2.6 Using PDGEADD for Data Paralization

Let present a way to achieve parallelization on data level (i.e. data sharing and distribution for processes) using function PDGEADD.

Purpose

PDGEADD adds a matrix to another sub(C) := beta*sub(C) + alpha*op(sub(A)), where sub(C) denotes C(IC:IC+M-1,JC:JC+N-1), and op(X) is one of op(X) = X, or op(X) = XT. Thus, op(sub(A)) denotes

- 1) A(IA:IA+M-1,JA:JA+N-1) if TRANS = 'N',
- 2) A(IA:IA+N-1,JA:JA+M-1)' if TRANS = 'T',
- 3) A(IA:IA+N-1,JA:JA+M-1)' if TRANS = 'C'.

Alpha and beta are scalars, sub(C) and op(sub(A)) are m by n submatrices.

In the other world: to distribute your matrix over process grid (block cyclic 2d distribution) we can do this by means of pdgeadd_ PBLAS routine. This routine computes sum of two matrices C, A: C:=alpha*A+beta*C. Matrices can have different distribution, in particular matrix A can be owned by only one process, thus, setting alpha=1, beta=0 you can simply copy your non-distributed matrix A into distributed matrix C.

Arguments

```
TRANS (global input) CHARACTER*1

On entry, TRANS specifies the form of op(sub(A)) to be used in the matrix addition as follows:

TRANS = 'N' \text{ or 'n'} \text{ op(sub(A))} = \text{sub(A)},
```

TRANS = 'T' or 't' op(sub(A)) = sub(A)',TRANS = 'C' or 'c' op(sub(A)) = sub(A)'.

M (global input) INTEGER

On entry, M specifies the number of rows of the submatrix sub(C) and the number of columns of the submatrix sub(A). M must be at least zero.

N (global input) INTEGER

On entry, N specifies the number of columns of the submatrix sub(C) and the number of rows of the submatrix sub(A). N must be at least zero.

ALPHA (global input) DOUBLE PRECISION

On entry, ALPHA specifies the scalar alpha. When ALPHA is supplied as zero then the local entries of the array A corresponding to the entries of the submatrix sub(A) need not be set on input.

A (local input) DOUBLE PRECISION array

On entry, A is an array of dimension (LLD_A, Ka), where Ka is at least Lc(1, JA+M-1). Before entry, this array contains the local entries of the matrix A.

IA (global input) INTEGER

On entry, IA specifies A's global row index, which points to the beginning of the submatrix sub(A).

JA (global input) INTEGER

On entry, JA specifies A's global column index, which points to the beginning of the submatrix sub(A).

DESCA (global and local input) INTEGER array

On entry, DESCA is an integer array of dimension DLEN_. This s the array

```
descriptor for the matrix A.
BETA
          (global input) DOUBLE PRECISION
        On entry, BETA specifies the scalar beta.
                                                  When BETA is supplied as zero
        then the local entries of the array C corresponding to the entries of the
        submatrix sub(C) neednot be set on input.
C
          (local input/local output) DOUBLE PRECISION array
        On entry, C is an array of dimension (LLD_C, Kc), where Kc is at least Lc( 1,
        JC+N-1). Before entry, this array contains the local entries of the matrix C. On
        exit, the entries of this array corresponding to the local entries of the submatrix
        sub(C) are overwritten by the local entries of the m by n updated submatrix.
IC
          (global input) INTEGER
        On entry, IC specifies C's global row index, which points to the beginning of the
        submatrix sub( C).
JC
          (global input) INTEGER
        On entry, JC specifies C's global column index, which points to the beginning of
        the submatrix sub( C ).
DESCC
          (global and local input) INTEGER array
        On entry, DESCC is an integer array of dimension DLEN_. This is the array
        descriptor for the matrix C.
```

The following example should demonstrate how the pdgeadd routine is used.

Example 2.6.1. This example illustrates the modalities of using the BLACS functions for carrying out the operations of parallelization on data level.

