

Galactic dynamics with



part 1: foundations

Eugene Vasiliev

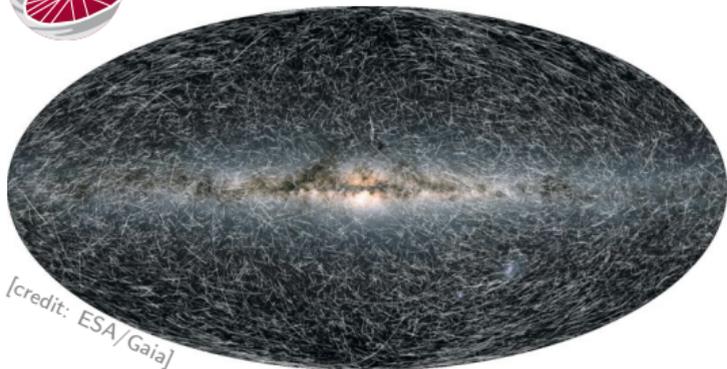
Galaxy Modelling & Galactic centre workshop

University of Surrey, December 2024

Observational and theoretical context



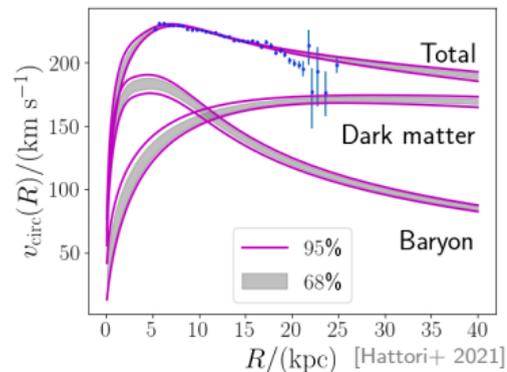
gaia



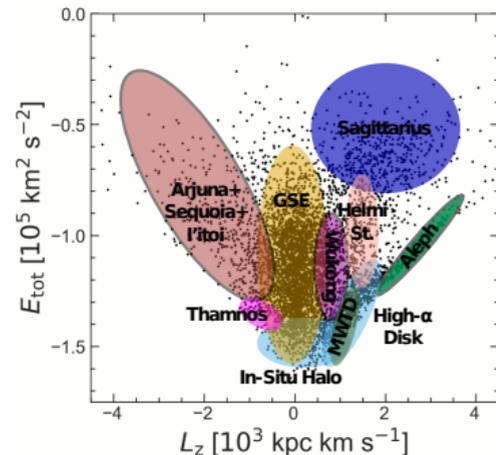
[credit: ESA/Gaia]



Milky Way rotation curve

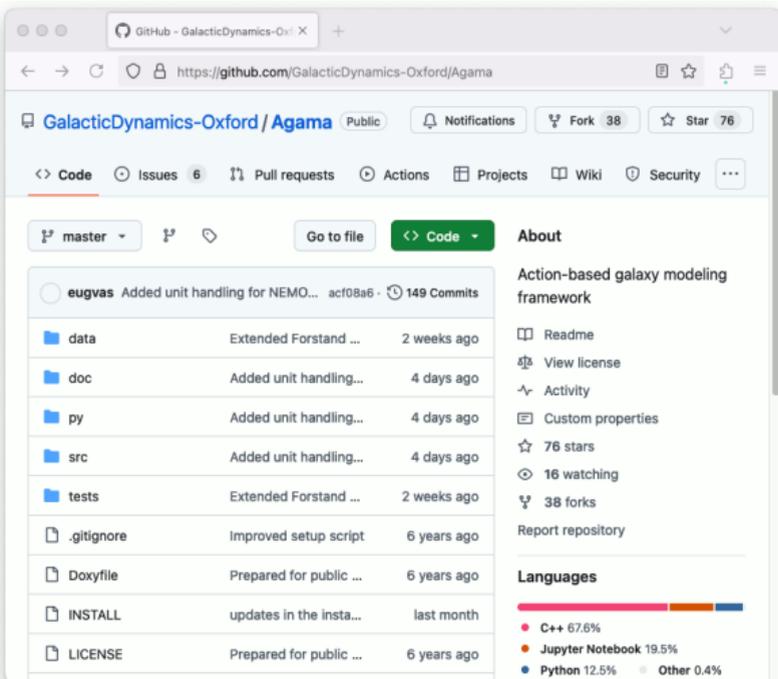


Classification of accreted stars



[Naidu+ 2020]

Agama software package



Monthly Notices

of the
ROYAL ASTRONOMICAL SOCIETY

MNRAS **482**, 1525–1544 (2019)

Advance Access publication 2018 October 3

AGAMA: action-based galaxy modelling architecture

Eugene Vasiliev ^{1,2,3}★

development started in 2015;
code paper published in 2018;
used in $\gtrsim 350$ publications.

Main features:

- ▶ core library written in C++;
- ▶ OpenMP parallelization;
- ▶ hand-made Python interface;
- ▶ extensible with user-defined functions;
- ▶ several dozen tests and example programs in C++ and Python;
- ▶ detailed documentation (~ 140 pages);
- ▶ $\sim 80\,000$ lines of code.

Structure of the Agama library

some maths stuff (C & Fortran)

GSL (C) & Eigen (C++) math libs

CVXOPT quad. opt. solver (Python)

C++ core \implies shared library `agama.so`

computationally heavy parts (potentials, actions, orbit integration, etc.);
built-in density, potential and DF models

Python interface (C & C++)

vectorization (operation on arrays);
OpenMP parallelization

external gravity
source interfaces
for N -body codes

C, Fortran
and Julia
interfaces

additional Python routines

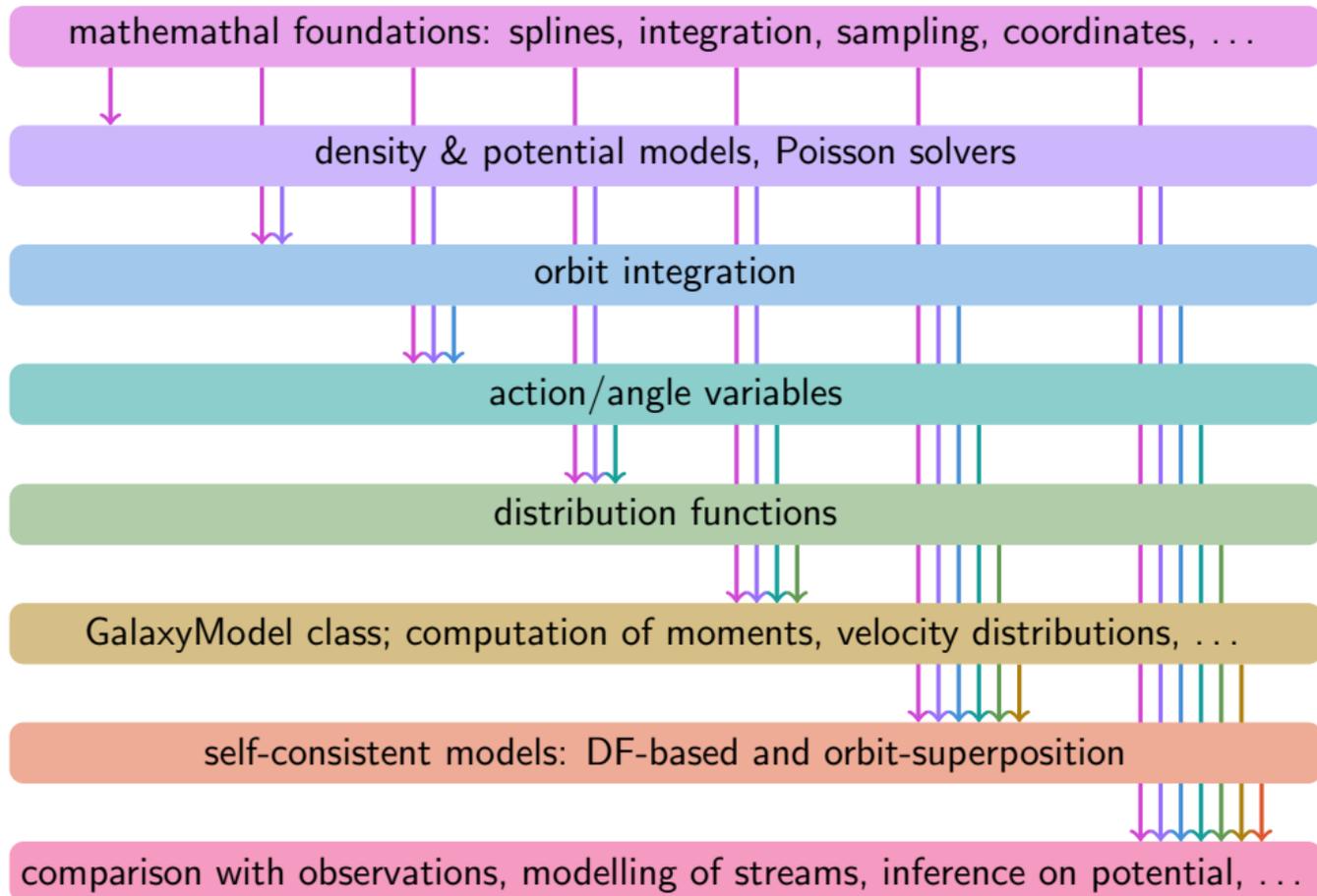
math & coord utility functions;
Forstand orbit-superposition code;
interfaces for GALPY and GALA

