

# Parallel Algorithms On Some Numerical Techniques Using PVM Platform On A Cluster Of Workstations

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## Abstract

In this paper, a few parallel algorithms are explained in solving one dimensional heat model problem using Parallel Virtual Machine (PVM). This research focuses on two iteration methods, Iterative Alternating Decomposition Explicit Method (IADE) and Alternating Group Explicit Scheme (AGE). Conjugate Gradient (CG) is selected as an alternative method to accelerate the convergent and efficiency of these two iteration methods.

## keywords

Parallel Virtual Machine(PVM), Iterative Alternating Decomposition Explicit Method(IADE), Alternating Group Explicit Scheme(AGE), Conjugate Gradient(CG).

## 1.0 Introduction

Six strategies of parallel algorithms are implemented to exploit the convergence of IADE\_CG. In the domain decomposition strategy  $\Omega$ , the IADE\_CG Michell-Fairweather which is fully explicit, is derived to produce the approximation of grid- $i$  and not totally dependent on the grid  $(i - 1)$  and  $(i + 1)$ . In IADE\_CG Red Black and IADE\_CG SOR strategies, the domain  $\Omega$  is decomposed into two different subdomains  $\Omega^H$  and  $\Omega^M$ . The concept of multidomain is observed in the IADE\_CG Multicoloring method. The decomposition of domain  $\Omega$  into  $w$  different groups of domain. The Domains for colors  $1, 2, 3, \dots, w$  are noted as  $\Omega^{w_1}, \Omega^{w_2}, \dots, \Omega^{w_w}$ . For the computational grid for domain  $\Omega$ , its execution started with level  $\Omega^{w_1}$ , followed by level  $\Omega^{w_2}$  and ends with level  $\Omega^{w_w}$ . On the vector iteration strategy, parallel IADE\_CG is run in two sections. This method converges if the inner convergence criterion is achieved for each section.

On the strategy of Incomplete Block LU preconditioners on slightly overlapping subdomains, the domain  $\Omega$  is decomposed into  $p$  processors with incomplete subdomain  $\bar{\Omega}$ . This strategy used a preconditioners, the incomplete factorization with certain parameters of algebraic boundary condition. Thus, AGE\_CG algorithm is shown to be extremely straightforward as implemented in parallel algorithms using PVM. CG is chosen as the alternative parallel algorithm because it does not increase the communication time between the processors. The application of CG is a correction to the parallel and sequential algorithms of IADE\_CG and AGE\_CG.

## 2.0 IADE\_CG Algorithms

IADE with Mitchell-Fairweather is introduced by Evans and Sahimi (1992) for solving the one dimensional heat problems. A generalized finite different approximation to the different equation at point  $(x_i, t_{j+\frac{1}{2}})$  is given by,

$$\begin{aligned} -\lambda\theta u_{i-1,j+1} + (1 + 2\lambda\theta)u_{i,j+1} - \lambda\theta u_{i+1,j+1} &= \lambda(1 - \theta)u_{i-1,j} + [1 - 2\lambda(1 - \theta)]u_{i,j} \\ &+ \lambda(1 - \theta)u_{i+1,j}, \quad i = 1, 2, 3, \dots, m, \end{aligned} \quad (1)$$

which leads to the three-point formulae, displayed in matrix form as

$$\mathbf{A}\mathbf{u} = \mathbf{f} \quad (2)$$

The IADE iterative employs the fractional splitting strategy,

$$(rI + \mathbf{G}_1)u^{(k+\frac{1}{2})} = (rI - g\mathbf{G}_2)u^{(k)} + \mathbf{f} \quad (3)$$

$$(rI + \mathbf{G}_2)u^{(k+1)} = (rI - g\mathbf{G}_1)u^{(k+\frac{1}{2})} + g\mathbf{f} \quad (4)$$

where the coefficient matrix  $A$  can be decomposed into the matrices  $\mathbf{G}_1$  and  $\mathbf{G}_2$ , as

$$\mathbf{A} = \mathbf{G}_1 + \mathbf{G}_2 - \frac{\mathbf{G}_1\mathbf{G}_2}{g}$$

