

# Using pseudopotentials to study strongly correlated phases

Gareth Conduit

Pascal Bugnion

Richard Needs

Pablo Lopez Rios

Jonathan Lloyd-Williams

Thomas Whitehead

Lars Schonenberg

TCM Group, Department of Physics

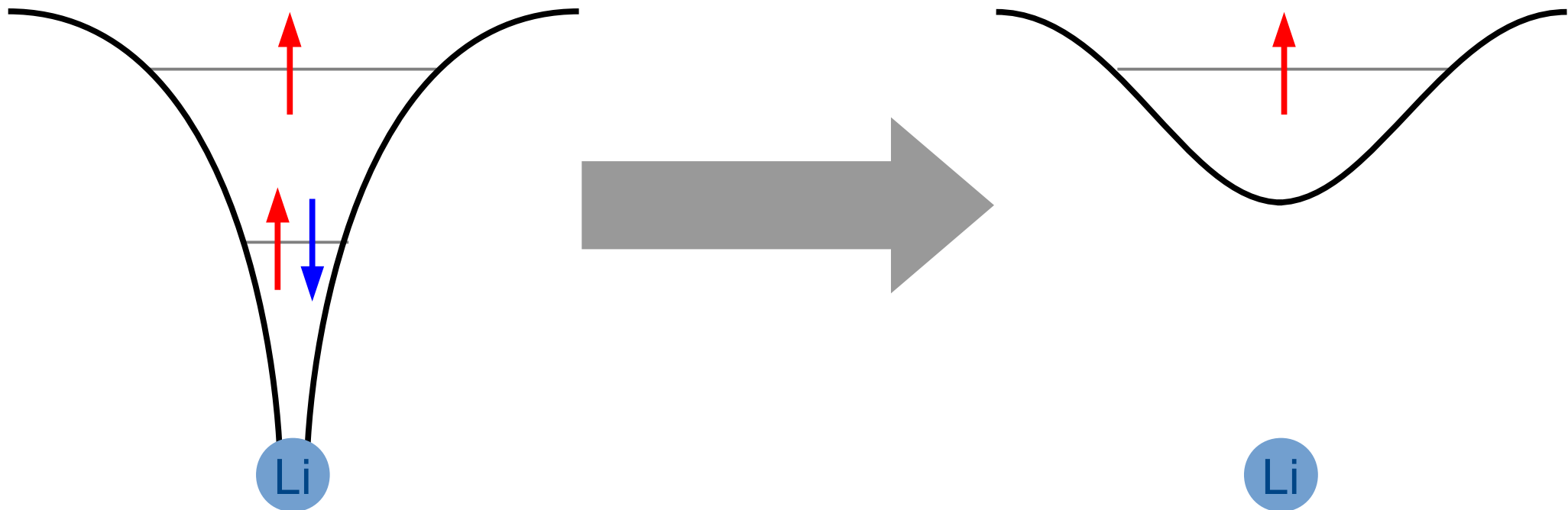
# QMC study of the ground state

$$H = KE + V_{e-i} + V_{e-e}$$

$$E = \frac{\int \bar{\psi} H \psi d\mathbf{r}}{\int \bar{\psi} \psi d\mathbf{r}}$$

