

The logo for CEA (Commissariat à l'énergie atomique et aux énergies alternatives) is displayed in white lowercase letters on a red square background. A thin green horizontal line is positioned below the letters.

# The Great Merge

## Progress towards a unified RAMSES

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# Motivation

RAMSES has a large variety of users

Many versions of the code exist

## Possible issues:

- Bug fixes are not always applied to every version
- Duplicated work
- Incompatible physics features
- Complicates comparison of results

**SOLUTION: Merge everything back into 1 central repository**

## Advantages:

- Overview of past/ongoing development
- More available physics
- Code reviews by other users
- Less bugs!
- Easier maintenance
- ...

# Merging process

## The case of ramses-ism

Fork of  
rteyssie/ramses

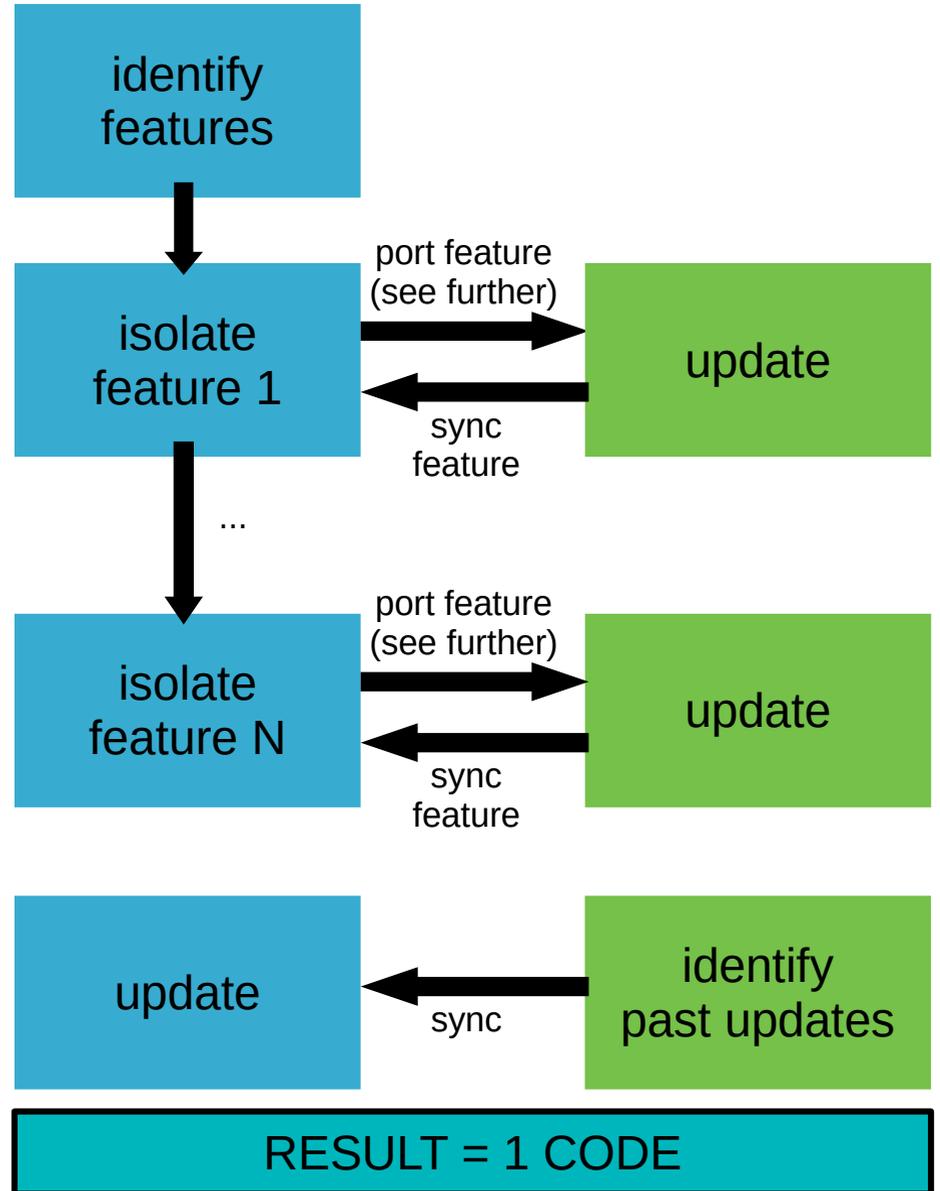
Sync (1542 commits behind)

Branch	Behind	Ahead	Updated
master MAIN DEVELOPMENT			BC 2017-11-09
ism	0	438	2 hours ago

Not possible to do a normal merge

# RAMSES-ISM

# RAMSES master



Big features:

Non-ideal MHD

Dust

Turbulence driving

Flux-limited-diffusion (FLD)

