



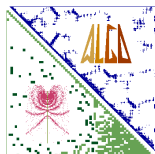
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## Preconditioners based on Strong Components

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## Abstract

This paper proposes an approach for obtaining block diagonal and block triangular preconditioners that can be used for solving a linear system  $\mathbf{Ax} = \mathbf{b}$ , where  $\mathbf{A}$  is a large, nonsingular, real,  $n \times n$  sparse matrix. The proposed approach uses Tarjan's algorithm for hierarchically decomposing a digraph into its strong subgraphs [22, 23]. To the best of our knowledge, this is the first work that uses this algorithm of Tarjan for preconditioning purposes. We describe the method, analyse its performance, and compare it with preconditioners from the literature such as ILUT [18, 19] and XPABLO [12, 13] and show that it is better than XPABLO and competitive with ILUT for many matrices with the advantage that a version of our preconditioner is fully parallelizable.

## 1 Introduction

Given a linear system

$$\mathbf{Ax} = \mathbf{b}, \tag{1}$$

where  $\mathbf{A}$  is a real, large, sparse square matrix, we propose a method to construct a preconditioning matrix  $\mathbf{M}$  to accelerate the solution of the system when using Krylov methods. The proposed method is based on a hierarchical decomposition of the associated digraph into its strong subgraphs. This decomposition can be used to find a permutation of the rows and columns of the original matrix  $\mathbf{A}$  and a block structure where the maximum size of a diagonal block is smaller than a desired value. After that, by dropping a set of nonzeros that are not in the block diagonal, we can generate a preconditioning matrix  $\mathbf{M}$ . This set of nonzeros depends on the desired structure of the preconditioning matrix: if the desired structure is block diagonal, the set contains all nonzeros which are not in any of the blocks on the diagonal. Note that a block diagonal preconditioning matrix is very suitable for parallel computation. Although less parallelizable, a block triangular structure may also be suitable as a preconditioner since, if  $\mathbf{M}$  is block triangular, it is relatively easy to solve a linear system with  $\mathbf{M}$  as the coefficient matrix. If this is the desired structure, our proposed method first applies an additional block permutation and then drops the nonzeros which are not in the block upper-triangular part.

The algorithm we use to create the blocks on the diagonal of  $\mathbf{M}$  is a modified version of Tarjan's algorithm **HD** that decomposes a digraph into its strong subgraphs hierarchically [23]. Tarjan assumed that the edges of the digraph are weighted and **HD** uses this weight information to create the hierarchical decomposition. However, **HD** requires distinct edge weights if it is implemented as given in [23]. In this paper, we propose a slight modification of **HD** which allows us to handle digraphs whose edge weights are not necessarily distinct. We made further modifications to the algorithm to use it for preconditioning purposes. The strong subgraphs formed by the modified version of **HD** correspond to the blocks on the diagonal of  $\mathbf{M}$ . To the best of our knowledge, this is the first work that uses Tarjan's hierarchical decomposition algorithm for preconditioning purposes. We call our modified version **HDPRE**.

We should emphasize at this point that this algorithm of Tarjan is different from the much better known algorithm for obtaining the strong components of a reducible matrix. This earlier algorithm [21], which we call **SCC**, is used widely in the solution of reducible systems and is also called by **HD** and **HDPRE** which in a sense extend the earlier work to irreducible matrices.

We have conducted several experiments to see the efficiency of the proposed algorithm. We compare the number of iterations for convergence and the memory requirement of the **GMRES** [20] iterative solver when the proposed approach and a set of **ILUT** preconditioners [18, 19] are used. We are aware that block based preconditioning techniques have been studied before and successful preconditioners such as **PABLO** and its derivatives have been proposed [12, 13]. These preconditioners were successfully used for several matrices [3, 5, 9]. In this paper, we compare our results also with **XPABLO** [12, 13].

For the experiments, we used several circuit simulation, device simulation and electromagnetics matrices from the University of Florida Sparse Matrix Collection (<http://www.cise.ufl.edu/research/sparse/matrices/>). Experimental results show that the performance of the proposed algorithm, called **SCPRE**, is better than **ILUT** and **XPABLO** preconditioners especially for device and circuit simulation matrices. We also show that when both **XPABLO** and **SCPRE** are used to generate parallel preconditioners, **SCPRE** again outperforms **XPABLO**.

Section 2 gives the notation used in the paper and background on Tarjan's algorithm **HD**. The proposed algorithm is described in Section 3 and the implementation details are given in Section 4. Section 5 gives the experimental results and Section 6 concludes the paper.

## 2 Background

Let  $\mathbf{A}$  be a large, nonsingular,  $n \times n$  sparse matrix with  $m$  nonzeros. The digraph  $G = (V, E)$ , associated with  $\mathbf{A}$ , has  $n$  vertices in its vertex set  $V$  where  $v_i$  corresponds to the  $i$ th row/column of  $\mathbf{A}$  for  $1 \leq i \leq n$  and  $v_i v_j$  is in the edge set  $E$  iff  $\mathbf{A}_{ij}$  is nonzero, for  $1 \leq i \neq j \leq n$ . Figure 1(b) shows a simple  $6 \times 6$  matrix with 13 nonzeros and its associated digraph.

A *path* is sequence of vertices such that there exists an edge between every two consecutive vertices. A path is called *closed* if its first and last vertex are the same. A vertex  $u \in V$  is *connected* to  $v \in V$  if there is path from  $u$  to  $v$  in  $G$ . A directed graph  $G$  is *strongly connected* if  $u$  is connected to  $v$  for all  $u, v \in V$ . Note that a digraph with a single vertex  $u$  is strongly

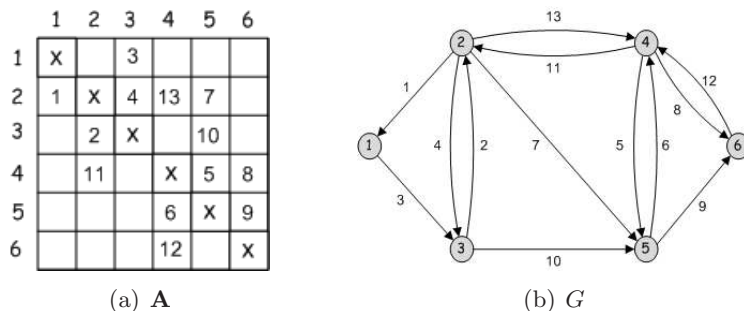


Figure 1: A  $6 \times 6$  matrix **A** with 13 nonzeros on the left and its associated digraph  $G$  on the right. The nonzeros on the diagonal of **A** are shown with  $\times$ . Except for these entries, there is an edge in the associated digraph  $G$  for each nonzero of **A**.

connected. A digraph  $G' = (V', E')$  is a subgraph of  $G$  if  $V' \subset V$  and  $E' \subseteq E \cap (V' \times V')$ . Furthermore, if  $G'$  is maximally strongly connected, i.e., if there is no strongly connected subgraph  $G''$  of  $G$  such that  $G'$  is a subgraph of  $G''$ , it is called a strong component (or a strongly connected component) of  $G$ .

Let  $G = (V, E)$  be a digraph and  $\mathcal{P}(V) = \{V_1, V_2, \dots, V_k\}$  define a partition of  $V$  into disjoint sets, i.e.,  $V_i \cap V_j = \emptyset$  for  $i \neq j$  and  $\cup_{i=1}^k V_i = V$ . Let  $\mathcal{V} = \{\mathcal{V}_1, \mathcal{V}_2\}$  be a set of two vertex partitions such that  $\mathcal{V}_1 = \mathcal{P}(V)$  and

$$\mathcal{V}_2 = \bigcup_{V_i \in \mathcal{V}_1} \mathcal{P}(V_i),$$

i.e.,  $\mathcal{V}_2$  is a finer partition obtained from partitioning the parts in  $\mathcal{V}_1$ . Hence, if  $\mathcal{V}_1 = \{\{1, 2, 3\}, \{4, 5, 6\}\}$  then  $\mathcal{V}_2$  can be  $\{\{1\}, \{2, 3\}, \{4, 5\}, \{6\}\}$  but cannot be  $\{\{1, 2\}, \{3, 4\}$

$\}, \{5, 6\}\}$ . Let  $\text{no}_1(v)$  and  $\text{no}_2(v)$  denote the index of the part containing vertex  $v \in V$  for  $\mathcal{V}_1$  and  $\mathcal{V}_2$ , respectively.

Let **condense** be an operation which takes  $G$  and  $\mathcal{V}$  as inputs and returns a condensed digraph  $\text{condense}(G, \mathcal{V}) = G^\mathcal{V} = (V^\mathcal{V}_2, E^\mathcal{V}_1)$  where each vertex set  $V_i \in \mathcal{V}_2$  is condensed into a single vertex  $\nu_i \in V^\mathcal{V}_2$ . For all  $uv \in E$ , with  $\text{no}_2(u) = i$  and  $\text{no}_2(v) = j$  there exists  $\nu_i \nu_j \in E^\mathcal{V}_1$  if and only if  $\text{no}_1(u) \neq \text{no}_1(v)$ , i.e.,  $u$  and  $v$  are in different coarse parts. Note that even though  $G$  is a simple digraph,  $G^\mathcal{V}$  can be a directed multigraph, i.e., there can be multiple edges between two vertices. The definitions of connectivity and strong connectivity in directed multigraphs are the same as those in digraphs. An example for the **condense** operation is given in Figure 2.

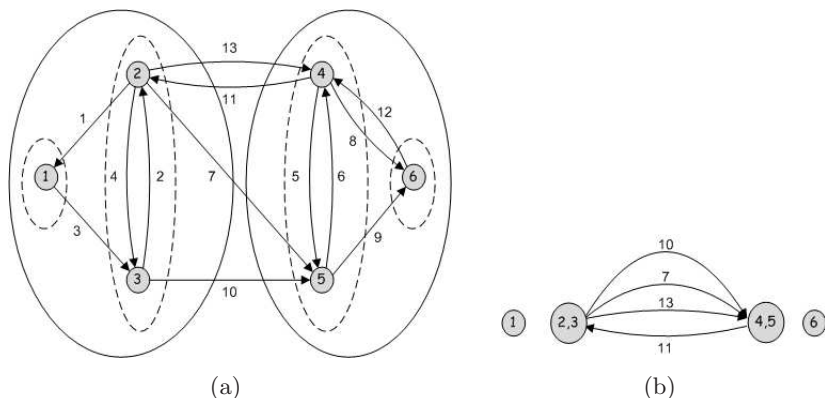


Figure 2: An example for the `condense` operation on the digraph in Figure 1(b). The partitions  $\mathcal{V}_1 = \{\{1, 2, 3\}, \{4, 5, 6\}\}$  and  $\mathcal{V}_2 = \{\{1\}, \{2, 3\}, \{4, 5\}, \{6\}\}$  are shown in 2(a). The condensed graph is shown in 2(b).

