

# Creating mixed FLASH/Chombo applications

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## 1 Accessing the code

The Chombo-dev svn branch is at the following location:

```
svn+ssh://USERNAME@flash.uchicago.edu/home/svn/repos/FLASH3/branches/Chombo-dev
```

There is also a svn vendor branch containing Chombo-3.0-Oct2009 at the following location:

```
svn+ssh://USERNAME@flash.uchicago.edu/home/svn/repos/FLASH3/vendor/chomboForFlash
```

It is convenient to use this vendor branch because it contains Chombo Makefiles for various Flash Center machines and a small patch to build applications on Intrepid BG/P.

## 2 Creating FLASH-Chombo applications

The standard FLASH build system is used to create a mixed FLASH-Chombo application. The FLASH Makefile will link FLASH against the appropriate dimensionality Chombo libraries and so Chombo must be pre-built in 1D, 2D and 3D before building a mixed FLASH-Chombo application.

A FLASH Config file registers Chombo as an external library to the setup script and so the generated FLASH Makefile depends on the macros CFLAGS\_CHOMBO and LIB\_CHOMBO. The macro CFLAGS\_CHOMBO is used when compiling the FLASH C++ code that interacts with Chombo library and LIB\_CHOMBO is used to link against the Chombo libraries. Both macros must be defined by the user in Makefile.h.chombo. Figure 1 shows the macro values used on code.uchicago.edu.

The setup shortcuts +chombo\_ug and +chombo\_amr are available to build UG and AMR FLASH applications respectively. They correspond to:

```
chombo_ug:- unit=Grid/GridMain/Chombo/UG:- index-reorder: Grid=Chombo:- maxblocks=1:  
-nofbs:- makefile=chombo:chomboCompatibleHydro=True  
  
chombo_amr:- unit=Grid/GridMain/Chombo/AMR:- index-reorder: Grid=Chombo:  
-nofbs:- makefile=chombo:chomboCompatibleHydro=True
```

## 3 Available software stacks

- GNU

```

CHOMBO_PATH = /home/cdaley/flash/chomboForFlash/current

_DIM = $(shell grep 'define NDIM' Flash.h | cut -d ' ' -f 3)
_CPP = $(shell make vars DIM=${_DIM} -C ${CHOMBO_PATH}/lib | \
    awk -F 'CPPFLAGS=' '/^CPPFLAGS/{print $$2}')
_LIB = $(shell make vars DIM=${_DIM} -C ${CHOMBO_PATH}/lib | \
    awk -F 'config=' '/^config/{print $$2}')

CFLAGS_CHOMBO = -I${CHOMBO_PATH}/lib/include ${_CPP} -DCH_LANG_CC
LIB_CHOMBO = -L${CHOMBO_PATH}/lib \
-lamrtimependent${_LIB} \
-lamrtools${_LIB} \
-lboxtools${_LIB} \
-lbasetools${_LIB} \
-lstdc++

```

Figure 1: FLASH Makefile.h.chombo variables

- lenovolaptop
  - \* sites/lenovolaptop/Makefile.h
  - \* lib/chombo/lib/mk/local/Make.defs.lenovolaptop
- archimedes
  - \* sites/archimedes.uchicago.edu/Makefile.h
  - \* lib/chombo/lib/mk/local/Make.defs.archimedes
- Intel
  - code
    - \* sites/code.uchicago.edu/Makefile.h
    - \* lib/chombo/lib/mk/local/Make.defs.code
  - mongchi
    - \* sites/mongchi.uchicago.edu/Makefile.h
    - \* lib/chombo/lib/mk/local/Make.defs.mongchi
- XL
  - intrepid
    - \* sites/intrepid.alcf.anl.gov/Makefile.h
    - \* lib/chombo/lib/mk/local/Make.defs.intrepid

NOTE: gfortran >= 4.4.2 is required because there is a pointer interoperability bug in earlier versions of gfortran. See **GCC bugzilla 40962**.

## 4 Runtime parameters

### 4.1 Chombo UG

- **iGridSize, jGridSize, kGridSize:** The total number of grid points along each dimension of the global domain.
- **iProcs, jProcs, kProcs:** The total number of processors along each dimension of the global domain.