# Electron-ion pseudopotential

$$H = KE + V_{e-i} + V_{e-e}$$

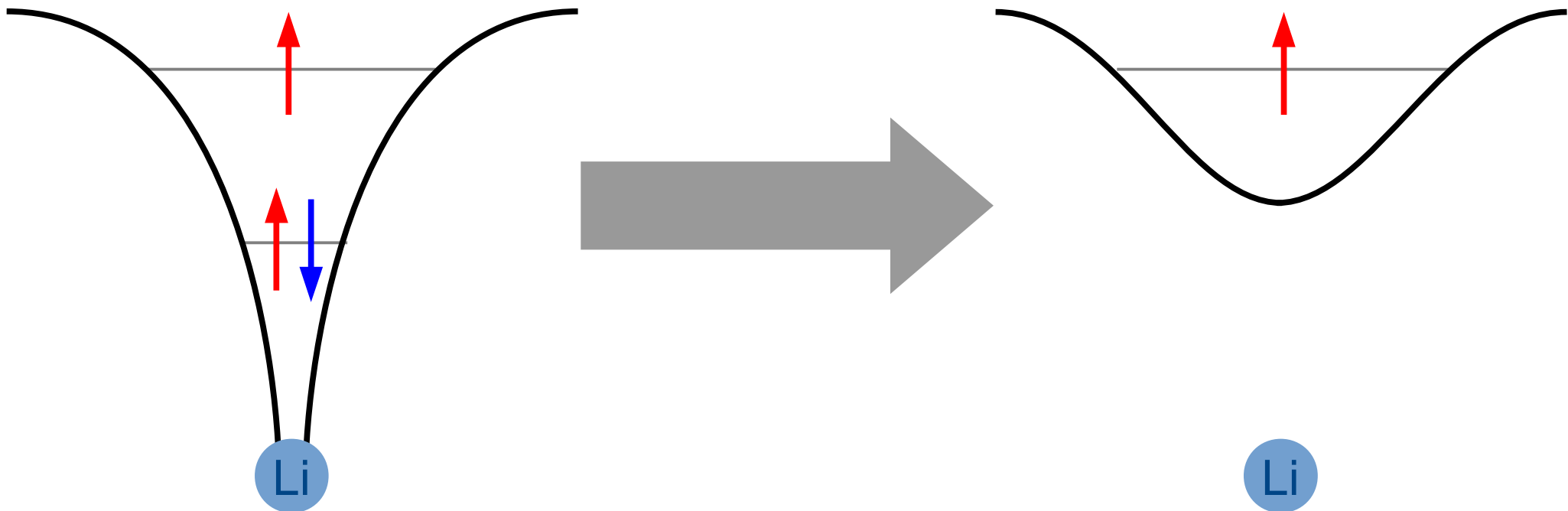


# Electron-ion pseudopotential

$$H = KE + V_{e-i} + V_{e-e}$$

Smooth background

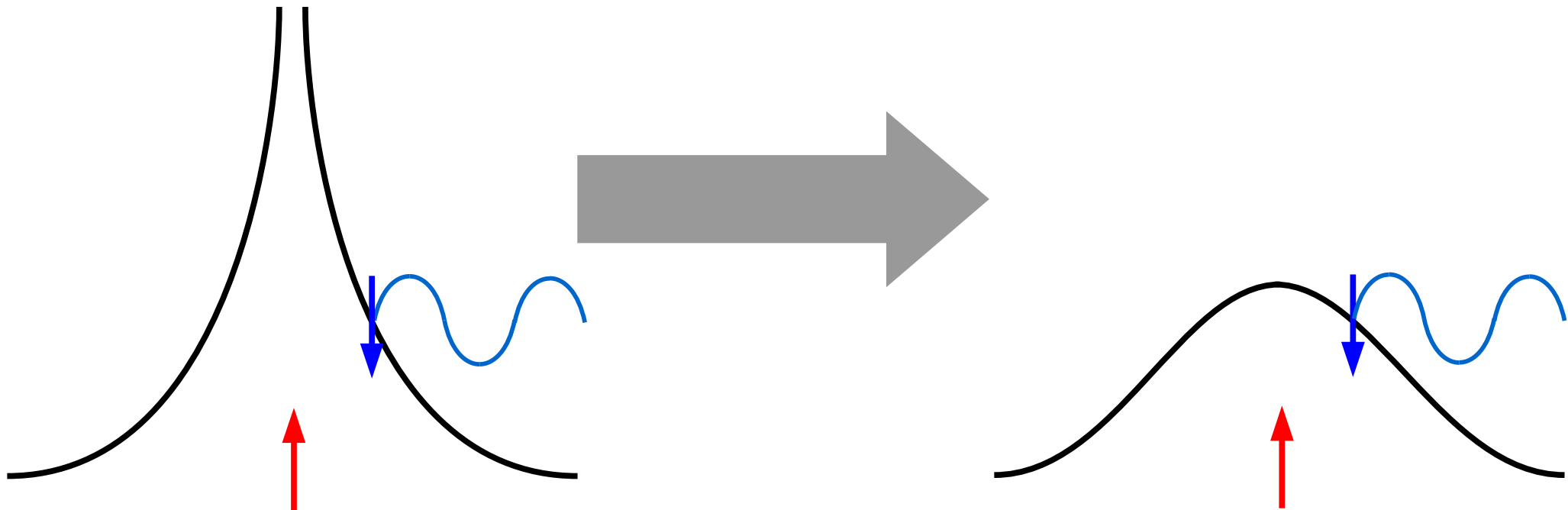
Fewer electrons



# Electron-electron pseudopotential

$$H = KE + V_{e-i} + V_{e-e}$$

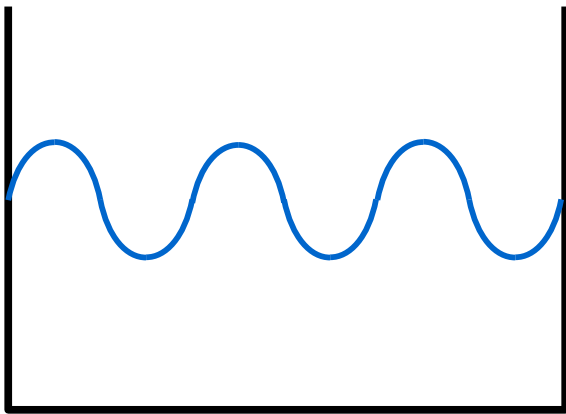
Smooth background



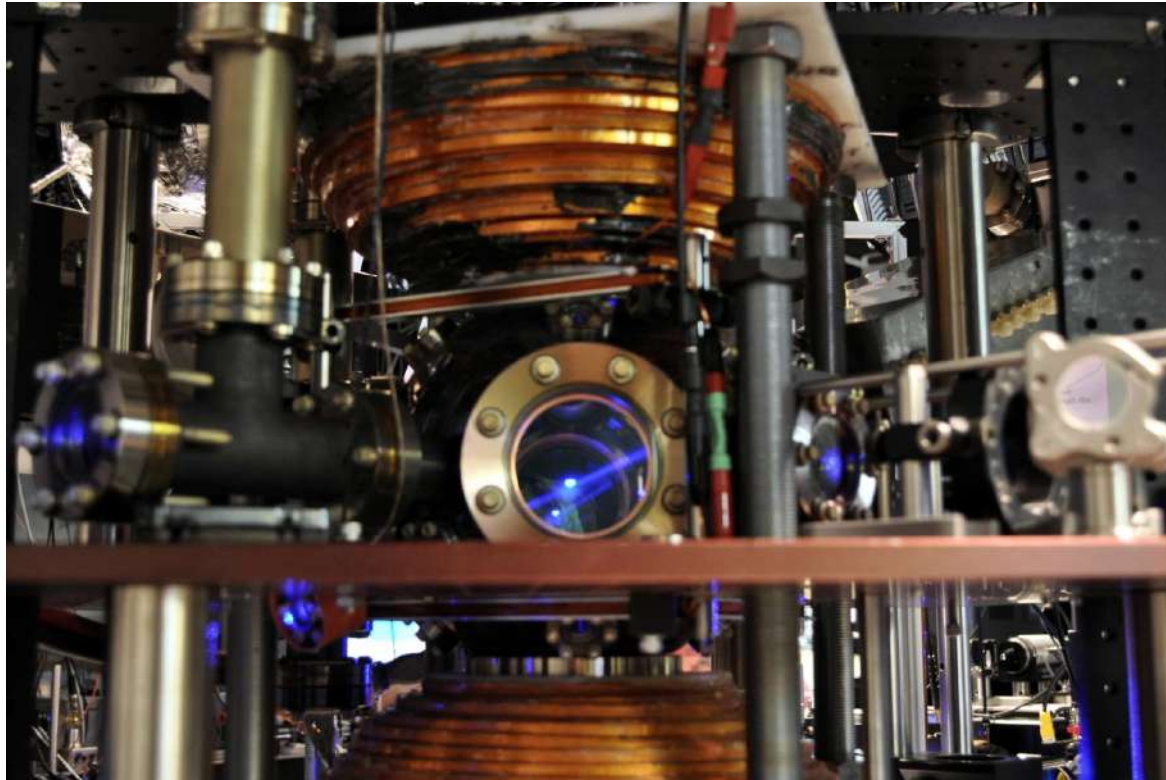
# Pseudo kinetic energy

$$H = \text{KE} + V_{e-i} + V_{e-e}$$