Attention! If the grid of processes is not a square, then the distributions do not correspond to 2D-ciclic algorithm.

```
#include <stdio.h>
#include <string.h>
#include <stdlib.h>
#include "mpi.h"
#include <iostream>
#include <iomanip>
#include <fstream>
#include <sstream>
using namespace std;
#define A(i,j) A[(i)*n+(j)]
#define A_distr(i,j) A_distr[(i)*n+(j)]
#define B(i,j) B[(i)*n+(j]
#define B_distr(i,j) B_distr[(i)*n+(j]
static int MAX( int a, int b ){
       if (a>b) return(a); else return(b);
extern "C"
void Cblacs pinfo( int* mypnum, int* nprocs);
void Cblacs get( int context, int request, int* value);
int Cblacs gridinit( int* context, char * order, int np row, int np col);
void Cblacs gridinfo( int context, int* np row, int* np col, int* my row, int*
my col);
void Cblacs_gridexit( int context);
void Cblacs_barrier(int, const char*);
void Cblacs_exit( int error_code);
void Cblacs pcoord(int, int, int*, int*);
int numroc_( int *n, int *nb, int *iproc, int *isrcproc, int *nprocs);
int indxl2g_(int*, int*, int*, int*);
void descinit (int *desc, int *m, int *n, int *mb, int *nb, int *irsrc, int *icsrc, int
*ictxt, int *lld, int *info);
void pdgeadd (char *TRANS,int *M, int *N, double * ALPHA, double *A,int *IA,int *JA,int
*DESCA, double *BETA, double *C,
              int *IC,int *JC,int *DESCC);
```

```
} // extern "C"
int main(int argc, char **argv) {
// Useful constants
int i one = 1, i zero = 0;
double zero=0.0E+0, one=1.0E+0;
int descA[9], descA distr[9], descB[9], descB distr[9];
int iam, nprocs, nprow, npcol, myrow, mycol;
int m,n,mb,nb,mp,nq,nqrhs, nrhs;
int i, j,mypnum;
int lld, lld distr;
int ictxt, info, lwork;
m=9; n=9;
/*
Where m-number of columns in matrix A, n-number of rows in matrix A
*/
mb=2; nb=2;
nrhs=1;
nprow=2; npcol=2; // Astfel, programul se executa pe 4 procese
double *A, *A distr, *B, *B distr, *work, *tau;
// Part with invoking of ScaLAPACK routines. Initialize process grid, first
Cblacs pinfo(&iam, &nprocs);
Cblacs_get( -1, 0, &ictxt );
Cblacs_gridinit(&ictxt, "R", nprow, npcol );
Cblacs_gridinfo(ictxt, &nprow, &npcol, &myrow, &mycol);
// Matricea A se initializeaza numai pentru procesul cu rankul 0
if ( iam==0 ) {
A = (double*) malloc(m*n*sizeof(double));
B = (double*) malloc(m*nrhs*sizeof(double));
//input matrix A abd B
for(i=0;i<m;i++)
     for(j=0;j<n;j++)
             A[i*n+j] = (10*i+j);
for(i=0;i<m;i++)
for (j=0; j< nrhs; j++)
     B[i*nrhs+j]=i+j;
else{
A = NULL;
B = NULL;
//other processes don't contain parts of A
if (iam==0)
 printf("======= REZULT OF THE PROGRAM %s \n", argv[0]);
  cout << "Global matrix A:\n";</pre>
               for (i = 0; i < m; ++i) {
                  for (j = 0; j < n; ++j) {
                       cout << setw(3) << *(A + n*i + j) << "";
                   cout << "\n";
               }
              cout << endl;
cout << "Global vector B:\n";</pre>
              for (i = 0; i < m; ++i) {
             for (j = 0; j < nrhs; ++j)
                   cout << setw(3) << *(B + i*nrhs+j ) << "";
                   cout << "\n";
               cout << endl;
Cblacs barrier(ictxt, "All");
// Compute dimensions of local part of distributed matrix A distr and B distr
mp = numroc_( &m, &mb, &myrow, &i_zero, &nprow );
nq = numroc_( &n, &nb, &mycol, &i_zero, &npcol );
ngrhs = numroc ( &nrhs, &mb, &mycol, &i zero, &npcol );
A_distr = (double*) malloc( mp*nq*sizeof(double));
```

```
B distr = (double*) malloc( mp*ngrhs*sizeof(double));
// Initialize discriptors (local matrix A is considered as distributed with blocking
parameters
// m, n, i.e. there is only one block - whole matrix A - which is located on process
(0,0)
11d = MAX( numroc ( &m, &n, &myrow, &i zero, &nprow ), 1 );
descinit_(descA, &m, &n, &m, &n, &i_zero, &i_zero, &ictxt, &lld, &info );
descinit (descB, &n, &nrhs, &n, &nrhs, &i zero, &i zero, &ictxt, &lld, &info);
lld distr = MAX( mp, 1 );//lld distr = MAX( nq, 1 );
descinit_(descA_distr, &m, &n, &mb, &nb, &i_zero, &i_zero, &ictxt, &lld_distr, &info);
descinit (descB distr, &m, &nrhs, &mb, &nb, &i zero, &i zero, &ictxt, &lld distr, &info
);
// Call pdgeadd to distribute matrix (i.e. copy A into A_distr and B into B_dist)
pdgeadd ("N", &m, &n, &one, A, &i one, &i one, descA, &zero, A distr, &i one, &i one,
descA distr );
pdgeadd ( "N", &m, &nrhs, &one, B, &i one, &i one, descB, &zero, B distr, &i one,
&i_one, descB_distr);
Here m-number of rows in A distr, n- number of columns in A distr
// print A distr and B distr
for (int id = 0; id < nprocs; ++id)
Cblacs barrier(ictxt, "All");
if (id == iam) {
printf("Local A(%d*%d) on node %d (%d,%d) \n", mp,nq,iam,myrow,mycol);
//cout << "A distr on node " << iam << endl;
for (i = 0; i < mp; i++)
 //for (j = 0; j < nq; j++)
 for (j = 0; j < nq; j++)
 //for (i = 0; i < mp; i++)
 cout << setw(3) << *(A distr+nq*i+j) << " ";//cout << setw(3) << *(A distr+mp*j+i) <<</pre>
" ";
 cout << endl;
}
if (nqrhs > 0) {
//if (mycol==0) {
printf("Local B(%d*%d) on node %d (%d,%d) \n", mp,nqrhs,iam,myrow,mycol);
//cout << "B distr on node " << iam << endl;</pre>
for (i = 0; i < mp; ++i)
for (j = 0; j < nqrhs; ++j)
cout << setw(3) << *(B distr+nqrhs*i+j) << " ";</pre>
cout << endl;
}
cout << endl;
Cblacs barrier(ictxt, "All");
if( iam==0 ){
for(i=0;i<m;i++)
    for(j=0;j<n;j++)
             A[i*m+j]=0;
cout << "Global matrix A (pana la restabilire):\n";</pre>
              for (i = 0; i < m; ++i) {
                  for (j = 0; j < n; ++j) {
                       cout << setw(3) << *(A + n*i + j) << " ";
                  cout << "\n";
               }
              cout << endl;
for(i=0;i<m;i++)
     for(j=0;j<nrhs;j++)</pre>
             B[i*nrhs+j]=-10;
```

```
cout << "Global matrix B (pana la restabilire):\n";</pre>
               for (i = 0; i < m; ++i) {
                   for (j = 0; j < nrhs; ++j) {
                       cout << setw(3) << *(B + nrhs*i + j) << " ";
                   cout << "\n";
               }
               cout << endl;
// Copy result into local matrix (adica "restabilirea matrice A si B)
pdgeadd ( "N", &m, &n, &one, A distr, &i one, &i one, descA distr, &zero, A, &i one,
&i one, descA);
pdgeadd ( "N", &m, &nrhs, &one, B distr, &i one, &i one, descB distr, &zero, B, &i one,
&i one, descB);
// Tipar A si B
if (iam==0)
  {
   cout << "Global matrix A (dupa restabilire):\n";</pre>
               for (i = 0; i < m; ++i) {
                   for (j = 0; j < n; ++j) {
                       cout << setw(3) << *(A + n*i + j) << " ";
                   cout << "\n";
               cout << endl;</pre>
cout << "Global vector B (dupa restabilire):\n";</pre>
               for (i = 0; i < m; ++i) {
              for (j = 0; j < nrhs; ++j)
                   cout << setw(3) << *(B + i*nrhs+j ) << " ";</pre>
                   cout << "\n";
               }
               cout << endl;
free ( A distr ); free ( B distr );
if ( myrow==0 && mycol==0 ) {
free(A);
free(B);
// End of ScaLAPACK part. Exit process grid.
Cblacs_gridexit(ictxt);
Cblacs exit(0);
```

