# Action-angle variables in galactic dynamics

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Summer School on Galactic Dynamics, Shanghai, June 2019

# Hamiltonian mechanics

Consider a particle moving in a potential  $\Phi(\mathbf{x})$ .

 $\mathbf{x}(t), \mathbf{v}(t)$  are "ordinary" *D*-dimensional position/velocity coordinates;  $H(\mathbf{x}, \mathbf{v}) = \Phi(\mathbf{x}) + \frac{1}{2} |\mathbf{v}|^2$  is the Hamiltonian.

The equations of motion are

$$\frac{d\mathbf{x}}{dt} \equiv \dot{\mathbf{x}} = \mathbf{v}, \quad \frac{d\mathbf{v}}{dt} \equiv \dot{\mathbf{v}} = -\frac{\partial\Phi}{\partial\mathbf{x}}$$

One may consider a general class of Hamiltonian systems defined by  $H(\mathbf{q}, \mathbf{p})$  as a function of generalized phase-space coordinates, which satisfy the Hamilton's equations of motion:

$$\dot{\mathbf{q}} = \frac{\partial H}{\partial \mathbf{p}}, \qquad \dot{\mathbf{p}} = -\frac{\partial H}{\partial \mathbf{q}}$$

#### **Poisson brackets**

Define the commutator operator for two functions of phase-space coordinates  $A(\mathbf{q}, \mathbf{p})$  and  $B(\mathbf{q}, \mathbf{p})$  as

$$[A, B] \equiv \frac{\partial A}{\partial \mathbf{q}} \frac{\partial B}{\partial \mathbf{p}} - \frac{\partial A}{\partial \mathbf{p}} \frac{\partial B}{\partial \mathbf{q}}$$

It follows immediately that

$$[A, A] = 0, \quad [A, B] = -[B, A], \qquad (\text{antisymmetry})$$

$$[[A, B], C] + [[B, C], A] + [[C, A], B] = 0, \qquad (\text{Jacobi identity})$$

$$[q_i, q_j] = 0, \quad [p_i, p_j] = 0, \quad [q_i, p_j] = \delta_{ij}, \qquad i, j = 1..D,$$

and the Hamilton equations can be written as

$$\dot{q}_i = [q_i, H], \quad \dot{p}_i = [p_i, H]$$

# **Integrals of motion**

If [A, B] = 0, we say that A commutes with B.

If a function  $A(\mathbf{q}, \mathbf{p})$  commutes with the Hamiltonian, it is conserved along the particle's trajectory – we call it an integral of motion:

$$\frac{dA}{dt} = \frac{\partial A}{\partial \mathbf{q}} \frac{d\mathbf{q}}{dt} + \frac{\partial A}{\partial \mathbf{p}} \frac{d\mathbf{p}}{dt}$$
$$= \frac{\partial A}{\partial \mathbf{q}} \frac{\partial H}{\partial \mathbf{p}} - \frac{\partial A}{\partial \mathbf{p}} \frac{\partial H}{\partial \mathbf{q}}$$
$$= [A, H] = 0$$

Obviously, the Hamiltonian itself is an integral of motion.

Phase-space distribution function  $f(\mathbf{q}, \mathbf{p})$  satisfies the collisionless Boltzmann equation and hence is also conserved along the trajectory of any particle.

# **Canonical transformations**

Consider a change of variables from  $\mathbf{p}, \mathbf{q}$  to  $\mathbf{P}, \mathbf{Q}$ , and express the Hamiltonian  $H(\mathbf{P}, \mathbf{Q})$  or any other function in phase space in terms of the new variables.

If the new variables satisfy the canonical commutatation relations  $[Q_i, Q_j] = 0$ ,  $[P_i, P_j] = 0$ ,  $[Q_i, P_j] = \delta_{ij}$ , such transformation is called canonical (or symplectic).