Smooth integrand



# Ultracold atom gases



Pristine

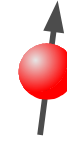
Contact interaction

High level of control

# Scattering in ultracold atom gases



$$|F = 1/2, m_F = 1/2\rangle$$

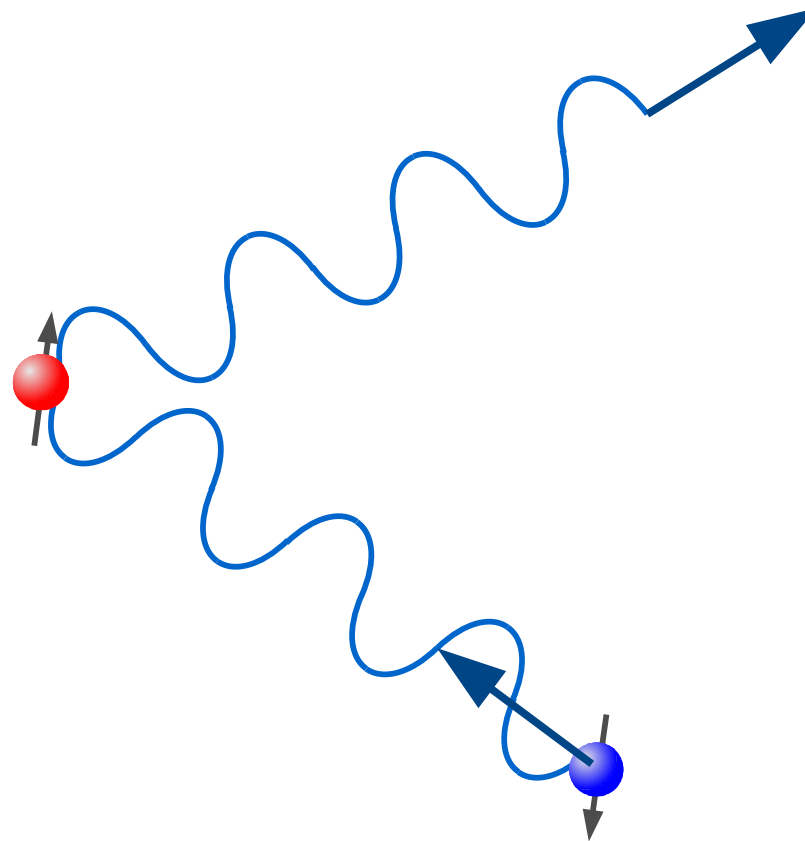


Up spin particle

$$|F = 1/2, m_F = -1/2\rangle$$

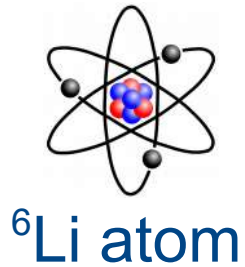


Down spin particle





# Scattering in ultracold atom gases



$$|F = 1/2, m_F = 1/2\rangle$$



