Setting up a new simulation

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Terminology



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"Child Grid"

A cosmology simulation is only 1 choice...



Since Enzo models both gas and particles, it can be used for most astrophysical problems

When you run Enzo:



CosmologySimulationInitialize.C



NewSimulationInitialize.C

Read parameters

Set up grid levels



т

Grid_NewSimulation InitializeGrid.C





An example:

Rotating Cylinder

cp workshop/RotatingCylinderInitialize.C enzo/src/enzo/. cp workshop/Grid_RotatingCylinderInitializeGrid.C enzo/src/enzo/.



RotatingCylinderInitialize.C

(called by InitializeNew.C)

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Set up grid levels



RotatingCylinderInitialize.C

FLOAT RotatingCylinderCenterPosition[MAX_DIMENSION]; for(dim = 0; dim < MAX_DIMENSION; dim++) RotatingCylinderCenterPosition[dim] = 0.5*(DomainRightEdge[dim]+DomainLeftEdge[dim]); // middle of the box

float RotatingCylinderVelocity[3] = {0.0, 0.0, 0.0}; // gas initally at rest
FLOAT RotatingCylinderRadius = 0.3;
float RotatingCylinderLambda = 0.05;
float RotatingCylinderOverdensity = 20.0;
int RotatingCylinderRefineAtStart = 1;

/* read input from file */

while (fgets(line, MAX_LINE_LENGTH, fptr) != NULL) {

ret = 0;

- /* read parameters specifically for radiating shock problem*/
- ret += sscanf(line, "RotatingCylinderOverdensity = %"FSYM, &RotatingCylinderOverdensity);
- ret += sscanf(line, "RotatingCylinderLambda = %"FSYM, &RotatingCylinderLambda);
- ret += sscanf(line, "RotatingCylinderRefineAtStart = %"ISYM, &RotatingCylinderRefineAtStart);
- ret += sscanf(line, "RotatingCylinderRadius = %"PSYM, &RotatingCylinderRadius);
- ret += sscanf(line, "RotatingCylinderCenterPosition = %"PSYM" %"PSYM" %"PSYM, RotatingCylinderCenterPosition, RotatingCylinderCenterPosition+1, RotatingCylinderCenterPosition+2);

/* if the line is suspicious, issue a warning */

"*** warning: the following parameter line was not interpreted:\n%s\n", line);

Check we've not missed any

} // end input from parameter file

Tuesday, May 22, 12

Set defaults

Read problem-specific parameters

if (TopGrid.GridData->RotatingCylinderInitializeGrid(RotatingCylinderRadius, RotatingCylinderCenterPosition, RotatingCylinderLambda, RotatingCylinderOverdensity) == f

ENZO_FAIL("Error in RotatingCylinderInitializeGrid.");

/* Set up initial AMR levels */

if (RotatingCylinderRefineAtStart) {

/* Declare, initialize and fill out the LevelArray. */

LevelHierarchyEntry *LevelArray[MAX_DEPTH_OF_HIERARCHY]; for (level = 0; level < MAX_DEPTH_OF_HIERARCHY; level++) LevelArray[level] = NULL; AddLevel(LevelArray, &TopGrid, 0);

/* Add levels to the maximum depth or until no new levels are created, and re-initialize the level after it is created. */

```
for (level = 0; level < MaximumRefinementLevel; level++) {
    if (RebuildHierarchy(&MetaData, LevelArray, level) == FAIL) {
      ENZO_FAIL("Error in RebuildHierarchy.");
    }
}</pre>
```

```
if (LevelArray[level+1] == NULL)
    break;
```

```
LevelHierarchyEntry *Temp = LevelArray[level+1];
```

while (Temp != NULL) {

Temp = Temp->NextGridThisLevel;

} // end: loop over levels

Set ρ, e, \overline{v} cells in top grid

Create AMR hierarchy



Set ρ, e, \overline{v} cells in child grid

Largely identical for all problem types

Grid_RotatingCylinderInitializeGrid.C

(called by RotatingCylinderInitialize.C)