The consistent  $g = \frac{6+r}{6}$ ,  $r$  is acceleration parameter and the constituent matrices  $\mathbf{G}_1$  and  $\mathbf{G}_2$  take the bidiagonal form (lower and upper respectively)

i. at the  $(k + \frac{1}{2})^{th}$  iterate

$$u_i^{(k+\frac{1}{2})} = \frac{1}{d}(-k_{i-1}u_{i-1}^{(k+\frac{1}{2})} + q_i u_i^{(k)} + w u_{i+1}^{(k)} + f_i), \quad i = 1, 2, \dots, m, \quad k_0 = 0 \quad (5)$$

with  $q_i = \frac{r-ge_i}{d}$ ,  $\forall i \in [1, m]$

ii. at the  $(k + 1)^{th}$  iterate

$$u_{m+1-i}^{(k+1)} = \frac{1}{d_{m+1-i}}(-v_{m-i}u_{m-i}^{(k+\frac{1}{2})} + s u_{m+1-i}^{(k+\frac{1}{2})} + g f_{m+1-i} - h_{m+1-i} u_{m+2-i}^{(k+1)}), \quad i = 1, 2, \dots, m, \quad v_0 = 0 \quad (6)$$

with  $d_i = r + e_i$  and  $\forall i \in [1, m]$ . Sequential IADE shows that the approximation solution for  $u_i^{(k+\frac{1}{2})}$  is dependent on  $u_{i-1}^{(k+\frac{1}{2})}$  and approximation solution for  $u_{m+1-i}^{(k+1)}$  is dependent on  $u_{m+2-i}^{(k+1)}$ . The parallel algorithm strategies are implemented to avoid the sequential IADE situation.

### 3.0 AGE\_CG Algorithms

Through the Alternating Direction Implicit (ADI), AGE methods with Peaceman-Rachford variation is created to be more extremely powerful, flexible and it offers users many advantages. The accuracy of this method is comparable if not better than that of the GE class of problems as well as other existing schemes (Evans and Abdullah, 1983). This method employs the fractional splitting strategy and the implicit form is as follows,

$$\begin{aligned} u^{(k+\frac{1}{2})} &= (\mathbf{G}_1 + rI)^{-1}[(rI - \mathbf{G}_2)u^{(k)} + \mathbf{f}] \\ u^{(k+1)} &= (\mathbf{G}_2 + rI)^{-1}[(rI - \mathbf{G}_1)u^{(k+\frac{1}{2})} + \mathbf{f}], \end{aligned} \quad (7)$$

we have

$$\mathbf{A} = \mathbf{G}_1 + \mathbf{G}_2$$

If we assume  $m$  to be odd then  $\widehat{\mathbf{G}}$  could be written as,

$$\widehat{\mathbf{G}} = \begin{bmatrix} r_2 & b \\ c & r_2 \end{bmatrix}_{(2 \times 2)}$$

where,  $r_2 = r + \frac{a}{2}$ . The alternating implicit nature of the  $(2 \times 2)$  groups where the implicit and explicit values are given on the forward and backward levels for sweeps on the  $(k + \frac{1}{2})^{th}$  and  $(k + 1)^{th}$  levels, with  $r_1 = r - \frac{a}{2}$ ,  $r_2 = r + \frac{a}{2}$  end  $\Delta = r_2^2 - bc$

i. at the  $(k + \frac{1}{2})^{th}$  iterate

$$\begin{aligned} u_1^{(k+\frac{1}{2})} &= \frac{r_1 u_1^{(k)} - b u_2^{(k)} + f_1}{r_2} \\ u_i^{(k+\frac{1}{2})} &= \frac{A u_{i-1}^{(k)} + B u_i^{(k)} + C u_{i+1}^{(k)} + D u_{i+2}^{(k)} + E_i}{\Delta} \\ u_{i+1}^{(k+\frac{1}{2})} &= \frac{\tilde{A} u_{i-1}^{(k)} + \tilde{B} u_i^{(k)} + \tilde{C} u_{i+1}^{(k)} + \tilde{D} u_{i+2}^{(k)} + \tilde{E}_i}{\Delta} \end{aligned} \quad (8)$$

with  $i = 2, 4, 6, \dots, m-1$ ,  $A = -cr_2$ ,  $B = r_1 r_2$ ,  $C = -br_1$ ,  $E_i = r_2 f_i - b f_{i+1}$ ,

$$D = \{b^2, i \neq m-1\}$$

and  $\tilde{A} = -cr_2$ ,  $\tilde{B} = r_1 r_2$ ,  $\tilde{C} = -br_1$ ,  $\tilde{E}_i = r_2 f_i - b f_{i+1}$ ,