### 4.2 Chombo AMR

- **iGridSize, jGridSize, kGridSize:** The total number of grid points along each dimension of the coarsest level.
- **lrefine\_max:** The unit-based maximum refinement level. A value of 3 means that blocks can refine a maximum of 2 times.
- **flux\_correct:** Logical flag to turn on/off flux correction. Currently always off because flux correction is not yet implemented.
- **maxBlockSize:** The maximum FLASH block size. In Chombo terminology this is the maximum grid size.
- **BRMeshRefineFillRatio:** See `a_fillRatio` argument of `BRMeshRefine` define method in Chombo user-guide.
- **BRMeshRefineBufferSize:** See `a_bufferSize` argument of `BRMeshRefine` define method in Chombo user-guide.
- **BRMeshRefineBlockFactor:** See `a_blockFactor` argument of `BRMeshRefine` define method in Chombo user-guide. Note that setting this parameter equal to **maxBlockSize** leads to a mesh where all blocks have the same number of cells.
- **tagRadius:** The radius of cells around a tagged cell that should also be tagged, i.e. marked for refinement

## 5 Unit\_test

There are unit tests in sub-directories 1 and 2 at the following location in the Chombo-dev source tree:

`source/Grid/GridMain/Chombo/wrapper/unit_tests`

## 5.1 1

This test is independent of FLASH and Chombo. It simply tests interoperability of pointers in a serial, mixed Fortran/C++ code. Note that this test will fail with versions of gfortran  $< 4.4.2$  because the compiler introduces a pointer indexing bug (see GCC bugzilla 40962). Build with make [gnu, intel, ibm, sun].

## 5.2 2

This test depends on Chombo only. It tests usage of Chombo library in a parallel, mixed Fortran/C++ code. The main() function is written in Fortran and calls C++ functions that use Chombo to generate a distributed uniform grid. The application must be run with 4 MPI processes. Build with ./compile.sh [gnu, intel, ibm].

# 6 Test problems

## 6.1 Sedov

The standard FLASH Sedov explosion problem can be run with Chombo Grid implementation. The generated checkpoint files are in HDF5 format and can be visualized using visit. The following aliases may be useful.

```
alias visit_flash='visit -assume_format FLASH'
alias visit_chombo='visit -assume_format Chombo'
```

### 6.1.1 Uniform grid

Creates a uniform grid with 1 box per MPI process. The box size is determined from igrdsize,jgrdsize,kgrdsize and iprocs,jprocs,kprocs runtime parameters.

```
./setup Sedov -auto +chombo_ug -parfile=test_chombo_ug_2d.par
```

A grid size of  $512^2$  is created for FLASH UG and Chombo UG and the resulting application is run on 1024 cores of Intrepid. The final checkpoint file occurs at  $t=0.05$  seconds and is visualized with visit. Figures 2 and 3 show the generated pseudocolor plot of density mesh variable for FLASH UG and Chombo UG.

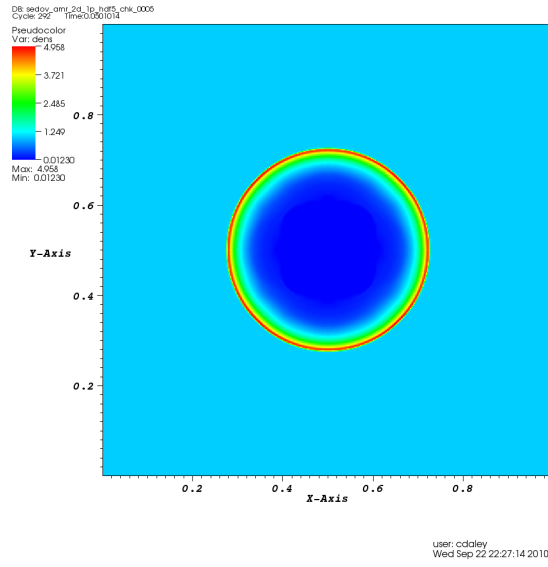


Figure 2: High resolution Sedov application with NOFBS uniform grid

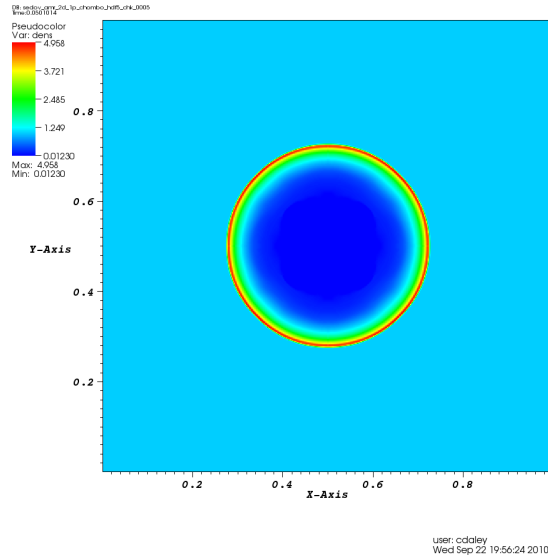


Figure 3: High resolution Sedov application with Chombo grid



### 6.1.2 Adaptive grid

The Sedov problem can also be run with a Chombo adaptive grid using the following setup line:

```
./setup Sedov -auto +chombo_amr -parfile=test_chombo_amr_2d.par
```

The density mesh variable and Chombo patches are shown in Figure 4 at 0.01 second intervals. It can be seen that the Chombo patches are clustered near the explosion front. This demonstrates that FLASH is able to communicate to Chombo the cells that need to be tagged for refinement, however, there are some anomalies with the clustering because the test problem is symmetrical. This non-symmetry is most likely an issue with my custom `gr_markRefineDerefine` which is a hacked cell-based version of the default block-based Paramesh `gr_markRefineDerefine`.