It also preserves

Hamilton's equations of motion:

 $\dot{\mathbf{Q}}_i = [Q_i, H], \ \dot{\mathbf{P}}_i = [P_i, H];$ 

• more generally, all Poisson brackets:  $[A(\mathbf{p}, \mathbf{q}), B(\mathbf{p}, \mathbf{q})] = [A(\mathbf{P}, \mathbf{Q}), B(\mathbf{P}, \mathbf{Q})];$ 

► all Poincaré invariants: ∮ p · dq

> 2*D*-dimensional phase volume element:  $d^D \mathbf{q} d^D \mathbf{p} = d^D \mathbf{Q} d^D \mathbf{P}$ 

## **Examples of canonical transformations**

1. Exchange:  $\mathbf{Q} = \mathbf{p}, \ \mathbf{P} = \mathbf{q}$ 

(i.e., there is no fundamental difference between coordinate and momentum variables).

2. Point transformation: define Q(q) in whatever way, and then P(q, p) is uniquely specified.

For instance, cartesian to polar coordinates:  $\mathbf{q} \equiv \{x, y\}$  to  $\mathbf{Q} \equiv \{r, \phi\}$ implies  $\mathbf{P} \equiv \{p_r, p_{\phi}\} = \{(xp_x + yp_y)/r, xp_y - yp_x\}.$ 

 Hamiltonian flow: integrate the equations of motion for some time τ, and let {Q, P}(q, p; τ) be the new coordinates and momenta of a point started from initial conditions q, p.

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One powerful way of constructing such transformations is to introduce

a generating function  $F(\mathbf{q}, \mathbf{P})$  such that  $\mathbf{p} = \frac{\partial F}{\partial \mathbf{q}}$ ,  $\mathbf{Q} = \frac{\partial F}{\partial \mathbf{P}}$ ;

F could also be expressed in terms of some other combination of old and new variables, e.g.,  $F(\mathbf{q}, \mathbf{Q})$ , etc.

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$$H(\mathbf{q},\mathbf{p})=rac{1}{2}|\mathbf{p}|^2 \implies p_i(t)=\mathrm{const}, \quad q_i(t)=p_i\,t+\mathrm{const}$$

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where **q** are treated  
as angle-like (periodic)  
variables,  $q + 2\pi \cong q$   
on a *D*-dimensional torus.  
These are action-angle variables  $q_1$ 

# Integrability and the Arnold–Liouville theorem

If  $I_1$  and  $I_2$  are two integrals of motion, then the Jacobi identity  $[[I_1, I_2], H] + [[I_2, H], I_1] + [[H, I_1], I_2] = 0$  implies that  $[I_1, I_2]$  is also an integral of motion.

(Example:  $I_1 = L_x, I_2 = L_y \Rightarrow [I_1, I_2] = L_z$ ).

If  $[I_1, I_2]$  is identically zero, the two integrals are said to be in involution.

A Hamiltonian system with D degrees of freedom is *integrable* if it has D independent integrals of motion  $I_1 ldots I_D$  (including the Hamiltonian itself) which are all in involution with each other.

The motion of any particle is restricted to a D-dimensional hypersurface of the 2D-dimensional phase space.

**Theorem:** this hypersurface is diffeomorphic to (i.e., could be smoothly deformed into) a *D*-torus, parametrized by *D* periodic variables  $\theta \in [0..2\pi)$ .

#### Action-angle variables for a 1d simple harmonic oscillator

The simplest possible Hamiltonian system:  $H(q, p) = \frac{1}{2}p^2 + \frac{1}{2}\omega^2 q^2$ .

The trajectory is  $q(t) = A \sin(\omega t + \phi_0)$ ,  $p(t) = A \omega \cos(\omega t + \phi_0)$ , and the energy is  $E = \frac{1}{2}\omega^2 A^2$ .

The motion is periodic with frequency  $\omega$  ( $\Leftrightarrow$  period  $2\pi/\omega$ ), so we define the angle  $\theta = \omega t + \phi_0$ .

The action J is  $\frac{1}{2\pi}$  × area enclosed by the trajectory:





#### Action-angle variables for a 2d simple harmonic oscillator

The same thing but in two dimensions:  $\mathbf{q} = \{x, y\}, \ \mathbf{p} = \{p_x, p_y\};$ Hamiltonian:

$$H(\mathbf{q}, \mathbf{p}) = \frac{1}{2} (p_x^2 + \omega_x^2 x^2) + \frac{1}{2} (p_y^2 + \omega_y^2 y^2)$$
  
$$\equiv H_x(x, p_x) + H_y(y, p_y)$$

Motion is separable in x, y - ttwo uncoupled simple harmonic oscillators, two integrals of motion  $E_x, E_y$ , actions are  $J_x = E_x/\omega_x, J_y = E_y/\omega_y$ .