$$\tilde{D} = \{-br_2, i \neq m-1\}$$

ii. at the  $(k + 1)^{th}$  iterate

$$\begin{aligned} u_i^{(k+1)} &= \frac{P u_{i-1}^{(k+\frac{1}{2})} + Q u_i^{(k+\frac{1}{2})} + R u_{i+1}^{(k+\frac{1}{2})} + S u_{i+2}^{(k+\frac{1}{2})} + T_i}{\Delta} \\ u_{i+1}^{(k+1)} &= \frac{\tilde{P} u_{i-1}^{(k)} + \tilde{Q} u_i^{(k)} + \tilde{R} u_{i+1}^{(k)} + \tilde{S} u_{i+2}^{(k)} + \tilde{T}_i}{\Delta} \\ u_m^{(k+1)} &= \frac{-c u_{m-1}^{(k+\frac{1}{2})} + r_1 u_m^{(k+\frac{1}{2})} + f_m}{r_2} \end{aligned} \quad (9)$$

with  $Q = r_1 r_2$ ,  $R = -br_1$ ,  $S = b^2$ ,  $T_i = r_2 f_i - b f_{i+1}$ ,

$$\tilde{P} = \{c^2, i \neq 1\}$$

and  $\tilde{Q} = -cr_1$ ,  $\tilde{R} = \tilde{Q} = r_1 r_2$ ,  $\tilde{S} = -br_2$ ,  $\tilde{T}_i = -c f_i + r_2 f_{i+1}$ ,

All the equations are not dependent on every point  $i-1$ ,  $i$  and  $i+1$  event for every time step. This is the advantages to create the parallel algorithm for AGE.

Furthermore, Conjugate Gradient algorithm (CG) is exploited to accelerate the convergence and efficiency of of IADE and AGE iterations. If  $\mathbf{r}_k = \mathbf{b} - \mathbf{A}\mathbf{x}_k$  is residual vector at level  $k$  and if  $\mathbf{A}$  is symmetric matrix ( $m \times m$ ) end positively defined, the CG algorithm is as follows,

$$\begin{aligned} \text{i. } \mathbf{p}_k &\leftarrow \mathbf{A}\mathbf{d}_k \\ \text{ii. } \mathbf{x}_{k+1} &\leftarrow \mathbf{x}_k + \alpha_k \mathbf{d}_k, \quad \alpha_k = \frac{\mathbf{r}_k^T \mathbf{r}_k}{\mathbf{d}_k^T \mathbf{p}_k} \\ \text{iii. } \mathbf{r}_{k+1} &\leftarrow \mathbf{r}_k - \alpha_k \mathbf{p}_k, \\ \text{iv. } \mathbf{d}_{k+1} &\leftarrow \mathbf{r}_{k+1} + \beta_k \mathbf{d}_k, \quad \beta_k = \frac{\mathbf{r}_{k+1}^T \mathbf{r}_{k+1}}{\mathbf{r}_k^T \mathbf{r}_k} \end{aligned} \quad (10)$$

if  $\mathbf{x}_0$  is initial value  $\mathbf{x}$  and  $\mathbf{d}_0 = \mathbf{r}_0 = \mathbf{b} - \mathbf{A}\mathbf{x}_0$  is initial value  $\mathbf{r}$  and  $\mathbf{d}$ . A few type of CG algorithms is created by chosen the value of acceleration parameter  $\alpha_k$  and  $\beta_k$ . The chosen normal value of acceleration parameter [Reid(1971)] is,

$$\alpha_k = \frac{\mathbf{d}_k^T \mathbf{r}_k}{\mathbf{p}_k^T \mathbf{d}_k} \quad \text{end} \quad \beta_k = \frac{-\mathbf{p}_k^T \mathbf{r}_{k+1}}{\mathbf{p}_k^T \mathbf{d}_k}$$

The first step, the IADE\_CG Algorithms is executed by using the equation (5) and sweeps along the domain  $\Omega$ . The second step, executes CG algorithms by using the equation (10)( $i - iv$ ). These two steps are continued until a specified convergence criterion is satisfied, when the requirement  $|u_{i,j}^{(k-1)} - u_{i,j}^{(k)}| \leq \varepsilon$  is met and  $\varepsilon$  becomes the convergence criterion. The accuracy of the solution at these grid points was determined by computing its root means square error. For an introduction to conjugate gradient acceleration of iterative methods, see Hageman and Young (1981).