NEMO, AMUSE, Gadget4, Arepo

AGAMA potentials can be used in
GALPY and GALA

user scripts in Python (including custom density, potential & DF models)

Structure of the Agama library



Gravitational potential

Task: given the density profile $\rho(\mathbf{x})$, determine the potential $\Phi(\mathbf{x})$ from the Poisson equation: $\nabla^2\Phi = 4\pi G \rho$.

Example 1: spherical Plummer model

$$\rho(r) = \frac{3M}{4\pi a^3 (1 + r^2/a^2)^{5/2}} \implies \Phi(r) = -\frac{GM}{\sqrt{r^2 + a^2}}.$$

Example 2: triaxial Hernquist model

$$\rho(x, y, z) = \frac{M}{2\pi abc s (1 + s)^3}, \quad s \equiv \sqrt{\frac{x^2}{a^2} + \frac{y^2}{b^2} + \frac{z^2}{c^2}} \implies$$
$$\Phi(x, y, z) = -GM \int_0^\infty d\tau \frac{\left(1 + \sqrt{\frac{x^2}{a^2 + \tau} + \frac{y^2}{b^2 + \tau} + \frac{z^2}{c^2 + \tau}}\right)^{-2}}{2\sqrt{(a^2 + \tau)(b^2 + \tau)(c^2 + \tau)}}.$$

It gets very complicated very quickly!

Gravitational potential

Commonly used analytic potential–density pairs: Plummer, NFW, MiyamotoNagai, Dehnen, Ferrers ...

If one needs more flexibility, there are three general-purpose Poisson solvers:

0. Direct integration:

$$\Phi(\mathbf{x}) = - \iiint d^3x' \rho(\mathbf{x}') \times \frac{G}{|\mathbf{x} - \mathbf{x}'|}. \quad (\text{impractical})$$

1. Azimuthal-harmonic expansion (CylSpline):

$$\Phi(R, z, \phi) = \sum_{m=-\infty}^{\infty} \Phi_m(R, z) e^{im\phi}.$$

2. Spherical-harmonic expansion (Multipole):

$$\Phi(r, \theta, \phi) = \sum_{l=0}^{\infty} \sum_{m=-l}^l \Phi_{lm}(r) Y_l^m(\theta, \phi).$$

interpolated functions

3. BasisSet expansion (a.k.a. self-consistent field method of Hernquist&Ostriker 1992):

$$\Phi(r, \theta, \phi) = \sum_{n=0}^{\infty} \sum_{l=0}^{\infty} \sum_{m=-l}^l \Phi_{nlm} A_{nl}(r) Y_l^m(\theta, \phi).$$

Gravitational potential

Workflow for the Multipole potential expansion:

original $\rho(r, \theta, \phi)$

\rightsquigarrow

approximate $\sum_{\ell=0}^{\ell_{\max}} \sum_{m=-\ell}^{\ell} \rho_{\ell m}(r) Y_{\ell}^m(\theta, \phi)$

solve Poisson eqn for each term

$$\Phi_{\ell m}(r) = -\frac{4\pi G}{2\ell + 1} \left[r^{-\ell-1} \int_0^r \rho_{\ell m}(s) s^{\ell+2} ds + r^{\ell} \int_r^{\infty} \rho_{\ell m}(s) s^{1-\ell} ds \right]$$

$\sum_{\ell=0}^{\ell_{\max}} \sum_{m=-\ell}^{\ell} \Phi_{\ell m}(r) Y_{\ell}^m(\theta, \phi)$

$\Phi(r, \theta, \phi)$

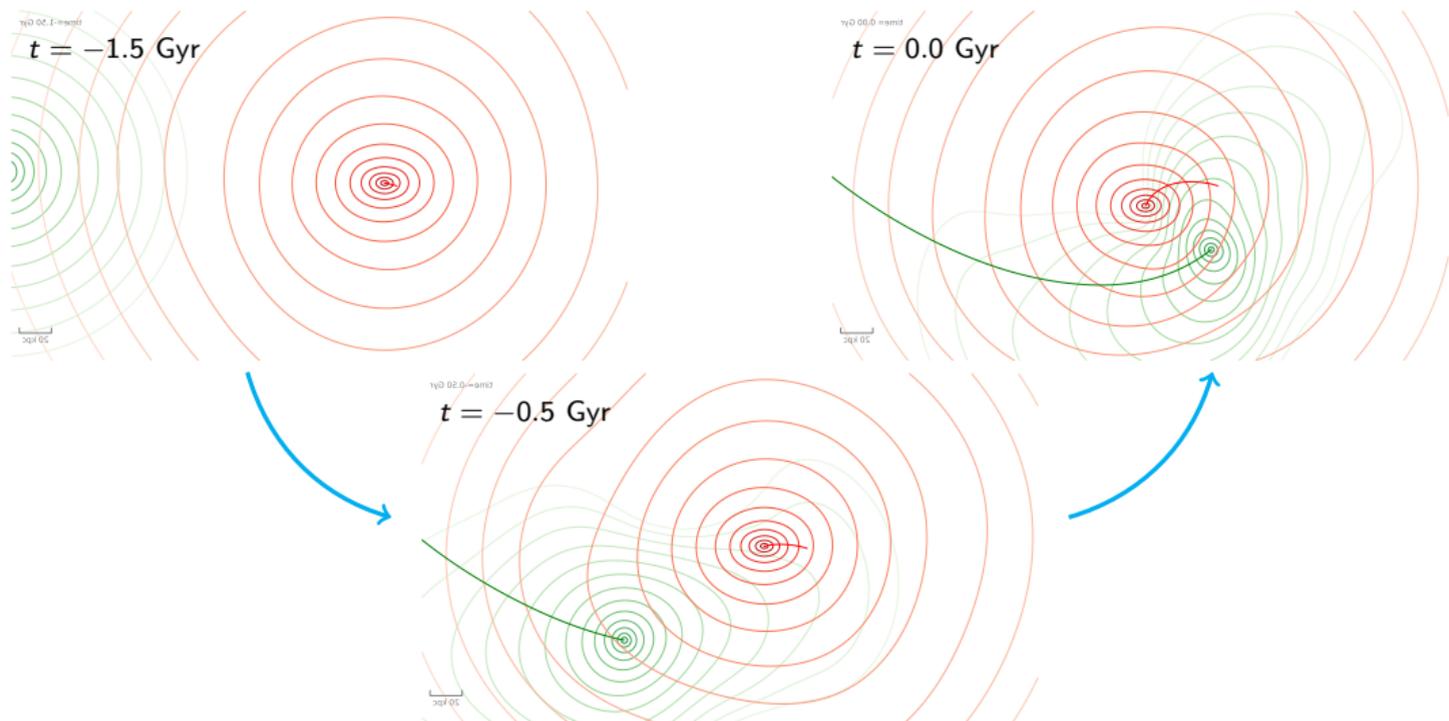
\parallel

Multipole and BasisSet potentials are well-suited for extended but not too flattened profiles; CylSpline is ideal for disk profiles, but has a finite spatial extent and doesn't like cusps.

Composite and time-dependent potentials

Potentials can be added, scaled or interpolated with time, rotated (e.g., bar or spiral arms), shifted along a time-dependent trajectory, etc.

Example: potentials of Milky Way and LMC extracted from an N -body simulation.



Gravitational potential: example 1

User-defined density model: a boxy bar $\rho(x, y, z) = \rho_0 \exp(-s^{1/n})$, where $s \equiv [(x/a)^k + (y/b)^k + (z/c)^k]^{1/k}$ is the generalized ellipsoidal radius (an ordinary ellipsoid has $k = 2$), and n is the Einasto index.

