

Demystifying the fusion mechanism in heavy ion collisions within full Langevin dissipative dynamics

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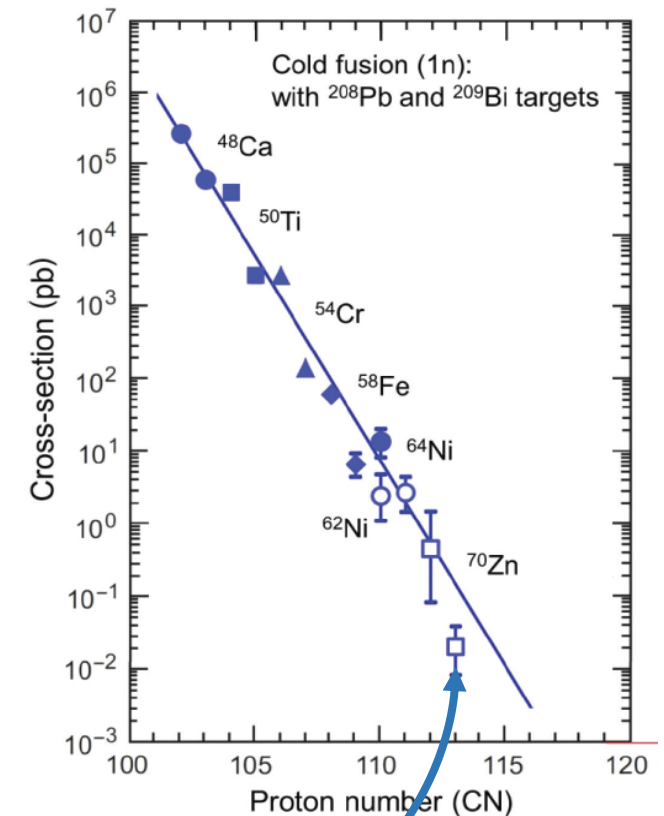
Description of the fusion mechanism within Langevin dynamics

Super-heavy elements with $Z > 103$ do not occur in nature.

They can only be produced in the laboratory by fusing two lighter nuclei.

Our primary goal is to gain insights into the **fusion reaction mechanisms in the domain of cold synthesis reactions** ($Z < 113$, $E^* \approx 10 - 20$ MeV), in particular on the **understanding of the hindrance mechanism which prevents the formation of super-heavy nuclei.**

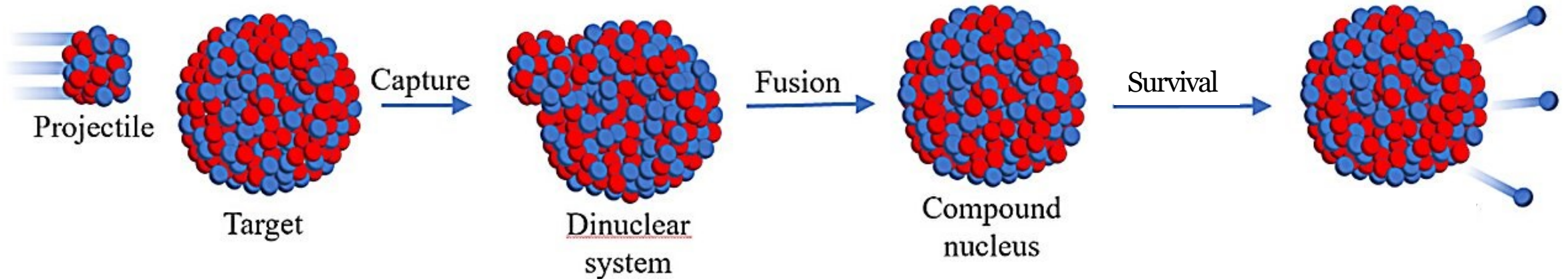
We propose a comprehensive dissipative dynamics **Langevin**-based formalism to describe the unrestricted motion of the systems in terms of **elongation**, **neck** and **asymmetry** variables.



**$Z = 113$, 22fb
(Only 3 atoms in
576 days of irradiation)**

Credits: T. Cap, M. Kowal

The fusion process (Schematic view)



Micro-macroscopic description of fusion:

- ▶ Complexity/impossibility of tracking all internal degrees of freedom
- ▶ Identification of **slow collective degrees of freedom** immersed in a **bath** of faster dynamics
- ▶ Emergence of the mechanisms of **friction** and **random forces**

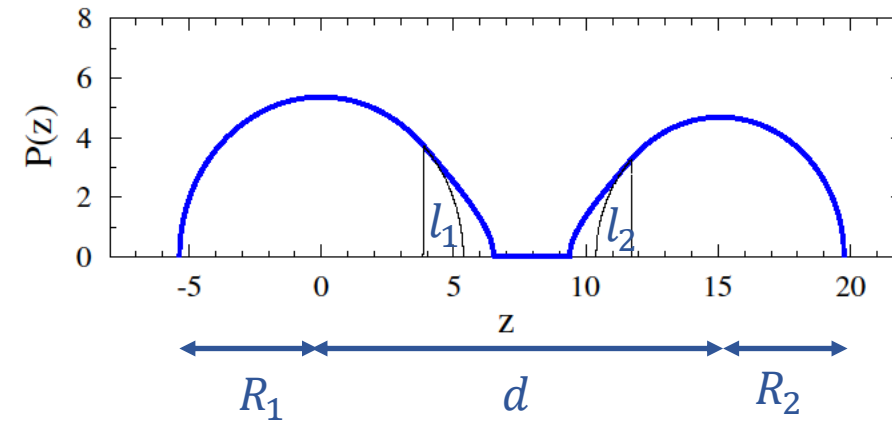
Collective variables adapted to fusion/fission – Shape variables

- ▶ Axially symmetric shapes
- ▶ **Spherical** cups connected by quadratic surfaces^[1]

▶ **Shape collective/slow variables:**

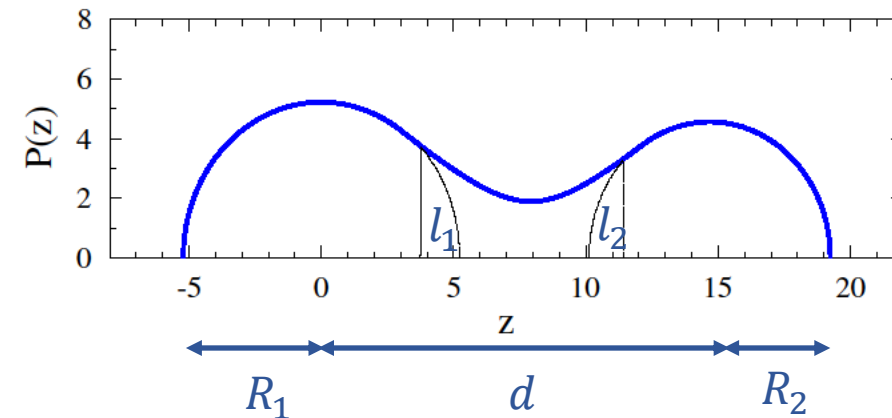
- ▶ **Distance/elongation:** $\rho = \frac{d}{R_1 + R_2}$
- ▶ **Neck/deformation:** $\lambda = \frac{l_1 + l_2}{R_1 + R_2}$
- ▶ **Asymmetry:** $\Delta = \frac{R_1 - R_2}{R_1 + R_2}$

$^{92}\text{Zr} + ^{64}\text{Ni}$



Bipartite

$$\begin{aligned} \rho &= 1.5 \\ \lambda &= 0.3 \\ \Delta &= \Delta_0 \end{aligned}$$



Monopartite

$$\begin{aligned} \rho &= 1.5 \\ \lambda &= 0.4 \\ \Delta &= \Delta_0 \end{aligned}$$