### 3 A Strong Subgraph based Preconditioner

Our proposed algorithm, `SCPRES`, generates a preconditioner  $\mathbf{M}$  with a block diagonal or block upper-triangular structure where the size of each block is smaller than a requested maximum block size  $mbs$ . For the experiments, we scale and permute  $\mathbf{A}$  from (1) by Duff and Koster's `MC64` with the option that uses the *maximum product transversal* [10]. The idea used by `MC64` is due to Olschowska and Neumaier [17] who propose an algorithm which permutes and scales the matrix in such a way that the magnitudes of the diagonal entries are one and the magnitudes of the off-diagonal entries are all less than or equal to one. Such a matrix is called an *I-matrix*. For direct methods, it has been observed that the more dominant the diagonal of a matrix, the higher the chance that diagonal entries are stable enough to serve as pivots for elimination. For iterative methods, as previous experiments have shown, such a scaling is also of interest [4, 10]. Additionally, in our experiments, we observed that permuting the matrix with respect to the maximum product transversal is the most promising one for the preconditioned iterative solver in terms of the iteration count. From now on, we will assume that the diagonal of  $\mathbf{A}$  is nonzero since this is the case after scaling and permutation.

In the first phase, with a symmetric permutation of the rows/columns, `SCPRES` finds the diagonal blocks of the preconditioner  $\mathbf{M}$  by using the algorithm `HDPRES` which is a modified version of Tarjan's algorithm `HD` which was

originally designed to hierarchically decompose a digraph into its strong subgraphs [23]. We then combine some of these blocks if the combination has fewer than  $mbs$  rows/columns and the combination is not block diagonal. The diagonal blocks of the resulting matrix can then be used to precondition the iterative solver using block Jacobi and can exploit parallel architectures as the blocks are independent. If a block diagonal structure is desired, SCPRE terminates. Otherwise, in the additional second phase, SCPRE extends the nonzero set contained in  $\mathbf{M}$  to possibly reduce the number of iterations required for convergence. To extend the preconditioner and put more nonzeros in it, SCPRE permutes the blocks obtained in the first phase and outputs the block upper-triangular part as the preconditioner  $\mathbf{M}$ .

If  $\mathbf{A}$  cannot be permuted into a block triangular form (BTF)[11] by simultaneous row/column permutations we say that it is irreducible. Otherwise, we call it reducible. If  $\mathbf{A}$  is reducible and the maximum block size in the BTF of  $\mathbf{A}$  is not larger than  $mbs$ , then the first phase of SCPRE just obtains the diagonal blocks of the BTF. The idea of permuting a matrix into its block triangular form is a well known technique which has been recently and successfully used by direct and iterative solvers for circuit simulation matrices [8, 24], which they state can often be permuted to a non-trivial BTF. For some applications, such as DC operating point analysis, the block triangular form has many but small blocks [24]. Such a matrix is usually easy to factorize if we initially permute it to BTF, so that a direct solver like KLU [8] only needs to factorize the diagonal blocks. However, even if the matrix is reducible, its BTF may sometimes not be sufficient since for applications like transient simulation or for circuit matrices with feedbacks, we may have one or more very large blocks. In this case HDPRE used in the first phase of SCPRE, decomposes this block by using the ideas developed for Tarjan’s HD algorithm [23]. In our experiments, we only use matrices that are irreducible or have a large block in their BTF. The details of SCPRE and the algorithms it uses are given in the next section. Note that since we use a combinatorial algorithm from graph theory for preconditioning purposes, while describing the phases, the terms *row/column* and *vertex* can be used interchangeably, as well as the terms *nonzero* and *edge*.

### 3.1 SCPRE: Obtaining the Block Diagonal Preconditioner

To obtain a block diagonal preconditioner, SCPRE uses HDPRE and then combines some of these blocks if the size of the combined block is at most  $mbs$  and the combined block is not block diagonal. In this section, we give the details of these algorithms. First, we describe Tarjan’s hierarchical decom-

position algorithm in detail.

### 3.1.1 Tarjan's Algorithm for Hierarchical Clustering

The digraph  $G = (V, E)$ , associated with  $\mathbf{A}$ , has  $n$  vertices in  $V$  where  $v_i$  corresponds to the  $i$ th row/column of  $\mathbf{A}$  for  $1 \leq i \leq n$  and  $v_i v_j \in E$  iff  $\mathbf{A}_{ij}$  is nonzero, for  $1 \leq i \neq j \leq n$ . The weight of an edge is set to the absolute value of the corresponding nonzero. Hence, there are  $m$  edges and all of the edges have positive weights. A hierarchical decomposition of  $G$  into its strong subgraphs can be defined in the following way. Let  $\sigma_0$  be a permutation of the edges. For  $1 \leq i \leq m$ , let  $\sigma_0(i)$  be the  $i$ th edge in  $\sigma_0$  and  $\sigma_0^{-1}(uv)$  be the index of the edge  $uv$  in the permutation for all  $uv \in E$ . Let  $G_0 = (V, \emptyset)$  be the graph obtained by removing all the edges from  $G$ . Consider that edges are added one by one to  $G_0$  in the order determined by  $\sigma_0$ . Let  $G_i = (V, \{\sigma(j) : 1 \leq j \leq i\})$  be the digraph obtained after the addition of the first  $i$  edges. Initially in  $G_0$ , there are  $n$  strong components, one for each vertex, and during the edge addition process, the strong components gradually coalesce until there is only one, as we are assuming that  $\mathbf{A}$  is irreducible. Note that if this is not the case, the algorithm will be used for the large irreducible blocks in  $\mathbf{A}$ . The hierarchical decomposition of  $G$  into its strong subgraphs with respect to the edge permutation  $\sigma_0$  shows which strong components are formed in this process hierarchically. Note that a strong component formed in this edge addition process is indeed a strong component of some digraph  $G_i$  but not of  $G$ . For  $G$  all except the last are just strong subgraphs.

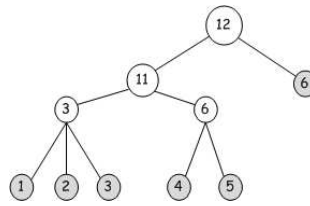


Figure 3: The hierarchical decomposition tree for the digraph  $G$  and the permutation in Figure 1(b).

A hierarchical decomposition can be represented by a hierarchical decomposition tree  $T$ , whose leaf nodes correspond to the vertices in  $V$ , non-leaf nodes correspond to the edges in  $E$ , and subtrees correspond to the decomposition trees of the strong components that form as the process proceeds. Note that only the edges that create strong components during the process

have corresponding internal nodes in  $T$ . If  $\sigma_0$  is the ordering determined by the edge numbers, the hierarchical decomposition tree for the digraph in Figure 1(b) is given in Figure 3. As the figure shows, during the edge addition process, after the addition of the 3rd and 6th edges in  $\sigma_0$ , the sets of vertices  $\{1, 2, 3\}$  and  $\{4, 5\}$  form a strong component of  $G_3$  and  $G_6$ , respectively. These strong components are then combined and form a larger one after the addition of the 11th edge. In Figure 3, the root of the tree is labelled with 12. Hence the first 12 edges in  $\sigma_0$  are sufficient to construct a strongly connected digraph. For the figures in this paper, we use the labels of the corresponding vertices and the  $\sigma_0^{-1}$  values of the corresponding edges to label each leaf and non-leaf node of a hierarchical decomposition tree, respectively.

Given a digraph  $G = (V, E)$  with  $n$  vertices and  $m$  edges, and a permutation  $\sigma_0$ , the hierarchical decomposition tree  $T$  can be obtained by first constructing  $G_0$  and executing SCC for each internal digraph  $G_i$  obtained during the edge addition process. Note that this is an  $\mathcal{O}(mn + m^2)$  algorithm since  $1 \leq i \leq m$  and the cost of SCC is  $\mathcal{O}(n + m)$  due to another algorithm by Tarjan [21]. To obtain  $T$  in a more efficient way, Tarjan first proposed an  $\mathcal{O}(m \log^2 n)$  recursive algorithm [22] and later improved his algorithm and reduced the complexity to  $\mathcal{O}(m \log n)$  [23]. He assumed that the weights of the edges in the digraph are distinct, i.e.,  $w(uv) \neq w(u'v')$  for two distinct edges  $uv$  and  $u'v'$ . Here we modify the description of the algorithm slightly so that it also works for the case when some edges have equal weights. Note that the edge weights do not play a role in the connectivity of the digraph. In Tarjan's algorithm HD, they are used in a preprocessing step which defines a permutation  $\sigma_0$  of the edges. In addition, they are also used for comparison purposes during the course of the algorithm. We eliminate the necessity of the latter by using the indices of the edges with respect to  $\sigma_0$  for the comparisons. With this slight modification, the algorithm remains correct even when some edges have the same weight.

HD uses a recursive approach and for every recursive call, it gets a digraph  $G = (V, E)$ , a permutation  $\sigma$  of the edges, and a parameter  $i$  as inputs such that  $G$  is strongly connected and  $G_i$  is known to be acyclic, i.e., every vertex is a separate strong component [23]. For the initial call,  $i$  is set to 0 and the initial permutation is set to  $\sigma_0$  which is a permutation of all the edges in the original digraph.

For a call of  $\text{HD}(G = (V, E), \sigma, i)$ , the *size* of the subproblem is set to  $|E| - i$ , the number of edges that remain to be investigated (Tarjan used the term *rank* to denote the size of a problem). Note that in the first step, HD knows that  $G_i$  is acyclic, i.e., there are  $|V|$  strong components of  $G_i$ , one for

each vertex. If the problem size is one, since  $G$  is strongly connected and  $G_i$  is acyclic, the vertices in  $V$  are combined with the addition of the  $|E|$ th edge in  $\sigma$ . Hence the algorithm HD returns a tree  $T$  having a root labelled with  $\sigma_0^{-1}(\sigma(|E|))$  and  $|V|$  leaves. If the problem size is not one, HD checks if  $G_j$ ,  $j = \lceil (i + |E|)/2 \rceil$  is strongly connected. If  $G_j$  is strongly connected, then all of the strong components should have been combined before the addition of the  $(j + 1)$ th edge. Hence the algorithm calls HD( $G_j, \sigma, i$ ). Otherwise, a recursive call is made for each strong component of size larger than one. A detailed pseudo-code of HD is given in Algorithm 1.