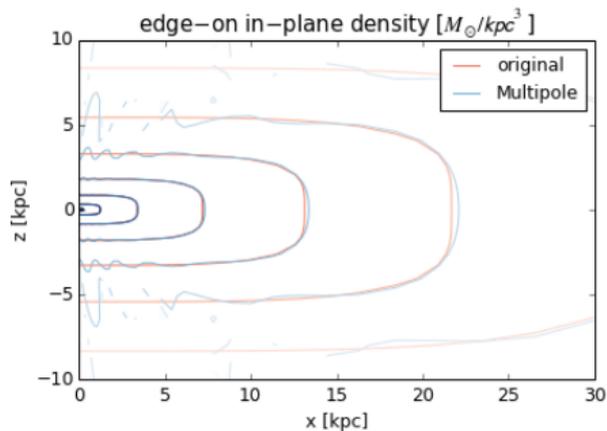
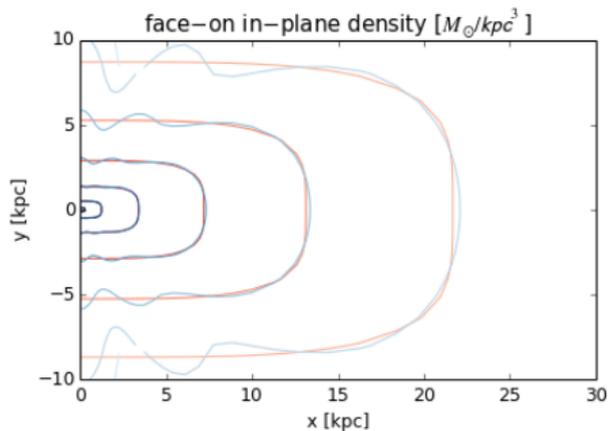
```
def dens_bar(xyz):
```

```
    x, y, z = abs(xyz).T
```

```
    s = ((x/a)**k + (y/b)**k + (z/c)**k)**(1./k)
```

```
    return rho0 * numpy.exp(-s**(1./n))
```

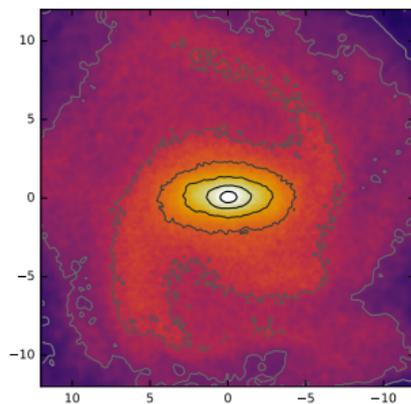
```
pot_bar = agama.Potential(type='Multipole', density=dens_bar,  
                           lmax=20, mmax=10, symmetry='triaxial')
```



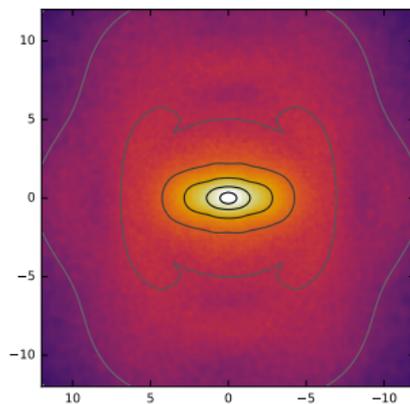
Gravitational potential: example 2

One may construct these potential expansions either from an analytic density profile (including any user-defined Python function for ρ or Φ) or from an N -body snapshot, specifying the desired level of symmetry.

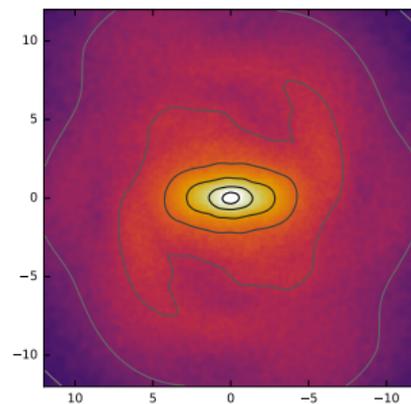
```
pot_nbody = agama.Potential(type='CylSpline',  
                             particles=(pos,mass), symmetry='bisymmetric')
```



original snapshot



triaxial



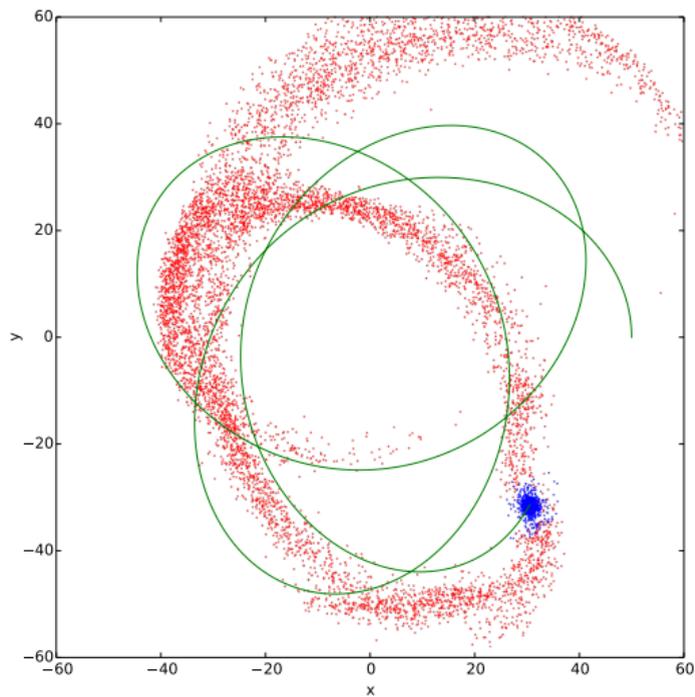
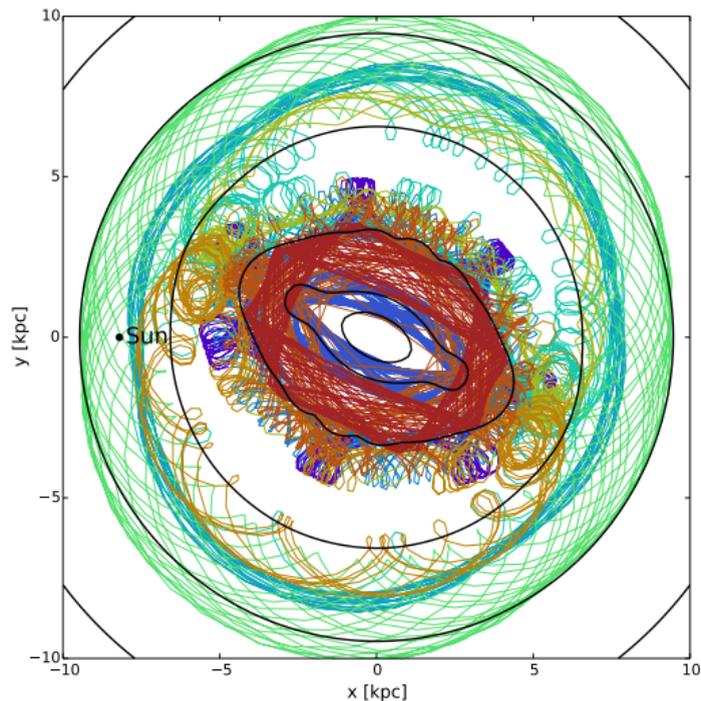
bisymmetric

Numerical integration of orbits

One of the most common tasks in galactic dynamics.

