temporary assignment of v_i.
12.
13
             Compute the total energy \Gamma_{G'}.
14.
             if \Gamma_{G'} > \Gamma_G then
                Remove v_i from processing unit b and assign it back to its original processing unit.
15.
16.
17.
                Permanently assign v_i to this new empty slot on processing unit b.
18.
             end if
           end if
19.
20.
        end for
21.
        if v_k is not the first task in \mathbf{g}_a then
22.
           Let v_k be a new task found after traversing list \mathbf{g}_a upward starting from the empty slot E.
23.
         end if
24.
      end while
25. end for
```

To illustrate how these four algorithms work, the algorithms are applied to schedule all tasks in the dependent task graph shown in Fig. 1. Each vertex represents a task with its name. A parenthesis to the right of each vertex v_i is a set of processing units capable of processing task v_i . For example, task v_1 can be processed by processing units a and c. The values of all variables in Eq. (2) are hypothetically given in Tables 2–4. Table 2 provides the amount of processing energy per unit time (α_a), amount of waiting energy per unit time (β_a), and transmission energy per unit amount of data (λ_a) of each processing unit a. In this example, there are four processing units which are a, b, c, and d. Table 3 defines the transmission speed of the link between each processing unit pair in unit amount of data per unit time. These constants are used to estimate data transmission time among processing units. Table 4 shows the amount of transmitted data of each task v_i and its estimated transmission time of each processing unit. To show how the energy consumption is estimated, the activities of task v_1 will be used as an example.

The amount of time to process task v_1 on a processing unit a is $t_a(v_1) = 260$ with $\alpha_a = 0.067$ as defined in Table 2. The estimated execution energy is equal to $\alpha_a t_a(v_1) = 0.067 \times 260 = 17.42$. The energy consumption of the other tasks can be similarly estimated. Table 5 summarizes the estimated energy consumption of all tasks in the given dependent task graph. Suppose v_1 is processed at processing unit a. After finishing task v_1 , the results must be transmitted to its descendant tasks which are v_4 , v_5 , v_6 , and v_7 . These four tasks can be executed by a set of processing units (a, b) for v_4 , (c, d) for v_5 , (b, d) for v_6 , and (a, c) for v_7 . Thus, the amount of transmission energy between processing unit pairs $\{a, b\}$, $\{a, c\}$, $\{a, d\}$ must

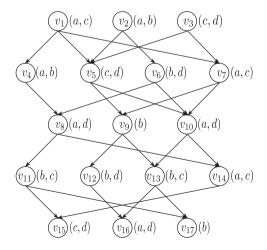


Fig. 1. An example of dependent task graph G. The parentheses to the right of each vertex denote candidate processing units to which the tasks can be executed. For example, vertex $v_1(a, c)$ denotes task v_1 that can be executed on either processing unit a or c.

Table 2 A list of given energy consumption constants for each processing unit $i \in \{a, b, c, d\}$. α_i is the execution energy at peak state per unit time, β_i is the waiting energy at idle state per unit time, λ_i is the transmission energy per unit data.

Energy constants	Processing units			
	а	b	С	d
$\alpha_{i\in\{a,\ b,\ c,\ d\}}$	0.067	0.065	0.064	0.062
	0.029	0.031	0.029	0.029
$eta_{i\in\{a,\ b,\ c,\ d\}}$ $\lambda_{i\in\{a,\ b,\ c,\ d\}}$	0.045	0.047	0.050	0.048

Table 3 Estimated data transmission rate $r_{a,b}$ between any processing unit pair a and b in unit amount of data per unit time.

Processing unit	а	ь	С	d
a	=	1	0.5	1
b	1	=	0.5	1.25
С	0.5	0.5	_	1.25
d	1	1.25	1.25	_

Table 4The initial values of output data size and the estimated execution time of each task processed by each processing unit.

Task	Amount of transmitted data	Amount of ex	ecution time				
v_i	$d_a(v_i)$	$t_a(v_i)$	$t_b(v_i)$	$t_c(v_i)$	$t_d(v_i)$		
1	60	260	-	760	_		
2	80	350	160	_	_		
3	10	_	_	890	820		
4	60	420	150	_	_		
5	90	_	_	940	430		
6	30	_	400	_	560		
7	50	670	_	850	_		
8	40	300	_	_	630		
9	50	_	470	_	_		
10	30	920	_	_	950		
11	80	_	550	120	_		
12	10	_	200	_	130		
13	90	_	100	790	_		
14	40	940	_	580	_		
15	50	_	_	350	310		
16	40	930	_	_	750		
17	70	-	370	=	-		

Table 5Estimated execution energy to execute each task on each processing unit for the first case.

