



1 Hydrodynamic electrons?

In clean, **two-dimensional materials**, the electrons can reach a **hydrodynamic** transport regime, flowing around obstacles like water.



- The lower carrier densities in 2D means a much lower Fermi temperature than in 3D. The electron-electron decay rate, increasing as $f(T/T_F)$, can therefore become large without the material melting. The electron fluid can then reach local equilibrium and behave hydrodynamically.
- This makes perturbative expansions in T/T_F poor. Hence **there is a need** for an **exact solution** to the governing Fermi liquid equations describing the electrons.

3 Methods

We study the linearized kinetic equation describing the electrons semi-classically:

$$\text{Decay rate} \quad \gamma \chi = \mathcal{L}[\chi]$$

Linearized collision operator

Fermi surface deformation

We evaluate matrix elements $\langle \chi | \mathcal{L} | \psi \rangle = \frac{m^* \lambda_T^2}{16\pi \hbar^3} \int \frac{d\mathbf{P} d\mathbf{q} d\Omega}{(2\pi)^4} |V|^2 F_{121'2} [\bar{\chi}(\mathbf{p}_1) + \bar{\chi}(\mathbf{p}_2) - \bar{\chi}(\mathbf{p}'_1) - \bar{\chi}(\mathbf{p}'_2)] [\bar{\chi} \rightarrow \psi]$