## 4.0 Parallel IADE\_CG Algorithms

### 1. IADE\_CG Michell-Fairweather

The IADE\_CG Michell-Fairweather which is fully explicit is derived to produce the approximation of grid- $i$  and which is not totally dependent on the grid  $(i - 1)$  and  $(i + 1)$ . The approximation at the first and second intermediate levels are computed directly by inverting  $(rI + \mathbf{G}_1)$  and  $(rI + \mathbf{G}_2)$ . The explicit form of equation (3) and (4) is given by

$$u^{(k+\frac{1}{2})} = (rI + \mathbf{G}_1)^{-1}(rI - g\mathbf{G}_2)u^{(k)} + \mathbf{f} \quad (11)$$

$$u^{(k+1)} = (rI + \mathbf{G}_2)^{-1}(rI - g\mathbf{G}_1)u^{(k+\frac{1}{2})} + g\mathbf{f} \quad (12)$$

In the first step, IADE\_CG Michell-Fairweather implements the equations (11) and (12) and the second step, exploits the CG algorithm as in equation (10)( $i - iv$ )

The parallel IADE\_CG Michell-Fairweather is executed by decomposing the  $m$  grid into  $\frac{m}{p}$  groups of grid and assigning the groups into  $p$  processors. For example if  $G_1, G_2, \dots, G_{\frac{m}{p}}$  are the groups of grid and  $p_1, p_2, \dots, p_p$  are the processors, the  $G_i, i = 1, 2, \dots, \frac{m}{p}$  are assigned to  $p_j, j = 1, 2, \dots, p$ .

### 2. IADE\_CG Red Black

In IADE\_CG Red Black strategy, the domain  $\Omega$  is decomposed into two different subdomains  $\Omega^H$  and  $\Omega^M$ .  $\Omega^H$  is the approximate solution on the odd grids and  $\Omega^M$  is the approximate solution on the even grids. Computation on  $\Omega^H$  is executed followed by  $\Omega^M$ . These two subdomains are not dependent on each other.  $\Omega^H$  is decomposed into groups,  $H_1, H_2, \dots, H_{\frac{m}{p}}$  and  $\Omega^M$  is decomposed into groups,  $M_1, M_2, \dots, M_{\frac{m}{p}}$ . Every group of  $H_i$  and  $M_i, i = 1, 2, \dots, \frac{m}{p}$  is assigned to processors  $p$ . IADE\_CG Red Black is run in parallel for each subdomain in alternating way on time steps  $(k + \frac{1}{2})$  and  $(k + 1)$ . The parallel IADE\_CG Red Black formulae for (3) and (4) are as follows,

i. at the  $(k + \frac{1}{2})^{th}$  iterate

$$\begin{aligned} a.u_i^{k+\frac{1}{2}} &= -\frac{q_i u_i^k}{d} + \frac{\omega}{d}(-k_{i-1}u_{i-1}^k + q_i u_i^k + w_i u_{i+1}^k + f_i), \\ i &= 1, 3, 5, \dots, m \\ b.u_i^{k+\frac{1}{2}} &= -\frac{q_i u_i^k}{d} + \frac{\omega}{d}(-k_{i-1}u_{i-1}^{k+\frac{1}{2}} + q_i u_i^k + w_i u_{i+1}^k + f_i), \\ i &= 2, 4, 6, \dots, m-1, k_0 = w_0 = 0 \end{aligned}$$

ii. at the  $(k + 1)^{th}$  iterate

$$\begin{aligned}
a. u_{m+1-i}^{k+1} &= -\frac{su_{m+1-i}^{k+\frac{1}{2}}}{d_{m+1-i}} + \frac{\omega}{d_{m+1-i}}(-v_{m-i}u_{m-i}^{k+\frac{1}{2}} + su_{m+1-i}^{k+\frac{1}{2}} + gf_{m+1-i} - h_{m+1-i}u_{m+2-i}^k), \\
i &= 1, 3, 5, \dots, m, v_0 = 0 \\
b. u_{m+1-i}^{k+1} &= -\frac{su_{m+1-i}^{k+\frac{1}{2}}}{d_{m+1-i}} + \frac{\omega}{d_{m+1-i}}(-v_{m-i}u_{m-i}^{k+\frac{1}{2}} + su_{m+1-i}^{k+\frac{1}{2}} + gf_{m+1-i} - h_{m+1-i}u_{m+2-i}^{k+1}), \\
i &= 2, 4, 6, \dots, m-1
\end{aligned}$$

and acceleration parameter  $\omega = 1$ .