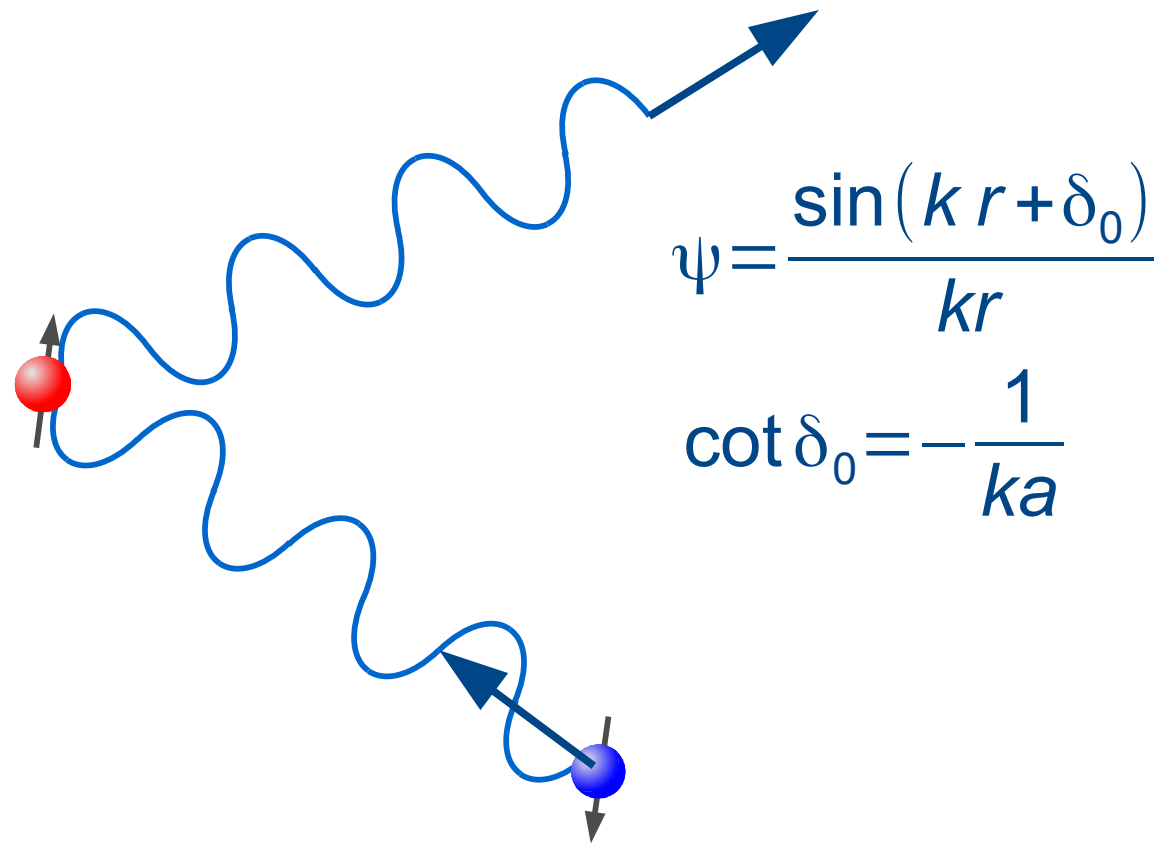
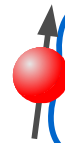
Up spin particle

$$|F = 1/2, m_F = -1/2\rangle$$



Down spin particle

$$V(r) = 4\pi a \delta(r) \frac{d}{dr} r$$



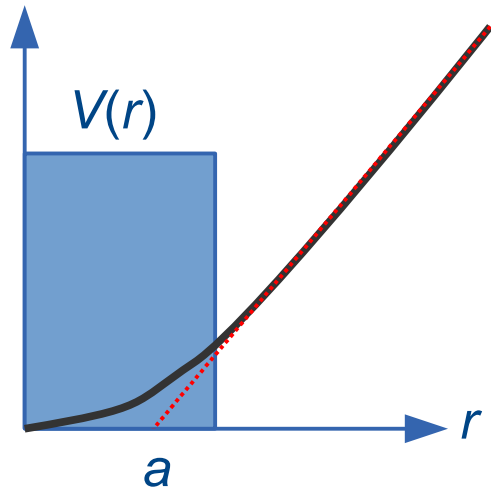
$$\psi = \frac{\sin(kr + \delta_0)}{kr}$$

$$\cot \delta_0 = -\frac{1}{ka}$$

# Scattering potentials

Underlying repulsive

Effective repulsive



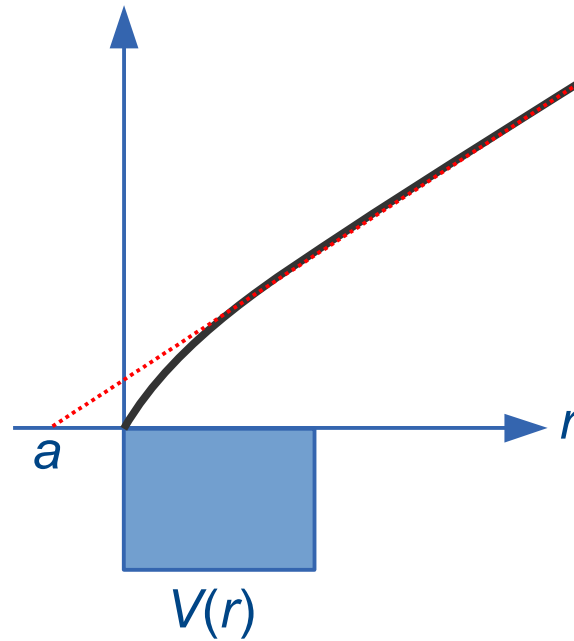
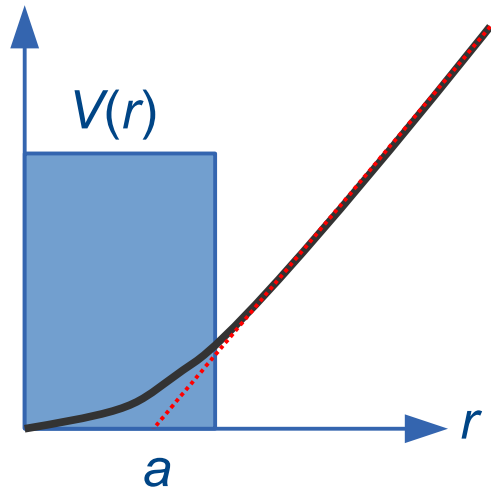
# Scattering potentials