[1] J. Błocki, H. Feldmeier and W. J. Świątecki, Nucl. Phys. A 459 (1986) 145

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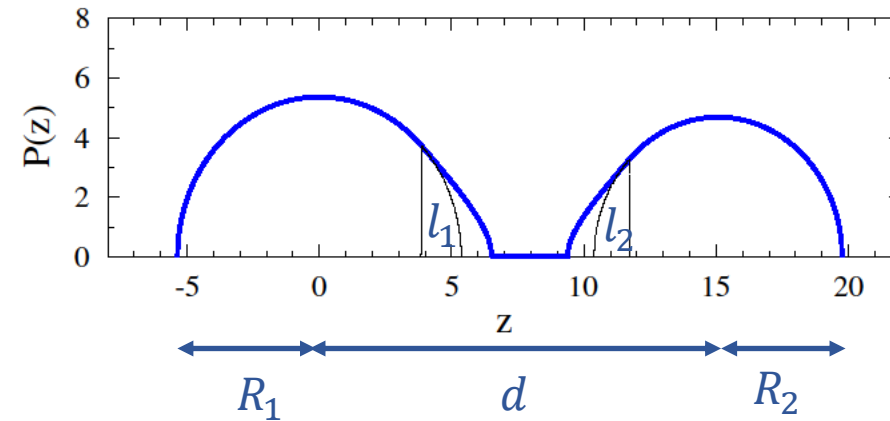
▶ **Neck/deformation:** $\lambda = \frac{l_1 + l_2}{R_1 + R_2}$

▶ **Asymmetry:** $\Delta = \frac{R_1 - R_2}{R_1 + R_2}$

▶ **Scission is well-defined:** $\lambda_{\text{scission}} = 1 - \frac{1}{\rho_{\text{scission}}}$

→ **Suited to describe fusion/fission**
(vs. multipole moments)

$^{92}\text{Zr} + ^{64}\text{Ni}$

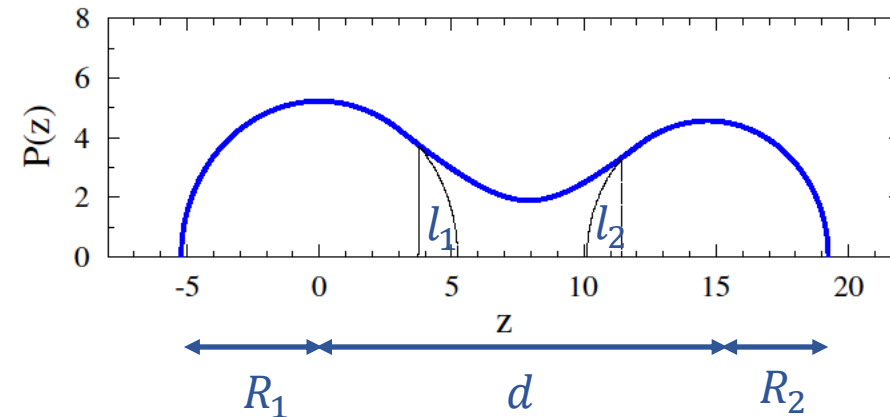


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Collective variables adapted to fusion/fission – Angle variables

- ▶ **Collective angle variables**

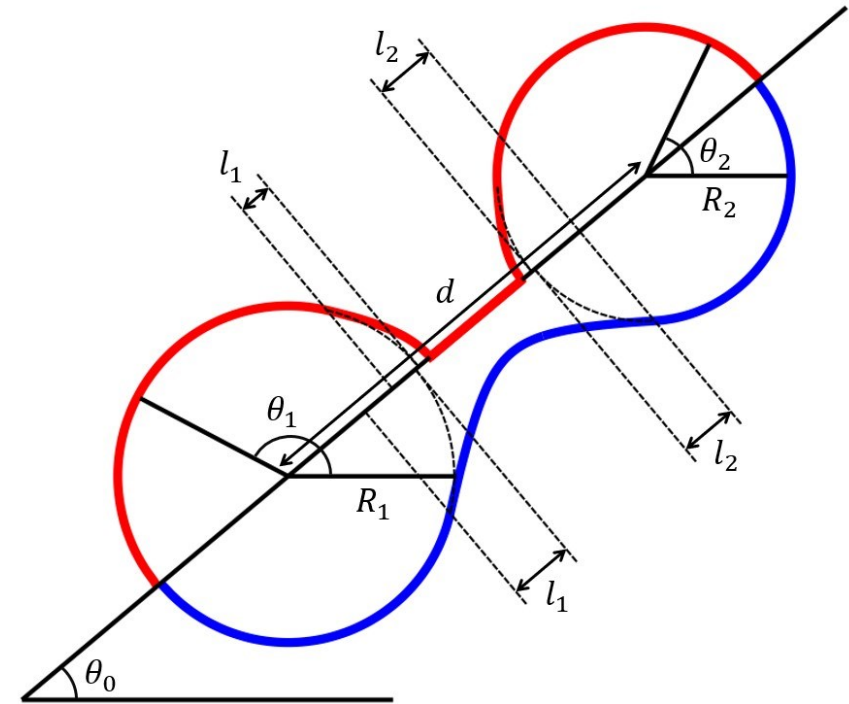
- ▶ Angle of the whole system θ_0
- ▶ Angle of the first sphere θ_1
- ▶ Angle of the second sphere θ_2

- ▶ Variations linked to angular momentum, in particular:

$$p_{\theta_0} + p_{\theta_1} + p_{\theta_2} = L_{init}$$

- ▶ **Exact treatment of angular momentum**

→ Full Langevin 6-dimensional dissipative dynamics



The Langevin system of equations

- ▶ Denoting **collective/slow variables** $q_i(t)$ and their **associated moments** $p_i(t)$, the **Langevin equations** read:

$$\dot{q}_i(t) = \sum_k (\mathcal{M}^{-1})_{ik} p_k \quad \Leftrightarrow (P = MV)$$

$$\dot{p}_i(t) = -\frac{\partial H}{\partial q_i} - \sum_k \gamma_{ik} \dot{q}_k + \sum_k g_{ik} \xi_k(t) \quad \Leftrightarrow \left(\frac{dP}{dt} = \Sigma F \right)$$

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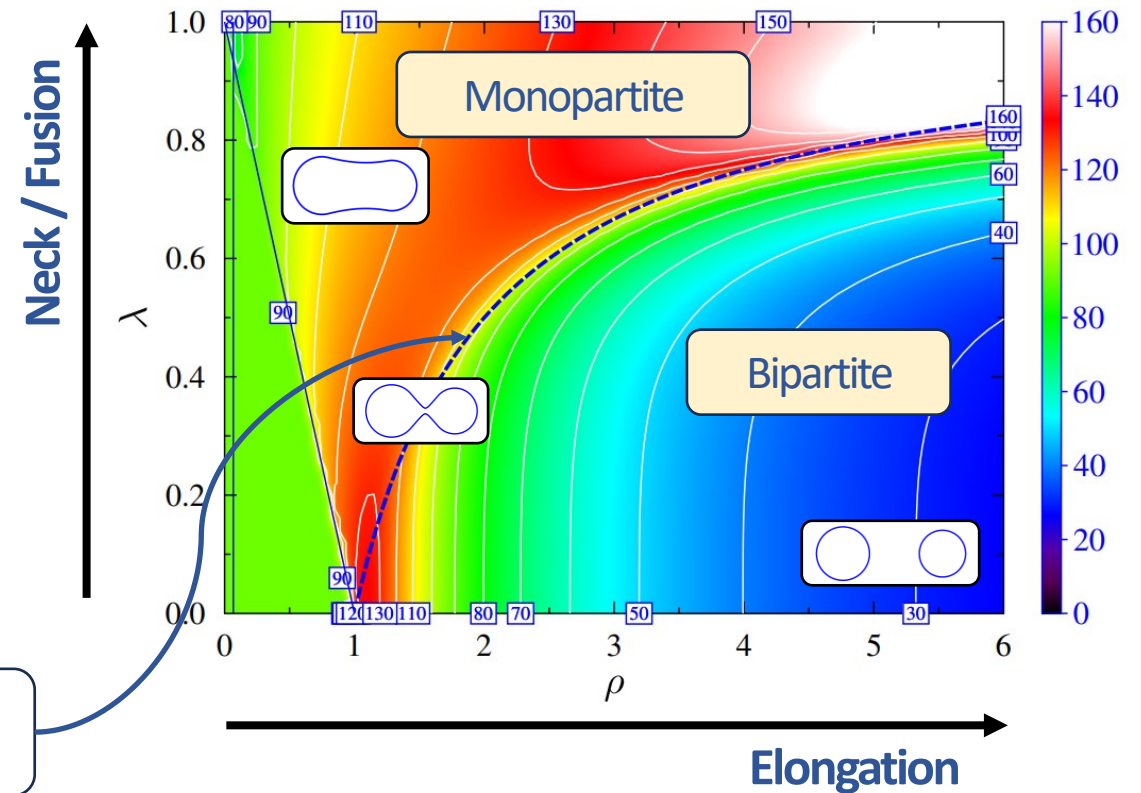
The diagram illustrates the Langevin system of equations. A light blue box labeled "Mass tensor" has an arrow pointing to the first equation. Three yellow boxes at the bottom are labeled "Conservative forces (H = T + V)", "Friction forces", and "Langevin/random forces". Arrows point from these boxes to the corresponding terms in the second equation.