Assign memory for ρ, e, \bar{v} fields

Set ρ, e, \overline{v} for each cell

Grid_RotatingCylinderInitializeGrid.C

/* create fields */

NumberOfBaryonFields = 0; FieldType[NumberOfBaryonFields++] = Density; FieldType[NumberOfBaryonFields++] = TotalEnergy; if (DualEnergyFormalism) FieldType[NumberOfBaryonFields++] = InternalEnergy; int vel = NumberOfBaryonFields; FieldType[NumberOfBaryonFields++] = Velocity1; FieldType[NumberOfBaryonFields++] = Velocity2; FieldType[NumberOfBaryonFields++] = Velocity2;

Create
$$\rho, e, \bar{v}$$
 fields

```
if (ProcessorNumber != MyProcessorNumber)
  return SUCCESS;
```

Only do this on I processor

/* declarations */

FLOAT x = 0, y = 0, z = 0, radius, z_distance, x_velocity = 0.0, y_velocity = 0.0, z_velocity = 0.0; float sintheta, costheta, omega; float outside_density = 1.0, outside_energy = 0.5, density = 1.0, energy = 0.5; int i, j, k, dim, cellindex;

/* compute size of fields */

int size = 1; for (dim = 0; dim < GridRank; dim++) size *= GridDimension[dim];

/* allocate fields */

```
int field;
for (field = 0; field < NumberOfBaryonFields; field++)
if (BaryonField[field] == NULL)
BaryonField[field] = new float[size];
```

int DensNum, GENum, TENum, Vel1Num, Vel2Num, Vel3Num, MetalNum; if (this->IdentifyPhysicalQuantities(DensNum, GENum, Vel1Num, Vel2Num, Vel3Num, TENum) == FAIL) { ENZO_FAIL("Error in IdentifyPhysicalQuantities.\n");

Assign memory for fields

Useful function for finding fields

```
for (k = 0; k < GridDimension[2]; k++)
for (j = 0; j < GridDimension[1]; j++)
for (i = 0; i < GridDimension[0]; i++) {
    cellindex = i + j*GridDimension[0] + k*GridDimension[0]*GridDimension[1];</pre>
```

energy = outside_energy; density = outside_density;

- x = CellLeftEdge[0][i] + 0.5*CellWidth[0][i];
- y = CellLeftEdge[1][j] + 0.5*CellWidth[1][j];z = CellLeftEdge[2][k] + 0.5*CellWidth[2][k];

/* Find distance from center. */

radius = POW(x-RotatingCylinderCenterPosition[0], 2.0) + POW(y-RotatingCylinderCenterPosition[1], 2.0);

radius = sqrt(radius); // ok, now it's just radius

z_distance = fabs(z-RotatingCylinderCenterPosition[2]);

if ((radius <= RotatingCylinderRadius) && (z_distance <= RotatingCylinderRadius)) {
 // inside the cylinder</pre>

density = outside_density * RotatingCylinderOverdensity;

sintheta = (y-RotatingCylinderCenterPosition[1])/radius; costheta = (x-RotatingCylinderCenterPosition[0])/radius;

// x,y and z velocity.

x_velocity = -1.0*sintheta*omega*radius; y_velocity = costheta*omega*radius; z_velocity = 0.0;

energy = outside_energy / RotatingCylinderOverdensity;

} // if (r <= RotatingCylinderRadius)

BaryonField[DensNum][cellindex] = density;

BaryonField[Vel1Num][cellindex] = x_velocity; BaryonField[Vel2Num][cellindex] = y_velocity; BaryonField[Vel3Num][cellindex] = z_velocity;

BaryonField[TENum][cellindex] = energy;

Loop over cells

set background values

Cell position

$ho, e, ar{v}$ inside cylinder

Set final field value

Make the 2 new files (or copy from another test problem): MyProblemInitialize.C Grid_MyProblemInitializeGrid.C