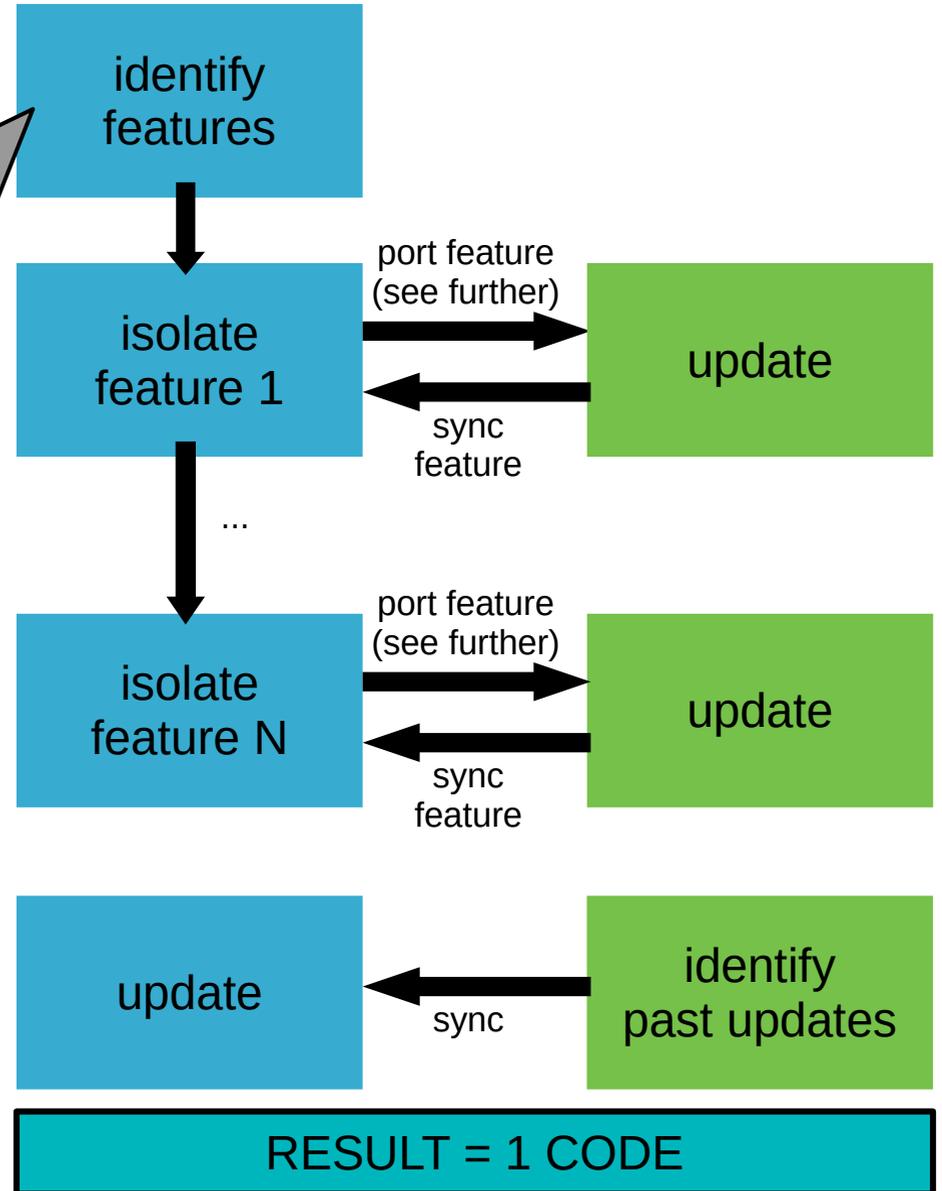
Stellar feedback (SN, HII, jets)

