hypre Reference Manual

— Version 2.11.1 —

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. 1

Struct System Interface

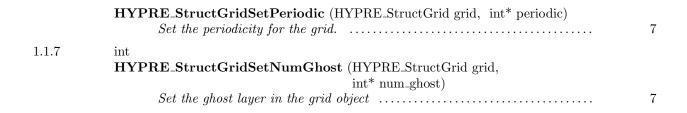
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This interface represents a structured-grid conceptual view of a linear system.

_ 1.1 _

Struct Grids

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typedef struct hypre_StructGrid_struct *HYPRE_StructGrid

A grid object is constructed out of several "boxes", defined on a global abstract index space

1.1.2 _

HYPRE_StructGridCreate (MPI_Comm comm, int ndim, HYPRE_StructGrid* grid)

Create an ndim-dimensional grid object

1.1.3

int HYPRE_StructGridDestroy (HYPRE_StructGrid grid)

Destroy a grid object. An object should be explicitly destroyed using this destructor when the user's code no longer needs direct access to it. Once destroyed, the object must not be referenced again. Note that the object may not be deallocated at the completion of this call, since there may be internal package references to the object. The object will then be destroyed when all internal reference counts go to zero.

1.1.4

int **HYPRE_StructGridSetExtents** (HYPRE_StructGrid grid, int* ilower, int* iupper)

Set the extents for a box on the grid

_ 1.1.5 _

int HYPRE_StructGridAssemble (HYPRE_StructGrid grid)

Finalize the construction of the grid before using

_ 1.1.6 _

int HYPRE_StructGridSetPeriodic (HYPRE_StructGrid grid, int* periodic)

Set the periodicity for the grid.

The argument periodic is an ndim-dimensional integer array that contains the periodicity for each dimension. A zero value for a dimension means non-periodic, while a nonzero value means periodic and contains the actual period. For example, periodicity in the first and third dimensions for a 10x11x12 grid is indicated by the array [10,0,12].

NOTE: Some of the solvers in hypre have power-of-two restrictions on the size of the periodic dimensions.

_ 1.1.7 _

int HYPRE_StructGridSetNumGhost (HYPRE_StructGrid grid, int* num_ghost)

Set the ghost layer in the grid object

1.2

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	Set a stencil entry.	9

1.2.1

 $typedef\ struct\ hypre_StructStencil_struct\ *HYPRE_StructStencil$

The stencil object

1.2.2

int

HYPRE_StructStencilCreate (int ndim, int size, HYPRE_StructStencil* stencil)

Create a stencil object for the specified number of spatial dimensions and stencil entries

1.2.3

int HYPRE_StructStencilDestroy (HYPRE_StructStencil stencil)

Destroy a stencil object

1.2.4

int

HYPRE_StructStencilSetElement (HYPRE_StructStencil stencil, int entry, int* offset)

Set a stencil entry.

NOTE: The name of this routine will eventually be changed to ${\tt HYPRE_StructStencilSetEntry}.$

1.3

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	HYPRE_StructVector x,	
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 $typedef\ struct\ hypre_StructMatrix_struct\ *HYPRE_StructMatrix$

The matrix object

_ 1.3.2 _

int

HYPRE_StructMatrixCreate (MPI_Comm comm, HYPRE_StructGrid grid, HYPRE_StructStencil stencil, HYPRE_StructMatrix* matrix)

Create a matrix object

_ 1.3.3 _

int HYPRE_StructMatrixDestroy (HYPRE_StructMatrix matrix)

Destroy a matrix object

int **HYPRE_StructMatrixInitialize** (HYPRE_StructMatrix matrix)

Prepare a matrix object for setting coefficient values

_ 1.3.5 __

int

HYPRE_StructMatrixSetValues (HYPRE_StructMatrix matrix, int* index, int nentries, int* entries, HYPRE_Complex* values)

Set matrix coefficients index by index. The values array is of length nentries.

NOTE: For better efficiency, use HYPRE_StructMatrixSetBoxValues to set coefficients a box at a time.

1.3.6

int

HYPRE_StructMatrixAddToValues (HYPRE_StructMatrix matrix, int* index, int nentries, int* entries, HYPRE_Complex* values)

Add to matrix coefficients index by index. The values array is of length nentries.

NOTE: For better efficiency, use HYPRE_StructMatrixAddToBoxValues to set coefficients a box at a time.

_ 1.3.7 _

int

HYPRE_StructMatrixSetConstantValues (HYPRE_StructMatrix matrix, int nentries, int* entries, HYPRE_Complex* values)

Set matrix coefficients which are constant over the grid. The values array is of length nentries.

int

 $\label{lem:hypre_structMatrix} \begin{tabular}{ll} HYPRE_StructMatrix Matrix, int nentries, int* entries, HYPRE_Complex* values) \end{tabular}$

Add to matrix coefficients which are constant over the grid. The values array is of length nentries.

__ 1.3.9 __

int

HYPRE_StructMatrixSetBoxValues (HYPRE_StructMatrix matrix, int* ilower, int* iupper, int nentries, int* entries, HYPRE_Complex* values)

Set matrix coefficients a box at a time. The data in values is ordered as follows:

```
m = 0;
for (k = ilower[2]; k <= iupper[2]; k++)
    for (j = ilower[1]; j <= iupper[1]; j++)
        for (i = ilower[0]; i <= iupper[0]; i++)
            for (entry = 0; entry < nentries; entry++)
        {
            values[m] = ...;
            m++;
        }
}</pre>
```

1.3.10

int

HYPRE_StructMatrixAddToBoxValues (HYPRE_StructMatrix matrix, int* ilower, int* iupper, int nentries, int* entries, HYPRE_Complex* values)

Add to matrix coefficients a box at a time. The data in values is ordered as in HYPRE_StructMatrixSetBoxValues.

int HYPRE_StructMatrixAssemble (HYPRE_StructMatrix matrix)

Finalize the construction of the matrix before using

___ 1.3.12 ____

int

HYPRE_StructMatrixGetValues (HYPRE_StructMatrix matrix, int* index, int nentries, int* entries, HYPRE_Complex* values)

Get matrix coefficients index by index. The values array is of length nentries.

NOTE: For better efficiency, use HYPRE_StructMatrixGetBoxValues to get coefficients a box at a time.

_ 1.3.13 _____

int

HYPRE_StructMatrixGetBoxValues (HYPRE_StructMatrix matrix, int* ilower, int* iupper, int nentries, int* entries, HYPRE_Complex* values)

Get matrix coefficients a box at a time. The data in values is ordered as in HYPRE_StructMatrixSetBoxValues.

1.3.14

int

 $\label{eq:hypre_structMatrixSetSymmetric} \textbf{(HYPRE_StructMatrix matrix, int symmetric)}$

Define symmetry properties of the matrix. By default, matrices are assumed to be nonsymmetric. Significant storage savings can be made if the matrix is symmetric.

int

 $\label{thm:hypre_structMatrixSetConstantEntries} \mbox{ (HYPRE_StructMatrix matrix, int nentries, int* entries)}$

Specify which stencil entries are constant over the grid. Declaring entries to be "constant over the grid" yields significant memory savings because the value for each declared entry will only be stored once. However, not all solvers are able to utilize this feature.

Presently supported:

- no entries constant (this function need not be called)
- all entries constant
- all but the diagonal entry constant

_ 1.3.16 _

ınt

HYPRE_StructMatrixSetNumGhost (HYPRE_StructMatrix matrix, int* num_ghost)

Set the ghost layer in the matrix

_ 1.3.17 __

int

HYPRE_StructMatrixPrint (const char* filename, HYPRE_StructMatrix matrix, int all)

Print the matrix to file. This is mainly for debugging purposes.

int

HYPRE_StructMatrixMatvec (HYPRE_Complex alpha, HYPRE_StructMatrix A, HYPRE_StructVector x, HYPRE_Complex beta, HYPRE_StructVector y)

Matvec operator. This operation is $y = \alpha Ax + \beta y$. Note that you can do a simple matrix-vector multiply by setting $\alpha = 1$ and $\beta = 0$.

_ 1.4 _

Struct Vectors

Names		
1.4.1	typedef struct hypre_StructVector_struct *HYPRE_StructVector The vector object	17
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1.4.4	int	
	HYPRE_StructVectorInitialize (HYPRE_StructVector vector)	
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1.4.5	int	
	HYPRE_StructVectorSetValues (HYPRE_StructVector vector, int* index, HYPRE_Complex value)	
	Set vector coefficients index by index.	18
1.4.6	int	
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	Add to vector coefficients index by index	18
1.4.7	int	
	HYPRE_StructVectorSetBoxValues (HYPRE_StructVector vector, int* ilower, int* iupper, HYPRE_Complex* values)	
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1.4.9	int	

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1.4.10	int	
	HYPRE_StructVectorGetValues (HYPRE_StructVector vector, int* index,	
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	Get vector coefficients index by index.	19
1.4.11	int	
	HYPRE_StructVectorGetBoxValues (HYPRE_StructVector vector,	
	int* ilower, int* iupper,	
	HYPRE_Complex* values)	
	Get vector coefficients a box at a time.	20
1.4.12	int	
	HYPRE_StructVectorPrint (const char* filename,	
	HYPRE_StructVector vector, int all)	
	Print the vector to file	20

1.4.1

typedef struct hypre_StructVector_struct *HYPRE_StructVector

The vector object

1.4.2

HYPRE_StructVectorCreate (MPI_Comm comm, HYPRE_StructGrid grid, HYPRE_StructVector* vector)

Create a vector object

1.4.3

int HYPRE_StructVectorDestroy (HYPRE_StructVector vector)

Destroy a vector object

1.4.4

int HYPRE_StructVectorInitialize (HYPRE_StructVector vector)

Prepare a vector object for setting coefficient values

__ 1.4.5 ____

HYPRE_StructVectorSetValues (HYPRE_StructVector vector, int* index, HYPRE_Complex value)

Set vector coefficients index by index.

NOTE: For better efficiency, use HYPRE_StructVectorSetBoxValues to set coefficients a box at a time.

1.4.6

HYPRE_StructVectorAddToValues (HYPRE_StructVector vector, int* index, HYPRE_Complex value)

Add to vector coefficients index by index.

NOTE: For better efficiency, use HYPRE_StructVectorAddToBoxValues to set coefficients a box at a time.

_ 1.4.7 ____

HYPRE_StructVectorSetBoxValues (HYPRE_StructVector vector, int* ilower, int* iupper, HYPRE_Complex* values)

Set vector coefficients a box at a time. The data in values is ordered as follows:

```
m = 0;
for (k = ilower[2]; k <= iupper[2]; k++)
  for (j = ilower[1]; j <= iupper[1]; j++)
    for (i = ilower[0]; i <= iupper[0]; i++)
    {
      values[m] = ...;
      m++;
    }</pre>
```

_ 1.4.8 _

int

HYPRE_StructVectorAddToBoxValues (HYPRE_StructVector vector, int* ilower, int* iupper, HYPRE_Complex* values)

Add to vector coefficients a box at a time. The data in values is ordered as in HYPRE_StructVectorSetBoxValues.

1.4.9

int HYPRE_StructVectorAssemble (HYPRE_StructVector vector)

Finalize the construction of the vector before using

1.4.10

HYPRE_StructVectorGetValues (HYPRE_StructVector vector, int* index, HYPRE_Complex* value)

Get vector coefficients index by index.

NOTE: For better efficiency, use HYPRE_StructVectorGetBoxValues to get coefficients a box at a time.

1.4.11

HYPRE_StructVectorGetBoxValues (HYPRE_StructVector vector, int* ilower, int* iupper, HYPRE_Complex* values)

Get vector coefficients a box at a time. The data in values is ordered as in HYPRE_StructVectorSetBoxValues.

___ 1.4.12 _____

HYPRE_StructVectorPrint (const char* filename, HYPRE_StructVector vector, int all)

Print the vector to file. This is mainly for debugging purposes.

2

SStruct System Interface

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This interface represents a semi-structured-grid conceptual view of a linear system.

2.1

SStruct Grids

Names		
2.1.1	typedef struct hypre_SStructGrid_struct *HYPRE_SStructGrid A grid object is constructed out of several structured "parts" and an optional unstructured "part".	23
2.1.2	typedef int HYPRE_SStructVariable An enumerated type that supports cell centered, node centered, face centered, and edge centered variables	23
2.1.3	int HYPRE_SStructGridCreate (MPI_Comm comm, int ndim, int nparts,	
	HYPRE_SStructGrid* grid) Create an ndim-dimensional grid object with nparts structured parts	24
2.1.4	int HYPRE_SStructGridDestroy (HYPRE_SStructGrid grid) Destroy a grid object.	24
2.1.5	int HYPRE_SStructGridSetExtents (HYPRE_SStructGrid grid, int part, int* ilower, int* iupper) Set the extents for a box on a structured part of the grid	24
2.1.6	int	

	HYPRE_SStructGridSetVariables (HYPRE_SStructGrid grid, int part,	
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	HYPRE_SStructVariable* vartypes)	
	Describe the variables that live on a structured part of the grid	25
2.1.7	int	
	HYPRE_SStructGridAddVariables (HYPRE_SStructGrid grid, int part,	
	int* index, int nvars,	
	HYPRE_SStructVariable* vartypes)	
	Describe additional variables that live at a particular index	25
2.1.8	int	
	HYPRE_SStructGridSetFEMOrdering (HYPRE_SStructGrid grid, int part,	
	int* ordering)	
	Set the ordering of variables in a finite element problem	25
2.1.9	int	
2.1.0	HYPRE_SStructGridSetNeighborPart (HYPRE_SStructGrid grid, int part,	
	int* ilower, int* iupper,	
	int nbor_part, int* nbor_ilower,	
	int* nbor_iupper, int* index_map,	
	$int^* index_dir)$	
	Describe how regions just outside of a part relate to other parts	26
2.1.10	int	
	HYPRE_SStructGridSetSharedPart (HYPRE_SStructGrid grid, int part,	
	int* ilower, int* iupper, int* offset,	
	int shared_part, int* shared_ilower,	
	int* shared_iupper, int* shared_offset,	
	int* index_map, int* index_dir)	
	Describe how regions inside a part are shared with regions in other parts.	26
2.1.11	int	
	HYPRE_SStructGridAddUnstructuredPart (HYPRE_SStructGrid grid,	
	int ilower, int iupper)	
	Add an unstructured part to the grid	27
2.1.12	int	
2.1.12	HYPRE_SStructGridAssemble (HYPRE_SStructGrid grid)	
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	Set the periodicity on a particular part.	28
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2.1.1

typedef struct hypre_SStructGrid_struct *HYPRE_SStructGrid

A grid object is constructed out of several structured "parts" and an optional unstructured "part". Each structured part has its own abstract index space.

2.1.2

typedef int HYPRE_SStructVariable

An enumerated type that supports cell centered, node centered, face centered, and edge centered variables. Face centered variables are split into x-face, y-face, and z-face variables, and edge centered variables are split into x-edge, y-edge, and z-edge variables. The edge centered variable types are only used in 3D. In 2D, edge centered variables are handled by the face centered types.

Variables are referenced relative to an abstract (cell centered) index in the following way:

- cell centered variables are aligned with the index;
- node centered variables are aligned with the cell corner at relative index (1/2, 1/2, 1/2);
- x-face, y-face, and z-face centered variables are aligned with the faces at relative indexes (1/2, 0, 0), (0, 1/2, 0), and (0, 0, 1/2), respectively;
- x-edge, y-edge, and z-edge centered variables are aligned with the edges at relative indexes (0, 1/2, 1/2), (1/2, 0, 1/2), and (1/2, 1/2, 0), respectively.

The supported identifiers are:

- HYPRE_SSTRUCT_VARIABLE_CELL
- HYPRE_SSTRUCT_VARIABLE_NODE
- HYPRE_SSTRUCT_VARIABLE_XFACE
- HYPRE_SSTRUCT_VARIABLE_YFACE
- HYPRE_SSTRUCT_VARIABLE_ZFACE
- HYPRE_SSTRUCT_VARIABLE_XEDGE
- HYPRE_SSTRUCT_VARIABLE_YEDGE
- HYPRE_SSTRUCT_VARIABLE_ZEDGE

NOTE: Although variables are referenced relative to a unique abstract cell-centered index, some variables are associated with multiple grid cells. For example, node centered variables in 3D are associated with 8 cells (away from boundaries). Although grid cells are distributed uniquely to different processes, variables may be owned by multiple processes because they may be associated with multiple cells.

 $_$ 2.1.3 $_$

HYPRE_SStructGridCreate (MPI_Comm comm, int ndim, int nparts, HYPRE_SStructGrid* grid)

Create an ndim-dimensional grid object with nparts structured parts

 $_$ 2.1.4 $_$

int HYPRE_SStructGridDestroy (HYPRE_SStructGrid grid)

Destroy a grid object. An object should be explicitly destroyed using this destructor when the user's code no longer needs direct access to it. Once destroyed, the object must not be referenced again. Note that the object may not be deallocated at the completion of this call, since there may be internal package references to the object. The object will then be destroyed when all internal reference counts go to zero.

 $_{-}$ 2.1.5 $_{-}$

HYPRE_SStructGridSetExtents (HYPRE_SStructGrid grid, int part, int* ilower, int* iupper)

Set the extents for a box on a structured part of the grid

2.1.6

HYPRE_SStructGridSetVariables (HYPRE_SStructGrid grid, int part, int nvars, HYPRE_SStructVariable* vartypes)

Describe the variables that live on a structured part of the grid

2.1.7

HYPRE_SStructGridAddVariables (HYPRE_SStructGrid grid, int part, int* index, int nvars, HYPRE_SStructVariable* vartypes)

Describe additional variables that live at a particular index. These variables are appended to the array of variables set in HYPRE_SStructGridSetVariables, and are referenced as such.

NOTE: This routine is not yet supported.

2.1.8

int **HYPRE_SStructGridSetFEMOrdering** (HYPRE_SStructGrid grid, int part, int* ordering)

Set the ordering of variables in a finite element problem. This overrides the default ordering described below.

Array ordering is composed of blocks of size (1 + ndim). Each block indicates a specific variable in the element and the ordering of the blocks defines the ordering of the variables. A block contains a variable number followed by an offset direction relative to the element's center. For example, a block containing (2, 1, -1, 0) means variable 2 on the edge located in the (1, -1, 0) direction from the center of the element. Note that here variable 2 must be of type ZEDGE for this to make sense. The ordering array must account for all variables in the element. This routine can only be called after HYPRE_SStructGridSetVariables.

The default ordering for element variables (var, i, j, k) varies fastest in index i, followed by j, then k, then var. For example, if var 0, var 1, and var 2 are declared to be XFACE, YFACE, and NODE variables, respectively, then the default ordering (in 2D) would first list the two XFACE variables, then the two YFACE variables, then the four NODE variables as follows:

(0,-1,0), (0,1,0), (1,0,-1), (1,0,1), (2,-1,-1), (2,1,-1), (2,-1,1), (2,1,1)

2.1.9

int

HYPRE_SStructGridSetNeighborPart (HYPRE_SStructGrid grid, int part, int* ilower, int* iupper, int nbor_part, int* nbor_ilower, int* nbor_iupper, int* index_map, int* index_dir)

Describe how regions just outside of a part relate to other parts. This is done a box at a time.

Parts part and nbor_part must be different, except in the case where only cell-centered data is used.

Indexes should increase from ilower to iupper. It is not necessary that indexes increase from nbor_ilower to nbor_iupper.

The index_map describes the mapping of indexes 0, 1, and 2 on part part to the corresponding indexes on part nbor_part. For example, triple (1, 2, 0) means that indexes 0, 1, and 2 on part part map to indexes 1, 2, and 0 on part nbor_part, respectively.

The index_dir describes the direction of the mapping in index_map. For example, triple (1, 1, -1) means that for indexes 0 and 1, increasing values map to increasing values on nbor_part, while for index 2, decreasing values map to increasing values.

NOTE: All parts related to each other via this routine must have an identical list of variables and variable types. For example, if part 0 has only two variables on it, a cell centered variable and a node centered variable, and we declare part 1 to be a neighbor of part 0, then part 1 must also have only two variables on it, and they must be of type cell and node. In addition, variables associated with FACEs or EDGEs must be grouped together and listed in X, Y, Z order. This is to enable the code to correctly associate variables on one part with variables on its neighbor part when a coordinate transformation is specified. For example, an XFACE variable on one part may correspond to a YFACE variable on a neighbor part under a particular transformation, and the code determines this association by assuming that the variable lists are as noted here.

2.1.10

int

HYPRE_SStructGridSetSharedPart (HYPRE_SStructGrid grid, int part, int* ilower, int* iupper, int* offset, int shared_part, int* shared_ilower, int* shared_ilower, int* shared_offset, int* index_map, int* index_dir)

Describe how regions inside a part are shared with regions in other parts.

Parts part and shared_part must be different.

Indexes should increase from ilower to iupper. It is not necessary that indexes increase from shared_ilower to shared_iupper. This is to maintain consistency with the SetNeighborPart function, which is also able

to describe shared regions but in a more limited fashion.

The offset is a triple (in 3D) used to indicate the dimensionality of the shared set of data and its position with respect to the box extents ilower and iupper on part part. The dimensionality is given by the number of 0's in the triple, and the position is given by plus or minus 1's. For example: (0, 0, 0) indicates sharing of all data in the given box; (1, 0, 0) indicates sharing of data on the faces in the (1, 0, 0) direction; (1, 0, -1) indicates sharing of data on the edges in the (1, 0, -1) direction; and (1, -1, 1) indicates sharing of data on the nodes in the (1, -1, 1) direction. To ensure the dimensionality, it is required that for every nonzero entry, the corresponding extents of the box are the same. For example, if offset is (0, 1, 0), then (2, 1, 3) and (10, 1, 15) are valid box extents, whereas (2, 1, 3) and (10, 7, 15) are invalid (because 1 and 7 are not the same).

The shared_offset is used in the same way as offset, but with respect to the box extents shared_ilower and shared_iupper on part shared_part.

The index_map describes the mapping of indexes 0, 1, and 2 on part part to the corresponding indexes on part shared_part. For example, triple (1, 2, 0) means that indexes 0, 1, and 2 on part part map to indexes 1, 2, and 0 on part shared_part, respectively.

The index_dir describes the direction of the mapping in index_map. For example, triple (1, 1, -1) means that for indexes 0 and 1, increasing values map to increasing values on shared_part, while for index 2, decreasing values map to increasing values.

NOTE: All parts related to each other via this routine must have an identical list of variables and variable types. For example, if part 0 has only two variables on it, a cell centered variable and a node centered variable, and we declare part 1 to have shared regions with part 0, then part 1 must also have only two variables on it, and they must be of type cell and node. In addition, variables associated with FACEs or EDGEs must be grouped together and listed in X, Y, Z order. This is to enable the code to correctly associate variables on one part with variables on a shared part when a coordinate transformation is specified. For example, an XFACE variable on one part may correspond to a YFACE variable on a shared part under a particular transformation, and the code determines this association by assuming that the variable lists are as noted here.

2.1.11

HYPRE_SStructGridAddUnstructuredPart (HYPRE_SStructGrid grid, int ilower, int iupper)

Add an unstructured part to the grid. The variables in the unstructured part of the grid are referenced by a global rank between 0 and the total number of unstructured variables minus one. Each process owns some unique consecutive range of variables, defined by ilower and iupper.

NOTE: This is just a placeholder. This part of the interface is not finished.

2.1.12

int HYPRE_SStructGridAssemble (HYPRE_SStructGrid grid)

Finalize the construction of the grid before using

_ 2.1.13 ____

HYPRE_SStructGridSetPeriodic (HYPRE_SStructGrid grid, int part, int* periodic)

Set the periodicity on a particular part.

The argument periodic is an ndim-dimensional integer array that contains the periodicity for each dimension. A zero value for a dimension means non-periodic, while a nonzero value means periodic and contains the actual period. For example, periodicity in the first and third dimensions for a 10x11x12 part is indicated by the array [10,0,12].

NOTE: Some of the solvers in hypre have power-of-two restrictions on the size of the periodic dimensions.

2.1.14

int **HYPRE_SStructGridSetNumGhost** (HYPRE_SStructGrid grid, int* num_ghost)

Setting ghost in the sgrids

2.2

SStruct Stencils

Names

2.2.1 typedef struct hypre_SStructStencil_struct *HYPRE_SStructStencil

	The stencil object	29
2.2.2	int	
	HYPRE_SStructStencilCreate (int ndim, int size,	
	HYPRE_SStructStencil* stencil)	
	Create a stencil object for the specified number of spatial dimensions and	
	stencil entries	29
2.2.3	int	
	HYPRE_SStructStencilDestroy (HYPRE_SStructStencil stencil)	
	Destroy a stencil object	29
2.2.4	int	
	HYPRE_SStructStencilSetEntry (HYPRE_SStructStencil stencil, int entry,	
	int* offset, int var)	
	Set a stencil entru	30

2.2.1

typedef struct hypre_SStructStencil_struct *HYPRE_SStructStencil

The stencil object

2.2.2

HYPRE_SStructStencilCreate (int ndim, int size, HYPRE_SStructStencil* stencil)

Create a stencil object for the specified number of spatial dimensions and stencil entries

 $_$ 2.2.3 $_$

int HYPRE_SStructStencilDestroy (HYPRE_SStructStencil stencil)

 ${\bf Destroy}\ {\bf a}\ {\bf stencil}\ {\bf object}$

2.2.4

HYPRE_SStructStencilSetEntry (HYPRE_SStructStencil stencil, int entry, int* offset, int var)

Set a stencil entry

__ 2.3 _

SStruct Graphs

Names		
2.3.1	$typedef\ struct\ hypre_SStructGraph_struct\ *HYPRE_SStructGraph$	
	The graph object is used to describe the nonzero structure of a matrix	31
2.3.2	int	
	HYPRE_SStructGraphCreate (MPI_Comm comm,	
	${ m HYPRE_SStructGrid\ grid},$	
	HYPRE_SStructGraph* graph)	
	Create a graph object	31
2.3.3	int	
	HYPRE_SStructGraphDestroy (HYPRE_SStructGraph graph)	
	Destroy a graph object	31
2.3.4	int	
	$\mathbf{HYPRE_SStructGraphSetDomainGrid} \ (\mathbf{HYPRE_SStructGraph} \ \mathbf{graph},$	
	HYPRE_SStructGrid domain_grid)	
	Set the domain grid	32
2.3.5	int	
	HYPRE_SStructGraphSetStencil (HYPRE_SStructGraph graph, int part,	
	int var, HYPRE_SStructStencil stencil)	
	Set the stencil for a variable on a structured part of the grid	32
2.3.6	int	
	HYPRE_SStructGraphSetFEM (HYPRE_SStructGraph graph, int part)	
	Indicate that an FEM approach will be used to set matrix values on this part	
		32
2.3.7	int	
	HYPRE_SStructGraphSetFEMSparsity (HYPRE_SStructGraph graph,	
	int part, int nsparse,	
	$int^* sparsity)$	
	Set the finite element stiffness matrix sparsity.	32
2.3.8	int	

	HYPRE_SStructGraphAddEntries (HYPRE_SStructGraph graph, int part, int* index, int var, int to_part,	
	int* to_index, int to_var)	
	Add a non-stencil graph entry at a particular index	33
2.3.9	int	
	HYPRE_SStructGraphAssemble (HYPRE_SStructGraph graph)	
	Finalize the construction of the graph before using	33
2.3.10	int	
	HYPRE_SStructGraphSetObjectType (HYPRE_SStructGraph graph,	
	int type)	
	Set the storage type of the associated matrix object	33

 $typedef\ struct\ hypre_SStructGraph_struct\ *HYPRE_SStructGraph$

The graph object is used to describe the nonzero structure of a matrix

2.3.2

HYPRE_SStructGraphCreate (MPI_Comm comm, HYPRE_SStructGrid grid, HYPRE_SStructGraph* graph)

Create a graph object

 $_$ 2.3.3 $_$

int HYPRE_SStructGraphDestroy (HYPRE_SStructGraph graph)

Destroy a graph object

int
HYPRE_SStructGraphSetDomainGrid (HYPRE_SStructGraph graph,
HYPRE_SStructGrid domain_grid)

Set the domain grid

2.3.5

int

HYPRE_SStructGraphSetStencil (HYPRE_SStructGraph graph, int part, int var, HYPRE_SStructStencil stencil)

Set the stencil for a variable on a structured part of the grid

_ 2.3.6 _

int HYPRE_SStructGraphSetFEM (HYPRE_SStructGraph graph, int part)

Indicate that an FEM approach will be used to set matrix values on this part

2.3.7

HYPRE_SStructGraphSetFEMSparsity (HYPRE_SStructGraph graph, int part, int nsparse, int* sparsity)

Set the finite element stiffness matrix sparsity. This overrides the default full sparsity pattern described below.

Array sparsity contains nsparse row/column tuples (I,J) that indicate the nonzeroes of the local stiffness matrix. The layout of the values passed into the routine HYPRE_SStructMatrixAddFEMValues is determined here.

The default sparsity is full (each variable is coupled to all others), and the values passed into the routine HYPRE_SStructMatrixAddFEMValues are assumed to be by rows (that is, column indices vary fastest).

2.3.8

int
HYPRE_SStructGraphAddEntries (HYPRE_SStructGraph graph, int part, int* index, int var, int to_part, int* to_index, int to_var)

Add a non-stencil graph entry at a particular index. This graph entry is appended to the existing graph entries, and is referenced as such.

NOTE: Users are required to set graph entries on all processes that own the associated variables. This means that some data will be multiply defined.

 $_$ 2.3.9 $_$

int HYPRE_SStructGraphAssemble (HYPRE_SStructGraph graph)

Finalize the construction of the graph before using

2.3.10

mt HYPRE_SStructGraphSetObjectType (HYPRE_SStructGraph graph, int type)

Set the storage type of the associated matrix object. It is used before AddEntries and Assemble to compute the right ranks in the graph.

NOTE: This routine is only necessary for implementation reasons, and will eventually be removed.

See Also:

HYPRE_SStructMatrixSetObjectType ($\rightarrow 2.4.16$, page 40)

2 4

SStruct Matrices

Names		
2.4.1	typedef struct hypre_SStructMatrix_struct *HYPRE_SStructMatrix The matrix object	36
2.4.2	int	
	HYPRE_SStructMatrixCreate (MPI_Comm comm,	
	HYPRE_SStructGraph graph,	
	HYPRE_SStructMatrix* matrix)	
	Create a matrix object	36
2.4.3	int	
	HYPRE_SStructMatrixDestroy (HYPRE_SStructMatrix matrix)	
	Destroy a matrix object	36
2.4.4	int	
2.4.4	HYPRE_SStructMatrixInitialize (HYPRE_SStructMatrix matrix)	
	Prepare a matrix object for setting coefficient values	36
		0.
2.4.5	int	
	HYPRE_SStructMatrixSetValues (HYPRE_SStructMatrix matrix, int part,	
	int* index, int var, int nentries,	
	int* entries, HYPRE_Complex* values)	37
	Set matrix coefficients index by index	3
2.4.6	int	
	$\mathbf{HYPRE_SStructMatrixAddToValues} \ (\mathbf{HYPRE_SStructMatrix} \ \mathbf{matrix},$	
	int part, int* index, int var,	
	int nentries, int* entries,	
	HYPRE_Complex* values)	0.5
	Add to matrix coefficients index by index	37
2.4.7	int	
	$\mathbf{HYPRE_SStructMatrixAddFEMValues} \ (\mathbf{HYPRE_SStructMatrix} \ \mathbf{matrix},$	
	int part, int* index,	
	HYPRE_Complex* values)	
	Add finite element stiffness matrix coefficients index by index	3
2.4.8	int	
	HYPRE_SStructMatrixGetValues (HYPRE_SStructMatrix matrix, int part,	
	int* index, int var, int nentries,	
	int* entries, HYPRE_Complex* values)	
	Get matrix coefficients index by index	38
2.4.9	int	
	HYPRE_SStructMatrixGetFEMValues (HYPRE_SStructMatrix matrix,	
	int part, int* index,	
	HYPRE_Complex* values)	
	Get finite element stiffness matrix coefficients index by index	38
2.4.10	int	
4.4.10	1110	

	HYPRE_SStructMatrixSetBoxValues (HYPRE_SStructMatrix matrix,	
	int part, int* ilower, int* iupper,	
	int var, int nentries, HYPRE_Inte	
	entiies HYRRE_oompeex vvalees)	
	Set matrix coefficients a box at a time.	38
2.4.11	int	
2.4.11	HYPRE_SStructMatrixAddToBoxValues (HYPRE_SStructMatrix matrix,	
	int part, int* ilower, int* iupper,	
	int var, int nentries, int* entries,	
	HYPRE_Complex* values)	
	Add to matrix coefficients a box at a time	39
	•	0.0
2.4.12	int	
	HYPRE_SStructMatrixGetBoxValues (HYPRE_SStructMatrix matrix,	
	int part, int* ilower, int* iupper,	
	int var, int nentries, int* entries,	
	HYPRE_Complex* values)	
	Get matrix coefficients a box at a time	39
2.4.13	int	
	HYPRE_SStructMatrixAssemble (HYPRE_SStructMatrix matrix)	
	Finalize the construction of the matrix before using	40
2.4.14	int	
2.4.14	HYPRE_SStructMatrixSetSymmetric (HYPRE_SStructMatrix matrix,	
	int part, int var, int to_var,	
	int symmetric)	
	Define symmetry properties for the stencil entries in the matrix	40
		40
2.4.15	int	
	HYPRE_SStructMatrixSetNSSymmetric (HYPRE_SStructMatrix matrix,	
	int symmetric)	
	Define symmetry properties for all non-stencil matrix entries	40
2.4.16	int	
	HYPRE_SStructMatrixSetObjectType (HYPRE_SStructMatrix matrix,	
	int type)	
	Set the storage type of the matrix object to be constructed	40
2.4.17	int	
2.4.17	HYPRE_SStructMatrixGetObject (HYPRE_SStructMatrix matrix,	
	void** object)	
	Get a reference to the constructed matrix object.	11
	Get a rejevence to the constructed matrix object.	41
2.4.18	int	
	HYPRE_SStructMatrixPrint (const char* filename,	
	HYPRE_SStructMatrix matrix, int all)	
	Print the matrix to file.	41

_ 2.4.1 _____

 $typedef\ struct\ hypre_SStructMatrix_struct\ \textbf{*HYPRE_SStructMatrix}$

The matrix object

2.4.2

int
HYPRE_SStructMatrixCreate (MPI_Comm comm, HYPRE_SStructGraph
graph, HYPRE_SStructMatrix* matrix)

Create a matrix object

 $_$ 2.4.3 $_$

int HYPRE_SStructMatrixDestroy (HYPRE_SStructMatrix matrix)

Destroy a matrix object

_ 2.4.4 _

int HYPRE_SStructMatrixInitialize (HYPRE_SStructMatrix matrix)

Prepare a matrix object for setting coefficient values

2.4.5

int
HYPRE_SStructMatrixSetValues (HYPRE_SStructMatrix matrix, int part, int* index, int var, int nentries, int* entries, HYPRE_Complex* values)

Set matrix coefficients index by index. The values array is of length nentries.

NOTE: For better efficiency, use HYPRE_SStructMatrixSetBoxValues to set coefficients a box at a time.

NOTE: Users are required to set values on all processes that own the associated variables. This means that some data will be multiply defined.

NOTE: The entries in this routine must all be of the same type: either stencil or non-stencil, but not both. Also, if they are stencil entries, they must all represent couplings to the same variable type (there are no such restrictions for non-stencil entries).

2.4.6

HYPRE_SStructMatrixAddToValues (HYPRE_SStructMatrix matrix, int part, int* index, int var, int nentries, int* entries, HYPRE_Complex* values)

Add to matrix coefficients index by index. The values array is of length nentries.

NOTE: For better efficiency, use HYPRE_SStructMatrixAddToBoxValues to set coefficients a box at a time.

NOTE: Users are required to set values on all processes that own the associated variables. This means that some data will be multiply defined.

NOTE: The entries in this routine must all be of the same type: either stencil or non-stencil, but not both. Also, if they are stencil entries, they must all represent couplings to the same variable type.

2.4.7

int
HYPRE_SStructMatrixAddFEMValues (HYPRE_SStructMatrix matrix, int
part, int* index, HYPRE_Complex* values)

Add finite element stiffness matrix coefficients index by index. The layout of the data in values is determined by the routines HYPRE_SStructGridSetFEMOrdering and HYPRE_SStructGraphSetFEMSparsity ($\rightarrow 2.3.6$, page 32).

2.4.8

int

HYPRE_SStructMatrixGetValues (HYPRE_SStructMatrix matrix, int part, int* index, int var, int nentries, int* entries, HYPRE_Complex* values)

Get matrix coefficients index by index. The values array is of length nentries.

NOTE: For better efficiency, use HYPRE_SStructMatrixGetBoxValues to get coefficients a box at a time.

NOTE: Users may get values on any process that owns the associated variables.

NOTE: The entries in this routine must all be of the same type: either stencil or non-stencil, but not both. Also, if they are stencil entries, they must all represent couplings to the same variable type (there are no such restrictions for non-stencil entries).

2.4.9

HYPRE_SStructMatrixGetFEMValues (HYPRE_SStructMatrix matrix, int part, int* index, HYPRE_Complex* values)

Get finite element stiffness matrix coefficients index by index. The layout of the data in values is determined by the routines HYPRE_SStructGridSetFEMOrdering and HYPRE_SStructGraphSetFEMSparsity ($\rightarrow 2.3.6$, page 32).

2.4.10

HYPRE_SStructMatrixSetBoxValues (HYPRE_SStructMatrix matrix, int part, int* ilower, int* iupper, int var, int nentries, HYPRE_Inte entiies HYRRE_oompeex vvalees)

Set matrix coefficients a box at a time. The data in values is ordered as follows:

```
m = 0;
for (k = ilower[2]; k <= iupper[2]; k++)
  for (j = ilower[1]; j <= iupper[1]; j++)
    for (i = ilower[0]; i <= iupper[0]; i++)
    for (entry = 0; entry < nentries; entry++)</pre>
```

```
{
   values[m] = ...;
   m++;
}
```

NOTE: Users are required to set values on all processes that own the associated variables. This means that some data will be multiply defined.

NOTE: The entries in this routine must all be of the same type: either stencil or non-stencil, but not both. Also, if they are stencil entries, they must all represent couplings to the same variable type (there are no such restrictions for non-stencil entries).

2.4.11

int

HYPRE_SStructMatrixAddToBoxValues (HYPRE_SStructMatrix matrix, int part, int* ilower, int* iupper, int var, int nentries, int* entries, HYPRE_Complex* values)

Add to matrix coefficients a box at a time. The data in values is ordered as in HYPRE_SStructMatrixSetBoxValues.

NOTE: Users are required to set values on all processes that own the associated variables. This means that some data will be multiply defined.

NOTE: The entries in this routine must all be of stencil type. Also, they must all represent couplings to the same variable type.

2.4.12

int

HYPRE_SStructMatrixGetBoxValues (HYPRE_SStructMatrix matrix, int part, int* ilower, int* iupper, int var, int nentries, int* entries, HYPRE_Complex* values)

Get matrix coefficients a box at a time. The data in values is ordered as in HYPRE_SStructMatrixSetBoxValues.

NOTE: Users may get values on any process that owns the associated variables.

NOTE: The entries in this routine must all be of stencil type. Also, they must all represent couplings to the same variable type.

2.4.13

int HYPRE_SStructMatrixAssemble (HYPRE_SStructMatrix matrix)

Finalize the construction of the matrix before using

_ 2.4.14 _

int

HYPRE_SStructMatrixSetSymmetric (HYPRE_SStructMatrix matrix, int part, int var, int to_var, int symmetric)

Define symmetry properties for the stencil entries in the matrix. The boolean argument symmetric is applied to stencil entries on part part that couple variable var to variable to_var. A value of -1 may be used for part, var, or to_var to specify "all". For example, if part and to_var are set to -1, then the boolean is applied to stencil entries on all parts that couple variable var to all other variables.