Add the 2 new files to Make.config.objects (so they get compiled) /src/enzo/Make.config.objects

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```
rotate2d.o \
rotate3d.o \
RotatingCylinderInitialize.o \
s66_st1.o \
s90_st1.o \
```

```
Grid_RHIonizationClumpInitializeGrid.o \
Grid_RHIonizationSteepInitializeGrid.o \
Grid_RHIonizationTestInitializeGrid.o \
Grid_RotatingCylinderInitialize.o \
Grid_SedovBlastInitializeGrid.o \
Grid_SedovBlastInitializeGrid3D.o \
```

Make the 2 new files (or copy from another test problem): MyProblemInitialize.C Grid_MyProblemInitializeGrid.C

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Add definition of Grid initialization routine to Grid.h /src/enzo/Grid.h

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Add definition of Grid initialization routine to Grid.h /src/enzo/Grid.h

FLOAT RadiatingShockCenterPosition[MAX_DIMENSION]);

/* Initialize a grid for a rotating cylinder collapse */ int RotatingCylinderInitializeGrid(FLOAT RotatingCylinderRadius, FLOAT RotatingCylinderCenterPosition[MAX_DIMENSION], float RotatingCylinderLambda, float RotatingCylinderOverdensity);

int RotatingSphereInitializeGrid(FLOAT RotatingSphereRadius, FLOAT RotatingSphereCenterPosition[MAX_DIMENSION], float RotatingSphereI ambda.

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Add definition of Grid initialization routine to Grid.h /src/enzo/Grid.h

Add prototype and call for non-Grid initialization routine in InitializeNew (make sure it has a unique ProblemType #).

/src/enzo/InitializeNew.C

Make the 2 new files (or copy from another test problem): MyProblemInitialize.C Grid_MyProblemInitializeGrid.C

Add the 2 new files to Make.config.objects (so they get compiled)

Grid.h

Src int ShockInABoxInitialize(FILE *fptr, FILE *Outfptr, HierarchyEntry &TopGrid, TopGridData &MetaData, ExternalBoundary &Exterior); int ImplosionInitialize(FILE *fptr, FILE *Outfptr, HierarchyEntry &TopGrid, TopGridData &MetaData); int RotatingCylinderInitialize(FILE *fptr, FILE *Outfptr, HierarchyEntry &TopGrid, TopGridData &MetaData); int ConductionTestInitialize(FILE *fptr, FILE *Outfptr, HierarchyEntry &TopGrid, TopGridData &MetaData);

Sr(// 10) RotatingCylinder

if (ProblemType == 10)
ret = RotatingCylinderInitialize(fptr, Outfptr, TopGrid, MetaData);

Add prototype and call for non-Grid initialization routine in InitializeNew (make sure it has a unique ProblemType #).

/src/enzo/InitializeNew.C

Points to watch

<u>Using Zeus</u> (HydroMethod = 2)

Zeus uses a face-centered velocity



/* Loop over dims if using Zeus (since vel's face-centered). */

for (dim = 0; dim < 1+(HydroMethod == Zeus_Hydro ? GridRank : 0); dim++) {

/* Compute position. */

xpos = x-DiskPosition[0] (dim == 1 ? 0.5*CellWidth[0][0] : 0.0);
ypos = y-DiskPosition[1] (dim == 2 ? 0.5*CellWidth[1][0] : 0.0);
zpos = z-DiskPosition[2] (dim == 3 ? 0.5*CellWidth[2][0] : 0.0);

/* Compute velocty: L x r_perp. */

Velocity[2] = DiskVelocityMag*(AngularMomentum[0]*xhat[1] -AngularMomentum[1]*xhat[0]);

It also uses internal energy, not total energy

Points to watch

<u>Energy</u>

BaryonField[TENum] is energy/mass

Particle Mass

ParticleMass[i] is particle mass / cell volume

GravitationalConstant

GravitationalConstant = 4 pi G

Must be in code units if SelfGravity = I