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→ A comprehensive understanding of the dynamics process (in comparison to the random walk f. eg.).

Potential Energy

- ▶ **Yukawa-plus-exponential folding potential + Coulomb**
- ▶ Parameters taken from a previous fit to experimental masses and fusion barrier heights [1]
- ▶ No shell effects at the moment.



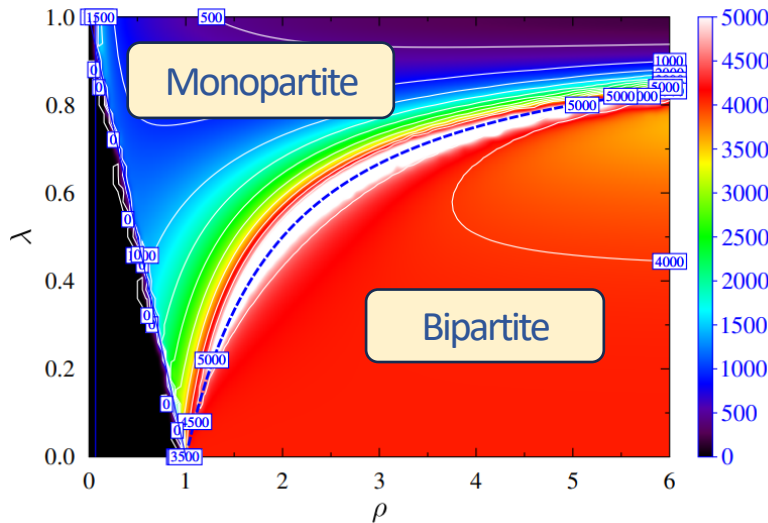
Scission line $\lambda = 1 - \frac{1}{\rho}$

Deformation potential of $^{92}\text{Zr} + ^{64}\text{Ni}$ in MeV

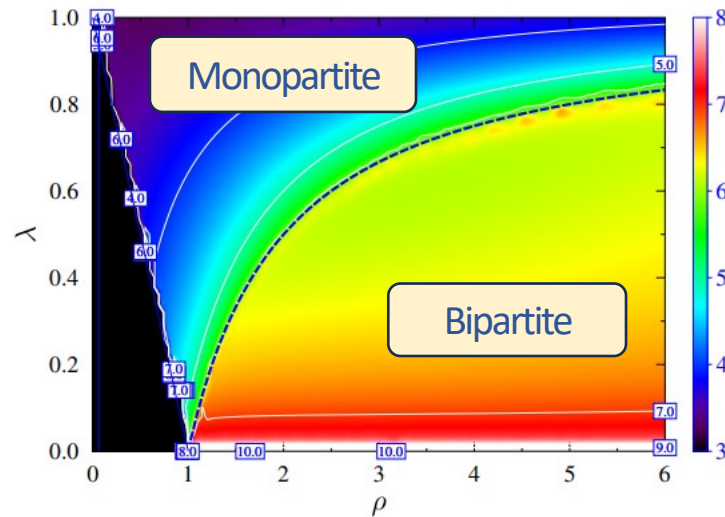
[1] H. J. Krappe *et al.*, Phys. Rev. C 20 (1979) 992–1013

Mass tensor / Kinetic Energy

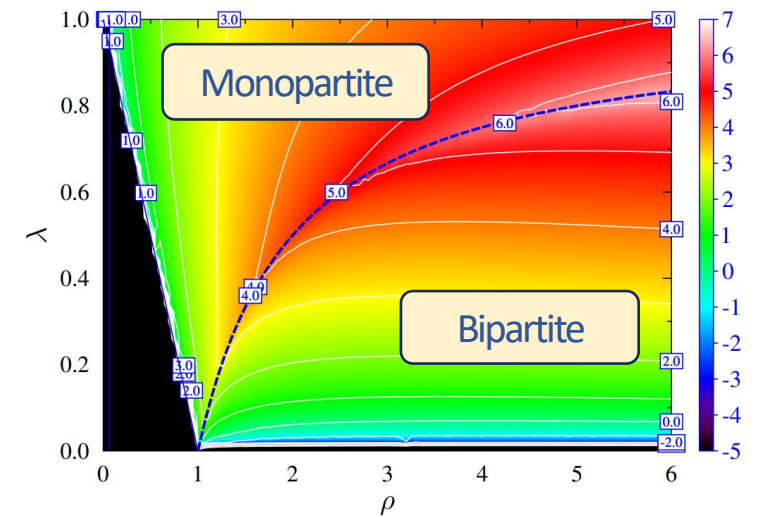
- ▶ **Werner-Wheeler flow approximation:**
 - ▶ **Incompressibility** (matter density is uniformly distributed)
 - ▶ The flow is **irrotational** (the moving planes remain plane)



$\mathcal{M}_{\rho\rho}$



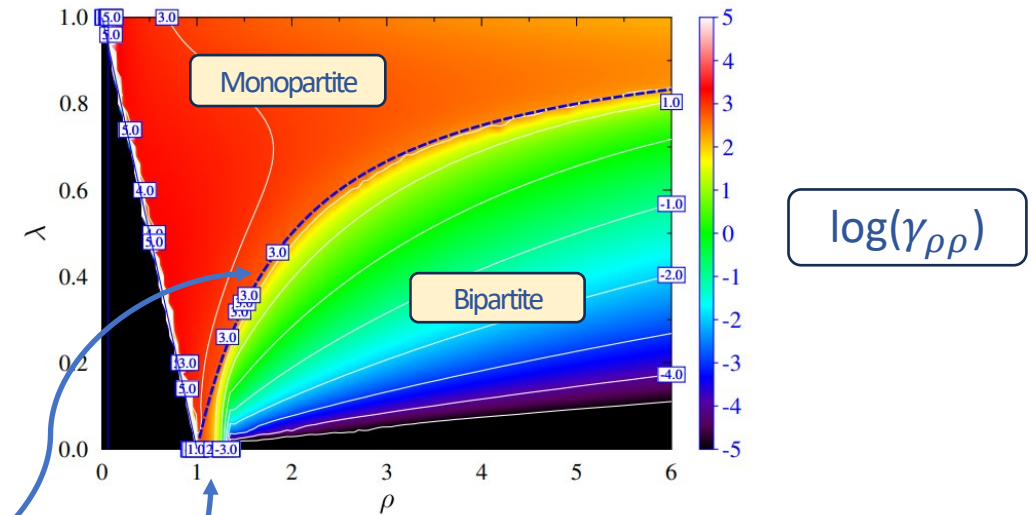
$\log(\mathcal{M}_{\Delta\Delta})$



$\log(\mathcal{M}_{\lambda\lambda})$

Friction forces

- ▶ **Proximity formalism** (to account for some quantum effects):
Possible matter flow/friction before contact ($d = 3.2$ fm)
- ▶ **Shape friction**:
 - ▶ **Wall friction** (collisions nucleons \leftrightarrow nuclear surface)
 - ▶ + **Wall-plus-window friction** (between the two fragments)