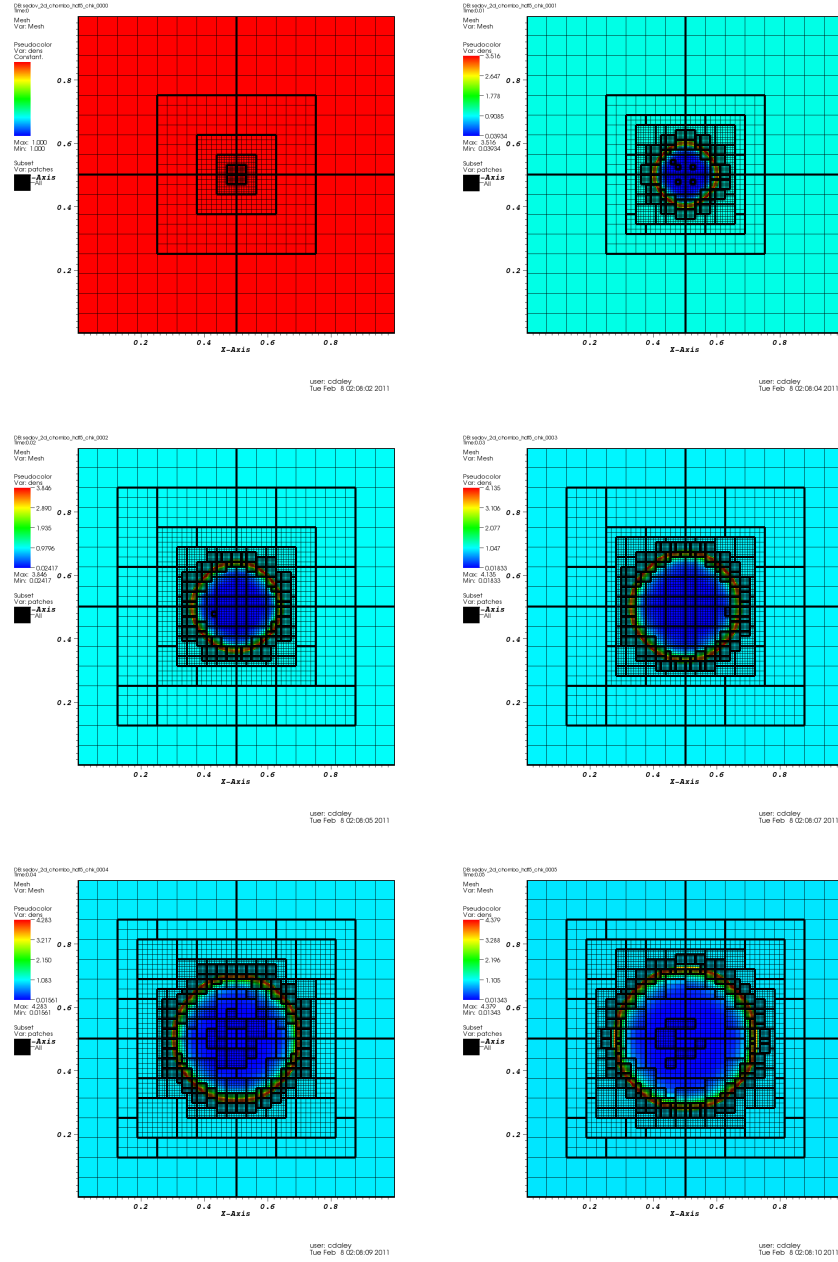


Figure 4: Density at 0.01 second intervals in a Sedov simulation using AMR capabilities of Chombo

## 6.2 Sod

The FLASH Sod problem can be setup with split or unsplit Hydro using:

```
./setup Sod -auto +chombo_amr -parfile=test_chombo_amr_2d.par  
./setup Sod -auto +chombo_amr +uhd -parfile=test_chombo_amr_2d.par
```

This problem tests the code’s ability to capture shocks. Originally we had problems with non-convergence in Riemann solver. We believe the non-convergence was because of fine-coarse interpolation in a region of the domain where there were steep gradients.