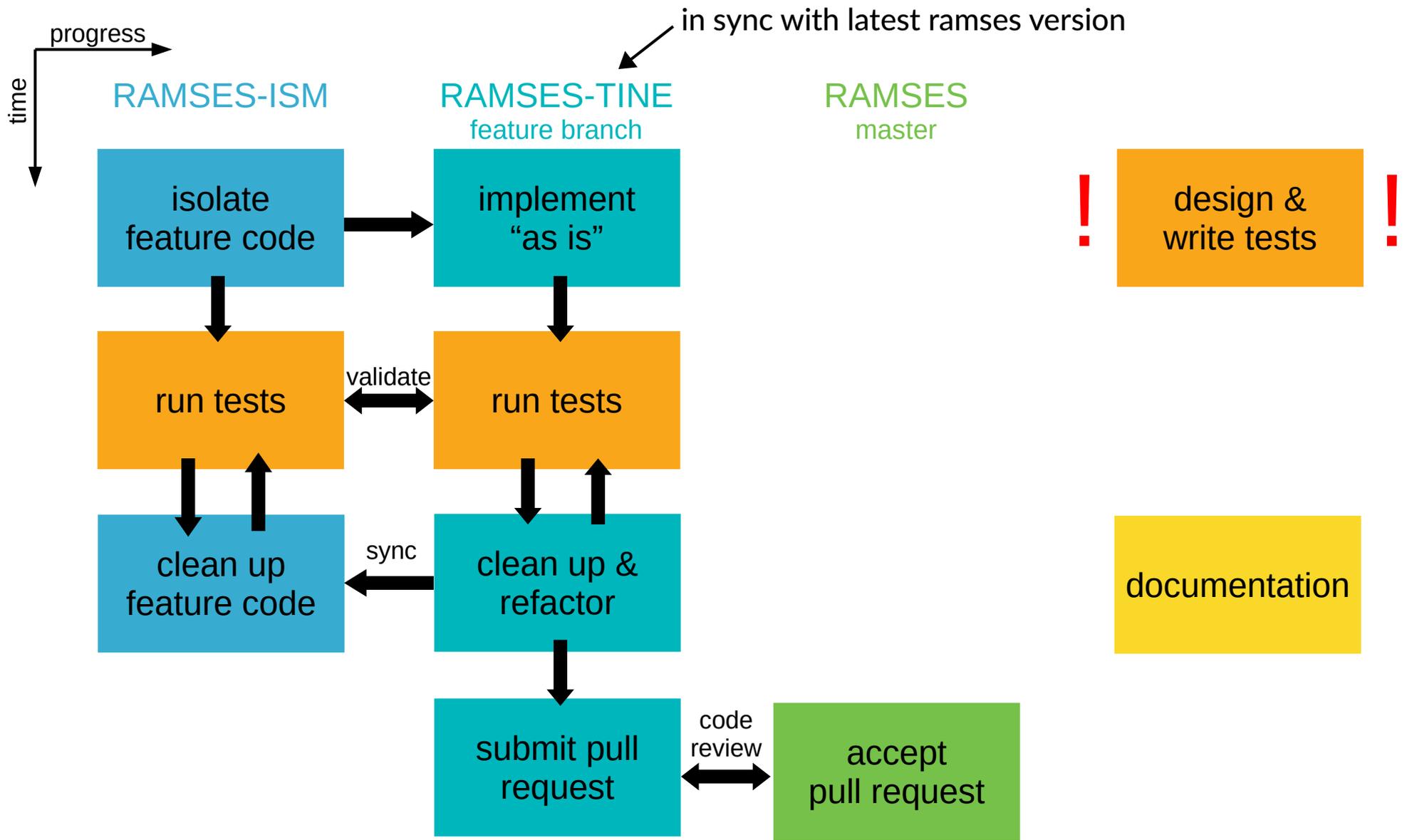
ISM cooling

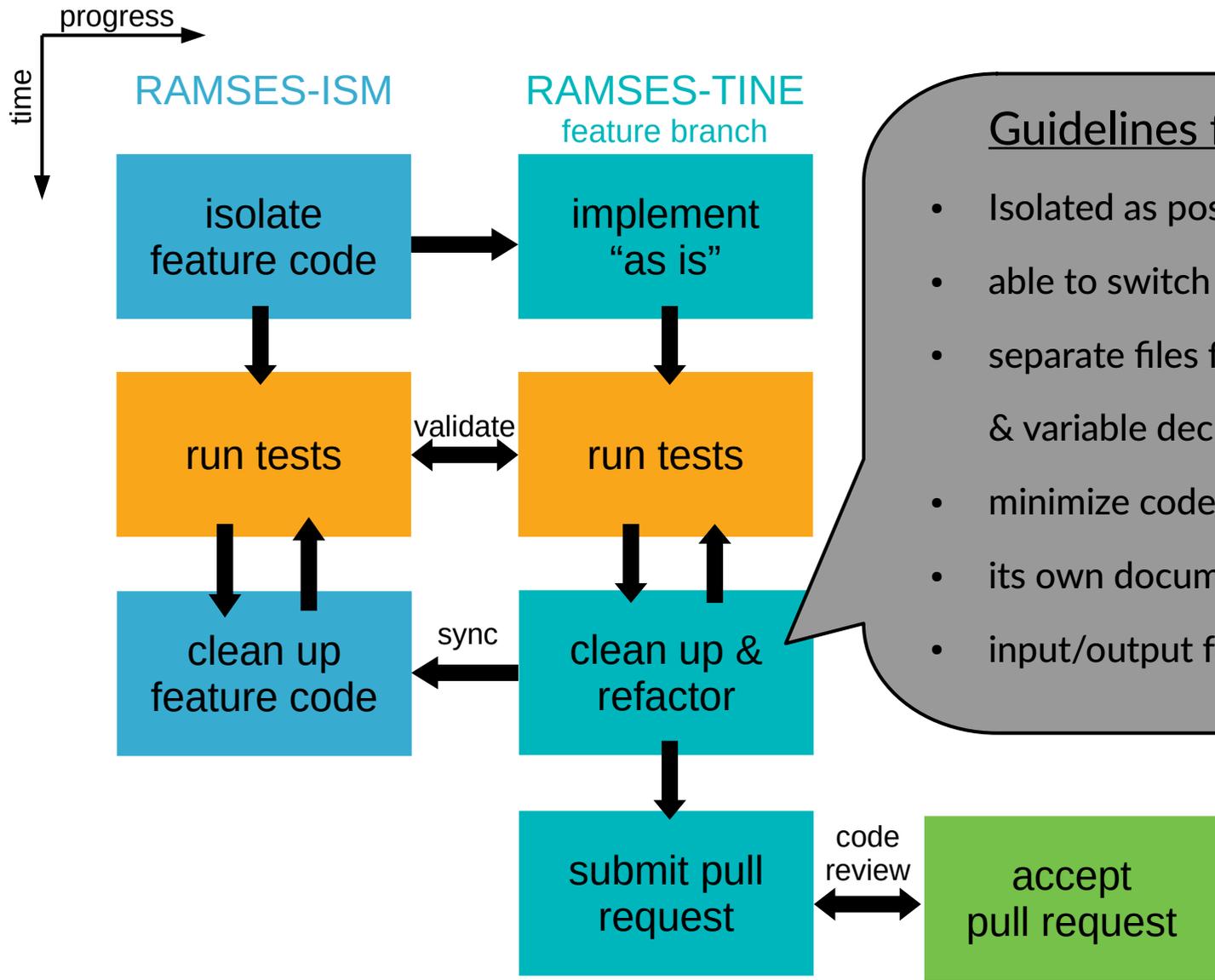
Cosmic ray anisotropic diffusion

RAMSES-ISM

RAMSES  
master

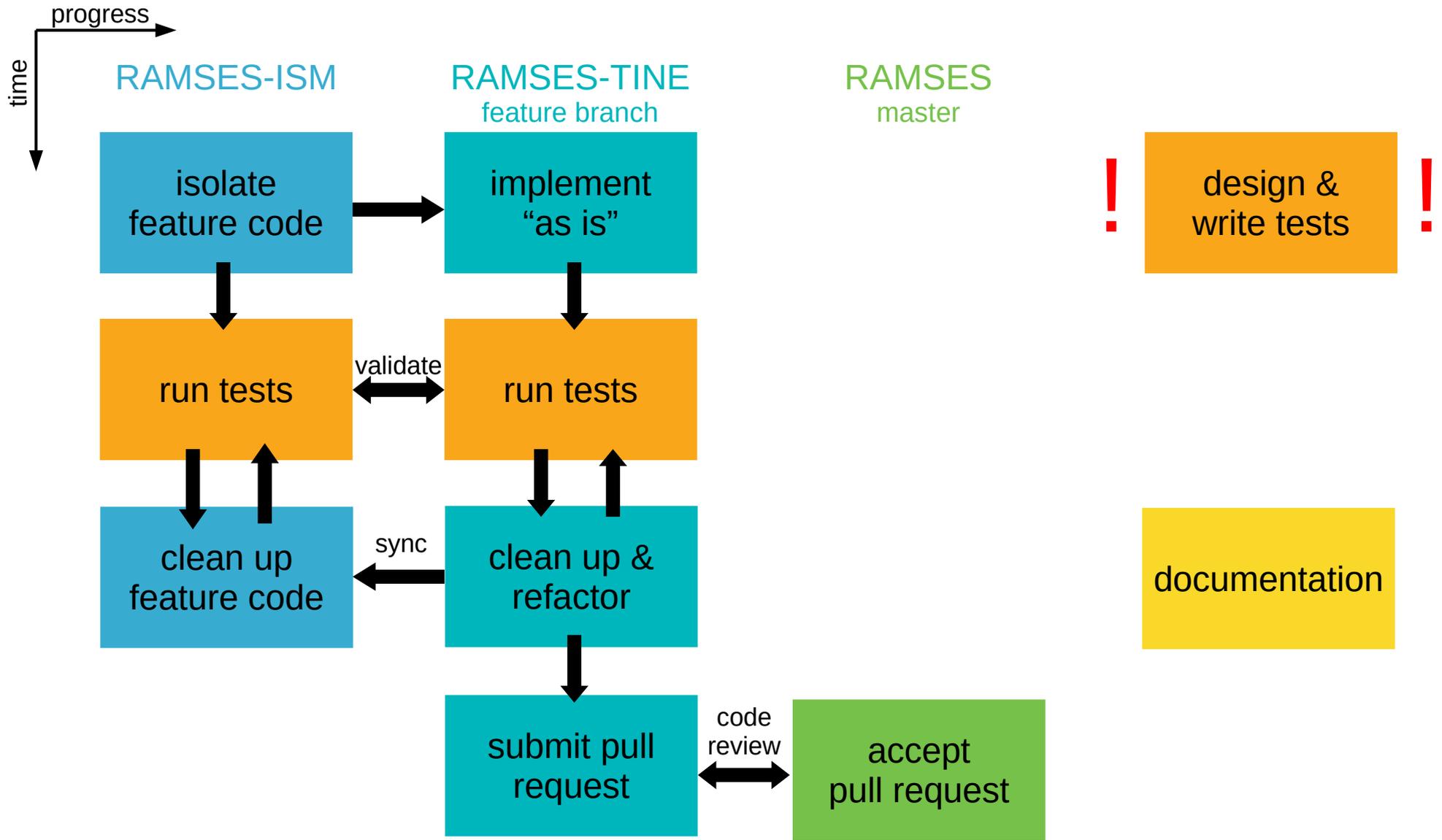






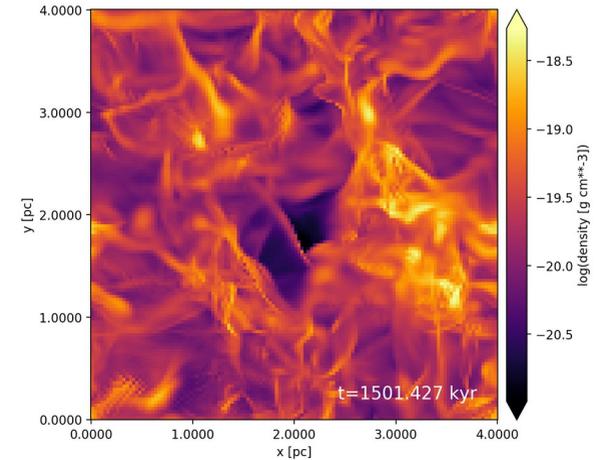
### Guidelines for features

- Isolated as possible
- able to switch on/off
- separate files for feature routines & variable declarations
- minimize code duplication
- its own documentation & tests
- input/output format compatibility