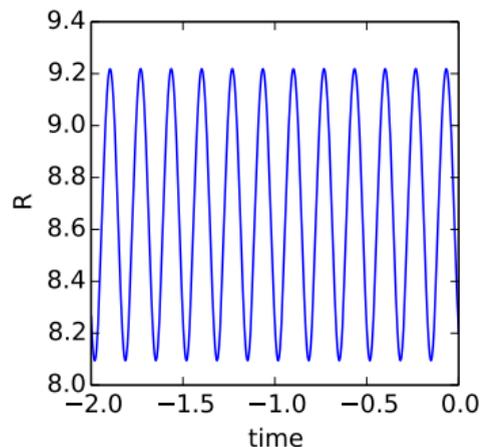
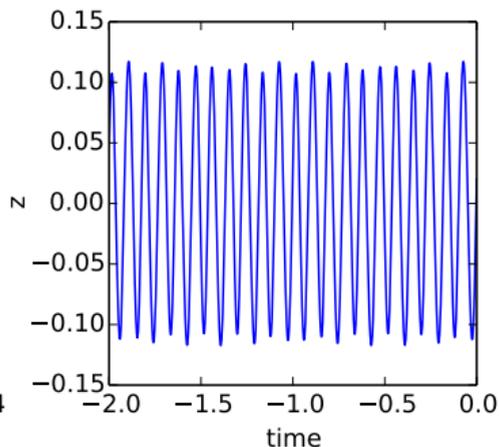
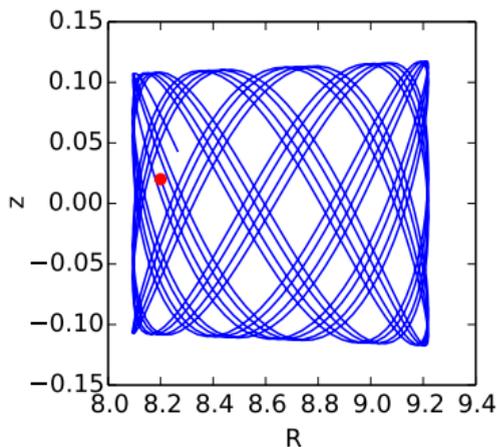
Examples: orbits in the MW bar;

test-particle simulations of a tidal stream



Numerical integration of orbits: example

```
agama.setUnits(length=1, velocity=1, mass=1)
#Note: distances are in kpc, velocities in km/s  $\Rightarrow$  time in kpc/(km/s)=0.978 Gyr
pot_mw = agama.Potential('McMillan17.ini')
time, traj = agama.orbit(potential=pot_mw,
    ic=[-8.2, 0, 0.02, 13, 245, 8], time=-2.0, trajsize=1001)
R = (traj[:,0]**2 + traj[:,1]**2)**0.5
z = traj[:,2]
```



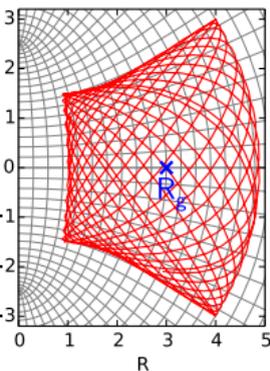
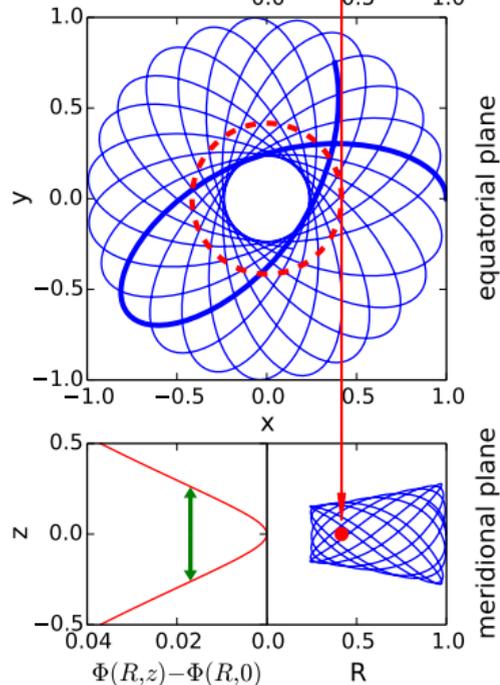
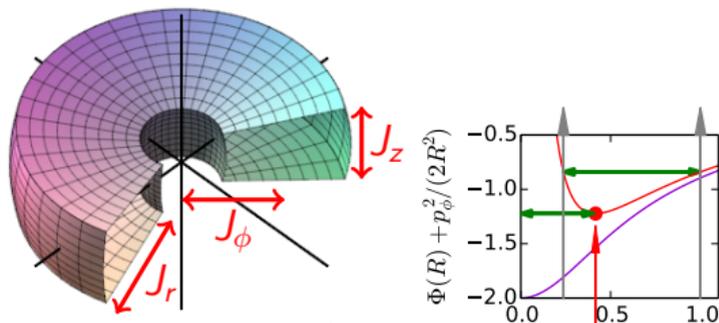
see `tutorial_potential_orbits.ipynb` for much more detail

Action-angle variables

Most orbits in axisymmetric potentials look like "rectangular tori" with three parameters defining the shape:

$J_\phi \equiv L_z = R_g v_{\text{circ}}(R_g)$ determines the overall size of the orbit ("guiding radius" R_g);
 J_R determines the extent of radial oscillations;
 J_z does the same for vertical oscillations.

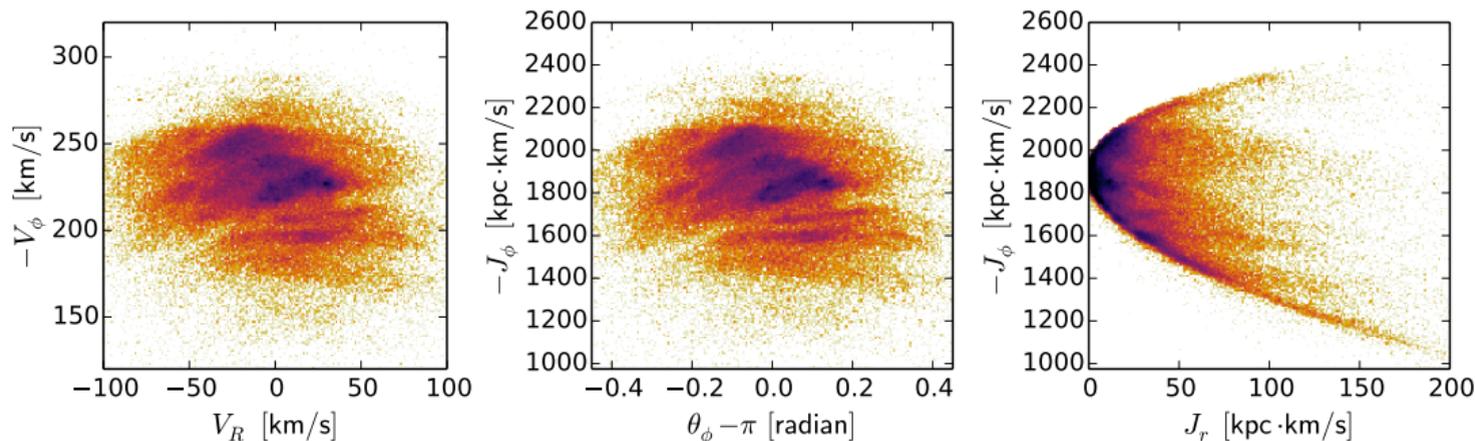
Corresponding phase angles $\theta_{\phi,R,z}$ determine the location on the orbit.



Actions are defined as $J_i \equiv \frac{1}{2\pi} \oint p_i dx_i$, and are computed in the Stäckel approximation [Binney 2012], using spheroidal coordinates for x_i . Inverse transformation $\{\mathbf{J}, \boldsymbol{\theta}\} \Rightarrow \{\mathbf{x}, \mathbf{v}\}$ is provided by the Torus code [Binney & McMillan 2017].

Action-angle variables: example

```
af = agama.ActionFinder(pot_mw)
#posvel is an array of 6d phase-space coords in Cartesian frame
actions, angles = af(posvel, angles=True, frequencies=False)
J_R, J_z, J_phi = actions.T
theta_R, theta_z, theta_phi = angles.T
```



Gaia sample of nearby stars ($D < 100$ pc)

other applications: clustering in the integrals space; study of resonances, ...

Distribution functions

DF $f(\mathbf{x}, \mathbf{v})$ offers a complete description of a stellar population.

Fundamental principle of stellar dynamics (Jeans's theorem):

in a steady state, DF must be a function of integrals of motion $f(\mathcal{I}(\mathbf{x}, \mathbf{v}; \Phi))$, and it is often convenient to use actions \mathbf{J} as integrals \mathcal{I} .