In Algorithm 1, for the  $\ell$ th strong component  $SC_\ell = (V_\ell, E_\ell)$ , lines 11–14 find the integer  $i_\ell$  such that the subgraph of  $SC_\ell$  containing only its first  $i_\ell$  edges is acyclic. Since  $G_i$ , the graph containing the first  $i$  edges of  $G$  in  $\sigma$ , is known to be acyclic, for  $SC_\ell$ ,  $i_\ell$  is set to the index of the last edge  $uv$  in  $E_\ell$  such that  $\sigma^{-1}(uv) \leq i$ . If no such edge exists, i.e., all the edges in  $E_\ell$  come after the  $i$ th edge in  $\sigma$ ,  $i_\ell$  is set to 0. Since  $G_j$  has more than one strong component and  $G$  is known to be strongly connected, with the addition of some edge(s) after the  $j$ th one at least two strong components of  $G_j$  should have been combined. To find this edge, another recursive call, HD( $G^\mathcal{V}, \sigma^\mathcal{V}, i^\mathcal{V}$ ), is made for the condensed graph  $G^\mathcal{V}$ . The  $i^\mathcal{V}$  value is set in a similar fashion to  $i_\ell$  as described above. But this time instead of  $i$ , we use  $j$  since we know that the graph  $G_j^\mathcal{V} = (V^\mathcal{V}, \{\sigma^\mathcal{V}(k) : 1 \leq k \leq j\})$  is acyclic.

At line 6 of Algorithm 1, the size of the problem becomes at most  $j - i$  and for lines 15 and 26, there will be smaller subproblems with size at most  $j - i$  and  $|E| - j$ , respectively. By definition of  $j$ , every subproblem has a size at most  $\frac{2}{3}$  of the original problem size (consider the case when  $i = 0$  and  $|E| = 3$ ). Note that every edge in the original problem corresponds to an edge in at most one subproblem and, if we do not count the recursive calls, the rest of the algorithm takes  $\mathcal{O}(|E|)$ . Let  $|E| = m$ ,  $\mathfrak{t}(m, r)$  be the total complexity of a problem with  $m$  edges and  $r$  problem size, and  $k$  be the number of recursive calls. Then

$$\mathfrak{t}(m, r) = \mathcal{O}(m) + \sum_{i=1}^k \mathfrak{t}(m_i, r_i).$$

Since  $\sum_{i=1}^k m_i \leq m$  and  $r_i \leq 2r/3$  for  $1 \leq i \leq k$ , an easy induction shows that  $\mathfrak{t}(m, r) = \mathcal{O}(m \log r)$ . Hence the total complexity of the algorithm is  $\mathcal{O}(m \log m)$  which is actually  $\mathcal{O}(m \log n)$  since the original graph is a simple digraph (not a directed multigraph).

Let us sketch the algorithm for the digraph  $G = (V, E)$  in Figure 1(b). Assume that  $\sigma_0$  is the ordering described in the figure. In the initial call,

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**Algorithm 1**  $T = \text{HD}(G = (V, E), \sigma, i)$  . For the initial call,  $\sigma = \sigma_0$  and  $i = 0$ .

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1: if  $|E| - i = 1$  then
2:   Let  $T$  be a tree with  $V$  leaves. Root is labelled with  $\sigma_0^{-1}(\sigma(|E|))$ 
3:   return  $T$ 
4:  $j = \lceil (i + |E|)/2 \rceil$ 
5: if  $G_j = (V, \{\sigma(k) : 1 \leq k \leq j\})$  is strongly connected then
6:   return  $T = \text{HD}(G_j, \sigma, i)$ 
7: else
8:   for each strong component  $SC_\ell = (V_\ell, E_\ell)$  of  $G_j$  do
9:     if  $|V_\ell| > 1$  then
10:       $\sigma_\ell =$  the permutation of  $E_\ell$  ordered with respect to  $\sigma$ 
11:      if  $i = 0$  or  $(\sigma^{-1}(uv) > i, \forall uv \in E_\ell)$  then
12:         $i_\ell = 0$ 
13:      else
14:         $i_\ell = \max\{k : \sigma^{-1}(\sigma_\ell(k)) \leq i\}$ 
15:         $T_\ell = \text{HD}(SC_\ell, \sigma_\ell, i_\ell)$ 
16:      else
17:         $T_\ell = (V_\ell, \emptyset)$ 
18:       $\mathcal{V}_1 = \mathcal{V}_2 = \{V_\ell : SC_\ell \text{ is a strong component of } G_j\}$ 
19:       $\mathcal{V} = \{\mathcal{V}_1, \mathcal{V}_2\}$ 
20:       $G^\mathcal{V} = \text{condense}(G, \mathcal{V}) = (V^{\mathcal{V}_2}, E^{\mathcal{V}_1})$ 
21:       $\sigma^\mathcal{V} =$  the permutation of  $E^{\mathcal{V}_1}$  ordered with respect to  $\sigma$ 
22:      if  $(\sigma^{-1}(uv) > j, \forall uv \in E^{\mathcal{V}_1})$  then
23:         $i^\mathcal{V} = 0$ 
24:      else
25:         $i^\mathcal{V} = \max\{k : \sigma^{-1}(\sigma^\mathcal{V}(k)) \leq j\}$ 
26:         $T^\mathcal{V} = \text{HD}(G^\mathcal{V}, \sigma^\mathcal{V}, i^\mathcal{V})$ 
27:      replace the leaves of  $T^\mathcal{V}$  with the corresponding trees  $T_\ell$ 
28:      return  $T^\mathcal{V}$ 

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step 4 of Algorithm 1 computes  $j = 7$  and checks if  $G_7$  is strongly connected. As Figure 4 shows  $G_7$  has three strong components where the first and second are the new subproblems solved recursively. Since the third strong component contains only one vertex, HD does not make a recursive call for it. An additional recursive call is made for the condensed graph. Figure 5 shows the graphs for the recursive calls and the returned trees. The number of edges in Figs. 5(a), 5(b) and 5(c) are 4, 2 and 7, whereas the corresponding problem sizes are 4, 2 and 6 respectively. Note that  $i_1$  and  $i_2$  are 0 for the first two calls and  $i^{\mathcal{V}} = 1$  for the last one with  $G^{\mathcal{V}}$  since  $G_1^{\mathcal{V}}$  is known to be acyclic because  $j = 7$  and  $\sigma^{\mathcal{V}}(1) = \sigma(7)$ .

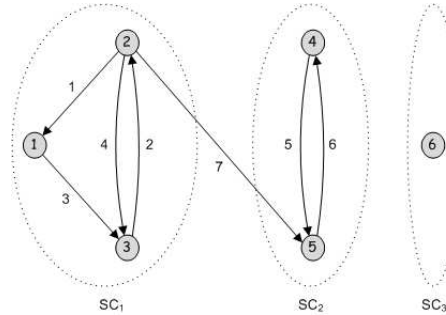


Figure 4: Strong components of  $G_7$  for the digraph  $G$  given in Figure 1(b)

Because of the multiple edges between two vertices, the condensed graph in Figure 5(c) has 7 edges. However, the algorithm still works if we sparsify the edges of  $G^{\mathcal{V}} = (V^{\mathcal{V}_2}, E^{\mathcal{V}_1})$  and obtain a simple digraph as follows: For a  $uv \in E$  such that  $u \in V_i$  and  $v \in V_j$  and  $i \neq j$ , there exists  $v_i v_j \in E^{\mathcal{V}_1}$  if no other  $u'v' \in E$  exists such that  $u' \in V_i$  and  $v' \in V_j$  and  $\sigma^{-1}(u'v') < \sigma^{-1}(uv)$ . That is, for multiple edges between  $u$  and  $v$ , we delete all but the first one in the permutation  $\sigma$ . In Figure 5(c), these edges,  $\sigma(7)$  and  $\sigma(8)$ , are shown in bold. In [23], Tarjan states that although having less edges in the condensed graphs with this modification is desirable, in practice the added simplicity does not compensate for the cost of the reduction of multigraphs to simple digraphs. This is also validated by our preliminary experiments.

### 3.1.2 HDPRE: Obtaining the Initial Block Structure

As mentioned in Section 3.1.1, Tarjan proposed HD for hierarchical clustering purposes and sorted the edges with respect to increasing edge weights. That is, for the permutation  $\sigma_0$  used for hierarchical clustering, if  $i < j$  then  $w(\sigma_0(i)) \leq w(\sigma_0(j))$ . In this work, we propose using two different

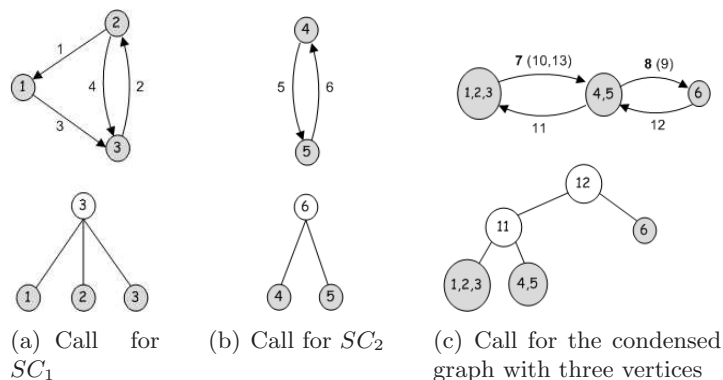


Figure 5: Three recursive calls for the digraph  $G$  and  $\sigma_0$  in Figure 1(b). Internal nodes in trees are labelled with the  $\sigma_0^{-1}$  value of the corresponding edge. Note that the overall hierarchical decomposition tree is already given in Figure 3.

approaches to obtain the permutation: the first one solely depends on the weights of the edges and sorts them in the order of decreasing edge weights, i.e., we define the permutation  $\sigma$  such as  $w(\sigma(i)) \geq w(\sigma(j))$  if  $i < j$ . The second one uses the sparsity pattern of the matrix by using the reverse Cuthill-McKee (RCM) ordering [6, 16] to find a symmetric row/column permutation and relabels the vertices of the digraph accordingly. After that the edges are ordered in a natural, row-wise order. That is, an edge  $ij$  always comes before  $k\ell$  if  $i < k$  or,  $i = k$  and  $j < \ell$ .