Program results.

```
[MI_gr_TPS1@hpc]$ ./mpiCC ScL -o Example2.6.1.exe Example2.6.1.cpp
[{\tt MI\_gr\_TPS1@hpc}] \$ / opt/openmpi/bin/mpirun -n \ 4 \ -host \ compute-0-0, compute-0-1 \ -host \ compute-0-1, compute
Example2.6.1.exe
Global matrix A:
     0 1 2 3
                                                                                                5
                                                                                                                   6 7
                                                                               4
    10 11 12 13 14 15 16 17 18
    20 21 22 23 24 25 26 27 28
    30 31 32 33 34 35 36 37 38
    40 41 42 43 44 45 46 47 48
    50 51 52 53 54 55 56 57 58
    60 61 62 63 64 65 66 67 68
    70 71 72 73 74 75 76 77 78
    80 81 82 83 84 85 86 87 88
Global vector B:
       0
         1
         2
        3
         4
         5
```

```
6
  7
Local A(5*5) on node 0 (0,0)
0 1 4 5 8
 10 11 14 15 18
 40 41 44 45 48

    50
    51
    54
    55
    58

    80
    81
    84
    85
    88

Local B(5*1) on node 0 (0,0)
 0
 1
 4
 5
 8
Local A(5*4) on node 1 (0,1)
20 21 24 25
 28 30 31 34
 35 38 60 61
 64 65 68 70
 71 74 75 78
Local A(4*5) on node 2 (1,0)
2 3 6 7 12
13 16 17 42 43
46 47 52 53 56
57 82 83 86 87
Local B(4*1) on node 2 (1,0)
 2
 3
 6
 7
Local A(4*4) on node 3 (1,1)
22 23 26 27
 32 33 36 37
 62 63 66 67
 72 73 76 77
Global matrix A (before restoring):

    0
    0
    0
    0
    0
    0
    0
    0
    0

    0
    0
    0
    0
    0
    0
    0
    0
    0

         0 0 0 0 0 0
 0 0
                                      0
 0 0 0 0 0 0 0
                                    0
 0 0 0 0 0 0 0
                                      0
  0 0 0 0 0 0 0
                                      0
  0 0 0 0 0 0 0
                                      0
  0 0 0 0 0 0
                           0 0
                                      0
  0 0
                       0
         0 0 0
                            0 0
Global matrix B (before restoring):
-10
-10
-10
-10
-10
-10
-10
-10
-10
Global matrix A (after restoring):
0 1 2 3 4 5 6 7 8
 10 11 12 13 14 15 16 17 18

    20
    21
    22
    23
    24
    25
    26
    27
    28

    30
    31
    32
    33
    34
    35
    36
    37
    38
```

```
40 41 42 43 44 45 46 47 48
50 51 52 53 54 55 56 57 58
60 61 62 63 64 65 66 67 68
70 71 72 73 74 75 76 77 78
80 81 82 83 84 85 86 87 88

Global vector B (after restoring):
0
1
2
3
4
5
6
7
7
8
```