Two ways of constructing a DF in AGAMA:

- ▶ Using the Cuddeford–Eddington [anisotropic] inversion formula to obtain the DF of the form $f(E, L) = \hat{f}(E + L^2/(2r_a^2)) L^{-2\beta_0}$ corresponding to a given density $\rho(r)$ and potential $\Phi(r)$:

```
df_qs = agama.DistributionFunction(type='QuasiSpherical',  
    density=dens, potential=pot, beta0=0.3, r_a=10)
```

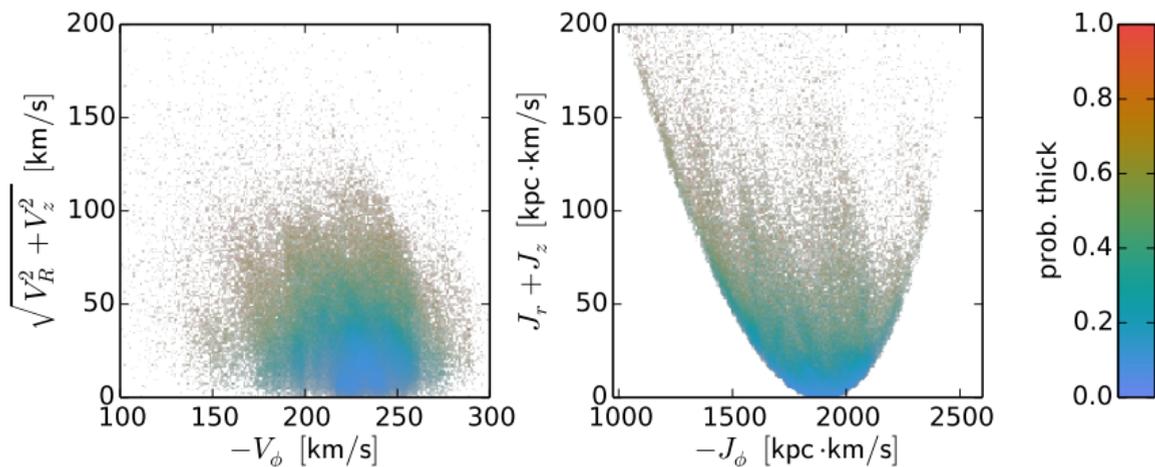
This DF is internally converted to the form $f(\mathbf{J})$.

- ▶ Using an *ad hoc* expression for the DF $f(\mathbf{J})$, either one of built-in models (DoublePowerLaw, QuasiIsothermal, Exponential) or a user-defined Python function:

```
df_dp = agama.DistributionFunction(type='DoublePowerLaw',  
    J0=J0, slopeIn=Gamma, slopeOut=Beta, mass=1)
```

Distribution function applications: classification

```
df_thin = agama.DistributionFunction(type='QuasiIsothermal',  
    potential=pot_mw, mass=0.7, Rdisk=2.5, Hdisk=0.15,  
    sigmar0=50, Rsigmar=10)  
df_thick = agama.DistributionFunction(type='QuasiIsothermal',  
    potential=pot_mw, mass=0.3, Rdisk=2.5, Hdisk=0.40,  
    sigmar0=150, Rsigmar=10)  
df_total = agama.DistributionFunction(df_thin, df_thick)  
prob_thick = df_thick(actions) / df_total(actions)
```



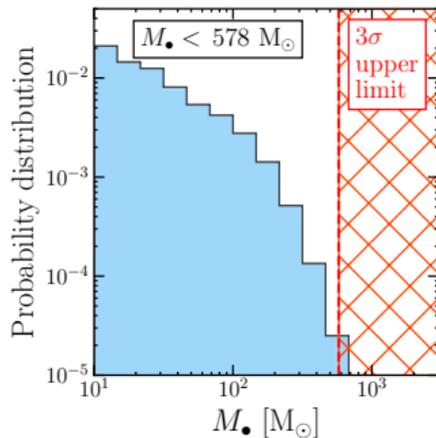
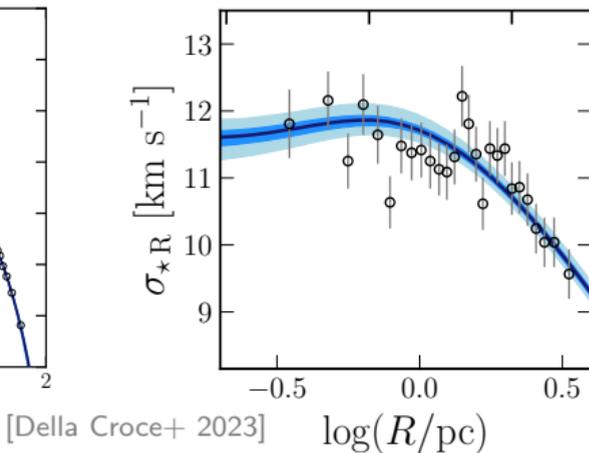
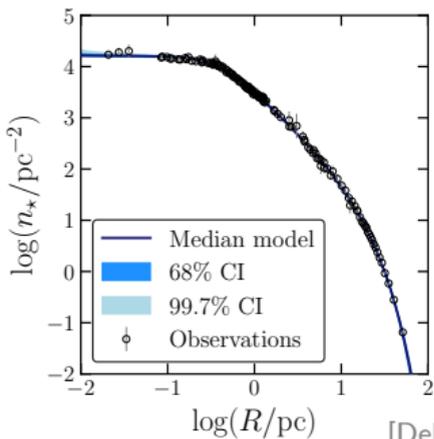
Distribution function applications: potential inference

DF is a *probability distribution* for finding a star with a given position and velocity, and it also depends on the potential Φ via the integrals of motion \mathcal{I} .

By maximising the likelihood of the observed dataset, one can determine the best-fit parameters α , β of the stellar system, including its mass distribution.

$$\ln \mathcal{L} = \sum_{i=1}^{N_{\text{stars}}} \ln f\left(\mathcal{I}(\mathbf{x}, \mathbf{v}; \Phi(\mathbf{x}; \alpha)); \beta\right)$$

Example: dynamical modelling of the globular cluster NGC 104 with the goal of constraining the mass of the central IMBH.



[Della Croce+ 2023]

Distribution function moments

The 6d DF $f(\mathbf{x}, \mathbf{v})$ can be reduced to more “easy to grasp” quantities:

▶ density $\rho(\mathbf{x}) = \int f(\mathbf{x}, \mathbf{v}) d^3 v,$

▶ mean velocity $\bar{\mathbf{v}}(\mathbf{x}) = \frac{1}{\rho(\mathbf{x})} \int \mathbf{v} f(\mathbf{x}, \mathbf{v}) d^3 v,$

▶ second moment of velocity $\overline{v_{ij}^2}(\mathbf{x}) = \frac{1}{\rho(\mathbf{x})} \int v_i v_j f(\mathbf{x}, \mathbf{v}) d^3 v,$

▶ more generally, velocity distribution at a given point

$$f(v_1; \mathbf{x}) = \frac{1}{\rho(\mathbf{x})} \int f(\mathbf{x}, \mathbf{v}) dv_2 dv_3 \quad (\text{it can be strongly non-Gaussian!}).$$

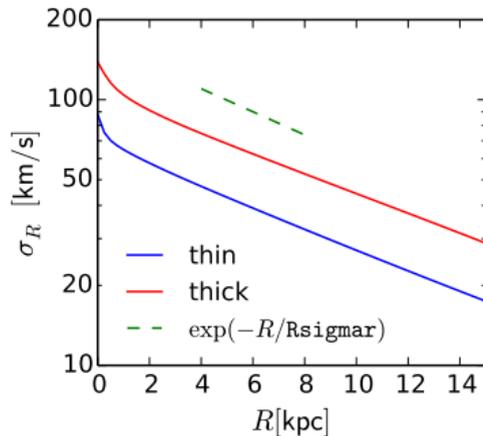
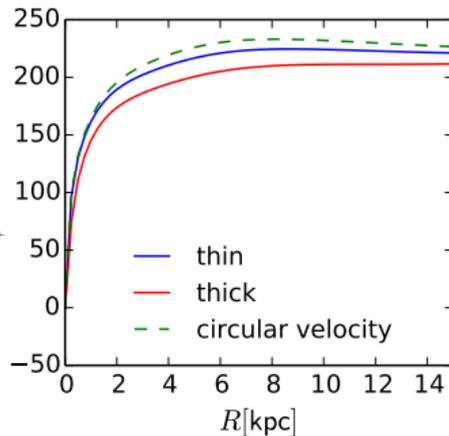
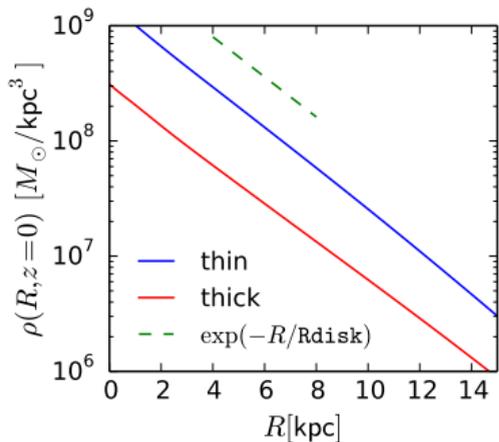
In most cases, we need $f(\mathbf{x}, \mathbf{v})$, but the DF gives us $f(\mathbf{J}) \implies$
use an action finder to convert from phase-space to action-space.