Underlying repulsive

Effective repulsive

Underlying attractive

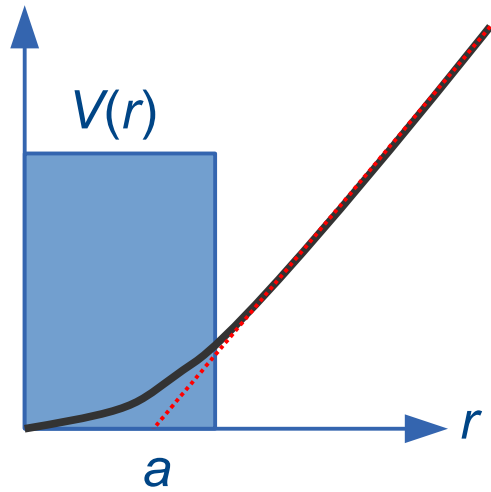
Effective attractive



# Scattering potentials

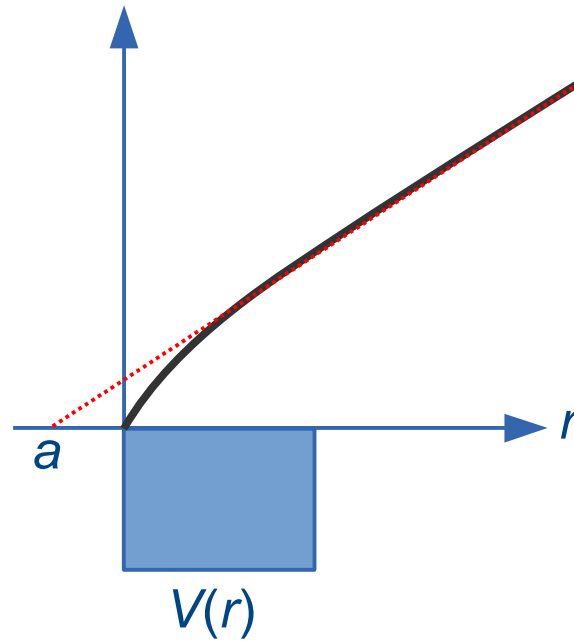
Underlying repulsive

Effective repulsive