## Action-angle variables for a 2d planar axisymmetric potential

A slightly more complicated system: two degrees of freedom, motion in an axisymmetric potential  $\Phi(x, y) = \Phi(R)$ , where  $R \equiv \sqrt{x^2 + y^2}$ .

Canonical coordinates:  $\mathbf{q} = \{R, \phi\}, \mathbf{p} = \{p_R, p_\phi\}$ 

Hamiltonian: 
$$H = \Phi(R) + \frac{1}{2}\left(p_R^2 + \frac{p_{\phi}^2}{R^2}\right) \equiv \Phi_{\text{eff}}(R) + \frac{1}{2}p_R^2$$

equations of motion:  $\dot{R} = p_R$ ,  $\dot{\phi} = \frac{p_{\phi}}{R^2}$ ,  $\dot{p}_R = -\frac{d\Phi_{\text{eff}}}{dR}$ ,  $\dot{p}_{\phi} = 0$ 

integrals of motion: E and  $p_{\phi}$ 

Motion in R is described by a 1d effective potential  $\Phi_{\rm eff}(R)\equiv \Phi(R)+p_{\phi}^2/R^2$ 

The radial action is

$$egin{split} J_R &= rac{1}{\pi} \int_{R_-}^{R_+} p_R(R;\,E,p_\phi)\,dR \ &= rac{1}{\pi} \int_{R_-}^{R_+} \sqrt{2ig[(E-\Phi_{
m eff}(R)ig]}\,\,dR \end{split}$$



#### Action-angle variables for a 2d planar axisymmetric potential

Motion in  $\phi$ :  $\dot{p}_{\phi} = 0 \Rightarrow p_{\phi} = \text{const}$ ,

hence the azimuthal action is

$$J_{\phi}=rac{1}{2\pi}\int_{0}^{2\pi} p_{\phi}\,d\phi=p_{\phi}.$$

The actions  $J_R$ ,  $J_\phi$  describe the extent of the orbit in two complementary dimensions:

 $J_{\phi}$  corresponds to the "guiding radius" (the radius of a circular orbit with the given angular momentum  $J_{\phi}$ ),

 $J_R$  gives the extent of radial oscillation about this guiding radius.

They can be varied independently, and any possible choice (provided that  $J_R \ge 0$ ) corresponds to some trajectory.



# **Angles and frequencies**

Note that  $\dot{\phi} = p_{\phi}/R^2(t) \neq \text{const}$ , so  $\phi$  is not a canonically conjugate angle variable to  $p_{\phi}$ !

Such variable is  $\theta_{\phi}$  defined to increase linearly with time, and similarly the radial phase angle  $\theta_R$  also increases linearly with time:

 $\theta_R = \Omega_R t, \quad \theta_\phi = \Omega_\phi t, \text{ where }$  $\Omega_R \equiv \frac{\partial H(J_R, J_{\phi})}{\partial J_{\Sigma}}, \quad \Omega_{\phi} \equiv \frac{\partial H(J_R, J_{\phi})}{\partial L} \quad \text{are orbital frequencies.}$  $\theta_R(R; E, p_\phi) = \Omega_R \int_{R}^{R} \frac{dt}{dR} dR = \Omega_R \int_{R}^{R} \frac{dR}{p_R(R; E, p_\phi)}$ Radial orbital period  $T_R \equiv \frac{2\pi}{\Omega_R} = 2 \int_{R_-}^{R_+} \frac{dR}{p_R} = 2 \int_{R_-}^{R_+} \frac{dR}{\sqrt{2[E - \Phi(R)] - \frac{p_{\phi}^2}{R^2}}}$ Azimuthal period  $T_{\phi} \equiv \frac{2\pi}{\Omega_{\phi}} = \frac{2\pi}{p_{\phi}} \int_{\Gamma_{e}}^{R_{+}} dR/p_{R}}{p_{\phi}}$ 