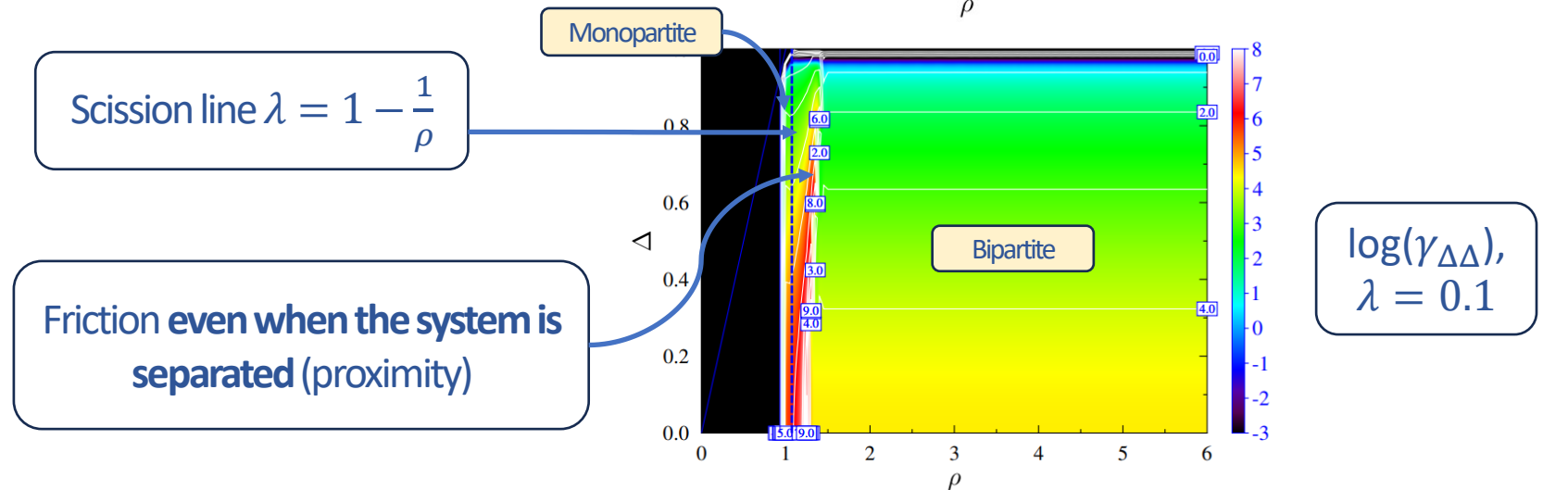
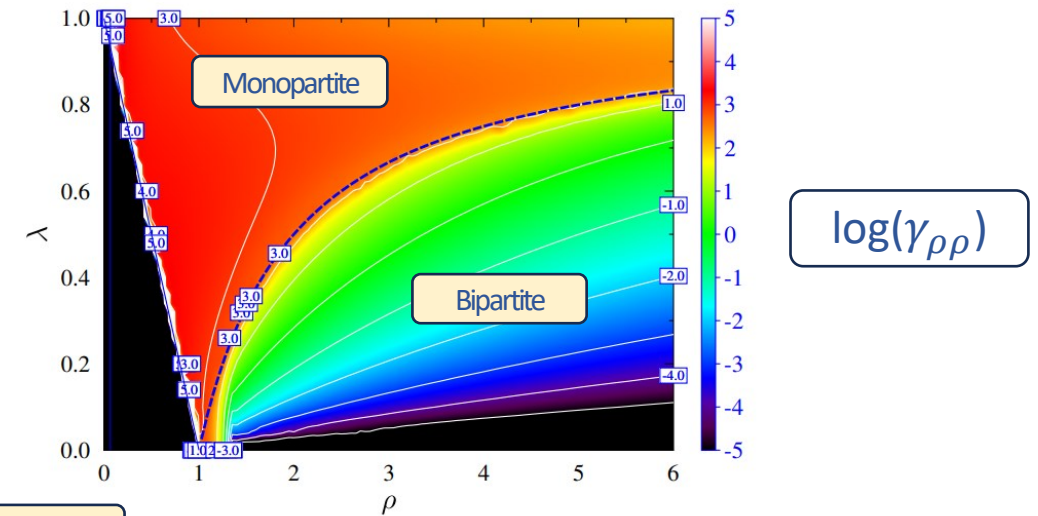


Scission line $\lambda = 1 - \frac{1}{\rho}$

Friction even when the system is separated (proximity)

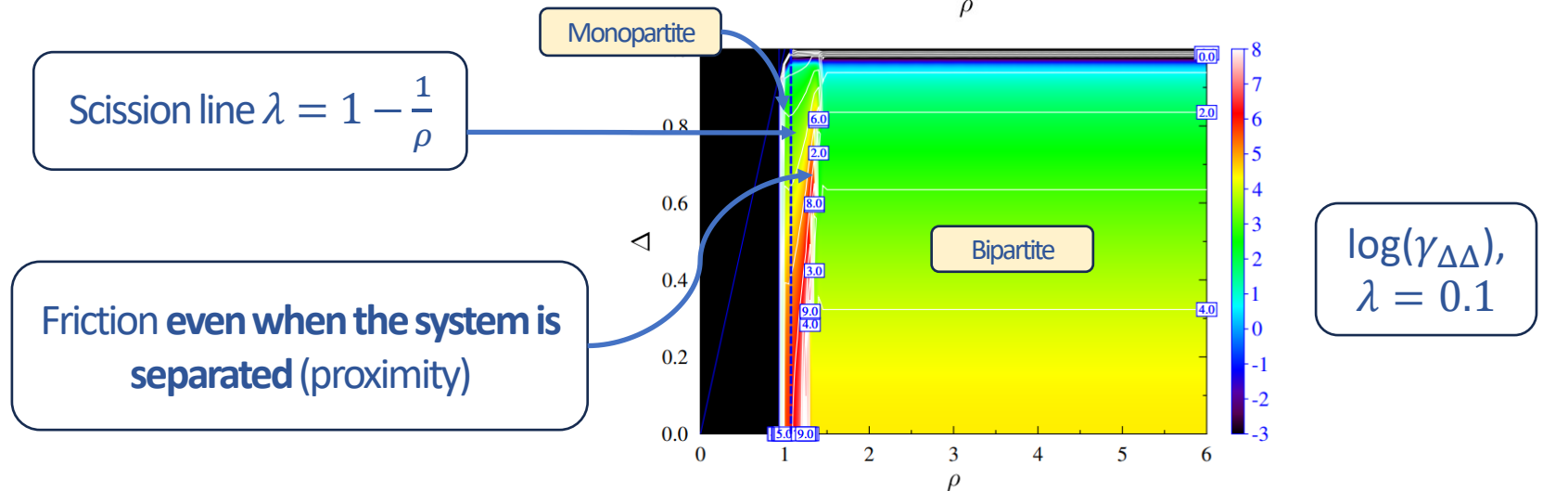
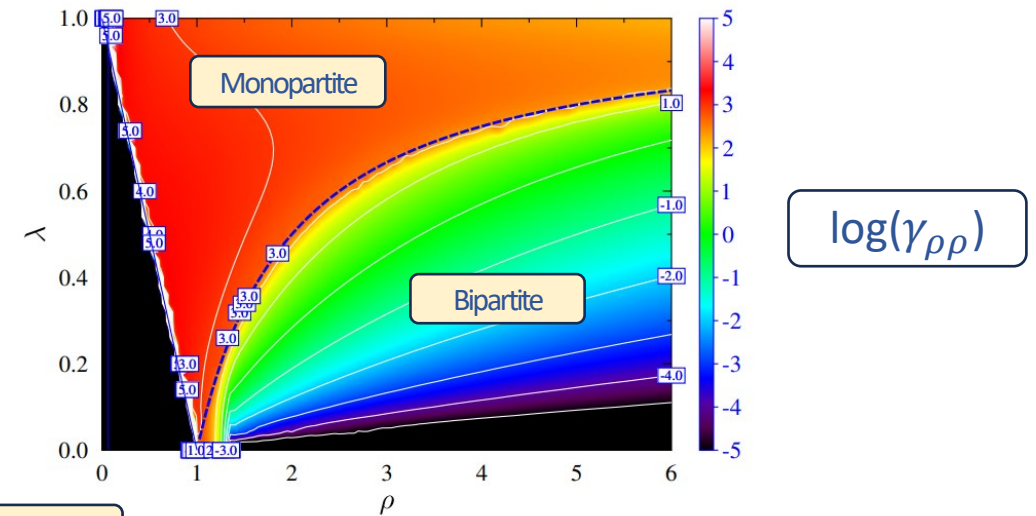
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- ▶ **Angular friction**:
 - ▶ **Sliding friction**
 - ▶ **No rolling friction**