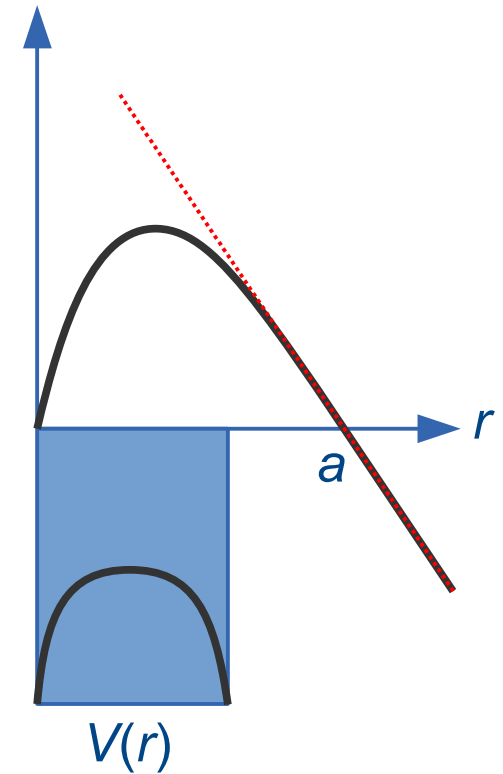
Underlying attractive

Effective attractive



Underlying attractive

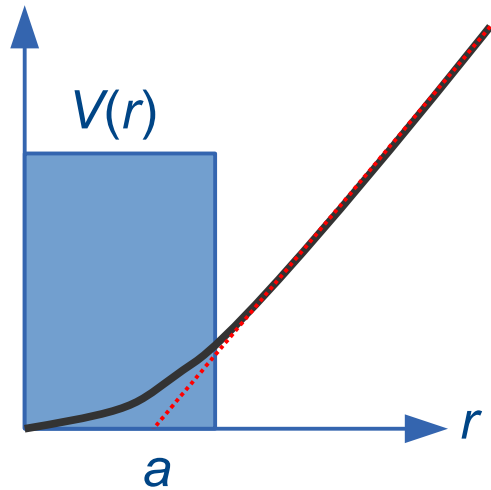
Effective repulsive



# Scattering potentials

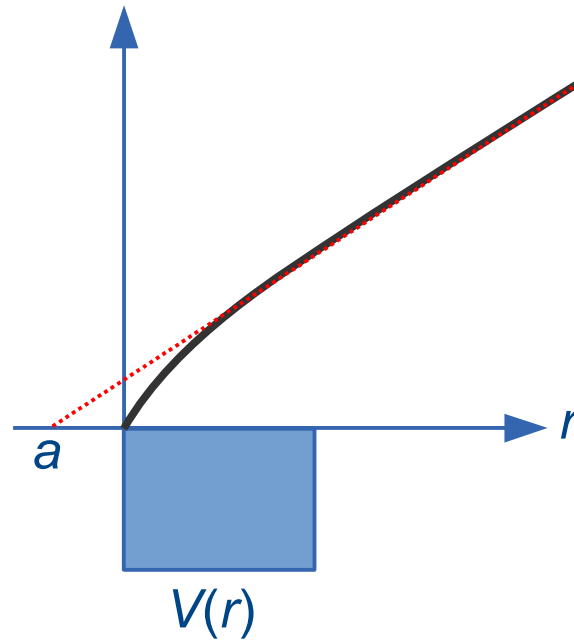
Underlying repulsive