We overcame the non-convergence by tagging cells for refinement more liberally; whenever we tagged a cell we also tagged nearest neighbor cells up to a set distance. This allows us to refine a region of the domain (rather than a single cell) in preparation for an approaching shock, and seems to avoid the fine-coarse interpolation issue. The default tag distance is 2 and is controlled by the parameter, tagRadius. The Sedov images above were obtained before we changed the tagging code and so had a tag distance of 0.

The density mesh variable is shown at 0.04 second intervals in Figure 5.

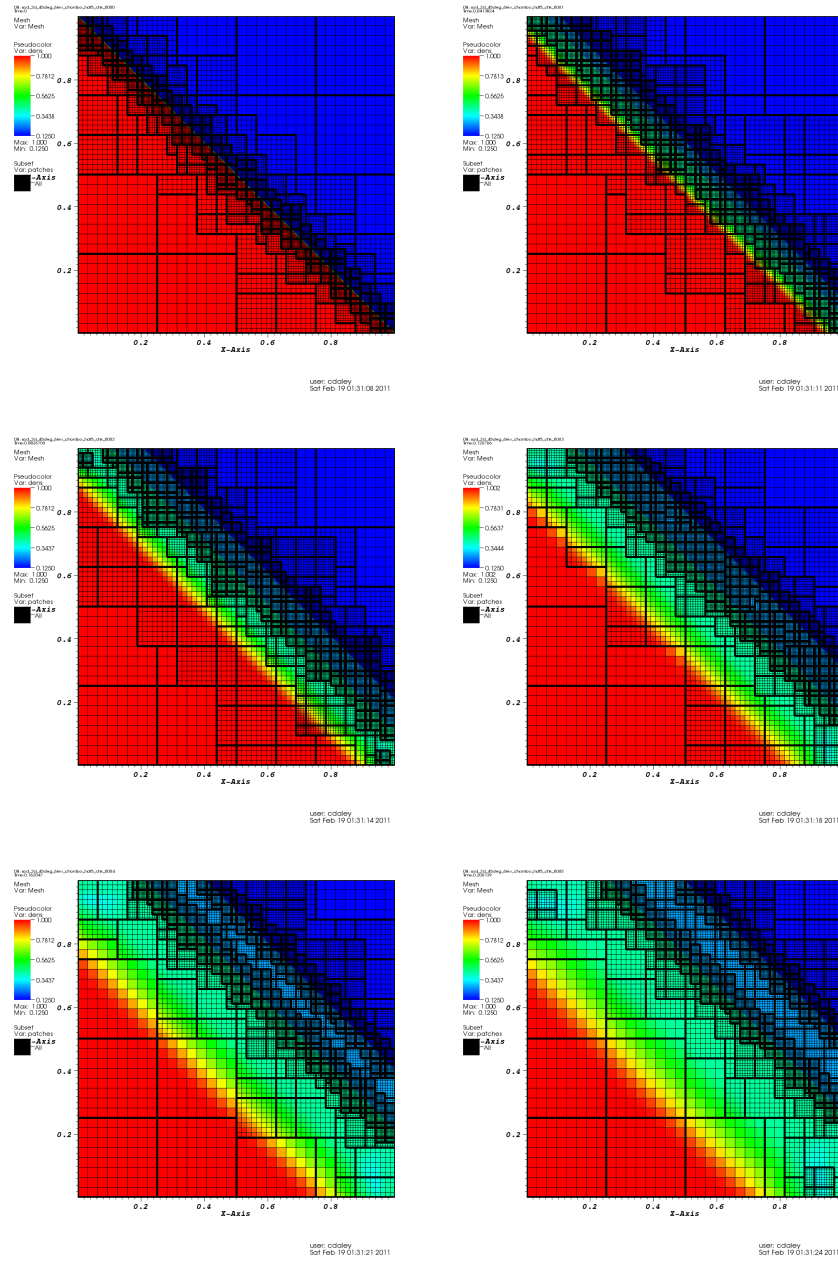


Figure 5: Density at 0.04 second intervals in a Sod simulation using AMR capabilities of Chombo

## 6.3 Others

The following unit tests work as expected:

```
setupName: unitTest/Eos/Multigamma
setupOptions: -auto -3d +chombo_amr +noio -debug
numProcs: 1
parfiles: <pathToSimulations>/unitTest/Eos/Multigamma/chombo_3d.par

setupName: unitTest/Eos/Multigamma
setupOptions: -auto -3d +chombo_ug +noio -debug
numProcs: 1
parfiles: <pathToSimulations>/unitTest/Eos/Multigamma/chombo_3d.par

setupName: unitTest/PFFT_PoissonFD
setupOptions: -auto -3d +chombo_ug +noio -debug
numProcs: 4
pathToParfiles: <pathToSimulations>/unitTest/PFFT_PoissonFD
parfiles: <pathToParfiles>/test_UG_4p_3d_16cube.par
          <pathToParfiles>/test_UG_4p_3d_32cube.par
          <pathToParfiles>/test_UG_4p_3d_64cube.par
          <pathToParfiles>/test_UG_4p_3d_128cube.par
```