By default, matrices are assumed to be nonsymmetric. Significant storage savings can be made if the matrix is symmetric.

2.4.15

HYPRE_SStructMatrixSetNSSymmetric (HYPRE_SStructMatrix matrix, int symmetric)

Define symmetry properties for all non-stencil matrix entries

_ 2.4.16 ___

HYPRE_SStructMatrixSetObjectType (HYPRE_SStructMatrix matrix, int type)

Set the storage type of the matrix object to be constructed. Currently, type can be either HYPRE_SSTRUCT (the default), HYPRE_STRUCT, or HYPRE_PARCSR.

See Also:

HYPRE_SStructMatrixGetObject (\rightarrow 2.4.17, page 41)

2.4.17

HYPRE_SStructMatrixGetObject (HYPRE_SStructMatrix matrix, void** object)

Get a reference to the constructed matrix object.

See Also:

HYPRE_SStructMatrixSetObjectType (\rightarrow 2.4.16, page 40)

2.4.18

HYPRE_SStructMatrixPrint (const char* filename, HYPRE_SStructMatrix matrix, int all)

Print the matrix to file. This is mainly for debugging purposes.

__ 2.5 ____

SStruct Vectors

Names		
2.5.1	typedef struct hypre_SStructVector_struct *HYPRE_SStructVector The vector object	43
2.5.2	int HYPRE_SStructVectorCreate (MPI_Comm comm, HYPRE_SStructGrid grid, HYPRE_SStructVector* vector)	
	Create a vector object	43
2.5.3	int HYPRE_SStructVectorDestroy (HYPRE_SStructVector vector) Destroy a vector object	43
2.5.4	int	

	HYPRE_SStructVectorInitialize (HYPRE_SStructVector vector) Prepare a vector object for setting coefficient values	44
2.5.5	int	
	HYPRE_SStructVectorSetValues (HYPRE_SStructVector vector, int part,	
	int* index, int var, HYPRE_Complex* value)	
	Set vector coefficients index by index	44
0.5.6		
2.5.6	int HYPRE_SStructVectorAddToValues (HYPRE_SStructVector vector,	
	int part, int* index, int var,	
	HYPRE_Complex* value)	
	Add to vector coefficients index by index	44
2.5.7	int	
2.0.1	HYPRE_SStructVectorAddFEMValues (HYPRE_SStructVector vector,	
	int part, int* index,	
	HYPRE_Complex* values)	
	Add finite element vector coefficients index by index	44
2.5.8	int	
	HYPRE_SStructVectorGetValues (HYPRE_SStructVector vector, int part,	
	int* index, int var,	
	HYPRE_Complex* value)	
	Get vector coefficients index by index	45
2.5.9	int	
	${\bf HYPRE_SStructVectorGetFEMValues}~({\bf HYPRE_SStructVector~vector},$	
	int part, int* index,	
	HYPRE_Complex* values)	45
	Get finite element vector coefficients index by index	45
2.5.10	int	
	HYPRE_SStructVectorSetBoxValues (HYPRE_SStructVector vector,	
	int part, int* ilower, int* iupper, int var, HYPRE_Complex* values)	
	Set vector coefficients a box at a time.	45
0 7 11		10
2.5.11	int HYPRE_SStructVectorAddToBoxValues (HYPRE_SStructVector vector,	
	int part, int* ilower, int* iupper,	
	int var,	
	HYPRE_Complex* values)	
	Add to vector coefficients a box at a time.	46
2.5.12	int	
	HYPRE_SStructVectorGetBoxValues (HYPRE_SStructVector vector,	
	int part, int* ilower, int* iupper,	
	int var, HYPRE_Complex* values)	
	Get vector coefficients a box at a time.	46
2.5.13	int	
	${\bf HYPRE_SStructVectorAssemble} \ ({\bf HYPRE_SStructVector\ vector})$	
	Finalize the construction of the vector before using	46
2.5.14	int	

	HYPRE_SStructVectorGather (HYPRE_SStructVector vector)	
	Gather vector data so that efficient GetValues can be done	47
2.5.15	int	
	HYPRE_SStructVectorSetObjectType (HYPRE_SStructVector vector,	
	int type)	
	Set the storage type of the vector object to be constructed	47
2.5.16	int	
	HYPRE_SStructVectorGetObject (HYPRE_SStructVector vector,	
	void** object)	
	Get a reference to the constructed vector object	47
2.5.17	int	
	HYPRE_SStructVectorPrint (const char* filename,	
	HYPRE_SStructVector vector, int all)	
	Print the vector to file	47

2.5.1

 $typedef \ struct \ hypre_SStructVector_struct \ *HYPRE_SStructVector$

The vector object

2.5.2

int

HYPRE_SStructVectorCreate (MPI_Comm comm, HYPRE_SStructGrid grid, HYPRE_SStructVector* vector)

Create a vector object

 $_$ 2.5.3 $_$

int HYPRE_SStructVectorDestroy (HYPRE_SStructVector vector)

Destroy a vector object

254

int HYPRE_SStructVectorInitialize (HYPRE_SStructVector vector)

Prepare a vector object for setting coefficient values

2.5.5

int

HYPRE_SStructVectorSetValues (HYPRE_SStructVector vector, int part, int* index, int var, HYPRE_Complex* value)

Set vector coefficients index by index.

NOTE: For better efficiency, use HYPRE_SStructVectorSetBoxValues to set coefficients a box at a time.

NOTE: Users are required to set values on all processes that own the associated variables. This means that some data will be multiply defined.

2.5.6

int

HYPRE_SStructVectorAddToValues (HYPRE_SStructVector vector, int part, int* index, int var, HYPRE_Complex* value)

Add to vector coefficients index by index.

NOTE: For better efficiency, use HYPRE_SStructVectorAddToBoxValues to set coefficients a box at a time.

NOTE: Users are required to set values on all processes that own the associated variables. This means that some data will be multiply defined.

 $_$ 2.5.7 $_$

int

HYPRE_SStructVectorAddFEMValues (HYPRE_SStructVector vector, int part, int* index, HYPRE_Complex* values)

Add finite element vector coefficients index by index. The layout of the data in values is determined by the routine HYPRE_SStructGridSetFEMOrdering.

2.5.8

int

HYPRE_SStructVectorGetValues (HYPRE_SStructVector vector, int part, int* index, int var, HYPRE_Complex* value)

Get vector coefficients index by index. Users must first call the routine HYPRE_SStructVectorGather to ensure that data owned by multiple processes is correct.

NOTE: For better efficiency, use HYPRE_SStructVectorGetBoxValues to get coefficients a box at a time.

NOTE: Users may only get values on processes that own the associated variables.

_ 2.5.9 _

int

HYPRE_SStructVectorGetFEMValues (HYPRE_SStructVector vector, int part, int* index, HYPRE_Complex* values)

Get finite element vector coefficients index by index. The layout of the data in values is determined by the routine HYPRE_SStructGridSetFEMOrdering. Users must first call the routine HYPRE_SStructVectorGather to ensure that data owned by multiple processes is correct.

 $_$ 2.5.10 $_$

int

HYPRE_SStructVectorSetBoxValues (HYPRE_SStructVector vector, int part, int* ilower, int* iupper, int var, HYPRE_Complex* values)

Set vector coefficients a box at a time. The data in values is ordered as follows:

```
m = 0;
for (k = ilower[2]; k <= iupper[2]; k++)
    for (j = ilower[1]; j <= iupper[1]; j++)
        for (i = ilower[0]; i <= iupper[0]; i++)</pre>
```

```
{
   values[m] = ...;
   m++;
}
```

NOTE: Users are required to set values on all processes that own the associated variables. This means that some data will be multiply defined.

2.5.11

int

HYPRE_SStructVectorAddToBoxValues (HYPRE_SStructVector vector, int part, int* ilower, int* iupper, int var, HYPRE_Complex* values)

Add to vector coefficients a box at a time. The data in values is ordered as in HYPRE_SStructVectorSetBoxValues.

NOTE: Users are required to set values on all processes that own the associated variables. This means that some data will be multiply defined.

_ 2.5.12 _

int

HYPRE_SStructVectorGetBoxValues (HYPRE_SStructVector vector, int part, int* ilower, int* iupper, int var, HYPRE_Complex* values)

Get vector coefficients a box at a time. The data in values is ordered as in HYPRE_SStructVectorSetBoxValues. Users must first call the routine HYPRE_SStructVectorGather to ensure that data owned by multiple processes is correct.

NOTE: Users may only get values on processes that own the associated variables.

2.5.13

int HYPRE_SStructVectorAssemble (HYPRE_SStructVector vector)

Finalize the construction of the vector before using

2.5.14

int HYPRE_SStructVectorGather (HYPRE_SStructVector vector)

Gather vector data so that efficient GetValues can be done. This routine must be called prior to calling GetValues to ensure that correct and consistent values are returned, especially for non cell-centered data that is shared between more than one processor.

__ 2.5.15 _____

int **HYPRE_SStructVectorSetObjectType** (HYPRE_SStructVector vector, int type)

Set the storage type of the vector object to be constructed. Currently, type can be either HYPRE_SSTRUCT (the default), HYPRE_STRUCT, or HYPRE_PARCSR.

See Also:

HYPRE_SStructVectorGetObject ($\rightarrow 2.5.16$, page 47)

2.5.16

HYPRE_SStructVectorGetObject (HYPRE_SStructVector vector, void** object)

Get a reference to the constructed vector object.

See Also:

HYPRE_SStructVectorSetObjectType ($\rightarrow 2.5.15$, page 47)

_ 2.5.17 __

HYPRE_SStructVectorPrint (const char* filename, HYPRE_SStructVector vector, int all)

Print the vector to file. This is mainly for debugging purposes.

3

IJ System Interface

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This interface represents a linear-algebraic conceptual view of a linear system. The 'I' and 'J' in the name are meant to be mnemonic for the traditional matrix notation A(I,J).

3.1

IJ Matrices

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	Print the matrix to file.	55

3.1.1

typedef struct hypre_IJMatrix_struct *HYPRE_IJMatrix

The matrix object

3.1.2

int

HYPRE_IJMatrixCreate (MPI_Comm comm, int ilower, int iupper, int jlower, int jupper, HYPRE_IJMatrix* matrix)

Create a matrix object. Each process owns some unique consecutive range of rows, indicated by the global row indices ilower and iupper. The row data is required to be such that the value of ilower on any process p be exactly one more than the value of iupper on process p-1. Note that the first row of the global matrix may start with any integer value. In particular, one may use zero- or one-based indexing.

For square matrices, jlower and jupper typically should match ilower and iupper, respectively. For rectangular matrices, jlower and jupper should define a partitioning of the columns. This partitioning must be used for any vector v that will be used in matrix-vector products with the rectangular matrix. The matrix data structure may use jlower and jupper to store the diagonal blocks (rectangular in general) of the matrix separately from the rest of the matrix.

Collective.

3.1.3

int HYPRE_IJMatrixDestroy (HYPRE_IJMatrix matrix)

Destroy a matrix object. An object should be explicitly destroyed using this destructor when the user's code no longer needs direct access to it. Once destroyed, the object must not be referenced again. Note that the object may not be deallocated at the completion of this call, since there may be internal package references to the object. The object will then be destroyed when all internal reference counts go to zero.

3.1.4

int HYPRE_IJMatrixInitialize (HYPRE_IJMatrix matrix)

Prepare a matrix object for setting coefficient values. This routine will also re-initialize an already assembled matrix, allowing users to modify coefficient values.

3.1.5

HYPRE_IJMatrixSetValues (HYPRE_IJMatrix matrix, int nrows, int* ncols, const int* rows, const int* cols, const HYPRE_Complex* values)

Sets values for nrows rows or partial rows of the matrix. The arrays ncols and rows are of dimension nrows and contain the number of columns in each row and the row indices, respectively. The array cols contains the column indices for each of the rows, and is ordered by rows. The data in the values array corresponds directly to the column entries in cols. Erases any previous values at the specified locations and replaces them with new ones, or, if there was no value there before, inserts a new one if set locally. Note that it is not possible to set values on other processors. If one tries to set a value from proc i on proc j, proc i will erase all previous occurrences of this value in its stack (including values generated with AddToValues), and treat it like a zero value. The actual value needs to be set on proc j.

Note that a threaded version (threaded over the number of rows) will be called if $HYPRE_IJMatrixSetOMPFlag$ is set to a value !=0. This requires that rows[i] != rows[j] for i!=j and is only efficient if a large number of rows is set in one call to $HYPRE_IJMatrixSetValues$.

Not collective.

3.1.6

int

HYPRE_IJMatrixAddToValues (HYPRE_IJMatrix matrix, int nrows, int* ncols, const int* rows, const int* cols, const HYPRE_Complex* values)

Adds to values for nrows rows or partial rows of the matrix. Usage details are analogous to HYPRE_IJMatrixSetValues. Adds to any previous values at the specified locations, or, if there was no value there before, inserts a new one. AddToValues can be used to add to values on other processors.

Note that a threaded version (threaded over the number of rows) will be called if HYPRE_IJMatrixSetOMPFlag is set to a value != 0. This requires that rows[i] != rows[j] for i!= j and is only efficient if a large number of rows is added in one call to HYPRE_IJMatrixAddToValues.

Not collective.

_ 3.1.7 _

int **HYPRE_IJMatrixAssemble** (HYPRE_IJMatrix matrix)

Finalize the construction of the matrix before using

_ 3.1.8 _

int

HYPRE_IJMatrixGetRowCounts (HYPRE_IJMatrix matrix, int nrows, int* rows, int* ncols)

Gets number of nonzeros elements for nrows rows specified in rows and returns them in ncols, which needs to be allocated by the user

__ 3.1.9 ____

int

HYPRE_IJMatrixGetValues (HYPRE_IJMatrix matrix, int nrows, int* ncols, int* rows, int* cols, HYPRE_Complex* values)

Gets values for nrows rows or partial rows of the matrix. Usage details are analogous to HYPRE_IJMatrixSetValues.

__ 3.1.10 ___

int HYPRE_IJMatrixSetObjectType (HYPRE_IJMatrix matrix, int type)

Set the storage type of the matrix object to be constructed. Currently, type can only be HYPRE_PARCSR.

Not collective, but must be the same on all processes.

See Also:

<code>HYPRE_IJMatrixGetObject</code> ($\rightarrow 3.1.13, \; page \; 53)$

_ 3.1.11 ____

int **HYPRE_IJMatrixGetObjectType** (HYPRE_IJMatrix matrix, int* type)

Get the storage type of the constructed matrix object

_ 3.1.12 _

HYPRE_IJMatrixGetLocalRange (HYPRE_IJMatrix matrix, int* ilower, int* iupper, int* jlower, int* jupper)

Gets range of rows owned by this processor and range of column partitioning for this processor

_ 3.1.13 _

int HYPRE_IJMatrixGetObject (HYPRE_IJMatrix matrix, void** object)

Get a reference to the constructed matrix object.

See Also:

HYPRE_IJMatrixSetObjectType ($\rightarrow 3.1.10$, page 53)

3.1.14

int HYPRE_IJMatrixSetRowSizes (HYPRE_IJMatrix matrix, const int* sizes)

(Optional) Set the max number of nonzeros to expect in each row. The array sizes contains estimated sizes for each row on this process. This call can significantly improve the efficiency of matrix construction, and should always be utilized if possible.

Not collective.

3.1.15

HYPRE_IJMatrixSetDiagOffdSizes (HYPRE_IJMatrix matrix, const int* diag_sizes, const int* offdiag_sizes)

(Optional) Sets the exact number of nonzeros in each row of the diagonal and off-diagonal blocks. The diagonal block is the submatrix whose column numbers correspond to rows owned by this process, and the off-diagonal block is everything else. The arrays diag_sizes and offdiag_sizes contain estimated sizes for each row of the diagonal and off-diagonal blocks, respectively. This routine can significantly improve the efficiency of matrix construction, and should always be utilized if possible.

Not collective.

3.1.16

int

 $\label{eq:hypre_interpolation} \textbf{HYPRE_IJMatrix SetMaxOffProcElmts} \ (\textbf{HYPRE_IJMatrix matrix}, \ \textbf{int} \\ \text{max_off_proc_elmts})$

(Optional) Sets the maximum number of elements that are expected to be set (or added) on other processors from this processor This routine can significantly improve the efficiency of matrix construction, and should always be utilized if possible.

Not collective.

3 1 17

int HYPRE_IJMatrixSetPrintLevel (HYPRE_IJMatrix matrix, int print_level)

(Optional) Sets the print level, if the user wants to print error messages. The default is 0, i.e. no error messages are printed.

3.1.18

int HYPRE_IJMatrixSetOMPFlag (HYPRE_IJMatrix matrix, int omp_flag)

(Optional) if set, will use a threaded version of HYPRE_IJMatrixSetValues and HYPRE_IJMatrixAddToValues. This is only useful if a large number of rows is set or added to at once.

NOTE that the values in the rows array of HYPRE_IJMatrixSetValues or HYPRE_IJMatrixAddToValues must be different from each other !!!

This option is VERY inefficient if only a small number of rows is set or added at once and/or if reallocation of storage is required and/or if values are added to off processor values.

3.1.19

HYPRE_IJMatrixRead (const char* filename, MPI_Comm comm, int type, HYPRE_IJMatrix* matrix)

Read the matrix from file. This is mainly for debugging purposes.

_ 3.1.20 __

int **HYPRE_IJMatrixPrint** (HYPRE_IJMatrix matrix, const char* filename)

Print the matrix to file. This is mainly for debugging purposes.

_ 3.2 _

IJ Vectors

Names		
3.2.1	typedef struct hypre_IJVector_struct *HYPRE_IJVector	
	The vector object	57
3.2.2	int HYPRE_IJVectorCreate (MPI_Comm comm, int jlower, int jupper, HYPRE_IJVector* vector)	
	Create a vector object.	57
3.2.3	int HYPRE_IJVectorDestroy (HYPRE_IJVector vector) Destroy a vector object.	58
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0.2.4	HYPRE_IJVectorInitialize (HYPRE_IJVector vector)	
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3.2.5	int HYPRE_IJVectorSetMaxOffProcElmts (HYPRE_IJVector vector, int max_off_proc_elmts)	
	(Optional) Sets the maximum number of elements that are expected to be set (or added) on other processors from this processor This routine can significantly improve the efficiency of matrix construction, and should always be utilized if possible.	58
3.2.6	int	
	HYPRE_IJVectorSetValues (HYPRE_IJVector vector, int nvalues, const int* indices, const HYPRE_Complex* values)	
	Sets values in vector.	59
3.2.7	int HYPRE_IJVectorAddToValues (HYPRE_IJVector vector, int nvalues, const int* indices, const HYPRE_Complex* values)	
	Adds to values in vector.	59
3.2.8	int HYPRE_IJVectorAssemble (HYPRE_IJVector vector) Finalize the construction of the vector before using	59
3.2.9	int HYPRE_IJVectorGetValues (HYPRE_IJVector vector, int nvalues,	
	const int* indices, HYPRE_Complex* values)	F0
2.0.10	Gets values in vector.	59
3.2.10	int	

	HYPRE_IJVectorSetObjectType (HYPRE_IJVector vector, int type) Set the storage type of the vector object to be constructed	60
3.2.11	int	
	HYPRE_IJVectorGetObjectType (HYPRE_IJVector vector, int* type)	
	Get the storage type of the constructed vector object	60
3.2.12	int	
	HYPRE_IJVectorGetLocalRange (HYPRE_IJVector vector, int* jlower, int* jupper)	
	Returns range of the part of the vector owned by this processor	60
3.2.13	int	
	HYPRE_IJVectorGetObject (HYPRE_IJVector vector, void** object)	
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3.2.14	int	
	HYPRE_IJVectorSetPrintLevel (HYPRE_IJVector vector, int print_level)	
	(Optional) Sets the print level, if the user wants to print error messages.	61
3.2.15	int	
	HYPRE_IJVector* (const char* filename, MPI_Comm comm, int type, HYPRE_IJVector* vector)	
	Read the vector from file.	61
3.2.16	int	
	HYPRE_IJVectorPrint (HYPRE_IJVector vector, const char* filename)	
	Print the vector to file.	61

 $_$ 3.2.1 $_$

typedef struct hypre_IJVector_struct *HYPRE_IJVector

The vector object

3.2.2

HYPRE_IJVectorCreate (MPI_Comm comm, int jlower, int jupper, HYPRE_IJVector* vector)

Create a vector object. Each process owns some unique consecutive range of vector unknowns, indicated by the global indices jlower and jupper. The data is required to be such that the value of jlower on any process p be exactly one more than the value of jupper on process p-1. Note that the first index of the global vector may start with any integer value. In particular, one may use zero- or one-based indexing.

Collective.

3.2.3

int HYPRE_IJVectorDestroy (HYPRE_IJVector vector)

Destroy a vector object. An object should be explicitly destroyed using this destructor when the user's code no longer needs direct access to it. Once destroyed, the object must not be referenced again. Note that the object may not be deallocated at the completion of this call, since there may be internal package references to the object. The object will then be destroyed when all internal reference counts go to zero.

 $_{-}$ 3.2.4 $_{-}$

int HYPRE_IJVectorInitialize (HYPRE_IJVector vector)

Prepare a vector object for setting coefficient values. This routine will also re-initialize an already assembled vector, allowing users to modify coefficient values.

_ 3.2.5 _

int **HYPRE_IJVectorSetMaxOffProcElmts** (HYPRE_IJVector vector, int max_off_proc_elmts)

(Optional) Sets the maximum number of elements that are expected to be set (or added) on other processors from this processor This routine can significantly improve the efficiency of matrix construction, and should always be utilized if possible.

Not collective.

3.2.6

int

HYPRE_IJVectorSetValues (HYPRE_IJVector vector, int nvalues, const int* indices, const HYPRE_Complex* values)

Sets values in vector. The arrays values and indices are of dimension nvalues and contain the vector values to be set and the corresponding global vector indices, respectively. Erases any previous values at the specified locations and replaces them with new ones. Note that it is not possible to set values on other processors. If one tries to set a value from proc i on proc j, proc i will erase all previous occurrences of this value in its stack (including values generated with AddToValues), and treat it like a zero value. The actual value needs to be set on proc j.

Not collective.

3.2.7

HYPRE_IJVectorAddToValues (HYPRE_IJVector vector, int nvalues, const int* indices, const HYPRE_Complex* values)

Adds to values in vector. Usage details are analogous to HYPRE_IJVectorSetValues. Adds to any previous values at the specified locations, or, if there was no value there before, inserts a new one. AddToValues can be used to add to values on other processors.

Not collective.

3.2.8

int HYPRE_IJVectorAssemble (HYPRE_IJVector vector)

Finalize the construction of the vector before using

 $_$ 3.2.9 $_$

int

HYPRE_IJVectorGetValues (HYPRE_IJVector vector, int nvalues, const int* indices, HYPRE_Complex* values)

Gets values in vector. Usage details are analogous to HYPRE_IJVectorSetValues.

Not collective.

 $_{-}$ 3.2.10 $_{-}$

int HYPRE_IJVectorSetObjectType (HYPRE_IJVector vector, int type)

Set the storage type of the vector object to be constructed. Currently, type can only be HYPRE_PARCSR.

Not collective, but must be the same on all processes.

See Also:

HYPRE_IJVectorGetObject (\rightarrow 3.2.13, page 60)

_ 3.2.11 _

int **HYPRE_IJVectorGetObjectType** (HYPRE_IJVector vector, int* type)

Get the storage type of the constructed vector object

 $_$ 3.2.12 $_$

int **HYPRE_IJVectorGetLocalRange** (HYPRE_IJVector vector, int* jlower, int* jupper)

Returns range of the part of the vector owned by this processor

_ 3.2.13 ____

int HYPRE_IJVectorGetObject (HYPRE_IJVector vector, void** object)

Get a reference to the constructed vector object.

See Also:

HYPRE_IJVectorSetObjectType (\rightarrow 3.2.10, page 60)

3.2.14

int HYPRE_IJVectorSetPrintLevel (HYPRE_IJVector vector, int print_level)

(Optional) Sets the print level, if the user wants to print error messages. The default is 0, i.e. no error messages are printed.

3.2.15

int **HYPRE_IJVectorRead** (const char* filename, MPI_Comm comm, int type, HYPRE_IJVector* vector)

Read the vector from file. This is mainly for debugging purposes.

_ 3.2.16 ____

int HYPRE_IJVectorPrint (HYPRE_IJVector vector, const char* filename)

Print the vector to file. This is mainly for debugging purposes.

4

Struct Solvers

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4.11	Struct Hybrid Solver	85
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These solvers use matrix/vector storage schemes that are tailored to structured grid problems.

_ 4.1 ____

Struct Solvers

Names

4.1.1	typedef struct hypre_StructSolver_struct *HYPRE_StructSolver	
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_ 4.1.1 _

 $typedef\ struct\ hypre_StructSolver_struct\ *HYPRE_StructSolver$

The solver object

_ 4.2 _

Struct Jacobi Solver

Names		
4.2.1	int	
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	HYPRE_StructSolver* solver)	
	Create a solver object	64
4.2.2	int	
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4.2.3	int	
	HYPRE_StructJacobiSetup (HYPRE_StructSolver solver,	
	HYPRE_StructMatrix A,	
	HYPRE_StructVector b,	
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	Prepare to solve the system	64
4.2.4	int	
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	HYPRE_StructMatrix A,	
	HYPRE_StructVector b,	
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4.2.7	int	
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	, - , , , , , , , , , , , , , , , , , ,	Je
4.2.8	int	

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	,	iterations)	
	Return the number of iterations taken	,	66
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		(HYPRE_StructSolver	
		solver,	
		HYPRE_Real*	
		norm)	
	Return the norm of the final relative residual		66

 $_{-}$ 4.2.1 $_{-}$

HYPRE_StructJacobiCreate (MPI_Comm comm, HYPRE_StructSolver* solver)

Create a solver object

 $_$ 4.2.2 $_$

int HYPRE_StructJacobiDestroy (HYPRE_StructSolver solver)

Destroy a solver object. An object should be explicitly destroyed using this destructor when the user's code no longer needs direct access to it. Once destroyed, the object must not be referenced again. Note that the object may not be deallocated at the completion of this call, since there may be internal package references to the object. The object will then be destroyed when all internal reference counts go to zero.

4.2.3

HYPRE_StructJacobiSetup (HYPRE_StructSolver solver, HYPRE_StructMatrix A, HYPRE_StructVector b, HYPRE_StructVector x)

Prepare to solve the system. The coefficient data in b and x is ignored here, but information about the layout of the data may be used.

4.2.4

int **HYPRE_StructJacobiSolve** (HYPRE_StructSolver solver, HYPRE_StructMatrix A, HYPRE_StructVector b, HYPRE_StructVector x)

Solve the system

4.2.5

HYPRE_StructJacobiSetTol (HYPRE_StructSolver solver, HYPRE_Real tol)

(Optional) Set the convergence tolerance

__ 4.2.6 _____

Int
HYPRE_StructJacobiSetMaxIter (HYPRE_StructSolver solver, int max_iter)

(Optional) Set maximum number of iterations

_ 4.2.7 _

int HYPRE_StructJacobiSetZeroGuess (HYPRE_StructSolver solver)

(Optional) Use a zero initial guess. This allows the solver to cut corners in the case where a zero initial guess is needed (e.g., for preconditioning) to reduce computational cost.

4.2.8

int HYPRE_StructJacobiSetNonZeroGuess (HYPRE_StructSolver solver)

(Optional) Use a nonzero initial guess. This is the default behavior, but this routine allows the user to switch back after using SetZeroGuess.

 $_$ 4.2.9 $_$

int

HYPRE_StructJacobiGetNumIterations (HYPRE_StructSolver solver, int* num_iterations)

Return the number of iterations taken

__ 4.2.10 _____

int

HYPRE_StructJacobiGetFinalRelativeResidualNorm (HYPRE_StructSolver solver, HYPRE_Real* norm)

Return the norm of the final relative residual

4.3

Struct PFMG Solver

Names

4.3.1 int

HYPRE_StructPFMGCreate (MPI_Comm comm,

HYPRE_StructSolver* solver)

4.3.2 int

HYPRE_StructPFMGDestroy (HYPRE_StructSolver solver)

4.3.3 int

	HYPRE_StructPFMGSetup (HYPRE_StructSolver solver,	
	HYPRE_StructMatrix A,	
	HYPRE_StructVector b,	
	HYPRE_StructVector x)	
	Prepare to solve the system	69
4.0.4		
4.3.4	int	
	HYPRE_StructPFMGSolve (HYPRE_StructSolver solver,	
	HYPRE_StructMatrix A,	
	HYPRE_StructVector b,	
	$HYPRE_StructVector x)$	
	Solve the system	69
4.3.5	int	
	HYPRE_StructPFMGSetTol (HYPRE_StructSolver solver, HYPRE_Real tol)	
	(Optional) Set the convergence tolerance	69
	· - /	
4.3.6	int	
	HYPRE_StructPFMGSetMaxIter (HYPRE_StructSolver solver,	
	int max_iter)	
	(Optional) Set maximum number of iterations	69
4.3.7	int	
	HYPRE_StructPFMGSetMaxLevels (HYPRE_StructSolver solver,	
	int max_levels)	
	(Optional) Set maximum number of multigrid grid levels	70
4.0.0	, <u> </u>	
4.3.8	int	
	HYPRE_StructPFMGSetRelChange (HYPRE_StructSolver solver,	
	int rel_change)	
	(Optional) Additionally require that the relative difference in successive it-	
	erates be small	70
4.3.9	int	
	HYPRE_StructPFMGSetZeroGuess (HYPRE_StructSolver solver)	
	(Optional) Use a zero initial guess	70
4.3.10	· - /	
4.5.10	int	
	HYPRE_StructPFMGSetNonZeroGuess (HYPRE_StructSolver solver)	70
	(Optional) Use a nonzero initial guess	70
4.3.11	int	
	HYPRE_StructPFMGSetRelaxType (HYPRE_StructSolver solver,	
	int relax_type)	
	(Optional) Set relaxation type	71
4.3.12	int	
1.0.12	HYPRE_StructPFMGSetRAPType (HYPRE_StructSolver solver,	
	int rap_type)	
	(Optional) Set type of coarse-grid operator to use	71
	(Optional) Set type of course-grid operator to use	11
4.3.13	int	
	${\bf HYPRE_StructPFMGSetNumPreRelax}~({\bf HYPRE_StructSolver}~solver,$	
	$int num_pre_relax)$	
	$(Optional)\ Set\ number\ of\ relaxation\ sweeps\ before\ coarse-grid\ correction$.	71
4 3 14	int	

	HYPRE_StructPFMGSetNumPostRelax (HYPRE_StructSolver solver,	
	int num_post_relax)	
	(Optional) Set number of relaxation sweeps after coarse-grid correction	72
4.3.15	int	
	HYPRE_StructPFMGSetSkipRelax (HYPRE_StructSolver solver,	
	int skip_relax)	
	(Optional) Skip relaxation on certain grids for isotropic problems	72
4.3.16	int	
	HYPRE_StructPFMGSetLogging (HYPRE_StructSolver solver, int logging)	
	(Optional) Set the amount of logging to do	72
4.3.17	int	
	HYPRE_StructPFMGSetPrintLevel (HYPRE_StructSolver solver,	
	int print_level)	
	(Optional) Set the amount of printing to do to the screen	72
4.3.18	int	
	HYPRE_StructPFMGGetNumIterations (HYPRE_StructSolver solver,	
	int* num_iterations)	
	Return the number of iterations taken	73
4.3.19	int	
	$HYPRE_StructPFMGGetFinalRelativeResidualNorm$	
	(HYPRE_StructSolver	
	solver,	
	HYPRE_Real*	
	norm)	
	Return the norm of the final relative residual	73

PFMG is a semicoarsening multigrid solver that uses pointwise relaxation. For periodic problems, users should try to set the grid size in periodic dimensions to be as close to a power-of-two as possible. That is, if the grid size in a periodic dimension is given by $N = 2^m * M$ where M is not a power-of-two, then M should be as small as possible. Large values of M will generally result in slower convergence rates.

 $_$ 4.3.1 $_$

HYPRE_StructPFMGCreate (MPI_Comm comm, HYPRE_StructSolver* solver)

Create a solver object

 $_{-}$ 4.3.2 $_{-}$

int HYPRE_StructPFMGDestroy (HYPRE_StructSolver solver)

Destroy a solver object

___ 4.3.3 _____

int

HYPRE_StructPFMGSetup (HYPRE_StructSolver solver, HYPRE_StructMatrix A, HYPRE_StructVector b, HYPRE_StructVector x)

Prepare to solve the system. The coefficient data in b and x is ignored here, but information about the layout of the data may be used.

_ 4.3.4 ___

int

HYPRE_StructPFMGSolve (HYPRE_StructSolver solver, HYPRE_StructMatrix A, HYPRE_StructVector b, HYPRE_StructVector x)

Solve the system

4.3.5

int

HYPRE_StructPFMGSetTol (HYPRE_StructSolver solver, HYPRE_Real tol)

(Optional) Set the convergence tolerance

_ 4.3.6 _____

ınt

HYPRE_StructPFMGSetMaxIter (HYPRE_StructSolver solver, int max_iter)

(Optional) Set maximum number of iterations

137

int **HYPRE_StructPFMGSetMaxLevels** (HYPRE_StructSolver solver, int max_levels)

(Optional) Set maximum number of multigrid grid levels

___ 4.3.8 _____

HYPRE_StructPFMGSetRelChange (HYPRE_StructSolver solver, int rel_change)

(Optional) Additionally require that the relative difference in successive iterates be small

__ 4.3.9 _____

int HYPRE_StructPFMGSetZeroGuess (HYPRE_StructSolver solver)

(Optional) Use a zero initial guess. This allows the solver to cut corners in the case where a zero initial guess is needed (e.g., for preconditioning) to reduce computational cost.

4.3.10

int HYPRE_StructPFMGSetNonZeroGuess (HYPRE_StructSolver solver)

(Optional) Use a nonzero initial guess. This is the default behavior, but this routine allows the user to switch back after using SetZeroGuess.

4.3.11

HYPRE_StructPFMGSetRelaxType (HYPRE_StructSolver solver, int relax_type)

(Optional) Set relaxation type.

Current relaxation methods set by relax_type are:

- 0 & Jacobi
- 1 & Weighted Jacobi (default)
- $2~\&~{\rm Red/Black}$ Gauss-Seidel (symmetric: RB pre-relaxation, BR post-relaxation) -
- 3 & Red/Black Gauss-Seidel (nonsymmetric: RB pre- and post-relaxation)

_ 4.3.12 __

HYPRE_StructPFMGSetRAPType (HYPRE_StructSolver solver, int rap_type)

(Optional) Set type of coarse-grid operator to use.

Current operators set by rap_type are:

- 0 Galerkin (default)
- 1 non-Galerkin 5-pt or 7-pt stencils

Both operators are constructed algebraically. The non-Galerkin option maintains a 5-pt stencil in 2D and a 7-pt stencil in 3D on all grid levels. The stencil coefficients are computed by averaging techniques.

4.3.13

HYPRE_StructPFMGSetNumPreRelax (HYPRE_StructSolver solver, int num_pre_relax)

(Optional) Set number of relaxation sweeps before coarse-grid correction

int HYPRE_StructPFMGSetNumPostRelax (HYPRE_StructSolver solver, int num_post_relax)

(Optional) Set number of relaxation sweeps after coarse-grid correction

HYPRE_StructPFMGSetSkipRelax (HYPRE_StructSolver solver, int skip_relax)

(Optional) Skip relaxation on certain grids for isotropic problems. This can greatly improve efficiency by eliminating unnecessary relaxations when the underlying problem is isotropic.

___ 4.3.16 _____

int HYPRE_StructPFMGSetLogging (HYPRE_StructSolver solver, int logging)

(Optional) Set the amount of logging to do

HYPRE_StructPFMGSetPrintLevel (HYPRE_StructSolver solver, int print_level)

(Optional) Set the amount of printing to do to the screen

4.3.18

int

HYPRE_StructPFMGGetNumIterations (HYPRE_StructSolver solver, int* num_iterations)

Return the number of iterations taken

4.3.19

int

$HYPRE_StructPFMGGetFinalRelativeResidualNorm$

(HYPRE_StructSolver solver, HYPRE_Real* norm)

Return the norm of the final relative residual

_ 4.4 _____

Struct SMG Solver

Names

4.4.1	int	
	HYPRE_StructSMGCreate (MPI_Comm comm,	
	HYPRE_StructSolver* solver)	
	Create a solver object	75
4.4.2	int	
	HYPRE_StructSMGDestroy (HYPRE_StructSolver solver)	
	Destroy a solver object	75
4.4.3	int	
	HYPRE_StructSMGSetup (HYPRE_StructSolver solver,	
	HYPRE_StructMatrix A,	
	HYPRE_StructVector b, HYPRE_StructVector x)	
	Prepare to solve the system.	75
4.4.4	int	
	HYPRE_StructSMGSolve (HYPRE_StructSolver solver,	
	HYPRE_StructMatrix A, HYPRE_StructVector b,	
	HYPRE_StructVector x)	
	Solve the system	75
4.4.5	int	

	HYPRE_StructSMGSetTol (HYPRE_StructSolver solver, HYPRE_Real tol) (Optional) Set the convergence tolerance	76
4.4.6	int HYPRE_StructSMGSetMaxIter (HYPRE_StructSolver solver, int max_iter)	
	(Optional) Set maximum number of iterations	76
4.4.7	int	
	HYPRE_StructSMGSetRelChange (HYPRE_StructSolver solver, int rel_change)	
	(Optional) Additionally require that the relative difference in successive iterates be small	76
4.4.8	int	
1.1.0	HYPRE_StructSMGSetZeroGuess (HYPRE_StructSolver solver) (Optional) Use a zero initial guess	76
4.4.9	int	
	HYPRE_StructSMGSetNonZeroGuess (HYPRE_StructSolver solver) (Optional) Use a nonzero initial guess.	77
4.4.10	int	
	HYPRE_StructSMGSetNumPreRelax (HYPRE_StructSolver solver, int num_pre_relax)	
	(Optional) Set number of relaxation sweeps before coarse-grid correction .	77
4.4.11	int	
	HYPRE_StructSMGSetNumPostRelax (HYPRE_StructSolver solver, int num_post_relax)	
	(Optional) Set number of relaxation sweeps after coarse-grid correction	77
4.4.12	int	
	HYPRE_StructSMGSetLogging (HYPRE_StructSolver solver, int logging) (Optional) Set the amount of logging to do	77
4.4.13	int	
	HYPRE_StructSMGSetPrintLevel (HYPRE_StructSolver solver, int print_level)	
	(Optional) Set the amount of printing to do to the screen	78
4.4.14	int	
	HYPRE_StructSMGGetNumIterations (HYPRE_StructSolver solver, int* num_iterations)	
	Return the number of iterations taken	78
4.4.15	int	
-	$\label{eq:hypre_struct} \begin{aligned} \textbf{HYPRE_StructSMGGetFinalRelativeResidualNorm} & \text{ (HYPRE_StructSolver} \\ & \text{solver}, \end{aligned}$	
	HYPRE_Real*	
	norm)	
	Return the norm of the final relative residual	78

SMG is a semicoarsening multigrid solver that uses plane smoothing (in 3D). The plane smoother calls a 2D SMG algorithm with line smoothing, and the line smoother is cyclic reduction (1D SMG). For periodic problems, the grid size in periodic dimensions currently must be a power-of-two.

HYPRE_StructSMGCreate (MPI_Comm comm, HYPRE_StructSolver* solver)

Create a solver object

 $_$ 4.4.2 $_$

int HYPRE_StructSMGDestroy (HYPRE_StructSolver solver)

Destroy a solver object

4.4.3

HYPRE_StructSMGSetup (HYPRE_StructSolver solver, HYPRE_StructMatrix A, HYPRE_StructVector b, HYPRE_StructVector x)

Prepare to solve the system. The coefficient data in b and x is ignored here, but information about the layout of the data may be used.