#### Action-angle variables for a 3d spherical potential

Spherical coordinates:  $r, \theta, \phi, p_r, p_{\theta}, p_{\phi}$ 

Hamiltonian: 
$$H = \Phi(r) + \frac{1}{2} \left( p_r^2 + \frac{p_\theta^2}{r^2} + \frac{p_\phi^2}{r^2 \sin^2 \theta} \right)$$

Integrals of motion:  $E, L_x, L_y, L_z \left[, L \equiv \sqrt{L_x^2 + L_y^2 + L_z^2}\right]$ 

Radial action: 
$$J_r = \frac{1}{\pi} \int_{r_-}^{r_+} \sqrt{2[E - \Phi(r)] - \frac{L^2}{r^2}} \ge 0$$

Azimuthal action:  $J_{\phi} = L_z$  (any sign)

Vertical action:  $J_{\theta} \equiv J_z = L - |L_z| \ge 0$ 

In general, actions, angles, frequencies, or  $H(\mathbf{J})$  do not have analytic expressions. One exception is the isochrone potential [Hénon 1959]:

$$\Phi(r) = -\frac{G M}{b + \sqrt{b^2 + r^2}} \quad \text{(includes Kepler and harmonic osclilator as limiting cases)}$$
$$H(\mathbf{J}) = -\frac{2 (G M)^2}{\left(2J_r + L + \sqrt{L^2 + 4 G M b}\right)^2}$$

## Action-angle variables for a 3d axisymmetric potential

For nearly-circular orbits close to the equatorial plane, one may use the **epicyclic approximation**:

 $\Phi(R,z) \approx \Phi_R(R) + \Phi_z(z)$ ,

motion in  $R, \phi$  as in the planar axisymmetric problem with effective potential  $\Phi_{\text{eff}} = \Phi_R(R) + \frac{1}{2}L^2/r^2$ , and independent, nearly harmonic motion in z.

However, it becomes increasingly inaccurate for orbits with high eccentricity and/or inclination.



## State of the art: Stäckel fudge

Fact: orbits in realistic axisymmetric galactic potentials are much better aligned with prolate spheroidal coordinates.



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One may explore the assumption that the motion is separable in these coordinates  $(\lambda, \nu)$ .



#### Stäckel fudge [Binney 2012]

The most general form of potential that satisfies the separability condition is the Stäckel potential<sup>1</sup>:  $\Phi(\lambda, \nu) = -\frac{f_1(\lambda) - f_2(\nu)}{\lambda - \nu}$ .

The motion in  $\lambda$  and  $\nu$  directions, with canonical momenta  $p_{\lambda}, p_{\nu}$ , is governed by two separate equations:

$$2(\lambda - \Delta^2) \lambda p_{\lambda}^2 = \left[ E - \frac{L_z^2}{2(\lambda - \Delta^2)} \right] \lambda - [I_3 + (\lambda - \nu)\Phi(\lambda, \nu)],$$
  
$$2(\nu - \Delta^2) \nu p_{\nu}^2 = \left[ E - \frac{L_z^2}{2(\nu - \Delta^2)} \right] \nu - [I_3 + (\nu - \lambda)\Phi(\lambda, \nu)].$$

Under the approximation that the separation constant  $I_3$  is indeed conserved along the orbit, actions are computed as

$$J_{\lambda} = rac{1}{\pi} \int_{\lambda_{\min}}^{\lambda_{\max}} p_{\lambda} \, d\lambda, \quad J_{
u} = rac{1}{\pi} \int_{
u_{\min}}^{
u_{\max}} p_{
u} \, d
u.$$

<sup>1</sup>Note that the potential of the Perfect Ellipsoid [de Zeeuw 1985] is of the Stäckel form, but it is only one example of a much wider class of potentials.