The following test problems run, but have been looked at less:

```
setupName: Blast2
setupOptions: -auto -1d +chombo_amr
numProcs: 4
parfiles: <pathToSimulations>/Blast2/test_chombo_amr_1d.par

setupName: IsentropicVortex
setupOptions: -auto -2d +chombo_amr
numProcs: 2
parfiles: <pathToSimulations>/IsentropicVortex/test_chombo_amr_2d.par

setupName: StirTurb
setupOptions: -auto -3d +chombo_amr
numProcs: 4
parfiles: <pathToSimulations>/StirTurb/test_chombo_amr_3d.par
```

## 7 Design overview

The FLASH Grid API subroutines call “C” linkage functions in `chombo_f.c_api` which in turn call methods on an object that interacts with Chombo library. There are two different objects for interacting with Chombo

library, either a `chombo_uniform_grid` (for UG) or a `chombo_adaptive_grid` (for AMR).

The `chombo_uniform_grid` object is much simpler and only contains a `DisjointBoxLayout` Chombo object and a `LevelData<FArrayBox>` Chombo object. It provides identical functionality to a single box per MPI process FLASH uniform grid.

The `chombo_adaptive_grid` object is much more complicated and contains partial usage of Chombo’s AMR framework. As described in the Chombo user-guide, the Chombo AMR framework consists an AMR object calling methods on a user’s object (which is instantiated from a class that inherits from `AMRLevel` class). In this framework the user’s `AMRLevel` class is expected to contain the data representation for one level and also implement the algorithms for advancing one level in time.

The derived `AMRLevel` class in FLASH does not contain any code for advancing one level in time. This does fit with the intended usage of AMR framework in a Chombo application, but doing things this way allows FLASH’s Hydro unit to perform time advancement. It is an abuse of Chombo’s AMR framework that allows re-use of nearly 1000 lines of code from the example `AMRLevel` class named `AMRLevelPolytropicGas`.

The core methods of AMR and `AMRLevel` class are compared against FLASH subroutines in Tables 1 and 2, respectively.

Chombo method	FLASH subroutines	Operations
<code>setupForNewAMRRun</code>	<code>Driver_initFlash</code>	Initialize the simulation
<code>run</code>	<code>Driver_evolveFlash</code>	Evolve the simulation
<code>conclude</code>	<code>Driver_finalizeFlash</code>	Finalize the simulation
<code>initialGrid</code>	<code>Grid_initDomain</code>	Initialize the domain
<code>regrid</code>	<code>Grid_updateRefinement</code>	Update refinement
<code>makeBaseLevelMesh</code>	<code>gr_createDomain</code>	Create base level

Table 1: Chombo AMR class

The AMR class performs setup, evolution and conclusion of a custom simulation. These same steps are performed by FLASH’s Driver unit and specifically the subroutines `Driver_initFlash`, `Driver_evolveFlash` and `Driver_finalizeFlash`. As such, we cannot use the AMR class because we want FLASH to “drive” the simulation and not Chombo. It is possible to view FLASH and `chombo_adaptive_grid` class as collectively providing the same functionality as AMR class.

The `AMRLevelPolytropicGas` class (derived from `AMRLevel` class) contains most of the functionality required by a FLASH Grid API implementation. It is therefore a good starting point for FLASH’s custom `AMRLevel` class named `AMRLevelFlash`. Some functionality is not required by FLASH such as the `advance`, `computeDt`, and `initialGrid` methods which are already provided by FLASH’s Hydro, Driver and Simulation units. As such, FLASH’s custom `AMRLevel` class only provides implementations for methods corresponding to FLASH subroutines prefixed with `Grid_` or `gr_` as listed in Table 2.

Chombo method	FLASH subroutines	Operations
advance	Hydro	Advance this level by one time step
computeDt	Driver_computeDt	Return the maximum stable time step
initialData	Simulation_initBlock	Initialize grid data
postInitialize	N/A	If there is a finer level then replace coarse data with the average of fine data
postTimeStep	Grid_conserveFluxes & Grid_restrictAllLevels	If there is a finer level then reflux and replace coarse data with the average of fine data
tagCells	Grid_markRefineDerefine	If there is a coarser level then interpolate undefined ghost cells. Exchange guardcells. Tag cells according to the relative gradient
regrid	amr_morton_order (called from within gr_updateRefinement)	Copy Unew to Uold. Load balance Unew. If there is a coarser level then interpolate internal data into Unew. Copy overlapping sections from Uold to Unew
postRegrid	N/A	No implementation, but can perform any post-regridding operations