4.4.4

HYPRE_StructSMGSolve (HYPRE_StructSolver solver, HYPRE_StructMatrix A, HYPRE_StructVector b, HYPRE_StructVector x)

Solve the system

int HYPRE_StructSMGSetTol (HYPRE_StructSolver solver, HYPRE_Real tol)

(Optional) Set the convergence tolerance

__ 4.4.6 _____

int HYPRE_StructSMGSetMaxIter (HYPRE_StructSolver solver, int max_iter)

(Optional) Set maximum number of iterations

4.4.7

HYPRE_StructSMGSetRelChange (HYPRE_StructSolver solver, int rel_change)

(Optional) Additionally require that the relative difference in successive iterates be small

_ 4.4.8 _

int HYPRE_StructSMGSetZeroGuess (HYPRE_StructSolver solver)

(Optional) Use a zero initial guess. This allows the solver to cut corners in the case where a zero initial guess is needed (e.g., for preconditioning) to reduce computational cost.

 $int \ \mathbf{HYPRE_StructSMGSetNonZeroGuess} \ (\mathbf{HYPRE_StructSolver} \ solver)$

(Optional) Use a nonzero initial guess. This is the default behavior, but this routine allows the user to switch back after using SetZeroGuess.

_ 4.4.10 __

int

 $\label{eq:hypre_struct_solver} \mathbf{HYPRE_StructSolver} \ \text{solver}, \ \mathbf{int} \ \mathbf{num_pre_relax})$

(Optional) Set number of relaxation sweeps before coarse-grid correction

_ 4.4.11 _____

int **HYPRE_StructSMGSetNumPostRelax** (HYPRE_StructSolver solver, int num_post_relax)

(Optional) Set number of relaxation sweeps after coarse-grid correction

4.4.12

int HYPRE_StructSMGSetLogging (HYPRE_StructSolver solver, int logging)

(Optional) Set the amount of logging to do

int

 ${\bf HYPRE_StructSMGSetPrintLevel}~({\bf HYPRE_StructSolver}~solver, int~print_level)$

(Optional) Set the amount of printing to do to the screen

_ 4.4.14 ____

int

HYPRE_StructSMGGetNumIterations (HYPRE_StructSolver solver, int* num_iterations)

Return the number of iterations taken

 $_{-}$ 4.4.15 $_{---}$

HYPRE_StructSMGGetFinalRelativeResidualNorm (HYPRE_StructSolver solver, HYPRE_Real* norm)

Return the norm of the final relative residual

4.5

Struct CycRed Solver

Names

	${\bf HYPRE_StructCycRedSetup}~({\tt HYPRE_StructSolver}~solver,$	
	$HYPRE_StructMatrix A,$	
	HYPRE_StructVector b,	
	$HYPRE_StructVector x)$	
	Prepare to solve the system.	80
4.5.4	int	
	HYPRE_StructCycRedSolve (HYPRE_StructSolver solver,	
	HYPRE_StructMatrix A,	
	HYPRE_StructVector b,	
	HYPRE_StructVector x)	
	Solve the system	80
4.5.5	int	
	$\mathbf{HYPRE_StructCycRedSetTDim}\ (\mathbf{HYPRE_StructSolver}\ solver,\ \mathrm{int}\ \mathrm{tdim})\ldots.$	80
4.5.6	int	
	HYPRE_StructCycRedSetBase (HYPRE_StructSolver solver, int ndim,	
	int* base_index, int* base_stride)	
	(Optional) Set the base index and stride for the embedded 1D systems	80

CycRed is a cyclic reduction solver that simultaneously solves a collection of 1D tridiagonal systems embedded in a d-dimensional grid.

$_$ 4.5.1 $_$

HYPRE_StructCycRedCreate (MPI_Comm comm, HYPRE_StructSolver* solver)

Create a solver object

 $_$ 4.5.2 $_$

int HYPRE_StructCycRedDestroy (HYPRE_StructSolver solver)

Destroy a solver object

4.5.3

HYPRE_StructCycRedSetup (HYPRE_StructSolver solver,
HYPRE_StructMatrix A, HYPRE_StructVector b, HYPRE_StructVector x)

Prepare to solve the system. The coefficient data in b and x is ignored here, but information about the layout of the data may be used.

__ 4.5.4 _____

int
HYPRE_StructCycRedSolve (HYPRE_StructSolver solver,
HYPRE_StructMatrix A, HYPRE_StructVector b, HYPRE_StructVector x)

Solve the system

4.5.5

int HYPRE_StructCycRedSetTDim (HYPRE_StructSolver solver, int tdim)

(Optional) Set the dimension number for the embedded 1D tridiagonal systems. The default is tdim = 0.

 $_$ 4.5.6 $_$

HYPRE_StructCycRedSetBase (HYPRE_StructSolver solver, int ndim, int* base_index, int* base_stride)

(Optional) Set the base index and stride for the embedded 1D systems. The stride must be equal one in the dimension corresponding to the 1D systems (see HYPRE_StructCycRedSetTDim).

46

Struct PCG Solver

int	
HYPRE_StructPCGCreate (MPI_Comm comm,	
HYPRE_StructSolver* solver)	
Create a solver object	81
int	
HYPRE_StructPCGDestroy (HYPRE_StructSolver solver)	
Destroy a solver object	81
int	
HYPRE_StructDiagScaleSetup (HYPRE_StructSolver solver,	
HYPRE_StructMatrix A,	
HYPRE_StructVector y,	
$HYPRE_StructVector x)$	
Setup routine for diagonal preconditioning	82
int	
HYPRE_StructDiagScale (HYPRE_StructSolver solver,	
HYPRE_StructMatrix HA,	
HYPRE_StructVector Hy,	
HYPRE_StructVector Hx)	
Solve routine for diagonal preconditioning	82
	HYPRE_StructPCGCreate (MPI_Comm comm,

These routines should be used in conjunction with the generic interface in PCG Solver.

__ 4.6.1 ____

int
HYPRE_StructPCGCreate (MPI_Comm comm, HYPRE_StructSolver* solver)

Create a solver object

4.6.2

int HYPRE_StructPCGDestroy (HYPRE_StructSolver solver)

Destroy a solver object

4.6.3

int

HYPRE_StructDiagScaleSetup (HYPRE_StructSolver solver, HYPRE_StructMatrix A, HYPRE_StructVector y, HYPRE_StructVector x)

Setup routine for diagonal preconditioning

4.6.4

int

HYPRE_StructDiagScale (HYPRE_StructSolver solver, HYPRE_StructMatrix HA, HYPRE_StructVector Hy, HYPRE_StructVector Hx)

Solve routine for diagonal preconditioning

_ 4.7 _

Struct GMRES Solver

Names

4.7.1 int

HYPRE_StructGMRESCreate (MPI_Comm comm,

HYPRE_StructSolver* solver)

4.7.2 int

 ${\bf HYPRE_StructGMRESDestroy}~({\tt HYPRE_StructSolver~solver})$

These routines should be used in conjunction with the generic interface in GMRES Solver.

_ 4.7.1 _____

int

HYPRE_StructGMRESCreate (MPI_Comm comm, HYPRE_StructSolver* solver)

Create a solver object

4.7.2

int HYPRE_StructGMRESDestroy (HYPRE_StructSolver solver)

Destroy a solver object

4.8 .

Struct FlexGMRES Solver

Names

These routines should be used in conjunction with the generic interface in FlexGMRES Solver.

_ 4.8.1 _

int

HYPRE_StructFlexGMRESCreate (MPI_Comm comm, HYPRE_StructSolver* solver)

Create a solver object

4.8.2

int HYPRE_StructFlexGMRESDestroy (HYPRE_StructSolver solver)

Destroy a solver object

__ 4.9 _____

Struct LGMRES Solver

Names

These routines should be used in conjunction with the generic interface in LGMRES Solver.

__ 4.9.1 _____

HYPRE_StructLGMRESCreate (MPI_Comm comm, HYPRE_StructSolver* solver)

Create a solver object

 $_$ 4.9.2 $_$

int HYPRE_StructLGMRESDestroy (HYPRE_StructSolver solver)

Destroy a solver object

4.10

Struct BiCGSTAB Solver

These routines should be used in conjunction with the generic interface in BiCGSTAB Solver.

__ 4.10.1 _____

HYPRE_StructBiCGSTABCreate (MPI_Comm comm, HYPRE_StructSolver* solver)

Create a solver object

 $_$ 4.10.2 $_$

int HYPRE_StructBiCGSTABDestroy (HYPRE_StructSolver solver)

Destroy a solver object

_ 4.11 _____

Struct Hybrid Solver

Names

4.11.1 int

	HYPRE_StructHybridCreate (MPI_Comm comm, HYPRE_StructSolver* solver)	
	Create a solver object	8'
4 1 1 0	•	O
4.11.2	int HVDDE Struct Hybrid Destroy (HVDDE Struct Selven gelven)	
	HYPRE_StructHybridDestroy (HYPRE_StructSolver solver) Destroy a solver object	88
	· · ·	0
4.11.3	int	
	HYPRE_StructHybridSetup (HYPRE_StructSolver solver,	
	HYPRE_StructMatrix A, HYPRE_StructVector b,	
	HYPRE_StructVector x)	
	Prepare to solve the system	88
		00
4.11.4	int	
	HYPRE_StructHybridSolve (HYPRE_StructSolver solver,	
	HYPRE_StructMatrix A,	
	HYPRE_StructVector b,	
	HYPRE_StructVector x) Solve the system	88
		00
4.11.5	int	
	HYPRE_StructHybridSetTol (HYPRE_StructSolver solver, HYPRE_Real tol)	
	(Optional) Set the convergence tolerance	88
4.11.6	int	
	$\mathbf{HYPRE_StructHybridSetConvergenceTol} \ (\mathbf{HYPRE_StructSolver} \ solver,$	
	HYPRE_Real cf_tol)	
	(Optional) Set an accepted convergence tolerance for diagonal scaling (DS).	
		89
4.11.7	int	
	HYPRE_StructHybridSetDSCGMaxIter (HYPRE_StructSolver solver,	
	$int ds_max_its)$	
	(Optional) Set maximum number of iterations for diagonal scaling (DS)	89
4.11.8	int	
	HYPRE_StructHybridSetPCGMaxIter (HYPRE_StructSolver solver,	
	int pre_max_its)	
	(Optional) Set maximum number of iterations for general preconditioner	
	(PRE)	89
4.11.9	int	
4.11.0	HYPRE_StructHybridSetTwoNorm (HYPRE_StructSolver solver,	
	int two_norm)	
	(Optional) Use the two-norm in stopping criteria	89
4 11 10		
4.11.10	int HYPRE_StructHybridSetRelChange (HYPRE_StructSolver solver,	
	int rel_change)	
	(Optional) Additionally require that the relative difference in successive it-	
	erates be small	90
4 4 4 4 4 4		90
4.11.11	int	
	HYPRE_StructHybridSetSolverType (HYPRE_StructSolver solver,	
	int solver_type) (Ontional) Set the type of Kryley selven to use	00
	(Optional) Set the type of Krylov solver to use	90
4.11.12	int	

	HYPRE_StructHybridSetKDim (HYPRE_StructSolver solver, int k_dim) (Optional) Set the maximum size of the Krylov space when using GMRES	90
4.11.13	int	
	${\bf HYPRE_StructHybridSetPrecond}~({\bf HYPRE_StructSolver}~solver,$	
	HYPRE_PtrToStructSolverFcn precond,	
	HYPRE_PtrToStructSolverFcn	
	precond_setup,	
	HYPRE_StructSolver precond_solver)	00
	(Optional) Set the preconditioner to use	90
4.11.14	int	
	HYPRE_StructHybridSetLogging (HYPRE_StructSolver solver, int logging)	
	(Optional) Set the amount of logging to do	91
4.11.15	int	
	HYPRE_StructHybridSetPrintLevel (HYPRE_StructSolver solver,	
	int print_level)	
	(Optional) Set the amount of printing to do to the screen	91
4.11.16	int	
	${\bf HYPRE_StructHybridGetNumIterations}~({\bf HYPRE_StructSolver}~solver,$	
	int* num_its)	
	Return the number of iterations taken	91
4.11.17	int	
	$\mathbf{HYPRE_StructHybridGetDSCGNumIterations} \ (\mathbf{HYPRE_StructSolver}$	
	solver, int* ds_num_its)	
	Return the number of diagonal scaling iterations taken	91
4.11.18	int	
	HYPRE_StructHybridGetPCGNumIterations (HYPRE_StructSolver	
	solver, int* pre_num_its)	
	Return the number of general preconditioning iterations taken	92
4.11.19	int	
	$HYPRE_StructHybridGetFinalRelativeResidualNorm$	
	(HYPRE_StructSolver	
	solver,	
	HYPRE_Real*	
	$\operatorname{norm})$ Return the norm of the final relative residual	92
	recount one norm of the final relative restaud	34

_ 4.11.1 _

 $\begin{array}{l} \text{int} \\ \textbf{HYPRE_StructHybridCreate} \text{ (MPI_Comm comm, HYPRE_StructSolver*} \\ \text{solver)} \end{array}$

Create a solver object

int HYPRE_StructHybridDestroy (HYPRE_StructSolver solver)

Destroy a solver object

__ 4.11.3 _____

int

HYPRE_StructHybridSetup (HYPRE_StructSolver solver, HYPRE_StructMatrix A, HYPRE_StructVector b, HYPRE_StructVector x)

Prepare to solve the system. The coefficient data in ${\tt b}$ and ${\tt x}$ is ignored here, but information about the layout of the data may be used.

_ 4.11.4 _____

int **HYPRE_StructHybridSolve** (HYPRE_StructSolver solver,
HYPRE_StructMatrix A, HYPRE_StructVector b, HYPRE_StructVector x)

Solve the system

4.11.5

int

HYPRE_StructHybridSetTol (HYPRE_StructSolver solver, HYPRE_Real tol)

(Optional) Set the convergence tolerance

HYPRE_StructHybridSetConvergenceTol (HYPRE_StructSolver solver, HYPRE_Real cf_tol)

(Optional) Set an accepted convergence tolerance for diagonal scaling (DS). The solver will switch preconditioners if the convergence of DS is slower than cf_tol.

4.11.7

int

 $\label{lem:hypre_struct} \begin{tabular}{ll} HYPRE_StructSolver solver, int ds_max_its) \end{tabular}$

(Optional) Set maximum number of iterations for diagonal scaling (DS). The solver will switch preconditioners if DS reaches ds_max_its.

4.11.8

int

HYPRE_StructHybridSetPCGMaxIter (HYPRE_StructSolver solver, int pre_max_its)

(Optional) Set maximum number of iterations for general preconditioner (PRE). The solver will stop if PRE reaches pre_max_its.

4.11.9

int

HYPRE_StructHybridSetTwoNorm (HYPRE_StructSolver solver, int two_norm)

(Optional) Use the two-norm in stopping criteria

int

 $\label{lem:hyprid} \textbf{HYPRE_StructHybridSetRelChange} \ (\textbf{HYPRE_StructSolver solver}, \ \textbf{int rel_change})$

(Optional) Additionally require that the relative difference in successive iterates be small

4.11.11

HYPRE_StructHybridSetSolverType (HYPRE_StructSolver solver, int solver_type)

(Optional) Set the type of Krylov solver to use.

Current krylov methods set by solver_type are:

- 0 PCG (default)
- 1 GMRES
- 2 BiCGSTAB

__ 4.11.12 ____

int HYPRE_StructHybridSetKDim (HYPRE_StructSolver solver, int k_dim)

(Optional) Set the maximum size of the Krylov space when using GMRES

_ 4.11.13 _____

int

HYPRE_StructHybridSetPrecond (HYPRE_StructSolver solver, HYPRE_PtrToStructSolverFcn precond, HYPRE_PtrToStructSolverFcn precond_solver)

(Optional) Set the preconditioner to use

4 11 14

int HYPRE_StructHybridSetLogging (HYPRE_StructSolver solver, int logging)

(Optional) Set the amount of logging to do

__ 4.11.15 _____

HYPRE_StructHybridSetPrintLevel (HYPRE_StructSolver solver, int print_level)

(Optional) Set the amount of printing to do to the screen

_ 4.11.16 __

int

HYPRE_StructHybridGetNumIterations (HYPRE_StructSolver solver, int* num_its)

Return the number of iterations taken

4.11.17

HYPRE_StructHybridGetDSCGNumIterations (HYPRE_StructSolver solver, int* ds_num_its)

Return the number of diagonal scaling iterations taken

HYPRE_StructHybridGetPCGNumIterations (HYPRE_StructSolver solver, int* pre_num_its)

Return the number of general preconditioning iterations taken

_ 4.11.19 _____

int

$HYPRE_StructHybridGetFinalRelativeResidualNorm$

(HYPRE_StructSolver solver, HYPRE_Real* norm)

Return the norm of the final relative residual

4.12

Struct LOBPCG Eigensolver

Names

4.12.1 int

HYPRE_StructSetupInterpreter (mv_InterfaceInterpreter* i)

Load interface interpreter. 92

4.12.2 int

HYPRE_StructSetupMatvec (HYPRE_MatvecFunctions* mv)

Load Matvec interpreter with hypre_StructKrylov functions 93

These routines should be used in conjunction with the generic interface in LOBPCG Eigensolver.

 $_$ 4.12.1 $_$

int HYPRE_StructSetupInterpreter (mv_InterfaceInterpreter* i)

 $Load\ interface\ interpreter.\ Vector\ part\ loaded\ with\ hypre_StructKrylov\ functions\ and\ multivector\ part\ loaded\ with\ mv_TempMultiVector\ functions.$

4.12.2

 $int \ \mathbf{HYPRE_StructSetupMatvec} \ (HYPRE_MatvecFunctions*\ mv)$

 ${\bf Load\ Matvec\ interpreter\ with\ hypre_StructKrylov\ functions}$

5

SStruct Solvers

Names		
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		120

These solvers use matrix/vector storage schemes that are taylored to semi-structured grid problems.

_ 5.1 _

SStruct Solvers

Names

5.1.1	typedef struct	hypre_SStructSolver_struct *HYPRE_SStructSolver	
	The so	lver object	95

_ 5.1.1 _

 $typedef\ struct\ hypre_SStructSolver_struct\ \textbf{*HYPRE_SStructSolver}$

The solver object

_ 5.2 .

SStruct SysPFMG Solver

Names		
5.2.1	int	
	HYPRE_SStructSysPFMGCreate (MPI_Comm comm,	
	HYPRE_SStructSolver* solver)	
	Create a solver object	97
5.2.2	int	
J	HYPRE_SStructSysPFMGDestroy (HYPRE_SStructSolver solver)	
	Destroy a solver object.	97
5.2.3	·	
3.2.3	int HYPRE_SStructSysPFMGSetup (HYPRE_SStructSolver solver,	
	HYPRE_StructSolver solver, HYPRE_StructMatrix A,	
	HYPRE_SStructVector b,	
	· ·	
	HYPRE_SStructVector x)	07
	Prepare to solve the system.	97
5.2.4	int	
	HYPRE_SStructSysPFMGSolve (HYPRE_SStructSolver solver,	
	HYPRE_SStructMatrix A,	
	HYPRE_SStructVector b,	
	HYPRE_SStructVector x)	
	Solve the system	97
5.2.5	int	
0.2.0	HYPRE_SStructSysPFMGSetTol (HYPRE_SStructSolver solver,	
	HYPRE_Real tol)	
	(Optional) Set the convergence tolerance	98
		30
5.2.6	int	
	HYPRE_SStructSysPFMGSetMaxIter (HYPRE_SStructSolver solver,	
	int max_iter)	
	(Optional) Set maximum number of iterations	98
5.2.7	int	
	HYPRE_SStructSysPFMGSetRelChange (HYPRE_SStructSolver solver,	
	int rel_change)	
	(Optional) Additionally require that the relative difference in successive it-	
	erates be small	98
TO 0		
5.2.8	int	

	HYPRE_SStructSysPFMGSetZeroGuess (HYPRE_SStructSolver solver) (Optional) Use a zero initial guess.	98
5.2.9	int HYPRE_SStructSysPFMGSetNonZeroGuess (HYPRE_SStructSolver	
	solver)	
	(Optional) Use a nonzero initial guess	99
5.2.10	int	
0.2.10	HYPRE_SStructSysPFMGSetRelaxType (HYPRE_SStructSolver solver, int relax_type)	
	(Optional) Set relaxation type.	99
F 0 11	,	00
5.2.11	int HYPRE_SStructSysPFMGSetJacobiWeight (HYPRE_SStructSolver solver, HYPRE_Real weight)	
	(Optional) Set Jacobi Weight	99
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5.2.12	int HYPRE_SStructSysPFMGSetNumPreRelax (HYPRE_SStructSolver solver, int num_pre_relax)	
	(Optional) Set number of relaxation sweeps before coarse-grid correction .	99
5.2.13	int	
0.2.10	HYPRE_SStructSysPFMGSetNumPostRelax (HYPRE_SStructSolver solver, int num_post_relax)	
	(Optional) Set number of relaxation sweeps after coarse-grid correction	100
		100
5.2.14	int	
	HYPRE_SStructSysPFMGSetSkipRelax (HYPRE_SStructSolver solver,	
	int skip_relax) (Optional) Skip relaxation on certain grids for isotropic problems	100
		100
5.2.15	int	
	HYPRE_SStructSysPFMGSetLogging (HYPRE_SStructSolver solver,	
	int logging)	100
	(Optional) Set the amount of logging to do	100
5.2.16	int	
	HYPRE_SStructSysPFMGSetPrintLevel (HYPRE_SStructSolver solver,	
	int print_level)	100
	(Optional) Set the amount of printing to do to the screen	100
5.2.17	int	
	HYPRE_SStructSysPFMGGetNumIterations (HYPRE_SStructSolver	
	solver, int* num_iterations)	101
	Return the number of iterations taken	101
5.2.18	int	
	HYPRE_SStructSysPFMGGetFinalRelativeResidualNorm	~ .
	(HYPRE_SStruct	Solver
	solver,	
	HYPRE_Real*	
	$\operatorname{norm})$ Return the norm of the final relative residual	101
	1000 and 100 m of one form 100000 100 and	101

SysPFMG is a semicoarsening multigrid solver similar to PFMG, but for systems of PDEs. For periodic problems, users should try to set the grid size in periodic dimensions to be as close to a power-of-two as possible (for more details, see Struct PFMG Solver).

HYPRE_SStructSysPFMGCreate (MPI_Comm comm, HYPRE_SStructSolver* solver)

Create a solver object

__ 5.2.2 _

int HYPRE_SStructSysPFMGDestroy (HYPRE_SStructSolver solver)

Destroy a solver object. An object should be explicitly destroyed using this destructor when the user's code no longer needs direct access to it. Once destroyed, the object must not be referenced again. Note that the object may not be deallocated at the completion of this call, since there may be internal package references to the object. The object will then be destroyed when all internal reference counts go to zero.

__ 5.2.3 ____

HYPRE_SStructSysPFMGSetup (HYPRE_SStructSolver solver, HYPRE_SStructMatrix A, HYPRE_SStructVector b, HYPRE_SStructVector x)

Prepare to solve the system. The coefficient data in b and x is ignored here, but information about the layout of the data may be used.

_ 5.2.4 _

int

HYPRE_SStructSysPFMGSolve (HYPRE_SStructSolver solver, HYPRE_SStructMatrix A, HYPRE_SStructVector b, HYPRE_SStructVector x)

Solve the system

HYPRE_SStructSysPFMGSetTol (HYPRE_SStructSolver solver, HYPRE_Real tol)

(Optional) Set the convergence tolerance

5.2.6

HYPRE_SStructSysPFMGSetMaxIter (HYPRE_SStructSolver solver, int max_iter)

(Optional) Set maximum number of iterations

__ 5.2.7 _____

int **HYPRE_SStructSysPFMGSetRelChange** (HYPRE_SStructSolver solver, int rel_change)

(Optional) Additionally require that the relative difference in successive iterates be small

__ 5.2.8 _____

int HYPRE_SStructSysPFMGSetZeroGuess (HYPRE_SStructSolver solver)

(Optional) Use a zero initial guess. This allows the solver to cut corners in the case where a zero initial guess is needed (e.g., for preconditioning) to reduce computational cost.

int

HYPRE_SStructSysPFMGSetNonZeroGuess (HYPRE_SStructSolver solver)

(Optional) Use a nonzero initial guess. This is the default behavior, but this routine allows the user to switch back after using SetZeroGuess.

_ 5.2.10 __

int

HYPRE_SStructSysPFMGSetRelaxType (HYPRE_SStructSolver solver, int relax_type)

(Optional) Set relaxation type.

Current relaxation methods set by relax_type are:

- 0 Jacobi
- 1 Weighted Jacobi (default)
- 2 Red/Black Gauss-Seidel (symmetric: RB pre-relaxation, BR post-relaxation)

__ 5.2.11 _____

int

 $\label{lem:hypre_structSysPFMGSetJacobiWeight} \begin{tabular}{l} HYPRE_SStructSolver solver, \\ HYPRE_Real weight) \end{tabular}$

(Optional) Set Jacobi Weight

5.2.12

int

HYPRE_SStructSysPFMGSetNumPreRelax (HYPRE_SStructSolver solver, int num_pre_relax)

(Optional) Set number of relaxation sweeps before coarse-grid correction

int HYPRE_SStructSysPFMGSetNumPostRelax (HYPRE_SStructSolver solver, int num_post_relax)

(Optional) Set number of relaxation sweeps after coarse-grid correction

5.2.14

int

 $\label{eq:hypre_structSysPFMGSetSkipRelax} \mbox{(HYPRE_SStructSolver solver, int skip_relax)}$

(Optional) Skip relaxation on certain grids for isotropic problems. This can greatly improve efficiency by eliminating unnecessary relaxations when the underlying problem is isotropic.

___ 5.2.15 _____

int

 $\label{eq:hypre_structSysPFMGSetLogging} \ (\mbox{HYPRE_SStructSolver solver}, \mbox{ int logging})$

(Optional) Set the amount of logging to do

_ 5.2.16 __

ınt

HYPRE_SStructSysPFMGSetPrintLevel (HYPRE_SStructSolver solver, int print_level)

(Optional) Set the amount of printing to do to the screen

HYPRE_SStructSysPFMGGetNumIterations (HYPRE_SStructSolver solver, int* num_iterations)

Return the number of iterations taken

5.2.18

int

$HYPRE_SStructSysPFMGGetFinalRelativeResidualNorm$

(HYPRE_SStructSolver solver, HYPRE_Real* norm)

Return the norm of the final relative residual

5.3

SStruct Split Solver

int HYPRE_SStructSplitCreate (MPI_Comm comm,	
HYPRE_SStructSolver* solver)	
Create a solver object	102
int	
HYPRE_SStructSplitDestroy (HYPRE_SStructSolver solver)	
Destroy a solver object.	102
int	
HYPRE_SStructSplitSetup (HYPRE_SStructSolver solver,	
HYPRE_SStructMatrix A,	
HYPRE_SStructVector b,	
HYPRE_SStructVector x)	
Prepare to solve the system	103
int	
HYPRE_SStructSplitSolve (HYPRE_SStructSolver solver,	
HYPRE_SStructMatrix A,	
HYPRE_SStructVector b,	
$HYPRE_SStructVector x)$	
Solve the system	103
int	
	HYPRE_SStructSplitCreate (MPI_Comm comm,

	HYPRE_SStructSplitSetTol (HYPRE_SStructSolver solver, HYPRE_Real tol) (Optional) Set the convergence tolerance	103
5.3.6	int	
	HYPRE_SStructSplitSetMaxIter (HYPRE_SStructSolver solver,	
	int max_iter)	
	(Optional) Set maximum number of iterations	103
5.3.7	int	
	HYPRE_SStructSplitSetZeroGuess (HYPRE_SStructSolver solver)	
	(Optional) Use a zero initial guess.	104
5.3.8	int	
	HYPRE_SStructSplitSetNonZeroGuess (HYPRE_SStructSolver solver)	
	(Optional) Use a nonzero initial guess	104
5.3.9	int	
	HYPRE_SStructSplitSetStructSolver (HYPRE_SStructSolver solver,	
	int ssolver)	
	(Optional) Set up the type of diagonal struct solver	104
5.3.10	int	
	HYPRE_SStructSplitGetNumIterations (HYPRE_SStructSolver solver,	
	$int* num_iterations)$	
	Return the number of iterations taken	104
5.3.11	int	
	$HYPRE_SStructSplitGetFinalRelativeResidualNorm$	
	$(HYPRE_SStructSolver)$	
	solver,	
	$\mathrm{HYPRE} ext{-}\mathrm{Real}^*$	
	norm)	
	Return the norm of the final relative residual	105

5.3.1

int **HYPRE_SStructSplitCreate** (MPI_Comm comm, HYPRE_SStructSolver* solver)

Create a solver object

 $_$ 5.3.2 $_$

int HYPRE_SStructSplitDestroy (HYPRE_SStructSolver solver)

Destroy a solver object. An object should be explicitly destroyed using this destructor when the user's code no longer needs direct access to it. Once destroyed, the object must not be referenced again. Note that the object may not be deallocated at the completion of this call, since there may be internal package references to the object. The object will then be destroyed when all internal reference counts go to zero.

_ 5.3.3 _

int
HYPRE_SStructSplitSetup (HYPRE_SStructSolver solver,
HYPRE_SStructMatrix A, HYPRE_SStructVector b, HYPRE_SStructVector x)

Prepare to solve the system. The coefficient data in b and x is ignored here, but information about the layout of the data may be used.

_ 5.3.4 _

HYPRE_SStructSplitSolve (HYPRE_SStructSolver solver,
HYPRE_SStructMatrix A, HYPRE_SStructVector b, HYPRE_SStructVector x)

Solve the system

5.3.5

int
HYPRE_SStructSplitSetTol (HYPRE_SStructSolver solver, HYPRE_Real tol)

(Optional) Set the convergence tolerance

 $_{-}$ 5.3.6 $_{-}$

int
HYPRE_SStructSplitSetMaxIter (HYPRE_SStructSolver solver, int max_iter)

(Optional) Set maximum number of iterations

5.3.7

int HYPRE_SStructSplitSetZeroGuess (HYPRE_SStructSolver solver)

(Optional) Use a zero initial guess. This allows the solver to cut corners in the case where a zero initial guess is needed (e.g., for preconditioning) to reduce computational cost.

 $_$ 5.3.8 $_$

int HYPRE_SStructSplitSetNonZeroGuess (HYPRE_SStructSolver solver)

(Optional) Use a nonzero initial guess. This is the default behavior, but this routine allows the user to switch back after using SetZeroGuess.

5.3.9

 $\mbox{\bf HYPRE_SStructSplitSetStructSolver}$ (HYPRE_SStructSolver solver, int ssolver)

(Optional) Set up the type of diagonal struct solver. Either ssolver is set to HYPRE_SMG or HYPRE_PFMG.

5.3.10

HYPRE_SStructSplitGetNumIterations (HYPRE_SStructSolver solver, int* num_iterations)

Return the number of iterations taken

5.3.11

int

${\bf HYPRE_SStructSplitGetFinalRelativeResidualNorm}$

(HYPRE_SStructSolver solver, HYPRE_Real* norm)

Return the norm of the final relative residual

___ 5.4 _____

SStruct FAC Solver

Names		
5.4.1	int	
	HYPRE_SStructFACCreate (MPI_Comm comm,	
	HYPRE_SStructSolver* solver)	
	Create a solver object	107
5.4.2	int	
	HYPRE_SStructFACDestroy2 (HYPRE_SStructSolver solver)	
	Destroy a solver object.	108
5.4.3	int	
	HYPRE_SStructFACAMR_RAP (HYPRE_SStructMatrix A,	
	int (*rfactors)[HYPRE_MAXDIM],	
	HYPRE_SStructMatrix* fac_A)	
	Re-distribute the composite matrix so that the amr hierarchy is approximately	
	nested.	108
5.4.4	int	
	HYPRE_SStructFACSetup2 (HYPRE_SStructSolver solver,	
	HYPRE_SStructMatrix A,	
	HYPRE_SStructVector b,	
	HYPRE_SStructVector x)	
	Set up the FAC solver structure	108
5.4.5	int	
	HYPRE_SStructFACSolve3 (HYPRE_SStructSolver solver,	
	HYPRE_SStructMatrix A,	
	HYPRE_SStructVector b,	
	HYPRE_SStructVector x)	
	Solve the system	108
5.4.6	int	
	HYPRE_SStructFACSetPLevels (HYPRE_SStructSolver solver, int nparts,	
	int* plevels)	
	Set up amr structure	109
5.4.7	int	

	HYPRE_SStructFACSetPRefinements (HYPRE_SStructSolver solver,	
	int nparts,	
	int (*rfactors)[HYPRE_MAXDIM])	
	Set up amr refinement factors	10
5.4.8	int	
0.4.0	HYPRE_SStructFACZeroCFSten (HYPRE_SStructMatrix A,	
	HYPRE_SStructGrid grid, int part,	
	int rfactors[HYPRE_MAXDIM])	
	(Optional, but user must make sure that they do this function otherwise)	
	Zero off the coarse level stencils reaching into a fine level grid	10
		10
5.4.9	int	
	HYPRE_SStructFACZeroFCSten (HYPRE_SStructMatrix A,	
	HYPRE_SStructGrid grid, int part)	
	(Optional, but user must make sure that they do this function otherwise)	
	Zero off the fine level stencils reaching into a coarse level grid	10
5.4.10	int	
0.1.10	HYPRE_SStructFACZeroAMRMatrixData (HYPRE_SStructMatrix A,	
	int part_crse, int	
	rfactors[HYPRE_MAXDIM])	
	(Optional, but user must make sure that they do this function otherwise)	
	Places the identity in the coarse grid matrix underlying the fine patches	11
	· · · · · · · · · · · · · · · · · · ·	
5.4.11	int	
	HYPRE_SStructFACZeroAMRVectorData (HYPRE_SStructVector b,	
	int* plevels, int	
	(*rfactors)[HYPRE_MAXDIM]	
)	
	(Optional, but user must make sure that they do this function otherwise)	
	Places zeros in the coarse grid vector underlying the fine patches	11
5.4.12	int	
	HYPRE_SStructFACSetMaxLevels (HYPRE_SStructSolver solver,	
	int max_levels)	
	(Optional) Set maximum number of FAC levels	11
F 4 19	· ·	
5.4.13	int	
	HYPRE_SStructFACSetTol (HYPRE_SStructSolver solver, HYPRE_Real tol)	11
	(Optional) Set the convergence tolerance	11
5.4.14	int	
	HYPRE_SStructFACSetMaxIter (HYPRE_SStructSolver solver,	
	$\operatorname{int} \ \operatorname{max_iter})$	
	(Optional) Set maximum number of iterations	11
5.4.15	int	
0.4.10	HYPRE_SStructFACSetRelChange (HYPRE_SStructSolver solver,	
	int rel_change)	
	(Optional) Additionally require that the relative difference in successive it-	
	erates be small	11
		11
5.4.16	int	
	HYPRE_SStructFACSetZeroGuess (HYPRE_SStructSolver solver)	
	(Optional) Use a zero initial guess	11
5.4.17	int	

	HYPRE_SStructFACSetNonZeroGuess (HYPRE_SStructSolver solver) (Optional) Use a nonzero initial guess.	111
5.4.18	int	
	HYPRE_SStructFACSetRelaxType (HYPRE_SStructSolver solver, int relax_type)	
	(Optional) Set relaxation type.	112
5.4.19	int	
	HYPRE_SStructFACSetJacobiWeight (HYPRE_SStructSolver solver, HYPRE_Real weight)	
	(Optional) Set Jacobi weight if weighted Jacobi is used	112
5.4.20	int	
	HYPRE_SStructFACSetNumPreRelax (HYPRE_SStructSolver solver, int num_pre_relax)	
	(Optional) Set number of relaxation sweeps before coarse-grid correction .	112
5.4.21	int	
	HYPRE_SStructFACSetNumPostRelax (HYPRE_SStructSolver solver, int num_post_relax)	
	(Optional) Set number of relaxation sweeps after coarse-grid correction	112
5.4.22	int HYPRE_SStructFACSetCoarseSolverType (HYPRE_SStructSolver solver, int csolver_type)	
	(Optional) Set coarsest solver type.	113
5.4.23	int HYPRE_SStructFACSetLogging (HYPRE_SStructSolver solver, int logging)	
	(Optional) Set the amount of logging to do	113
5.4.24	int	
	HYPRE_SStructFACGetNumIterations (HYPRE_SStructSolver solver, int* num_iterations)	
	Return the number of iterations taken	113
5.4.25	$\operatorname{int} \\ \mathbf{HYPRE_SStructFACGetFinalRelativeResidualNorm}$	
	(HYPRE_SStructSolver	
	solver,	
	$\mathrm{HYPRE}_{-}\mathrm{Real}^*$	
	norm)	113
	Return the norm of the final relative residual	113

 $\begin{array}{l} \operatorname{int} \\ \mathbf{HYPRE_SStructFACCreate} \ (\operatorname{MPI_Comm} \ \operatorname{comm}, \ \operatorname{HYPRE_SStructSolver*} \ \operatorname{solver}) \end{array}$

Create a solver object

int HYPRE_SStructFACDestroy2 (HYPRE_SStructSolver solver)

Destroy a solver object. An object should be explicitly destroyed using this destructor when the user's code no longer needs direct access to it. Once destroyed, the object must not be referenced again. Note that the object may not be deallocated at the completion of this call, since there may be internal package references to the object. The object will then be destroyed when all internal reference counts go to zero.

5.4.3

int

HYPRE_SStructFACAMR_RAP (HYPRE_SStructMatrix A, int (*rfactors)[HYPRE_MAXDIM], HYPRE_SStructMatrix* fac_A)

Re-distribute the composite matrix so that the amr hierarchy is approximately nested. Coarse underlying operators are also formed.

5.4.4

int

HYPRE_SStructFACSetup2 (HYPRE_SStructSolver solver, HYPRE_SStructMatrix A, HYPRE_SStructVector b, HYPRE_SStructVector x)

Set up the FAC solver structure

_ 5.4.5 _

inf

HYPRE_SStructFACSolve3 (HYPRE_SStructSolver solver, HYPRE_SStructMatrix A, HYPRE_SStructVector b, HYPRE_SStructVector x)

Solve the system

HYPRE_SStructFACSetPLevels (HYPRE_SStructSolver solver, int nparts, int* plevels)

Set up amr structure

5.4.7

int

 $\label{lem:hypre_struct} \textbf{HYPRE_SStructSolver solver}, int \\ nparts, int (*rfactors)[HYPRE_MAXDIM])$

Set up amr refinement factors

__ 5.4.8 _____

int
HYPRE_SStructFACZeroCFSten (HYPRE_SStructMatrix A,
HYPRE_SStructGrid grid, int part, int rfactors[HYPRE_MAXDIM])

(Optional, but user must make sure that they do this function otherwise) Zero off the coarse level stencils reaching into a fine level grid

5.4.9

HYPRE_SStructFACZeroFCSten (HYPRE_SStructMatrix A, HYPRE_SStructGrid grid, int part)

(Optional, but user must make sure that they do this function otherwise) Zero off the fine level stencils reaching into a coarse level grid

int
HYPRE_SStructFACZeroAMRMatrixData (HYPRE_SStructMatrix A, int
part_crse, int rfactors[HYPRE_MAXDIM])

(Optional, but user must make sure that they do this function otherwise) Places the identity in the coarse grid matrix underlying the fine patches. Required between each pair of amr levels.

5.4.11

HYPRE_SStructFACZeroAMRVectorData (HYPRE_SStructVector b, int* plevels, int (*rfactors)[HYPRE_MAXDIM])

(Optional, but user must make sure that they do this function otherwise) Places zeros in the coarse grid vector underlying the fine patches. Required between each pair of amr levels.