# Example: Turbulence driving

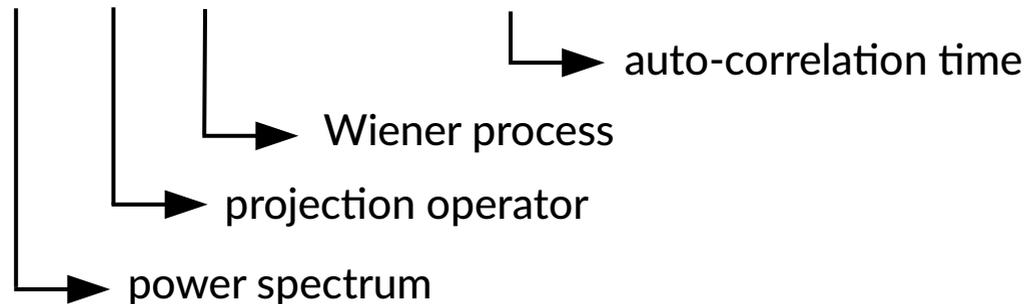
- Adds forcing to box to generate/maintain turbulence
- Andrew McLeod, in 2014
- Was easy because standalone module
- Dependency: fftw3
- Implementation:



generalised Ornstein-Uhlenbeck process

random growth      decay

$$dF(k, t) = \underbrace{F_0(k)P(k)}_{\text{random growth}} dW_t - \underbrace{F(k, t)}_{\text{decay}} dt/T$$



```

amr/amr_step.f90 MODIFIED
...
10 10 use coolrates_module, only: update_coolrates_tables
11 11 use rt_cooling_module, only: update_UVrates
12 12 #endif
13 13 +#if USE_TURB==1 preprocessor directive
14 14 + use turb_commons
15 15 + #endif
13 16 use mpi_mod
14 17 implicit none
15 18 #ifndef WITHOUTMPI
...
286 289 if(rt .and. rt_star) call update_star_RT_feedback(ilevel)
287 290 #endif
288 291
292 292 +#if USE_TURB==1
293 293 + ! Compute turbulent forcing
294 294 + call timer('turb','start')
295 295 + if (turb .and. turb_type/=3) then
296 296 + ! Calculate turbulent acceleration on each cell in this level
297 297 + call calc_turb_forcing(ilevel)
298 298 + end if
299 299 + #endif
300 300 +
289 301 !-----
290 302 ! Compute new time step
291 303 !-----
...
384 396 call timer('poisson','start')
385 397 if(poisson)call synchro_hydro_fine(ilevel,+0.5*dtnew(ilevel))
386 398
387 399 +#if USE_TURB==1
398 400 + ! Compute turbulent forcing
399 400 + call timer('turb','start')
401 401 + if (turb .AND. turb_type/=3) then
402 402 + ! Euler step, adding turbulent acceleration
403 403 + call add_turb_forcing(ilevel,dtnew(ilevel))
404 404 + end if
405 405 +
406 406 +
407 407 + #endif
408 408 +
387 409 ! Restriction operator
388 410 call timer('hydro upload fine','start')
389 411 call upload_fine(ilevel)
...

```

## Files changed (23)

+6	-0	M	amr/adaptive_loop.f90	← minimal changes to existing files
+22	-0	M	amr/amr_step.f90	
+14	-0	M	amr/output_amr.f90	
+3	-0	M	amr/read_params.f90	
+16	-0	M	amr/update_time.f90	
+22	-1	M	bin/Makefile	
+13	-0	M	hydro/courant_fine.f90	
+11	-0	M	hydro/init_flow_fine.f90	
+13	-0	M	mhd/courant_fine.f90	
+11	-0	M	mhd/init_flow_fine.f90	
+10	-0	M	pm/newdt_fine.f90	
+11	-0	A	turb/README	← New directory for feature
+216	-0	A	turb/add_turb_forcing.f90	
+215	-0	A	turb/init_turb.f90	
+47	-0	A	turb/mpi_share_turb_fields.f90	
+57	-0	A	turb/read_turb_fields.f90	
+58	-0	A	turb/read_turb_params.f90	
+44	-0	A	turb/turb_check_time.f90	
+31	-0	A	turb/turb_commons.f90	
+770	-0	A	turb/turb_force_utils.f90	
+35	-0	A	turb/turb_next_field.f90	
+54	-0	A	turb/turb_parameters.f90	
+59	-0	A	turb/write_turb_fields.f90	

## Makefile:

```
# Use turbulence? 1=Yes, 0=No (requires fftw3)
USE_TURB=1
```

```
ifeq ($(USE_TURB),1)
  LIBS_TURB = -L$(HOME)/local/lib -lfftw3
  LIBS_OBJ_TURB = -I$(HOME)/local/include -lfftw3
endif
```

## Namelist:

```
&TURB_PARAMS
turb=.true.
turb_type=1          ! 1=forced evolving, 2=forced fix
turb_seed=1         ! -1 = random
comp_frac=0.33      ! compressive fraction
instant_turb=.true. ! Generate initial turbulence before start
forcing_power_spectrum='parabolic' !powerlaw, parabolic
turb_T=0.001        ! Turbulent velocity autocorrelation time
turb_Ndt=100        ! Number of timesteps per autocorrelation time
turb_rms=1.e4        ! rms turbulent forcing acceleration
turb_min_rho=1d-50 ! Minimum density for turbulence
```

## Wiki

[ramses / TurbulenceDriving](#)

### Turbulence driving

The block named `&TURB_PARAMS` contains the parameters related to the turbulence driving. Originally implemented by Andrew Mcleod

#### Overview of parameters

Variable name	Fortran type	Default value	Description
<code>turb</code>	boolean	<code>.false.</code>	Turn on or off driving
<code>turb_seed</code>	integer	<code>-1</code>	Random number generator seed. -1 = random
<code>turb_type</code>	integer	<code>1</code>	How the driving changes over time. 1=driven evolving, 3=decaying
<code>instant_turb</code>	boolean	<code>.true.</code>	Generate initial turbulence before start
<code>comp_frac</code>	float	<code>0.3333</code>	The weight of compressive over solenoidal modes
<code>turb_T</code>	float	<code>1</code>	Turbulent velocity auto-correlation time in code units.
<code>turb_Ndt</code>	integer	<code>100</code>	Number of timesteps per auto-correlation time
<code>turb_rms</code>	float	<code>1</code>	Root-mean-square turbulent forcing in code units.
<code>turb_min_rho</code>	float	<code>1d-50</code>	Minimum density for turbulence. Not forcing is added onto cells with a density less than this value.
<code>forcing_power_spectrum</code>	string	<code>parabolic</code>	Power spectrum type of the forcing, which describes the relative strength of individual modes. Options are: <code>power_law</code> , <code>parabolic</code> , <code>konstantin</code>