Task	Execution energy			
$ u_i$	$\alpha_a t_a(v_i)$	$\alpha_b t_b(v_i)$	$\alpha_c t_c(v_i)$	$\alpha_d t_d(v_i)$
1	17.42	_	48.64	=
2	23.45	10.40	=	_
3	=	=	56.96	50.84
4	28.14	9.75	=	_
5	-	_	60.16	26.66
6	-	26.00	_	34.72
7	44.89	=	54.40	_
8	20.10	=	=	39.06
9	-	30.55	-	_
10	61.64	_	_	58.90
11	-	35.75	7.68	_
12	-	13.00	_	8.06
13	-	6.5	50.56	_
14	62.98	=	37.12	_
15	-	-	22.40	19.22
16	62.31	_	_	46.50
17	_	24.05	_	_

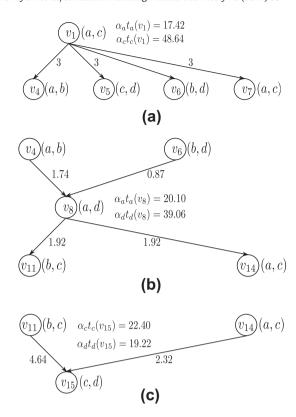


Fig. 2. Examples of applying the Algorithm 1 to identify preliminary candidate processing units for the tasks v_1 , v_8 , and v_{15} . The incoming edge of the considered task v_1 shows the waiting energy having maximum weight for each ancestor task. The outgoing edge shows the transmission energy having maximum weight for the transmission to each descendant task. The numbers next to the task denote the minimum execution energy utilized by the candidate processing units. (a) Task v_1 , (b) Task v_8 , and (c) Task v_{15} .

be estimated. Suppose we consider the pair $\{a, c\}$. From Tables 2 and 4, the amount of transmitted data is $d_a(v_1) = 60$ and the transmission energy per unit amount of data is $\lambda_a = 0.045$. Hence, the transmission energy is equal to $\lambda_a d_a(v_1) = 0.045 - 0.045$. The transmission energy of other processing unit pairs can be similarly estimated.

The waiting energy of any processing unit depends upon the data transmission rate between the sending unit and the receiving unit. Let us consider tasks v_6 and v_8 . Task v_6 can be processed by either processing units b or d and task v_8 can be processed by either processing units a or d. Assume that v_6 is processed by d and v_8 is by a. According to Table 4, the amount of data to be transmitted from d to a is equal to $d_d(v_6) = 30$ and the data transmission rate from d to a is equal to r_d , r_d = 1. Therefore, the data transmission time r_d is equal to r_d and r_d is estimated by r_d is esti

The results produced by each algorithm will be explained next. From Algorithm 1, three different types of tasks will be considered to illustrate how Algorithm 1 works. The first type is a task without any incoming degree such as task v_1 . The second type is a task with both incoming and outgoing degrees such as task v_8 . The last type is a task having only incoming degree such as task v_1 . For task v_1 , the value of minimum energy τ estimated from every processing unit in list P_{v_1} is equal to

$$\tau = \min_{k \in P_{\nu_1}} (\alpha_k t_k(\nu_1)) + \max_{k \in P_{\nu_1}} (o\kappa_k \lambda_k d_k(\nu_1) + \mu_k \beta_k w_k(\nu_1))$$

$$\tag{3}$$

$$= 17.42 + 3 \times 3 + 0 = 26.42 \tag{4}$$

Since v_1 has no incoming degree, the term $\mu_k \beta_k w_k(v_1)$ is set to zero. For task v_8 , the value of minimum energy τ estimated from every processing unit in list P_{v_8} is equal to

$$\tau = \min_{k \in P_{v_8}} (\alpha_k t_k(\nu_8)) + \max_{k \in P_{v_8}} (o\kappa_k \lambda_k d_k(\nu_8) + \mu_k \beta_k w_k(\nu_8))$$
(5)

$$=20.10+2\times1.92+1.74=25.68\tag{6}$$

For task v_{15} , the value of minimum energy au estimated from every processing unit in list $P_{v_{15}}$ is equal to

$$\tau = \min_{k \in P_{\nu_{15}}} (\alpha_k t_k(\nu_{15})) + \max_{k \in P_{\nu_{15}}} (\sigma_k \lambda_k d_k(\nu_{15}) + \mu_k \beta_k w_k(\nu_{15}))$$
(7)

$$= 19.22 + 0 + 4.64 = 23.86 \tag{8}$$

From Fig. 2, processing unit a is selected as a secondary candidate, while processing unit c is selected as a tertiary candidate for v_1 with respect to τ . Fig. 3 shows all secondary candidate processing units for each task after applying Algorithm 1.