# Stäckel fudge in practice

A rather flexible approximation: for each orbit, we have the freedom of using two functions  $f_1(\lambda)$ ,  $f_2(\nu)$  (directly evaluated from the actual potential  $\Phi(R, z)$ ) to describe the motion in two independent directions. These functions are different for each orbit (implicitly depend on  $E, L_z, I_3$ ). Moreover, we may choose the focal distance  $\Delta$  of the auxiliary prolate spheroidal coordinate system for each orbit independently.



#### Accuracy of the Stäckel fudge

Accuracy of action conservation using the Stäckel fudge:  $\leq 1\%$  for most disk orbits,  $\leq 10\%$  even for high-eccentricity orbits [except near resonances!]. Interpolation of  $J_r$ ,  $J_z$  on a 3d grid of E,  $L_z$ ,  $I_3$ : 10x speed-up at the expense of a moderate [not always acceptable!] decrease in accuracy.



# Other methods for action computation

The accuracy of the Stäckel approximation is "uncontrollable" (cannot be systematically improved), and it is mainly used in axisymmetric potentials.

However, actions offer the only **systematic** method for computing the integrals of motion in a **non-perturbative** way for an arbitrary potential:

Introduce a simple enough "toy" potential Φ<sup>t</sup> (e.g., isochrone), for which the mapping between position–velocity {x, v} and action–angle {J<sup>t</sup>, θ<sup>t</sup>} coordinates is known analytically.

We seek a canonical transformation between the true (yet unknown) {J, θ} and the "toy" {J<sup>t</sup>, θ<sup>t</sup>}. This transformation is described by a generating function S(J, θ<sup>t</sup>), which can be expanded into Fourier series in θ<sup>t</sup>: S(J, θ<sup>t</sup>) = J ⋅ θ<sup>t</sup> + ∑<sub>n</sub> S<sub>n</sub>(J) exp(in ⋅ θ<sup>t</sup>), where n are triplets of integers.

- Choose the Fourier coefficients S<sub>n</sub> up to some maximum order n to approximate the true Hamiltonian to any desired accuracy.
- The transformation is given by  $\mathbf{J}^t = \partial S / \partial \boldsymbol{\theta}^t$ ,  $\boldsymbol{\theta} = \partial S / \partial \mathbf{J}$ .

# Other methods for action computation



There are several variants of these methods, but we won't go into details.

- First numerically integrate the orbit, then obtain the coefficients S<sub>n</sub> to minimize the variation of the toy Hamiltonian across the real orbit [Sanders & Binney 2014; Bovy 2014]. This gives the transformation from {x, v} to {J, θ}.
- Reverse transformation (torus mapping) allows one to compute the position/velocity from action/angle without the need to integrate an orbit [McGill & Binney 1990; McMillan & Binney 2008].

 A variation of this approach also works for resonantly-trapped orbits [Kaasalainen 1994; Binney 2016, 2018].

## Advantages of action/angle variables

- Clear physical meaning (describe the extent of oscillations in each dimension).
- Most natural description of motion (angles change linearly with time).
- Possible range for each action variable is [0..∞) or (-∞..∞), independently of the other ones (unlike *E* and *L*, say).
- Canonical coordinates  $\Rightarrow$  the 6d phase-space volume element is  $d^3x \ d^3v = d^3J \ d^3\theta$ .
- Actions are adiabatic invariants (are conserved under slow variation of potential).
- Perturbation theory most naturally formulated in terms of actions.
- Efficient methods for conversion between  $\{x, v\}$  and  $\{J, \theta\}$  now exist.

## Fun facts / rules of thumb about actions

- Dimension of actions is length×velocity: if a star at a galactocentric distance r travels with velocity v, then [at least one of the actions] J ~ r v.
- Frequencies: Ω<sub>i</sub>(J) = ∂H/∂J<sub>i</sub> characteristic velocity v<sub>i</sub> ~ √Ω<sub>i</sub> J<sub>i</sub> e.g., for a circular orbit J<sub>φ</sub> = R v<sub>φ</sub>, Ω<sub>φ</sub> = v<sub>φ</sub>/R.
- Surfaces of constant energy  $H(\mathbf{J}) = E$  are approximately tetrahedra in the 3d action space, with  $E \approx E(\Omega_r J_r + \Omega_z J_z + \Omega_\phi J_\phi)$ .

