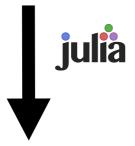


Existing Packages:

- Two Language problem: Python + C++
- Slow: e.g. Python slow in reading RAMSES data
- Fast, command line, difficult to extend
- use interpolated full grid —> huge memory consumption
- Effort to program personal extensions
- Not shared with community



Solution:

- -one language
- —high level
- —easy to use, extend and share with community
- -interactive
- -fast
- —memory leightweight (AMR)
- —easy to install
- Workflow enables to concentrate more on science
- MERA can be used on with other grid-based or N-body codes



Based on a database

Multiple dispatch

Predefined quantities

Macros

Fast projections

Compressed MERA files

Detailed documentation

Platform to easily share your notebooks

Little paper

[15]:	select	(gas.da	ta, (:	level,	:cx, :cy, :cz, :rho))
[15]:	Table level	with 50 cx	553133 cy	rows,	5 columns: rho
	6	1	1	1	1.47746e-8
	6	1	1	2	1.47746e-8
	6	1	1	3	1.47746e-8
	6	1	1	4	1.47746e-8
	6	1	1	5	1.47746e-8
	6	1	1	6	1.47746e-8
	6	1	1	7	1.47746e-8
	6	1	1	8	1.47746e-8
	6	1	1	9	1.47746e-8
	6	1	1	10	1.47746e-8
	6	1	1	11	1.47746e-8
	6	1	1	12	1.47746e-8
	:				
	14	13038	8855	8122	1.41561
	14	13038	8855	8123	1.29724
	14	13038	8855	8124	1.27011
	14	13038	8856	8117	1.89495
	14	13038	8856	8118	1.83967
	14	13038	8856	8119	1.67429
	14	13038	8856	8120	1.59913
	14	13038	8856	8121	1.45658
	14	13038	8856	8122	1.41468
	14	13038	8856	8123	1.3037
	14	13038	8856	8124	1.29413

Advantages e.g.:

- easy to filter data
- easy to parallize



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One single function name For different data-types

```
function projection(
    dataobject::HydroDataType, vars::Array{Symbol,1};
    units::Array{Symbol,1}=[:standard],
    #coordinates::Symbol=:cartesian,
    lmax::Number=dataobject.lmax,
    mask=[false],
    direction::Symbol=:perpendicular,
    weighting::Bool=true,
    mode::Symbol=:weighting,
    gamm::Number=1.4,
    xrange::Array{<:Any,1}=dataobject.ranges[1:2],
    yrange::Array{<:Any,1}=dataobject.ranges[3:4],
    zrange::Array{<:Number,1}=[0., 0., 0.],
    range_units::Symbol=:standard,
    data_center::Array{<:Number,1}=[0.5, 0.5, 0.5],
    data_center_units::Symbol=:standard,
    verbose::Bool=verbose_mode)</pre>
```

```
# exported
mutable struct HydroDataType <: DataSetType
    data::JuliaDB.AbstractIndexedTable #todo: check
    info::InfoType
    lmin::Int
    lmax::Int
    boxlen::Float64
    ranges::Array{Float64,1}
    selected_hydrovars::Array{Int,1}
    used_descriptors::Dict{Any,Any}
    smallr::Float64
    smallc::Float64
    scale::ScalesType
    HydroDataType() = new()
end</pre>
```

```
Particles
```

```
# exported
mutable struct PartDataType <: DataSetType
    data::JuliaDB.AbstractIndexedTable #todo:check
    info::InfoType
    lmin::Int
    lmax::Int
    boxlen::Float64
    ranges::Array{Float64,1}
    selected_partvars::Array{Symbol,1}
    scale::ScalesType
    PartDataType() = new()
end</pre>
```



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```
projection()
Predefined Quantities for Projections:
             =======[gas]:======
       -all the non derived hydro vars-
:cpu, :level, :rho, :cx, :cy, :cz, :vx, :vy, :vz, :p, var6,...
              -derived hydro vars-
:x, :y, :z
:sd or :Σ or :surfacedensity
:mass, :cellsize, :freefall_time
gamm needed => :cs, :mach, :jeanslength, :jeansnumber
 ===========[particles]:============
        all the non derived vars:
:cpu, :level, :id, :x, :y, :z, :vx, :vy, :vz, :mass, :age,....
=========[gas or particles]:=========
:v, :ekin
squared => :vx2, :vy2, :vz2
velocity dispersion \Rightarrow \sigma x, \sigma y, \sigma z, \sigma
related to a given center:
:vr_cylinder, vr_sphere
:vΦ, :vθ
squared => :vr_cylinder2, :v\u00f32
velocity dispersion \Rightarrow \sigma r_cylinder, \sigma \phi
:l, :lx, :ly, :lz :lr, :lφ, :lθ
2d maps:
:r_cylinder, :r_sphere
:φ, :θ
```

Quantities internally defined, And can be used by many functions In MERA or extended by the user.



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Macros