The next step, exploits the CG algorithm as in equation (10)( $i - iv$ )

### 3. IADE\_CG RBSOR

Using the well-known fact of the IADE\_CG Red Black, the parallel algorithm for IADE\_CG RBSOR takes the form similar to IADE\_CG Red Black but the acceleration parameter  $\omega$  was chosen to provide the most rapid convergence.

### 4. IADE\_CG Multicoloring

By the definition of multidomain, domain  $\Omega$  is decomposed into  $w$  different groups. IADE\_CG Multicoloring is an advanced concept of IADE\_CG Red Black. If  $w = 2$ , then IADE\_CG Multicoloring is equaled to IADE\_CG Red Black.

The Domains for colors  $1, 2, 3, \dots, w$  are noted as  $\Omega^{w_1}, \Omega^{w_2}, \dots, \Omega^{w_w}$ . The subdomain  $\Omega^{w_i}$  is distributed into different groups of grid  $W_{i1}, W_{i2}, \dots, W_{i\frac{m}{wp}}$ , where  $i = 1, 2, \dots, w$ . In the process of assignment,  $W_{ij}, i = 1, 2, \dots, w$  and  $j = 1, 2, \dots, \frac{m}{wp}$  are mapped into the processors  $p$  in the alternating way.

At each time step, the computational grid for domain  $\Omega$  started its execution with level  $\Omega^{w_1}$ , followed by level  $\Omega^{w_2}$  and ends with level  $\Omega^{w_w}$ . The next step, exploits the CG algorithm as in equation (10)( $i - iv$ )

### 5. IADE\_CG vector

On the vector iteration strategy, the parallel IADE\_CG is run in two convergence sections. The first section is at the  $(k + \frac{1}{2})$  time step and the second section is the  $(k + 1)$  time step. This method converges if the inner convergence criterion is achieved for each section. The inner convergence criterions  $\varepsilon^{(k+\frac{1}{2})}$  and  $\varepsilon^{(k+1)}$  are definite global convergence criterion  $\varepsilon$ .

The CG algorithm as in equation (10)( $i - iv$ ) is attained before the examination of global convergence criterion.

### 6. IADE\_CG Incomplete Block LU

On the strategy of Incomplete Block LU preconditioners on slightly overlapping subdomains, the domain  $\Omega$  is decomposed into  $p$  processors with incomplete subdomain  $\bar{\Omega}$ . This strategy used the incomplete factorization with parameter  $\beta$  and  $\delta$  of algebraic boundary condition as follows,

i. at the  $(k + \frac{1}{2})^{th}$  iterate

$$k_{i-2}u_{i-2}^Q + \beta k_{i-2}u_{i-1}^Q - \delta k_{i-2}u_i^Q + (1 + \beta)k_{i-2}u_{i-1}^Q + \delta k_{i-2}u_{i+1}^Q = f_{i-1}$$

ii. at the  $(k + 1)^{th}$  iterate

$$d_{i-1}u_{i-1}^R + h_{i-1}u_i^R = f_{i-1}$$

The CG algorithm as in equation (10) ( $i - iv$ ) is attained after the execution of IADE\_CG Incomplete Block LU.

## 5.0 Numerical experiments

The problem statement under consideration is based on one dimensional heat conductor equation.(Smith (1979))

$$\frac{\partial U}{\partial t} = \frac{\partial^2 U}{\partial x^2}, \quad 0 \leq x \leq 1, \quad 0 < t, \quad (13)$$

subject to the initial condition,  $U(x, 0) = \sin(\pi, x)$ ,  $0 \leq x \leq 1$   
and the boundary condition,  $U(0, t) = U(1, t) = 0$ ,  $0 < t \leq 1$

## 6.0 Experiment Result And Parallel Performances

This paper presents the numerical properties of the parallel solver on the homogeneous architecture of 12 PC systems with Linux operating, Intel Pentium IV processors, speedup 20GB HDD and connected with internal network Intel 10/100 NIC. The following definitions are used to measure the parallel algorithms in term of speedup, efficiency and effectiveness.