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## Natural coordinates to describe orbits

- The entire 6d phase space is foliated into non-intersecting 3d orbital tori.
- Actions tell you which orbit the star is on, angles – where it is located on this orbit.
- Angles change linearly with time,  $\theta_i = \Omega_i t$
- ► Torus construction provides the transformation
  J, θ → x, v, i.e., one can find the position-velocity at any time without the need to integrate the orbit.
- In a time-averaged sense, only actions are significant (distribution of stars is averaged over angles – phase mixed);

however, for an initially localized ensemble of stars (e.g., a stream from a disrupted cluster), the distribution over angles is not uniform.



- **1.** Numerically integrate the trajectory:  $x(t), y(t), p_x(t), p_y(t)$ .
- 2. Every time it passes through the axis y = 0 with  $\dot{y} > 0$ , put a point on the  $x, p_x$  plane.



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- 3. Each orbit corresponds to a closed loop in this plane.
- **4.** Repeat for many different initial conditions to get the "phase portrait" of the Hamiltonian.



- 3. Each orbit corresponds to a closed loop in this plane.
- **5.** The action of an orbit is just the area inside its Poincaré curve:  $J_x = \frac{1}{2\pi} \oint p_x \, dx$



A convenient tool for analyzing orbits in 2d Hamiltonian systems at a fixed E (e.g., motion in the equatorial plane, or in the meridional plane of an axisymmetric potential at a fixed  $L_z$ )

6. Now repeat this exercise for a different choice of energy E



- 7. This portrait may contain more than one orbit family!
- 8. The meaning of actions is different for each orbit family (e.g., it is  $J_x$  for a box orbit, and  $J_r$  for a loop orbit)



- 7. This portrait may contain more than one orbit family!
- 9. For high-order resonances the action *may* describe the width around parent orbit, and for chaotic orbits the actions are not defined at all..



#### Action-angles are canonical coordinates

A distribution function (DF) in phase space  $f(\mathbf{x}, \mathbf{v})$  can be expressed in any other coordinates, e.g., f(E, L) or  $f(\mathbf{J}, \theta)$ .

In a phase-mixed system, it is independent of  $\theta$ .

The mass in a given volume of phase space is  $\int f(\mathbf{x}, \mathbf{v}) d^3x d^3v = \int f(\mathbf{J}, \theta) d^3J d^3\theta = (2\pi)^3 \int f(\mathbf{J}) d^3J - does not depend on the potential and has no extra multiplicative factors! (e.g., compared to <math>\int f(E, L) 4\pi^2 T(E, L; \Phi) dL^2$  for a conventional DF).

The functional form of the DF is independent of the potential  $\Phi$ , but the correspondence between **x**, **v** and **J**,  $\theta$ , of course, does depend on  $\Phi$ .

# Adiabatic invariance of actions

Actions are conserved under slow variations of potential ( $\Omega \tau \gg 1$ ).

Examples:

- accreted satellite galaxies should stay localized in the action space;
- compression of the dark halo after the formation of the stellar disk:



#### Perturbation theory in action space

$$\begin{split} f(\mathbf{J}, \boldsymbol{\theta}, t) &= f_0(\mathbf{J}) + \epsilon f_1(\mathbf{J}, \boldsymbol{\theta}, t), \\ H(\mathbf{J}, \boldsymbol{\theta}, t) &= H_0(\mathbf{J}) + \epsilon H_1(\mathbf{J}, \boldsymbol{\theta}, t) = H(\mathbf{x}, \mathbf{v}, t) \equiv \Phi_0(\mathbf{x}) + \epsilon \Phi_1(\mathbf{x}, t) + \frac{1}{2}v^2. \\ \text{Linearized Vlasov / collisionless Boltzmann equation:} \end{split}$$

$$0 = \frac{\partial f}{\partial t} + \left[H, f\right] \approx \frac{\partial f_1}{\partial t} + \frac{\partial f_1}{\partial \theta} \frac{\partial H_0}{\partial J} - \frac{\partial f_0}{\partial J} \frac{\partial \Phi_1}{\partial \theta}.$$

 $\Phi_1(\mathbf{x}, t)$  is the external perturbation augmented with internal self-gravity (diverges at resonances!).