The following table contains a list of all functions BLACS:

BLACS Routine List

	C Name	Fortran Name	Date	Description
_			type v	
	Cblacs_pinfo	BLACS_PINFO		Get initial system information that is required before BLACS is set up
Initialization				
	Cblacs_setup	BLACS_SETUP		Functionally equivalent to blas_pinfo
	Cblacs_get	BLACS_GET		Returns values BLACS is using for internal defaults
	Cblacs_set	BLACS_SET		Sets BLACS internal defaults
	Cblacs_gridinit	BLACS_GRIDINIT		Assigns processors to BLACS process grid
	Cblacs_gridmap	BLACS_GRIDMAP		Assigns processors to BLACS process grid in arbitrary manner
Destruction	Cblacs_freebuff	BLACS_FREEBUFF		Releases BLACS buffer
	Cblacs_gridexit	BLACS_GRIDEXIT		Frees a BLACS context
	Cblacs_abort	BLACS_ABORT		Aborts all BLACS processes
	Cblacs_exit	BLACS_EXIT		Frees all BLACS contexts and allocated memory
Sending	Cvgesd2d	vGESD2D	SDCZI	General send 2-d
	Cvgebs2d	vGEBS2D	SDCZI	General broadcast send 2-d
	Cvtrsd2d	vTRSD2D	SDCZI	Trapezoidal send 2-d
	Cvtrbs2d	vTRBS2D	SDCZI	Trapezoidal broadcast send 2-d
Receiving	Cvgerv2d	vGERV2D	SDCZI	General receive
	Cvgebr2d	vGEBR2D	SDCZI	General broadcast receive
	Cvtrrv2d	vTRRV2D	SDCZI	Trapezoidal receive
	Cvtrbr2d	vTRBR2D	SDCZI	Trapezoidal broadcast receive

	Cvgamx2d	vGAMX2D	SDCZI	General element-wise absolute value maximum
Combine				value maximum
	Cvgamn2d	vGAMN2D	SDCZI	General element-wise absolute value minimum
	Cvgsum2d	vGSUM2D	SDCZI	General element-wise summation
	Cblacs_gridinfo	BLACS_GRIDINFO		Returns information on BLACS grid
Information and Miscellaneo				
us	Cblacs_pnum	BLACS_PNUM		Returns system process number
	Cblacs_pcoord	BLACS_PCOORD		Returns row and col in BLACS process grid
	Cblacs_barrier	BLACS_BARRIER		Holds up execution of all processes till all processes call this routine
	Csetpvmtids	SETPVMTIDS		PVM routine, not used
Non- Standard				
	Cdcputime00	DCPUTIME00		Returns CPU seconds since arbitrary starting point
	Cdwalltime00	DWALLTIME00		Returns wall clock seconds since arbitrary starting point
	Cksendid	SENDID		Returns BLACS message ID
	Cdrecvid	DRECVID		Returns BLACS message ID for receive
	Ckbsid	KBSID		Returns BLACS message ID for source
	Ckbrid	KBRID		Returns BLACS message ID for destination in broadcast