The combination of a potential $\Phi(\mathbf{x})$, action finder $\mathbf{J}(\mathbf{x}, \mathbf{v} \mid \Phi)$ and DF $f(\mathbf{J})$ is called GalaxyModel.

Distribution function moments

```
gm = agama.GalaxyModel(pot_mw, df_total)
radii = numpy.linspace(0, 15, 61)
rho, vel, vel2 = gm.moments(
    numpy.column_stack([radii, radii*0, radii*0]),
    dens=True, vel=True, vel2=True, separate=True)
ax[0].plot(radii, rho[:,0], color='b', label='rho, thin')
ax[0].plot(radii, rho[:,1], color='r', label='rho, thick')
ax[1].plot(radii, vel[:,0,1], color='b', label='mean vphi, thin')
ax[2].plot(radii, vel2[:,0,0]**0.5, color='b', label='sigma_r, thin')
```

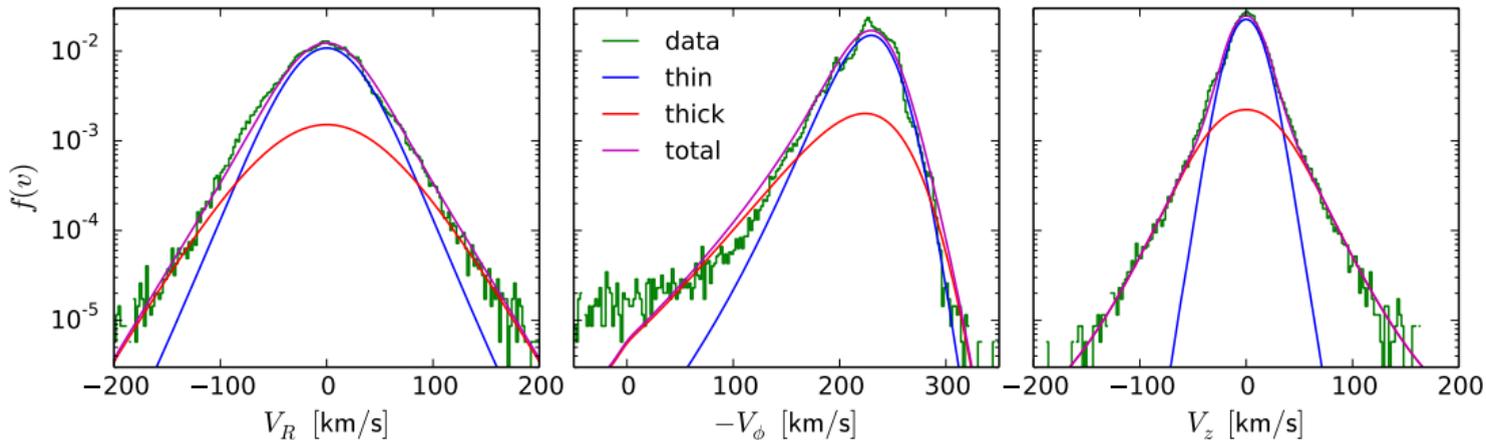
index of DF component ↗ ↖ index of velocity dimension: x, y, ...



Distribution functions in velocity (1d projections)

```
gridv = numpy.linspace(-200, 350)
vdf_vx, vdf_vy, vdf_vz, norm = gm.vdf([8.2, 0, 0],
    gridv=gridv, separate=True, dens=True)
frac0 = norm[0] / sum(norm)
frac1 = 1 - frac0
ax[0].plot(gridv, vdf_vx[0](gridv) * frac0, label='f(v_R), thin')
ax[0].plot(gridv, vdf_vx[1](gridv) * frac1, label='f(v_R), thick')
ax[0].plot(gridv, vdf_vy[0](gridv) * frac0, label='f(v_phi), thin')
ax[0].plot(gridv, vdf_vz[0](gridv) * frac0, label='f(v_z), thin')
```

index of DF component ↗

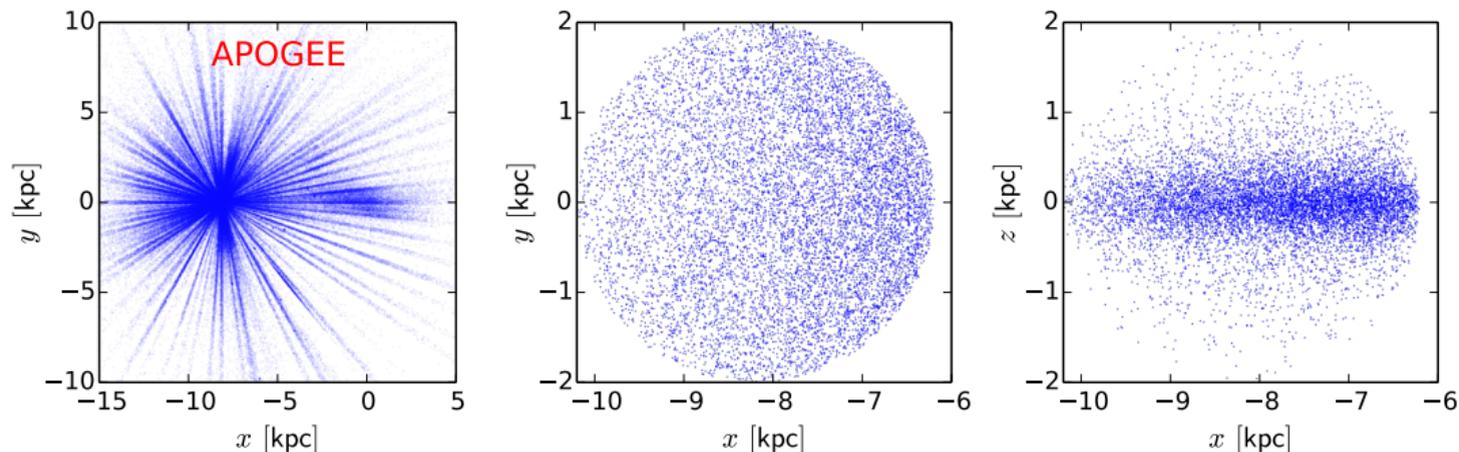


Sampling from a distribution function

One can use a selection function (usually spatial) to restrict the GalaxyModel to a limited volume, essentially using $DF \times SF$: $f(\mathbf{J}(\mathbf{x}, \mathbf{v})) \times S(\mathbf{x})$.

```
sf = agama.SelectionFunction([8.2, 0, 0], radius=2.0)
gm1 = agama.GalaxyModel(pot_mw, df_total, sf=sf)
posvel, mass = gm1.sample(1000000)
ax[0].scatter(posvel[:,0], posvel[:,1]) # x,y
ax[0].scatter(posvel[:,0], posvel[:,2]) # x,z
```

Or sample from the entire model (e.g., to create an N -body representation of it)...



Construction of self-consistent equilibrium models

Distribution function of stars $f(\mathbf{x}, \mathbf{v}, t)$

satisfies [sometimes] the collisionless Boltzmann equation:

$$\frac{\partial f(\mathbf{x}, \mathbf{v}, t)}{\partial t} + \mathbf{v} \cdot \frac{\partial f(\mathbf{x}, \mathbf{v}, t)}{\partial \mathbf{x}} - \frac{\partial \Phi(\mathbf{x}, t)}{\partial \mathbf{x}} \cdot \frac{\partial f(\mathbf{x}, \mathbf{v}, t)}{\partial \mathbf{v}} = 0.$$

Potential \Leftrightarrow mass distribution

not measured directly on human timescales

In order to infer anything about the potential from a time-dependent DF, need to make further assumptions about the initial state of the system, e.g., that the stars belong to a single stream or were perturbed from an equilibrium configuration in a specific way, etc.

