

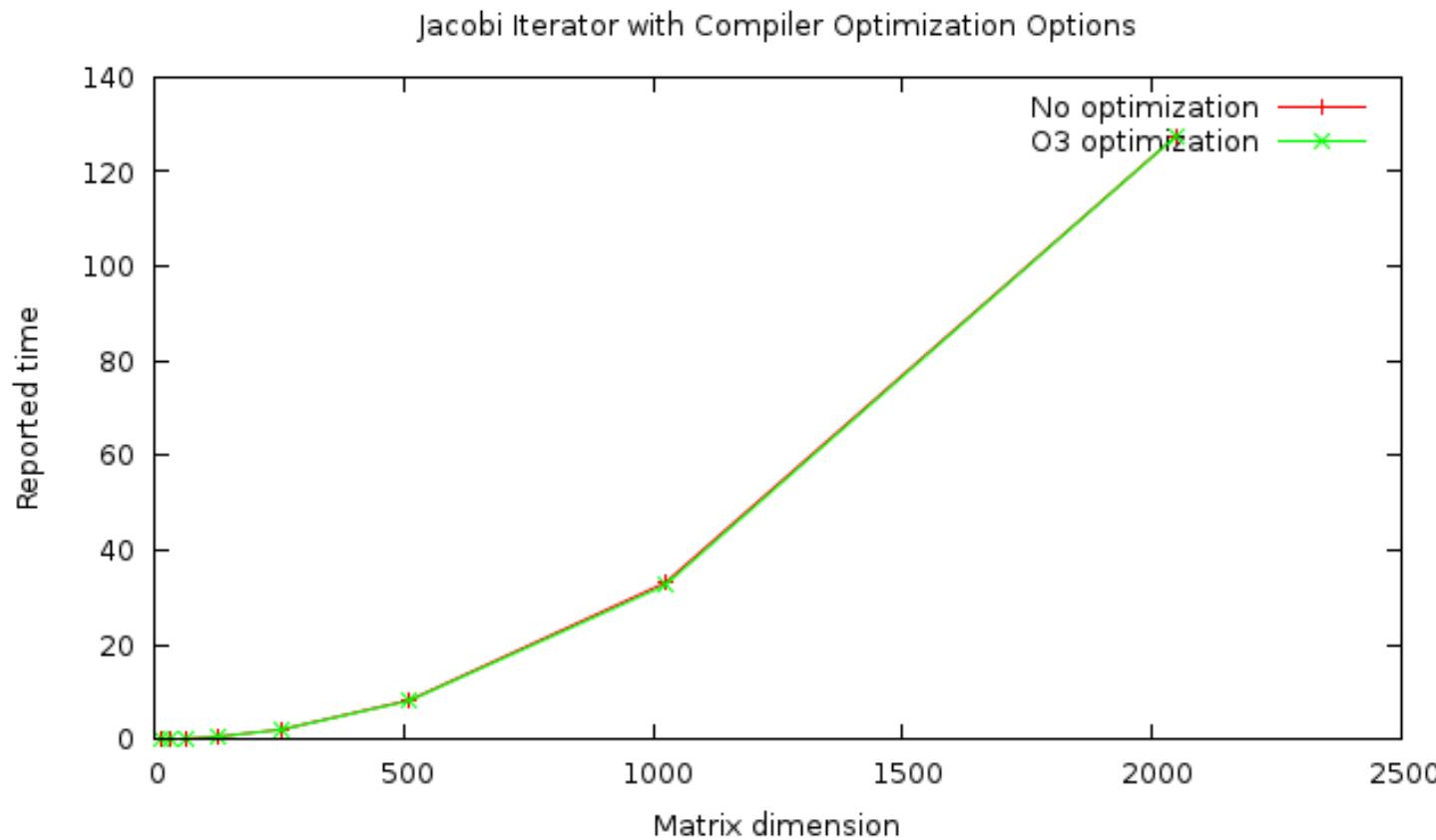
The Jacobi Iterator

(to solve Laplace's equation)