In the second step, Algorithm 2 is used to identify the actual candidate processing units. The input is a dependent graph *G* obtained from the first step as shown in Fig. 3. Algorithm 2 counts the number of ancestors as well as descendants of each task with respect to each secondary candidate processing unit. Any secondary processing unit of a task is set as a primary candidate processing unit if the task has the largest number of counts. Then the algorithm computes the actual energy consumption for each primary and secondary candidate processing units. An example is shown in Fig. 4 along with procedural computations below.

For task v_{10} in Fig. 3, both \mathbf{g}_a and \mathbf{g}_d have the same counts of 2. As the execution energy of task v_{10} on processing unit d is less than the execution energy on processing unit a, the processing unit d is marked as a primary candidate processing unit. Similarly, for task v_{15} in Fig. 3, group \mathbf{g}_c has the maximum counts of 3. Thus, the processing unit c is marked as a primary candidate processing unit.

After identifying the primary candidate processing units, the algorithm determines the energy consumption of primary and secondary processing units of each task. For example, consider task v_{10} . This task has both incoming and outgoing degrees. The energy consumption of processing units a and d for task v_{10} are shown in Fig. 4(a) and (b), respectively. The energy consumption of task v_{10} by processing unit a is equal to

$$e_a(v_{10}) = \alpha_a t_a(v_{10}) + \kappa_a \lambda_a d_a(v_{10}) + \mu_a \beta_a w_a(v_{10}) \tag{9}$$

$$= 61.64 + 2 \times 1.35 + 2.61 \tag{10}$$

$$=66.95$$
 (11)

and by processing unit d is equal to

$$e_d(v_{10}) = \alpha_d t_d(v_{10}) + \kappa_d \lambda_d d_d(v_{10}) + \mu_d \beta_d w_d(v_{10}) \tag{12}$$

$$= 58.90 + 2 \times 1.44 + 1.45 \tag{13}$$

$$=63.23$$
 (14)

Since $e_d(v_{10})$ is less than $e_a(v_{10})$, the actually selected candidate processing unit for task v_{10} is d. For a task without any outgoing degree such as v_{15} , a similar computation for energy consumption can be estimated. Task v_{15} has two candidate processing units which are c and d. Fig. 4(c) and 4(d) show the energy consumption of each processing unit for task v_{15} . The energy consumption of task v_{15} by processing unit c is equal to

$$e_c(v_{15}) = \alpha_c t_c(v_{15}) + \kappa_c \lambda_c d_c(v_{15}) + \mu_c \beta_c w_c(v_{15}) \tag{15}$$

$$= 22.40 + 0 + 0 \tag{16}$$

$$= 22.40$$
 (17)

and by processing unit d is equal to

$$e_d(v_{15}) = \alpha_d t_d(v_{15}) + \kappa_d \lambda_d d_d(v_{15}) + \mu_d \beta_d w_d(v_{15})$$
(18)

$$=19.22+0+1.856$$
 (19)

$$=21.076$$
 (20)

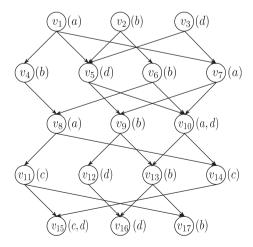


Fig. 3. Results from the first step after applying Algorithm 1 to the dependent task graph G of Fig. 1. Note that some tasks, such as v_{10} and v_{15} still have more than one secondary candidate processing unit.