Construction of self-consistent equilibrium models

Distribution function of stars $f(\mathbf{x}, \mathbf{v}, t)$

satisfies [sometimes] the collisionless Boltzmann equation:

$$\mathbf{v} \frac{\partial f(\mathbf{x}, \mathbf{v})}{\partial \mathbf{x}} - \frac{\partial \Phi(\mathbf{x})}{\partial \mathbf{x}} \frac{\partial f(\mathbf{x}, \mathbf{v})}{\partial \mathbf{v}} = 0.$$

Steady-state assumption \implies Jeans theorem:

$$f(\mathbf{x}, \mathbf{v}) = f(\mathcal{I}(\mathbf{x}, \mathbf{v}; \Phi))$$

3D - 6D
(observed)

integrals of motion ($\leq 3D?$), e.g., $\mathcal{I} = \{E, L, \dots\}$

3D
(want to infer)

Construction of **self-consistent** equilibrium models

Definition: a stellar system described by a time-independent DF $f(\mathcal{I}(\mathbf{x}, \mathbf{v}; \Phi))$ and potential $\Phi(\mathbf{x})$, which are related by the Poisson equation:

$$\nabla^2 \Phi(\mathbf{x}) = 4\pi G \rho(\mathbf{x}), \quad \text{where } \rho(\mathbf{x}) = \iiint d^3\mathbf{v} f(\mathcal{I}(\mathbf{x}, \mathbf{v})).$$

Applications:

- ▶ inference on gravitational potential from stellar kinematics (so-called *dynamical modelling*)
- ▶ creation of initial conditions for isolated galaxy simulations

Methods: (non-exhaustive list)

- ▶ DF₁: $\Phi + \rho \implies f$ (Eddington–Ossipkov–Merritt–Cuddeford inversion) only in spherical systems; **QuasiSpherical** DF with two free params β_0, r_a
- ▶ DF₂: $f \implies \Phi + \rho$ (iterative method)
- ▶ orbit-superposition: $\Phi + \rho + f_{i=1..N} \implies w_i$ (Schwarzschild method)

Iterative construction of DF-based self-consistent models

1. assume $f(\mathcal{I})$ and an initial guess for Φ

2. repeat
establish $\mathcal{I}(\mathbf{x}, \mathbf{v}; \Phi)$

$$\text{compute } \rho(\mathbf{x}) = \iiint d^3\mathbf{v} f(\mathcal{I}(\mathbf{x}, \mathbf{v}))$$

converged?

no

yes

update $\Phi(\mathbf{x})$ from the Poisson equation

3. enjoy!

```
scm_params = dict(
    rminSph=0.01, rmaxSph=100., sizeRadialSph=25, lmaxAngularSph=4)
df = agama.DistributionFunction(**df_params)
comp = agama.Component(df=df, disklike=False, **scm_params)
scm = agama.SelfConsistentModel(components=[comp], **scm_params)
scm.potential = init_potential
for i in range(5):
    scm.iterate()
comp.density.export('final_density.ini')
scm.potential.export('final_potential.ini')
```

Orbit-superposition method for self-consistent models

Introduced by Schwarzschild (1979) as a practical approach for constructing self-consistent triaxial models with prescribed $\rho(\mathbf{x}) \Leftrightarrow \Phi(\mathbf{x})$.

To invert the equation $\rho(\mathbf{x}) = \iiint f(\mathcal{I}(\mathbf{x}, \mathbf{v}; \Phi)) d^3\mathbf{v}$,
integrals of motion

discretize both the density profile and the distribution function:

$\rho(\mathbf{x}) \implies$ cells of a spatial grid;

mass of each cell is $M_c = \iiint_{\mathbf{x} \in V_c} \rho(\mathbf{x}) d^3\mathbf{x}$;

$f(\mathcal{I}) \implies$ collection of orbits with unknown weights:

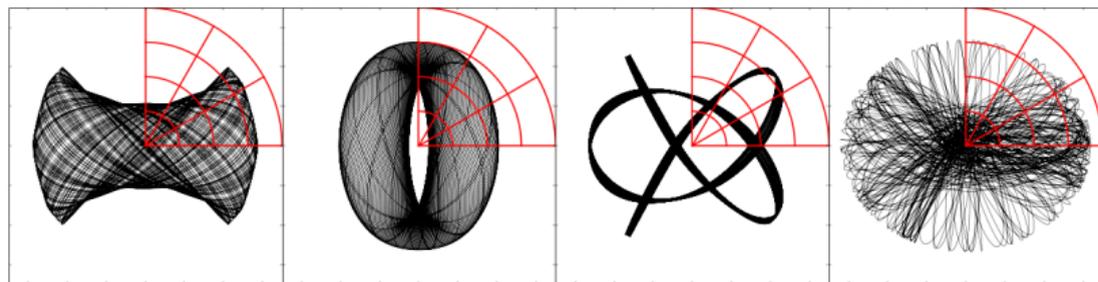
$$f(\mathcal{I}) = \sum_{k=1}^{N_{\text{orb}}} w_k \delta(\mathcal{I} - \mathcal{I}_k)$$

each orbit is a delta-function in the space of integrals of motion

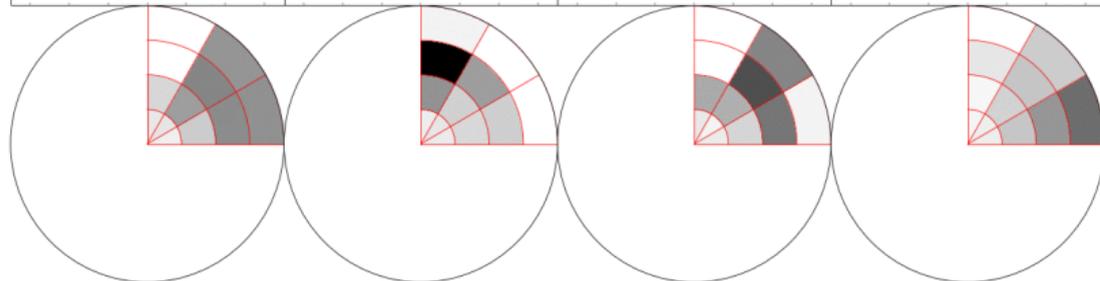
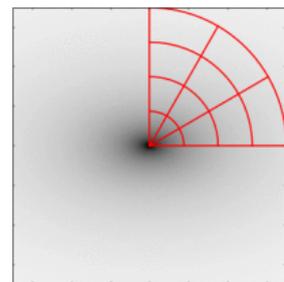
adjustable weight of each orbit [to be determined]

Orbit-superposition method for self-consistent models

orbits in the model



target density



discretized orbit density

(fraction of time t_{kc} that k -th orbit spends in c -th cell)

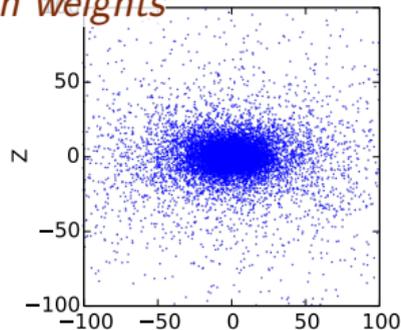
discretized density

(mass M_c in grid cells)

For each c -th cell we require $\sum_k w_k t_{kc} = M_c$, where $w_k \geq 0$ is orbit weight
system of linear equations with nonnegativity constraints \Rightarrow optimization problem