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Compiler Optimization

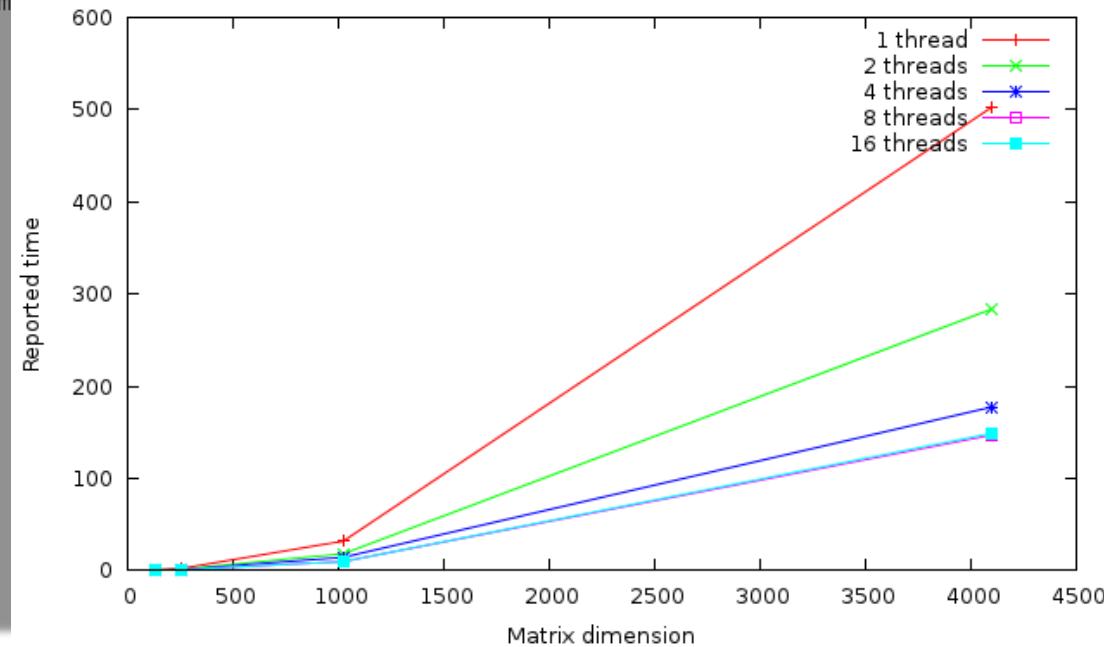


Open MP

```

73 // Iteration
74 double TimeStart = WallTime();
75
76 #pragma omp parallel for
77 for(int iCount = 1; iCount <=Iterations; iCount++){
78
79     //This will be a row dominant program.
80
81     for(int i=1;i<=Dimension;i++)
82         for(int j=1;j<=Dimension;j++)
83             SurfaceMatrix_t[i][j] = (0.25)*(SurfaceMatrix[i-1][j] +
84                                         SurfaceMatrix[i][j+1] +
85                                         SurfaceMatrix[i+1][j] +
86                                         SurfaceMatrix[i][j-1]);
87
88     PrintSurfaceMatrix(SurfaceMatrix_t,Dim);
89     double ** tmp;
90     tmp = SurfaceMatrix;
91     SurfaceMatrix = SurfaceMatrix_t;
92     SurfaceMatrix_t = tmp;
93 }
94 double TimeEnd = WallTime();
95