Table 2: Chombo AMRLevelPolytropicGas class (derived from AMRLevel class)

Both AMRLevelPolytropicGas and AMRLevelFlash contain cell centered solution data at an old and new state (m\_UOld and m\_UNew). In AMRLevelFlash there is also additional storage for temporary solution data at a single state (m\_scratchCtr). This has identical size to the solution data in m\_UOld and m\_UNew and is required by certain FLASH units such as Unsplit Hydro. The data is “scratch” space and is updated only by FLASH; ghost cells are not exchanged and the data is not written to file. All solution data is of type LevelData<FArrayBox>.

The tagCells implementation in AMRLevelFlash simply iterates over all cells in the scratchCtr array for the ‘tagc’ mesh variable. Cell values can be either FLASH\_TRUE or FLASH\_FALSE depending on whether or not a cell should be tagged. This means that the tagCells implementation simply translates from one data representation (array) to another data representation (IntVectSet). The actual tagging happens in a customized gr\_markRefineDerefine; this version is a cell-based version of the default block-based gr\_markRefineDerefine. The design strategy of communicating information through the scratchCtr array avoids the complexity of callbacks from Chombo to FLASH.

Since there is no time advancement code in AMRLevelFlash, new methods are added to the class so that an outsider (i.e. FLASH) can interact with the level data. This includes being able to exchange guardcells, interpolate and restrict grid data, retrieve and modify grid data, access grid metadata, regrid and tag cells. These new methods are:

- **fillGuardCells** Exchange guardcells for m\_UNew. If there is a fine-coarse interface then fill fine guardcells using interpolated coarse values. Note: My combination of PiecewiseLinearFillPatch and QuadCFInterp may be silly - it may be better to just accept the *everywhere* first order accurate Piece-

wiseLinearFillPatch. In AMRLevelPolytropic the exchange and interpolated exchange only happen in tagCells. In FLASH Grid\_markRefineDerefine calls fillGuardcells to do the exchange and then calls gr.markRefineDerefine to do the custom cell tagging.

- **coarseAverage** Replaces coarse data with the average of fine data. In AMRLevelPolytropic a coarse average happens in postInitialize and postTimeStep.
- **makeBoxLookupTable** Saves a vector of DataIndexes so that a FLASH integer block ID can be directly translated to a Chombo grid data index.
- **getBoxInfo** Obtains box metadata for FLASH
- **getDataPtr** Obtains a pointer to grid data for FLASH
- **getDataIndex** Helper method for getBoxInfo and getDataPtr
- **getDisjointBoxLayout** Helper method for makeBoxLookupTable and getBoxInfo
- **getNumBoxes** Obtains the number of boxes on the current level

The flow of control for Chombo and FLASH-Chombo applications can be broken down into initialization, evolution and finalization sections. Figure 6 shows Chombo flow on the left and FLASH-Chombo flow on the right. The amr object is an instance of AMR class for Chombo and an instance of Chombo\_Adaptive\_Grid class for FLASH-Chombo. The subroutines / methods being called are indented.

## 8 Key functions/subroutines

- **chombo\_f\_c\_interfaces.F90** The interface of all C++ functions called by Fortran.
- **chombo\_f\_c\_api.C** The implementation of all C++ function called by Fortran. All functions have C linkage and operate on a Chombo grid object (either chombo\_uniform\_grid or chombo\_adaptive\_grid).
- **flash\_ftypes.F90** The Fortran declaration of the interoperable types. This includes flash\_amr\_info.t, flash\_amr\_info.t for defining the initial grid and box.t to retrieve box metadata from Chombo to FLASH.
- **flash\_ctypes.h** The C++ declaration of the interoperable types mentioned in flash\_ftypes.
- **chombo\_uniform\_grid.C** A Chombo grid object that provides a simple uniform grid (1 box per MPI process).
- **chombo\_adaptive\_grid.C** A Chombo grid object that provides a AMR mesh.
- **AMRLevelFlash.C** A subclass of AMRLevel that is used by AMRLevelFlashFactory. Based on example/AMRGodunov/srcPolytropic/AMRLevelPolytropicGas.cpp.
- **AMRLevelFlashFactory.C** A subclass of AMRLevelFactory that is used by chombo\_adaptive\_grid. Based on example/AMRGodunov/srcPolytropic/AMRLevelPolytropicGasFactory.cpp.