Langevin/random forces

- ▶ We assume a simple memoryless Langevin force (white noise):

$$F_i = \sum_k g_{ik} \xi_k(t)$$

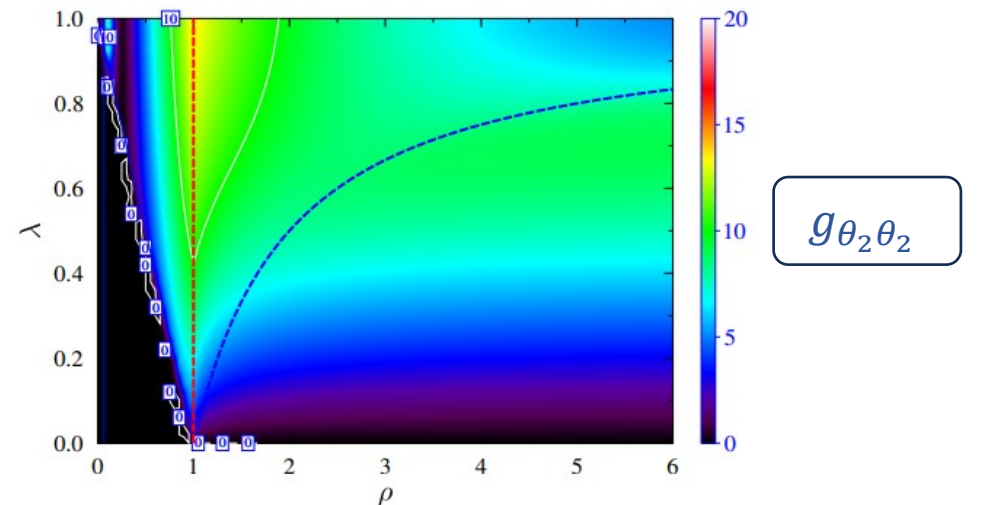
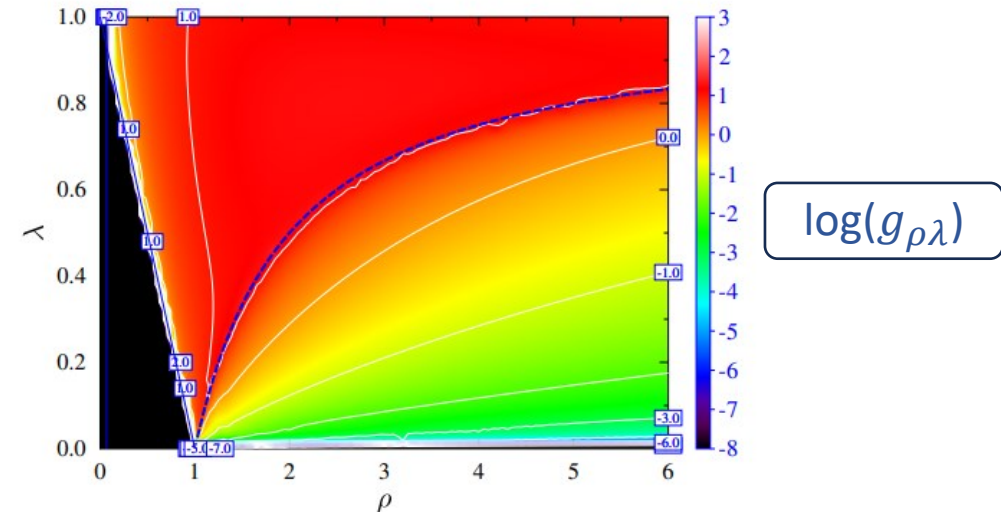
$\xi_k(t)$ are **time-dependent Gaussian random variables**:

$$\begin{aligned} \langle \xi_k(t) \rangle &= 0 \\ \langle \xi_k(t), \xi_{k'}(t') \rangle &= 2\delta_{kk'}\delta(t - t') \end{aligned}$$

- ▶ The diffusion tensor is given by the **Einstein relation**:

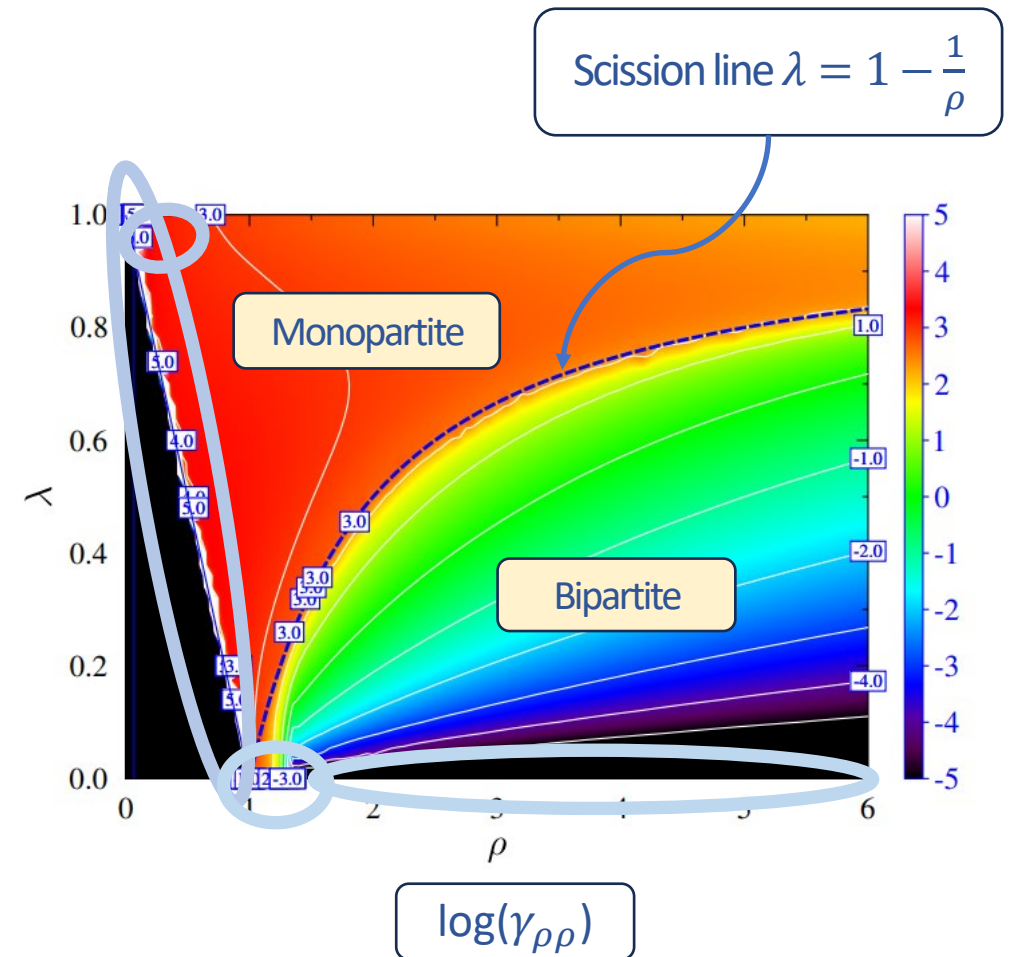
$$\sum_k g_{ik} g_{kj} = D_{ij} = k_B T^* \gamma_{ij}, \quad T^* = \frac{E_0}{\tanh(E_0/T)}, \quad T = \sqrt{E^*/a}$$

$E_0 = 2$ MeV is the zero-point collective energy of the heat bath oscillators
 E^* is the **dissipated energy**, $a = A/8$ MeV is the **level density parameter**.



