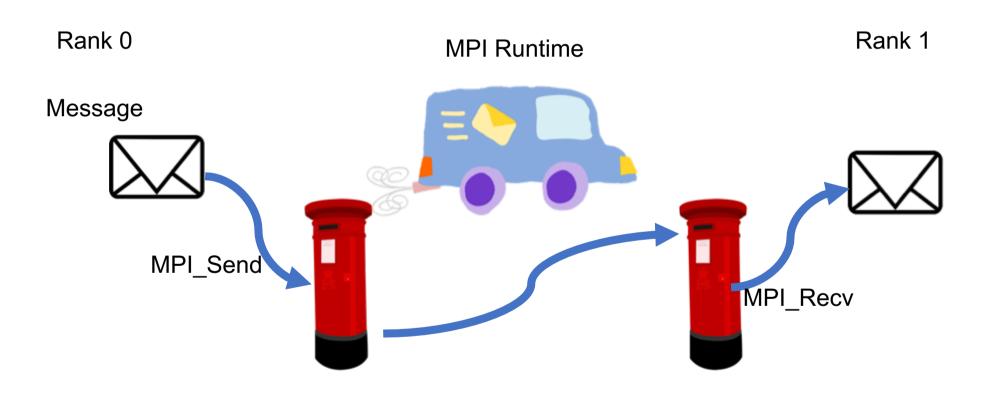
# Message Passing



#### 6 Basic Functions

```
MPI_Init ()
MPI_Comm_rank (Comm)
MPI_Comm_size (Comm)
MPI_Send (Data, Num, Type, Dest, Tag, Comm)
MPI_Recv (Data, Num, Type, Src, Tag, Comm, Stat)
MPI_Finalize ()
```

## Communication type

Point-to-Point communication

Collective communication

Blocking communication

Non-blocking communication

#### Non-blocking communication

```
MPI_Isend (Data, Num, Type, Dest, Tag, Comm, &Req)
MPI_Irecv (Data, Num, Type, Src, Tag, Comm, &Req)
MPI_Wait (&Req, &Status)
MPI_Waitall (Num, Req[], Stat[])
```

#### Collective Communication

```
MPI_Bcast MPI_Ibcast
```

MPI\_Scatter MPI\_Gather

MPI\_Allscatter MPI\_Alltoall

#### **Collective Communication**

rank	send buf		recv buf
0	a,b,c	MPI_Allgather	a,b,c,A,B,C,#,@,%
1	A,B,C	>	a,b,c,A,B,C,#,@,%
2	#,@,%		a,b,c,A,B,C,#,@,%

	and but						
rank	send buf	recv buf					
0	a,b,c MPI_Alltoall	a,A,#					
1	A,B,C>	b,B,@					
2	#,@,%	c, C, %					
_	,	3, 3, 3					
(a more	more elaborate case with two elements per process)						
rank	send buf	recv buf					
0	a,b,c,d,e,f MPI_Alltoall	a,b,A,B,#,@					
1	A,B,C,D,E,F>	c,d,C,D,%,\$					
2	#,@,%,\$,&,*	e,f,E,F,&,*					
_	161-171-1	01.1-1.1-1.					

Rank i wants to send message i to all other ranks

Implementation 1 : Blocking one-to-one:

```
for (int irank = 0; irank < size; ++irank) {
   if (irank == rank) continue;
   int stag = irank + size * rank;
   MPI_Send (&sendbuff, 1, MPI_INT, irank, stag, MPI_COMM_WORLD);
   std::cout << "rank " << rank << " sending message " << sendbuff << " to rank " << irank << std::endl;
   int rtag = rank + size * irank;
   MPI_Recv (&(recvbuff[irank]), 1, MPI_INT, irank, rtag, MPI_COMM_WORLD, MPI_STATUS_IGNORE);
}</pre>
```

Notice encoding of sender and receiver in tag!

**DANGER: Deadlock!** 

Rank i wants to send message i to all other ranks

Implementation 2 : Non-Blocking one-to-one:

```
for (int irank = 0; irank < size; ++irank) {
   if (irank == rank) continue;
   int stag = irank + size * rank;
   MPI_Isend (&sendbuff, 1, MPI_INT, irank, stag, MPI_COMM_WORLD, &(reqs[numreq++]));
   std::cout << "rank " << rank << " sending message " << sendbuff << " to rank " << irank << std::endl;
   int rtag = rank + size * irank;
   MPI_Irecv (&(recvbuff[irank]), 1, MPI_INT, irank, rtag, MPI_COMM_WORLD, &(reqs[numreq++]));
}
MPI_Waitall (numreq, &(reqs[0]), MPI_STATUSES_IGNORE);</pre>
```

Notice encoding of sender and receiver in tag!

Notice Use of Waitall!

Rank i wants to send message i to all other ranks

Implementation 3 : Blocking Collective (All-to-all)

```
recvbuff.assign (recvbuff.size (), rank);
MPI_Alltoall (MPI_IN_PLACE, 1, MPI_INT, &(recvbuff[0]), 1, MPI_INT, MPI_COMM_WORLD);
```

Notice MPI\_IN\_PLACE

Very convenient when messages of same known size are to be sent / received!

Rank i wants to send message i to all other ranks

Implementation 4 : Blocking Collective (All-gather)

```
recvbuff.assign (recvbuff.size (), rank);
MPI_Allgather (&sendbuff, 1, MPI_INT, &(recvbuff[0]), 1, MPI_INT, MPI_COMM_WORLD);
```

Very convenient when messages of same known size are to be sent / received!

#### Advanced Example

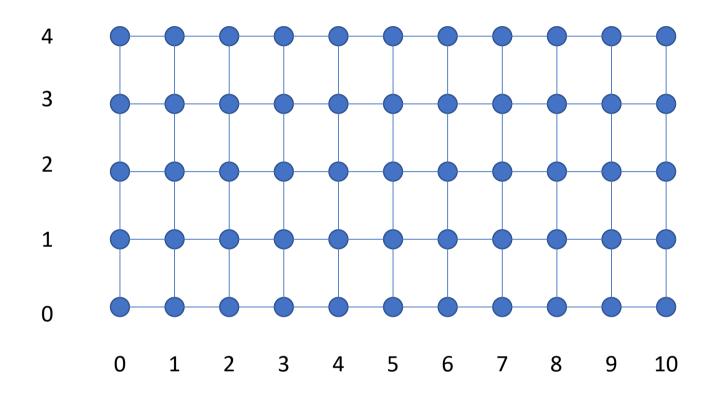
#### Distributed assembly of sparse\_matrix:

- Each rank owns a contiguous range of rows
- Each rank can assign elements within its ownership range or outside its ownership range
- If multiple ranks assign one matrix element, after assembly the actual value must be the sum of local contributions

This is the most ommon situation encountered in **Finite Element** problems

- Each rank owns a range of mesh elements and a range of DOFs
- For low order continuous formulations **DOFs** correspond to mesh **vertices**
- The same vertex touches multiple **elements**
- The matrix entries related to a given DOF are the sum of contributions from all elements touching the corresponding vertex

## Exercise: assemble FEM Laplacian



	10	J1	J2	J3
10	2	-1	-1	
I1	-1	2		-1
12	-1		2	-1
13		-1	-1	2

## Exercise: assemble FEM Laplacian