# Test case

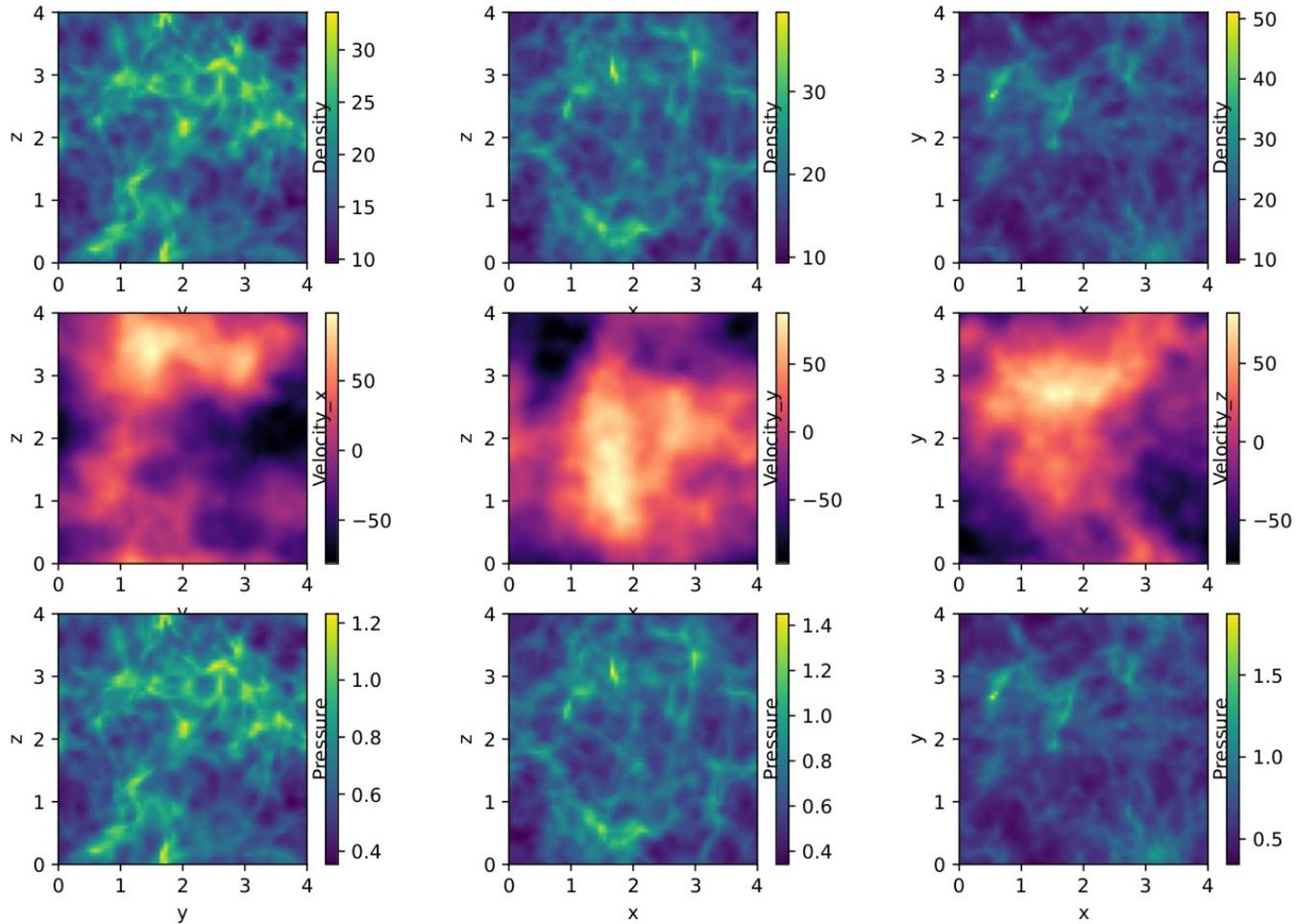


Figure 1: turb/driving test

amr	
adaptive_loop.f90	+6
amr_step.f90	+24 -3
output_amr.f90	+14
read_params.f90	+3
update_time.f90	+16
bin	
Makefile	+25 -3
hydro	
courant_fine.f90	+13
init_flow_fine.f90	+11
synchro_hydro_fine.f90	+42 -10
mhd	
courant_fine.f90	+13
init_flow_fine.f90	+11
synchro_hydro_fine.f90	-106
pm	
newdt_fine.f90	+10
tests (5)	
turb	
+ README	+11
+ add_turb_forcing.f90	+99
+ init_turb.f90	+202
+ mpi_share_turb_fields.f90	+43
+ read_turb_fields.f90	+55
+ read_turb_params.f90	+62
+ turb_check_time.f90	+42
+ turb_commons.f90	+31
+ turb_force_utils.f90	+726
+ turb_next_field.f90	+33
+ turb_parameters.f90	+54
+ write_turb_fields.f90	+57

← Final PR

# Turbulence driving update

<https://bitbucket.org/rteyssie/ramses/pull-requests/447>

Now available in main ramses!