Physical vs. practical collective variables

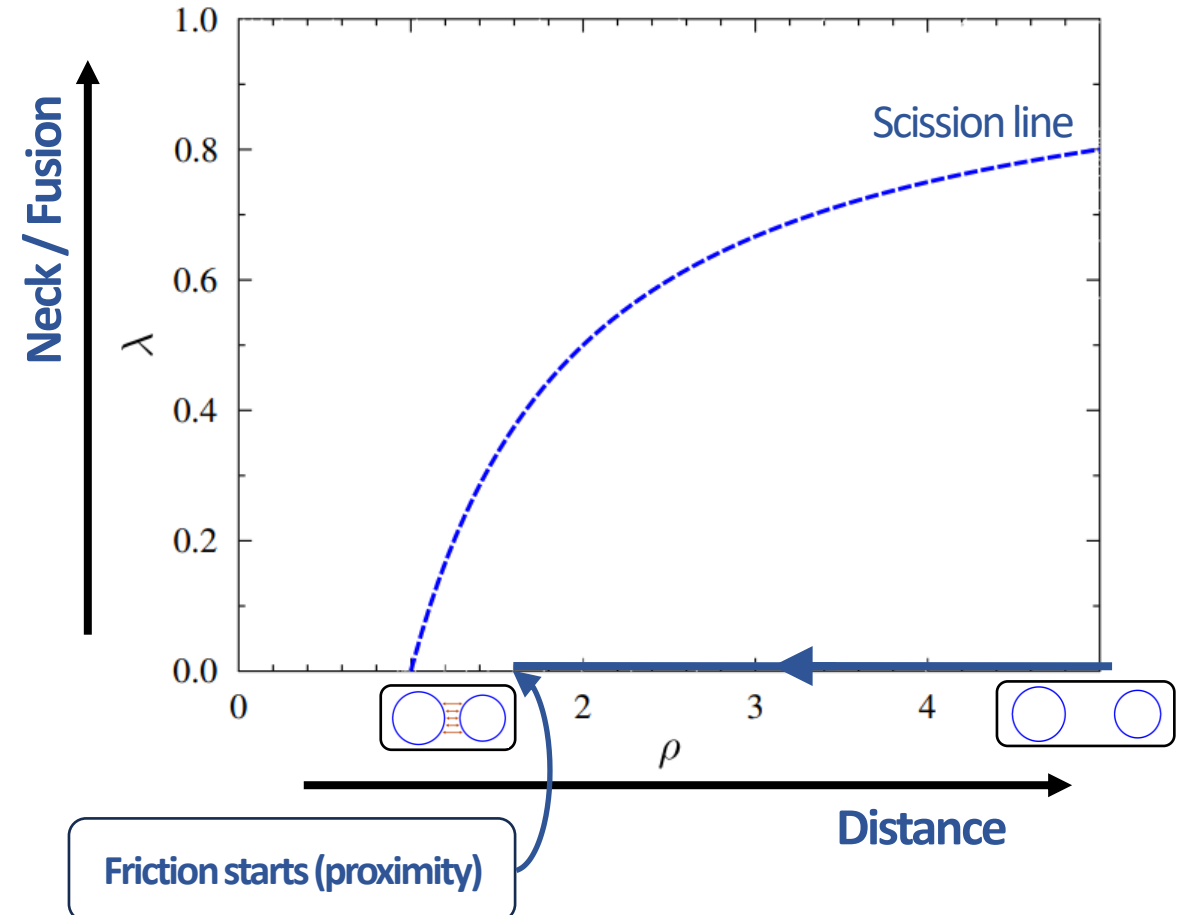
- ▶ Unlike the multipole moments, the (ρ, λ, σ) variables are **extremely irregular**:
 - ▶ Many borders,
 - ▶ Small proximity region,
 - ▶ Regime change.
- ▶ **A correct treatment of numerical precision is needed.**
- ▶ **Extrapolation** is needed for the calculations of the physical quantities and their derivatives.
The regions of extrapolation should **avoid the borders.**
- ▶ The process of fusion itself is not fully tractable numerically.



The fusion process

The three main stages of the collision:

1. **A first violent deceleration** during which:
 - The system loses most of its kinetic energy
 - There is **almost no deformation of the nuclei** as **no interaction has happened yet**
$$(\dot{\lambda} = (\mathcal{M}_{\rho\lambda})^{-1} p_\rho + (\mathcal{M}_{\lambda\lambda})^{-1} p_\lambda = -\infty + \infty)$$
 - **Unstable balance close to the lambda border**
 - **Treated exactly** (conservative forces).

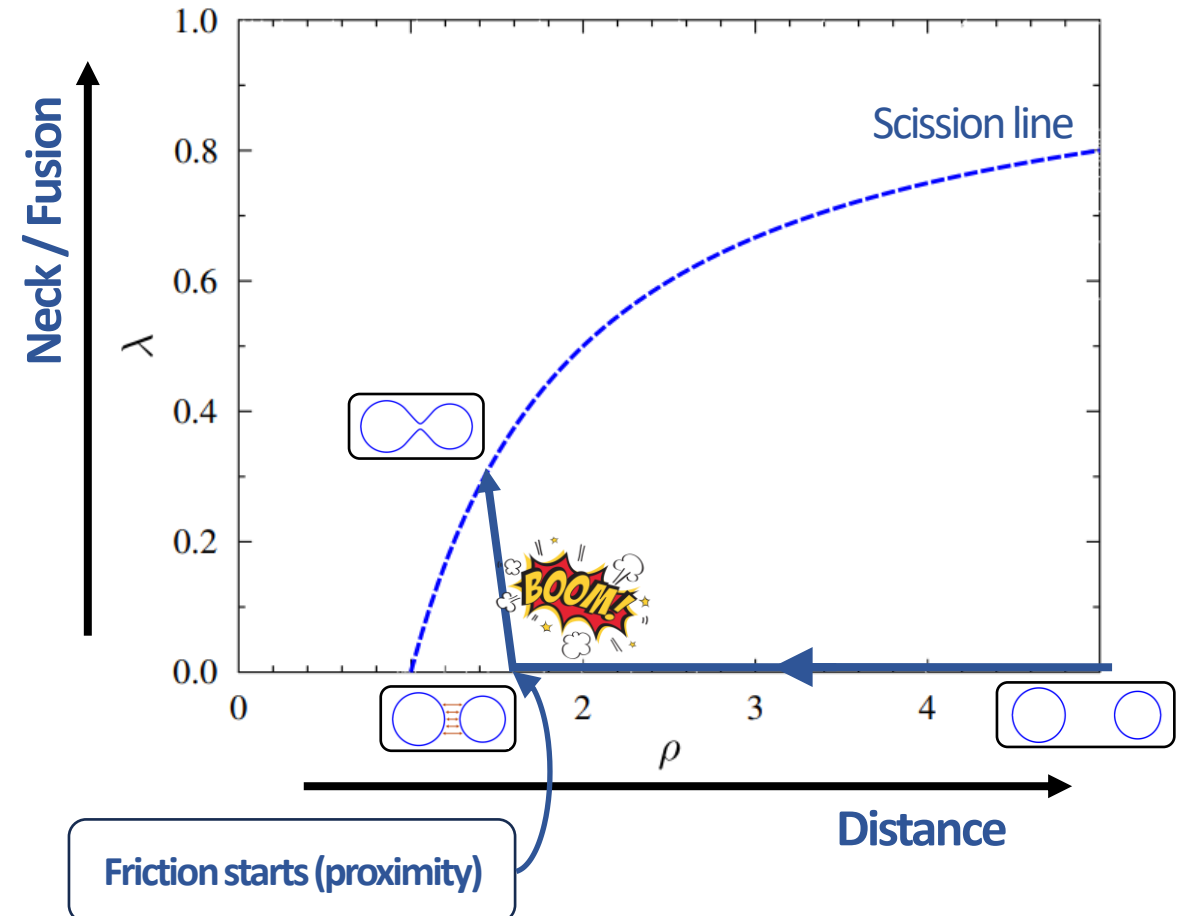


The fusion process

2. The "Kiss of death" when friction starts

- A little sudden change in $p_\rho \rightarrow$ infinite push to the scission line $(\dot{\lambda} = (\mathcal{M}_{\rho\lambda})^{-1} p_\rho + (\mathcal{M}_{\lambda\lambda})^{-1} p_\lambda = -\infty + \infty)$

- Deformation starts and remains.