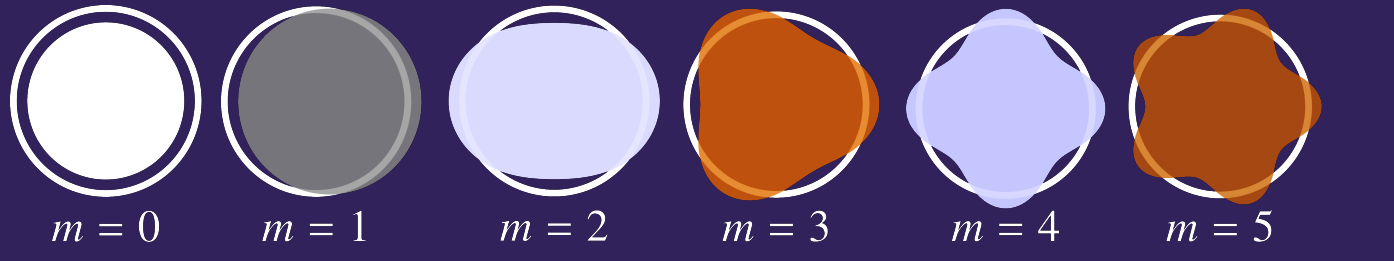
Eigenvalues γ_m

Product of Fermi factors, sharply peaked at low T

Screened Coulomb interaction

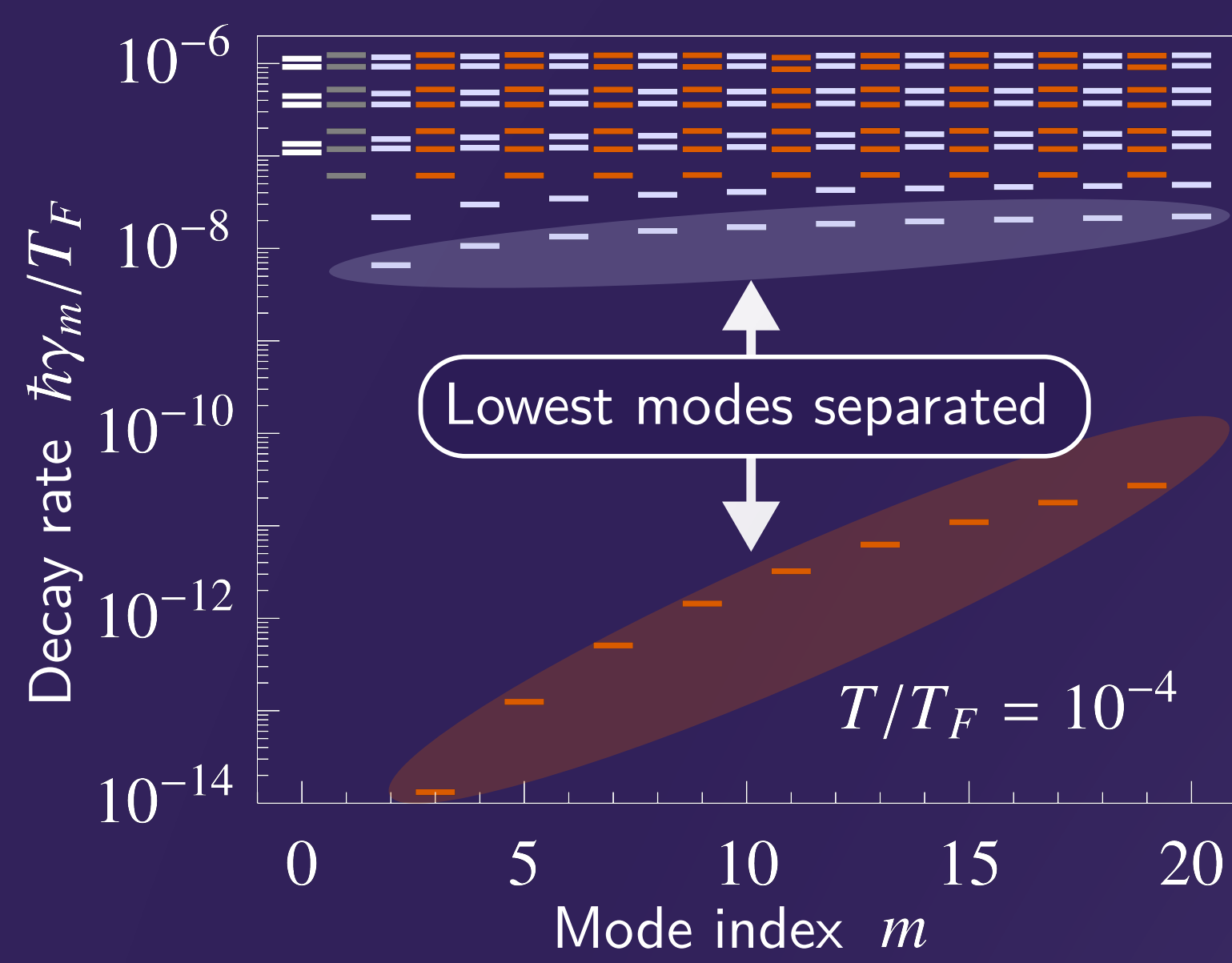
$$V(q; r_s) = \frac{2\pi e^2}{q + \sqrt{2} k_F r_s}$$