For the given  $f_0$  and  $\Phi_1$ , one may compute the perturbed DF  $f_1(\mathbf{J}, \boldsymbol{\theta}, t)$  [e.g., Monari+ 2016–2018] is possible.



### **Distribution functions**

The pair  $f(\mathbf{J})$ ,  $\Phi(\mathbf{x})$  provides the complete description of the system:

Φ determines the transformation {x, v} ↔ {J, θ};
density is ρ(x) = ∫ f [J(x, v; Φ)] d<sup>3</sup>v;

► velocity moments are 
$$\overline{\mathbf{v}} = \frac{1}{\rho} \int f[\mathbf{J}(\mathbf{x}, \mathbf{v}; \Phi)] \mathbf{v} d^3 v;$$
  
 $\overline{v_i v_j} = \frac{1}{\rho} \int f[\mathbf{J}(\mathbf{x}, \mathbf{v}; \Phi)] v_i v_j d^3 v$ , etc.

Now two questions remain:

- **1.** How to choose a sensible  $f(\mathbf{J})$
- **2.** How to find  $\Phi(\mathbf{x})$  consistent with this DF

#### Distribution functions for spheroidal systems

Recall that surfaces of constant energy are approximately tetrahedra in action space,  $E \approx E(\Omega_r J_r + \Omega_z J_z + \Omega_\phi J_\phi).$ 

So if we consider  $f(\mathbf{J}) = f_0[h(\mathbf{J})]$ , where  $h(\mathbf{J}) = k_r J_r + k_z J_z + k_{\phi} |J_{\phi}|$  is a linear combination of three actions with the above coefficients, it will be approximately isotropic (dependent on energy only).

We may construct tailored anisotropic systems by changing the coefficients  $k_i$ . [Binney 2014, Posti+ 2015, Williams & Evans 2015].

The function  $f_0$  is responsible for the overall density profile of the system; a reasonable choice is a double-power-law model:

$$f_0(h) \propto rac{h^{\Gamma}}{ig(1+[h/h_0]^\etaig)^{(\Gamma-{
m B})/\eta}}$$

Γ: inner slope B: outer slope  $h_0$ : scale action

pola

prograde

#### Distribution functions for disky systems

In the epicyclic approximation, the motion is separable in R, z, and the DF in each dimension has a nearly Boltzmann form:  $f(E, L_z, E_z) \propto f_0(L_z) \exp(-E_R/\sigma_R^2) \exp(-E_z/\sigma_z^2).$ 

One may construct a similarly behaving DF expressed in terms of actions, replacing  $E_R \rightarrow \Omega_R J_R$ ,  $E_z \rightarrow \Omega_z J_z$  – a quasi-isothermal DF [Binney & McMillan 2011]:

$$\begin{split} f(\mathbf{J}) &= \frac{\tilde{\Sigma}\,\Omega}{2\pi^2\,\kappa^2} \times \frac{\kappa}{\tilde{\sigma}_r^2} \exp\left(-\frac{\kappa\,J_r}{\tilde{\sigma}_r^2}\right) \times \frac{\nu}{\tilde{\sigma}_z^2} \exp\left(-\frac{\nu\,J_z}{\tilde{\sigma}_z^2}\right) \times \begin{cases} 1 & \text{if } J_\phi \geq 0, \\ \exp\left(\frac{2\Omega\,J_\phi}{\tilde{\sigma}_r^2}\right) & \text{if } J_\phi < 0, \end{cases} \\ \tilde{\Sigma}(R_c) &\equiv \Sigma_0 \exp\left(-\frac{R_c}{R_{\text{disk}}}\right), \quad \tilde{\sigma}_r^2(R_c) \equiv \sigma_{r,0}^2 \exp\left(-\frac{2R_c}{R_{\sigma,r}}\right), \quad \tilde{\sigma}_z^2(R_c) \equiv 2\,h_{\text{disk}}^2\,\nu^2(R_c). \end{split}$$

It produces nearly exponential radial profiles of surface density and nearly isothermal vertical density profiles:  $\rho(R, z) \propto \exp\left(-\frac{R}{R_{\text{disk}}}\right) \operatorname{sech}^2\left(\frac{z}{h}\right)$ .