The fusion process

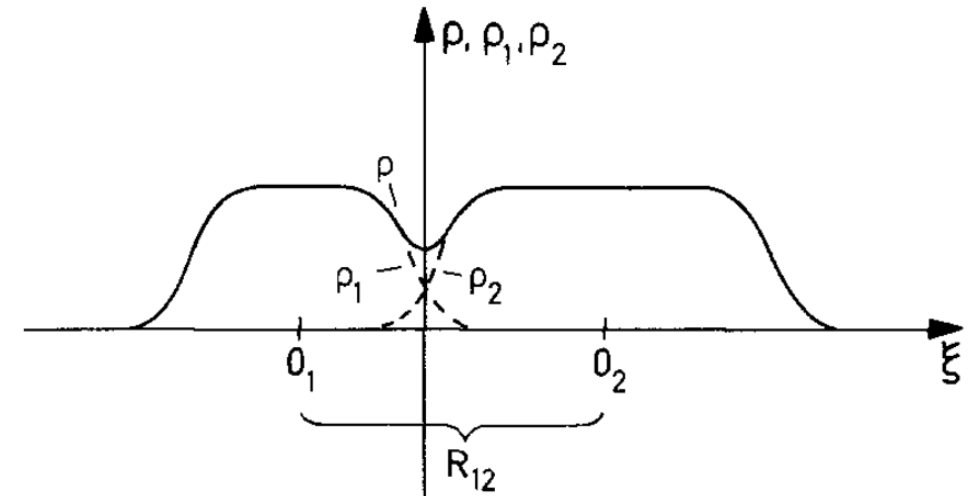
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- Deformation starts and remains.

- **Inherent instability of non-saturated nucleonic densities** (Sudden approximation in HF when the nuclear tails touch)

- **This step is NOT treated numerically:**
We start the calculation from the touching point



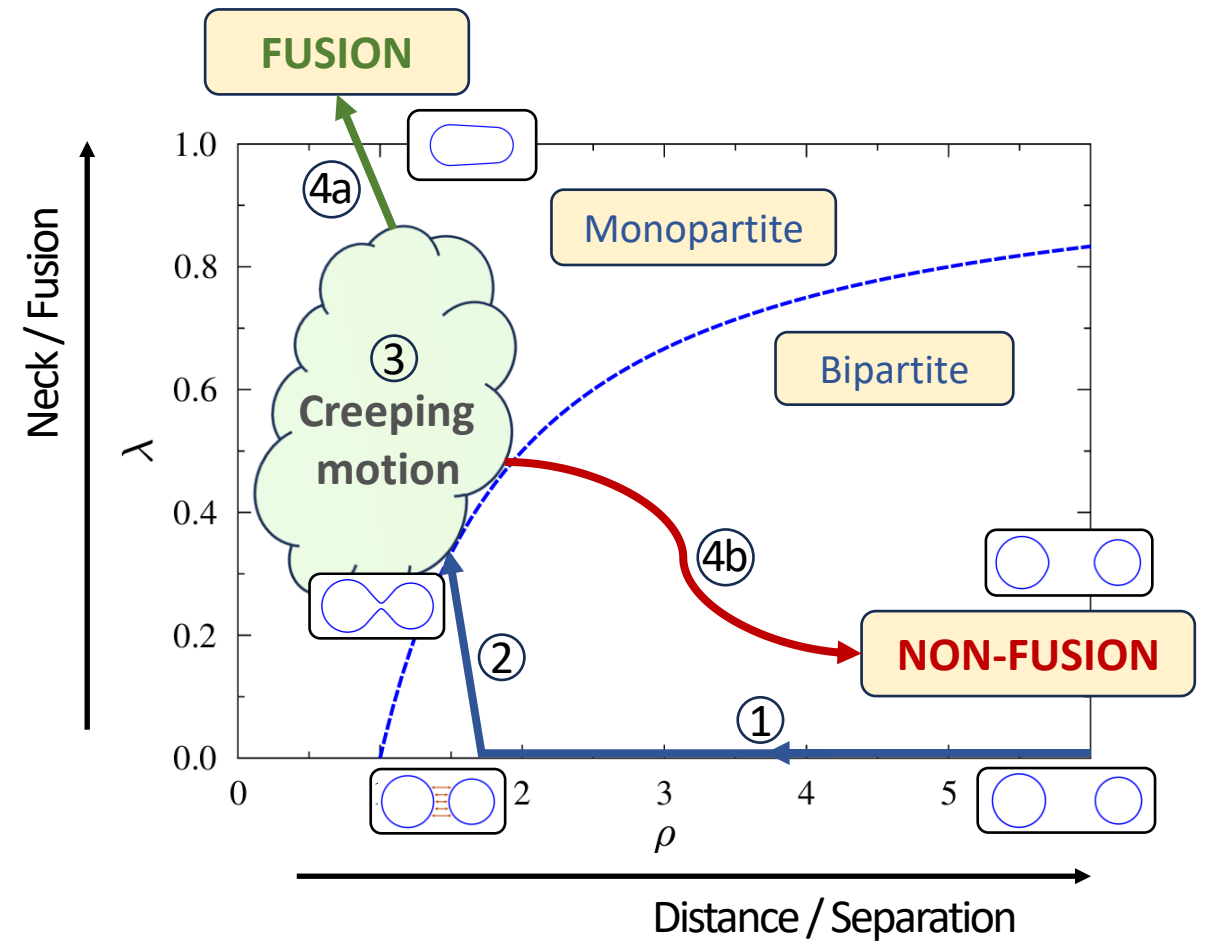
"Sudden approximation" – K. Pomorski, K. Dietrich
Z. Phys. A, 295 (1980) 335

The fusion process

The three main stages of the collision:

3. A **long creeping motion**

- which leads to fusion or separation
- Solved numerically



The observables

- ▶ The resolution of the Langevin equations generates a **distribution of trajectories due to the fluctuation force**.
- ▶ We use **500,000 - 1,000,000 trajectories**.
- ▶ **Asymmetry is free to change**.
- ▶ Calculations performed on the **CiŚ cluster** (Świerk/Warsaw).

- ▶ The **spin distribution** is calculated as a Monte-Carlo integral on a given bin $i \equiv \ell_i$:

$$\sigma_\ell = \left(\frac{d\sigma_{fus}}{d\ell} \right)_{\ell_i} = \frac{2\pi}{k^2} \ell_i^2 \frac{N_i^{fus}}{N_i^{tot}}$$

where $\ell_{init} = \ell_{max}\sqrt{x}$, x a random number in $[0,1]$ (for easy derived formulas).

- ▶ From the spin distribution, one can calculate:
 - ▶ The **total cross section / probability for the formation of the compound nucleus**
 - ▶ $\langle \ell \rangle, \langle \ell^2 \rangle$
 - ▶ **Excitation functions.**

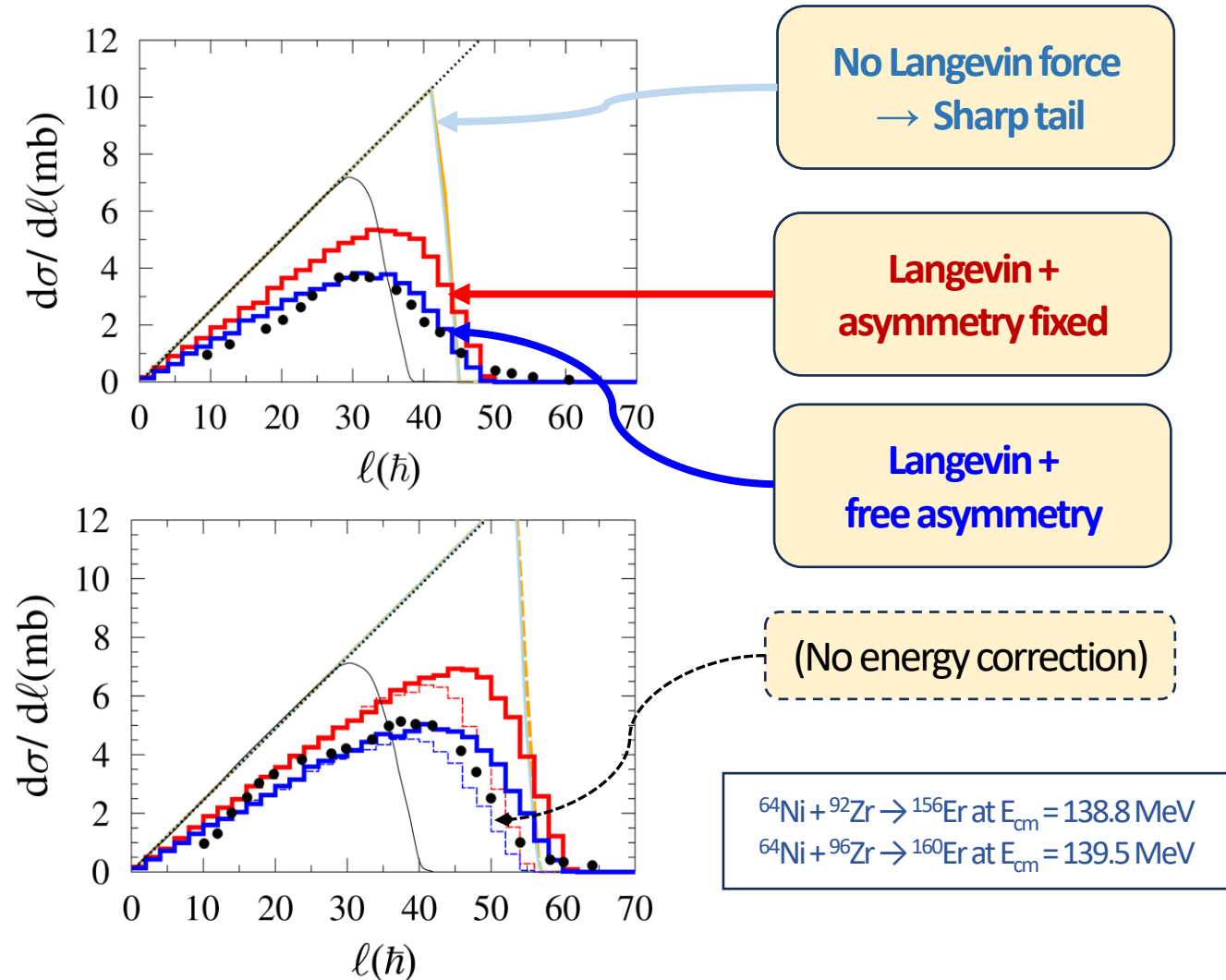
$^{64}\text{Ni} + ^{92,96}\text{Zr} \rightarrow ^{156,160}\text{Er}$ – Effect of the Asymmetry