```

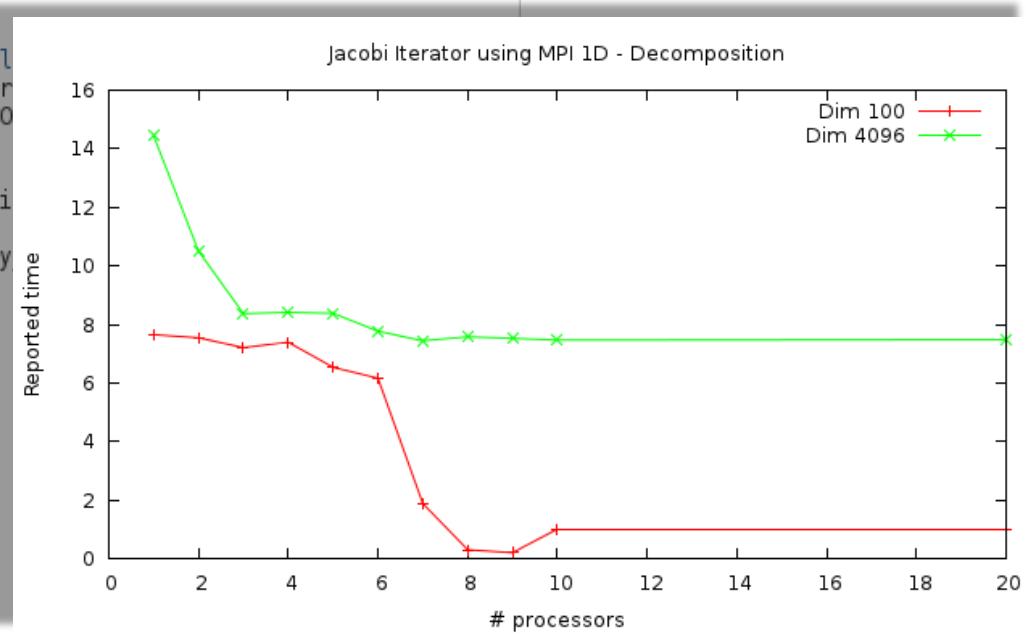


MPI: 1D - Decomposition

```

92 // Once each process has finished, sends a message to the next with the corresponden
t ghost cell
93 int next, prev;
94 next = (my_rank+1)%size;
95 prev = (my_rank-1+size)%size;
96 if(DEBUG)printf("P%d: Sending row %d to P%d\n",my_rank,load[my_rank]+1,next);
97 MPI_Isend(&SurfaceMatrix[load[my_rank]*(Dimension+2)+1], Dimension+2, MPI_DOUBLE, ne
xt, 0, MPI_COMM_WORLD, &request);
98
99 // Receive the message sent by the previous process
100 MPI_Recv(&SurfaceMatrix_t[0], Dimension+2, MPI_DOUBLE, prev, 0, MPI_COMM_WORLD, MPI_
STATUS_IGNORE);
101 if(DEBUG)printf("P%d: Received row %d from P%d\n",my_rank,load[my_rank]+1,prev);
102 // Wait until the message has been delivered and continue to calculate
103 MPI_Wait(&request, MPI_STATUS_IGNORE);
104
105
106
107 // Send a message to the previous process to fil
108 if(DEBUG)printf("P%d: Sending row %d to P%d\n",my_r
109 MPI_Isend(&SurfaceMatrix[0], Dimension+2, MPI_DO
est);
110
111 MPI_Recv(&SurfaceMatrix_t[load[my_rank]*(Dimensi
ext, 0,MPI_COMM_WORLD, MPI_STATUS_IGNORE);
112 if(DEBUG)printf("P%d: Received row 0 from P%d\n",my
113
114 MPI_Wait(&request, MPI_STATUS_IGNORE);
115

```



MPI: 1D - Decomposition

```
143 // ****
144 // *** Master process' section ***
145 if(!my_rank){
146
147     double *CompleteMatrix;
148     CompleteMatrix = (double *) malloc((Dimension+2)*Dimension*sizeof(double));
149
150     for(i=0;i<load[0];i++){
151         for(j=0;j<Dimension+2;j++)
152             CompleteMatrix[i*(Dimension+2)+j]=SurfaceMatrix[(i+1)*(Dimension+2)+j];
153     }
154
155     // Collect the information from everybody but myself
156     for(i=1;i<size;i++){
157         MPI_Recv(&CompleteMatrix[i*load[i]*(Dimension+2)], load[i]*(Dimension+2),
158                  MPI_DOUBLE, i, 1, MPI_COMM_WORLD, MPI_STATUS_IGNORE);
158     }
159
160
161     FILE *outfile;
162     outfile = fopen("jacOut.dat","w");
163
164     fprintf(outfile, "\t");
165     for(i=0;i<Dimension+2;i++){
166         fprintf(outfile, "%d\t",i);
167     }
168     fprintf(outfile, "\n");
169
```