 $_$ 5.4.12 $_$

int ${\bf HYPRE_SStructFACSetMaxLevels}$ (${\bf HYPRE_SStructSolver}$ solver, int max_levels)

(Optional) Set maximum number of FAC levels

_ 5.4.13 _

HYPRE_SStructFACSetTol (HYPRE_SStructSolver solver, HYPRE_Real tol)

(Optional) Set the convergence tolerance

int

HYPRE_SStructFACSetMaxIter (HYPRE_SStructSolver solver, int max_iter)

(Optional) Set maximum number of iterations

 $_$ 5.4.15 $_$

HYPRE_SStructFACSetRelChange (HYPRE_SStructSolver solver, int rel_change)

(Optional) Additionally require that the relative difference in successive iterates be small

 $_$ 5.4.16 $_$

int HYPRE_SStructFACSetZeroGuess (HYPRE_SStructSolver solver)

(Optional) Use a zero initial guess. This allows the solver to cut corners in the case where a zero initial guess is needed (e.g., for preconditioning) to reduce computational cost.

5.4.17

int HYPRE_SStructFACSetNonZeroGuess (HYPRE_SStructSolver solver)

(Optional) Use a nonzero initial guess. This is the default behavior, but this routine allows the user to switch back after using SetZeroGuess.

HYPRE_SStructFACSetRelaxType (HYPRE_SStructSolver solver, int relax_type)

(Optional) Set relaxation type. See HYPRE_SStructSysPFMGSetRelaxType for appropriate values of $relax_type$.

5.4.19

HYPRE_SStructFACSetJacobiWeight (HYPRE_SStructSolver solver, HYPRE_Real weight)

(Optional) Set Jacobi weight if weighted Jacobi is used

___ 5.4.20 _____

int **HYPRE_SStructFACSetNumPreRelax** (HYPRE_SStructSolver solver, int num_pre_relax)

(Optional) Set number of relaxation sweeps before coarse-grid correction

5.4.21

int **HYPRE_SStructFACSetNumPostRelax** (HYPRE_SStructSolver solver, int num_post_relax)

(Optional) Set number of relaxation sweeps after coarse-grid correction

int

 $\label{lem:hypre_struct} \begin{tabular}{ll} HYPRE_SStructSolver solver, int csolver_type \\ \end{tabular}$

(Optional) Set coarsest solver type.

Current solver types set by csolver_type are:

- 1 SysPFMG-PCG (default)
- 2 SysPFMG

5.4.23

int HYPRE_SStructFACSetLogging (HYPRE_SStructSolver solver, int logging)

(Optional) Set the amount of logging to do

 $_$ 5.4.24 $_$

int

HYPRE_SStructFACGetNumIterations (HYPRE_SStructSolver solver, int* num_iterations)

Return the number of iterations taken

 $_$ 5.4.25 $__$

int

 $HYPRE_SStructFACGetFinalRelativeResidualNorm$

(HYPRE_SStructSolver solver, HYPRE_Real* norm)

Return the norm of the final relative residual

_ 5.5 _

SStruct Maxwell Solver

\mathbf{Names}		
5.5.1	int	
	HYPRE_SStructMaxwellCreate (MPI_Comm comm,	
	HYPRE_SStructSolver* solver)	
	Create a solver object	116
5.5.2	int	
	HYPRE_SStructMaxwellDestroy (HYPRE_SStructSolver solver)	
	Destroy a solver object	116
5.5.3	int	
0.0.0	HYPRE_SStructMaxwellSetup (HYPRE_SStructSolver solver,	
	HYPRE_SStructMatrix A,	
	HYPRE_SStructVector b,	
	HYPRE_SStructVector x)	
	Prepare to solve the system	116
5.5.4	int	
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	HYPRE_SStructMatrix A,	
	HYPRE_SStructVector b,	
	HYPRE_SStructVector x)	
	Solve the system.	117
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5.5.5	int HYPRE_SStructMaxwellSolve2 (HYPRE_SStructSolver solver,	
	HYPRE_SStructMatrix A,	
	HYPRE_SStructVector b,	
	HYPRE_SStructVector x)	
	Solve the system.	117
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5.5.0	int HYPRE_SStructMaxwellSetGrad (HYPRE_SStructSolver solver,	
	HYPRE_ParCSRMatrix T)	
	Sets the gradient operator in the Maxwell solver	117
		11.
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	Sets the coarsening factor	117
		111
5.5.8	int	
	HYPRE_SStructMaxwellPhysBdy (HYPRE_SStructGrid* grid_l,	
	int num_levels,	
	int rfactors[HYPRE_MAXDIM], int*** BdryRanks_ptr,	
	int** BdryRanks_ptr,	
	Finds the physical boundary row ranks on all levels	118
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5.5.9	int	

	$\mathbf{HYPRE_SStructMaxwellEliminateRowsCols} \ (\mathbf{HYPRE_ParCSRMatrix}$
	parA, int nrows, int* rows)
	Eliminates the rows and cols corresponding to the physical boundary in a parcsr matrix
5.5.10	int
0.0.10	HYPRE_SStructMaxwellZeroVector (HYPRE_ParVector b, int* rows, int nrows)
	Zeros the rows corresponding to the physical boundary in a par vector
5.5.11	int
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	(Optional) Set the constant coefficient flag- Nedelec interpolation used
E E 10	
5.5.12	int HYPRE_SStructMaxwellGrad (HYPRE_SStructGrid grid, HYPRE_ParCSRMatrix* T)
	(Optional) Creates a gradient matrix from the grid
5.5.13	int
0.0.10	HYPRE_SStructMaxwellSetTol (HYPRE_SStructSolver solver, HYPRE_Real tol)
	(Optional) Set the convergence tolerance
5 5 1 4	
5.5.14	int HYPRE_SStructMaxwellSetMaxIter (HYPRE_SStructSolver solver, int max_iter)
	(Optional) Set maximum number of iterations
F F 1 F	
5.5.15	int HYPRE_SStructMaxwellSetRelChange (HYPRE_SStructSolver solver, int rel_change)
	(Optional) Additionally require that the relative difference in successive iterates be small
5.5.16	int HYPRE_SStructMaxwellSetNumPreRelax (HYPRE_SStructSolver solver, int num_pre_relax)
	(Optional) Set number of relaxation sweeps before coarse-grid correction .
E E 17	
5.5.17	int HYPRE_SStructMaxwellSetNumPostRelax (HYPRE_SStructSolver solver, int num_post_relax)
	(Optional) Set number of relaxation sweeps after coarse-grid correction
5.5.18	int
	HYPRE_SStructMaxwellSetLogging (HYPRE_SStructSolver solver,
	int logging) (Ontional) Set the amount of logging to do
	(Optional) Set the amount of logging to do
5.5.19	int HYPRE_SStructMaxwellGetNumIterations (HYPRE_SStructSolver solver,
	int* num_iterations)
	Return the number of iterations taken
5.5.20	int

$HYPRE_SStructMaxwellGetFinalRelativeResidualNorm$

(HYPRE_SStructSolver solver, HYPRE_Real* norm)

5.5.1

int **HYPRE_SStructMaxwellCreate** (MPI_Comm comm, HYPRE_SStructSolver* solver)

Create a solver object

5.5.2

int HYPRE_SStructMaxwellDestroy (HYPRE_SStructSolver solver)

Destroy a solver object. An object should be explicitly destroyed using this destructor when the user's code no longer needs direct access to it. Once destroyed, the object must not be referenced again. Note that the object may not be deallocated at the completion of this call, since there may be internal package references to the object. The object will then be destroyed when all internal reference counts go to zero.

_ 5.5.3 _

ınt

HYPRE_SStructMaxwellSetup (HYPRE_SStructSolver solver, HYPRE_SStructMatrix A, HYPRE_SStructVector b, HYPRE_SStructVector x)

Prepare to solve the system. The coefficient data in b and x is ignored here, but information about the layout of the data may be used.

int

HYPRE_SStructMaxwellSolve (HYPRE_SStructSolver solver, HYPRE_SStructMatrix A, HYPRE_SStructVector b, HYPRE_SStructVector x)

Solve the system. Full coupling of the augmented system used throughout the multigrid hierarchy.

5.5.5

int

HYPRE_SStructMaxwellSolve2 (HYPRE_SStructSolver solver, HYPRE_SStructMatrix A, HYPRE_SStructVector b, HYPRE_SStructVector x)

Solve the system. Full coupling of the augmented system used only on the finest level, i.e., the node and edge multigrid cycles are coupled only on the finest level.

__ 5.5.6 __

int **HYPRE_SStructMaxwellSetGrad** (HYPRE_SStructSolver solver,
HYPRE_ParCSRMatrix T)

Sets the gradient operator in the Maxwell solver

_ 5.5.7 _

int

 $\label{lem:hypre_sstructMaxwellSetRfactors} \ (\mbox{HYPRE_SStructSolver solver, int rfactors} \ (\mbox{HYPRE_MAXDIM}))$

Sets the coarsening factor

int

HYPRE_SStructMaxwellPhysBdy (HYPRE_SStructGrid* grid_l, int num_levels, int rfactors[HYPRE_MAXDIM], int*** BdryRanks_ptr, int*** BdryRanksCnt_ptr)

Finds the physical boundary row ranks on all levels

__ 5.5.9 _____

 int

HYPRE_SStructMaxwellEliminateRowsCols (HYPRE_ParCSRMatrix parA, int nrows, int* rows)

Eliminates the rows and cols corresponding to the physical boundary in a parcsr matrix

5.5.10

int

 $\label{eq:hypre_structMaxwellZeroVector} \textbf{(HYPRE_ParVector b, int* rows, int nrows)}$

Zeros the rows corresponding to the physical boundary in a par vector

_ 5.5.11 _

int

HYPRE_SStructMaxwellSetSetConstantCoef (HYPRE_SStructSolver solver, int flag)

(Optional) Set the constant coefficient flag- Nedelec interpolation used

int **HYPRE_SStructMaxwellGrad** (HYPRE_SStructGrid grid,
HYPRE_ParCSRMatrix* T)

(Optional) Creates a gradient matrix from the grid. This presupposes a particular orientation of the edge elements.

5.5.13

HYPRE_SStructMaxwellSetTol (HYPRE_SStructSolver solver, HYPRE_Real tol)

(Optional) Set the convergence tolerance

___ 5.5.14 _____

HYPRE_SStructMaxwellSetMaxIter (HYPRE_SStructSolver solver, int max_iter)

(Optional) Set maximum number of iterations

5.5.15

HYPRE_SStructMaxwellSetRelChange (HYPRE_SStructSolver solver, int rel_change)

(Optional) Additionally require that the relative difference in successive iterates be small

int

 $\label{lem:hypre_struct} \textbf{HYPRE_SStructSolver solver}, \ int \\ num_pre_relax)$

(Optional) Set number of relaxation sweeps before coarse-grid correction

5.5.17

int

 $\label{lem:hypre_sstruct} \textbf{HYPRE_SStructSolver solver}, \\ \textbf{int } num_post_relax)$

(Optional) Set number of relaxation sweeps after coarse-grid correction

__ 5.5.18 _____

int

HYPRE_SStructMaxwellSetLogging (HYPRE_SStructSolver solver, int logging)

(Optional) Set the amount of logging to do

5.5.19

int

HYPRE_SStructMaxwellGetNumIterations (HYPRE_SStructSolver solver, int* num_iterations)

Return the number of iterations taken

int

$HYPRE_SStruct Maxwell GetFinal Relative Residual Norm$

 $({\it HYPRE_SStructSolver solver}, \, {\it HYPRE_Real* norm})$

Return the norm of the final relative residual

___ 5.6 _____

SStruct PCG Solver

Names

5.6.1	int HYPRE_SStructPCGCreate (MPI_Comm comm, HYPRE_SStructSolver* solver) Create a solver object	121
5.6.2	int	
	HYPRE_SStructPCGDestroy (HYPRE_SStructSolver solver)	
	Destroy a solver object.	122
5.6.3	int	
	HYPRE_SStructDiagScaleSetup (HYPRE_SStructSolver solver,	
	HYPRE_SStructMatrix A,	
	HYPRE_SStructVector y,	
	HYPRE_SStructVector x)	
	Setup routine for diagonal preconditioning	122
5.6.4	int	
	HYPRE_SStructDiagScale (HYPRE_SStructSolver solver,	
	HYPRE_SStructMatrix A,	
	HYPRE_SStructVector y,	
	HYPRE_SStructVector x)	
	Solve routine for diagonal preconditioning	122

These routines should be used in conjunction with the generic interface in PCG Solver.

5.6.1

int
HYPRE_SStructPCGCreate (MPI_Comm comm, HYPRE_SStructSolver*
solver)

Create a solver object

5.6.2

int HYPRE_SStructPCGDestroy (HYPRE_SStructSolver solver)

Destroy a solver object. An object should be explicitly destroyed using this destructor when the user's code no longer needs direct access to it. Once destroyed, the object must not be referenced again. Note that the object may not be deallocated at the completion of this call, since there may be internal package references to the object. The object will then be destroyed when all internal reference counts go to zero.

__ 5.6.3 ____

HYPRE_SStructDiagScaleSetup (HYPRE_SStructSolver solver, HYPRE_SStructMatrix A, HYPRE_SStructVector y, HYPRE_SStructVector x)

Setup routine for diagonal preconditioning

_ 5.6.4 _

int
HYPRE_SStructDiagScale (HYPRE_SStructSolver solver,
HYPRE_SStructMatrix A, HYPRE_SStructVector y, HYPRE_SStructVector x)

Solve routine for diagonal preconditioning

_ 5.7 _

SStruct GMRES Solver

Names

5.7.1 int

	HYPRE_SStructGMRESCreate (MPI_Comm comm,	
	HYPRE_SStructSolver* solver)	
	Create a solver object	123
5.7.2	int	
	HYPRE_SStructGMRESDestroy (HYPRE_SStructSolver solver)	
	Destroy a solver object.	123

These routines should be used in conjunction with the generic interface in GMRES Solver.

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HYPRE_SStructGMRESCreate (MPI_Comm comm, HYPRE_SStructSolver* solver)

Create a solver object

5.7.2

int HYPRE_SStructGMRESDestroy (HYPRE_SStructSolver solver)

Destroy a solver object. An object should be explicitly destroyed using this destructor when the user's code no longer needs direct access to it. Once destroyed, the object must not be referenced again. Note that the object may not be deallocated at the completion of this call, since there may be internal package references to the object. The object will then be destroyed when all internal reference counts go to zero.

5.8

SStruct FlexGMRES Solver

Names

5.8.1 int

 ${\bf HYPRE_SStructFlexGMRESCreate}~({\rm MPI_Comm}~{\rm comm},$

HYPRE_SStructSolver* solver)

5.8.2 int

These routines should be used in conjunction with the generic interface in FlexGMRES Solver.

__ 5.8.1 ___

int
HYPRE_SStructFlexGMRESCreate (MPI_Comm comm,
HYPRE_SStructSolver* solver)

Create a solver object

5.8.2

int HYPRE_SStructFlexGMRESDestroy (HYPRE_SStructSolver solver)

Destroy a solver object. An object should be explicitly destroyed using this destructor when the user's code no longer needs direct access to it. Once destroyed, the object must not be referenced again. Note that the object may not be deallocated at the completion of this call, since there may be internal package references to the object. The object will then be destroyed when all internal reference counts go to zero.

_ 5.9 _

SStruct LGMRES Solver

Names

5.9.1 int

HYPRE_SStructLGMRESCreate (MPI_Comm comm,

HYPRE_SStructSolver* solver)

5.9.2 int

 ${\bf HYPRE_SStructLGMRESDestroy} \ ({\bf HYPRE_SStructSolver} \ solver)$

Destroy a solver object. 125

These routines should be used in conjunction with the generic interface in LGMRES Solver.

5.9.1

HYPRE_SStructLGMRESCreate (MPI_Comm comm, HYPRE_SStructSolver* solver)

Create a solver object

__ 5.9.2 __

int HYPRE_SStructLGMRESDestroy (HYPRE_SStructSolver solver)

Destroy a solver object. An object should be explicitly destroyed using this destructor when the user's code no longer needs direct access to it. Once destroyed, the object must not be referenced again. Note that the object may not be deallocated at the completion of this call, since there may be internal package references to the object. The object will then be destroyed when all internal reference counts go to zero.

5.10

SStruct BiCGSTAB Solver

Names

These routines should be used in conjunction with the generic interface in BiCGSTAB Solver.

5.10.1

int
HYPRE_SStructBiCGSTABCreate (MPI_Comm comm,
HYPRE_SStructSolver* solver)

Create a solver object

__ 5.10.2 ____

int HYPRE_SStructBiCGSTABDestroy (HYPRE_SStructSolver solver)

Destroy a solver object. An object should be explicitly destroyed using this destructor when the user's code no longer needs direct access to it. Once destroyed, the object must not be referenced again. Note that the object may not be deallocated at the completion of this call, since there may be internal package references to the object. The object will then be destroyed when all internal reference counts go to zero.

5.11

SStruct LOBPCG Eigensolver

Names

These routines should be used in conjunction with the generic interface in LOBPCG Eigensolver.

_ 5.11.1 _

int HYPRE_SStructSetupInterpreter (mv_InterfaceInterpreter* i)

 $Load\ interface\ interpreter.\ Vector\ part\ loaded\ with\ hypre_SStructKrylov\ functions\ and\ multivector\ part\ loaded\ with\ mv_TempMultiVector\ functions.$

_ 5.11.2 _

int HYPRE_SStructSetupMatvec (HYPRE_MatvecFunctions* mv)

 ${\bf Load\ Matvec\ interpreter\ with\ hypre_SStructKrylov\ functions}$

6

ParCSR Solvers

Names		
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		199
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		216

These solvers use matrix/vector storage schemes that are taylored for general sparse matrix systems.

_ 6.1 _

ParCSR Solvers

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6.1.1	#define HYPRE_SOLVER_STRUCT	
	The solver object	129

__ 6.1.1 _____

#define HYPRE_SOLVER_STRUCT

The solver object

6.2

ParCSR BoomerAMG Solver and Preconditioner

Names		
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6.2.2	int HYPRE_BoomerAMGDestroy (HYPRE_Solver solver) Destroy a solver object	138
6.2.3	int HYPRE_BoomerAMGSetup (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x) Set up the BoomerAMG solver or preconditioner.	138
6.2.4	int HYPRE_BoomerAMGSolve (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x) Solve the system or apply AMG as a preconditioner.	139
6.2.5	int	

	HYPRE_BoomerAMGSolveT (HYPRE_Solver solver, HYPRE_ParCSRMatrix A,
	HYPRE_ParVector b, HYPRE_ParVector x) Solve the transpose system $A^Tx = b$ or apply AMG as a preconditioner to the transpose system.
C O C	
6.2.6	int HYPRE_BoomerAMGSetOldDefault (HYPRE_Solver solver) Recovers old default for coarsening and interpolation, ie Falgout coarsening and untruncated modified classical interpolation
6.2.7	int
	HYPRE_BoomerAMGGetResidual (HYPRE_Solver solver, HYPRE_ParVector* residual)
	Returns the residual
6.2.8	int HYPRE_BoomerAMGGetNumIterations (HYPRE_Solver solver,
	int* num_iterations) Returns the number of iterations taken
600	•
6.2.9	$\begin{array}{c} \operatorname{int} \\ \mathbf{HYPRE_BoomerAMGGetFinalRelativeResidualNorm} \ (\operatorname{HYPRE_Solver} \\ \operatorname{solver}, \end{array}$
	HYPRE_Real*
	rel_resid_norm)
	Returns the norm of the final relative residual
6.2.10	int HYPRE_BoomerAMGSetNumFunctions (HYPRE_Solver solver, int num_functions) (Optional) Sets the size of the system of PDEs, if using the systems version.
6.2.11	int HYPRE_BoomerAMGSetDofFunc (HYPRE_Solver solver, int* dof_func) (Optional) Sets the mapping that assigns the function to each variable, if using the systems version.
6.2.12	int HYPRE_BoomerAMGSetTol (HYPRE_Solver solver, HYPRE_Real tol) (Optional) Set the convergence tolerance, if BoomerAMG is used as a solver.
6.2.13	int HYPRE_BoomerAMGSetMaxIter (HYPRE_Solver solver, int max_iter) (Optional) Sets maximum number of iterations, if BoomerAMG is used as a solver.
6.2.14	int HYPRE_BoomerAMGSetMinIter (HYPRE_Solver solver, int min_iter) (Optional)
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6.2.15	int HYPRE_BoomerAMGSetMaxCoarseSize (HYPRE_Solver solver, int max_coarse_size)
	(Optional) Sets maximum size of coarsest grid
6.2.16	int

int min_coarse_size) (Optional) Sets minimum size of coarsest grid
HYPRE_BoomerAMGSetMaxLevels (HYPRE_Solver solver, int max_levels) (Optional) Sets maximum number of multigrid levels
int HYPRE_BoomerAMGSetStrongThreshold (HYPRE_Solver solver, HYPRE_Real strong_threshold) (Optional) Sets AMG strength threshold. (Optional) Sets AMG strength threshold. HYPRE_BoomerAMGSetSCommPkgSwitch (HYPRE_Solver solver, HYPRE_Real S_commpkg_switch) (Optional) Defines the largest strength threshold for which the strength matrix S uses the communication package of the operator A. 143 int HYPRE_BoomerAMGSetMaxRowSum (HYPRE_Solver solver, HYPRE_Real max_row_sum) (Optional) Sets a parameter to modify the definition of strength for diagonal dominant portions of the matrix. 143 int HYPRE_BoomerAMGSetCoarsenType (HYPRE_Solver solver,
HYPRE_BoomerAMGSetStrongThreshold (HYPRE_Solver solver, HYPRE_Real strong_threshold) (Optional) Sets AMG strength threshold
HYPRE_Real strong_threshold) (Optional) Sets AMG strength threshold
int HYPRE_BoomerAMGSetSCommPkgSwitch (HYPRE_Solver solver, HYPRE_Real S_commpkg_switch) (Optional) Defines the largest strength threshold for which the strength matrix S uses the communication package of the operator A
HYPRE_BoomerAMGSetSCommPkgSwitch (HYPRE_Solver solver, HYPRE_Real S_commpkg_switch) (Optional) Defines the largest strength threshold for which the strength matrix S uses the communication package of the operator A
HYPRE_Real S_commpkg_switch) (Optional) Defines the largest strength threshold for which the strength matrix S uses the communication package of the operator A
S_commpkg_switch) (Optional) Defines the largest strength threshold for which the strength matrix S uses the communication package of the operator A
(Optional) Defines the largest strength threshold for which the strength matrix S uses the communication package of the operator A
int HYPRE_BoomerAMGSetMaxRowSum (HYPRE_Solver solver, HYPRE_Real max_row_sum) (Optional) Sets a parameter to modify the definition of strength for diagonal dominant portions of the matrix. 143 int HYPRE_BoomerAMGSetCoarsenType (HYPRE_Solver solver,
HYPRE_BoomerAMGSetMaxRowSum (HYPRE_Solver solver, HYPRE_Real max_row_sum) (Optional) Sets a parameter to modify the definition of strength for diagonal dominant portions of the matrix. 143 int HYPRE_BoomerAMGSetCoarsenType (HYPRE_Solver solver,
HYPRE_Real max_row_sum) (Optional) Sets a parameter to modify the definition of strength for diagonal dominant portions of the matrix
(Optional) Sets a parameter to modify the definition of strength for diagonal dominant portions of the matrix
int HYPRE_BoomerAMGSetCoarsenType (HYPRE_Solver solver,
int HYPRE_BoomerAMGSetCoarsenType (HYPRE_Solver solver,
${\bf HYPRE_BoomerAMGSetCoarsenType}~({\tt HYPRE_Solver}~solver,$
int coarsen_type)
(Optional) Defines which parallel coarsening algorithm is used
int
· · · · · · · · · · · · · · · · · · ·
· · · · · · · · · · · · · · · · · · ·
nongalerkin_tol, int level)
(Optional) Defines the level specific non-Galerkin drop-tolerances for spar-
sifying coarse grid operators and thus reducing communication
int
V = (
int measure_tvpe)
Ÿ - /
(Optional) Defines whether local or global measures are used
(Optional) Defines whether local or global measures are used
(Optional) Defines whether local or global measures are used
(Optional) Defines whether local or global measures are used
int HYPRE_BoomerAMGSetAggNumLevels (HYPRE_Solver solver, int agg_num_levels) (Optional) Defines the number of levels of aggressive coarsening. 145
int HYPRE_BoomerAMGSetAggNumLevels (HYPRE_Solver solver,
int HYPRE_BoomerAMGSetAggNumLevels (HYPRE_Solver solver, int agg_num_levels) (Optional) Defines the number of levels of aggressive coarsening. 145
int HYPRE_BoomerAMGSetNonGalerkinTol (HYPRE_Solver solver, HYPRE_Real nongalerkin_tol) (Optional) Defines the non-Galerkin drop-tolerance for sparsifying coarse grid operators and thus reducing communication. int HYPRE_BoomerAMGSetLevelNonGalerkinTol (HYPRE_Solver solver, HYPRE_Real nongalerkin_tol, int level) (Optional) Defines the level specific non-Galerkin drop-tolerances for sparsifying coarse grid operators and thus reducing communication. 14

	HYPRE_BoomerAMGSetCGCIts (HYPRE_Solver solver, int its) (optional) Defines the number of pathes for CGC-coarsening	146
6.2.28	int HYPRE_BoomerAMGSetNodal (HYPRE_Solver solver, int nodal)	
	(Optional) Sets whether to use the nodal systems coarsening	146
6.2.29	int HYPRE_BoomerAMGSetNodalDiag (HYPRE_Solver solver, int nodal_diag) (Optional) Sets whether to give special treatment to diagonal elements in the nodal systems version.	146
6.2.30	int HYPRE_BoomerAMGSetInterpType (HYPRE_Solver solver,	
	int interp_type) (Optional) Defines which parallel interpolation operator is used	146
6.2.31	int HYPRE_BoomerAMGSetTruncFactor (HYPRE_Solver solver, HYPRE_Real trunc_factor)	
	(Optional) Defines a truncation factor for the interpolation	147
6.2.32	int HYPRE_BoomerAMGSetPMaxElmts (HYPRE_Solver solver, int P_max_elmts)	
	(Optional) Defines the maximal number of elements per row for the interpolation.	147
6.2.33	int HYPRE_BoomerAMGSetSepWeight (HYPRE_Solver solver, int sep_weight) (Optional) Defines whether separation of weights is used when defining strength for standard interpolation or multipass interpolation	147
6.2.34	int HYPRE_BoomerAMGSetAggInterpType (HYPRE_Solver solver, int agg_interp_type)	
	(Optional) Defines the interpolation used on levels of aggressive coarsening The default is 4, ie.	148
6.2.35	int HYPRE_BoomerAMGSetAggTruncFactor (HYPRE_Solver solver, HYPRE_Real agg_trunc_factor)	
	(Optional) Defines the truncation factor for the interpolation used for aggressive coarsening.	148
6.2.36	int HYPRE_BoomerAMGSetAggP12TruncFactor (HYPRE_Solver solver,	
	HYPRE_Real agg_P12_trunc_factor)	
	(Optional) Defines the truncation factor for the matrices P1 and P2 which are used to build 2-stage interpolation.	148
6.2.37	int HYPRE_BoomerAMGSetAggPMaxElmts (HYPRE_Solver solver, int agg_P_max_elmts)	
	(Optional) Defines the maximal number of elements per row for the interpolation used for aggressive coarsening.	149
6.2.38	int	

	HYPRE_BoomerAMGSetAggP12MaxElmts (HYPRE_Solver solver,	
	int agg_P12_max_elmts)	
	(Optional) Defines the maximal number of elements per row for the matrices P1 and P2 which are used to build 2-stage interpolation	149
6.2.39	int	
	HYPRE_BoomerAMGSetInterpVectors (HYPRE_Solver solver,	
	int num_vectors,	
	HYPRE_ParVector* interp_vectors	
	(Optional) Allows the user to incorporate additional vectors into the inter-	
	polation for systems AMG, eg	149
6.2.40	int	
	HYPRE_BoomerAMGSetInterpVecVariant (HYPRE_Solver solver, int var)	
	(Optional) Defines the interpolation variant	
	$used \hspace{1cm} for \hspace{1cm} HYPRE_BoomerAMGSetInterp Vectors:$	
	1 GM approach 1	
	2 GM approach 2 (to be preferred over 1)	149
	$3 \mid LN \ approach$	
6.2.41	int	
	HYPRE_BoomerAMGSetInterpVecQMax (HYPRE_Solver solver,	
	int q_max)	
	(Optional) Defines the maximal elements per row for Q , the additional columns added to the original interpolation matrix P , to reduce complexity.	
	columns daded to the original interpolation matrix F, to reduce complexity.	150
		100
6.2.42	int	
	HYPRE_BoomerAMGSetInterpVecAbsQTrunc (HYPRE_Solver solver,	
	HYPRE_Real q_trunc)	
	(Optional) Defines a truncation factor for Q, the additional columns added	150
	to the original interpolation matrix P , to reduce complexity	150
6.2.43	int	
	HYPRE_BoomerAMGSetGSMG (HYPRE_Solver solver, int gsmg)	
	(Optional) Specifies the use of GSMG - geometrically smooth coarsening and	
	interpolation.	150
6.2.44	int	
	HYPRE_BoomerAMGSetNumSamples (HYPRE_Solver solver,	
	int num_samples)	
	(Optional) Defines the number of sample vectors used in GSMG or LS in-	
	terpolation	150
6.2.45	int	
	HYPRE_BoomerAMGSetCycleType (HYPRE_Solver solver, int cycle_type)	
	(Optional) Defines the type of cycle	151
6.2.46	int	
0.4.40	HYPRE_BoomerAMGSetAdditive (HYPRE_Solver solver, int addlvl)	
	(Optional) Defines use of an additive $V(1, 1)$ -cycle using the classical addi-	
	tive method starting at level 'addlvl'.	151
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6.2.57	(Optional) Sets the number of sweeps at a specified cycle	1			

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0.2.00	HYPRE_BoomerAMGSetRelaxType (HYPRE_Solver solver, int relax_type) (Optional) Defines the smoother to be used
6.2.59	int
	HYPRE_BoomerAMGSetCycleRelaxType (HYPRE_Solver solver, int relax_type, int k)
	(Optional) Defines the smoother at a given cycle
6.2.60	int
0.2.00	HYPRE_BoomerAMGSetRelaxOrder (HYPRE_Solver solver, int relax_order)
	(Optional) Defines in which order the points are relaxed
6.2.61	int
0.2.01	HYPRE_BoomerAMGSetRelaxWt (HYPRE_Solver solver, HYPRE_Real relax_weight)
	(Optional) Defines the relaxation weight for smoothed Jacobi and hybrid SOR on all levels.
6.2.62	int
0.2.02	HYPRE_BoomerAMGSetLevelRelaxWt (HYPRE_Solver solver, HYPRE_Real relax_weight, int level)
	(Optional) Defines the relaxation weight for smoothed Jacobi and hybrid SOR on the user defined level.
6.2.63	int
0.2.00	HYPRE_BoomerAMGSetOmega (HYPRE_Solver solver,
	HYPRE_Real* omega) (Ontional) Defines the outer relaxation weight for hybrid SOP
	(Optional) Defines the outer relaxation weight for hybrid SOR
6.2.64	int HYPRE_BoomerAMGSetOuterWt (HYPRE_Solver solver, HYPRE_Real omega)
	(Optional) Defines the outer relaxation weight for hybrid SOR and SSOR
0.0.05	
6.2.65	int HYPRE_BoomerAMGSetLevelOuterWt (HYPRE_Solver solver,
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6.2.67	int
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	(Optional) Fraction of the spectrum to use for the Chebyshev smoother
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	int** grid_relax_type_ptr,	
	int*** grid_relax_points_ptr,	
	int coarsen_type,	
	HYPRE_Real**	
	$relax_weights_ptr,$	
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6.2.91	int	

 $\begin{array}{c} \textbf{HYPRE_BoomerAMGSetKeepTranspose} \text{ (HYPRE_Solver solver,} \\ \text{int keepTranspose)} \end{array}$

(Optional) If set to 1, the local interpolation transposes will be saved to use more efficient matvecs instead of matvecTs

Parallel unstructured algebraic multigrid solver and preconditioner

 $_{-}$ 6.2.1 $_{-}$

int HYPRE_BoomerAMGCreate (HYPRE_Solver* solver)

Create a solver object

6.2.2

int HYPRE_BoomerAMGDestroy (HYPRE_Solver solver)

Destroy a solver object

_ 6.2.3 _

int

HYPRE_BoomerAMGSetup (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x)

Set up the BoomerAMG solver or preconditioner. If used as a preconditioner, this function should be passed to the iterative solver SetPrecond function.

Parameters:

solver [IN] object to be set up.

A [IN] ParCSR matrix used to construct the solver/preconditioner.

b Ignored by this function.

x Ignored by this function.

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6.2.4

int **HYPRE_BoomerAMGSolve** (HYPRE_Solver solver, HYPRE_ParCSRMatrix A,
HYPRE_ParVector b, HYPRE_ParVector x)

Solve the system or apply AMG as a preconditioner. If used as a preconditioner, this function should be passed to the iterative solver SetPrecond function.

Parameters:	solver	[IN] solver or preconditioner object to be applied.
	Α	[IN] ParCSR matrix, matrix of the linear system to be
		solved
	b	[IN] right hand side of the linear system to be solved
	x	[OUT] approximated solution of the linear system to
		be solved

6.2.5

int **HYPRE_BoomerAMGSolveT** (HYPRE_Solver solver, HYPRE_ParCSRMatrix
A, HYPRE_ParVector b, HYPRE_ParVector x)

Solve the transpose system $A^Tx=b$ or apply AMG as a preconditioner to the transpose system . Note that this function should only be used when preconditioning CGNR with BoomerAMG. It can only be used with Jacobi smoothing (relax_type 0 or 7) and without CF smoothing, i.e relax_order needs to be set to 0. If used as a preconditioner, this function should be passed to the iterative solver SetPrecond function.

Parameters:	solver	[IN] solver or preconditioner object to be applied.
	A	[IN] ParCSR matrix
	b	[IN] right hand side of the linear system to be solved
	x	[OUT] approximated solution of the linear system to
		be solved

6.2.6

int HYPRE_BoomerAMGSetOldDefault (HYPRE_Solver solver)

Recovers old default for coarsening and interpolation, ie Falgout coarsening and untruncated modified classical interpolation. This option might be preferred for 2 dimensional problems.

6.2.7

int

HYPRE_BoomerAMGGetResidual (HYPRE_Solver solver, HYPRE_ParVector* residual)

Returns the residual

6.2.8

int

 $\label{eq:hypre_bound} \begin{aligned} \mathbf{HYPRE_BoomerAMGGetNumIterations} & \text{ (HYPRE_Solver solver, int* } \\ \text{num_iterations)} \end{aligned}$

Returns the number of iterations taken

__ 6.2.9 ____

int

HYPRE_BoomerAMGGetFinalRelativeResidualNorm (HYPRE_Solver solver, HYPRE_Real* rel_resid_norm)

Returns the norm of the final relative residual

6.2.10

int

HYPRE_BoomerAMGSetNumFunctions (HYPRE_Solver solver, int num_functions)

(Optional) Sets the size of the system of PDEs, if using the systems version. The default is 1, i.e. a scalar system.

6211

int HYPRE_BoomerAMGSetDofFunc (HYPRE_Solver solver, int* dof_func)

(Optional) Sets the mapping that assigns the function to each variable, if using the systems version. If no assignment is made and the number of functions is k > 1, the mapping generated is (0,1,...,k-1,0,1,...,k-1,...).

6.2.12

int HYPRE_BoomerAMGSetTol (HYPRE_Solver solver, HYPRE_Real tol)

(Optional) Set the convergence tolerance, if BoomerAMG is used as a solver. If it is used as a preconditioner, it should be set to 0. The default is 1.e-7.

_ 6.2.13 _

int HYPRE_BoomerAMGSetMaxIter (HYPRE_Solver solver, int max_iter)

(Optional) Sets maximum number of iterations, if BoomerAMG is used as a solver. If it is used as a preconditioner, it should be set to 1. The default is 20.

6.2.14

int HYPRE_BoomerAMGSetMinIter (HYPRE_Solver solver, int min_iter)

(Optional)

6.2.15

int

 $\label{lem:hypre_bound} \mathbf{HYPRE_BoomerAMGSetMaxCoarseSize} \ (\mathbf{HYPRE_Solver} \ solver, \ int \\ \mathbf{max_coarse_size})$

(Optional) Sets maximum size of coarsest grid. The default is 9.

___ 6.2.16 _____

int

 $\label{lem:hypre_bound} \textbf{HYPRE_BoomerAMGSetMinCoarseSize} \ (\textbf{HYPRE_Solver solver}, \ \textbf{int min_coarse_size})$

(Optional) Sets minimum size of coarsest grid. The default is 1.

__ 6.2.17 _____

int

HYPRE_BoomerAMGSetMaxLevels (HYPRE_Solver solver, int max_levels)

(Optional) Sets maximum number of multigrid levels. The default is 25.

__ 6.2.18 _____

int

HYPRE_BoomerAMGSetStrongThreshold (HYPRE_Solver solver, HYPRE_Real strong_threshold)

(Optional) Sets AMG strength threshold. The default is 0.25. For 2d Laplace operators, 0.25 is a good value, for 3d Laplace operators, 0.5 or 0.6 is a better value. For elasticity problems, a large strength threshold, such as 0.9, is often better.

6.2.19

int

HYPRE_BoomerAMGSetSCommPkgSwitch (HYPRE_Solver solver, HYPRE_Real S_commpkg_switch)

(Optional) Defines the largest strength threshold for which the strength matrix S uses the communication package of the operator A. If the strength threshold is larger than this values, a communication package is generated for S. This can save memory and decrease the amount of data that needs to be communicated, if S is substantially sparser than A. The default is 1.0.

_ 6.2.20 __

int **HYPRE_BoomerAMGSetMaxRowSum** (HYPRE_Solver solver, HYPRE_Real max_row_sum)

(Optional) Sets a parameter to modify the definition of strength for diagonal dominant portions of the matrix. The default is 0.9. If max_row_sum is 1, no checking for diagonally dominant rows is performed.

6.2.21

int **HYPRE_BoomerAMGSetCoarsenType** (HYPRE_Solver solver, int coarsen_type)

(Optional) Defines which parallel coarsening algorithm is used. There are the following options for coarsen_type:

- 0 CLJP-coarsening (a parallel coarsening algorithm using independent sets.
- 1 | classical Ruge-Stueben coarsening on each processor, no boundary treatment (not recommended!)
- 3 classical Ruge-Stueben coarsening on each processor, followed by a third pass, which adds coarse points on the boundaries
- Falgout coarsening (uses 1 first, followed by CLJP using the interior coarse points generated by 1 as its first independent set)
- 7 | CLJP-coarsening (using a fixed random vector, for debugging purposes only)
- 8 PMIS-coarsening (a parallel coarsening algorithm using independent sets, generating lower complexities than CLJP, might also lead to slower convergence)
- 9 PMIS-coarsening (using a fixed random vector, for debugging purposes only)
- HMIS-coarsening (uses one pass Ruge-Stueben on each processor independently, followed by PMIS using the interior C-points generated as its first independent set)
- 11 one-pass Ruge-Stueben coarsening on each processor, no boundary treatment (not recommended!)
- 21 | CGC coarsening by M. Griebel, B. Metsch and A. Schweitzer
- 22 CGC-E coarsening by M. Griebel, B. Metsch and A.Schweitzer

The default is 6.

6.2.22

int

HYPRE_BoomerAMGSetNonGalerkinTol (HYPRE_Solver solver, HYPRE_Real nongalerkin_tol)

(Optional) Defines the non-Galerkin drop-tolerance for sparsifying coarse grid operators and thus reducing communication. Value specified here is set on all levels. This routine should be used before HYPRE_BoomerAMGSetLevelNonGalerkinTol, which then can be used to change individual levels if desired

6.2.23 $_{-}$

int

HYPRE_BoomerAMGSetLevelNonGalerkinTol (HYPRE_Solver solver, HYPRE_Real nongalerkin_tol, int level)

(Optional) Defines the level specific non-Galerkin drop-tolerances for sparsifying coarse grid operators and thus reducing communication. A drop-tolerance of 0.0 means to skip doing non-Galerkin on that level. The maximum drop tolerance for a level is 1.0, although much smaller values such as 0.03 or 0.01 are recommended.