Effective repulsive



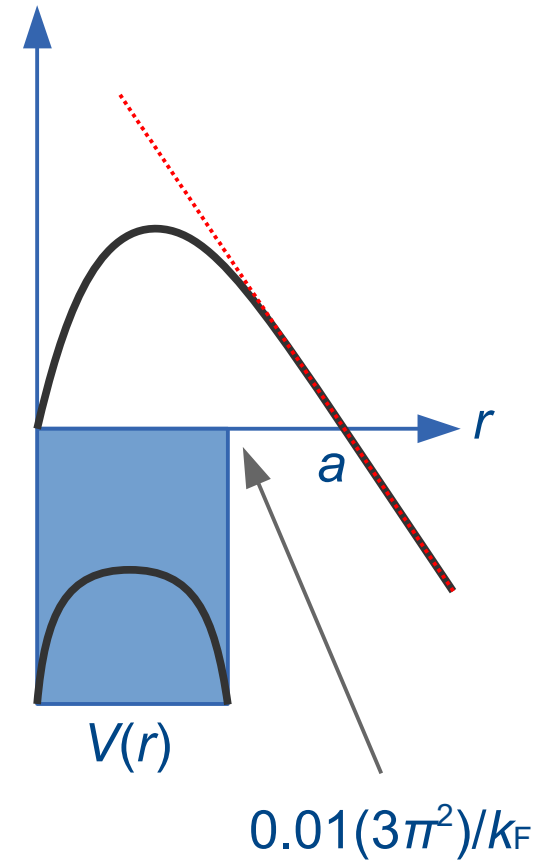
Underlying attractive

Effective attractive

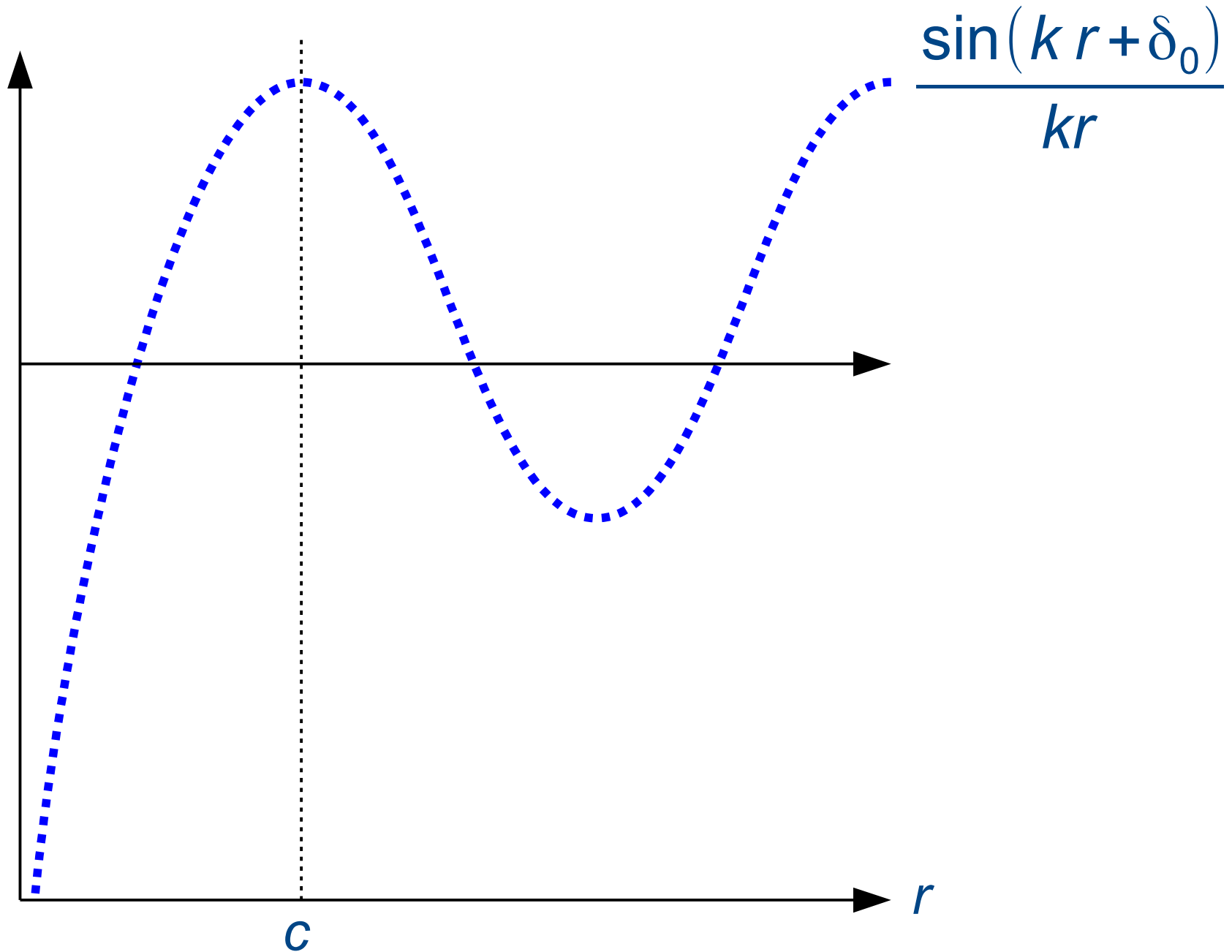


Underlying attractive

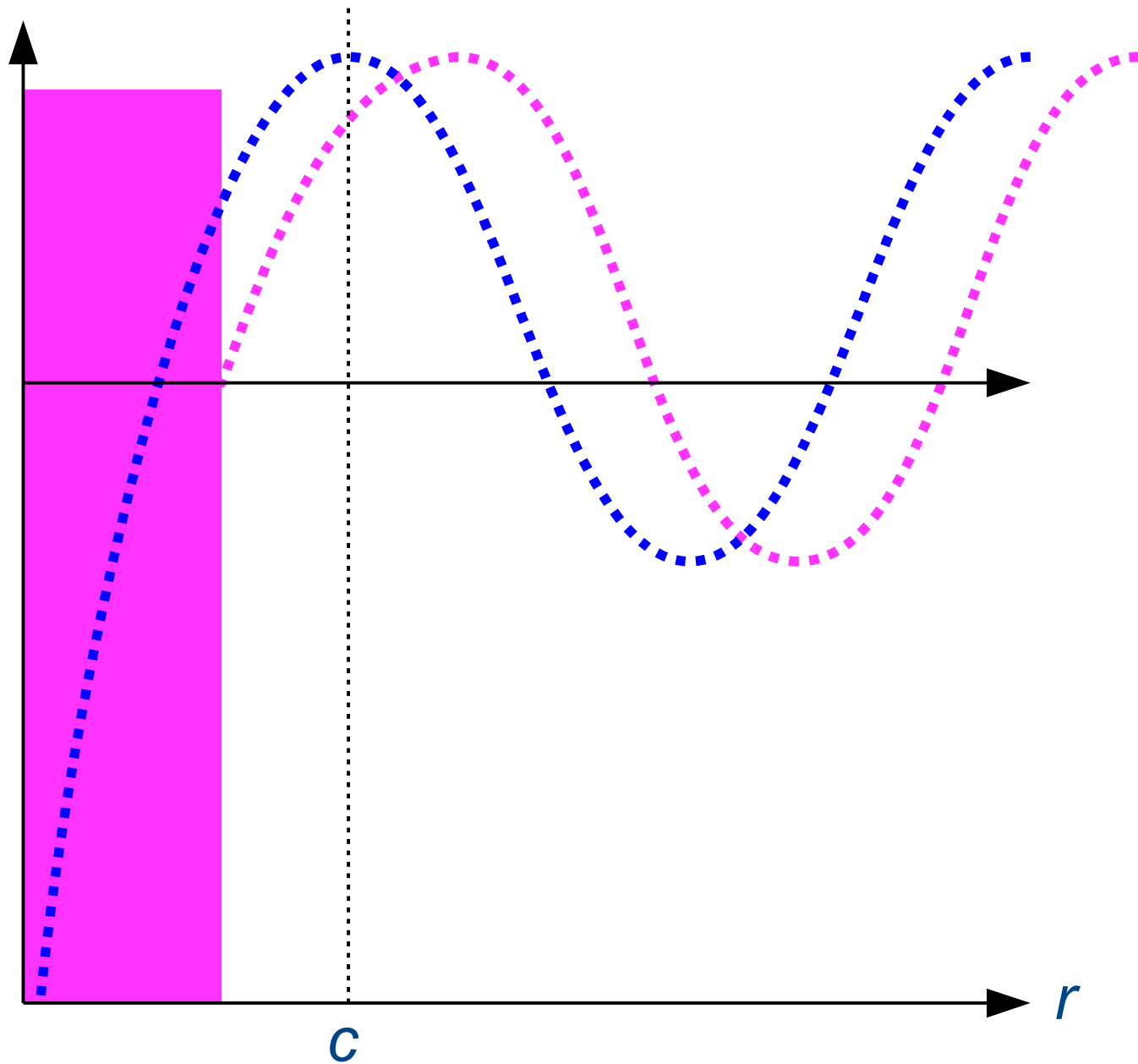
Effective repulsive



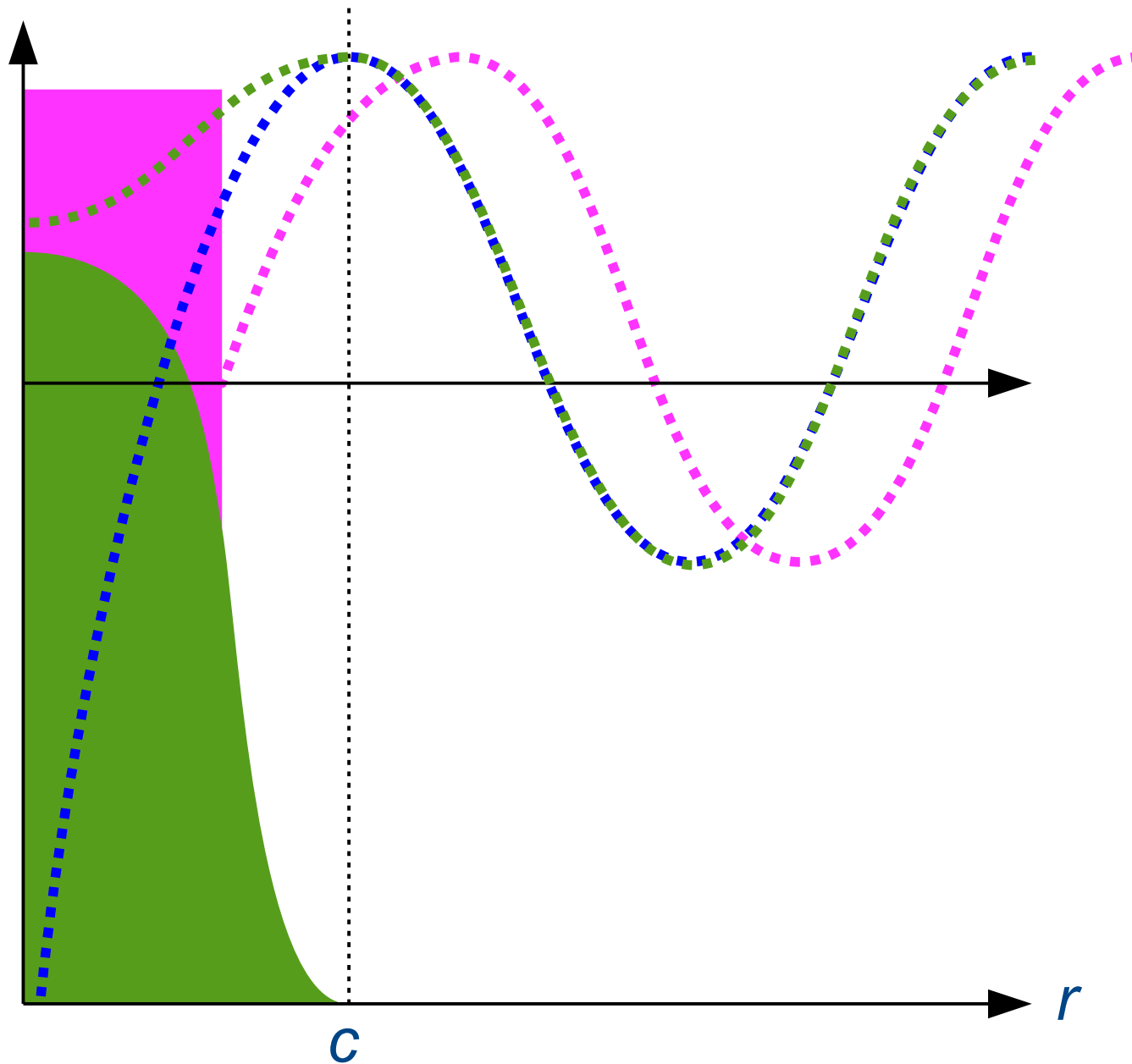