# **Self-consistent models**



(Assumption: a galaxy is a collisionless system in a steady state)

# **Self-consistent models**

# distribution function integrals of motion gravitational potential

1. Collisionless Boltzmann equation:

$$\mathbf{v} \frac{\partial f}{\partial \mathbf{x}} - \frac{\partial \Phi}{\partial \mathbf{x}} \frac{\partial f}{\partial \mathbf{v}} = 0 \implies f = f\left(\mathcal{I}(\mathbf{x}, \mathbf{v}; \Phi)\right).$$

(Assumption: a galaxy is a collisionless system in a steady state)

2. Poisson equation:

$$\nabla^2 \Phi(\mathbf{x}) = 4\pi \ G \ \rho(\mathbf{x}).$$
 total density

(Assumption: Newtonian gravity)

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**3.** The link:

$$\rho(\mathbf{x}) = \iiint d^3 v \, f(\mathbf{x}, \mathbf{v}).$$

(Assumption: self-consistency)



#### Self-consistent models – iterative approach

- **1.** Assume a particular distribution function  $f(\mathcal{I})$ ;
- **2.** Adopt an initial guess for  $\Phi(\mathbf{x})$ ;
- 3. Establish the integrals of motion I(x, v) in this potential;
  4. Compute the density ρ(x) = ∫∫∫ d<sup>3</sup>v f(I(x, v));
  5. Solve the Poisson equation to find the new potential Φ(x);
  6. Repeat until convergence.

**Origin**: Prendergast & Tomer 1970;

used in Kuijken & Dubinski 1995, Widrow+ 2008, Taranu+ 2017 (GalactICs),

Piffl+ 2014, Cole & Binney 2016, Sanders & Evans 2016 (action-based formalism).

# Advantages of models based on distribution function

#### Clear physical meaning

(localized structures in the space of integrals of motion);

Easy to compare different models

(how to compare two *N*-body or *N*-orbit models?);

- Easy to compare models to discrete observational data;
- Easy to sample particles from the distribution function (convert to an N-body model);
- Stability analysis

(perturbation theory most naturally formulated in terms of actions);

#### **Caveats:**

- Implicitly rely on the integrability of the potential, ignore the presence of resonant orbit families (but see Binney 2016, 2018);
- So far implemented only for axisymmetric models (not a fundamental limitation).

# Galactic modelling tasks

- $\checkmark \checkmark \checkmark \checkmark \checkmark \checkmark \checkmark$  For Gravitational potentials and forces
- $\checkmark \checkmark \checkmark \checkmark \checkmark \checkmark \checkmark$  **>** Orbit integration and analysis
- $\checkmark \checkmark \checkmark \checkmark \checkmark \checkmark \checkmark$  Econversion between position–velocity and action–angle variables
- $\checkmark$   $\checkmark$   $\checkmark$   $\checkmark$  **>** Distribution functions
  - $\checkmark$   $\checkmark$  **Streams modelling** 
    - ✓ ► DF-based self-consistent models
    - ✓ ► Orbit-superposition (Schwarzschild) models
  - $\checkmark$   $\checkmark$   $\blacktriangleright$  Jeans models

# Galactic modelling software

Torus Mapper [McMillan & Binney 2008; Binney & McMillan 2016]

TACT (the Action Computation Toolbox) [Sanders & Binney 2012–2016]

Galpy [Bovy 2015]

Gala [Price-Whelan 2017]

Agama [Vasiliev 2019]

# AGAM - All-purpose galaxy modeling architecture

- Extensive collection of gravitational potential models (analytic profiles, azimuthal- and spherical-harmonic expansions) constructed from smooth density profiles or N-body snapshots;
- 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: 1d,2d,3d spline interpolation, penalized spline fitting and density estimation, multidimensional sampling;
- Efficient and carefully designed C++ implementation, examples, Python and Fortran interfaces, plugins for Galpy, NEMO, AMUSE.

arXiv:1802.08239, 1802.08255 https://github.com/GalacticDynamics-Oxford/Agama