Note that if the user wants to set a specific tolerance on all levels, HYPRE_BooemrAMGSetNonGalerkinTol should be used. Individual levels can then be changed using this routine.

In general, it is safer to drop more aggressively on coarser levels. For instance, one could use 0.0 on the finest level, 0.01 on the second level and then using 0.05 on all remaining levels. The best way to achieve this is to set 0.05 on all levels with HYPRE_BoomerAMGSetNonGalerkinTol and then change the tolerance on level 0 to 0.0 and the tolerance on level 1 to 0.01 with HYPRE_BoomerAMGSetLevelNonGalerkinTol. Like many AMG parameters, these drop tolerances can be tuned. It is also common to delay the start of the non-Galerkin process further to a later level than level 1.

Parameters:

solver

[IN] solver or preconditioner object to be applied.

 $nongalerkin_tol$

[IN] level specific drop tolerance

level

[IN] level on which drop tolerance is used

6.2.24

int

HYPRE_BoomerAMGSetMeasureType (HYPRE_Solver solver, int measure_type)

(Optional) Defines whether local or global measures are used

6.2.25 $_{-}$

HYPRE_BoomerAMGSetAggNumLevels (HYPRE_Solver solver, int agg_num_levels)

(Optional) Defines the number of levels of aggressive coarsening. The default is 0, i.e. no aggressive coarsening.

6.2.26

int

HYPRE_BoomerAMGSetNumPaths (HYPRE_Solver solver, int num_paths)

(Optional) Defines the degree of aggressive coarsening. The default is 1. Larger numbers lead to less aggressive coarsening.

int HYPRE_BoomerAMGSetCGCIts (HYPRE_Solver solver, int its)

(optional) Defines the number of pathes for CGC-coarsening

_ 6.2.28 _

int HYPRE_BoomerAMGSetNodal (HYPRE_Solver solver, int nodal)

(Optional) Sets whether to use the nodal systems coarsening. Should be used for linear systems generated from systems of PDEs. The default is 0 (unknown-based coarsening, only coarsens within same function). For the remaining options a nodal matrix is generated by applying a norm to the nodal blocks and applying

2

Frobenius norm
sum of absolute values of elements in each block

- 3 largest element in each block (not absolute value)
- 4 row-sum norm
- 6 sum of all values in each block

 $_$ 6.2.29 $_$

the coarsening algorithm to this matrix.

int

HYPRE_BoomerAMGSetNodalDiag (HYPRE_Solver solver, int nodal_diag)

(Optional) Sets whether to give special treatment to diagonal elements in the nodal systems version. The default is 0. If set to 1, the diagonal entry is set to the negative sum of all off diagonal entries. If set to 2, the signs of all diagonal entries are inverted.

6.2.30

1111U

HYPRE_BoomerAMGSetInterpType (HYPRE_Solver solver, int interp_type)

(Optional) Defines which parallel interpolation operator is used. There are the following options for interp_type:

- 0 classical modified interpolation
- 1 LS interpolation (for use with GSMG)
- 2 classical modified interpolation for hyperbolic PDEs
- 3 direct interpolation (with separation of weights)
- 4 multipass interpolation
- 5 | multipass interpolation (with separation of weights)
- 6 extended+i interpolation
- 7 extended+i (if no common C neighbor) interpolation
- 8 standard interpolation
- 9 standard interpolation (with separation of weights)
- 10 classical block interpolation (for use with nodal systems version only)
- classical block interpolation (for use with nodal systems version only) with diagonalized diagonal blocks
- 12 | FF interpolation
- 13 | FF1 interpolation
- 14 extended interpolation

The default is 0.

6.2.31

int

HYPRE_BoomerAMGSetTruncFactor (HYPRE_Solver solver, HYPRE_Real trunc_factor)

(Optional) Defines a truncation factor for the interpolation. The default is 0.

_ 6.2.32 _

ınt

HYPRE_BoomerAMGSetPMaxElmts (HYPRE_Solver solver, int P_max_elmts)

(Optional) Defines the maximal number of elements per row for the interpolation. The default is 0.

6.2.33

ınt

HYPRE_BoomerAMGSetSepWeight (HYPRE_Solver solver, int sep_weight)

(Optional) Defines whether separation of weights is used when defining strength for standard interpolation or multipass interpolation. Default: 0, i.e. no separation of weights used.

6.2.34

int **HYPRE_BoomerAMGSetAggInterpType** (HYPRE_Solver solver, int agg_interp_type)

(Optional) Defines the interpolation used on levels of aggressive coarsening The default is 4, ie. multipass interpolation. The following options exist:

- 2-stage extended+i interpolation
- 2 | 2-stage standard interpolation
- 3 2-stage extended interpolation
- 4 | multipass interpolation

6.2.35

HYPRE_BoomerAMGSetAggTruncFactor (HYPRE_Solver solver, HYPRE_Real agg_trunc_factor)

(Optional) Defines the truncation factor for the interpolation used for aggressive coarsening. The default is 0.

6.2.36

HYPRE_BoomerAMGSetAggP12TruncFactor (HYPRE_Solver solver, HYPRE_Real agg_P12_trunc_factor)

(Optional) Defines the truncation factor for the matrices P1 and P2 which are used to build 2-stage interpolation. The default is 0.

int **HYPRE_BoomerAMGSetAggPMaxElmts** (HYPRE_Solver solver, int agg_P_max_elmts)

(Optional) Defines the maximal number of elements per row for the interpolation used for aggressive coarsening. The default is 0.

_ 6.2.38 _

HYPRE_BoomerAMGSetAggP12MaxElmts (HYPRE_Solver solver, int agg_P12_max_elmts)

(Optional) Defines the maximal number of elements per row for the matrices P1 and P2 which are used to build 2-stage interpolation. The default is 0.

6.2.39

int

HYPRE_BoomerAMGSetInterpVectors (HYPRE_Solver solver, int num_vectors, HYPRE_ParVector* interp_vectors)

(Optional) Allows the user to incorporate additional vectors into the interpolation for systems AMG, eg. rigid body modes for linear elasticity problems. This can only be used in context with nodal coarsening and still requires the user to choose an interpolation.

6.2.40

HYPRE_BoomerAMGSetInterpVecVariant (HYPRE_Solver solver, int var)

(Optional) Defines the interpolation variant used for HYPRE_BoomerAMGSetInterpVectors:

- 1 GM approach 1
- 2 GM approach 2 (to be preferred over 1)
- 3 | LN approach

int HYPRE_BoomerAMGSetInterpVecQMax (HYPRE_Solver solver, int q_max)

(Optional) Defines the maximal elements per row for Q, the additional columns added to the original interpolation matrix P, to reduce complexity. The default is no truncation.

__ 6.2.42 _____

int **HYPRE_BoomerAMGSetInterpVecAbsQTrunc** (HYPRE_Solver solver, HYPRE_Real q_trunc)

(Optional) Defines a truncation factor for Q, the additional columns added to the original interpolation matrix P, to reduce complexity. The default is no truncation.

 $_{-}$ 6.2.43 $_{-}$

int HYPRE_BoomerAMGSetGSMG (HYPRE_Solver solver, int gsmg)

(Optional) Specifies the use of GSMG - geometrically smooth coarsening and interpolation. Currently any nonzero value for gsmg will lead to the use of GSMG. The default is 0, i.e. (GSMG is not used)

6.2.44

HYPRE_BoomerAMGSetNumSamples (HYPRE_Solver solver, int num_samples)

(Optional) Defines the number of sample vectors used in GSMG or LS interpolation

HYPRE_BoomerAMGSetCycleType (HYPRE_Solver solver, int cycle_type)

(Optional) Defines the type of cycle. For a V-cycle, set cycle_type to 1, for a W-cycle set cycle_type to 2. The default is 1.

6.2.46

int HYPRE_BoomerAMGSetAdditive (HYPRE_Solver solver, int addlvl)

(Optional) Defines use of an additive V(1,1)-cycle using the classical additive method starting at level 'addlvl'. The multiplicative approach is used on levels 0, ...'addlvl+1'. 'addlvl' needs to be > -1 for this to have an effect. Can only be used with weighted Jacobi and l1-Jacobi(default).

Can only be used when AMG is used as a preconditioner!!!

6.2.47

int HYPRE_BoomerAMGSetMultAdditive (HYPRE_Solver solver, int addlvl)

(Optional) Defines use of an additive V(1,1)-cycle using the mult-additive method starting at level 'addlvl'. The multiplicative approach is used on levels 0, ...'addlvl+1'. 'addlvl' needs to be > -1 for this to have an effect. Can only be used with weighted Jacobi and l1-Jacobi(default).

Can only be used when AMG is used as a preconditioner!!!

6.2.48

int HYPRE_BoomerAMGSetSimple (HYPRE_Solver solver, int addlvl)

(Optional) Defines use of an additive V(1,1)-cycle using the simplified mult-additive method starting at level 'addlvl'. The multiplicative approach is used on levels 0, ...'addlvl+1'. 'addlvl' needs to be > -1 for this to have an effect. Can only be used with weighted Jacobi and l1-Jacobi(default).

Can only be used when AMG is used as a preconditioner!!!

_ 6.2.49 _

HYPRE_BoomerAMGSetMultAddTruncFactor (HYPRE_Solver solver, HYPRE_Real add_trunc_factor)

(Optional) Defines the truncation factor for the smoothed interpolation used for mult-additive or simple method. The default is 0.

6.2.50

HYPRE_BoomerAMGSetMultAddPMaxElmts (HYPRE_Solver solver, int add_P_max_elmts)

(Optional) Defines the maximal number of elements per row for the smoothed interpolation used for mult-additive or simple method. The default is 0.

 $_$ 6.2.51 $_$

HYPRE_BoomerAMGSetAddRelaxType (HYPRE_Solver solver, int add_rlx_type)

(Optional) Defines the relaxation type used in the (mult)additive cycle portion (also affects simple method) The default is 18 (L1-Jacobi). Currently the only other option allowed is 0 (Jacobi) which should be used in combination with HYPRE_BoomerAMGSetAddRelaxWt.

HYPRE_BoomerAMGSetAddRelaxWt (HYPRE_Solver solver, HYPRE_Real add_rlx_wt)

(Optional) Defines the relaxation weight used for Jacobi within the (mult)additive or simple cycle portion. The default is 1. The weight only affects the Jacobi method, and has no effect on L1-Jacobi

6.2.53

int **HYPRE_BoomerAMGSetSeqThreshold** (HYPRE_Solver solver, int seq_threshold)

(Optional) Sets maximal size for agglomeration or redundant coarse grid solve. When the system is smaller than this threshold, sequential AMG is used on process 0 or on all remaining active processes (if redundant = 1).

6.2.54

int **HYPRE_BoomerAMGSetRedundant** (HYPRE_Solver solver, int redundant)

(Optional) operates switch for redundancy. Needs to be used with HYPRE_BoomerAMGSetSeqThreshold. Default is 0, i.e. no redundancy.

6.2.55

int

HYPRE_BoomerAMGSetNumSweeps (HYPRE_Solver solver, int num_sweeps)

(Optional) Sets the number of sweeps. On the finest level, the up and the down cycle the number of sweeps are set to num_sweeps and on the coarsest level to 1. The default is 1.

int HYPRE_BoomerAMGSetCycleNumSweeps (HYPRE_Solver solver, int num_sweeps, int k)

(Optional) Sets the number of sweeps at a specified cycle. There are the following options for k:

the down cycle	if k=1
the up cycle	if $k=2$
the coarsest level	if k=3.

 $_$ 6.2.57 $_$

HYPRE_BoomerAMGSetGridRelaxType (HYPRE_Solver solver, int* grid_relax_type)

(Optional) Defines which smoother is used on the fine and coarse grid, the up and down cycle.

Note: This routine will be phased out!!!! Use HYPRE_BoomerAMGSetRelaxType or HYPRE_BoomerAMGSetCycleRelaxType instead.

6.2.58

HYPRE_BoomerAMGSetRelaxType (HYPRE_Solver solver, int relax_type)

(Optional) Defines the smoother to be used. It uses the given smoother on the fine grid, the up and the down cycle and sets the solver on the coarsest level to Gaussian elimination (9). The default is Gauss-Seidel (3).

There are the following options for relax_type:

0	Jaco	bi
---	------	----

- 1 Gauss-Seidel, sequential (very slow!)
- 2 Gauss-Seidel, interior points in parallel, boundary sequential (slow!)
- 3 hybrid Gauss-Seidel or SOR, forward solve
- 4 hybrid Gauss-Seidel or SOR, backward solve
- 5 | hybrid chaotic Gauss-Seidel (works only with OpenMP)
- 6 hybrid symmetric Gauss-Seidel or SSOR
- 8 ℓ_1 -scaled hybrid symmetric Gauss-Seidel
- 9 Gaussian elimination (only on coarsest level)
- 15 | CG (warning not a fixed smoother may require FGMRES)
- 16 Chebyshev
- 17 FCF-Jacobi
- 18 ℓ_1 -scaled jacobi

int

 $\label{eq:hypre_bound} \mathbf{HYPRE_BoomerAMGSetCycleRelaxType} \ (\mathbf{HYPRE_Solver} \ \mathbf{solver}, \ \mathbf{int} \ \mathbf{relax_type}, \ \mathbf{int} \ \mathbf{k})$

(Optional) Defines the smoother at a given cycle. For options of relax_type see description of HYPRE_BoomerAMGSetRelaxType). Options for k are

the down cycle	if k=1
the up cycle	if $k=2$
the coarsest level	if k=3.

6.2.60 $\overline{}$

int

HYPRE_BoomerAMGSetRelaxOrder (HYPRE_Solver solver, int relax_order)

(Optional) Defines in which order the points are relaxed. There are the following options for relax_order:

- 0 the points are relaxed in natural or lexicographic order on each processor
- 1 CF-relaxation is used, i.e on the fine grid and the down cycle the coarse points are relaxed first, followed by the fine points; on the up cycle the F-points are relaxed first, followed by the C-points. On the coarsest level, if an iterative scheme is used, the points are relaxed in lexicographic order.

The default is 1 (CF-relaxation).

int

 $\label{eq:hypre_bound} \mathbf{HYPRE_BoomerAMGSetRelaxWt} \ (\mathbf{HYPRE_Solver} \ \mathbf{solver}, \ \mathbf{HYPRE_Real} \ \mathbf{relax_weight})$

(Optional) Defines the relaxation weight for smoothed Jacobi and hybrid SOR on all levels.

re	elax_weight > 0	this assigns the given relaxation weight on all levels
re	$elax_weight = 0$	the weight is determined on each level with the estimate $\frac{3}{4\ D^{-1/2}AD^{-1/2}\ }$,
		where D is the diagonal matrix of A (this should only be used with Jacobi)
re	$elax_weight = -k$	the relaxation weight is determined with at most k CG steps on each level
		this should only be used for symmetric positive definite problems)

The default is 1.

$_{-}$ 6.2.62 $_{-}$

HYPRE_BoomerAMGSetLevelRelaxWt (HYPRE_Solver solver, HYPRE_Real relax_weight, int level)

(Optional) Defines the relaxation weight for smoothed Jacobi and hybrid SOR on the user defined level. Note that the finest level is denoted 0, the next coarser level 1, etc. For nonpositive relax_weight, the parameter is determined on the given level as described for HYPRE_BoomerAMGSetRelaxWt. The default is 1

6.2.63 _____

int

HYPRE_BoomerAMGSetOmega (HYPRE_Solver solver, HYPRE_Real* omega)

(Optional) Defines the outer relaxation weight for hybrid SOR. Note: This routine will be phased out!!!! Use HYPRE_BoomerAMGSetOuterWt or HYPRE_BoomerAMGSetLevelOuterWt instead.

HYPRE_BoomerAMGSetOuterWt (HYPRE_Solver solver, HYPRE_Real omega)

(Optional) Defines the outer relaxation weight for hybrid SOR and SSOR on all levels.

omega > 0	this assigns the same outer relaxation weight omega on each level
omega = -k	an outer relaxation weight is determined with at most k CG steps on each level
	(this only makes sense for symmetric positive definite problems and smoothers, e.g. SSOR)

The default is 1.

 $_{-}$ 6.2.65 $_{-}$

int **HYPRE_BoomerAMGSetLevelOuterWt** (HYPRE_Solver solver, HYPRE_Real omega, int level)

(Optional) Defines the outer relaxation weight for hybrid SOR or SSOR on the user defined level. Note that the finest level is denoted 0, the next coarser level 1, etc. For nonpositive omega, the parameter is determined on the given level as described for HYPRE_BoomerAMGSetOuterWt. The default is 1.

_ 6.2.66 ____

int HYPRE_BoomerAMGSetChebyOrder (HYPRE_Solver solver, int order)

(Optional) Defines the Order for Chebyshev smoother. The default is 2 (valid options are 1-4).

_ 6.2.67 ___

HYPRE_BoomerAMGSetChebyFraction (HYPRE_Solver solver, HYPRE_Real ratio)

(Optional) Fraction of the spectrum to use for the Chebyshev smoother. The default is .3 (i.e., damp on upper 30% of the spectrum).

_ 6.2.68 _

int **HYPRE_BoomerAMGSetSmoothType** (HYPRE_Solver solver, int smooth_type)

(Optional) Enables the use of more complex smoothers. The following options exist for smooth_type:

value	smoother	routines needed to set smoother parameters
6	Schwarz smoothers	HYPRE_BoomerAMGSetDomainType, HYPRE_BoomerAMGSetOverlap,
		HYPRE_BoomerAMGSetVariant, HYPRE_BoomerAMGSetSchwarzRlxWeight
7	Pilut	HYPRE_BoomerAMGSetDropTol, HYPRE_BoomerAMGSetMaxNzPerRow
8	ParaSails	HYPRE_BoomerAMGSetSym, HYPRE_BoomerAMGSetLevel,
		HYPRE_BoomerAMGSetFilter, HYPRE_BoomerAMGSetThreshold
9	Euclid	HYPRE_BoomerAMGSetEuclidFile

The default is 6. Also, if no smoother parameters are set via the routines mentioned in the table above, default values are used.

6.2.69

HYPRE_BoomerAMGSetSmoothNumLevels (HYPRE_Solver solver, int smooth_num_levels)

(Optional) Sets the number of levels for more complex smoothers. The smoothers, as defined by HYPRE_BoomerAMGSetSmoothType, will be used on level 0 (the finest level) through level smooth_num_levels-1. The default is 0, i.e. no complex smoothers are used.

6.2.70

HYPRE_BoomerAMGSetSmoothNumSweeps (HYPRE_Solver solver, int smooth_num_sweeps)

(Optional) Sets the number of sweeps for more complex smoothers. The default is 1.

int HYPRE_BoomerAMGSetVariant (HYPRE_Solver solver, int variant)

(Optional) Defines which variant of the Schwarz method is used. The following options exist for variant:

- 0 hybrid multiplicative Schwarz method (no overlap across processor boundaries)
- 1 | hybrid additive Schwarz method (no overlap across processor boundaries)
- 2 | additive Schwarz method
- 3 | hybrid multiplicative Schwarz method (with overlap across processor boundaries)

The default is 0.

6.2.72

int HYPRE_BoomerAMGSetOverlap (HYPRE_Solver solver, int overlap)

(Optional) Defines the overlap for the Schwarz method. The following options exist for overlap:

- 0 no overlap
- minimal overlap (default)
- 2 | overlap generated by including all neighbors of domain boundaries

6.2.73 _

HYPRE_BoomerAMGSetDomainType (HYPRE_Solver solver, int domain_type)

(Optional) Defines the type of domain used for the Schwarz method. The following options exist for domain_type:

- 0 each point is a domain
- 1 | each node is a domain (only of interest in "systems" AMG)
- 2 each domain is generated by agglomeration (default)

HYPRE_BoomerAMGSetSchwarzRlxWeight (HYPRE_Solver solver, HYPRE_Real schwarz_rlx_weight)

(Optional) Defines a smoothing parameter for the additive Schwarz method

___ 6.2.75 ____

int

 $\label{lem:hypre_bound} \mbox{\bf HYPRE_BoomerAMGSetSchwarzUseNonSymm (HYPRE_Solver solver, intuse_nonsymm)}$

(Optional) Indicates that the aggregates may not be SPD for the Schwarz method. The following options exist for use_nonsymm:

- 0 assume SPD (default)
 1 assume non-symmetric
 - _ 6.2.76 _

int HYPRE_BoomerAMGSetSym (HYPRE_Solver solver, int sym)

(Optional) Defines symmetry for ParaSAILS. For further explanation see description of ParaSAILS.

 $_$ 6.2.77 $_$

int HYPRE_BoomerAMGSetLevel (HYPRE_Solver solver, int level)

(Optional) Defines number of levels for ParaSAILS. For further explanation see description of ParaSAILS.

HYPRE_BoomerAMGSetThreshold (HYPRE_Solver solver, HYPRE_Real threshold)

(Optional) Defines threshold for ParaSAILS. For further explanation see description of ParaSAILS.

___ 6.2.79 _____

int **HYPRE_BoomerAMGSetFilter** (HYPRE_Solver solver, HYPRE_Real filter)

(Optional) Defines filter for ParaSAILS. For further explanation see description of ParaSAILS.

_ 6.2.80 _

int

 $\label{eq:hypre_bound} \begin{aligned} \mathbf{HYPRE_BoomerAMGSetDropTol} \ (\mathbf{HYPRE_Solver} \ solver, \ \mathbf{HYPRE_Real} \\ \mathbf{drop_tol}) \end{aligned}$

(Optional) Defines drop tolerance for PILUT. For further explanation see description of PILUT.

6.2.81

int **HYPRE_BoomerAMGSetMaxNzPerRow** (HYPRE_Solver solver, int max_nz_per_row)

(Optional) Defines maximal number of nonzeros for PILUT. For further explanation see description of PILUT.

Int HYPRE_BoomerAMGSetEuclidFile (HYPRE_Solver solver, char* euclidfile)

(Optional) Defines name of an input file for Euclid parameters. For further explanation see description of Euclid.

__ 6.2.83 _____

int HYPRE_BoomerAMGSetEuLevel (HYPRE_Solver solver, int eu_level)

(Optional) Defines number of levels for ILU(k) in Euclid. For further explanation see description of Euclid.

6.2.84

int **HYPRE_BoomerAMGSetEuSparseA** (HYPRE_Solver solver, HYPRE_Real eu_sparse_A)

(Optional) Defines filter for ILU(k) for Euclid. For further explanation see description of Euclid.

6.2.85

int HYPRE_BoomerAMGSetEuBJ (HYPRE_Solver solver, int eu_bj)

(Optional) Defines use of block jacobi ILUT for Euclid. For further explanation see description of Euclid.

int HYPRE_BoomerAMGSetPrintLevel (HYPRE_Solver solver, int print_level)

(Optional) Requests automatic printing of setup and solve information.

- 0 no printout (default)
- 1 print setup information
- 2 print solve information
- 3 | print both setup and solve information

Note, that if one desires to print information and uses BoomerAMG as a preconditioner, suggested print_level is 1 to avoid excessive output, and use print_level of solver for solve phase information.

6.2.87

int HYPRE_BoomerAMGSetLogging (HYPRE_Solver solver, int logging)

(Optional) Requests additional computations for diagnostic and similar data to be logged by the user. Default to 0 for do nothing. The latest residual will be available if logging > 1.

__ 6.2.88 ____

HYPRE_BoomerAMGSetDebugFlag (HYPRE_Solver solver, int debug_flag)

(Optional)

6.2.89

int

HYPRE_BoomerAMGInitGridRelaxation (int** num_grid_sweeps_ptr, int** grid_relax_type_ptr, int*** grid_relax_points_ptr, int coarsen_type, HYPRE_Real** relax_weights_ptr, int max_levels)

(Optional) This routine will be eliminated in the future

6 2 90

int HYPRE_BoomerAMGSetRAP2 (HYPRE_Solver solver, int rap2)

(Optional) If rap2 not equal 0, the triple matrix product RAP is replaced by two matrix products

__ 6.2.91 _____

int

 $\label{lem:hypre_bound} \mbox{\bf HYPRE_BoomerAMGSetKeepTranspose} \ (\mbox{HYPRE_Solver solver}, \mbox{ int keepTranspose})$

(Optional) If set to 1, the local interpolation transposes will be saved to use more efficient matvecs instead of matvecTs $^{\circ}$

6.3

ParCSR ParaSails Preconditioner

Names		
6.3.1	int HYPRE_ParaSailsCreate (MPI_Comm comm, HYPRE_Solver* solver) Create a ParaSails preconditioner	165
6.3.2	int	
	HYPRE_ParaSailsDestroy (HYPRE_Solver solver)	
	Destroy a ParaSails preconditioner	165
6.3.3	int	
	HYPRE_ParaSailsSetup (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x)	
	Set up the ParaSails preconditioner.	165
6.3.4	int	
	HYPRE_ParaSailsSolve (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x)	
	Apply the ParaSails preconditioner.	166
6.3.5	int	
	HYPRE_ParaSailsSetParams (HYPRE_Solver solver, HYPRE_Real thresh, int nlevels)	
	Set the threshold and levels parameter for the ParaSails preconditioner	166
6.3.6	int	

	$\label{lem:hypre_paraSailsSetFilter} \textbf{HYPRE_Solver solver}, \ \ \textbf{HYPRE_Real filter})$	
	Set the filter parameter for the ParaSails preconditioner	167
6.3.7	int	
	HYPRE_ParaSailsSetSym (HYPRE_Solver solver, int sym)	
	Set the symmetry parameter for the ParaSails preconditioner	167
6.3.8	int	
	HYPRE_ParaSailsSetLoadbal (HYPRE_Solver solver, HYPRE_Real loadbal)	
	Set the load balance parameter for the ParaSails preconditioner	167
6.3.9	int	
	HYPRE_ParaSailsSetReuse (HYPRE_Solver solver, int reuse)	
	Set the pattern reuse parameter for the ParaSails preconditioner	168
6.3.10	int	
	HYPRE_ParaSailsSetLogging (HYPRE_Solver solver, int logging)	
	Set the logging parameter for the ParaSails preconditioner	168
6.3.11	int	
	HYPRE_ParaSailsBuildIJMatrix (HYPRE_Solver solver,	
	HYPRE_IJMatrix* pij_A)	
	Build IJ Matrix of the sparse approximate inverse (factor)	169

Parallel sparse approximate inverse preconditioner for the ParCSR matrix format.

_ 6.3.1 _

int HYPRE_ParaSailsCreate (MPI_Comm comm, HYPRE_Solver* solver)

Create a ParaSails preconditioner

_ 6.3.2 _

int HYPRE_ParaSailsDestroy (HYPRE_Solver solver)

Destroy a ParaSails preconditioner

__ 6.3.3 __

HYPRE_ParaSailsSetup (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x)

Set up the ParaSails preconditioner. This function should be passed to the iterative solver SetPrecond function.

Parameters: solver [IN] Preconditioner object to set up.

A [IN] ParCSR matrix used to construct the precondi-

tioner.

b Ignored by this function.x Ignored by this function.

6.3.4

HYPRE_ParaSailsSolve (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x)

Apply the ParaSails preconditioner. This function should be passed to the iterative solver SetPrecond function.

Parameters: solver [IN] Preconditioner object to apply.

A Ignored by this function.

b [IN] Vector to precondition.

x [OUT] Preconditioned vector.

6.3.5 _

int **HYPRE_ParaSailsSetParams** (HYPRE_Solver solver, HYPRE_Real thresh, int nlevels)

Set the threshold and levels parameter for the ParaSails preconditioner. The accuracy and cost of ParaSails are parameterized by these two parameters. Lower values of the threshold parameter and higher values of levels parameter lead to more accurate, but more expensive preconditioners.

Parameters: solver [IN] Preconditioner object for which to set parameters.

thresh [IN] Value of threshold parameter, $0 \le \text{thresh} \le 1$. The

deault valuue is 0.1.

nlevels [IN] Value of levels parameter, $0 \le$ nlevels. The dfault

valuue is 1.

6.3.6

int HYPRE_ParaSailsSetFilter (HYPRE_Solver solver, HYPRE_Real filter)

Set the filter parameter for the ParaSails preconditioner.

Parameters: solver [IN] Preconditioner object for which to set filter pa-

rameter.

filter [IN] Value of filter parameter. The filter parameter

is used to drop small nonzeros in the preconditioner, to reduce the oost of applying the preconditioner. Values from .055 to 0.1 are recommended. The efault value

is 0.1.

6.3.7

int HYPRE_ParaSailsSetSym (HYPRE_Solver solver, int sym)

Set the symmetry parameter for the ParaSails preconditioner.

value	meaning
0	nonsymmetric and/or indefinite problem, and nonsymmetric preconditioner
1	SPD problem, and SPD (factored) preconditioner
2	nonsymmetric, definite problem, and SPD (factored) preconditioner

Parameters: solver [IN] Preconditioner object for which to set symmetry

parameter.

sym [IN] Symmetry parameter.

6.3.8

int HYPRE_ParaSailsSetLoadbal (HYPRE_Solver solver, HYPRE_Real loadbal)

Set the load balance parameter for the ParaSails preconditioner.

Parameters:

solver

[IN] Preconditioner object for which to set the load balance parameter.

loadbal

[IN] Value of the load balance parameter, $0 \le \text{loadbl} \le 1$. A zero value indicates that no load alance is attepted; a value of unity indicatestat perfect llooad balance will e attempted. The recommended value is 0.9 to balance the overhad ofdata exchhanes for load balancing. No load balancings need if the preeconditioner is very sparse and fast to construct. The default value when this parameter is not tsse iis 0.

6.3.9

int HYPRE_ParaSailsSetReuse (HYPRE_Solver solver, int reuse)

Set the pattern reuse parameter for the ParaSails preconditioner.

Parameters:

solver

[IN] Preconditioner object for which to set the pattern

reuse parameter.

reuse

[IN] Value of the pattern reuse parameter. A nonzero value indicaes that the pattern of the preconditioner should be reused for ubsequent constructions of the productioner... A zero value inicates that the peconditioner should be constructed from scrath. The default value when this parameter is not se iis 0.

.

int HYPRE_ParaSailsSetLogging (HYPRE_Solver solver, int logging)

Set the logging parameter for the ParaSails preconditioner.

Parameters:

solver

[IN] Preconditioner object for which to set the logging-

parameter.

logging

[IN] Value of the logging parameter. A nonzero valueseds statistics of the setup procedure to stdout. The default value when this parameter is not set is 0.

6.3.11

HYPRE_ParaSailsBuildIJMatrix (HYPRE_Solver solver, HYPRE_IJMatrix* pij_A)

Build IJ Matrix of the sparse approximate inverse (factor). This function explicitly creates the IJ Matrix corresponding to the sparse approximate inverse or the inverse factor. Example: HYPRE_IJMatrix ij_A; HYPRE_ParaSailsBuildIJMatrix(solver, &ij_A);

Parameters:

solver [IN] Preconditioner object.
pij_A [OUT] Pointer to the IJ Matrix.

_ 6.4

ParCSR Euclid Preconditioner

Names		
6.4.1	int HYPRE_EuclidCreate (MPI_Comm comm, HYPRE_Solver* solver) Create a Euclid object	170
6.4.2	int HYPRE_EuclidDestroy (HYPRE_Solver solver) Destroy a Euclid object	171
6.4.3	int HYPRE_EuclidSetup (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x) Set up the Euclid preconditioner.	171
6.4.4	int HYPRE_EuclidSolve (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x) Apply the Euclid preconditioner.	171
6.4.5	int HYPRE_EuclidSetParams (HYPRE_Solver solver, int argc, char* argv[]) Insert (name, value) pairs in Euclid's options database by passing Euclid the command line (or an array of strings).	172
6.4.6	int HYPRE_EuclidSetParamsFromFile (HYPRE_Solver solver, char* filename) Insert (name, value) pairs in Euclid's options database	172
6.4.7	int HYPRE_EuclidSetLevel (HYPRE_Solver solver, int level) Set level k for ILU(k) factorization, default: 1	172
6.4.8	int	

	HYPRE_EuclidSetBJ (HYPRE_Solver solver, int bj)	
	Use block Jacobi ILU preconditioning instead of PILU	173
6.4.9	int	
	HYPRE_EuclidSetStats (HYPRE_Solver solver, int eu_stats)	
	If eu_stats not equal 0, a summary of runtime settings and timing information is printed to stdout	173
6.4.10	int	
	HYPRE_EuclidSetMem (HYPRE_Solver solver, int eu_mem)	
	If eu_mem not equal 0, a summary of Euclid's memory usage is printed to stdout	173
6.4.11	int	
	HYPRE_EuclidSetSparseA (HYPRE_Solver solver, HYPRE_Real sparse_A)	
	Defines a drop tolerance for $ILU(k)$	173
6.4.12	int	
	HYPRE_EuclidSetRowScale (HYPRE_Solver solver, int row_scale)	
	If row_scale not equal 0, values are scaled prior to factorization so that	
	largest value in any row is +1 or -1.	173
6.4.13	int	
	HYPRE_EuclidSetILUT (HYPRE_Solver solver, HYPRE_Real drop_tol)	
	uses ILUT and defines a drop tolerance relative to the largest absolute value	
	of any entry in the row being factored	174

MPI Parallel ILU preconditioner

Options summary:

Option	Default	Synopsis
-level	1	ILU(k) factorization level
-bj	0 (false)	Use Block Jacobi ILU instead of PILU
-eu_stats	0 (false)	Print internal timing and statistics
-eu_mem	0 (false)	Print internal memory usage

6.4.1

 $int \ \mathbf{HYPRE_EuclidCreate} \ (\mathrm{MPI_Comm} \ \mathrm{comm}, \ \mathrm{HYPRE_Solver*} \ \mathrm{solver})$

Create a Euclid object

6.4.2

int HYPRE_EuclidDestroy (HYPRE_Solver solver)

Destroy a Euclid object

 $_$ 6.4.3 $_$

int

HYPRE_EuclidSetup (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x)

Set up the Euclid preconditioner. This function should be passed to the iterative solver SetPrecond function.

Parameters: solver [IN] Preconditioner object to set up.

A [IN] ParCSR matrix used to construct the precondi-

tioner.

b Ignored by this function.x Ignored by this function.

_ 6.4.4 _

int

HYPRE_EuclidSolve (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x)

Apply the Euclid preconditioner. This function should be passed to the iterative solver SetPrecond function.

Parameters: solver [IN] Preconditioner object to apply.

A Ignored by this function.
b [IN] Vector to precondition.
x [OUT] Preconditioned vector.

6.4.5

int HYPRE_EuclidSetParams (HYPRE_Solver solver, int argc, char* argv[])

Insert (name, value) pairs in Euclid's options database by passing Euclid the command line (or an array of strings). All Euclid options (e.g, level, drop-tolerance) are stored in this database. If a (name, value) pair already exists, this call updates the value. See also: HYPRE_EuclidSetParamsFromFile.

Parameters:

argc [IN] Length of argv array argv [IN] Array of strings

6.4.6

int HYPRE_EuclidSetParamsFromFile (HYPRE_Solver solver, char* filename)

Insert (name, value) pairs in Euclid's options database. Each line of the file should either begin with a "#," indicating a comment line, or contain a (name value) pair, e.g.:

>cat optionsFile #sample runtime parameter file -blockJacobi 3 -matFile /home/hysom/myfile.euclid -doSomething true -xx_coeff -1.0

See also: $HYPRE_EuclidSetParams$.

Parameters: filename [IN] Pathname/filename to read

6.4.7 _

int HYPRE_EuclidSetLevel (HYPRE_Solver solver, int level)

Set level k for ILU(k) factorization, default: 1

__ 6.4.8 ____

int HYPRE_EuclidSetBJ (HYPRE_Solver solver, int bj)

Use block Jacobi ILU preconditioning instead of PILU

__ 6.4.9 _____

int HYPRE_EuclidSetStats (HYPRE_Solver solver, int eu_stats)

If eu_stats not equal 0, a summary of runtime settings and timing information is printed to stdout

6.4.10

int HYPRE_EuclidSetMem (HYPRE_Solver solver, int eu_mem)

If eu_mem not equal 0, a summary of Euclid's memory usage is printed to stdout

_ 6.4.11 ____

int **HYPRE_EuclidSetSparseA** (HYPRE_Solver solver, HYPRE_Real sparse_A)

Defines a drop tolerance for $\mathrm{ILU}(k)$. Default: 0 Use with HYPRE_EuclidSetRowScale. Note that this can destroy symmetry in a matrix.

6.4.12

int HYPRE_EuclidSetRowScale (HYPRE_Solver solver, int row_scale)

If row_scale not equal 0, values are scaled prior to factorization so that largest value in any row is +1 or -1. Note that this can destroy symmetry in a matrix.