```
Use Pipeline Macros
Fast project
               boxlen = info.boxlen
               cv = boxlen/2.
Compress
               density = 3. /gas.scale.Msol_pc3
               radius = 3. /gas.scale.kpc
               height = 2. /gas.scale.kpc
Detailed d
               filtered_db = @apply gas.data begin
                    @where :rho >= density
Platform to
                    @where sqrt( (:cx * boxlen/2^:level - cv)^2 + (:cy * boxlen/2^:level - cv)^2 ) <= radius
                    @where abs(:cz * boxlen/2^:level -cv) <= height</pre>
Little pape
               var_filtered = getvar(gas, :mass, filtered_db=filtered_db, unit=:Msol)
               sum(var_filtered) # [Msol]
```



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Is fast due to:

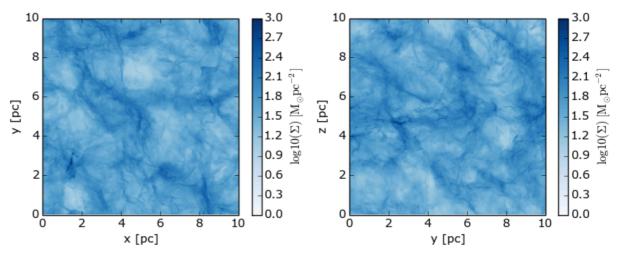
- AMR data only
- Julia
- and a trick

Fast high resolution projections

```
onto a (2^14)^2 grid !!!
```

```
proj z = projection(gas, vars=[:sd], direction=:z);
[Mera]: 2018-12-29T16:54:18.666
domain
xmin::xmax: 0.0 :: 1.0
                                ==> 0.0 [pc] :: 10.0 [pc]
ymin::ymax: 0.0 :: 1.0
                                ==> 0.0 [pc] :: 10.0 [pc]
                                ==> 0.0 [pc] :: 10.0 [pc]
zmin::zmax: 0.0 :: 1.0
Selected vars: Symbol[:sd]
                                                          Time: 0:02:08
100%
proj_x = projection(gas, vars=[:sd], direction=:x);
[Mera]: 2018-12-29T16:56:29.763
domain
xmin::xmax: 0.0 :: 1.0
                                ==> 0.0 [pc] :: 10.0 [pc]
                                ==> 0.0 [pc] :: 10.0 [pc]
ymin::ymax: 0.0 :: 1.0
zmin::zmax: 0.0 :: 1.0
                                ==> 0.0 [pc] :: 10.0 [pc]
Selected vars: Symbol[:sd]
                                                          Time: 0:02:10
```

Moeckels et al. - turbulent box





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Reduces significantly

- disc space: factor 5 (my simulations)
- data loading: factor 17 (2048 cores sim)

-> good for time-sequence Post-processing



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Detailed documentation

Mera.jl Search docs Home First Steps Simulation Overview Fields **Physical Units Physical Constants** Disc Space 1-Data Inspection Hydro **Particles** Clumps 2-Load by Selection Hydro **Particles** Clumps 3-Get Subregions

» First Steps Clit on GitHub

First Steps

Simulation Overview

The first call of MERA will compile the package.

Get information about the simulation for a given output number and assign it to an object (composite type) here: "info". By default, the RAMSES output folders are assumed to be in the current working directory, and the relative or absolute path can be given. The information is read from several files: info-file, the header-file, from the header of the Fortran binary files of the first CPU (hydro, grav, part, clump, if they exist), etc. Many familiar names and acronyms known from RAMSES are maintained. The printed units are automatically adapted to a human-readable representation (see getinfo):

```
info = getinfo(); # default: output=1 in current folder,
```

Documentation contains detailed explanations, many examples and Jupyter notebooks to try it on your simulation data.

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