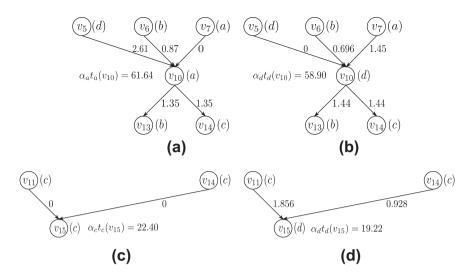


Fig. 4. Examples of candidate processing units for tasks v_{10} and v_{15} . The tasks still contain secondary candidate processing units at the end of the first step and all the ancestors and descendants are primary candidate processing units. The incoming and outgoing edges are actual waiting and transmission energy, respectively. (a) Task v_{10} on processing unit a, (b) Task v_{10} on processing unit d, (c) Task v_{15} on processing unit d, and (d) Task v_{15} on processing unit d.

Because $e_d(v_{15})$ is less than $e_c(v_{15})$, the actually selected candidate processing unit for task v_{15} is d. The actual processing unit of each task in the dependent task graph G as the result of Algorithm 2 is shown in Fig. 5.

Scheduling is performed by Algorithm 3. For example, from Fig. 5, there are three tasks in the first level, namely, v_1 , v_2 and v_3 , with the assigned primary candidate processing units a, b, and d, having the execution time of 260, 160, and 820, respectively. Thus, task v_2 will be scheduled first, followed by v_1 , and v_3 . Next, the unassigned task v_6 whose ancestor task v_2 is considered. Task v_6 has the latest finishing time comparing with all of its ancestors, but has the earliest finishing time comparing with the ancestors of other tasks. Thus, it is assigned to primary candidate processing unit b. Task v_4 , the descendant task of v_1 , has the latest finishing time comparing with all of its ancestors, but has the earliest finishing time comparing with the ancestors of other tasks. Thus, it is assigned to primary candidate processing unit b. Task b_4 , the descendant of b_4 , is scheduled next. It has the latest finishing time comparing with all of its ancestors, but has the earliest finishing time comparing with the ancestors of other tasks. Therefore, it is assigned to primary candidate processing unit b_4 . Tasks b_4 and b_4 have a common ancestor task b_4 has the latest finishing time comparing with all of its ancestors, but has the earliest finishing time comparing with the ancestors of other tasks. However, task b_4 has a shorter execution time. Hence, it is assigned to primary candidate processing unit b_4 . The algorithms will continue until all tasks are assigned. Fig. b_4 illustrates the list of scheduled tasks assigned by Algorithm 3.

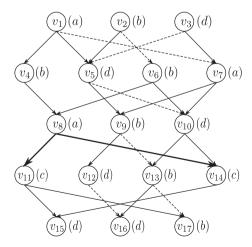


Fig. 5. Results of Algorithm 2. A solid line represents data transmission between different processing units. A dashed line represents data transmission within the same processing unit. A thick line represents data transmission from a task to its descendant tasks which are executed by the same processing unit, e.g., v_8 to v_{11} and v_{14} .

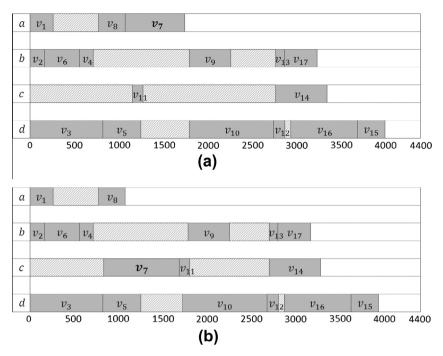


Fig. 6. The scheduling results of Algorithms 3 and 4. (a) The result of Algorithm 3, (b) The result of Algorithm 4. In this example, only task v_7 is removed from processing unit a and re-assigned to processing unit c.

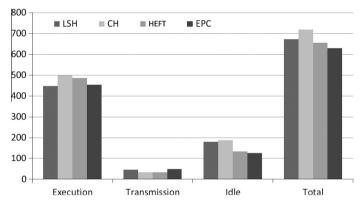


Fig. 7. Results of energy consumption for the first case.

Algorithm 4 is applied further to reduce the empty slots so as to minimize the idle energy of the scheduled tasks. From Fig. 6(a), the latest finished task is v_{15} on processing unit d. The algorithm traverses the time slots of processing unit d upwards until an empty slot is found. In this case, v_{16} is the task next to this first idle slot and its ancestors are v_{12} and v_{13} . A re-assignment trial of v_{12} and v_{13} to other processing units is proceeded. Although there are empty slots and no data dependency violation for v_{12} and v_{13} , the re-assignment of v_{12} and v_{13} increases the amount of total energy. Therefore, no re-assignment for v_{12} and v_{13} is done. The next empty slot when traversing the time slot towards the beginning is a slot before v_{10} . Task v_{10} has three ancestors which are v_{5} , v_{6} , and v_{7} . In this case, v_{7} can be re-assigned to processing unit c which decreases the idle energy and total energy. Other tasks in each processing unit are re-assigned in a similar fashion. Fig. 6(b) depicts the results from Algorithm 4.