$$\text{Speedup ratio } S_p = \frac{T_1}{T_p} \quad (14)$$

$$\text{Efficiency } C_p = \frac{S_p}{p} \quad (15)$$

$$\text{Effectiveness } F_p = \frac{S_p}{C_p} \quad (16)$$

where  $C_p = pT_p$ ,  $T_1$  is the execution time on a serial machine and  $T_p$  is the computing time on a parallel machine with  $p$  processors. Gauss Seidel Red Black is chosen as the control scheme. The number of iterations, the *rmse*, the maximum *rmse*, the maximum error and the execution time cost in sequential algorithms are shown in Figure 1. As shown in the figure 1, the sequential IADE\_CG, AGE\_CG and IADE\_CG Michell-Fairweather take the same number of relaxations to reach the convergence criterion  $\varepsilon$  with the minimum execution time. With domain decomposition, the implementation of parallel strategies, IADE\_CG Incomplete block LU, IADE\_CG Red Black, IADE\_CG Red Black SOR, IADE\_CG Multicoloring and IADE\_CG vector are increased in the number of relaxations and the execution time. The convergence rates, the *rmse*, the maximum *rmse* and the maximum error of all the sequential algorithms are slightly similar.

Figure 2 shows the execution time vs. number of processors. As expected, the computation time decreases with the increasing  $p$ . The 3 categories of the parallel algorithms indicated the best execution time in the following arrangement,

1. AGE\_CG and IADE\_CG Incomplete Block,
2. IADE\_CG Red Black SOR and IADE\_CG Red Black,
3. IADE\_CG Multicoloring and IADE\_CG Vector

Using this control scheme, IADE\_CG Michell-Fairweather and IADE\_CG vector are not suitable to be implemented as the parallel algorithms scheme for IADE\_CG. It involves high computational cost and communication cost. All the parallel algorithms for these categories have a

better execution time compared to their sequential algorithms as the number of processors are increased.

Figure 3 shows the speedup plotted against the number of processors  $p$ . As observed from the experiment, a nice speedup can be obtained for all applications with 12 processors except for IADE.CG Michell-Fairweather. The reductions in execution time often becomes smaller when a large number of processors is used. This phenomenon as stated in Amdahl's law, indicates that the number of processors increases, the communication cost (e.g., the latency for message passing) and the cost for global operations will eventually become dominant over local computation cost after a certain stage.

The IADE.CG Multicoloring, IADE.CG Red Black and IADE.CG Red Black SOR are good in terms of speedup and efficiency where data decomposition is run asynchronously and concurrently at every time step. AGE.CG and IADE.CG Incomplete Block allow for inconsistencies due to load balancing when the extra computation cost is needed for boundary condition. From equation (14) to equation (16),

$$F_p = \frac{S_p}{pT_p} = \frac{E_p}{T_p} = \frac{E_p S_p}{T_1}$$

which shows that  $F_p$  measure both speedup and efficiency. Therefore, a parallel algorithm is said to be effective when it maximizes  $F_p$  hence,  $F_p T_1 (= S_p E_p)$ . The optimal performance of the effectiveness could be obtained when more than 12 processors are used. Figure 4 shows the best effectiveness in the following order,

1. IADE.CG Incomplete Block,
2. AGE.CG
3. IADE.CG Red Black
4. IADE.CG Red Black SOR
5. IADE.CG Vector
6. IADE.CG Multicoloring

## 7.0 Conclusion

Conjugate Gradient (CG) can be regarded as a very efficient technique to increase the convergent and efficiency of these two iteration methods. The stable and high accuracy of IADE.CG Incomplete Block is found to be an alternative parallel algorithm for IADE.CG on PVM. All the parallel algorithms maintained the accuracy of the sequential IADE.CG algorithms. Parallel AGE.CG is inherently explicit, the domain decomposition strategy is efficiently utilized, straightforward to implement on PVM. These parallel algorithms is available to be implemented on heterogeneous architecture cluster of workstations. The experiment proved that the communication and computing times affect the speedup ratio, efficiency and effectiveness for the number of processors  $p$ . Therefore, we reach the conclusion that these parallel algorithms gets better efficiencies when it is implemented on a cluster workstations for solving a large-scale problems.

## 8.0 Acknowledgment

We wish to express our gratitude and indebtedness to the Universiti Kebangsaan Malaysia, Universiti Tenaga Nasional and Malaysian Government for providing the moral and financial support under IRPA grant for the successful completion of this project.