amr.setupForNewAMRRun	call Driver_initFlash
amr.initialGrid	call Grid_initDomain
amr.makeBaseLevelMesh	call gr_createDomain
	amr.makeBaseLevelMesh
	call gr_expandDomain
<b>do</b>	<b>do</b>
amrlevels.initialGrid	amr.BuildInitialGrid
amrlevels.initialData	amrlevels.initialGrid
amrlevels.tagCellsInit	call Simulation_initBlock
	call Grid_markRefineDerefine
	call Grid_fillGuardCells
	amr.FillGuardCells
	amrlevels.fillGuardCells
	call gr_markRefineDerefine
	amr.RefineInitialGrid
	amrlevels.tagCells
	mesh_refine.regrid
mesh_refine.regrid	amr.BuildInitialGrid
	amrlevels.initialGrid
amrlevels.initialGrid	call Simulation_initBlock
amrlevels.initialData	amr.FinalizeInitialGrid
	amrlevels.postInitialize
amrlevels.postInitialize	call Driver_verifyInitDt
amrlevels.computeInitialDt	
	call Driver_evolveFlash
amr.run	<b>do</b>
amr.timeStep	
<b>do</b>	
amr.regrid	amr.PreAdvance
amrlevels.tagCells	amrlevels.preAdvance
mesh_refine.regrid	call Hydro
amrlevels.preRegrid	call Driver_computeDt
amrlevels.regrid	amr.PostTimeStep
amrlevels.postRegrid	amrlevels.postTimeStep
	call Grid_updateRefinement
	call Grid_markRefineDerefine
	amr.regrid
	amrlevels.tagCells
	mesh_refine.regrid
	amrlevels.preRegrid
	amrlevels.regrid
	amrlevels.postRegrid
amrlevels.advance	
amrlevels.computeDt	
amrlevels.postTimeStep	
amr.conclude	call Driver_finalizeFlash

Figure 6: Flow of control for Chombo and FLASH-Chombo applications

## 9 Limitations / Known issues

### 9.1 Memory tracking

The AMR implementation crashes when Chombo library is compiled with memory tracking. The crash only happens when using more than 1 MPI process. My current thinking is that this is a problem with Chombo memory tracking rather than a FLASH memory bug because valgrind does not detect any heap management problems. A setup that experiences this problem and the resulting back trace is shown below:

```
./setup Sedov -auto +chombo_amr -parfile=test_chombo_amr_2d.par

(gdb) bt
#0 0x000000397e0332c0 in exit () from /lib64/libc.so.6
#1 0x00002aaaaadb2206 in MPIU_Exit (exit_code=255) at exit.c:22
#2 0x00002aaaaadd2f7e in MPID_Abort (comm=0xff, mpi_errno=0,
    exit_code=2117413264,
    error_msg=0xffffffffffffffff <Address 0xffffffffffffffff out of bounds>)
    at mpid_abort.c:104
#3 0x00002aaaaad39e1b in PMPI_Abort (comm=255, errorcode=0) at abort.c:119
#4 0x0000000000c20b25 in MayDay::Error (
    a_msg=0xff <Address 0xff out of bounds>, exit_code=0,
    $02=<value optimized out>, $03=<value optimized out>) at MayDay.cpp:76
#5 0x0000000000c23b0e in freep (a_p=0xff) at memtrack.cpp:785
#6 0x0000000000bbbc8b in Copier::~Copier (this=0x7fffffffcb0,
    $n2=<value optimized out>) at Copier.cpp:811
#7 0x0000000000b20c4b in PiecewiseLinearFillPatch::~PiecewiseLinearFillPatch (
    this=0x7fffffff650, $=<value optimized out>)
    at PiecewiseLinearFillPatch.cpp:52
#8 0x000000000041dc4b in AMRLevelFlash::fillGuardCells (this=0x17a9360,
    $&7=<value optimized out>) at AMRLevelFlash.C:1140
#9 0x0000000000cf99a in Chombo_Adaptive_Grid::FillGuardCells (
    this=0x179caf0, $;7=<value optimized out>) at chombo_adaptive_grid.C:597
#10 0x00000000006fb87d in ch_fill_guardcells () at chombo_f_c_api.C:89
#11 0x000000000056e289 in grid_fillguardcells (mype=0, griddatastruct=380,
    idir=-1, minlayers=Cannot access memory at address 0x0
    ) at Grid_fillGuardCells.F90:152
```

### 9.2 VisIt

VisIt only supports 2D and 3D Chombo output files. The following error appears when opening a 1D Chombo output file in a VisIt session launched with `visit -debug`.

```
A.mdserver.1.vlog:ERROR: Reader only supports 2D and 3D data sets.
```

```
A.mdserver.1.vlog-Exception: (InvalidFilesException)
/home/cdaley/build/gnu/visit-2.1.2/visit2.1.2/src/databases/Chombo/
avtChomboFileFormat.C, line 460: There was an error opening
/home/cdaley/flash/Chombo-dev/sod_1d_chombo_reg/run/sod_1d_64pt_chombo.hdf5_chk_0001.
It may be an invalid file.
```