6.4.13

int HYPRE_EuclidSetILUT (HYPRE_Solver solver, HYPRE_Real drop_tol)

uses ILUT and defines a drop tolerance relative to the largest absolute value of any entry in the row being factored

_ 6.5 _

ParCSR Pilut Preconditioner

Names		
6.5.1	int HYPRE_ParCSRPilutCreate (MPI_Comm comm, HYPRE_Solver* solver) Create a preconditioner object	175
6.5.2	int HYPRE_ParCSRPilutDestroy (HYPRE_Solver solver) Destroy a preconditioner object	175
6.5.3	int HYPRE_ParCSRPilutSetup (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x)	175
6.5.4	int HYPRE_ParCSRPilutSolve (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x) Precondition the system	175
6.5.5	int HYPRE_ParCSRPilutSetMaxIter (HYPRE_Solver solver, int max_iter) (Optional) Set maximum number of iterations	175
6.5.6	int HYPRE_ParCSRPilutSetDropTolerance (HYPRE_Solver solver, HYPRE_Real tol) (Optional)	176
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int HYPRE_ParCSRPilutCreate (MPI_Comm comm, HYPRE_Solver* solver)

Create a preconditioner object

__ 6.5.2 ____

int HYPRE_ParCSRPilutDestroy (HYPRE_Solver solver)

Destroy a preconditioner object

_ 6.5.3 _

int

HYPRE_ParCSRPilutSetup (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x)

6.5.4

HYPRE_ParCSRPilutSolve (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x)

Precondition the system

6.5.5

int HYPRE_ParCSRPilutSetMaxIter (HYPRE_Solver solver, int max_iter)

(Optional) Set maximum number of iterations

6.5.6

HYPRE_ParCSRPilutSetDropTolerance (HYPRE_Solver solver, HYPRE_Real tol)

(Optional)

6.5.7

int HYPRE_ParCSRPilutSetFactorRowSize (HYPRE_Solver solver, int size)

(Optional)

6.6

ParCSR AMS Solver and Preconditioner

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6.6.3	int	
	HYPRE_AMSSetup (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x)	
	Set up the AMS solver or preconditioner.	179
6.6.4	int	
	HYPRE_AMSSolve (HYPRE_Solver solver, HYPRE_ParCSRMatrix A,	
	HYPRE_ParVector b, HYPRE_ParVector x)	
	Solve the system or apply AMS as a preconditioner	180
6.6.5	int	
	HYPRE_AMSSetDimension (HYPRE_Solver solver, int dim)	
	(Optional) Sets the problem dimension (2 or 3)	180
6.6.6	int	
	HYPRE_AMSSetDiscreteGradient (HYPRE_Solver solver,	
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	Sets the discrete gradient matrix G.	180
6.6.7	int	

	HYPRE_AMSSetCoordinateVectors (HYPRE_Solver solver,	
	HYPRE_ParVector x,	
	HYPRE_ParVector y,	
	HYPRE_ParVector z)	
	Sets the x , y and z coordinates of the vertices in the mesh	181
6.6.8	int	
	HYPRE_AMSSetEdgeConstantVectors (HYPRE_Solver solver,	
	HYPRE_ParVector Gx,	
	HYPRE_ParVector Gy,	
	HYPRE_ParVector Gz)	
	Sets the vectors Gx , Gy and Gz which give the representations of the con-	
	stant vector fields $(1, 0, 0)$, $(0, 1, 0)$ and $(0, 0, 1)$ in the edge element	
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6.6.9	int	
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	HYPRE_ParCSRMatrix Pi,	
	HYPRE_ParCSRMatrix Pix,	
	HYPRE_ParCSRMatrix Piy,	
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6.6.10	int	
	${\bf HYPRE_AMSSetAlphaPoissonMatrix} \ ({\bf HYPRE_Solver} \ solver,$	
	HYPRE_ParCSRMatrix A_alpha)	
	(Optional) Sets the matrix A_{α} corresponding to the Poisson problem with	
	coefficient α (the curl-curl term coefficient in the Maxwell problem)	182
6.6.11	int	
	HYPRE_AMSSetBetaPoissonMatrix (HYPRE_Solver solver,	
	HYPRE_ParCSRMatrix A_beta)	
	(Optional) Sets the matrix A_{β} corresponding to the Poisson problem with	
	coefficient β (the mass term coefficient in the Maxwell problem)	182
6.6.12	int	
	HYPRE_AMSSetInteriorNodes (HYPRE_Solver solver,	
	HYPRE_ParVector interior_nodes)	
	(Optional) Set the list of nodes which are interior to a zero-conductivity	
	region.	182
6.6.13	int	
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	(Optional) Set the frequency at which a projection onto the compatible sub-	
	space for problems with zero-conductivity regions is performed	183
6.6.14	int	
0.0.14	HYPRE_AMSSetMaxIter (HYPRE_Solver solver, int maxit)	
	(Optional) Sets maximum number of iterations, if AMS is used as a solver.	
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6615		
6.6.15	int HYPRE_AMSSetTol (HYPRE_Solver solver, HYPRE_Real tol)	
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0.010		100
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	HYPRE_AMSSetCycleType (HYPRE_Solver solver, int cycle_type) (Optional) Choose which three-level solver to use	183
6.6.17	int HYPRE_AMSSetPrintLevel (HYPRE_Solver solver, int print_level) (Optional) Control how much information is printed during the solution iterations.	184
6.6.18	int	
	HYPRE_AMSSetSmoothingOptions (HYPRE_Solver solver, int relax_type, int relax_times,	
	HYPRE_Real relax_weight,	
	HYPRE_Real omega)	
	(Optional) Sets relaxation parameters for A	184
6.6.19	int	
	HYPRE_AMSSetAlphaAMGOptions (HYPRE_Solver solver,	
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	int alpha_agg_levels,	
	int alpha_relax_type, HYPRE_Real	
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	int alpha_interp_type,	
	$\operatorname{int\ alpha_Pmax})$ (Optional) Sets AMG parameters for B_{Π}	185
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6.6.20	int HVDDE AMSSet Alpha AMCCeanag Delev Tyme (HVDDE Selven gelven int	
	HYPRE_AMSSetAlphaAMGCoarseRelaxType (HYPRE_Solver solver, int alpha_coarse_relax_type)	
	(Optional) Sets the coarsest level relaxation in the AMG solver for B_{Π}	185
6.6.21	int	
	HYPRE_AMSSetBetaAMGOptions (HYPRE_Solver solver,	
	int beta_coarsen_type,	
	int beta_agg_levels, int beta_relax_type,	
	HYPRE_Real beta_strength_threshold, int beta_interp_type, int beta_Pmax)	
	(Optional) Sets AMG parameters for B_G	185
C C 00	,	100
6.6.22	int HYPRE_AMSSetBetaAMGCoarseRelaxType (HYPRE_Solver solver,	
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	(Optional) Sets the coarsest level relaxation in the AMG solver for B_G .	185
6.6.23	int	
	HYPRE_AMSGetNumIterations (HYPRE_Solver solver, int* num_iterations)	
	Returns the number of iterations taken	186
0.004	·	100
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	HYPRE_AMSGetFinalRelativeResidualNorm (HYPRE_Solver solver, HYPRE_Real*	
	rel_resid_norm)	
	Returns the norm of the final relative residual	186
6.6.25	int	- •
0.0.20	1110	

HYPRE_AMSProjectOutGradients (HYPRE_Solver solver, HYPRE_ParVector x) For problems with zero-conductivity regions, project the vector onto the compatible subspace: $x = (I - G_0(G_0^t G_0)^{-1} G_0^T)x$, where G_0 is the discrete gradient restricted to the interior nodes of the regions with zero conductivity. 186 6.6.26 int HYPRE_AMSConstructDiscreteGradient (HYPRE_ParCSRMatrix A. HYPRE_ParVector x_coord, int* edge_vertex, int edge_orientation, HYPRE_ParCSRMatrix* G) $Construct\ and\ return\ the\ lowest-order\ discrete\ gradient\ matrix\ G\ using\ some$ 186

Parallel auxiliary space Maxwell solver and preconditioner

6.6.1

int HYPRE_AMSCreate (HYPRE_Solver* solver)

Create an AMS solver object

6.6.2

int HYPRE_AMSDestroy (HYPRE_Solver solver)

Destroy an AMS solver object

_ 6.6.3 ___

HYPRE_AMSSetup (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x)

Set up the AMS solver or preconditioner. If used as a preconditioner, this function should be passed to the iterative solver SetPrecond function.

Parameters:	solver A	[IN] object to be set up.[IN] ParCSR matrix used to construct the solver/preconditioner.
	b	Ignored by this function.
	x	Ignored by this function.

6.6.4

HYPRE_AMSSolve (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x)

Solve the system or apply AMS as a preconditioner. If used as a preconditioner, this function should be passed to the iterative solver SetPrecond function.

Parameters:	solver A	[IN] solver or preconditioner object to be applied.[IN] ParCSR matrix, matrix of the linear system to be solved
	b	[IN] right hand side of the linear system to be solved
	х	[OUT] approximated solution of the linear system to
		be solved

 $_$ 6.6.5 $_$

int HYPRE_AMSSetDimension (HYPRE_Solver solver, int dim)

(Optional) Sets the problem dimension (2 or 3). The default is 3.

6.6.6

int
HYPRE_AMSSetDiscreteGradient (HYPRE_Solver solver,
HYPRE_ParCSRMatrix G)

Sets the discrete gradient matrix G. This function should be called before HYPRE_AMSSetup()!

6.6.7

int
HYPRE_AMSSetCoordinateVectors (HYPRE_Solver solver,
HYPRE_ParVector x, HYPRE_ParVector y, HYPRE_ParVector z)

Sets the x, y and z coordinates of the vertices in the mesh.

Either HYPRE_AMSSetCoordinateVectors() or HYPRE_AMSSetEdgeConstantVectors() should be called before HYPRE_AMSSetup()!

6.6.8

int
HYPRE_AMSSetEdgeConstantVectors (HYPRE_Solver solver,
HYPRE_ParVector Gx, HYPRE_ParVector Gy, HYPRE_ParVector Gz)

Sets the vectors Gx, Gy and Gz which give the representations of the constant vector fields (1,0,0), (0,1,0) and (0,0,1) in the edge element basis.

Either HYPRE_AMSSetCoordinateVectors() or HYPRE_AMSSetEdgeConstantVectors() should be called before HYPRE_AMSSetup()!

_ 6.6.9 _

int
HYPRE_AMSSetInterpolations (HYPRE_Solver solver,
HYPRE_ParCSRMatrix Pi, HYPRE_ParCSRMatrix Pix, HYPRE_ParCSRMatrix Piy, HYPRE_ParCSRMatrix Piz)

(Optional) Set the (components of) the Nedelec interpolation matrix $\Pi = [\Pi^x, \Pi^y, \Pi^z]$.

This function is generally intended to be used only for high-order Nedelec discretizations (in the lowest order case, Π is constructed internally in AMS from the discreet gradient matrix and the coordinates of the vertices), though it can also be used in the lowest-order case or for other types of discretizations (e.g. ones based on the second family of Nedelec elements).

By definition, Π is the matrix representation of the linear operator that interpolates (high-order) vector nodal finite elements into the (high-order) Nedelec space. The component matrices are defined as $\Pi^x \varphi = \Pi(\varphi, 0, 0)$ and similarly for Π^y and Π^z . Note that all these operators depend on the choice of the basis and degrees of freedom in the high-order spaces.

The column numbering of Pi should be node-based, i.e. the x/y/z components of the first node (vertex or high-order dof) should be listed first, followed by the x/y/z components of the second node and so on (see the documentation of HYPRE_BoomerAMGSetDofFunc).

If used, this function should be called before HYPRE_AMSSetup() and there is no need to provide the vertex coordinates. Furthermore, only one of the sets $\{\Pi\}$ and $\{\Pi^x, \Pi^y, \Pi^z\}$ needs to be specified (though it is OK to provide both). If Pix is NULL, then scalar Π -based AMS cycles, i.e. those with cycle_type > 10, will be unavailable. Similarly, AMS cycles based on monolithic Π (cycle_type < 10) require that Pi is not NULL.

6.6.10

HYPRE_AMSSetAlphaPoissonMatrix (HYPRE_Solver solver, HYPRE_ParCSRMatrix A_alpha)

(Optional) Sets the matrix A_{α} corresponding to the Poisson problem with coefficient α (the curl-curl term coefficient in the Maxwell problem).

If this function is called, the coarse space solver on the range of Π^T is a block-diagonal version of A_{Π} . If this function is not called, the coarse space solver on the range of Π^T is constructed as $\Pi^T A \Pi$ in HYPRE_AMSSetup(). See the user's manual for more details.

_ 6.6.11 __

int **HYPRE_AMSSetBetaPoissonMatrix** (HYPRE_Solver solver,
HYPRE_ParCSRMatrix A_beta)

(Optional) Sets the matrix A_{β} corresponding to the Poisson problem with coefficient β (the mass term coefficient in the Maxwell problem).

If not given, the Poisson matrix will be computed in HYPRE_AMSSetup(). If the given matrix is NULL, we assume that β is identically 0 and use two-level (instead of three-level) methods. See the user's manual for more details.

 $_{-}$ 6.6.12 $_{-}$

HYPRE_AMSSetInteriorNodes (HYPRE_Solver solver, HYPRE_ParVector interior_nodes)

(Optional) Set the list of nodes which are interior to a zero-conductivity region. This way, a more robust solver is constructed, that can be iterated to lower tolerance levels. A node is interior if its entry in the array is 1.0. This function should be called before HYPRE_AMSSetup()!

6.6.13

int **HYPRE_AMSSetProjectionFrequency** (HYPRE_Solver solver, int projection_frequency)

(Optional) Set the frequency at which a projection onto the compatible subspace for problems with zero-conductivity regions is performed. The default value is 5.

6.6.14

int HYPRE_AMSSetMaxIter (HYPRE_Solver solver, int maxit)

(Optional) Sets maximum number of iterations, if AMS is used as a solver. To use AMS as a preconditioner, set the maximum number of iterations to 1. The default is 20.

 $_{-}$ 6.6.15 $_{-}$

int HYPRE_AMSSetTol (HYPRE_Solver solver, HYPRE_Real tol)

(Optional) Set the convergence tolerance, if AMS is used as a solver. When using AMS as a preconditioner, set the tolerance to 0.0. The default is 10^{-6} .

6.6.16

int HYPRE_AMSSetCycleType (HYPRE_Solver solver, int cycle_type)

(Optional) Choose which three-level solver to use. Possible values are:

- 1 3-level multiplicative solver (01210)
- 2 | 3-level additive solver (0+1+2)
- 3 | 3-level multiplicative solver (02120)
- 4 | 3-level additive solver (010+2)
- 5 | 3-level multiplicative solver (0102010)
- 6 3-level additive solver (1+020)
- 7 | 3-level multiplicative solver (0201020)
- 8 | 3-level additive solver (0(1+2)0)
- 11 | 5-level multiplicative solver (013454310)
- 12 | 5-level additive solver (0+1+3+4+5)
- 13 | 5-level multiplicative solver (034515430)
- 14 | 5-level additive solver (01(3+4+5)10)

The default is 1. See the user's manual for more details.

6.6.17

int HYPRE_AMSSetPrintLevel (HYPRE_Solver solver, int print_level)

(Optional) Control how much information is printed during the solution iterations. The default is 1 (print residual norm at each step).

6.6.18

int

HYPRE_AMSSetSmoothingOptions (HYPRE_Solver solver, int relax_type, int relax_times, HYPRE_Real relax_weight, HYPRE_Real omega)

(Optional) Sets relaxation parameters for A. The defaults are 2, 1, 1.0, 1.0.

The available options for relax_type are:

- 1 ℓ_1 -scaled Jacobi
- 2 ℓ_1 -scaled block symmetric Gauss-Seidel/SSOR
- 3 | Kaczmarz
- 4 | truncated version of ℓ_1 -scaled block symmetric Gauss-Seidel/SSOR
- 16 Chebyshev

6.6.19

int

HYPRE_AMSSetAlphaAMGOptions (HYPRE_Solver solver, int alpha_coarsen_type, int alpha_agg_levels, int alpha_relax_type, HYPRE_Real alpha_strength_threshold, int alpha_interp_type, int alpha_Pmax)

(Optional) Sets AMG parameters for B_{Π} . The defaults are 10, 1, 3, 0.25, 0, 0. See the user's manual for more details.

6.6.20

int

HYPRE_AMSSetAlphaAMGCoarseRelaxType (HYPRE_Solver solver, int alpha_coarse_relax_type)

(Optional) Sets the coarsest level relaxation in the AMG solver for B_{Π} . The default is 8 (l1-GS). Use 9, 19, 29 or 99 for a direct solver.

6.6.21

int

HYPRE_AMSSetBetaAMGOptions (HYPRE_Solver solver, int beta_coarsen_type, int beta_agg_levels, int beta_relax_type, HYPRE_Real beta_strength_threshold, int beta_interp_type, int beta_Pmax)

(Optional) Sets AMG parameters for B_G . The defaults are 10, 1, 3, 0.25, 0, 0. See the user's manual for more details.

 $_$ 6.6.22 $_$

int

 $\label{lem:hypre_ams} \textbf{HYPRE_AMSSetBetaAMGCoarseRelaxType} \ (\textbf{HYPRE_Solver solver}, \ \textbf{int beta_coarse_relax_type})$

(Optional) Sets the coarsest level relaxation in the AMG solver for B_G . The default is 8 (l1-GS). Use 9, 19, 29 or 99 for a direct solver.

6.6.23

int
HYPRE_AMSGetNumIterations (HYPRE_Solver solver, int* num_iterations)

Returns the number of iterations taken

 $_$ 6.6.24 $_$

int **HYPRE_AMSGetFinalRelativeResidualNorm** (HYPRE_Solver solver, HYPRE_Real* rel_resid_norm)

Returns the norm of the final relative residual

 $_{-}$ 6.6.25 $_{--}$

HYPRE_AMSProjectOutGradients (HYPRE_Solver solver, HYPRE_ParVector x)

For problems with zero-conductivity regions, project the vector onto the compatible subspace: $x = (I - G_0(G_0^t G_0)^{-1} G_0^T)x$, where G_0 is the discrete gradient restricted to the interior nodes of the regions with zero conductivity. This ensures that x is orthogonal to the gradients in the range of G_0 .

This function is typically called after the solution iteration is complete, in order to facilitate the visualization of the computed field. Without it the values in the zero-conductivity regions contain kernel components.

6.6.26

int
HYPRE_AMSConstructDiscreteGradient (HYPRE_ParCSRMatrix A,
HYPRE_ParVector x_coord, int* edge_vertex, int edge_orientation,
HYPRE_ParCSRMatrix* G)

Construct and return the lowest-order discrete gradient matrix G using some edge and vertex information. We assume that edge_vertex lists the edge vertices consecutively, and that the orientation of all edges is consistent.

If edge_orientation = 1, the edges are already oriented.

If edge_orientation = 2, the orientation of edge i depends only on the sign of edge_vertex[2*i+1] - edge_vertex[2*i].

_ 6.7 _

ParCSR ADS Solver and Preconditioner

Names		
6.7.1	int	
	HYPRE_ADSCreate (HYPRE_Solver* solver)	
	Create an ADS solver object	189
6.7.2	int	
	HYPRE_ADSDestroy (HYPRE_Solver solver)	
	Destroy an ADS solver object	189
6.7.3	int	
	HYPRE_ADSSetup (HYPRE_Solver solver, HYPRE_ParCSRMatrix A,	
	HYPRE_ParVector b, HYPRE_ParVector x)	
	Set up the ADS solver or preconditioner.	189
6.7.4	int	
	HYPRE_ADSSolve (HYPRE_Solver solver, HYPRE_ParCSRMatrix A,	
	HYPRE_ParVector b, HYPRE_ParVector x)	
	Solve the system or apply ADS as a preconditioner.	190
6.7.5	int	
	HYPRE_ADSSetDiscreteCurl (HYPRE_Solver solver,	
	HYPRE_ParCSRMatrix C)	
	Sets the discrete curl matrix C.	190
6.7.6	int	
	HYPRE_ADSSetDiscreteGradient (HYPRE_Solver solver,	
	HYPRE_ParCSRMatrix G)	
	Sets the discrete gradient matrix G.	190
6.7.7	int	
	HYPRE_ADSSetCoordinateVectors (HYPRE_Solver solver,	
	HYPRE_ParVector x,	
	HYPRE_ParVector y,	
	HYPRE_ParVector z)	
	Sets the x , y and z coordinates of the vertices in the mesh	191
6.7.8	int	

	HYPRE_ADSSetInterpolations (HYPRE_Solver solver, HYPRE_ParCSRMatrix RT_Pi, HYPRE_ParCSRMatrix RT_Pix, HYPRE_ParCSRMatrix RT_Piy,	
	HYPRE_ParCSRMatrix RT_Piz,	
	HYPRE_ParCSRMatrix ND_Pi,	
	HYPRE_ParCSRMatrix ND_Pix,	
	HYPRE_ParCSRMatrix ND_Piy,	
	HYPRE_ParCSRMatrix ND_Piz)	
	(Optional) Set the (components of) the Raviart-Thomas (Π_{RT}) and the Ned-	
	elec (Π_{ND}) interpolation matrices.	191
6.7.9	int	
	HYPRE_ADSSetMaxIter (HYPRE_Solver solver, int maxit) (Optional) Sets maximum number of iterations, if ADS is used as a solver.	192
0.7.10		132
6.7.10	int HYPRE_ADSSetTol (HYPRE_Solver solver, HYPRE_Real tol)	
	(Optional) Set the convergence tolerance, if ADS is used as a solver	192
0 = 44		132
6.7.11	int HVDDE ADSS at Cycle Type (HVDDE Selven selven, intervals type)	
	HYPRE_ADSSetCycleType (HYPRE_Solver solver, int cycle_type) (Optional) Choose which auxiliary-space solver to use	192
		132
6.7.12	int	
	HYPRE_ADSSetPrintLevel (HYPRE_Solver solver, int print_level) (Optional) Control how much information is printed during the solution	
	iterations.	193
<i>c</i> 7 19		100
6.7.13	int HYPRE_ADSSetSmoothingOptions (HYPRE_Solver solver, int relax_type,	
	int relax_times,	
	HYPRE_Real relax_weight,	
	HYPRE_Real omega)	
	(Optional) Sets relaxation parameters for A	193
6.7.14	int	
0.7.14	HYPRE_ADSSetChebySmoothingOptions (HYPRE_Solver solver,	
	int cheby_order,	
	int cheby_fraction)	
	(Optional) Sets parameters for Chebyshev relaxation	193
6715		
6.7.15	int HYPRE_ADSSetAMSOptions (HYPRE_Solver solver, int cycle_type,	
	int coarsen_type, int agg_levels,	
	int relax_type, int agg_levels,	
	HYPRE_Real strength_threshold,	
	int interp_type, int Pmax)	
	(Optional) Sets AMS parameters for B_C	193
6.7.16		
0.7.10	int HYPRE_ADSSetAMGOptions (HYPRE_Solver solver, int coarsen_type,	
	int agg_levels, int relax_type,	
	HYPRE_Real strength_threshold,	
	int interp_type, int Pmax)	
	(Optional) Sets AMG parameters for B_{Π}	194
6.7.17	int	
V. L. L.	1110	

Parallel auxiliary space divergence solver and preconditioner

__ 6.7.1 __

int HYPRE_ADSCreate (HYPRE_Solver* solver)

Create an ADS solver object

_ 6.7.2 _

int HYPRE_ADSDestroy (HYPRE_Solver solver)

Destroy an ADS solver object

 $_$ 6.7.3 $_$

int

HYPRE_ADSSetup (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x)

Set up the ADS solver or preconditioner. If used as a preconditioner, this function should be passed to the iterative solver SetPrecond function.

Parameters:

solver [IN] object to be set up.

A [IN] ParCSR matrix used to construct the solver/preconditioner.

b Ignored by this function.

x Ignored by this function.

HYPRE_ADSSolve (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x)

Solve the system or apply ADS as a preconditioner. If used as a preconditioner, this function should be passed to the iterative solver SetPrecond function.

Parameters:

solver [IN] solver or preconditioner object to be applied.

A [IN] ParCSR matrix, matrix of the linear system to be solved

b [IN] right hand side of the linear system to be solved

x [OUT] approximated solution of the linear system to be solved

6.7.5

int **HYPRE_ADSSetDiscreteCurl** (HYPRE_Solver solver, HYPRE_ParCSRMatrix C)

Sets the discrete curl matrix C. This function should be called before HYPRE_ADSSetup()!

__ 6.7.6 _____

HYPRE_ADSSetDiscreteGradient (HYPRE_Solver solver, HYPRE_ParCSRMatrix G)

Sets the discrete gradient matrix G. This function should be called before HYPRE_ADSSetup()!

int **HYPRE_ADSSetCoordinateVectors** (HYPRE_Solver solver,
HYPRE_ParVector x, HYPRE_ParVector y, HYPRE_ParVector z)

Sets the x, y and z coordinates of the vertices in the mesh. This function should be called before $HYPRE_ADSSetup()!$

 $_{-}$ 6.7.8

int
HYPRE_ADSSetInterpolations (HYPRE_Solver solver, HYPRE_ParCSRMatrix RT_Pi, HYPRE_ParCSRMatrix RT_Pix, HYPRE_ParCSRMatrix RT_Piy,
HYPRE_ParCSRMatrix RT_Piz, HYPRE_ParCSRMatrix ND_Pi,
HYPRE_ParCSRMatrix ND_Pix, HYPRE_ParCSRMatrix ND_Piy,
HYPRE_ParCSRMatrix ND_Piz)

(Optional) Set the (components of) the Raviart-Thomas (Π_{RT}) and the Nedelec (Π_{ND}) interpolation matrices.

This function is generally intended to be used only for high-order H(div) discretizations (in the lowest order case, these matrices are constructed internally in ADS from the discreet gradient and curl matrices and the coordinates of the vertices), though it can also be used in the lowest-order case or for other types of discretizations.

By definition, RT_Pi and ND_Pi are the matrix representations of the linear operators Π_{RT} and Π_{ND} that interpolate (high-order) vector nodal finite elements into the (high-order) Raviart-Thomas and Nedelec spaces. The component matrices are defined in both cases as $\Pi^x \varphi = \Pi(\varphi, 0, 0)$ and similarly for Π^y and Π^z . Note that all these operators depend on the choice of the basis and degrees of freedom in the high-order spaces.

The column numbering of RT_Pi and ND_Pi should be node-based, i.e. the x/y/z components of the first node (vertex or high-order dof) should be listed first, followed by the x/y/z components of the second node and so on (see the documentation of HYPRE_BoomerAMGSetDofFunc).

If used, this function should be called before hypre_ADSSetup() and there is no need to provide the vertex coordinates. Furthermore, only one of the sets $\{\Pi_{RT}\}$ and $\{\Pi_{RT}^x, \Pi_{RT}^y, \Pi_{RT}^z\}$ needs to be specified (though it is OK to provide both). If RT_Pix is NULL, then scalar Π -based ADS cycles, i.e. those with cycle_type > 10, will be unavailable. Similarly, ADS cycles based on monolithic Π (cycle_type < 10) require that RT_Pi is not NULL. The same restrictions hold for the sets $\{\Pi_{ND}\}$ and $\{\Pi_{ND}^x, \Pi_{ND}^y, \Pi_{ND}^z\}$ – only one of them needs to be specified, and the availability of each enables different AMS cycle type options.

int HYPRE_ADSSetMaxIter (HYPRE_Solver solver, int maxit)

(Optional) Sets maximum number of iterations, if ADS is used as a solver. To use ADS as a preconditioner, set the maximum number of iterations to 1. The default is 20.

_ 6.7.10 _

int HYPRE_ADSSetTol (HYPRE_Solver solver, HYPRE_Real tol)

(Optional) Set the convergence tolerance, if ADS is used as a solver. When using ADS as a preconditioner, set the tolerance to 0.0. The default is 10^{-6} .

6.7.11

int HYPRE_ADSSetCycleType (HYPRE_Solver solver, int cycle_type)

(Optional) Choose which auxiliary-space solver to use. Possible values are:

- 1 3-level multiplicative solver (01210)
- 2 | 3-level additive solver (0+1+2)
- 3 | 3-level multiplicative solver (02120)
- 4 3-level additive solver (010+2)
- 5 | 3-level multiplicative solver (0102010)
- 6 3-level additive solver (1+020)
- 7 | 3-level multiplicative solver (0201020)
- 8 3-level additive solver (0(1+2)0)
- 11 | 5-level multiplicative solver (013454310)
- 12 | 5-level additive solver (0+1+3+4+5)
- 13 | 5-level multiplicative solver (034515430)
- 14 | 5-level additive solver (01(3+4+5)10)

The default is 1. See the user's manual for more details.

int HYPRE_ADSSetPrintLevel (HYPRE_Solver solver, int print_level)

(Optional) Control how much information is printed during the solution iterations. The default is 1 (print residual norm at each step).

 $_{-}$ 6.7.13 $_{-}$

int

HYPRE_ADSSetSmoothingOptions (HYPRE_Solver solver, int relax_type, int relax_times, HYPRE_Real relax_weight, HYPRE_Real omega)

(Optional) Sets relaxation parameters for A. The defaults are 2, 1, 1.0, 1.0.

The available options for relax_type are:

- 1 ℓ_1 -scaled Jacobi
- 2 ℓ_1 -scaled block symmetric Gauss-Seidel/SSOR
- 3 Kaczmarz
- 4 truncated version of ℓ_1 -scaled block symmetric Gauss-Seidel/SSOR
- 16 Chebyshev

_ 6.7.14 _

int

HYPRE_ADSSetChebySmoothingOptions (HYPRE_Solver solver, int cheby_order, int cheby_fraction)

(Optional) Sets parameters for Chebyshev relaxation. The defaults are 2, 0.3.

6.7.15

int

HYPRE_ADSSetAMSOptions (HYPRE_Solver solver, int cycle_type, int coarsen_type, int agg_levels, int relax_type, HYPRE_Real strength_threshold, int interp_type, int Pmax)

(Optional) Sets AMS parameters for B_C . The defaults are 11, 10, 1, 3, 0.25, 0, 0. Note that cycle_type should be greater than 10, unless the high-order interface of HYPRE_ADSSetInterpolations is being used! See the user's manual for more details.

6.7.16

int **HYPRE_ADSSetAMGOptions** (HYPRE_Solver solver, int coarsen_type, int agg_levels, int relax_type, HYPRE_Real strength_threshold, int interp_type, int Pmax)

(Optional) Sets AMG parameters for B_{Π} . The defaults are 10, 1, 3, 0.25, 0, 0. See the user's manual for more details.

6.7.17

HYPRE_ADSGetNumIterations (HYPRE_Solver solver, int* num_iterations)

Returns the number of iterations taken

6.7.18

HYPRE_ADSGetFinalRelativeResidualNorm (HYPRE_Solver solver, HYPRE_Real* rel_resid_norm)

Returns the norm of the final relative residual

_ 6.8 _____

ParCSR PCG Solver

Names

6.8.1 int

	HYPRE_ParCSRPCGCreate (MPI_Comm comm, HYPRE_Solver* solver)	
	Create a solver object	195
6.8.2	int	
	HYPRE_ParCSRPCGDestroy (HYPRE_Solver solver)	
	Destroy a solver object	195
6.8.3	int	
	HYPRE_ParCSRDiagScaleSetup (HYPRE_Solver solver,	
	HYPRE_ParCSRMatrix A,	
	HYPRE_ParVector y,	
	HYPRE_ParVector x)	
	Setup routine for diagonal preconditioning	195
6.8.4	int	
	HYPRE_ParCSRDiagScale (HYPRE_Solver solver,	
	HYPRE_ParCSRMatrix HA,	
	HYPRE_ParVector Hy, HYPRE_ParVector Hx)	
	Solve routine for diagonal preconditioning	196

These routines should be used in conjunction with the generic interface in PCG Solver.

__ 6.8.1 __

 $int \ \mathbf{HYPRE_ParCSRPCGCreate} \ (MPI_Comm \ comm, \ HYPRE_Solver* \ solver)$

Create a solver object

_ 6.8.2 _

int HYPRE_ParCSRPCGDestroy (HYPRE_Solver solver)

Destroy a solver object

_ 6.8.3 _

int
HYPRE_ParCSRDiagScaleSetup (HYPRE_Solver solver,
HYPRE_ParCSRMatrix A, HYPRE_ParVector y, HYPRE_ParVector x)

Setup routine for diagonal preconditioning

6.8.4

HYPRE_ParCSRDiagScale (HYPRE_Solver solver, HYPRE_ParCSRMatrix HA, HYPRE_ParVector Hy, HYPRE_ParVector Hx)

Solve routine for diagonal preconditioning

_ 6.9 _

ParCSR GMRES Solver

Names

These routines should be used in conjunction with the generic interface in GMRES Solver.

_ 6.9.1 _

int
HYPRE_ParCSRGMRESCreate (MPI_Comm comm, HYPRE_Solver* solver)

Create a solver object

6.9.2

int HYPRE_ParCSRGMRESDestroy (HYPRE_Solver solver)

Destroy a solver object

___ 6.10 _____

ParCSR FlexGMRES Solver

Names

These routines should be used in conjunction with the generic interface in FlexGMRES Solver.

 $_{-}$ 6.10.1

HYPRE_ParCSRFlexGMRESCreate (MPI_Comm comm, HYPRE_Solver* solver)

Create a solver object

6.10.2

int HYPRE_ParCSRFlexGMRESDestroy (HYPRE_Solver solver)

Destroy a solver object

ParCSR LGMRES Solver

Names

6.11.1 int

HYPRE_ParCSRLGMRESCreate (MPI_Comm comm,

HYPRE_Solver* solver)

198

198

6.11.2int

HYPRE_ParCSRLGMRESDestroy (HYPRE_Solver solver)

Destroy a solver object

Create a solver object

These routines should be used in conjunction with the generic interface in LGMRES Solver.

__ 6.11.1 __

HYPRE_ParCSRLGMRESCreate (MPI_Comm comm, HYPRE_Solver* solver)

Create a solver object

-6.11.2 _____

int HYPRE_ParCSRLGMRESDestroy (HYPRE_Solver solver)

Destroy a solver object

-6.12

ParCSR BiCGSTAB Solver

Names

6.12.1

HYPRE_ParCSRBiCGSTABCreate (MPI_Comm comm, HYPRE_Solver* solver)

Create a solver object 199

6.12.2int

These routines should be used in conjunction with the generic interface in BiCGSTAB Solver.

_ 6.12.1 __

HYPRE_ParCSRBiCGSTABCreate (MPI_Comm comm, HYPRE_Solver* solver)

Create a solver object

6.12.2

int HYPRE_ParCSRBiCGSTABDestroy (HYPRE_Solver solver)

Destroy a solver object

_ 6.13 __

ParCSR Hybrid Solver

Names		
6.13.1	int	
	HYPRE_ParCSRHybridCreate (HYPRE_Solver* solver)	
	Create solver object	203
6.13.2	int	
	HYPRE_ParCSRHybridDestroy (HYPRE_Solver solver)	
	Destroy solver object	204
6.13.3	int	
	HYPRE_ParCSRHybridSetup (HYPRE_Solver solver,	
	HYPRE_ParCSRMatrix A,	
	HYPRE_ParVector b, HYPRE_ParVector x)	
	Setup the hybrid solver	204
6.13.4	int	

	HYPRE_ParCSRHybridSolve (HYPRE_Solver solver,
	HYPRE_ParCSRMatrix A,
	HYPRE_ParVector b, HYPRE_ParVector x)
	Solve linear system
6.13.5	int
	HYPRE_ParCSRHybridSetTol (HYPRE_Solver solver, HYPRE_Real tol)
	Set the convergence tolerance for the Krylov solver
6.13.6	int
0.10.0	HYPRE_ParCSRHybridSetAbsoluteTol (HYPRE_Solver solver,
	HYPRE_Real tol)
	Set the absolute convergence tolerance for the Krylov solver
6.13.7	int
0.13.7	HYPRE_ParCSRHybridSetConvergenceTol (HYPRE_Solver solver,
	HYPRE_Real cf_tol)
	Set the desired convergence factor
0.10.0	
6.13.8	int
	HYPRE_ParCSRHybridSetDSCGMaxIter (HYPRE_Solver solver, int dscg_max_its)
	Set the maximal number of iterations for the diagonally preconditioned solver
	See the maximum number of heracions for the anagonality preconditioned solver
C 19 O	
6.13.9	int HYPRE_ParCSRHybridSetPCGMaxIter (HYPRE_Solver solver,
	int pcg_max_its)
	Set the maximal number of iterations for the AMG preconditioned solver .
	* * * * * * * * * * * * * * * * * * * *
6.13.10	int
	HYPRE_ParCSRHybridSetSolverType (HYPRE_Solver solver,
	int solver_type)
	Set the desired solver type.
6.13.11	int
	HYPRE_ParCSRHybridSetKDim (HYPRE_Solver solver, int k_dim)
	Set the Krylov dimension for restarted GMRES
6.13.12	int
	HYPRE_ParCSRHybridSetTwoNorm (HYPRE_Solver solver, int two_norm)
	Set the type of norm for PCG
6.13.13	int
	HYPRE_ParCSRHybridSetPrecond (HYPRE_Solver solver,
	HYPRE_PtrToParSolverFcn precond,
	HYPRE_PtrToParSolverFcn
	$\operatorname{precond_setup},$
	HYPRE_Solver precond_solver)
	Set preconditioner if wanting to use one that is not set up by the hybrid
	solver
6.13.14	int
	HYPRE_ParCSRHybridSetLogging (HYPRE_Solver solver, int logging)
	Set logging parameter (default: 0, no logging)
6.13.15	int
0.10.10	1110

	HYPRE_ParCSRHybridSetPrintLevel (HYPRE_Solver solver,	
	int print_level)	207
	Set print level (default: 0, no printing)	207
6.13.16	int	
	HYPRE_ParCSRHybridSetStrongThreshold (HYPRE_Solver solver, HYPRE_Real	
	strong_threshold)	
	(Optional) Sets AMG strength threshold.	207
		201
6.13.17	int	
	HYPRE_ParCSRHybridSetMaxRowSum (HYPRE_Solver solver,	
	HYPRE_Real max_row_sum)	
	(Optional) Sets a parameter to modify the definition of strength for diagonal dominant portions of the matrix.	208
		200
6.13.18	int	
	HYPRE_ParCSRHybridSetTruncFactor (HYPRE_Solver solver,	
	HYPRE_Real trunc_factor)	200
	(Optional) Defines a truncation factor for the interpolation	208
6.13.19	int	
	HYPRE_ParCSRHybridSetPMaxElmts (HYPRE_Solver solver,	
	int P_max_elmts)	
	(Optional) Defines the maximal number of elements per row for the inter-	200
	polation.	208
6.13.20	int	
	HYPRE_ParCSRHybridSetMaxLevels (HYPRE_Solver solver,	
	int max_levels)	
	(Optional) Defines the maximal number of levels used for AMG	208
6.13.21	int	
	$\mathbf{HYPRE_ParCSRHybridSetMeasureType} \ (\mathbf{HYPRE_Solver} \ \mathbf{solver},$	
	int measure_type)	
	(Optional) Defines whether local or global measures are used	209
6.13.22	int	
	HYPRE_ParCSRHybridSetCoarsenType (HYPRE_Solver solver,	
	int coarsen_type)	
	(Optional) Defines which parallel coarsening algorithm is used	209
6.13.23	int	
	HYPRE_ParCSRHybridSetCycleType (HYPRE_Solver solver,	
	int cycle_type)	
	(Optional) Defines the type of cycle	209
6.13.24	int	
0.10.21	HYPRE_ParCSRHybridSetNumSweeps (HYPRE_Solver solver,	
	int num_sweeps)	
	(Optional) Sets the number of sweeps	210
6.13.25	int	
0.13.20	HYPRE_ParCSRHybridSetCycleNumSweeps (HYPRE_Solver solver,	
	int num_sweeps, int k)	
	(Optional) Sets the number of sweeps at a specified cycle	210
0.10.00		210
6.13.26	int	

	HYPRE_ParCSRHybridSetRelaxType (HYPRE_Solver solver, int relax_type)	
	(Optional) Defines the smoother to be used	210
6.13.27	int HYPRE_ParCSRHybridSetCycleRelaxType (HYPRE_Solver solver,	
	int relax_type, int k)	01:
£ 12 20	(Optional) Defines the smoother at a given cycle	21
6.13.28	int HYPRE_ParCSRHybridSetRelaxOrder (HYPRE_Solver solver,	
	int relax_order)	
	(Optional) Defines in which order the points are relaxed	21
6.13.29	int	
	HYPRE_ParCSRHybridSetRelaxWt (HYPRE_Solver solver, HYPRE_Real relax_wt)	
	(Optional) Defines the relaxation weight for smoothed Jacobi and hybrid	
	SOR on all levels.	21
6.13.30	int	
	HYPRE_ParCSRHybridSetLevelRelaxWt (HYPRE_Solver solver,	
	HYPRE_Real relax_wt, int level)	
	(Optional) Defines the relaxation weight for smoothed Jacobi and hybrid	
	SOR on the user defined level.	21
6.13.31	int	
	HYPRE_ParCSRHybridSetOuterWt (HYPRE_Solver solver, HYPRE_Real outer_wt)	
	(Optional) Defines the outer relaxation weight for hybrid SOR and SSOR	
	on all levels.	213
6.13.32	int	
	HYPRE_ParCSRHybridSetLevelOuterWt (HYPRE_Solver solver,	
	HYPRE_Real outer_wt, int level)	
	(Optional) Defines the outer relaxation weight for hybrid SOR or SSOR on	
	the user defined level.	213
6.13.33	int	
	HYPRE_ParCSRHybridSetMaxCoarseSize (HYPRE_Solver solver,	
	int max_coarse_size) (Optional) Defines the maximal coarse grid size	213
6.13.34	int	
0.10.04	HYPRE_ParCSRHybridSetMinCoarseSize (HYPRE_Solver solver,	
	int min_coarse_size)	
	(Optional) Defines the minimal coarse grid size	213
6.13.35	int	
	HYPRE_ParCSRHybridSetSeqThreshold (HYPRE_Solver solver, int seq_threshold)	
	(Optional) enables redundant coarse grid size	213
6.13.36	int	

	HYPRE_ParCSRHybridSetAggNumLevels (HYPRE_Solver solver,
	int agg_num_levels)
	(Optional) Defines the number of levels of aggressive coarsening, starting with the finest level.
6.13.37	int
	${\bf HYPRE_ParCSRHybridSetNumPaths} \ ({\bf HYPRE_Solver} \ solver,$
	int num_paths)
	(Optional) Defines the degree of aggressive coarsening
6.13.38	int
	${\bf HYPRE_ParCSRHybridSetNumFunctions}~({\bf HYPRE_Solver}~solver,$
	int num_functions)
	(Optional) Sets the size of the system of PDEs, if using the systems version.
6.13.39	int
	HYPRE_ParCSRHybridSetDofFunc (HYPRE_Solver solver, int* dof_func)
	(Optional) Sets the mapping that assigns the function to each variable, if
	using the systems version.
6.13.40	int
	HYPRE_ParCSRHybridSetNodal (HYPRE_Solver solver, int nodal)
	(Optional) Sets whether to use the nodal systems version
6.13.41	int
	HYPRE_ParCSRHybridGetNumIterations (HYPRE_Solver solver,
	int* num_its)
	Retrieves the total number of iterations
6.13.42	int
	HYPRE_ParCSRHybridGetDSCGNumIterations (HYPRE_Solver solver,
	$\operatorname{int}^*\operatorname{dscg_num_its})$ Retrieves the number of iterations used by the diagonally scaled solver
	Retrieves the number of tierations used by the diagonally scaled solver
6.13.43	int
	HYPRE_ParCSRHybridGetPCGNumIterations (HYPRE_Solver solver,
	int* pcg_num_its) Retrieves the number of iterations used by the AMG preconditioned solver
	v · ·
6.13.44	int
	HYPRE_ParCSRHybridGetFinalRelativeResidualNorm (HYPRE_Solver
	solver, HYPRE_Real*
	norm)
	Retrieves the final relative residual norm

_ 6.13.1 _

 ${\rm int}\; {\bf HYPRE_ParCSRHybridCreate}\; ({\rm HYPRE_Solver*\; solver})$

Create solver object

6.13.2

int HYPRE_ParCSRHybridDestroy (HYPRE_Solver solver)

Destroy solver object

6.13.3 _

int

HYPRE_ParCSRHybridSetup (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x)

Setup the hybrid solver

Parameters: solver [IN] object to be set up.