5. Experimental results and discussion

We conducted four experimental cases on LSH, CH, HEFT, and EPC algorithms to test and evaluate their performance using different scheduling scenarios. The dependent task graph in the first case was synthesized. The second case was taken directly from [25] as a validation benchmark. The third case employed the same scenario as that of the second case but

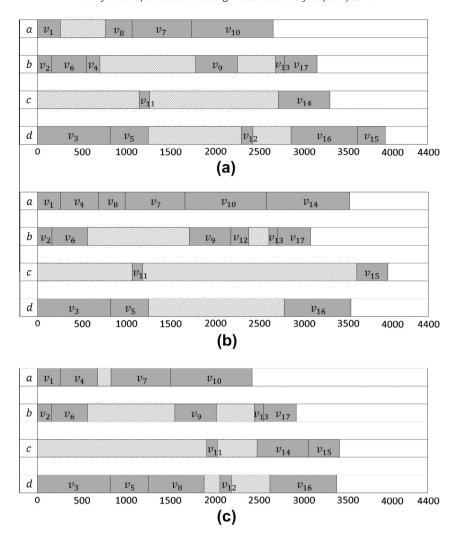


Fig. 8. A list of scheduled tasks in each processing unit of dependent task graph in Fig. 1. (a) List scheduling heuristic (LSH), (b) Clustering heuristic (CH), (c) Heterogeneous earliest finish time (HEFT).

 Table 6

 Energy consumption in each processing unit by LSH algorithm in the first case.

Processing unit	Execution energy	Transmission energy	Idle energy	Total energy
а	144.05	12.15	14.79	170.99
b	107.25	14.57	46.81	168.63
С	44.80	10.00	75.40	130.20
d	151.28	9.12	42.98	203.38
Total	447.38	45.84	179.98	673.20

imposed the processing unit limitation of the first case on the second case. The fourth case was extended from the first case to cover more complex task graphs and emphasize on execution intensive scenario rather than transmission load. Thus, consideration of the transmission time in the first and the fourth cases was not our main concern in this situation as the network capability is technologically increased.

The simulation environment for these experiments consists of four heterogeneous processing units. Each unit has unique competence but different energy consumption during peak or idle states. The amount of energy consumption for the four processing units were selected from the values of experimental energy consumption database for a number of CPU specifications [26,27]. The selected values are shown in Table 1. In this study, the energy consumption was converted to execution energy at peak state and waiting energy at idle state in watts per seconds. For example, for processing unit *a*, the execution

energy per unit time α_a = 240/3600 = 0.067 and the waiting energy per unit time β_a = 105/3600 = 0.029. As processing units are connected to one another, the transmission energy (λ) were assigned to each pair. The values were randomly generated within the range of energy consumption during the peak state and the idle state. The obtained values for the execution energy per unit time, waiting energy per unit time, and transmission energy per unit data are shown in Table 2. In order to calculate delay time due to data transmission, the network transmission speed for all the processing unit pairs were assigned as shown in Table 3.

For the first case, the dependent task graph is shown in Fig. 1. There are 17 given tasks, i.e., v_1 – v_{17} , and 4 processing units, i.e., a, b, c, and d. The tasks were created with different attributes and the processing units possessed various competency to handle individual task execution requirements. The initial assignment of processing units to their executable tasks are shown in Fig. 1. The size of output data and the estimated execution time of each task processed by each processing unit from the beginning of simulation are given in Table 4. The total execution energy consumed by each processing unit for each task is shown in Table 5.

Although other scheduling algorithms, namely, LSH, CH, HEFT, did not consider energy consumption as their cost functions, they are worthy to be deployed in performance comparison. Fig. 7 shows the comparison of energy consumption of LSH, CH, and HEFT scheduling algorithms with the proposed algorithm based on the first experimental case. This energy consumption was set up as a cost function during the scheduling process for each processing unit in terms of execution, transmission, and idle energy. The introduction of energy consumption was not deployed as a part of the cost function computation in those comparative algorithms. The resulting schedules are depicted in Fig. 8. The dark and striped slots denote execution time and idle time of each processing unit, respectively. The system completion time of LSH, CH, HEFT, and EPC were 3922, 3940, 3410, and 3924, respectively. At any rate, this number is of no concern for this study as this overhead could be allocated off-line, thus imposing no performance degradation for the proposed algorithm.