### 9.3 Isentropic Vortex

The Isentropic Vortex problem aborts at step 89, but only when run with more than 2 processors. The problem can be reproduced using r13336 of Chombo-dev on code.uchicago.edu with an intel software stack. The setup line, abort message and stack trace from a 4 processor-run are:

```
./setup IsentropicVortex -auto -2d +chombo_amr
-parfile=test_chombo_amr_2d.par

MayDay: TreeIntVectSet.cpp:1995: Assertion 'bxNumPts != 0' failed. !!!

#4 0x0000000000ce2703 in MayDay::Abort (
    a_msg=0xff <Address 0xff out of bounds>, $01=<value optimized
    out>)
    at MayDay.cpp:87
#5 0x0000000000c868f in TreeIntVectSet::numPts (this=0xff,
    $R0=<value optimized out>) at TreeIntVectSet.cpp:1995
#6 0x0000000000cb311b in IntVectSet::numPts (this=0xff,
    $W6=<value optimized out>) at IntVectSet.cpp:487
#7 0x0000000000c7690d in BRMeshRefine::makeBoxesParallel (this=0xff,
    a_mesh=..., a_tags=..., a_pnd=..., a_domain=...,
    a_maxBoxSize=1293968485,
    a_depth=0, a_totalBufferSize=3, a_minSize=100, a_procInterval=...,
    $ =<value optimized out>, $ =<value optimized out>,
    $ =<value optimized out>, $ =<value optimized out>,
    $ =<value optimized out>, $ =<value optimized out>,
    $ =<value optimized out>, $ =<value optimized out>)
    at BRMeshRefine.cpp:315
---Type <return> to continue, or q <return> to quit---
#8 0x0000000000c765bb in BRMeshRefine::makeBoxes (this=0xff,
    a_mesh=...,
    a_tags=..., a_pnd=..., a_domain=..., a_maxBoxSize=1293968485,
    a_totalBufferSize=26481840, $ =<value optimized out>,
    $ =<value optimized out>, $ =<value optimized out>,
    $ =<value optimized out>, $ =<value optimized out>,
    $ =<value optimized out>)
    at BRMeshRefine.cpp:199
#9 0x0000000000cbe2a4 in MeshRefine::regrid (this=0xff,
    a_newmeshes=...,
```

```

    a_tags=..., a_baseLevel=-1, a_topLevel=-1408463424,
    a_OldMeshes=...,
    $e1=<value optimized out>, $y4=<value optimized out>,
    $y5=<value optimized out>, $y6=<value optimized out>,
    $y7=<value optimized out>, $y8=<value optimized out>) at
    MeshRefine.cpp:647
#10 0x00000000007605e0 in Chombo_Adaptive_Grid::Regrid
    (this=0x1854460,
    a_base_level=0, $N1=<value optimized out>, $N2=<value optimized
    out>)
    at chombo_adaptive_grid.C:343
#11 0x000000000078e1ee in ch_regrid (baseLevel=0) at
    chombo_f_c_api.C:77
#12 0x0000000000614ba3 in grid_updaterefinement (mype=0, nstep=90,
    time=4.4601542468994086, gridchanged=.FALSE.)
    at Grid_updateRefinement.F90:94
#13 0x00000000004ede63 in driver_evolveflash () at
    Driver_evolveFlash.F90:239
#14 0x000000000053d866 in flash () at Flash.F90:49
#15 0x0000000000409ffc in main ()

```

## 9.4 TODO

- Flux conservation.
- Sfocu tool for comparison of Chombo layout HDF5 files.
- Investigate lack of symmetrical refinement in Sedov problem - is this simply because I have not run the problem with high enough resolution?
- Restarts.
- Add ability to specify refinement jumps for each level in flash.par.
- Remove all references to gr\_ilo, gr\_ihi e.t.c from Grid unit because blocks can have different size in the same MPI process. This has already been done in Chombo-dev but needs to be cleaned up. In some cases I simply commented out code that is only used for debugging purposes.
- Non-Cartesian geometries.
- Burn unit needs to be made NOFBS compliant.
- Multipole unit needs to be made compatible with Chombo.

```

The first problem found is:
gr_mpoleInit.F90:  call RuntimeParameters_get("Nblockx", nBlockx),

setup line is:

```

```
./setup unitTest/Gravity/Poisson3 -auto -3d -maxblocks=600 -debug  
+chombo_amr -noc  
—without—unit=physics/Gravity/GravityMain/Poisson/Multigrid  
—unit=physics/Gravity/GravityMain/Poisson/Multipole
```