The output of Tarjan's HD algorithm, the decomposition tree  $T$ , could be used for preconditioning without modifying the algorithm, but we post-process the tree to ensure that all leaf nodes are as large as they can be but still have fewer than  $mbs$  nodes. For the decomposition tree  $T$  in Figure 3, the cases for  $mbs = 2$  and  $mbs = 3$  are given in Figure 6. In  $T$ , for the case  $mbs = 2$ , vertices 1, 2, and 3 cannot be combined since the number of vertices in the combined component will be 3, greater than  $mbs$ . Hence, there will be 5 blocks after this phase. However, such a combination is possible for the case  $mbs = 3$  and the number of blocks will be 3. Note that for preconditioning, we do not need to construct the whole tree of HD. We only need to continue hierarchically decomposing the blocks until they contain at most  $mbs$  vertices. Hence, for efficiency we modify line 9 of HD to check if the current strong component has more than  $mbs$  vertices (instead of one vertex). Hence the modified algorithm will make a recursive call for a strong component if and only if the component has more than  $mbs$  vertices.

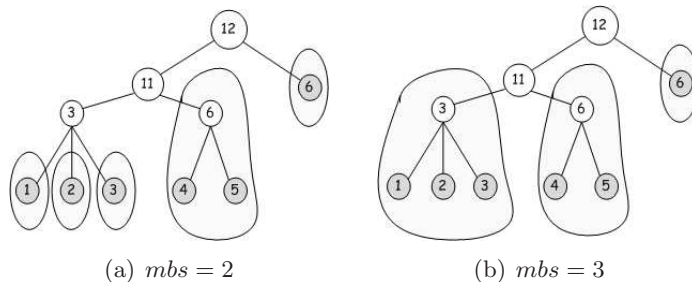


Figure 6: Using the output of HD algorithm. Two cases,  $mbs = 2$  and  $mbs = 3$ , are investigated for the decomposition tree in Figure 3.

To obtain denser and larger blocks, we incorporate some more modifications to HD as follows: first, we modify the definition of  $\mathcal{V}$ . Note that  $\mathcal{V} = \{\mathcal{V}_1, \mathcal{V}_2\}$  for HD, where the parts in  $\mathcal{V}_1 = \mathcal{V}_2$  are the vertex sets of the strong components of  $G_j$ . For preconditioning, we keep the definition of  $\mathcal{V}_1$  but we use a finer partition  $\mathcal{V}_2$  which contains the vertex sets of strong components obtained by hierarchically decomposing the strong components of size larger than  $mbs$ . For example, in Figure 4, we have 3 strong components of sizes 3, 2 and 1, respectively. Hence,  $\mathcal{V}_1 = \{\{1, 2, 3\}, \{4, 5\}, \{6\}\}$ . If  $mbs = 2$ ,  $SC_1$  will be further divided so  $\mathcal{V}_2 = \{\{1\}, \{2\}, \{3\}, \{4, 5\}, \{6\}\}$ . However, if  $mbs = 3$  no more decomposition will occur and  $\mathcal{V}_1$  will be equal to  $\mathcal{V}_2$ . With this modification, the algorithm will try to combine the smaller strong components and obtain larger ones with at most  $mbs$  vertices. Note that setting  $\mathcal{V} = \{\mathcal{V}_2, \mathcal{V}_2\}$  tries to do the same but it will fail since the only components that can be formed by this approach will be the same as those in  $\mathcal{V}_1$ . Hence, by deleting the edges within the vertex sets in  $\mathcal{V}_1$ , we eliminate the possibility of obtaining the same components.

A second modification is applied to the **condense** operation by deleting the edges between two vertices  $\nu_i, \nu_j \in V^{\mathcal{V}_1}$  in the condensed graph  $G^{\mathcal{V}}$ , if the total size of the corresponding parts  $V_i, V_j \in \mathcal{V}_2$  is larger than  $mbs$ . Note that if we were to retain these edges, they would only be used to form blocks of size more than  $mbs$ . We call this modified **condense** operation **pcondense**. An example of the difference between **condense** and **pcondense** is given in Figure 7.

As Figure 7 shows, with the last modification some of the graphs for the recursive calls may not be strongly connected. Hence, instead of a whole decomposition tree, we may obtain a forest such that each tree in the forest, which corresponds to a strong subgraph in the hierarchical decomposition,

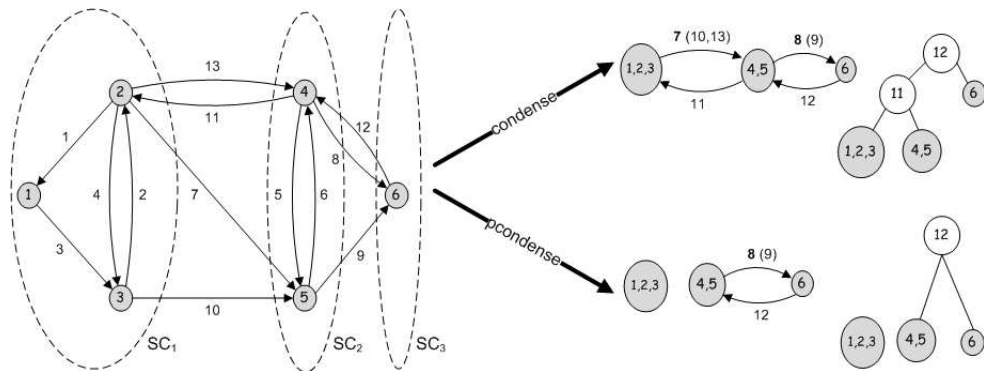


Figure 7: Difference between `condense` and `pcondense` operations for the strong components of  $G_7$  given in Figure 4. Let  $mbs = 3$  so all of the components have a desired number of vertices and  $\mathcal{V}_1 = \mathcal{V}_2 = \{\{1, 2, 3\}, \{4, 5\}, \{6\}\}$ . Note that the condensed graphs obtained by `condense` and `pcondense` are the same except that the latter does not have some of the edges that the former has. For this example, the edges 7, 10, 11 and 13 are missing since the total size size of  $SC_1$  and  $SC_2$  is 5, greater than  $mbs$ . As a result, for the `condense` graph, we obtain 3 blocks of sizes 3, 2 and 1, respectively, whereas for the `pcondense` graph, we have 2 blocks of size 3.

has less than  $mbs$  leaves. The modified algorithm **HDPRE**, described in Algorithm 2, also handles digraphs which are not strongly connected. Note that, for preconditioning, the only information we need is the block information for the rows/columns. That is, we need to know which vertex is in which tree in the forest after the modified hierarchical decomposition algorithm is performed. Instead of a tree (or a forest), **HDPRE** returns this information in the *scomp* array.

The structure of the algorithm **HDPRE** is similar to that of **HD**. In addition to  $G$ ,  $\sigma$  and  $i$ , **HDPRE** requires an additional input array *vsiz*e which stores the number of vertices condensed into each vertex of  $V$ . Note that for a simple vertex, this value is one. Hence, for the initial call with  $G = (V, E)$ , *vsiz*e is an array containing  $|V|$  ones. On the other hand, for the condensed vertices, this value will be equal to the sum of the *vsiz*e values condensed into that vertex. For the condensed digraph in Figure 5(c),  $vsiz$ e =  $\{3, 2, 1\}$  when its vertices are ordered from left to right. To be precise, for a recursive call with  $G = (V, E)$ , the total number of simple vertices is  $\sum_{v \in V} vsiz$ e( $v$ ) and this number is larger than  $mbs$  for all recursive calls because of the size check in line 14 of Algorithm 2.

At first, **HDPRE** checks if the problem size of the recursive call  $|E| - i$  is equal to one. If this is the case, it finds the strong components of  $G$ . If a strong component  $SC_\ell = (V_\ell, E_\ell)$  has  $\sum_{v \in V_\ell} vsiz$ e( $v$ )  $>$   $mbs$  vertices then **HDPRE** considers each vertex in  $V_\ell$  as a different strong component. Otherwise, i.e., if the size of a strong component is less than or equal to  $mbs$ , that component is not divided. **HDPRE** constructs the *scomp* array and returns. If the problem size,  $|E| - i$  is greater than 1, similar to **HD**, it constructs  $G_j$  for  $j = \lceil (i + |E|)/2 \rceil$  and if it is strongly connected the search for the combining edge among the first  $j$  edges starts with the call **HDPRE**( $G_j, \sigma, i, vsiz$ e). If not, for every strong component  $SC_\ell = (V_\ell, E_\ell)$  of  $G_j$  such that  $\sum_{v \in V_\ell} vsiz$ e( $v$ )  $>$   $mbs$ , it makes a recursive call **HDPRE**( $SC_\ell, \sigma_\ell, i_\ell, vsiz$ e $_\ell$ ) and updates the strong component information for the vertices in  $V_\ell$ . This update operation can be considered as further dividing the strong component  $SC_\ell$  hierarchically until all of the strong components obtained during this process contain at most  $mbs$  vertices.

Similarly to **HD**, at line 28, **HDPRE** makes one more recursive call for the condensed graph  $G^\mathcal{V}$  where the definition of the vertex partition  $\mathcal{V}$  (in line 22) is modified as described above. In **HD**, each vertex in the condensed graph corresponds to a strong component of  $G_j$  which defines a partition  $\mathcal{V}_1$ . In **HDPRE**, these components are further divided until all of them have a size no larger than  $mbs$ . A second partition,  $\mathcal{V}_2$  is obtained from these

---

**Algorithm 2**  $scomp = \text{HDPRE}(G = (V, E), \sigma, i, vsize)$  ( $mbs$  is global,  $i = 0$  for the initial call).