By expanding in angular harmonics and orthogonal polynomials

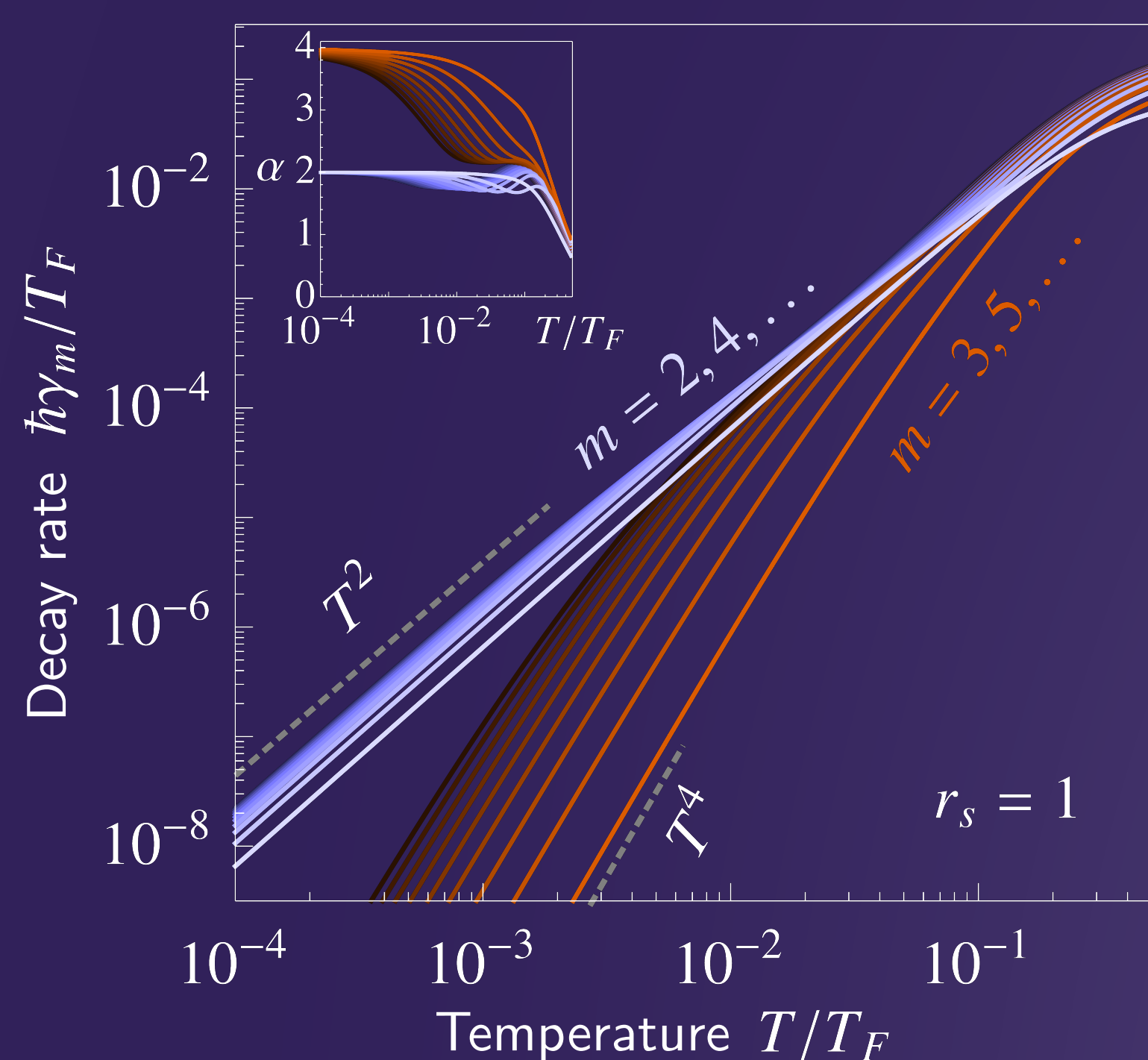
$$\chi(\mathbf{p}) = \sum_m \sum_{\ell=1}^N c_{\ell} u_{\ell}(\mathbf{p}) e^{im\theta}$$


Divonne algorithm
(CUBA library, c)