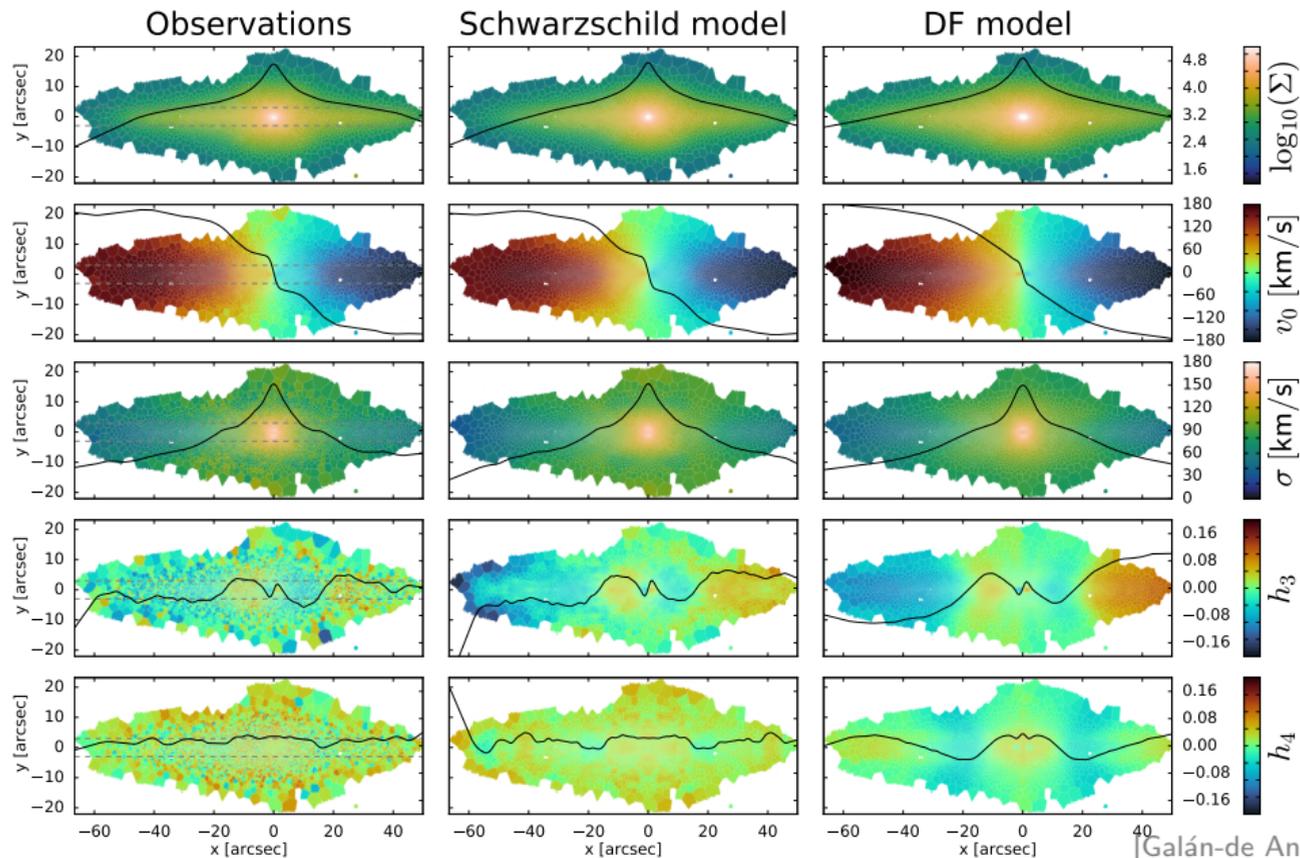
Orbit-superposition method: example

```
dens = agama.Density(type='Spheroid', axisRatioZ=0.5)
pot = agama.Potential(type='Multipole', density=dens)
target = agama.Target(type='DensityClassicLinear',
    gridr=numpy.logspace(-1.5, 2), stripsPerPane=2)
# prepare initial conditions for the orbit library
numOrbits = 10**4
ic,_ = dens.sample(numOrbits, potential=pot)
# compute the orbits and their contribution to the density
matrix, trajs = agama.orbit(potential=pot, ic=ic,
    time=100*pot.Tcirc(ic), dtype=object, targets=[target])
# solve the optimization problem to get the orbit weights
rhs = target(dens)
weights = agama.solveOpt(matrix=matrix.T, rhs=rhs)
# create N-body snapshot from recorded trajectories with weights
nbody = 10**5
_,(xv,m) = agama.sampleOrbitLibrary(
    nbody, trajs, weights)
plt.scatter(xv[:,0], xv[:,2])
# (omitting some non-essential but useful steps)
```



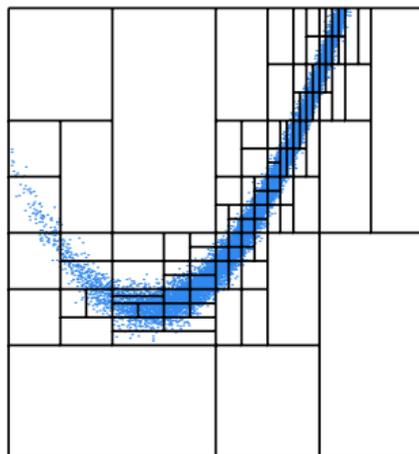
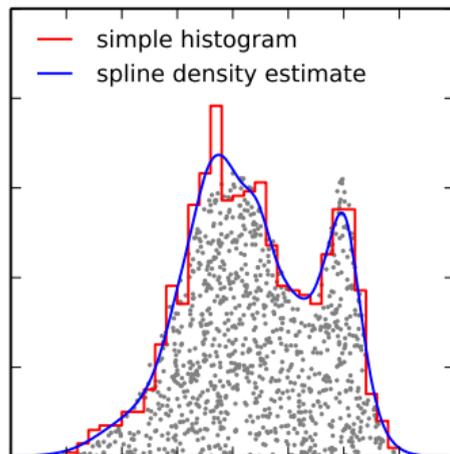
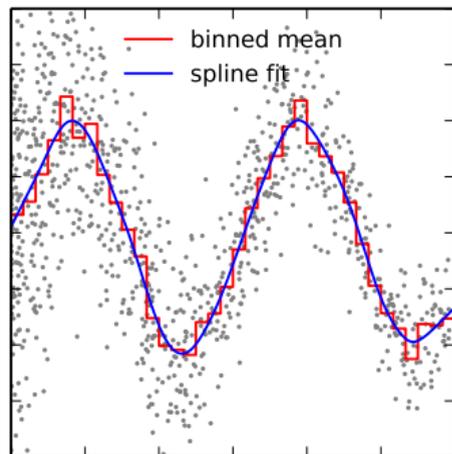
Example of DF- and orbit-based dynamical models

Model of an edge-on S0 galaxy FCC 170 constrained by MUSE IFU kinematics



Various mathematical methods

- ▶ spline interpolation (B-spline/cubic/quintic, 1d/2d/3d);
- ▶ penalized (i.e., with automatic optimal smoothing) spline fitting and density estimation;
- ▶ N-dimensional integration (using cubature or cuba libraries);
- ▶ drawing uniform-weight samples from an arbitrary N-dim probability function (rejection sampling with adaptive domain refinement).



Other features, caveats and limitations

- ▶ `AGAMA` is not a general-purpose N -body simulation code, but it can be used in this role to some extent, utilizing orbit integration coupled with the Multipole potential expansion (aka BFE).
- ▶ It can create initial conditions for galaxy simulations, provide external potential for several simulation codes (`NEMO` / `GYRFALCON`, `AMUSE`, `GADGET4`, `AREPO`), and assist in the analysis (e.g., extract potential from snapshots, integrate orbits, etc.)
- ▶ User-defined Python functions can serve as density, potential, DF and selection functions along with built-in C++ models, but are [much] less efficient; recommended approach is to approximate density or potential with C++-native expansions. A possible future development is to compile them as Cython code and use natively from the C++ core.
- ▶ No universal support for differentiable programming (`Jax autodiff` does not propagate into C++), but some functions can provide analytic derivatives (orbit integration, DFs, and in the future action-angle transformations).

Software for galactic dynamics

	GALPY	GALA	AGAMA
	[Bovy 2015]	[Price-Whelan 2017]	[Vasiliev 2019]
density and potential profiles:			
collection of analytic models	+	+	+
solution of the Poisson equation for an arbitrary $\rho(\mathbf{r})$ or an N -body snapshot	+	+	+
numerical integration of orbits	+	+	+
conversion between position/velocity and action/angle variables	+	+	+
distribution functions and their moments	+	-	+
construction of equilibrium models	-	-	+
modelling of tidal streams	+	+	+
N -body simulations with BFE	-	-	+
integration with <code>ASTROPY</code>	+	+	-
language	Python, C	Cython	C++, Python

AGAMA potentials can be used in GALA & GALPY, although not most efficiently

Summary

AGAMA is a versatile toolbox for stellar dynamics catering to many needs:

- ▶ Extensive collection of gravitational potential models
(analytic profiles, azimuthal- and spherical-harmonic expansions)
constructed from smooth density profiles or N -body snapshots;
 - ▶ Numerical orbit integration;
 - ▶ Conversion to/from action/angle variables;
 - ▶ Self-consistent multicomponent models with action-based DFs;
 - ▶ Schwarzschild orbit-superposition models;
 - ▶ Generation of initial conditions for N -body simulations;
 - ▶ Various math tools: spline-based interpolation, fitting and density estimation, multidimensional sampling;
 - ▶ Efficient and carefully designed C++ implementation, examples, Python and Fortran interfaces, plugins for Galpy, Gala, NEMO, AMUSE.
- <https://github.com/GalacticDynamics-Oxford/Agama>