---

```

1: if  $|E| - i = 1$  then
2:   find strong components of  $G$ 
3:   for each strong component  $SC_\ell = (V_\ell, E_\ell)$  of  $G$  do
4:     if  $\sum_{v \in V_\ell} vsize(v) > mbs$  then
5:       consider each  $v \in V_\ell$  as a strong component
6:     else
7:        $\forall v \in V_\ell, scomp(v) = \ell$ 
8:   return  $scomp$ 
9:  $j = \lceil (i + |E|)/2 \rceil$ 
10: if  $G_j = (V, \{\sigma(k) : 1 \leq k \leq j\})$  is strongly connected then
11:   return  $scomp = \text{HDPRE}(G_j, \sigma, i, vsize)$ 
12: else
13:   for each strong component  $SC_\ell = (V_\ell, E_\ell)$  of  $G_j$  do
14:     if  $\sum_{v \in V_\ell} vsize(v) > mbs$  then
15:        $\sigma_\ell =$  the permutation of  $E_\ell$  ordered with respect to  $\sigma$ 
16:       compute  $i_\ell$  as in Algorithm 1
17:        $vsize_\ell(v) = vsize(v), \forall v \in V_\ell$ 
18:        $scomp_\ell = \text{HDPRE}(SC_\ell, \sigma_\ell, i_\ell, vsize_\ell)$ 
19:       update  $scomp$  according to  $scomp_\ell$ 
20:    $\mathcal{V}_1 = \{V_\ell : SC_\ell \text{ is a strong component of } G_j\}$ 
21:    $\mathcal{V}_2 = \{V_{\ell'} : SC_{\ell'} = (V_{\ell'}, E_{\ell'}) \text{ is a strong component in } scomp\}$ 
22:    $\mathcal{V} = \{\mathcal{V}_1, \mathcal{V}_2\}$ 
23:    $G^\mathcal{V} = \text{pcondense}(G, \mathcal{V}, mbs) = (V^{\mathcal{V}_2}, E^{\mathcal{V}_1})$ 
24:    $\sigma^\mathcal{V} =$  the permutation of  $E^{\mathcal{V}_1}$  ordered with respect to  $\sigma$ 
25:   compute  $i^\mathcal{V}$  as in Algorithm 1
26:   if  $i^\mathcal{V} \neq |E^{\mathcal{V}_1}|$  then
27:      $vsize^\mathcal{V}(v_{\ell'}) = \sum_{v \in V_{\ell'}} vsize(v), \forall V_{\ell'} \in \mathcal{V}_2$ 
28:      $scomp^\mathcal{V} = \text{HDPRE}(G^\mathcal{V}, \sigma^\mathcal{V}, i^\mathcal{V}, vsize^\mathcal{V})$ 
29:     update  $scomp$  with respect to  $scomp^\mathcal{V}$ 
30:   return  $scomp$ 

```

---

smaller strong components and  $\mathcal{V} = \{\mathcal{V}_1, \mathcal{V}_2\}$  is defined. After obtaining the condensed graph  $G^\mathcal{V}$ , HDPRE checks if  $G^\mathcal{V}$  is known to be acyclic. Note that if  $i^\mathcal{V} = |E^{\mathcal{V}_1}|$ , no strong component with two or more vertices exists in  $G^\mathcal{V}$  and hence it is acyclic. If  $i^\mathcal{V} \neq |E^{\mathcal{V}_1}|$ , after obtaining  $scomp^\mathcal{V}$ , HDPRE updates  $scomp$  if a larger strong component is obtained.

For the matrix given in Figure 1(a), HDPRE generates the blocks for the cases  $mbs = 2$  and  $mbs = 3$  as shown in Figure 8(a) and Figure 8(b), respectively. For  $mbs = 2$ , the condensed graph has 5 vertices and no edges hence no combination will occur. For  $mbs = 3$ , as shown in Figure 8(b), the condensed graph has 3 vertices where 2 of them will combine with the 12th edge in  $\sigma_0$ .

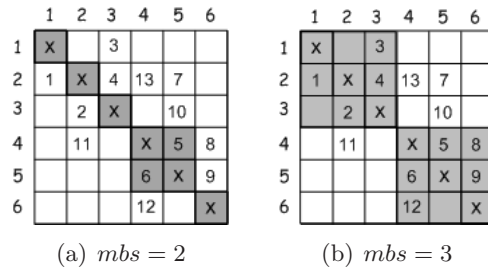


Figure 8: Initial block structure of the preconditioner after HDPRE algorithm. Two cases,  $mbs = 2$  and  $mbs = 3$ , are investigated for the matrix in Figure 1(a).

### 3.1.3 Combining the Blocks

After HDPRE obtains a block diagonal partition, SCPRE performs a loop on the nonzeros which are not contained in a block on the diagonal to see if it is possible to put more into the block diagonal by combining original blocks. To do this, SCPRE first constructs a condensed simple graph  $H$  where the vertices of  $H$  correspond to the diagonal blocks and inter-block edges of  $G$  in both directions are combined as a single edge with a weight that is the sum of the weights of the combined edges.

After  $H$  is obtained, its edges are visited in an order corresponding to a permutation  $\sigma_H$ . This permutation is consistent with the original permutation  $\sigma_0$ . That is, if the edges of the original digraph are sorted in descending order with respect to the edge weights,  $\sigma_H$  permutes the edges of  $H$  with respect to descending edge weights. On the other hand, if the initial permutation is based on the RCM ordering we compute the RCM ordering of  $H$ ,

relabel the vertices of  $H$  accordingly, and order the edges with respect to this RCM ordering. Let  $vsize(u)$  be the number of rows/columns in a block corresponding to the vertex  $u$ .

Assume that SCPRE constructs  $\sigma_0$  by sorting the edges with respect to decreasing weights. For the matrix given in Figure 8(a), if  $w(2) + w(4) > w(1)$  then vertices 2 and 3 are combined. Since  $mbs = 2$  and there is no edge between vertices 1 and 6, these two vertices remain as singletons.

### 3.2 SCPRE: Extending the Preconditioner

If the desired structure of  $\mathbf{M}$  is block diagonal SCPRE stops after the first phase. Otherwise, while preserving the blocks, it tries to extend the block diagonal preconditioner to a block upper-triangular one. Note that in this case the order of the blocks is important since it changes the nonzeros in the upper-triangular part of  $\mathbf{M}$ . Hence, before this extension, SCPRE tries to put more magnitude into the block upper-triangular part via a block permutation.

Let  $G = (V, E)$  be the digraph associated with the matrix and  $k$  be the number of its diagonal blocks after the first phase. Let  $\mathcal{V}_1 = \{V_1, V_2, \dots, V_k\}$  be a partition of  $V$  such that the vertices in  $V_i$  correspond to the rows/columns of  $i$ th block. Let  $\mathcal{V} = \{\mathcal{V}_1, \mathcal{V}_1\}$  and  $G^\mathcal{V} = \text{condense}(G, \mathcal{V})$  be the condensed multigraph. Note that if  $G^\mathcal{V}$  is acyclic, a topological sort in  $G^\mathcal{V}$  gives a symmetric block permutation such that all of the nonzeros in the matrix will be in the upper-triangular part of the permuted matrix. However, this only happens for a reducible matrix with blocks having no more than  $mbs$  rows/columns.

The problem of finding a good block permutation, which maximizes the number of nonzeros in the upper part of  $\mathbf{M}$ , can be reduced to the problem of finding the smallest edge set  $E'$  such that  $\overline{G}^\mathcal{V} = (V^\mathcal{V}_1, E^\mathcal{V}_1 \setminus E')$  is acyclic. For the weighted version of the problem, i.e., to maximize the total magnitude in the upper part, we need to find an edge set  $E'$  where  $\overline{G}^\mathcal{V}$  is acyclic and the sum  $\sum_{uv \in E'} |w(uv)|$  is minimum. In the literature, the first problem is called the *directed feedback arc set* problem and the second one is called the *directed weighted feedback arc set* problem. Both problems are NP-complete [14, 15].

Our simple heuristic proposed for this problem works as follows: Let  $G^\mathcal{V}$  be the condensed graph described above. For each vertex  $u \in V^\mathcal{V}$ , let  $tweight(u) = \sum_{w \in E^\mathcal{V}} w(uw)$ . The main body of the algorithm is a **for** loop where at the  $i$ th iteration, it chooses the vertex  $u$  with maximum  $tweight$  and assigns it as the  $i$ th vertex in the permutation. It then removes  $u$  from  $V^\mathcal{V}$ , its edges from  $E^\mathcal{V}$ , and continues with the next iteration.

## 4 Using SCPRE with an Iterative Solver

The iterative solver we use in our experiments is the right-preconditioned GMRES [20] with restarts. A template for this can be found in [2]. Let  $\mathbf{A} = \mathbf{D} + \mathbf{U} + \mathbf{L}$  be the scaled and permuted matrix such that  $\mathbf{D}$ ,  $\mathbf{U}$ , and  $\mathbf{L}$  are the block diagonal, upper, and lower parts, respectively.

Hence, if the desired structure is block diagonal, which is suitable for exploitation of parallelism,  $\mathbf{M} = \mathbf{D}$  is the preconditioner. If this is not the case,  $\mathbf{M} = \mathbf{D} + \mathbf{U}$  is the preconditioner for  $\mathbf{A}$ . For the latter case, the computation  $\mathbf{A}\mathbf{M}^{-1}\mathbf{x}$  becomes

$$\mathbf{A}\mathbf{M}^{-1}\mathbf{x} = (\mathbf{D} + \mathbf{U} + \mathbf{L})(\mathbf{D} + \mathbf{U})^{-1}\mathbf{x} = \mathbf{x} + \mathbf{L}((\mathbf{D} + \mathbf{U})^{-1}\mathbf{x}).$$

Note that SCPRE tries to maximize the total magnitude in  $\mathbf{D}$  and  $\mathbf{U}$ . As a consequence and as in experiments not included here show,  $\mathbf{L}$  usually contains many fewer nonzeros than  $\mathbf{A}$ . Hence computing the vector  $\mathbf{z} = \mathbf{L}\mathbf{y}$  usually takes very little time and the main operation is to compute  $\mathbf{y} = (\mathbf{D} + \mathbf{U})^{-1}\mathbf{x} = \mathbf{M}^{-1}\mathbf{x}$ .

In our implementation, in addition to  $\mathbf{A}$ , we store the LU factors of the diagonal blocks, i.e., the factors  $\mathbf{L}_i$  and  $\mathbf{U}_i$  such that  $\mathbf{D}_i = \mathbf{L}_i\mathbf{U}_i$  where  $\mathbf{D}_i$  is the  $i$ th diagonal block. We reduce the memory requirements for these factors by ordering the blocks using the approximate minimum degree (AMD) heuristic [1, 7] before the MATLAB sparse factorization. We then solve the upper block triangular system  $\mathbf{M}\mathbf{y} = \mathbf{x}$  using these factors, starting with the last block, so that the off-diagonal part  $\mathbf{U}$  is only used to multiply vectors.

### 4.1 Robustness

Although it is very rare, we observe that the preconditioner  $\mathbf{M}$  obtained by SCPRE is singular for some matrices in our test set. That is, some of the blocks on the diagonal of  $\mathbf{M}$  are singular. When the original matrix is diagonally dominant or close to an  $I$ -matrix, with high probability the diagonal blocks of  $\mathbf{M}$  are  $I$ -matrices as well. If this is not the case, scaling and permuting the matrix via MC64 will probably help us to convert it to an  $I$ -matrix.