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method	AGE	IADE	LU	RBSOR	RB	MULTI	MF	VECTOR	GSRB
me (ms)	49.743	40.2641	50.4327	124.5471	126.1464	205.9478	1896.65	261.0082	156.4432
erations	250	252	305	362	363	362	252	252	600
rmse	1.5921E-09	1.5921E-18	1.5921E-09	1.5921E-16	1.5921E-09	1.5921E-09	1.5921E-18	1.5921E-09	1.5921E-09
iax_error	1.1102E-16	3.3307E-16	4.4409E-16	2.2204E-16	3.7921E-16	4.2306E-16	3.3307E-16	3.3307E-16	1.11E-16
qr_error	1.9846E-07	1.9793E-07	1.9846E-07	1.9846E-07	1.9846E-07	1.9846E-07	1.9793E-07	1.9846E-07	1.9846E-07
ve_rmse	5.3374E-17	5.3374E-17	5.3374E-17	5.3374E-17	5.3374E-17	5.3374E-17	5.3374E-17	5.3374E-17	5.3374E-17
m	720003	720003	720003	720003	720003	720003	720003	720003	720003
dx	1.3889E-06	1.3889E-06	1.3889E-06	1.3889E-06	1.3889E-06	1.3889E-06	1.3889E-06	1.3889E-06	1.3889E-06
dt	9.6450E-13	9.6450E-13	9.6450E-13	9.6450E-13	9.6450E-13	9.6450E-13	9.6450E-13	9.6450E-13	9.6450E-13
tt	4.8225E-11	4.8225E-11	4.8225E-11	4.8225E-11	4.8225E-11	4.8225E-11	4.8225E-11	4.8225E-11	4.8225E-11
lamda	0.5	0.5	0.5	0.5	0.5	0.5	0.5	0.5	0.5
teta	1	1	1	1	1	1	1	1	1
round	50	50	50	50	50	50	50	50	50
r	0.8	0.8	1.1	1.3	1.29	0.98	0.8	0.9	-
ry	-	-	-	1.1	1	0.92	-	-	-
rz	-	-	-	1	1	1.09	-	-	-
tol	1.0000E-15	1.0000E-15	1.0000E-15	1.0000E-15	1.0000E-15	1.0000E-15	1.0000E-15	1.0000E-15	1.0000E-15
tol_y	-	-	-	-	-	-	-	1.0000E-15	-
tol_z	-	-	-	-	-	-	-	1.0000E-15	-