A [IN] ParCSR matrix used to construct the

solver/preconditioner.

b Ignored by this function.

x Ignored by this function.

6.13.4

int

HYPRE_ParCSRHybridSolve (HYPRE_Solver solver, HYPRE_ParCSRMatrix A, HYPRE_ParVector b, HYPRE_ParVector x)

Solve linear system

Parameters: solver [IN] solver or preconditioner object to be applied.

A [IN] ParCSR matrix, matrix of the linear system to be

 ${\rm solved}$

b [IN] right hand side of the linear system to be solved

x [OUT] approximated solution of the linear system to

be solved

int HYPRE_ParCSRHybridSetTol (HYPRE_Solver solver, HYPRE_Real tol)

Set the convergence tolerance for the Krylov solver. The default is 1.e-7.

__ 6.13.6 _____

HYPRE_ParCSRHybridSetAbsoluteTol (HYPRE_Solver solver, HYPRE_Real tol)

Set the absolute convergence tolerance for the Krylov solver. The default is 0.

_ 6.13.7 ___

int

 $\label{lem:hypre_parcsrhybridSetConvergenceTol} \begin{tabular}{l} HYPRE_ParcsRHybridSetConvergenceTol \end{tabular} \begin{tabular}{l} (HYPRE_Solver solver, HYPRE_Real cf_tol) \end{tabular}$

Set the desired convergence factor

6.13.8

HYPRE_ParCSRHybridSetDSCGMaxIter (HYPRE_Solver solver, int dscg_max_its)

Set the maximal number of iterations for the diagonally preconditioned solver

HYPRE_ParCSRHybridSetPCGMaxIter (HYPRE_Solver solver, int pcg_max_its)

Set the maximal number of iterations for the AMG preconditioned solver

___ 6.13.10 _____

HYPRE_ParCSRHybridSetSolverType (HYPRE_Solver solver, int solver_type)

Set the desired solver type. There are the following options: 2 GMRES

1 PCG (default)

3 BiCGSTAB

_ 6.13.11 __

int HYPRE_ParCSRHybridSetKDim (HYPRE_Solver solver, int k_dim)

Set the Krylov dimension for restarted GMRES. The default is 5.

HYPRE_ParCSRHybridSetTwoNorm (HYPRE_Solver solver, int two_norm)

Set the type of norm for PCG

int

HYPRE_ParCSRHybridSetPrecond (HYPRE_Solver solver, HYPRE_PtrToParSolverFcn precond, HYPRE_PtrToParSolverFcn precond_setup, HYPRE_Solver precond_solver)

Set preconditioner if wanting to use one that is not set up by the hybrid solver

_ 6.13.14 _

int HYPRE_ParCSRHybridSetLogging (HYPRE_Solver solver, int logging)

Set logging parameter (default: 0, no logging)

_ 6.13.15 ____

int

HYPRE_ParCSRHybridSetPrintLevel (HYPRE_Solver solver, int print_level)

Set print level (default: 0, no printing)

6.13.16

HYPRE_ParCSRHybridSetStrongThreshold (HYPRE_Solver solver, HYPRE_Real strong_threshold)

(Optional) Sets AMG strength threshold. The default is 0.25. For 2d Laplace operators, 0.25 is a good value, for 3d Laplace operators, 0.5 or 0.6 is a better value. For elasticity problems, a large strength threshold, such as 0.9, is often better.

HYPRE_ParCSRHybridSetMaxRowSum (HYPRE_Solver solver, HYPRE_Real max_row_sum)

(Optional) Sets a parameter to modify the definition of strength for diagonal dominant portions of the matrix. The default is 0.9. If max_row_sum is 1, no checking for diagonally dominant rows is performed.

6.13.18

HYPRE_ParCSRHybridSetTruncFactor (HYPRE_Solver solver, HYPRE_Real trunc_factor)

(Optional) Defines a truncation factor for the interpolation. The default is 0.

__ 6.13.19 __

int **HYPRE_ParCSRHybridSetPMaxElmts** (HYPRE_Solver solver, int P_max_elmts)

(Optional) Defines the maximal number of elements per row for the interpolation. The default is 0.

_ 6.13.20 ___

HYPRE_ParCSRHybridSetMaxLevels (HYPRE_Solver solver, int max_levels)

(Optional) Defines the maximal number of levels used for AMG. The default is 25.

int HYPRE_ParCSRHybridSetMeasureType (HYPRE_Solver solver, int measure_type)

(Optional) Defines whether local or global measures are used

 $_{-}$ 6.13.22 $_{-}$

HYPRE_ParCSRHybridSetCoarsenType (HYPRE_Solver solver, int coarsen_type)

(Optional) Defines which parallel coarsening algorithm is used. There are the following options for coarsen_type:

- 0 | CLJP-coarsening (a parallel coarsening algorithm using independent sets).
- 1 classical Ruge-Stueben coarsening on each processor, no boundary treatment
- 3 classical Ruge-Stueben coarsening on each processor, followed by a third pass, which adds coarse points on the boundaries
- Falgout coarsening (uses 1 first, followed by CLJP using the interior coarse points generated by 1 as its first independent set)
- 7 CLJP-coarsening (using a fixed random vector, for debugging purposes only)
- 8 PMIS-coarsening (a parallel coarsening algorithm using independent sets with lower complexities than CLJP, might also lead to slower convergence)
- 9 PMIS-coarsening (using a fixed random vector, for debugging purposes only)
- HMIS-coarsening (uses one pass Ruge-Stueben on each processor independently, followed by PMIS using the interior C-points as its first independent set)
- 11 one-pass Ruge-Stueben coarsening on each processor, no boundary treatment

The default is 6.

6.13.23

HYPRE_ParCSRHybridSetCycleType (HYPRE_Solver solver, int cycle_type)

(Optional) Defines the type of cycle. For a V-cycle, set cycle_type to 1, for a W-cycle set cycle_type to 2. The default is 1.

int

HYPRE_ParCSRHybridSetNumSweeps (HYPRE_Solver solver, int num_sweeps)

(Optional) Sets the number of sweeps. On the finest level, the up and the down cycle the number of sweeps are set to num_sweeps and on the coarsest level to 1. The default is 1.

 $_$ 6.13.25 $_$

int HYPRE_ParCSRHybridSetCycleNumSweeps (HYPRE_Solver solver, int num_sweeps, int k)

(Optional) Sets the number of sweeps at a specified cycle. There are the following options for k:

the down cycle	if k=1
the up cycle	if $k=2$
the coarsest level	if $k=3$.

HYPRE_ParCSRHybridSetRelaxType (HYPRE_Solver solver, int relax_type)

(Optional) Defines the smoother to be used. It uses the given smoother on the fine grid, the up and the down cycle and sets the solver on the coarsest level to Gaussian elimination (9). The default is Gauss-Seidel (3).

There are the following options for relax_type:

- 0 Jacobi Gauss-Seidel, sequential (very slow!) Gauss-Seidel, interior points in parallel, boundary sequential (slow!) hybrid Gauss-Seidel or SOR, forward solve 3 hybrid Gauss-Seidel or SOR, backward solve 4
- hybrid chaotic Gauss-Seidel (works only with OpenMP)
- hybrid symmetric Gauss-Seidel or SSOR
- Gaussian elimination (only on coarsest level)

int

 $\label{eq:hypre_parcsr} \textbf{HYPRE_ParCSRHybridSetCycleRelaxType} \ (\textbf{HYPRE_Solver solver}, \ \textbf{int} \\ \textbf{relax_type}, \ \textbf{int} \ \textbf{k})$

(Optional) Defines the smoother at a given cycle. For options of relax_type see description of HYPRE_BoomerAMGSetRelaxType). Options for k are

the down cycle	if k=1
the up cycle	if $k=2$
the coarsest level	if k=3.

_ 6.13.28 _

int

HYPRE_ParCSRHybridSetRelaxOrder (HYPRE_Solver solver, int relax_order)

(Optional) Defines in which order the points are relaxed. There are the following options for relax_order:

- 0 the points are relaxed in natural or lexicographic order on each processor
- 1 CF-relaxation is used, i.e on the fine grid and the down cycle the coarse points are relaxed first, followed by the fine points; on the up cycle the F-points are relaxed first, followed by the C-points. On the coarsest level, if an iterative scheme is used, the points are relaxed in lexicographic order.

The default is 1 (CF-relaxation).

6.13.29

HYPRE_ParCSRHybridSetRelaxWt (HYPRE_Solver solver, HYPRE_Real relax_wt)

(Optional) Defines the relaxation weight for smoothed Jacobi and hybrid SOR on all levels.

	this assigns the given relaxation weight on all levels
$relax_weight = 0$	the weight is determined on each level with the estimate $\frac{3}{4\ D^{-1/2}AD^{-1/2}\ }$,
	where D is the diagonal matrix of A (this should only be used with Jacobi)
$relax_weight = -k$	the relaxation weight is determined with at most k CG steps on each level
	this should only be used for symmetric positive definite problems)

The default is 1.

6.13.30

int

HYPRE_ParCSRHybridSetLevelRelaxWt (HYPRE_Solver solver, HYPRE_Real relax_wt, int level)

(Optional) Defines the relaxation weight for smoothed Jacobi and hybrid SOR on the user defined level. Note that the finest level is denoted 0, the next coarser level 1, etc. For nonpositive relax_weight, the parameter is determined on the given level as described for HYPRE_BoomerAMGSetRelaxWt. The default is 1.

6.13.31

HYPRE_ParCSRHybridSetOuterWt (HYPRE_Solver solver, HYPRE_Real outer_wt)

(Optional) Defines the outer relaxation weight for hybrid SOR and SSOR on all levels.

omega > 0	this assigns the same outer relaxation weight omega on each level
omega = -k	an outer relaxation weight is determined with at most k CG steps on each level
	(this only makes sense for symmetric positive definite problems and smoothers, e.g. SSOR)

The default is 1.

6.13.32

HYPRE_ParCSRHybridSetLevelOuterWt (HYPRE_Solver solver, HYPRE_Real outer_wt, int level)

(Optional) Defines the outer relaxation weight for hybrid SOR or SSOR on the user defined level. Note that the finest level is denoted 0, the next coarser level 1, etc. For nonpositive omega, the parameter is determined on the given level as described for HYPRE_BoomerAMGSetOuterWt. The default is 1.

int

 $\label{lem:hypre_parcsr} \mathbf{HYPRE_ParcSRHybridSetMaxCoarseSize} \ (\\ \mathbf{HYPRE_Solver} \ solver, \ int \\ \mathbf{max_coarse_size})$

(Optional) Defines the maximal coarse grid size. The default is 9.

6.13.34

int

 $\label{lem:hypre-parcsrhybridSetMinCoarseSize} \ (\mbox{HYPRE_Solver solver, int min_coarse_size})$

(Optional) Defines the minimal coarse grid size. The default is 0.

__ 6.13.35 ___

ınt

HYPRE_ParCSRHybridSetSeqThreshold (HYPRE_Solver solver, int seq_threshold)

(Optional) enables redundant coarse grid size. If the system size becomes smaller than seq_threshold, sequential AMG is used on all remaining processors. The default is 0.

6.13.36

ınt

HYPRE_ParCSRHybridSetAggNumLevels (HYPRE_Solver solver, int agg_num_levels)

(Optional) Defines the number of levels of aggressive coarsening, starting with the finest level. The default is 0, i.e. no aggressive coarsening.

HYPRE_ParCSRHybridSetNumPaths (HYPRE_Solver solver, int num_paths)

(Optional) Defines the degree of aggressive coarsening. The default is 1, which leads to the most aggressive coarsening. Setting num_paths to 2 will increase complexity somewhat, but can lead to better convergence.*

_ 6.13.38 _

int **HYPRE_ParCSRHybridSetNumFunctions** (HYPRE_Solver solver, int num_functions)

(Optional) Sets the size of the system of PDEs, if using the systems version. The default is 1.

 $_{-}$ 6.13.39 $_{-}$

int HYPRE_ParCSRHybridSetDofFunc (HYPRE_Solver solver, int* dof_func)

(Optional) Sets the mapping that assigns the function to each variable, if using the systems version. If no assignment is made and the number of functions is k > 1, the mapping generated is (0,1,...,k-1,0,1,...,k-1,...).

6.13.40

int HYPRE_ParCSRHybridSetNodal (HYPRE_Solver solver, int nodal)

(Optional) Sets whether to use the nodal systems version. The default is 0 (the unknown based approach).

HYPRE_ParCSRHybridGetNumIterations (HYPRE_Solver solver, int* num_its)

Retrieves the total number of iterations

6.13.42

HYPRE_ParCSRHybridGetDSCGNumIterations (HYPRE_Solver solver, int* dscg_num_its)

Retrieves the number of iterations used by the diagonally scaled solver

__ 6.13.43 _____

HYPRE_ParCSRHybridGetPCGNumIterations (HYPRE_Solver solver, int* pcg_num_its)

Retrieves the number of iterations used by the AMG preconditioned solver

__ 6.13.44 _____

HYPRE_ParCSRHybridGetFinalRelativeResidualNorm (HYPRE_Solver solver, HYPRE_Real* norm)

Retrieves the final relative residual norm

6 14

ParCSR LOBPCG Eigensolver

These routines should be used in conjunction with the generic interface in LOBPCG Eigensolver.

_ 6.14.1 _

int HYPRE_ParCSRSetupInterpreter (mv_InterfaceInterpreter* i)

 $\label{loaded} Load interface interpreter. \ Vector \ part\ loaded\ with\ hypre_ParKrylov\ functions\ and\ multivector\ part\ loaded\ with\ mv_TempMultiVector\ functions.$

_ 6.14.2 _

int HYPRE_ParCSRSetupMatvec (HYPRE_MatvecFunctions* mv)

 ${\bf Load\ Matvec\ interpreter\ with\ hypre_ParKrylov\ functions}$

7

Krylov Solvers

Names		
7.1	Krylov Solvers	
		217
7.2	PCG Solver	
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7.5	LGMRES Solver	
		239
7.6	BiCGSTAB Solver	
		245
7.7	CGNR Solver	
		249

These solvers support many of the matrix/vector storage schemes in hypre. They should be used in conjunction with the storage-specific interfaces, particularly the specific Create() and Destroy() functions.

7.1

Krylov Solvers

Names		
7.1.1	typedef struct hypre_Solver_struct *HYPRE_Solver The solver object	21
7.1.2	typedef struct hypre_Matrix_struct *HYPRE_Matrix The matrix object	21
7.1.3	typedef struct hypre_Vector_struct *HYPRE_Vector The vector object	21

__ 7.1.1 _____

 $typedef\ struct\ hypre_Solver_struct\ \textbf{*HYPRE_Solver}$

The solver object

7.1.2

 $typedef\ struct\ \ hypre_Matrix_struct\ \ \textbf{*HYPRE_Matrix}$

The matrix object

7.1.3

 $typedef \ struct \ hypre_Vector_struct \ *HYPRE_Vector$

The vector object

_ 7.2 _

PCG Solver

Names

7.2.1	int HYPRE_PCGSetup (HYPRE_Solver solver, HYPRE_Matrix A, HYPRE_Vector b, HYPRE_Vector x)	
	Prepare to solve the system.	220
7.2.2	int HYPRE_PCGSolve (HYPRE_Solver solver, HYPRE_Matrix A, HYPRE_Vector b, HYPRE_Vector x)	222
	Solve the system	220
7.2.3	int HYPRE_PCGSetTol (HYPRE_Solver solver, HYPRE_Real tol) (Optional) Set the relative convergence tolerance	221
7.2.4	int	

	HYPRE_PCGSetAbsoluteTol (HYPRE_Solver solver, HYPRE_Real a_tol) (Optional) Set the absolute convergence tolerance (default is 0)	
7.2.5	int	
	HYPRE_PCGSetResidualTol (HYPRE_Solver solver, HYPRE_Real rtol) (Optional) Set a residual-based convergence tolerance which checks if $ r_{old} - r_{new} < rtol b $.	
7.2.6	int	
	HYPRE_PCGSetMaxIter (HYPRE_Solver solver, int max_iter) (Optional) Set maximum number of iterations	
7.2.7	int HYPRE_PCGSetTwoNorm (HYPRE_Solver solver, int two_norm) (Optional) Use the two-norm in stopping criteria	
7.2.8	int	
1.2.0	HYPRE_PCGSetRelChange (HYPRE_Solver solver, int rel_change) (Optional) Additionally require that the relative difference in successive iterates be small	
7.2.9	int HYPRE_PCGSetRecomputeResidual (HYPRE_Solver solver,	
	int recompute_residual) (Optional) Recompute the residual at the end to double-check convergence	
7.2.10	int	
7.2.10	HYPRE_PCGSetRecomputeResidualP (HYPRE_Solver solver, int recompute_residual_p)	
	(Optional) Periodically recompute the residual while iterating	
7.2.11	int HYPRE_PCGSetPrecond (HYPRE_Solver solver, HYPRE_PtrToSolverFcn precond, HYPRE_PtrToSolverFcn precond_setup, HYPRE_Solver precond_solver)	
	(Optional) Set the preconditioner to use	
7.2.12	int HYPRE_PCGSetLogging (HYPRE_Solver solver, int logging) (Optional) Set the amount of logging to do	
7.2.13	int HYPRE_PCGSetPrintLevel (HYPRE_Solver solver, int level)	
	(Optional) Set the amount of printing to do to the screen	
7.2.14	int HYPRE_PCGGetNumIterations (HYPRE_Solver solver, int* num_iterations) Return the number of iterations taken	
7.2.15	int HYPRE_PCGGetFinalRelativeResidualNorm (HYPRE_Solver solver, HYPRE_Real* norm)	
	Return the norm of the final relative residual	
7.2.16	int HYPRE_PCGGetResidual (HYPRE_Solver solver, void** residual)	
	Return the residual	
7.2.17	int	

	HYPRE_PCGGetTol (HYPRE_Solver solver, HYPRE_Real* tol)	224
7.2.18	int HYPRE_PCGGetResidualTol (HYPRE_Solver solver, HYPRE_Real* rtol)	224
7.2.19	int HYPRE_PCGGetMaxIter (HYPRE_Solver solver, int* max_iter)	224
7.2.20	int HYPRE_PCGGetTwoNorm (HYPRE_Solver solver, int* two_norm)	225
7.2.21	int HYPRE_PCGGetRelChange (HYPRE_Solver solver, int* rel_change)	225
7.2.22	int HYPRE_GMRESGetSkipRealResidualCheck (HYPRE_Solver solver, int* skip_real_r_check)	225
7.2.23	int HYPRE_PCGGetPrecond (HYPRE_Solver solver, HYPRE_Solver* precond_data_ptr)	225
7.2.24	int HYPRE_PCGGetLogging (HYPRE_Solver solver, int* level)	225
7.2.25	int HYPRE_PCGGetPrintLevel (HYPRE_Solver solver, int* level)	226
7.2.26	int HYPRE_PCGGetConverged (HYPRE_Solver solver, int* converged)	226

7.2.1

int **HYPRE_PCGSetup** (HYPRE_Solver solver, HYPRE_Matrix A, HYPRE_Vector b, HYPRE_Vector x)

Prepare to solve the system. The coefficient data in ${\tt b}$ and ${\tt x}$ is ignored here, but information about the layout of the data may be used.

_ 7.2.2 _

int **HYPRE_PCGSolve** (HYPRE_Solver solver, HYPRE_Matrix A, HYPRE_Vector b, HYPRE_Vector x)

Solve the system

7.2.3

int HYPRE_PCGSetTol (HYPRE_Solver solver, HYPRE_Real tol)

(Optional) Set the relative convergence tolerance

 $_$ 7.2.4 $_$

int HYPRE_PCGSetAbsoluteTol (HYPRE_Solver solver, HYPRE_Real a_tol)

(Optional) Set the absolute convergence tolerance (default is 0). If one desires the convergence test to check the absolute convergence tolerance only, then set the relative convergence tolerance to 0.0. (The default convergence test is $< C*r, r > \le \max(\text{relative_tolerance}^2 * < C*b, b >$, absolute_tolerance²).)

 $_$ 7.2.5 $_$

int HYPRE_PCGSetResidualTol (HYPRE_Solver solver, HYPRE_Real rtol)

(Optional) Set a residual-based convergence tolerance which checks if $||r_{old} - r_{new}|| < rtol ||b||$. This is useful when trying to converge to very low relative and/or absolute tolerances, in order to bail-out before roundoff errors affect the approximation.

_ 7.2.6 _

int HYPRE_PCGSetMaxIter (HYPRE_Solver solver, int max_iter)

(Optional) Set maximum number of iterations

7.2.7

int HYPRE_PCGSetTwoNorm (HYPRE_Solver solver, int two_norm)

(Optional) Use the two-norm in stopping criteria

___ 7.2.8 _____

int HYPRE_PCGSetRelChange (HYPRE_Solver solver, int rel_change)

(Optional) Additionally require that the relative difference in successive iterates be small

___ 7.2.9 _____

HYPRE_PCGSetRecomputeResidual (HYPRE_Solver solver, int recompute_residual)

(Optional) Recompute the residual at the end to double-check convergence

_ 7.2.10 _____

int **HYPRE_PCGSetRecomputeResidualP** (HYPRE_Solver solver, int recompute_residual_p)

(Optional) Periodically recompute the residual while iterating

7 2 11

int

HYPRE_PCGSetPrecond (HYPRE_Solver solver, HYPRE_PtrToSolverFcn precond, HYPRE_PtrToSolverFcn precond_setup, HYPRE_Solver precond_solver)

(Optional) Set the preconditioner to use

___ 7.2.12 _____

int HYPRE_PCGSetLogging (HYPRE_Solver solver, int logging)

(Optional) Set the amount of logging to do

_ 7.2.13 __

int HYPRE_PCGSetPrintLevel (HYPRE_Solver solver, int level)

(Optional) Set the amount of printing to do to the screen

_ 7.2.14 __

HYPRE_PCGGetNumIterations (HYPRE_Solver solver, int* num_iterations)

Return the number of iterations taken

7.2.15

HYPRE_PCGGetFinalRelativeResidualNorm (HYPRE_Solver solver, HYPRE_Real* norm)

Return the norm of the final relative residual

7.2.16

int HYPRE_PCGGetResidual (HYPRE_Solver solver, void** residual)

Return the residual

7.2.17

int HYPRE_PCGGetTol (HYPRE_Solver solver, HYPRE_Real* tol)

7.2.18

int HYPRE_PCGGetResidualTol (HYPRE_Solver solver, HYPRE_Real* rtol)

7.2.19

int HYPRE_PCGGetMaxIter (HYPRE_Solver solver, int* max_iter)

7.2.20

int HYPRE_PCGGetTwoNorm (HYPRE_Solver solver, int* two_norm)

7.2.21

int HYPRE_PCGGetRelChange (HYPRE_Solver solver, int* rel_change)

 $_$ 7.2.22 $_$

 $\label{lem:hypre_gmres} \begin{array}{l} \operatorname{int} \\ \mathbf{HYPRE_GMRESGetSkipRealResidualCheck} \ (\operatorname{HYPRE_Solver} \ \operatorname{solver}, \ \operatorname{int}^* \\ \operatorname{skip_real_r_check}) \end{array}$

__ 7.2.23 _____

int **HYPRE_PCGGetPrecond** (HYPRE_Solver solver, HYPRE_Solver* precond_data_ptr)

7.2.24

int HYPRE_PCGGetLogging (HYPRE_Solver solver, int* level)

7 2 25

int HYPRE_PCGGetPrintLevel (HYPRE_Solver solver, int* level)

_ 7.2.26 _

int **HYPRE_PCGGetConverged** (HYPRE_Solver solver, int* converged)

7.3

GMRES Solver

Names		
7.3.1	int	
	HYPRE_GMRESSetup (HYPRE_Solver solver, HYPRE_Matrix A,	
	HYPRE_Vector b, HYPRE_Vector x)	
	Prepare to solve the system.	228
7.3.2	int	
	HYPRE_GMRESSolve (HYPRE_Solver solver, HYPRE_Matrix A, HYPRE_Vector b, HYPRE_Vector x)	
	Solve the system	228
7.3.3	int	
	HYPRE_GMRESSetTol (HYPRE_Solver solver, HYPRE_Real tol)	
	(Optional) Set the relative convergence tolerance	228
7.3.4	int	
	HYPRE_GMRESSetAbsoluteTol (HYPRE_Solver solver,	
	HYPRE_Real a_tol)	
	(Optional) Set the absolute convergence tolerance (default is 0)	229
7.3.5	int	
	HYPRE_GMRESSetMaxIter (HYPRE_Solver solver, int max_iter)	
	(Optional) Set maximum number of iterations	229
7.3.6	int	
	HYPRE_GMRESSetKDim (HYPRE_Solver solver, int k_dim)	
	(Optional) Set the maximum size of the Krylov space	229
7.3.7	int	
	HYPRE_GMRESSetRelChange (HYPRE_Solver solver, int rel_change)	
	(Optional) Additionally require that the relative difference in successive iterates be small	229
7.3.8	int	

	HYPRE_GMRESSetSkipRealResidualCheck (HYPRE_Solver solver, int skip_real_r_check)	
	(Optional) By default, hypre checks for convergence by evaluating the actual residual before returning from GMRES (with restart if the true residual does not indicate convergence).	230
7.0.0	,	230
7.3.9	int HYPRE_GMRESSetPrecond (HYPRE_Solver solver, HYPRE_PtrToSolverFcn precond, HYPRE_PtrToSolverFcn precond_setup,	
	HYPRE_Solver precond_solver) (Optional) Set the preconditioner to use	230
7.3.10	· - / -	230
7.3.10	int HYPRE_GMRESSetLogging (HYPRE_Solver solver, int logging) (Optional) Set the amount of logging to do	230
7.3.11	int	
	HYPRE_GMRESSetPrintLevel (HYPRE_Solver solver, int level) (Optional) Set the amount of printing to do to the screen	230
7.3.12	int HYPRE_GMRESGetNumIterations (HYPRE_Solver solver,	
	int* num_iterations) Return the number of iterations taken	231
7.3.13	· ·	201
7.3.13	int HYPRE_GMRESGetFinalRelativeResidualNorm (HYPRE_Solver solver, HYPRE_Real* norm)	
	Return the norm of the final relative residual	231
7.3.14	int HYPRE_GMRESGetResidual (HYPRE_Solver solver, void** residual) Return the residual	231
7.3.15	int HYPRE_GMRESGetTol (HYPRE_Solver solver, HYPRE_Real* tol)	231
7.3.16	int HYPRE_GMRESGetAbsoluteTol (HYPRE_Solver solver, HYPRE_Real* tol)	231
7.3.17	int	201
	HYPRE_GMRESGetMaxIter (HYPRE_Solver solver,	
	HYPRE_Intm mxx_ieer)	232
7.3.18	int HYPRE_GMRESGetKDim (HYPRE_Solver solver, int* k_dim)	232
7.3.19	int HYPRE_GMRESGetRelChange (HYPRE_Solver solver, int* rel_change)	232
7.3.20	int	
	HYPRE_GMRESGetPrecond (HYPRE_Solver solver, HYPRE_Solver* precond_data_ptr)	232
7.3.21	int HYPRE_GMRESGetLogging (HYPRE_Solver solver, int* level)	232
7.3.22	int	

	HYPRE_GMRESGetPrintLevel (HYPRE_Solver solver,	int* level)	233
7.3.23	int		
	HYPRE_GMRESGetConverged (HYPRE_Solver solver,	int^* converged)	233

int
HYPRE_GMRESSetup (HYPRE_Solver solver, HYPRE_Matrix A,
HYPRE_Vector b, HYPRE_Vector x)

Prepare to solve the system. The coefficient data in ${\tt b}$ and ${\tt x}$ is ignored here, but information about the layout of the data may be used.

7.3.2

int **HYPRE_GMRESSolve** (HYPRE_Solver solver, HYPRE_Matrix A, HYPRE_Vector b, HYPRE_Vector x)

Solve the system

7.3.3

int HYPRE_GMRESSetTol (HYPRE_Solver solver, HYPRE_Real tol)

(Optional) Set the relative convergence tolerance

Int
HYPRE_GMRESSetAbsoluteTol (HYPRE_Solver solver, HYPRE_Real a_tol)

(Optional) Set the absolute convergence tolerance (default is 0). If one desires the convergence test to check the absolute convergence tolerance only, then set the relative convergence tolerance to 0.0. (The convergence test is $||r|| \le \max(\text{relative_tolerance*}||b||)$, absolute_tolerance).)

____ 7.3.5 ______

int HYPRE_GMRESSetMaxIter (HYPRE_Solver solver, int max_iter)

(Optional) Set maximum number of iterations

_ 7.3.6 _

int HYPRE_GMRESSetKDim (HYPRE_Solver solver, int k_dim)

(Optional) Set the maximum size of the Krylov space

7.3.7

int HYPRE_GMRESSetRelChange (HYPRE_Solver solver, int rel_change)

(Optional) Additionally require that the relative difference in successive iterates be small

int **HYPRE_GMRESSetSkipRealResidualCheck** (HYPRE_Solver solver, int skip_real_r_check)

(Optional) By default, hypre checks for convergence by evaluating the actual residual before returning from GMRES (with restart if the true residual does not indicate convergence). This option allows users to skip the evaluation and the check of the actual residual for badly conditioned problems where restart is not expected to be beneficial.

7.3.9

int **HYPRE_GMRESSetPrecond** (HYPRE_Solver solver, HYPRE_PtrToSolverFcn precond, HYPRE_PtrToSolverFcn precond_setup, HYPRE_Solver precond_solver)

(Optional) Set the preconditioner to use

_ 7.3.10 _

int HYPRE_GMRESSetLogging (HYPRE_Solver solver, int logging)

(Optional) Set the amount of logging to do

_ 7.3.11 _

int HYPRE_GMRESSetPrintLevel (HYPRE_Solver solver, int level)

(Optional) Set the amount of printing to do to the screen

7 3 12

int **HYPRE_GMRESGetNumIterations** (HYPRE_Solver solver, int* num_iterations)

Return the number of iterations taken

7.3.13

HYPRE_GMRESGetFinalRelativeResidualNorm (HYPRE_Solver solver, HYPRE_Real* norm)

Return the norm of the final relative residual

7.3.14 _____

int HYPRE_GMRESGetResidual (HYPRE_Solver solver, void** residual)

Return the residual

__ 7.3.15 _____

int HYPRE_GMRESGetTol (HYPRE_Solver solver, HYPRE_Real* tol)

7.3.16

HYPRE_GMRESGetAbsoluteTol (HYPRE_Solver solver, HYPRE_Real* tol)

7317

Int
HYPRE_GMRESGetMaxIter (HYPRE_Solver solver, HYPRE_Intm mxx_ieer)

__ 7.3.18 _____

int HYPRE_GMRESGetKDim (HYPRE_Solver solver, int* k_dim)

_ 7.3.19 _____

int HYPRE_GMRESGetRelChange (HYPRE_Solver solver, int* rel_change)

7.3.20

HYPRE_GMRESGetPrecond (HYPRE_Solver solver, HYPRE_Solver* precond_data_ptr)

7.3.21

int HYPRE_GMRESGetLogging (HYPRE_Solver solver, int* level)

int HYPRE_GMRESGetPrintLevel (HYPRE_Solver solver, int* level)

7.3.23

int HYPRE_GMRESGetConverged (HYPRE_Solver solver, int* converged)

7.4

FlexGMRES Solver

Names		
7.4.1	int	
	HYPRE_FlexGMRESSetup (HYPRE_Solver solver, HYPRE_Matrix A,	
	HYPRE_Vector b, HYPRE_Vector x)	
	Prepare to solve the system.	235
7.4.2	int	
	HYPRE_FlexGMRESSolve (HYPRE_Solver solver, HYPRE_Matrix A, HYPRE_Vector b, HYPRE_Vector x)	
	Solve the system	235
7.4.3	int	
	HYPRE_FlexGMRESSetTol (HYPRE_Solver solver, HYPRE_Real tol) (Optional) Set the convergence tolerance	235
7.4.4	int	
1.4.4	HYPRE_FlexGMRESSetAbsoluteTol (HYPRE_Solver solver,	
	HYPRE_Real a_tol)	
	(Optional) Set the absolute convergence tolerance (default is 0)	235
7.4.5	int	
	HYPRE_FlexGMRESSetMaxIter (HYPRE_Solver solver, int max_iter) (Optional) Set maximum number of iterations	236
	· -	200
7.4.6	int	
	HYPRE_FlexGMRESSetKDim (HYPRE_Solver solver, int k_dim)	200
	(Optional) Set the maximum size of the Krylov space	236
7.4.7	int	

	HYPRE_FlexGMRESSetPrecond (HYPRE_Solver solver, HYPRE_PtrToSolverFcn precond, HYPRE_PtrToSolverFcn precond_setup,	
	HYPRE_Solver precond_solver)	224
	(Optional) Set the preconditioner to use	236
7.4.8	int HYPRE_FlexGMRESSetLogging (HYPRE_Solver solver, int logging) (Optional) Set the amount of logging to do	236
7.4.9	int HYPRE_FlexGMRESSetPrintLevel (HYPRE_Solver solver, int level) (Optional) Set the amount of printing to do to the screen	237
7.4.10	int HYPRE_FlexGMRESGetNumIterations (HYPRE_Solver solver, int* num_iterations)	
	Return the number of iterations taken	237
7.4.11	int HYPRE_FlexGMRESGetFinalRelativeResidualNorm (HYPRE_Solver solver, HYPRE_Real* norm)	
	Return the norm of the final relative residual	237
7.4.12	int HYPRE_FlexGMRESGetResidual (HYPRE_Solver solver, void** residual) Return the residual	237
7.4.13	int HYPRE_FlexGMRESGetTol (HYPRE_Solver solver, HYPRE_Real* tol)	237
7.4.14	int HYPRE_FlexGMRESGetMaxIter (HYPRE_Solver solver, int* max_iter)	238
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7.4.19	int HYPRE_FlexGMRESGetConverged (HYPRE_Solver solver, int* converged)	239
7.4.20	int HYPRE_FlexGMRESSetModifyPC (HYPRE_Solver solver, HYPRE_PtrToModifyPCFcn modify_pc)	
	(Optional) Set a user-defined function to modify solve-time preconditioner	239

_ 7.4.1 _

HYPRE_FlexGMRESSetup (HYPRE_Solver solver, HYPRE_Matrix A, HYPRE_Vector b, HYPRE_Vector x)

Prepare to solve the system. The coefficient data in b and x is ignored here, but information about the layout of the data may be used.

7.4.2

int **HYPRE_FlexGMRESSolve** (HYPRE_Solver solver, HYPRE_Matrix A, HYPRE_Vector b, HYPRE_Vector x)

Solve the system

7.4.3

int HYPRE_FlexGMRESSetTol (HYPRE_Solver solver, HYPRE_Real tol)

(Optional) Set the convergence tolerance

_ 7.4.4 _

HYPRE_FlexGMRESSetAbsoluteTol (HYPRE_Solver solver, HYPRE_Real a_tol)

(Optional) Set the absolute convergence tolerance (default is 0). If one desires the convergence test to check the absolute convergence tolerance only, then set the relative convergence tolerance to 0.0. (The convergence test is $||r|| \le \max(\text{relative_tolerance*}||b||)$, absolute_tolerance).)

 $_{-}$ 7.4.5 $_{-}$

int HYPRE_FlexGMRESSetMaxIter (HYPRE_Solver solver, int max_iter)

(Optional) Set maximum number of iterations

 $_$ 7.4.6 $_$

int HYPRE_FlexGMRESSetKDim (HYPRE_Solver solver, int k_dim)

(Optional) Set the maximum size of the Krylov space

__ 7.4.7 ____

int

HYPRE_FlexGMRESSetPrecond (HYPRE_Solver solver, HYPRE_PtrToSolverFcn precond_setup, HYPRE_Solver precond_solver)

(Optional) Set the preconditioner to use

_ 7.4.8 _

int HYPRE_FlexGMRESSetLogging (HYPRE_Solver solver, int logging)

(Optional) Set the amount of logging to do

7 4 9

int HYPRE_FlexGMRESSetPrintLevel (HYPRE_Solver solver, int level)

(Optional) Set the amount of printing to do to the screen

7.4.10

int

HYPRE_FlexGMRESGetNumIterations (HYPRE_Solver solver, int* num_iterations)

Return the number of iterations taken

____ 7.4.11 _____

int

HYPRE_FlexGMRESGetFinalRelativeResidualNorm (HYPRE_Solver solver, HYPRE_Real* norm)

Return the norm of the final relative residual

____ 7.4.12 _____

int HYPRE_FlexGMRESGetResidual (HYPRE_Solver solver, void** residual)

Return the residual

_ 7.4.13 _____

int HYPRE_FlexGMRESGetTol (HYPRE_Solver solver, HYPRE_Real* tol)

7.4.14

 $int~ \mathbf{HYPRE_FlexGMRESGetMaxIter}~(\mathrm{HYPRE_Solver}~solver,~int^*~max_iter)$

7.4.15

int **HYPRE_FlexGMRESGetKDim** (HYPRE_Solver solver, int* k_dim)

_ 7.4.16 _

int **HYPRE_FlexGMRESGetPrecond** (HYPRE_Solver solver, HYPRE_Solver* precond_data_ptr)

__ 7.4.17 _____

int HYPRE_FlexGMRESGetLogging (HYPRE_Solver solver, int* level)

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int HYPRE_FlexGMRESGetPrintLevel (HYPRE_Solver solver, int* level)

7.4.19

int
HYPRE_FlexGMRESGetConverged (HYPRE_Solver solver, int* converged)

___ 7.4.20 _____

int
HYPRE_FlexGMRESSetModifyPC (HYPRE_Solver solver,
HYPRE_PtrToModifyPCFcn modify_pc)

(Optional) Set a user-defined function to modify solve-time preconditioner attributes

__ 7.5 ____

LGMRES Solver

Names		
7.5.1	int	
	HYPRE_LGMRESSetup (HYPRE_Solver solver, HYPRE_Matrix A,	
	HYPRE_Vector b, HYPRE_Vector x)	
	Prepare to solve the system	241
7.5.2	int	
	HYPRE_LGMRESSolve (HYPRE_Solver solver, HYPRE_Matrix A,	
	HYPRE_Vector b, HYPRE_Vector x)	
	Solve the system.	241
7.5.3	int	
	HYPRE_LGMRESSetTol (HYPRE_Solver solver, HYPRE_Real tol)	
	(Optional) Set the convergence tolerance	241
7.5.4	int	
	HYPRE_LGMRESSetAbsoluteTol (HYPRE_Solver solver,	
	HYPRE_Real a_tol)	
	(Optional) Set the absolute convergence tolerance (default is 0)	241
7.5.5	int	
	HYPRE_LGMRESSetMaxIter (HYPRE_Solver solver, int max_iter)	
	(Optional) Set maximum number of iterations	242
7.5.6	int	
	HYPRE_LGMRESSetKDim (HYPRE_Solver solver, int k_dim)	
	(Optional) Set the maximum size of the approximation space (includes the augmentation vectors)	242
7.5.7	int	

	HYPRE_LGMRESSetAugDim (HYPRE_Solver solver, int aug_dim) (Optional) Set the number of augmentation vectors (default: 2)	242
7.5.8	int HYPRE_LGMRESSetPrecond (HYPRE_Solver solver, HYPRE_PtrToSolverFcn precond, HYPRE_PtrToSolverFcn precond_setup, HYPRE_Solver precond_solver)	
	(Optional) Set the preconditioner to use	242
7.5.9	int HYPRE_LGMRESSetLogging (HYPRE_Solver solver, int logging) (Optional) Set the amount of logging to do	242
7.5.10	int HYPRE_LGMRESSetPrintLevel (HYPRE_Solver solver, int level) (Optional) Set the amount of printing to do to the screen	243
7.5.11	int HYPRE_LGMRESGetNumIterations (HYPRE_Solver solver, int* num_iterations)	
	Return the number of iterations taken	243
7.5.12	int HYPRE_LGMRESGetFinalRelativeResidualNorm (HYPRE_Solver solver, HYPRE_Real* norm) Return the norm of the final relative residual	243
7.5.13	int HYPRE_LGMRESGetResidual (HYPRE_Solver solver, void** residual) Return the residual	243
7.5.14	int HYPRE_LGMRESGetTol (HYPRE_Solver solver, HYPRE_Real* tol)	243
7.5.15	int HYPRE_LGMRESGetMaxIter (HYPRE_Solver solver, int* max_iter)	244
7.5.16	int HYPRE_LGMRESGetKDim (HYPRE_Solver solver, int* k_dim)	244
7.5.17	int HYPRE_LGMRESGetAugDim (HYPRE_Solver solver, int* k_dim)	244
7.5.18	int HYPRE_LGMRESGetPrecond (HYPRE_Solver solver, HYPRE_Solver* precond_data_ptr)	244
7.5.19	int HYPRE_LGMRESGetLogging (HYPRE_Solver solver, int* level)	244
7.5.20	int HYPRE_LGMRESGetPrintLevel (HYPRE_Solver solver, int* level)	244
7.5.21	int HYPRE_LGMRESGetConverged (HYPRE_Solver solver, int* converged)	245

7.5.1

HYPRE_LGMRESSetup (HYPRE_Solver solver, HYPRE_Matrix A, HYPRE_Vector b, HYPRE_Vector x)

Prepare to solve the system. The coefficient data in b and x is ignored here, but information about the layout of the data may be used.