When using the MATLAB factorization, we guard against this potential problem by using the simple and cheap stability check proposed and used by XPABLO [12, 13]. That is, if  $n_i$  is the dimension of  $\mathbf{D}_i$ , after computing  $\mathbf{L}_i$  and  $\mathbf{U}_i$ , we check whether

$$\left| 1 - \frac{\|\mathbf{U}_i^{-1}\mathbf{L}_i^{-1}\mathbf{x}\|}{\|\mathbf{e}\|} \right| < \sqrt{\epsilon_M}, \quad (2)$$

where  $\mathbf{e} = (1, \dots, 1)^T$  is an  $n_i \times 1$  column vector,  $\mathbf{x} = \mathbf{D}_i \mathbf{e}$ , and  $\epsilon_M$  is machine epsilon. They suggest that, if a block does not satisfy (2),  $\mathbf{D}_i$  is replaced either by  $\mathbf{U}_i$  or  $\mathbf{L}_i$  according to whether they are solving a block upper- or lower-triangular system, respectively. For **SCP**, we always use the factor having the largest Frobenius norm to replace  $\mathbf{D}_i$ , where the Frobenius norm of an  $n \times n$  matrix  $\mathbf{B}$  is given by

$$\|\mathbf{B}\|_F = \sqrt{\sum_{1 \leq i, j \leq n} |\mathbf{B}_{ij}|^2}.$$

## 5 Experiments

All of the experiments are conducted on an Intel 2.4Ghz Quad Core computer, equipped with 4GB RAM with a Linux operating system. For the experiments, we use two sets of matrices from the University of Florida Sparse Matrix Collection<sup>1</sup>. Each set contains matrices arising from either a circuit simulation problem (CSP), or a semiconductor device problem (SDP), or an electromagnetics problem (EMP). The first set contains 45 matrices with  $m \leq 2 \times 10^6$  nonzeros. For this set, we use  $mbs = 2000$  in the experiments. The second set contains 13 relatively large matrices with  $2.5 \times 10^6 \leq m \leq 3 \times 10^6$  nonzeros. For this set, we use  $mbs = 5000$  since they are larger. To construct the sets, we first choose all matrices with these properties and then remove those whose largest blocks in their BTF form have less than  $mbs$  rows/columns. In the tables, we exclude the matrices on which none of our preconditioned iterative solvers converged. The lists of the remaining 37 matrices in the first set and 12 matrices in the second set are given in Table 1.

In our experiments, we restarted **GMRES** [20] after every 50 iterations. The desired error tolerance for **GMRES**(50) is set to  $\epsilon = 10^{-8}$  and the stopping criterion we use for **GMRES** is

$$\frac{\|\mathbf{A}\mathbf{M}^{-1}\bar{\mathbf{z}} - \mathbf{b}\|}{\|\mathbf{b}\|} < \epsilon$$

where  $\bar{\mathbf{z}} = \mathbf{M}\bar{\mathbf{x}}$ , with  $\bar{\mathbf{z}}$  the computed solution of the preconditioned system and  $\bar{\mathbf{x}}$  the computed solution of the original system. After obtaining the solution  $\bar{\mathbf{x}}$  to the original system, we compute the relative error  $\|\mathbf{A}\bar{\mathbf{x}} - \mathbf{b}\|/\|\mathbf{b}\|$  to the unpreconditioned system. For all cases, this error is smaller than  $10^{-7}$  and indeed, for most of the cases it is also smaller than  $\epsilon$ .

<sup>1</sup><http://www.cise.ufl.edu/research/sparse/matrices/>

Table 1: Properties of the matrices used for the experiments.  $n$  is the dimension of the matrix,  $m$  is the number of nonzeros, and  $n_1$  and  $n_2$  are the size of the largest and second largest blocks in the BTF form. Note that  $n_2 = 0$  means that the matrix is irreducible, i.e.,  $n_1 = n$ . The column type shows the application from which the matrix arises. The sets are sorted first according to the type of the problem and then their  $n_1$  values.

	Matrix	Group	$n$	$m$	$n_1$	$n_2$	Type	
SET 1	<i>Hamrle2</i>	Hamrle	5952	22162	5952	0	CSP	
	<i>rajat03</i>	Rajat	7602	32653	7500	1		
	<i>circuit_3</i>	Bomhof	12127	48137	7607	1		
	<i>coupled</i>	IBM_Austin	11341	97193	11293	1		
	<i>memplus</i>	Hamm	17758	99147	17736	1		
	<i>rajat22</i>	Rajat	39899	195429	26316	7672		
	<i>onetone2</i>	ATandT	36057	222596	32211	2		
	<i>onetone1</i>	ATandT	36057	335552	32211	2		
	<i>rajat15</i>	Rajat	37261	443573	37243	1		
	<i>ckt11752_tr_0</i>	IBM_EDA	49702	332807	49371	44		
	<i>circuit_4</i>	Bomhof	80209	307604	52005	7		
	<i>bcircuit</i>	Hamm	68902	375558	68902	0		
	<i>rajat18</i>	Rajat	94294	479151	84507	52		
	<i>hcircuit</i>	Hamm	105676	513072	92144	4927		
	<i>ASIC_100ks</i>	Sandia	99190	578890	98843	2		
	<i>ASIC_100k</i>	Sandia	99340	940621	98843	2		
	<i>ASIC_680ks</i>	Sandia	682712	1693767	98843	2		
	<i>rajat23</i>	Rajat	110355	555441	103024	216		
	<i>twotone</i>	ATandT	120750	1206265	105740	6		
	<i>trans5</i>	IBM_EDA	116835	749800	116817	1		
	<i>dc2</i>	IBM_EDA	116835	766396	116817	1		
	<i>G2_circuit</i>	AMD	150102	726674	150102	0		
	<i>scircuit</i>	Hamm	170998	958936	170493	216		
	<i>transient</i>	Freescall	178866	961368	178823	11		
	<i>Raj1</i>	Rajat	263743	1300261	263571	5		
	<i>ASIC_320ks</i>	Sandia	321671	1316085	320926	6		
	<i>ASIC_320k</i>	Sandia	321821	1931828	320926	6		
	<i>utm5940</i>	TOKAMAK	5940	83842	5794	1		EMP
	<i>dw4096</i>	Bai	8192	41746	8192	0		
	<i>Zhao1</i>	Zhao	33861	166453	33861	0		
<i>igbt3</i>	Schenk_ISEI	10938	130500	10938	0	SDP		
<i>wang3</i>	Wang	26064	177168	26064	0			
<i>wang4</i>	Wang	26068	177196	26068	0			
<i>ecl32</i>	Sanghavi	51993	380415	42341	1			
<i>ibm_matrix_2</i>	Schenk_IBMSDS	51448	537038	44822	1			
<i>matrix-new_3</i>	Schenk_IBMSDS	125329	893984	78672	1			
<i>matrix_9</i>	Schenk_IBMSDS	103430	1205518	99372	1			
SET 2	<i>ASIC_680k</i>	Sandia	682862	2638997	98843	2	CSP	
	<i>G3_circuit</i>	AMD	1585478	7660826	181343	0		
	<i>rajat29</i>	Rajat	643994	3760246	629328	71		
	<i>rajat30</i>	Rajat	643994	6175244	632151	0		
	<i>Hamrle3</i>	Hamrle	1447360	5514242	1447360	0		
	<i>memchip</i>	Freescall	2707524	13343948	2706851	0	EMP	
	<i>offshore</i>	Um	259789	4242673	259789	0		
	<i>tmt_sym</i>	CEMW	726713	5080961	726713	0		
	<i>t2em</i>	CEMW	921632	4590832	917300	1		
	<i>tmt_unsym</i>	CEMW	917825	4584801	917825	0		
<i>para-4</i>	Schenk_ISEI	153226	2930882	153226	0	SDP		
<i>ohne2</i>	Schenk_ISEI	181343	6869939	181343	0			

The maximum number of outer iterations is set to 20, hence the maximum number of inner iterations is 1000. In the tables, we give the inner iteration counts when this criterion is satisfied. Otherwise, if the criterion is not satisfied, we put “–” in the corresponding cell to denote that GMRES did not converge. Also, we put the lowest iteration count for each matrix in bold font.

To compare the efficiency of the preconditioner, we used a generic preconditioner, ILUT [18, 19], from MATLAB 7.11 with two drop tolerances,  $dtol = 10^{-3}$  and  $10^{-4}$ . In addition to ILUT, we also compared our results with those of XPABLO [12, 13]. For all of the preconditioners, we use MC64 and obtain a maximum product transversal by scaling and permuting the matrix as a preprocessing step.

In the MATLAB implementation of ILUT, for the  $j$ th column of the incomplete  $\mathbf{L}$  and  $\mathbf{U}$ , entries smaller in magnitude than  $dtol \times \|\mathbf{A}_{*j}\|$  are deleted from the factor where  $\|\mathbf{A}_{*j}\|$  is the norm of the  $j$ th column of  $\mathbf{A}$ . However, the diagonal entries of  $\mathbf{U}$  are always kept to avoid a singular factor. To use ILUT based preconditioners, we use AMD before computing the incomplete factorization of the matrix. For XPABLO preconditioners, we use the J variant for the block Jacobi iterations, and LX and UX variants for the forward and backward block Gauss-Seidel iterations, respectively, with the parameters given in [12, 13]. For the maximum block size of XPABLO, we used the same  $mbs$  as for SCPRE.

When the matrix is reducible, for GMRES with block Gauss-Seidel iterations, SCPRE has an advantage as it will automatically find the block triangular form and will then work within the diagonal blocks of the form. As the other algorithms do not detect this form automatically, for a fair comparison we used the reducibility information also for the ILUT and tt XPABLO preconditioners. While using ILUT (XPABLO) for reducible matrices, we first compute the BTF form and apply ILUT (XPABLO) only to the blocks with more than  $mbs$  rows and columns. For smaller blocks, we compute the complete factors. We then use these complete and incomplete factors together while computing a matrix vector product using  $\mathbf{M}^{-1}$ . Our experiments show that this approach is almost always better than using ILUT (XPABLO) in a straightforward manner in terms of iteration count. We also tried this approach while using the J variant of the XPABLO preconditioner. Surprisingly, even for the block Jacobi case, this approach slightly helps to reduce the iteration counts for most of the reducible matrices. We call this variant **J-red** in the tables below. Note that for the block Gauss-Seidel case, when we apply XPABLO (or ILUT) only to the large blocks in the BTF form of a reducible matrix, we keep all of the nonzeros in the preconditioner from the

off-diagonal blocks. However, for block Jacobi iterations, we automatically drop them from the preconditioning matrix  $\mathbf{M}$  since its desired structure is block diagonal, not block triangular.