The execution, transmission, and idle energy consumption of each processing unit obtained from LSH, CH, HEFT, and EPC are shown in Tables 6–9, respectively. Note that LSH algorithm has the least execution energy usage but ranks third in the total energy usage as the algorithm concerns mainly execution energy. The least transmission energy usage belongs to CH algorithm, although the algorithm yielded the most total energy consumption as it concerns mainly transmission energy. HEFT has the least execution time, but came in the second place for the total energy usage as the algorithm concerns mainly the idle and execution time. The EPC yielded good execution and total energy usages because the algorithm selected the

Table 7Energy consumption in each processing unit by CH algorithm in the first case.

Processing unit	Execution energy	Transmission energy	Idle energy	Total energy
а	235.17	9.90	0.00	245.07
b	110.50	9.87	42.47	162.84
С	30.08	4.00	100.63	134.71
d	124.00	9.12	44.14	177.26
Total	499.75	32.89	187.24	719.88

Table 8Energy consumption in each processing unit by HEFT algorithm in the first case.

Processing unit	Execution energy	Transmission energy	Idle energy	Total energy
а	152.09	10.35	4.35	166.79
b	97.50	13.16	44.02	154.68
c	67.20	4.00	68.44	139.64
d	171.12	6.24	17.75	195.12
Total	487.91	33.75	134.56	656.22

Table 9Energy consumption in each processing unit by EPC algorithm in the first case.

Processing unit	Execution energy	Transmission energy	Idle energy	Total energy
а	37.52	9.90	14.79	62.21
b	107.25	15.98	46.87	170.10
С	99.20	15.00	49.94	164.14
d	210.18	7.68	15.49	233.35
Total	454.15	48.56	127.09	629.80

Table 10 Execution cost of dependent task graph of Fig. 9 for the second case.

Task	$lpha_a t_a(v_i)$	$lpha_b t_b(v_i)$	$\alpha_c t_c(v_i)$
1	14	16	9
2	13	19	18
3	11	13	19
4	13	8	17
5	12	13	10
6	13	16	9
7	7	15	11
8	5	11	14
9	18	12	20
10	21	7	16

Table 11Execution cost of dependent task graph of Fig. 9 for the third case.

Task	$\alpha_a t_a(v_i)$	$\alpha_b t_b(v_i)$	$\alpha_c t_c(v_i)$
1	14	_	9
2	13	19	=
3	=	-	19
4	13	8	-
5	-	-	10
6	-	16	-
7	7	-	11
8	5	-	-
9	-	12	_
10	21	=	_

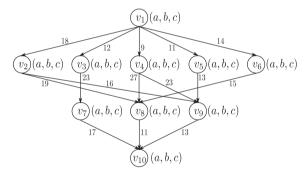


Fig. 9. The dependent task graph with similar transmission and execution costs from [25] used in the second case.

optimal server assignment for minimal execution and transmission energy consumption, and proceeded to reduce idle energy to attain the best total energy performance. In addition, the optimal time-line might not be reached as it is not the prime directive of EPC algorithm (see Tables 10 and 11).

The second case used a task graph taken from [25] as shown in Fig. 9. Each number at each edge of the graph denotes the transmission time. The transmission energy was determined in terms of delay, which was different for every task. The execution energy of each task depended on a processing unit. The waiting energy was set to 50% of execution energy. Scheduling of a task limited a processing unit to use only one unit of execution energy per one unit of execution time and one unit of transmission energy per one unit of transmission time. The execution energy, transmission energy, idle energy, and total energy used by each algorithm are shown in Fig. 10. The system completion time of LSH, CH, HEFT, and EPC were 80, 139, 80, and 127, respectively. The reason why LSH and HEFT ended up with the shortest computation time is because they focused primarily on execution time and did not take energy consumption issue into consideration. On the other hand, the transmission and idle energy consumption was equal to zero in EPC algorithm because all processing units were capable of executing every task and processing unit *a* required the minimum energy.