[WIP] More options to select modes

[WIP] User guide

- description implementation
- what parameters to choose

## TO DO



put back threshold sink accretion scheme

adapt test suite to allow restarts



add option to restart with different number of cpus

Port jets

Interaction tracer particles with sinks

Port FLD (ism)

Port Cosmic Ray diffusion (ism)

General EOS

+ Add a card



## In progress



Option to output electric current in ideal MHD

2D turbulence

options to choose turbulence driving modes



TC

Port FRIGG cooling & couple to H2 formation

Port stellar particles & sink feedback (ism)

Port non-ideal MHD (ism)

Port dust (Ugo)

+ Add a card



## Done



Port turbulence driving (ism)

Port tracer particles (Corentin)

Easier analytic equation of state

+ Add a card



TO DO

In progress

Done

Masson et al 2012  
Marchand et al 2018

Ambipolar diffusion  $\checkmark$

Ohmic dissipation  $\checkmark$

Hall effect  $\times$

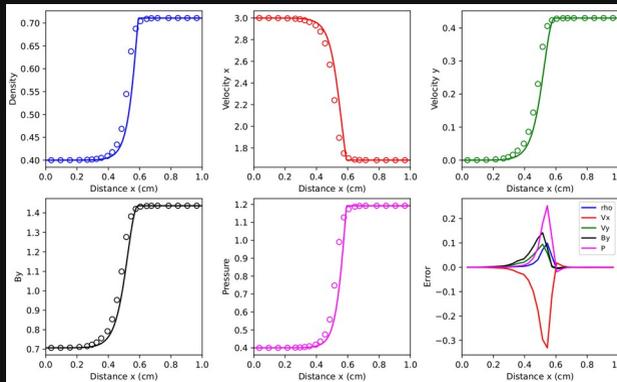


Figure 2: nimhd/nimhd-shock-ohm test

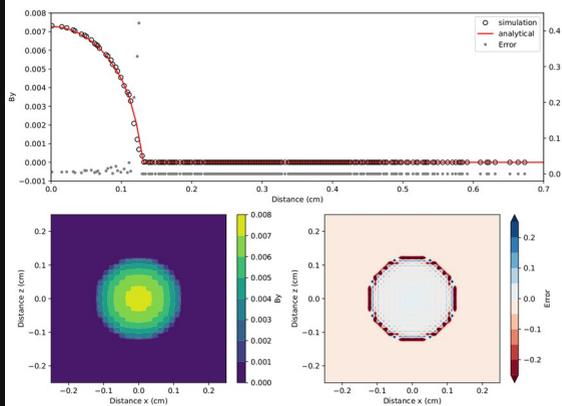


Figure 3: nimhd/nimhd-diffusion-ad test

Table 3: nimhd-diffusion-ad error summary

Variable	This run	Reference	Error	Tolerance
B x left	0.00000000000000e+00	0.00000000000000e+00	0.00000000000000e+00	2.99999999999998e-13
B x right	0.00000000000000e+00	0.00000000000000e+00	0.00000000000000e+00	2.99999999999998e-13
B y left	6.399999999999370e-01	6.40000000000001e+01	9.72555057173154e-14	2.99999999999998e-13
B y right	6.399999999999370e-01	6.40000000000001e+01	9.72555057173154e-14	2.99999999999998e-13
B z left	0.00000000000000e+00	0.00000000000000e+00	0.00000000000000e+00	2.99999999999998e-13
B z right	0.00000000000000e+00	0.00000000000000e+00	0.00000000000000e+00	2.99999999999998e-13
hoshen	1.00000000000000e+00	1.00000000000000e+00	0.00000000000000e+00	2.99999999999998e-13
current x	7.586446588798568e+02	7.592928512366497e+02	8.54408386974322e-04	2.99999999999998e-13
current y	0.00000000000000e+00	0.00000000000000e+00	0.00000000000000e+00	2.99999999999998e-13
current z	7.586446588798568e+02	7.592928512366497e+02	8.54408386974322e-04	2.99999999999998e-13
density	0.00000000000000e+00	0.00000000000000e+00	0.00000000000000e+00	2.99999999999998e-13
dx	1.21300000000000e-03	1.21300000000000e-03	0.00000000000000e+00	2.99999999999998e-13
level	3.40128000000000e+05	3.40128000000000e+05	0.00000000000000e+00	2.99999999999998e-13
ncells	5.94240000000000e+04	5.94240000000000e+04	0.00000000000000e+00	2.99999999999998e-13
pressure	-2.264719684267732e+04	0.00000000000000e+00	inf	2.99999999999998e-13
time	5.402496678193103e+01	5.021153116957440e+01	3.803220180117077e-03	2.99999999999998e-13
unit d	1.00000000000000e+00	1.00000000000000e+00	0.00000000000000e+00	2.99999999999998e-13
unit l	1.00000000000000e+00	1.00000000000000e+00	0.00000000000000e+00	2.99999999999998e-13
unit t	1.00000000000000e+00	1.00000000000000e+00	0.00000000000000e+00	2.99999999999998e-13
velocity x	0.00000000000000e+00	0.00000000000000e+00	0.00000000000000e+00	2.99999999999998e-13
velocity y	0.00000000000000e+00	0.00000000000000e+00	0.00000000000000e+00	2.99999999999998e-13
velocity z	0.00000000000000e+00	0.00000000000000e+00	0.00000000000000e+00	2.99999999999998e-13
x	2.98275000000000e+04	2.98275000000000e+04	0.00000000000000e+00	2.99999999999998e-13
y	2.97120000000000e+04	2.97120000000000e+04	0.00000000000000e+00	2.99999999999998e-13
z	2.98275000000000e+04	2.98275000000000e+04	0.00000000000000e+00	2.99999999999998e-13