In addition to the number of iterations required for convergence, we compare the performance of the preconditioners according to the relative memory requirement with respect to the number of nonzeros in  $\mathbf{A}$ . Let  $nz(\mathbf{B})$  be the number of nonzeros in a matrix  $\mathbf{B}$ . For ILUT, the relative memory requirement is equal to

$$mem_{\text{ILUT}} = \frac{nz(\mathbf{L}) + nz(\mathbf{U})}{nz(\mathbf{A})},$$

where  $\mathbf{L}$  and  $\mathbf{U}$  are the incomplete triangular factors of  $\mathbf{A}$ . On the other hand, the relative memory requirement for SCPRE and XPABLO is equal to

$$mem_{\text{SCPRE}} = mem_{\text{XPABLO}} = \frac{\sum_{i=1}^k (nz(\mathbf{L}_i) + nz(\mathbf{U}_i))}{nz(\mathbf{A})},$$

where  $k$  is the number of blocks in the block diagonal  $\mathbf{D}$  and  $\mathbf{L}_i$  and  $\mathbf{U}_i$  are the lower- and upper-triangular factors of the  $LU$  factorization of the  $i$ th block in  $\mathbf{D}$ . Note that the relative memory requirements of the preconditioners can give an idea for the cost of computing  $\mathbf{M}^{-1}\mathbf{x}$ . Assuming  $x$  is a dense vector, a preconditioned GMRES iteration will require approximately  $nz(\mathbf{A})(1 + mem_{\mathbf{X}})$  operations for the preconditioner generated by the algorithm  $\mathbf{X}$ .

There are two parameters for the proposed algorithm: the first is the maximum block size,  $mbs$ , the second is the permutation for the nonzeros, denoted by  $\sigma_0$ . As expected, our experiments (not reported on here) show that increasing  $mbs$  usually reduces the iteration counts and increases the relative memory requirements of the solver.

We conduct some experiments to show the effect of our choice of  $\sigma_0$  on the performance of our algorithm. Note that in HD the edges are sorted in increasing order with respect to their weights. In our implementation, we define the weight of an edge as the magnitude of the corresponding nonzero and sort the edges in decreasing order. We test our decision by comparing its effect with that of a random permutation. As Table 2 shows, our decision to sort the edges in decreasing order with respect to the edge weights makes the solver converge more quickly.

## 5.1 Experiments with Block Gauss-Seidel Iterations

Table 3 shows the performance of SCPRE, XPABLO and ILUT preconditioners for block Gauss-Seidel iterations. Note that both SCPRE(*dec*) and SCPRE(RCM)

Table 2: Effect of the permutation  $\sigma_0$  on the number of iterations. Two options are compared: decreasing order with respect to the edge weights and a random order. Maximum block size for SCPRE is set to 2000 where the structure of  $\mathbf{M}$  is block upper-triangular. For each case, the ratio of the total magnitude in  $\mathbf{M}$  to the total magnitude in  $\mathbf{A}$ , the relative memory requirement, and the number of inner iterations for preconditioned GMRES are given.

Matrix	Decreasing			Random		
	$\frac{\sum  \mathbf{M}_{ij} }{\sum  \mathbf{A}_{ij} }$	$mem_{\text{SCPRE}}$	$iters$	$\frac{\sum  \mathbf{M}_{ij} }{\sum  \mathbf{A}_{ij} }$	$mem_{\text{SCPRE}}$	$iters$
<i>Hamrle2</i>	0.998	2.03	<b>16</b>	0.993	2.05	157
<i>rajat03</i>	0.999	1.07	<b>2</b>	0.997	1.02	5
<i>circuit_3</i>	0.996	1.45	<b>9</b>	0.987	1.23	445
<i>coupled</i>	0.998	1.57	<b>11</b>	0.992	1.58	34
<i>memplus</i>	0.999	1.03	<b>5</b>	0.998	1.03	7
<i>rajat22</i>	0.973	1.20	<b>21</b>	0.962	1.15	-

are robust that is, the solvers converge for most of the matrices. Although there are a few matrices for which the SCPRE(RCM) preconditioned solver converges more quickly than that preconditioned with SCPRE(*dec*) (such as *ASIC\_680k*) and, amongst all preconditioners, only SCPRE(RCM) converges for matrices *onetone1* and *onetone2*, SCPRE(*dec*) is almost always better and is our preferred preconditioner.

In general, all the preconditioners work well for the matrices in the first set. However, SCPRE(*dec*) is the most robust since the preconditioned solver fails to converge only for 3 out of 37 matrices whereas the next best result is 9 by XPABLO variants. Thus SCPRE(*dec*) is the best block preconditioner on this set of matrices. When comparing SCPRE(*dec*) to ILUT( $10^{-4}$ ) on this set we see that they are comparable in terms of the number of best performances but ILUT( $10^{-4}$ ) is less robust failing to converge for 10 matrices in this set and requiring more memory than SCPRE(*dec*).

For the second set, ILUT( $10^{-4}$ ) is the best preconditioner in terms of robustness and iteration count. For the matrices in this set, the ILUT( $10^{-4}$ ) preconditioned solver fails to converge in only 2 out of 12 matrices whereas SCPRE(*dec*) does not converge on 4. Although ILUT( $10^{-4}$ ) is better than SCPRE(*dec*) for 10 out of 12 matrices in the second set, its average relative memory usage is 9.39 which is almost 3 times as much as the relative memory requirement of SCPRE(*dec*). Note that, for the second set, even ILUT( $10^{-3}$ ) uses slightly more memory than SCPRE(*dec*). However, it fails to converge on 7 matrices. Hence, if memory is the bottleneck, SCPRE(*dec*) may be a suitable choice for preconditioning.

The performance of the SCPRE based preconditioners changes with respect to the application. For example, as Table 3 shows, SCPRE(*dec*) preconditioned GMRES fails to converge in 3 out of 7 matrices from electromagnetics applications. On the other hand, it fails to converge on only 4 of the remaining 42 matrices. Hence its performance is much better for circuit and device simulation applications. Note that, even though some of these matrices are reducible, they have a large reducible block with size much larger than *mbs*. That is, we still have a large subproblem to deal with. On the circuit simulation and semiconductor device matrices, SCPRE’s performance is far better than that of XPABLO which is another block based preconditioner with a promising performance in practice for several matrix classes [3, 5, 9]. Note that we used the BTF forms of the reducible matrices for both the XPABLO and ILUT preconditioners. Hence, reducibility alone is not a reason for the good performance of SCPRE based preconditioners.

## 5.2 Experiments with Block Jacobi Iterations

Table 4 shows the performance of SCPRE and XPABLO preconditioners for block Jacobi iterations. Similar to the experiments with block Gauss-Seidel iterations, performance of SCPRE(*dec*) is much better than that of SCPRE(RCM) for the matrices in our sets. For XPABLO, applying the preconditioner only to the blocks in the BTF form, variant J-red, reduces the number of iterations on 11 matrices. Furthermore, for 5 of the matrices J-red converges whereas J does not. Note that there are 32 reducible matrices in the sets and J-red differs from J only for these matrices. Although J-red required more iterations for convergence for matrices *matrix\_new\_3* and *matrix\_9*, in general, for the matrices in our experiments, J-red performs better than J.

As Table 4 shows, SCPRE(*dec*) preconditioned GMRES converges for 36 matrices whereas XPABLO’s J-red variant converges for only 24 matrices. The XPABLO based preconditioner has the least number of iterations in only 8 cases whereas the SCPRE variants are better on 35 matrices. The difference in the performance is not due to the relative memory usage of SCPRE variants. For the first set, SCPRE(*dec*) uses only 8% more memory than XPABLO(J-red) on average and, for the second set, its memory usage is much less.

In the right-hand side of Table 4, the execution times of the GMRES solver are given. As the table shows, for most of the cases, the best solver in terms of iteration count has also the best execution time. Note that there are some exceptions such as *matrix\_9* for which the solver preconditioned by XPABLO(J) requires 49 iterations fewer than when preconditioned by SCPRE but its execution time is slightly more. This is because, for this matrix,

Table 3: Number of inner iterations for GMRES using XPABLO, ILUT and SCPRE preconditioners and block Gauss-Seidel iterations. For SCPRE and XPABLO,  $mbs$  is set to 2000 and 5000 for the first and second sets, respectively. For SCPRE, we give the results using two permutations for  $\sigma_0$ , based on descending order and RCM. For XPABLO, we give the results for the UX and LX variants which are used with the parameters suggested in [12]. For ILUT, the drop tolerance is set to  $10^{-3}$  and  $10^{-4}$ . A ‘-’ sign indicates that the preconditioned solver did not converge. Average relative memory requirements are computed by taking the averages over the cases when the solver converges.