For the third case using the task graph shown in Fig. 9, the same capabilities and conditions of the processing units as those in the second case were adopted. The results are shown in Fig. 11. The system completion time of LSH, CH, HEFT, and EPC were 102, 116, 102, and 116, respectively.

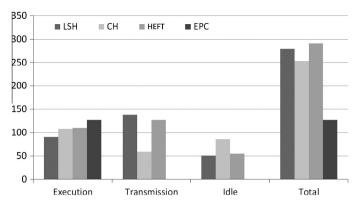


Fig. 10. Results of energy consumption for the second case. Note that in case of transmission and idle situations, the results from EPC were not shown because all tasks in EPC were executed only at processing unit *a*.

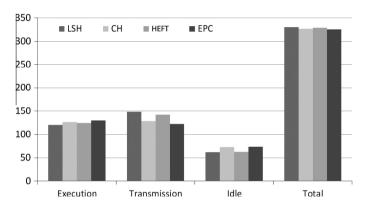


Fig. 11. Results of energy consumption for the third case.

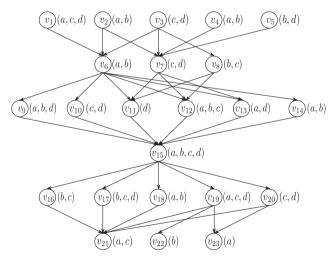


Fig. 12. The dependent task graph for the fourth case.

Similar results to those of the second case mean that the adopted capabilities and conditions have a little effect on the performance of each algorithm.

The fourth case uses a generated task graph shown in Fig. 12. The transmission time was reduced to 10% of the execution time. The execution time and data size of this case for each task are shown in Table 12. The same estimated energy consumption and data transmission coefficients of each processing unit used in the first case were used in this case as shown in

Table 12 A list of transmission data of each task $d_a(v_i)$ and execution time by each processing unit $t_a(v_i)$ for the fourth case.

Task	Amount of transmission data	Amount of ex	ecution time		
v_i	$d_a(v_i)$	$t_a(v_i)$	$t_b(v_i)$	$t_c(v_i)$	$t_d(v_i)$
1	25	350	-	570	430
2	30	430	270	_	_
3	20	_	_	510	450
4	45	590	710	_	_
5	60	_	920	_	790
6	70	360	480	_	_
7	30	_	_	840	720
8	55	_	710	380	_
9	20	390	670	_	380
10	25	_	_	550	580
11	15	_	_	_	620
12	65	870	790	720	_
13	20	560	-	-	540
14	10	430	410	_	_
15	85	280	260	270	250
16	35	_	520	480	_
17	20	_	380	390	420
18	15	610	690	-	_
19	30	250	_	270	480
20	10	-	-	620	710
21	70	760	-	890	-
22	30	-	540	-	-
23	35	580	_	_	_

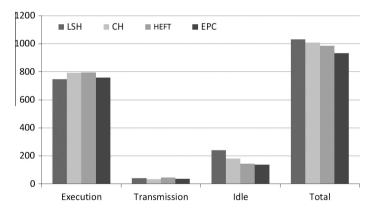


Fig. 13. Results of energy consumption for the fourth case.

Table 2. The execution energy, transmission energy, idle energy, and total energy used by each algorithm are shown in Fig. 13. The system completion time of LSH, CH, HEFT, and EPC were 6278, 6299, 5230, and 5450, respectively. HEFT algorithm gave the minimum scheduling length, while EPC algorithm gave the lowest total energy. As seen from Fig. 13, a significant reduction of the transmission energy from the total energy consumption was given by CH algorithm. In EPC algorithm, additional idle energy reduction helped shorten the scheduling length at the small expense of execution and transmission energy. However, EPC algorithm still yielded the minimum total energy.

6. Conclusion

A new set of task scheduling and assignment algorithms focusing on minimizing energy consumption, named Energy-Efficient Process Clustering Assignment algorithm or EPC algorithm, was proposed. In this study, multiple capabilities of each processing unit was involved during the task scheduling and enhance the assignment flexibility. The proposed Algorithm 4 significantly reduces the idle time and further minimizes the energy consumption which is not properly handled by the other compared algorithms. The experimental results showed that EPC algorithm spent less energy consumption than those from LSH, CH, and HEFT algorithms in different scenarios. Although the significant benefits of EPC algorithm was confirmed by the experimental results, the performance of the algorithm still depends upon the provided information on estimated transmission data of each task, execution time of each task, and the computational capability of each processing unit.

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