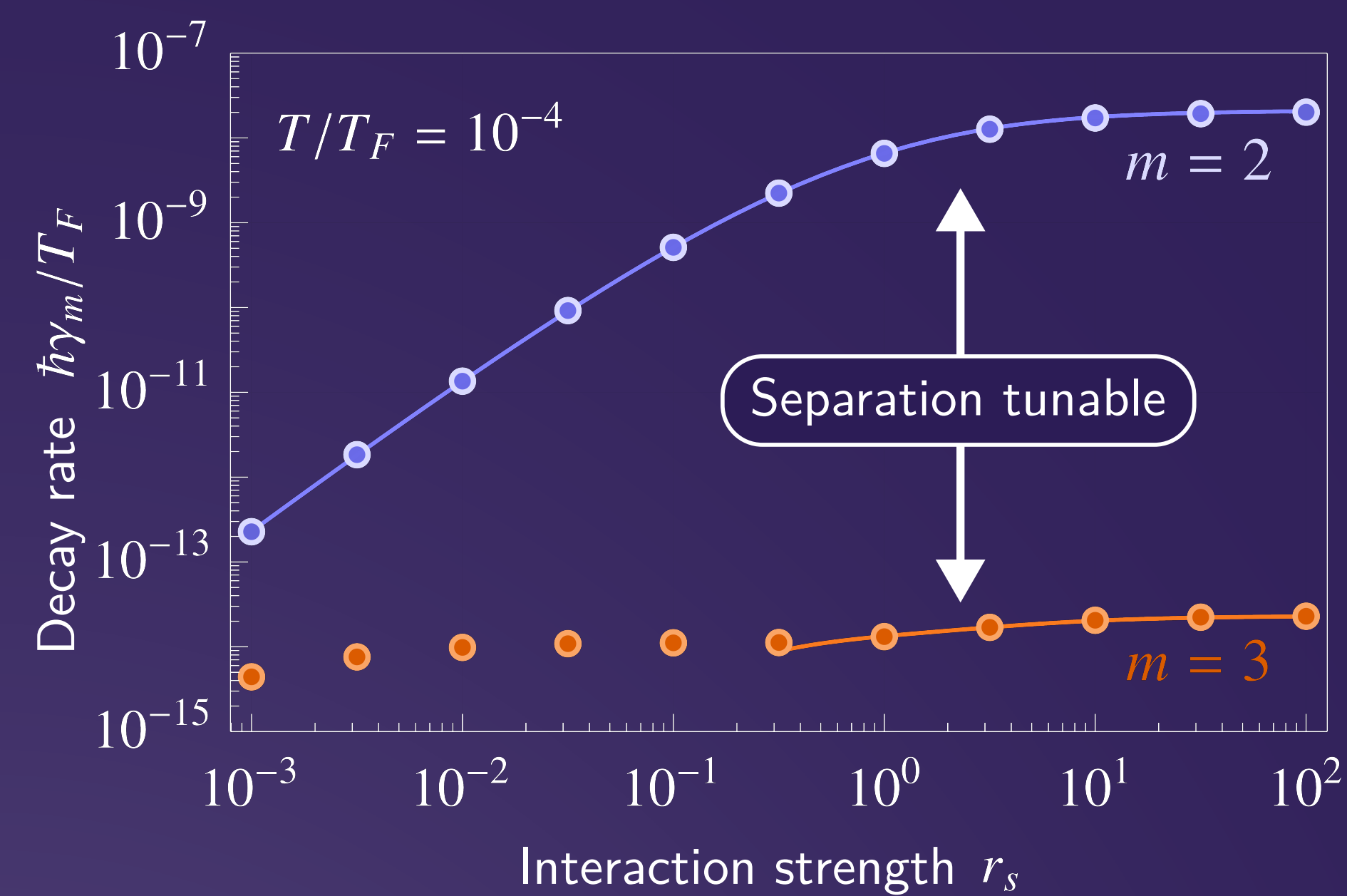
4 Results



We are able to obtain a **complete spectrum** of modes of the collision operator. There exists a set of **long-lived modes**, with a decay rate separated by several orders of magnitude from the rest.



The separation remains up to **experimentally realizable temperatures**. The onset of a regime depends on the angular mode index.



The separation between even and odd parity modes **can be tuned** by the interaction strength between the electrons. Experiments should choose the doping and substrate environment to make the effect as large as possible.

5 Conclusions

We have formulated a numerically exact method of describing the interacting 2D electron gas, going beyond the relaxation time approximation or perturbative expansions. This has enabled us to fully characterize a set of long-lived, odd-parity modes that play a role in a tomographic transport regime. The method also allows for the calculation of general linear-response transport coefficients, such as the shear viscosity [2].