 $_{-}$ 7.5.2 $_{-}$

int **HYPRE_LGMRESSolve** (HYPRE_Solver solver, HYPRE_Matrix A, HYPRE_Vector b, HYPRE_Vector x)

Solve the system. Details on LGMRES may be found in A. H. Baker, E.R. Jessup, and T.A. Manteuffel, "A technique for accelerating the convergence of restarted GMRES." SIAM Journal on Matrix Analysis and Applications, 26 (2005), pp. 962-984. LGMRES(m,k) in the paper corresponds to LGMRES(Kdim+AugDim, AugDim).

__ 7.5.3 _____

int HYPRE_LGMRESSetTol (HYPRE_Solver solver, HYPRE_Real tol)

(Optional) Set the convergence tolerance

 $_$ 7.5.4 $_$

HYPRE_LGMRESSetAbsoluteTol (HYPRE_Solver solver, HYPRE_Real a_tol)

(Optional) Set the absolute convergence tolerance (default is 0). If one desires the convergence test to check the absolute convergence tolerance only, then set the relative convergence tolerance to 0.0. (The convergence test is $||r|| \le \max(\text{relative_tolerance*}||b||)$, absolute_tolerance).)

int HYPRE_LGMRESSetMaxIter (HYPRE_Solver solver, int max_iter)

(Optional) Set maximum number of iterations

___ 7.5.6 _____

int HYPRE_LGMRESSetKDim (HYPRE_Solver solver, int k_dim)

(Optional) Set the maximum size of the approximation space (includes the augmentation vectors)

___ 7.5.7 _____

int HYPRE_LGMRESSetAugDim (HYPRE_Solver solver, int aug_dim)

(Optional) Set the number of augmentation vectors (default: 2)

7.5.8

int **HYPRE_LGMRESSetPrecond** (HYPRE_Solver solver, HYPRE_PtrToSolverFcn precond, HYPRE_PtrToSolverFcn precond_setup, HYPRE_Solver precond_solver)

(Optional) Set the preconditioner to use

7.5.9

int HYPRE_LGMRESSetLogging (HYPRE_Solver solver, int logging)

(Optional) Set the amount of logging to do

7 5 10

int HYPRE_LGMRESSetPrintLevel (HYPRE_Solver solver, int level)

(Optional) Set the amount of printing to do to the screen

7.5.11

int **HYPRE_LGMRESGetNumIterations** (HYPRE_Solver solver, int* num_iterations)

Return the number of iterations taken

____ 7.5.12 _____

int

HYPRE_LGMRESGetFinalRelativeResidualNorm (HYPRE_Solver solver, HYPRE_Real* norm)

Return the norm of the final relative residual

___ 7.5.13 _____

int HYPRE_LGMRESGetResidual (HYPRE_Solver solver, void** residual)

Return the residual

_ 7.5.14 __

int HYPRE_LGMRESGetTol (HYPRE_Solver solver, HYPRE_Real* tol)

7.5.15

int HYPRE_LGMRESGetMaxIter (HYPRE_Solver solver, int* max_iter)

7.5.16

int **HYPRE_LGMRESGetKDim** (HYPRE_Solver solver, int* k_dim)

7.5.17

int HYPRE_LGMRESGetAugDim (HYPRE_Solver solver, int* k_dim)

7.5.18

int **HYPRE_LGMRESGetPrecond** (HYPRE_Solver solver, HYPRE_Solver* precond_data_ptr)

7.5.19

int HYPRE_LGMRESGetLogging (HYPRE_Solver solver, int* level)

7.5.20

int HYPRE_LGMRESGetPrintLevel (HYPRE_Solver solver, int* level)

7.5.21

 $int \ \mathbf{HYPRE_LGMRESGetConverged} \ (HYPRE_Solver \ solver, \ int^* \ converged)$

__ 7.6 ____

BiCGSTAB Solver

Names		
7.6.1	int HYPRE_BiCGSTABSetup (HYPRE_Solver solver, HYPRE_Matrix A, HYPRE_Vector b, HYPRE_Vector x)	
	Prepare to solve the system	246
7.6.2	int HYPRE_BiCGSTABSolve (HYPRE_Solver solver, HYPRE_Matrix A, HYPRE_Vector b, HYPRE_Vector x) Solve the system	246
7.6.3	int HYPRE_BiCGSTABSetTol (HYPRE_Solver solver, HYPRE_Real tol) (Optional) Set the convergence tolerance	246
7.6.4	int HYPRE_BiCGSTABSetAbsoluteTol (HYPRE_Solver solver, HYPRE_Real a_tol) (Optional) Set the absolute convergence tolerance (default is 0)	247
7.6.5	int HYPRE_BiCGSTABSetMaxIter (HYPRE_Solver solver, int max_iter) (Optional) Set maximum number of iterations	247
7.6.6	int HYPRE_BiCGSTABSetPrecond (HYPRE_Solver solver, HYPRE_PtrToSolverFcn precond, HYPRE_PtrToSolverFcn precond_setup, HYPRE_Solver precond_solver) (Optional) Set the preconditioner to use	247
7.6.7	int HYPRE_BiCGSTABSetLogging (HYPRE_Solver solver, int logging) (Optional) Set the amount of logging to do	247
7.6.8	int HYPRE_BiCGSTABSetPrintLevel (HYPRE_Solver solver, int level) (Optional) Set the amount of printing to do to the screen	248
7.6.9	int HYPRE_BiCGSTABGetNumIterations (HYPRE_Solver solver,	248
7.6.10	int	

	${\bf HYPRE_BiCGSTABGetFinalRelativeResidualNorm}~({\tt HYPRE_Solver}$	
	solver,	
	$\mathrm{HYPRE}_{-}\mathrm{Real}^*$	
	norm)	
	Return the norm of the final relative residual	248
7.6.11	int HYPRE_BiCGSTABGetResidual (HYPRE_Solver solver, void** residual) Return the residual	248
7.6.12	int	240
7.0.12	HYPRE_BiCGSTABGetPrecond (HYPRE_Solver solver, HYPRE_Solver* precond_data_ptr)	
		249

7.6.1

HYPRE_BiCGSTABSetup (HYPRE_Solver solver, HYPRE_Matrix A, HYPRE_Vector b, HYPRE_Vector x)

Prepare to solve the system. The coefficient data in ${\tt b}$ and ${\tt x}$ is ignored here, but information about the layout of the data may be used.

7.6.2

HYPRE_BiCGSTABSolve (HYPRE_Solver solver, HYPRE_Matrix A, HYPRE_Vector b, HYPRE_Vector x)

Solve the system

7.6.3

int HYPRE_BiCGSTABSetTol (HYPRE_Solver solver, HYPRE_Real tol)

(Optional) Set the convergence tolerance

7.6.4

HYPRE_BiCGSTABSetAbsoluteTol (HYPRE_Solver solver, HYPRE_Real a_tol)

(Optional) Set the absolute convergence tolerance (default is 0). If one desires the convergence test to check the absolute convergence tolerance only, then set the relative convergence tolerance to 0.0. (The convergence test is $||r|| \le \max(\text{relative_tolerance} *||b||)$, absolute_tolerance).)

 $_$ 7.6.5 $_$

int HYPRE_BiCGSTABSetMaxIter (HYPRE_Solver solver, int max_iter)

(Optional) Set maximum number of iterations

7.6.6

HYPRE_BiCGSTABSetPrecond (HYPRE_Solver solver, HYPRE_PtrToSolverFcn precond, HYPRE_PtrToSolverFcn precond_setup, HYPRE_Solver precond_solver)

(Optional) Set the preconditioner to use

__ 7.6.7 __

int HYPRE_BiCGSTABSetLogging (HYPRE_Solver solver, int logging)

(Optional) Set the amount of logging to do

768

int HYPRE_BiCGSTABSetPrintLevel (HYPRE_Solver solver, int level)

(Optional) Set the amount of printing to do to the screen

7.6.9

int **HYPRE_BiCGSTABGetNumIterations** (HYPRE_Solver solver, int* num_iterations)

Return the number of iterations taken

7.6.10

int

 $\label{eq:hypre_bicgstab} \begin{aligned} \mathbf{HYPRE_BiCGSTABGetFinalRelativeResidualNorm} \ (\mathbf{HYPRE_Solver} \ solver, \\ \mathbf{HYPRE_Real*} \ norm) \end{aligned}$

Return the norm of the final relative residual

__ 7.6.11 _____

int HYPRE_BiCGSTABGetResidual (HYPRE_Solver solver, void** residual)

Return the residual

7.6.12

int **HYPRE_BiCGSTABGetPrecond** (HYPRE_Solver solver, HYPRE_Solver* precond_data_ptr)

7.7

CGNR Solver

Names		
7.7.1	int HYPRE_CGNRSetup (HYPRE_Solver solver, HYPRE_Matrix A, HYPRE_Vector b, HYPRE_Vector x) Prepare to solve the system.	25
7.7.2	int HYPRE_CGNRSolve (HYPRE_Solver solver, HYPRE_Matrix A, HYPRE_Vector b, HYPRE_Vector x)	or
7.7.3	int HYPRE_CGNRSetTol (HYPRE_Solver solver, HYPRE_Real tol) (Optional) Set the convergence tolerance	25 25
7.7.4	int HYPRE_CGNRSetMaxIter (HYPRE_Solver solver, int max_iter) (Optional) Set maximum number of iterations	25
7.7.5	int HYPRE_CGNRSetPrecond (HYPRE_Solver solver, HYPRE_PtrToSolverFcn precond, HYPRE_PtrToSolverFcn precondT, HYPRE_PtrToSolverFcn precond_setup, HYPRE_Solver precond_solver) (Optional) Set the preconditioner to use.	25
7.7.6	int HYPRE_CGNRSetLogging (HYPRE_Solver solver, int logging) (Optional) Set the amount of logging to do	25
7.7.7	int HYPRE_CGNRGetNumIterations (HYPRE_Solver solver,	25
7.7.8	int HYPRE_CGNRGetFinalRelativeResidualNorm (HYPRE_Solver solver, HYPRE_Real* norm) Return the norm of the final relative residual	25
7.7.9	int	

HYPRE_CGNRGetPrecond (HYPRE_Solver solver, HYPRE_Solver* precond_data_ptr) 252
202
7.7.1
int HYPRE_CGNRSetup (HYPRE_Solver solver, HYPRE_Matrix A, HYPRE_Vector b, HYPRE_Vector x)
Prepare to solve the system. The coefficient data in b and x is ignored here, but information about the layout of the data may be used.
7.7.2
int HYPRE_CGNRSolve (HYPRE_Solver solver, HYPRE_Matrix A, HYPRE_Vector b, HYPRE_Vector x)
Solve the system
7.7.3
int HYPRE_CGNRSetTol (HYPRE_Solver solver, HYPRE_Real tol)
(Optional) Set the convergence tolerance
7.7.4
int HYPRE_CGNRSetMaxIter (HYPRE_Solver solver, int max_iter)

(Optional) Set maximum number of iterations

7.7.5

int
HYPRE_CGNRSetPrecond (HYPRE_Solver solver, HYPRE_PtrToSolverFcn
precond, HYPRE_PtrToSolverFcn precondT, HYPRE_PtrToSolverFcn
precond_setup, HYPRE_Solver precond_solver)

(Optional) Set the preconditioner to use. Note that the only preconditioner available in hypre for use with CGNR is currently BoomerAMG. It requires to use Jacobi as a smoother without CF smoothing, i.e. relax_type needs to be set to 0 or 7 and relax_order needs to be set to 0 by the user, since these are not default values. It can be used with a relaxation weight for Jacobi, which can significantly improve convergence.

7.7.6

int HYPRE_CGNRSetLogging (HYPRE_Solver solver, int logging)

(Optional) Set the amount of logging to do

_ 7.7.7 _

HYPRE_CGNRGetNumIterations (HYPRE_Solver solver, int* num_iterations)

Return the number of iterations taken

_ 7.7.8 _

int

 $\label{lem:hypre_converse} \begin{aligned} \mathbf{HYPRE_CGNRGetFinalRelativeResidualNorm} \ (\mathbf{HYPRE_Solver} \ solver, \\ \mathbf{HYPRE_Real*} \ norm) \end{aligned}$

Return the norm of the final relative residual

__ 7.7.9 _

int **HYPRE_CGNRGetPrecond** (HYPRE_Solver solver, HYPRE_Solver* precond_data_ptr)

${\bf Eigensolvers}$

\mathbf{Names}		
8.1	EigenSolvers	
		253
8.2	LOBPCG Eigensolver	
		254

These eigensolvers support many of the matrix/vector storage schemes in hypre. They should be used in conjunction with the storage-specific interfaces.

8.1

EigenSolvers

Names		
8.1.1	typedef struct hypre_Solver_struct *HYPRE_Solver The solver object	253
8.1.2	typedef struct hypre_Matrix_struct *HYPRE_Matrix The matrix object	25
8.1.3	typedef struct hypre_Vector_struct *HYPRE_Vector The vector object	25

Q 1 1

 $typedef\ struct\ hypre_Solver_struct\ \textbf{*HYPRE_Solver}$

The solver object

__ 8.1.2 _____

 $typedef \ struct \ \ hypre_Matrix_struct \ \ *HYPRE_Matrix$

The matrix object

____ 8.1.3 _____

 $typedef \ struct \ \ hypre_Vector_struct \ \ \textbf{*HYPRE_Vector}$

The vector object

_ 8.2 _

LOBPCG Eigensolver

int	
HYPRE_LOBPCGCreate (mv_InterfaceInterpreter* interpreter,	
HYPRE_MatvecFunctions* mvfunctions,	
· ·	
LOBPCG constructor	255
int	
HYPRE_LOBPCGDestroy (HYPRE_Solver solver)	
LOBPCG destructor	255
int	
HYPRE_LOBPCGSetPrecond (HYPRE_Solver solver,	
HYPRE_PtrToSolverFcn precond,	
• '	
<u> </u>	
(Optional) Set the preconditioner to use.	256
int	
HYPRE_LOBPCGGetPrecond (HYPRE_Solver solver.	
	256
int	
- 1	
Set up A and the preconditioner (if there is one)	256
int	
	HYPRE_LOBPCGCreate (mv_InterfaceInterpreter* interpreter,

	HYPRE_LOBPCGSetupB (HYPRE_Solver solver, HYPRE_Matrix B, HYPRE_Vector x)	
	(Optional) Set up B	256
8.2.7	int	
	HYPRE_LOBPCGSetupT (HYPRE_Solver solver, HYPRE_Matrix T, HYPRE_Vector x)	
	(Optional) Set the preconditioning to be applied to $Tx = b$, not $Ax = b$.	257
8.2.8	int	
	HYPRE_LOBPCGSolve (HYPRE_Solver solver, mv_MultiVectorPtr y, mv_MultiVectorPtr x, HYPRE_Real* lambda) Solve $A = lambda B x$, $y'x = 0$	257
8.2.9	int	
0.2.9	HYPRE_LOBPCGSetTol (HYPRE_Solver solver, HYPRE_Real tol) (Optional) Set the absolute convergence tolerance	257
8.2.10	int	
	HYPRE_LOBPCGSetRTol (HYPRE_Solver solver, HYPRE_Real tol)	
	(Optional) Set the relative convergence tolerance	257
8.2.11	int	
	HYPRE_LOBPCGSetMaxIter (HYPRE_Solver solver, int max_iter)	
	(Optional) Set maximum number of iterations	258
8.2.12	int	
	HYPRE_LOBPCGSetPrecondUsageMode (HYPRE_Solver solver, int mode)	
	Define which initial guess for inner PCG iterations to use: $mode = 0$: use zero initial guess, otherwise use RHS	258
8.2.13	int	
	HYPRE_LOBPCGSetPrintLevel (HYPRE_Solver solver, int level)	
	(Optional) Set the amount of printing to do to the screen	258

_ 8.2.1 _

HYPRE_LOBPCGCreate (mv_InterfaceInterpreter* interpreter, HYPRE_MatvecFunctions* mvfunctions, HYPRE_Solver* solver)

LOBPCG constructor

_ 8.2.2 _

 $\operatorname{int} \ \mathbf{HYPRE_LOBPCGDestroy} \ (\operatorname{HYPRE_Solver} \ \operatorname{solver})$

LOBPCG destructor

8.2.3

int

HYPRE_LOBPCGSetPrecond (HYPRE_Solver solver, HYPRE_PtrToSolverFcn precond, HYPRE_PtrToSolverFcn precond_setup, HYPRE_Solver precond_solver)

(Optional) Set the preconditioner to use. If not called, preconditioning is not used.

__ 8.2.4 _____

int

HYPRE_LOBPCGGetPrecond (HYPRE_Solver solver, HYPRE_Solver* precond_data_ptr)

8.2.5

int

HYPRE_LOBPCGSetup (HYPRE_Solver solver, HYPRE_Matrix A, HYPRE_Vector b, HYPRE_Vector x)

Set up A and the preconditioner (if there is one)

_ 8.2.6 _

int **HYPRE_LOBPCGSetupB** (HYPRE_Solver solver, HYPRE_Matrix B, HYPRE_Vector x)

(Optional) Set up B. If not called, B = I.

8.2.7

HYPRE_LOBPCGSetupT (HYPRE_Solver solver, HYPRE_Matrix T, HYPRE_Vector x)

(Optional) Set the preconditioning to be applied to Tx = b, not Ax = b

8.2.8 ____

int

HYPRE_LOBPCGSolve (HYPRE_Solver solver, mv_MultiVectorPtr y, mv_MultiVectorPtr x, HYPRE_Real* lambda)

Solve A x = lambda B x, y'x = 0

__ 8.2.9 _____

int HYPRE_LOBPCGSetTol (HYPRE_Solver solver, HYPRE_Real tol)

(Optional) Set the absolute convergence tolerance

__ 8.2.10 ____

int HYPRE_LOBPCGSetRTol (HYPRE_Solver solver, HYPRE_Real tol)

(Optional) Set the relative convergence tolerance

8 2 11

int HYPRE_LOBPCGSetMaxIter (HYPRE_Solver solver, int max_iter)

(Optional) Set maximum number of iterations

_ 8.2.12 _

int

HYPRE_LOBPCGSetPrecondUsageMode (HYPRE_Solver solver, int mode)

Define which initial guess for inner PCG iterations to use: mode = 0: use zero initial guess, otherwise use RHS

8.2.13

int HYPRE_LOBPCGSetPrintLevel (HYPRE_Solver solver, int level)

(Optional) Set the amount of printing to do to the screen

9

Finite Element Interface

Names		
9.1	FEI Functions	
		25
9.2	FEI Solver Parameters	
		269

_ 9.1 _

FEI Functions

Names		
9.1.1	LLNL_FEI_Impl (MPI_Comm comm) Finite element interface constructor: this function creates an instantiation	
	of the HYPRE fei class.	261
9.1.2	~LLNL_FEI_Impl () Finite element interface destructor: this function destroys the object as well as its internal memory allocations	261
9.1.3	int parameters (int numParams, char** paramStrings) The parameter function is the single most important function to pass solver information (which solver, which preconditioner, tolerance, other solver parameters) to HYPRE.	262
9.1.4	int intFields (int numFields, int* fieldSizes, int* fieldIDs) Each node or element variable has one or more fields	262
9.1.5	int initElemBlock (int elemBlockID, int numElements, int numNodesPerElement, int* numFieldsPerNode, int** nodalFieldIDs, int numElemDOFFieldsPerElement, int* elemDOFFieldIDs, int interleaveStrategy) The whole finite element mesh can be broken down into a number of element blocks.	262
9.1.6	int initElem (int elemBlockID, int elemID, int* elemConn) This function initializes element connectivity (that is, the node identifiers associated with the current element) given an element block identifier and the element identifier with the element block.	263
9.1.7	int	

	initSharedNodes (int nShared, int* sharedIDs, int* sharedLengs, int** sharedProcs)	
	This function initializes the nodes that are shared between the current processor and its neighbors.	263
9.1.8	int	
0.1.0	initCRMult (int CRListLen, int* CRNodeList, int* CRFieldList, int* CRID) This function initializes the Lagrange multiplier constraints	263
9.1.9	int	
	initComplete () This function signals to the FEI that the initialization step has been completed	264
9.1.10	int	
	resetSystem (double s) This function resets the global matrix to be of the same sparsity pattern as before but with every entry set to s	264
9.1.11	int	
	resetMatrix (double s)	
	This function resets the global matrix to be of the same sparsity pattern as before but with every entry set to s.	264
9.1.12	int	
	resetRHSVector (double s) This function resets the right hand side vector to s	265
9.1.13	int	
	resetInitialGuess (double s) This function resets the solution vector to s.	265
9.1.14	int	
	loadNodeBCs (int nNodes, int* nodeIDs, int fieldID, double** alpha, double** beta, double** gamma)	
	This function loads the nodal boundary conditions	265
9.1.15	int	
	<pre>sumInElem (int elemBlockID, int elemID, int* elemConn, double** elemStiff,</pre>	
	and also the element load to the right hand side vector	266
9.1.16	int	
	sumInElemMatrix (int elemBlock, int elemID, int* elemConn, double** elemStiffness, int elemFormat)	
	This function differs from the sumInElem function in that the right hand load vector is not passed.	266
9.1.17	int	
	sumInElemRHS (int elemBlock, int elemID, int* elemConn, double* elemLoad)	200
	This function adds the element load to the right hand side vector	266
9.1.18	int	
	loadComplete () This function signals to the FFI that the loading phase has been completed	
	This function signals to the FEI that the loading phase has been completed.	267
9.1.19	int	• •

	getNumBlockActNodes (int elemBlockID, int* nNodes)	
	This function returns the number of nodes given the element block	267
9.1.20	int	
	getNumBlockActEqns (int elemBlockID, int* nEqns)	
	This function returns the number of unknowns given the element block	267
9.1.21	int	
	getBlockNodeIDList (int elemBlockID, int numNodes, int* nodeIDList)	
	This function returns the node identifiers given the element block	268
9.1.22	int	
	getBlockNodeSolution (int elemBlockID, int numNodes, int* nodeIDList,	
	int* solnOffsets, double* solnValues)	
	This function returns the nodal solutions given the element block number.	268
9.1.23	int	
	loadCRMult (int CRID, int CRListLen, int* CRNodeList, int* CRFieldList,	
	double* CRWeightList, double CRValue)	
	This function loads the Lagrange multiplier constraints	268

9.1.1

LLNL_FEI_Impl (MPI_Comm comm)

Finite element interface constructor: this function creates an instantiation of the HYPRE fei class.

Parameters: comm - an MPI communicator

9.1.2

$^{\sim}$ LLNL_FEI_Impl ()

Finite element interface destructor: this function destroys the object as well as its internal memory allocations.

Parameters: - no parameter needed

int parameters (int numParams, char** paramStrings)

The parameter function is the single most important function to pass solver information (which solver, which preconditioner, tolerance, other solver parameters) to HYPRE.

Parameters: numParams - number of command strings

paramStrings - the command strings

int initFields (int numFields, int* fieldSizes, int* fieldIDs)

Each node or element variable has one or more fields. The field information can be set up using this function.

Parameters: - total number of fields for all variable types numFields

fieldSizes - degree of freedom for each field type

- a list of field identifiers fieldIDs

int

initElemBlock (int elemBlockID, int numElements, int numNodesPerElement, int* numFieldsPerNode, int** nodalFieldIDs, int numElemDOFFieldsPerElement, int* elemDOFFieldIDs, int interleaveStrategy)

The whole finite element mesh can be broken down into a number of element blocks. The attributes for each element block are: an identifier, number of elements, number of nodes per elements, the number of fields in each element node, etc.

Parameters: - element block identifier elemblockID

> - number of element in this block numElements

- number of nodes per element in this block numNodesPerElement

numFieldsPerNode - number of fields for each node - field identifiers for the nodal unknowns nodalFieldIDs

- number of fields for the element numElemDOFFieldsPerElement

- field identifier for the element unknowns elemDOFFieldIDs

- indicates how unknowns are ordered interleaveStratety

9.1.6

int initElem (int elemBlockID, int elemID, int* elemConn)

This function initializes element connectivity (that is, the node identifiers associated with the current element) given an element block identifier and the element identifier with the element block.

Parameters: elemblockID - element block identifier

elemID - element identifier

elemConn - a list of node identifiers for this element

9.1.7

int initSharedNodes (int nShared, int* sharedIDs, int* sharedLengs, int** sharedProcs)

This function initializes the nodes that are shared between the current processor and its neighbors. The FEI will decide a unique processor each shared node will be assigned to.

Parameters: nShared - number of shared nodes

sharedIDs - shared node identifiers

sharedLengs - the number of processors each node shares withsharedProcs - the processor identifiers each node shares with

9.1.8

int initCRMult (int CRListLen, int* CRNodeList, int* CRFieldList, int* CRID)

This function initializes the Lagrange multiplier constraints

Parameters: CRListLen - the number of constraints

CRNodeList - node identifiers where constraints are applied
 CRFieldList - field identifiers within nodes where constraints are

applied

CRID - the constraint identifier

 $_$ 9.1.9 $_$

int initComplete ()

This function signals to the FEI that the initialization step has been completed. The loading step will follow.

Parameters: - no parameter needed

_ 9.1.10 __

int resetSystem (double s)

This function resets the global matrix to be of the same sparsity pattern as before but with every entry set to s. The right hand side is set to 0.

Parameters:

s - the value each matrix entry is set to.

9.1.11

int resetMatrix (double s)

This function resets the global matrix to be of the same sparsity pattern as before but with every entry set to s.

Parameters:

s - the value each matrix entry is set to.

9.1.12

int resetRHSVector (double s)

This function resets the right hand side vector to s.

Parameters:

s - the value each right hand side vector entry is set to.

9.1.13

int resetInitialGuess (double s)

This function resets the solution vector to s.

Parameters:

s - the value each solution vector entry is set to.

9.1.14

loadNodeBCs (int nNodes, int* nodeIDs, int fieldID, double** alpha, double** beta, double** gamma)

This function loads the nodal boundary conditions. The boundary conditions

Parameters: nNodes - number of nodes boundary conditions are imposed

nodeIDs - nodal identifiers

fieldID - field identifier with nodes where BC are imposed

alpha - the multipliers for the field

the multipliers for the normal derivative of the fieldthe boundary values on the right hand side of the

 ${\it equations}$

9.1.15

int **sumInElem** (int elemBlockID, int elemID, int* elemConn, double** elemStiff, double* elemLoad, int elemFormat)

This function adds the element contribution to the global stiffness matrix and also the element load to the right hand side vector

Parameters: elemBlockID - element block identifier

elemID - element identifier

elemConn - a list of node identifiers for this element

elemStiff - element stiffness matrix

elemLoad - right hand side (load) for this element- the format the unknowns are passed in

9.1.16

int sumInElemMatrix (int elemBlock, int elemID, int* elemConn, double** elemStiffness, int elemFormat)

This function differs from the sumInElem function in that the right hand load vector is not passed.

Parameters: elemBlockID - element block identifier

elemID - element identifier

elemConn - a list of node identifiers for this element

elemStiff - element stiffness matrix

elemFormat - the format the unknowns are passed in

9.1.17

sumInElemRHS (int elemBlock, int elemID, int* elemConn, double* elemLoad)

This function adds the element load to the right hand side vector

Parameters: elemBlockID - element block identifier

elemID - element identifier

elemConna list of node identifiers for this elementright hand side (load) for this element

9.1.18

int loadComplete ()

This function signals to the FEI that the loading phase has been completed.

Parameters: - no parameter needed

9.1.19

int getNumBlockActNodes (int elemBlockID, int* nNodes)

This function returns the number of nodes given the element block.

Parameters: elemBlockID - element block identifier

nNodes - the number of nodes to be returned

_ 9.1.20 _

int **getNumBlockActEqns** (int elemBlockID, int* nEqns)

This function returns the number of unknowns given the element block.

Parameters: elemBlockID - element block identifier

nEqns - the number of unknowns to be returned

9.1.21 _

int getBlockNodeIDList (int elemBlockID, int numNodes, int* nodeIDList)

This function returns the node identifiers given the element block.

Parameters: elemBlockID - element block identifier

numNodes - the number of nodesnodeIDList - the node identifiers

9.1.22 _

int

getBlockNodeSolution (int elemBlockID, int numNodes, int* nodeIDList, int* solnOffsets, double* solnValues)

This function returns the nodal solutions given the element block number.

Parameters: elemBlockID - element block identifier

numNodes - the number of nodesnodeIDList - the node identifiers

solnOffsets - the equation number for each nodal solution

solnValues - the nodal solution values

 $_{-}$ 9.1.23 $_{-}$

int

loadCRMult (int CRID, int CRListLen, int* CRNodeList, int* CRFieldList, double* CRWeightList, double CRValue)

This function loads the Lagrange multiplier constraints

Parameters: CRID - the constraint identifier

CRListLen - the number of constraints

CRNodeList - node identifiers where constraints are applied
 CRFieldList - field identifiers within nodes where constraints are

applied

CRWeightList - a list of weights applied to each specified field

CRValue - the constraint value (right hand side of the con-

straint)

9.2

FEI Solver Parameters

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9.2.1

Preconditioners and Solvers

Here the various options for solvers and preconditioners are defined.

solver xxx where xxx specifies one of cg, gmres, fgmres, bicgs, bicgstab, tfqmr, symqmr, superlu, or superlux. The default is gmres. The solver type can be followed by override to specify its priority when multiple solvers are declared at random order.

preconditioner xxx where xxx is one of diagonal, pilut, euclid, parasails, boomeramg, poly, or mli. The default is diagonal. Another option for xxx is reuse which allows the preconditioner to be reused (this should only be set after a preconditioner has been set up already). The preconditioner type can be followed by override to specify its priority when multiple preconditioners are declared at random order.

- maxIterations xxx where xxx is an integer specifying the maximum number of iterations permitted for the iterative solvers. The default value is 1000.
- **tolerance xxx** where **xxx** is a floating point number specifying the termination criterion for the iterative solvers. The default value is 1.0E-6.
- gmresDim xxx where xxx is an integer specifying the value of m in restarted GMRES(m). The default value is 100.
- **stopCrit xxx** where xxx is one of absolute or relative stopping criterion.
- **superluOrdering xxx** where xxx specifies one of natural or mmd (minimum degree ordering). This ordering is used to minimize the number of nonzeros generated in the LU decomposition. The default is natural ordering.
- **superluScale xxx** where xxx specifies one of y (perform row and column scalings before decomposition) or n. The default is no scaling.

9.2.2

BoomerAMG

Parameter options for the algebraic multigrid preconditioner BoomerAMG.

- **amgMaxLevels xxx** where xxx is an integer specifying the maximum number of levels to be used for the grid hierarchy.
- amgCoarsenType xxx where xxx specifies one of falgout or ruge, or default (CLJP) coarsening for Boomer-AMG.
- **amgMeasureType xxx** where xxx specifies one of local or or global. This parameter affects how coarsening is performed in parallel.
- amgRelaxType xxx where xxx is one of jacobi (Damped Jacobi), gs-slow (sequential Gauss-Seidel), gs-fast (Gauss-Seidel on interior nodes), or hybrid. The default is hybrid.
- **amgNumSweeps xxx** where xxx is an integer specifying the number of pre- and post-smoothing at each level of BoomerAMG. The default is two pre- and two post-smoothings.
- amgRelaxWeight xxx where xxx is a floating point number between 0 and 1 specifying the damping factor for BoomerAMG's damped Jacobi and GS smoothers. The default value is 1.0.
- amgRelaxOmega xxx where xxx is a floating point number between 0 and 1 specifying the damping factor for BoomerAMG's hybrid smoother for multiple processors. The default value is 1.0.
- amgStrongThreshold xxx where xxx is a floating point number between 0 and 1 specifying the threshold used to determine strong coupling in BoomerAMG's coasening. The default value is 0.25.
- amgSystemSize xxx where xxx is the degree of freedom per node.

amgMaxLevels xxx where xxx is an integer specifying the maximum number of iterations to be used during the solve phase.

amgUseGSMG - tells BoomerAMG to use a different coarsening called GSMG.

amgGSMGNumSamples where xxx is the number of samples to generate to determine how to coarsen for GSMG.

Parameter options for the smoothed aggregation preconditioner MLI.

outputLevel xxx where xxx is the output level for diagnostics.

method xxx where xxx is either AMGSA (default), AMGSAe, to indicate which MLI algorithm is to be used.

numLevels xxx where xxx is the maximum number of levels (default=30) used.

maxIterations xxx where xxx is the maximum number of iterations (default = 1 as preconditioner).

cycleType xxx where xxx is either 'V' or 'W' cycle (default = 'V').

strengthThreshold xxx strength threshold for coarsening (default = 0).

smoother xxx where xxx is either Jacobi, BJacobi, GS, SGS, HSGS (SSOR,default), BSGS, ParaSails, MLS, CGJacobi, CGBJacobi, or Chebyshev.

numSweeps xxx where xxx is the number of smoother sweeps (default = 2).

coarseSolver xxx where xxx is one of those in 'smoother' or SuperLU (default).

minCoarseSize xxx where xxx is the minimum coarse grid size to control the number of levels used (default = 3000).

Pweight xxx where xxx is the relaxation parameter for the prolongation smoother (default 0.0).

nodeDOF xxx where xxx is the degree of freedom for each node (default = 1).

nullSpaceDim $\mathbf{x}\mathbf{x}\mathbf{x}$ where $\mathbf{x}\mathbf{x}\mathbf{x}$ is the dimension of the null space for the coarse grid (default = 1).

useNodalCoord xxx where xxx is either 'on' or 'off' (default) to indicate whether the nodal coordinates are used to generate the initial null space.

saAMGCalibrationSize xxx where xxx is the additional null space vectors to be generated via calibration (default = 0).

 $numSmoothVecs\ xxx$ where xxx is the number of near null space vectors used to create the prolongation operator (default = 0).

smoothVecSteps xxx where xxx is the number of smoothing steps used to generate the smooth vectors (default = 0).

In addition, to use 'AMGSAe', the parameter 'haveSFEI' has to be sent into the FEI using the parameters function (this option is valid only for the Sandia FEI implementation).

Various

Parameter options for ILUT, ParaSails and polynomial preconditioners are defined.

euclidNlevels xxx where xxx is an non-negative integer specifying the desired sparsity of the incomplete factors. The default value is 0.

euclidThreshold xxx where xxx is a floating point number specifying the threshold used to sparsify the incomplete factors. The default value is 0.0.

parasailsThreshold xxx where xxx is a floating point number between 0 and 1 specifying the threshold used to prune small entries in setting up the sparse approximate inverse. The default value is 0.0.

parasailsNlevels xxx where xxx is an integer larger than 0 specifying the desired sparsity of the approximate inverse. The default value is 1.

parasailsFilter xxx where xxx is a floating point number between 0 and 1 specifying the threshold used to prune small entries in A. The default value is 0.0.

parasailsLoadbal xxx where xxx is a floating point number between 0 and 1 specifying how load balancing has to be done (Edmond, explain please). The default value is 0.0.

parasailsSymmetric sets Parasails to take A as symmetric.

parasailsUnSymmetric sets Parasails to take A as nonsymmetric (default).

parasailsReuse sets Parasails to reuse the sparsity pattern of A.

polyorder xxx where xxx is the order of the least-squares polynomial preconditioner.

9.2.5

Matrix Reduction

Parameters which define different reduction modes.

schurReduction turns on the Schur reduction mode.

slideReduction turns on the slide reduction mode.

slideReduction2 turns on the slide reduction mode version 2 (see section 2).

slideReduction3 turns on the slide reduction mode version 3 (see section 2).

9.2.6

Performance Tuning and **Diagnostics**

Parameters control diagnostic information, memory use, etc.

outputLevel xxx where xxx is an integer specifying the output level. An output level of 1 prints only the solver information such as number of iterations and timings. An output level of 2 prints debug information such as the functions visited and preconditioner information. An output level of 3 or higher prints more debug information such as the matrix and right hand side loaded via the LinearSystemCore functions to the standard output.

setDebug xxx where xxx is one of slideReduction1, slideReduction2, slideReduction3 (level 1,2,3 diagnostics in the slide surface reduction code), printMat (print the original matrix into a file), printReducedMat (print the reduced matrix into a file), printSol (print the solution into a file), ddilut (output diagnostic information for DDIlut preconditioner setup), and amgDebug (output diagnostic information for AMG).

optimizeMemory cleans up the matrix sparsity pattern after the matrix has been loaded. (It has been kept to allow matrix reuse.)

imposeNoBC turns off the boundary condition to allow diagnosing the matrix (for example, checking the null space.)

Miscellaneous

Parameters that are helpful for finite element information.

- **AConjugateProjection xxx** where xxx specifies the number of previous solution vectors to keep for the A-conjugate projection. The default is 0 (the projection is off).
- minResProjection xxx where xxx specifies the number of previous solution vectors to keep for projection. The default is 0 (the projection is off).
- haveFEData indicates that additional finite element information are available to assist in building more efficient solvers.
- have SFEI indicates that the simplified finite element information are available to assist in building more efficient solvers.