- ▶ $^{64}\text{Ni} + ^{92}\text{Zr}$: $Q_{\text{fus,calc}} = Q_{\text{fus,exp}} \rightarrow$ no correction needed
- ▶ $^{64}\text{Ni} + ^{96}\text{Zr}$: Difference of 3.5 MeV \rightarrow **correction needed**
- ▶ Spin distributions are more natural with the Langevin force (vs. Ref. [1])
- ▶ **Great agreement to the experimental data**^[2,3].
- ▶ In the $^{64}\text{Ni} + ^{96}\text{Zr}$ case, little discrepancies in the spin distributions (deformation?)
- ▶ **The total cross section (160 mb) is well reproduced.**

[1] W. Przystupa, K. Pomorski, Nucl. Phys. A 572(1) (1994) 153

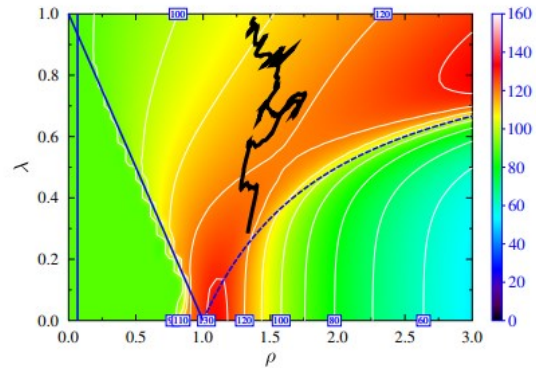
[2] W. Kuhn *et al.*, Phy. Rev. Lett. 62 (1989) 1103

[3] A. M. Stefanini *et al.*, Phys. Lett. B 252 (1990) 43

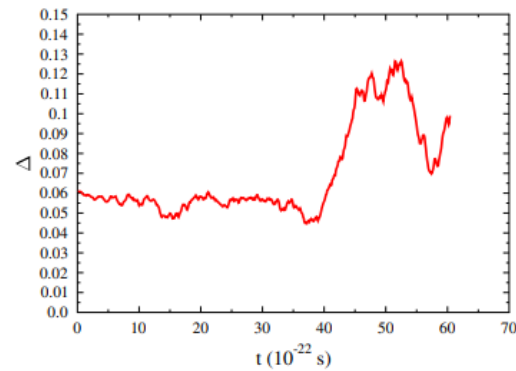


A deep understanding of the fusion process (Preliminary examples)

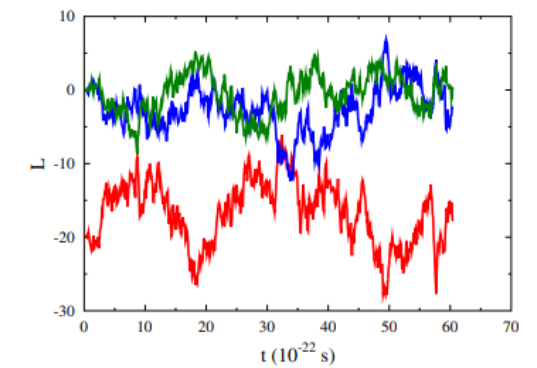
Paths



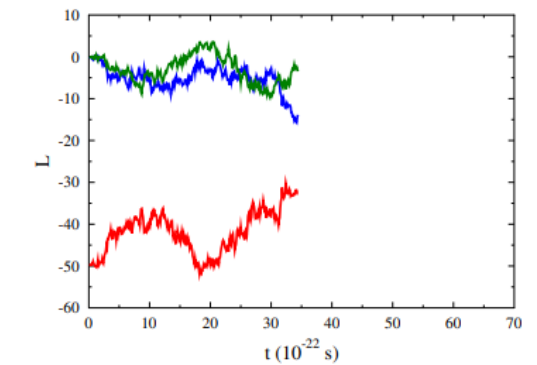
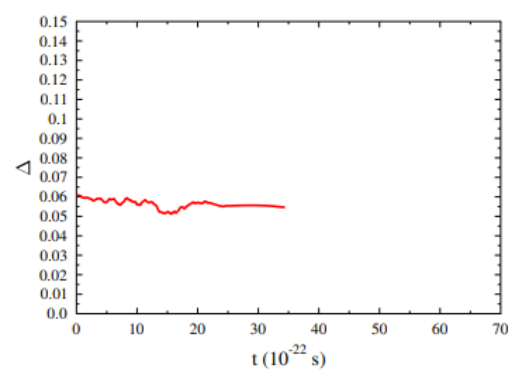
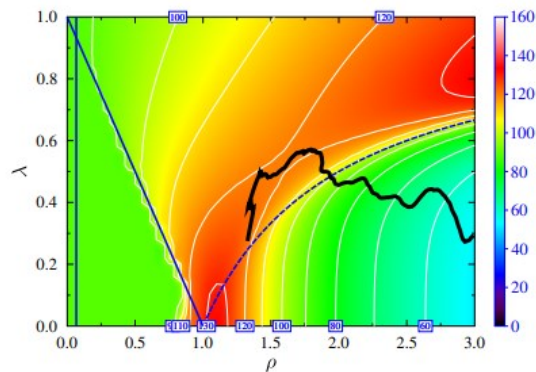
Asymmetry vs. time



Angular momenta



- ▶ $\ell = 20\hbar$,
Fusion case:



- ▶ $\ell = 50\hbar$,
No-fusion case:

- ▶ And various other observables: **diffusion tensor**, the **number of rotations the system undergoes before fusion etc.**

Summary and Perspectives

- ▶ We have derived a fully **6-dimensional dissipative dynamics Langevin**-based formalism to describe the **unrestricted motion of the systems** in terms of **elongation, neck** and **asymmetry** variables.
- ▶ Thanks to a **correct treatment of the different stages of fusion**, the spin distributions are now **in great agreement with experimental data**.
- ▶ The Langevin formalism allows for a deep understanding of the fusion process: evolution of the asymmetry, angular momentum/rotations of the fragments etc.
- ▶ In the very near future:
 - ▶ We will study the effect of the **asymmetry of entrance channel** on the formation of the compound system.
 - ▶ We will tackle the **hindrance problem** by comparing the $^{48}\text{Ca}/^{50}\text{Ti}/^{54}\text{Cr} + ^{208}\text{Pb}$ systems.
- ▶ We are planning to make the following improvements of the formalism:
 - ▶ The addition of **shell effects** for a **fully microscopic-macroscopic picture**
 - ▶ The testing of different forms of stochastic noises (**color noises**), which will allow us to explore **memory effects** as the process evolves.

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Thank you for
your attention!