Port turbulence driving (ism)

Port tracer particles (Corentin)

Easier analytic equation of state

+ Add a card

Port Cosmic Ray diffusion (ism)

General EOS

+ Add a card

Port non-ideal MHD (ism)

Port dust (Ugo)

+ Add a card

## TO DO

put back threshold sink accretion

## In progress

Option to output electric current in

## Done

Port turbulence driving (ism)

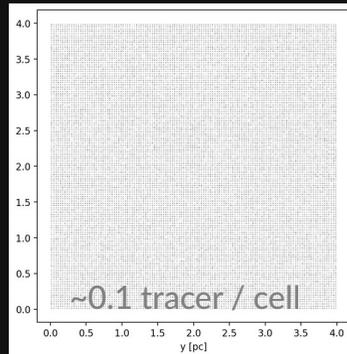
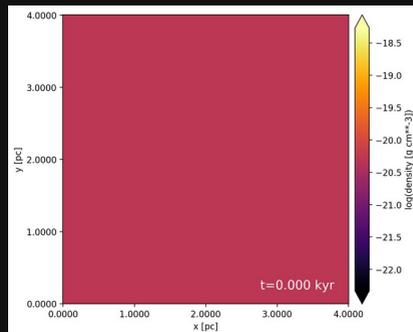
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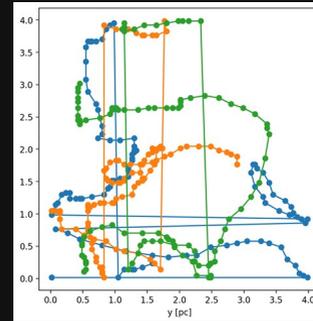
+ Add a card

### Cadiou et al 2019

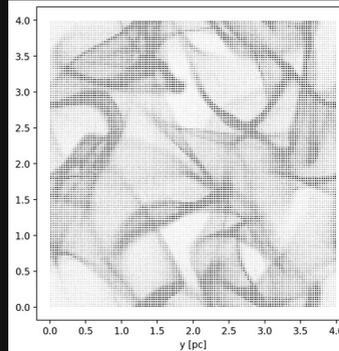
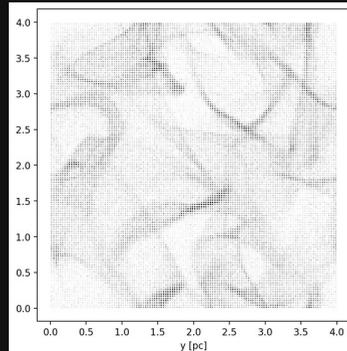
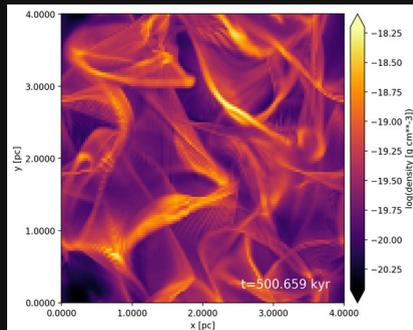
```
&TRACER_PARAMS
tracer=.true.
MC_tracer=.true.
tracer_feed_fmt='inplace' !can also initialize from file
tracer_mass=0.00001 !this gives about 1 particle per cell
/
```



### Example trajectories



~1 tracer / cell



## TO DO

put back threshold sink accretion scheme

adapt test suite to allow restarts

add option to restart with different number of cpus

Port jets

Interaction tracer particles with sinks

Port FLD (ism)

Port Cosmic Ray diffusion (ism)

General EOS

+ Add a card

## In progress

Option to output electric current in ideal MHD

2D turbulence

options to choose turbulence driving modes

Port F H2 for

Port s feedb

Port n

Port d

+ Add a card

## Done

Port turbulence driving (ism)

Port tracer particles (Corentin)

Easier analytic equation of state

### COOLING\_PARAMS

isothermal	logical	.false.	(deprecated) Enable equation of state for gas (heating and cooling disabled if <code>.true.</code> .)
barotropic_eos	logical	.false.	Enable barotropic equation of state for gas (heating and cooling disabled if <code>.true.</code> ). Replaced 'isothermal'
barotropic_eos_form	string	legacy	Type of barotropic EOS. Options: isothermal, polytrope, double_polytrope, custom, legacy
polytrope_rho	real	1.0d50	sets rho0 in EOS (density normalisation or knee-density), in g/cm3
polytrope_index	real	1.0d0	sets gamma in EOS (polytropic index)
T_eos	real	10	sets T0 in EOS (isothermal temperature or temperature normalisation), in K
mu_gas	real	1d0	molecular weight

+ hydro / eos.f90

# How can you help?

- Use git branches instead of patches for development
- Add your typical (low resolution) setup as a test case
- Merge your features into the main repo

Join RAMSES merge slack (or make a specific feature channel on the teamramses slack)

[https://join.slack.com/t/mergingramsesversions/shared\\_invite/zt-wat3j23a-~Q01jBM~9x7douyYJOX2wA](https://join.slack.com/t/mergingramsesversions/shared_invite/zt-wat3j23a-~Q01jBM~9x7douyYJOX2wA)