1038+890

Figure 1 Sequential algorithms for IADE

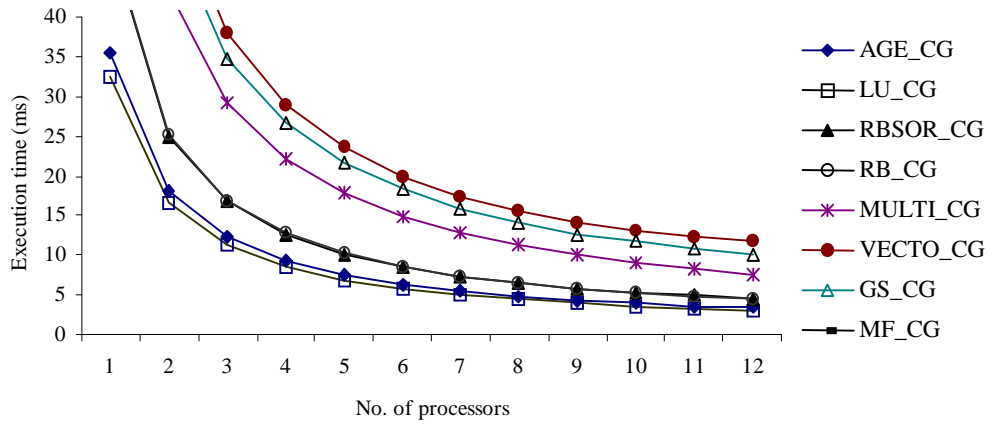


Figure 2 The execution time vs. number of processors

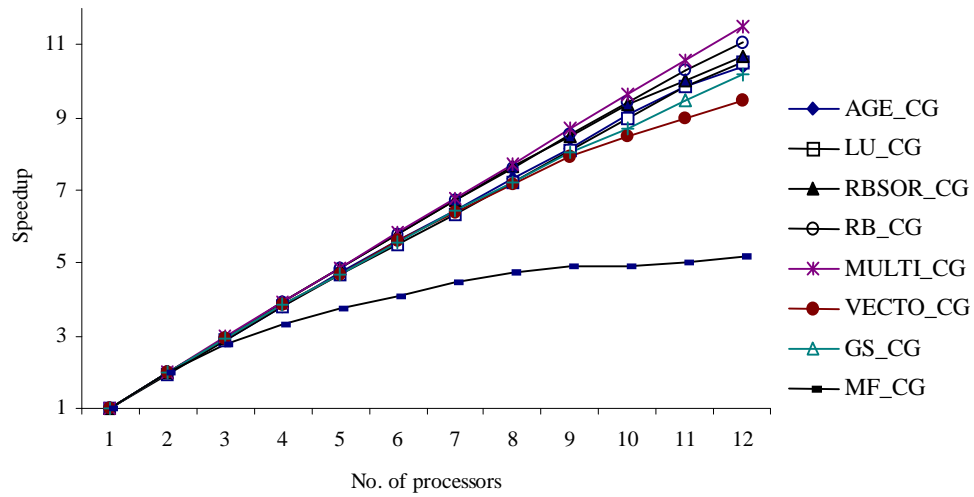


Figure 3 The speedup vs. number of processors

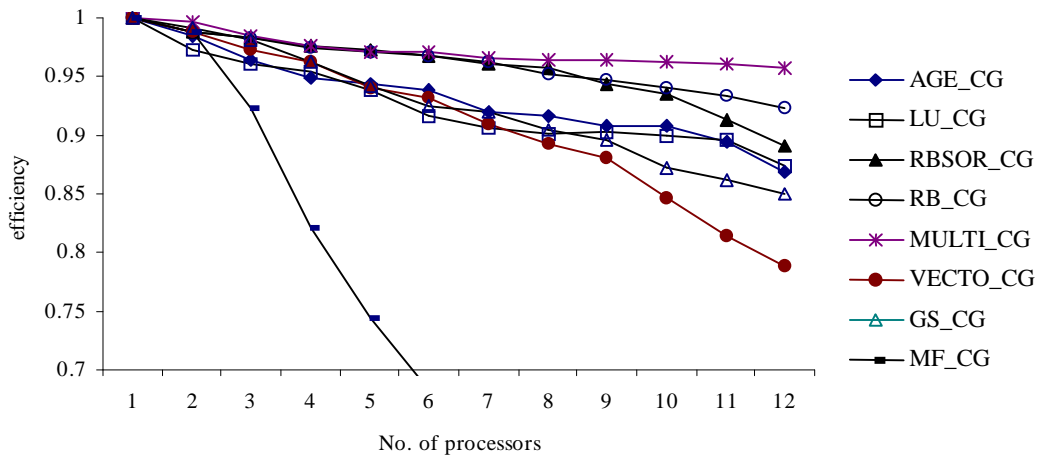


Figure 4 The efficiency vs. number of processors

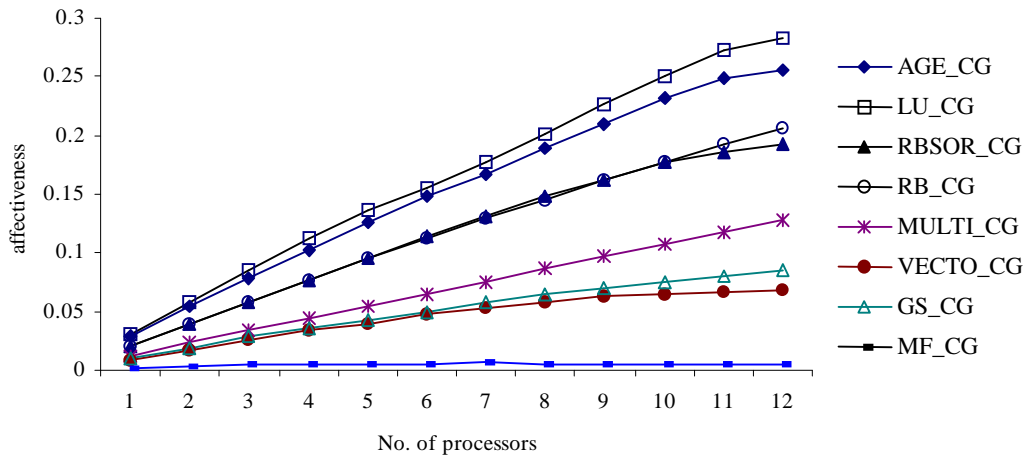


Figure 5 The effectiveness vs. number of processors