	Matrix	XPABLO		SCPRE		ILUT	
		UX	LX	dec	RCM	$10^{-3}$	$10^{-4}$
$mbs =$ 2000	<i>Hamrle2</i>	31	31	16	28	6	4
	<i>rajat03</i>	<b>2</b>	<b>2</b>	<b>2</b>	<b>2</b>	<b>2</b>	<b>2</b>
	<i>circuit_3</i>	135	137	<b>9</b>	61	-	-
	<i>coupled</i>	12	12	11	13	6	4
	<i>memplus</i>	9	9	<b>5</b>	18	15	9
	<i>rajat22</i>	36	37	21	61	36	<b>16</b>
	<i>onetone2</i>	-	-	-	<b>248</b>	-	-
	<i>onetone1</i>	-	-	-	<b>297</b>	-	-
	<i>rajat15</i>	-	-	120	467	-	<b>33</b>
	<i>ckt11752_tr_0</i>	197	188	<b>19</b>	323	-	-
	<i>circuit_4</i>	100	81	<b>39</b>	346	-	-
	<i>bcircuit</i>	-	-	<b>40</b>	620	568	93
	<i>rajat18</i>	-	-	<b>11</b>	-	393	54
	<i>hcircuit</i>	8	9	9	21	9	<b>5</b>
	<i>ASIC_100ks</i>	9	10	9	10	<b>4</b>	<b>4</b>
	<i>ASIC_100k</i>	9	9	10	10	<b>4</b>	<b>4</b>
	<i>ASIC_680ks</i>	<b>3</b>	4	<b>3</b>	4	4	4
	<i>rajat23</i>	40	41	<b>16</b>	88	47	18
	<i>twotone</i>	-	-	<b>25</b>	128	-	48
	<i>trans5</i>	9	9	<b>5</b>	7	7	6
	<i>dc2</i>	13	12	12	11	10	<b>6</b>
	<i>G2_circuit</i>	-	-	444	834	124	<b>30</b>
	<i>scircuit</i>	741	764	<b>317</b>	977	-	-
	<i>transient</i>	-	-	<b>33</b>	-	-	-
	<i>Raj1</i>	775	789	636	-	269	<b>39</b>
	<i>ASIC_320ks</i>	4	4	<b>1</b>	4	2	2
	<i>ASIC_320k</i>	5	5	<b>2</b>	3	3	3
	<i>utm5940</i>	-	-	-	-	-	<b>29</b>
	<i>dw4096</i>	881	798	13	141	24	<b>10</b>
	<i>Zhao1</i>	7	7	4	9	4	<b>3</b>
	<i>igbt3</i>	29	29	20	17	94	<b>12</b>
	<i>wang3</i>	107	105	54	58	18	<b>9</b>
<i>wang4</i>	39	38	21	36	11	<b>6</b>	
<i>ecl32</i>	99	99	30	32	32	<b>13</b>	
<i>ibm_matrix_2</i>	-	249	<b>10</b>	16	-	-	
<i>matrix-new_3</i>	85	86	<b>30</b>	41	-	-	
<i>matrix_9</i>	146	90	98	<b>88</b>	-	-	
Avg. relative memory		2.95	3.04	3.36	3.19	2.12	4.02
$mbs =$ 5000	<i>ASIC_680k</i>	<b>2</b>	<b>2</b>	27	<b>2</b>	3	3
	<i>G3_circuit</i>	-	-	357	422	212	<b>81</b>
	<i>rajat29</i>	-	-	<b>11</b>	-	-	-
	<i>rajat30</i>	12	12	14	15	7	<b>5</b>
	<i>Hamrle3</i>	-	-	-	-	-	<b>17</b>
	<i>memchip</i>	26	27	10	20	8	<b>5</b>
	<i>offshore</i>	330	327	488	451	-	<b>15</b>
	<i>tmt_sym</i>	-	-	-	-	-	<b>69</b>
	<i>t2em</i>	-	-	876	-	132	<b>38</b>
	<i>tmt_unsym</i>	-	-	-	-	-	<b>136</b>
<i>para-4</i>	-	-	-	-	-	<b>433</b>	
<i>ohne2</i>	-	-	<b>196</b>	-	-	-	
Avg. relative memory		3.58	3.58	3.23	2.51	3.36	9.39

$mem_{\text{XPABLO}(J)} = 8.43$  and  $mem_{\text{SCPRED}(dec)} = 3.69$ , and the cheaper cost of computing  $\mathbf{M}^{-1}\mathbf{x}$  more than compensates for the difference in iteration counts. For 39 matrices, a **SCPRED** variant has the best or very close to the best time. In summary, **SCPRED**(*dec*) performs much better than **XPABLO** variants in our block Jacobi experiments.

## 6 Conclusions and Future Work

Given a linear system  $\mathbf{Ax} = \mathbf{b}$ , we have proposed a method to construct generic block diagonal and block triangular preconditioners. The proposed approach is based on Tarjan’s algorithm HD for hierarchical decomposition of a digraph into its strong subgraphs. Although our preconditioner **SCPRED**, in common with the other block preconditioners, does not perform well for electromagnetics matrices, we obtain promising results for many device and circuit simulation matrices and we suggest using it with these types of problems. In future research, the structure of graphs for different classes of matrices can be analysed to try to understand the reason for the difference in performance.

There are two main parameters for the algorithm: the way that a permutation  $\sigma_0$  of the edges is obtained and the maximum block size *mbs*. For  $\sigma_0$ , we used two approaches: the first sorts the edges in the order of decreasing weights. With this approach, we wanted to include nonzeros with large magnitudes in our preconditioner. The second approach uses the well known reverse Cuthill-McKee ordering. We tested this approach since a sparsity structure with a small bandwidth may be useful for putting more nonzeros into the preconditioner. The permutation decisions are validated by the experiments which also show that the first approach is usually better than the second. In future work, other ways to generate  $\sigma_0$  can be investigated.

The second parameter, *mbs*, affects the memory requirement of the matrix significantly, and hence the number of iterations required for convergence. The experiments show that for the preconditioners **ILUT**, **SCPRED** and **XPABLO**, the memory requirement and the number of iterations are inversely correlated. For the proposed preconditioner **SCPRED**, *mbs* needs to be set by the user without knowing how much memory will be required by the solver. In future work, we will look for a self-tuning mechanism which enables **SCPRED** to determine *mbs* automatically given the memory available to store the preconditioner.

Table 4: Number of inner iterations and solver times (in seconds) for GMRES using XPABLO and SCPRE preconditioners and block Jacobi iterations. The maximum block size  $mbs$  is set to 2000 and 5000 for the first and second sets, respectively. For SCPRE, we give the results using two permutations for  $\sigma_0$ , based on descending order and RCM. For XPABLO, we give the results for the J variant which is used with the parameters suggested in [12]. The J-red variant, described in the text, also uses the same parameters. A ‘-’ sign indicates that the preconditioned solver did not converge. Average relative memory requirements are computed by taking the averages over the cases when the solver converges.

Matrix	# iterations				solver time (in secs.)			
	XPABLO		SCPRE		XPABLO		SCPRE	
	J	J-red	dec	RCM	J	J-red	dec	RCM
<i>Hamrle2</i>	99	99	<b>31</b>	96	0.32	0.32	<b>0.08</b>	0.30
<i>rajat03</i>	7	<b>3</b>	4	4	0.03	<b>0.01</b>	<b>0.01</b>	<b>0.01</b>
<i>circuit_3</i>	680	327	<b>19</b>	179	3.72	1.81	<b>0.07</b>	0.93
<i>coupled</i>	43	22	<b>21</b>	25	0.22	<b>0.09</b>	<b>0.08</b>	0.11
<i>memplus</i>	17	17	<b>8</b>	33	0.08	0.08	<b>0.03</b>	0.19
<i>rajat22</i>	190	77	<b>42</b>	124	3.03	1.61	<b>0.63</b>	1.88
<i>onetone2</i>	-	-	-	<b>627</b>	-	-	-	<b>9.62</b>
<i>onetone1</i>	-	-	-	<b>622</b>	-	-	-	<b>14.95</b>
<i>rajat15</i>	-	-	<b>265</b>	-	-	-	<b>4.85</b>	-
<i>ckt11752_tr_0</i>	-	776	<b>36</b>	-	-	20.28	<b>0.83</b>	-
<i>circuit_4</i>	-	864	<b>112</b>	-	-	27.48	<b>3.33</b>	-
<i>bcircuit</i>	-	-	<b>107</b>	-	-	-	<b>3.06</b>	-
<i>rajat18</i>	-	-	<b>16</b>	-	-	-	<b>0.39</b>	-
<i>hcircuit</i>	16	<b>15</b>	16	40	<b>0.43</b>	<b>0.47</b>	<b>0.43</b>	1.55
<i>ASIC_100ks</i>	17	17	<b>16</b>	18	<b>0.46</b>	0.50	<b>0.44</b>	0.50
<i>ASIC_100k</i>	17	<b>16</b>	17	18	<b>0.48</b>	0.52	<b>0.49</b>	0.51
<i>ASIC_680ks</i>	-	<b>8</b>	-	<b>8</b>	-	<b>0.72</b>	-	<b>0.74</b>
<i>rajat23</i>	203	140	<b>32</b>	208	9.29	7.65	<b>1.17</b>	9.09
<i>twotone</i>	-	-	<b>49</b>	322	-	-	<b>2.70</b>	17.11
<i>trans5</i>	23	16	<b>9</b>	13	0.75	0.47	<b>0.24</b>	0.36
<i>dc2</i>	76	21	<b>20</b>	<b>20</b>	3.32	<b>0.67</b>	<b>0.64</b>	<b>0.64</b>
<i>G2_circuit</i>	-	-	<b>833</b>	-	-	-	<b>56.55</b>	-
<i>scircuit</i>	-	-	<b>682</b>	-	-	-	<b>49.25</b>	-
<i>transient</i>	-	-	<b>186</b>	-	-	-	<b>13.60</b>	-
<i>Raj1</i>	-	-	-	-	-	-	-	-
<i>ASIC_320ks</i>	5	6	<b>1</b>	7	0.43	0.70	<b>0.19</b>	0.54
<i>ASIC_320k</i>	11	9	<b>3</b>	10	0.91	0.85	<b>0.42</b>	0.61
<i>utm5940</i>	-	-	-	-	-	-	-	-
<i>dw4096</i>	-	-	<b>24</b>	-	-	-	<b>0.11</b>	-
<i>Zhao1</i>	12	12	<b>7</b>	16	0.12	0.12	<b>0.08</b>	0.19
<i>igbt3</i>	60	60	32	<b>26</b>	0.52	0.52	<b>0.21</b>	0.17
<i>wang3</i>	263	263	140	<b>138</b>	3.26	3.26	1.96	<b>1.78</b>
<i>wang4</i>	91	91	<b>39</b>	79	1.24	1.24	<b>0.54</b>	1.02
<i>ecl32</i>	-	-	<b>79</b>	90	-	-	<b>2.50</b>	3.08
<i>ibm_matrix_2</i>	-	344	<b>22</b>	30	-	12.29	<b>0.65</b>	1.09
<i>matrix-new_3</i>	184	248	<b>71</b>	95	13.21	20.20	<b>4.64</b>	6.79
<i>matrix_9</i>	<b>208</b>	240	257	346	<b>14.85</b>	18.80	<b>14.41</b>	21.83
Avg. relative memory	2.67	3.10	3.35	3.32				
<i>ASIC_680k</i>	-	<b>3</b>	-	<b>3</b>	-	<b>0.54</b>	-	<b>0.54</b>
<i>G3_circuit</i>	-	-	<b>674</b>	-	-	-	<b>516.38</b>	-
<i>rajat29</i>	-	-	<b>18</b>	-	-	-	<b>3.03</b>	-
<i>rajat30</i>	43	<b>22</b>	24	27	11.84	<b>4.59</b>	5.12	6.12
<i>Hamrle3</i>	-	-	-	-	-	-	-	-
<i>memchip</i>	41	50	<b>17</b>	38	53.39	74.55	<b>14.60</b>	38.82
<i>offshore</i>	<b>883</b>	<b>883</b>	-	-	<b>189.98</b>	<b>189.98</b>	-	-
<i>tmt_sym</i>	-	-	-	-	-	-	-	-
<i>t2em</i>	-	-	-	-	-	-	-	-
<i>tmt_unsym</i>	-	-	-	-	-	-	-	-
<i>para-4</i>	-	-	-	-	-	-	-	-
<i>ohme2</i>	-	-	-	-	-	-	-	-
Avg. relative memory	3.58	2.84	1.81	0.92				

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