# Construction of a pseudopotential



# Construction of a pseudopotential



# Construction of a pseudopotential



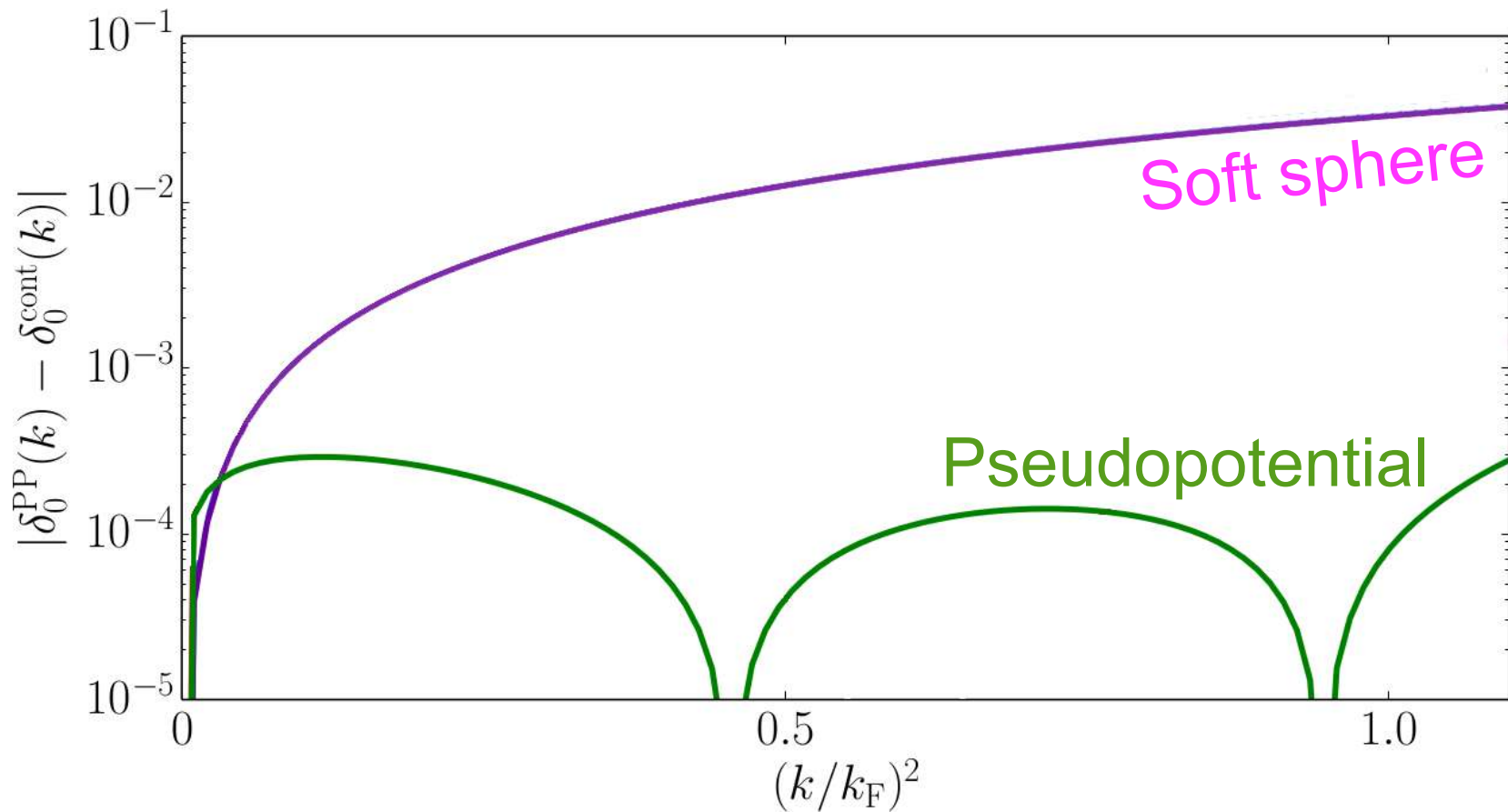
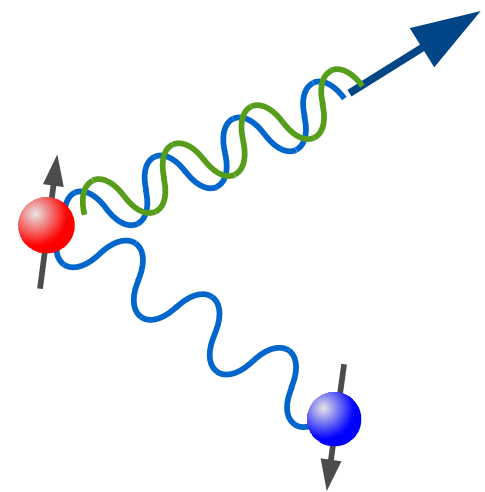


# Construction of a pseudopotential

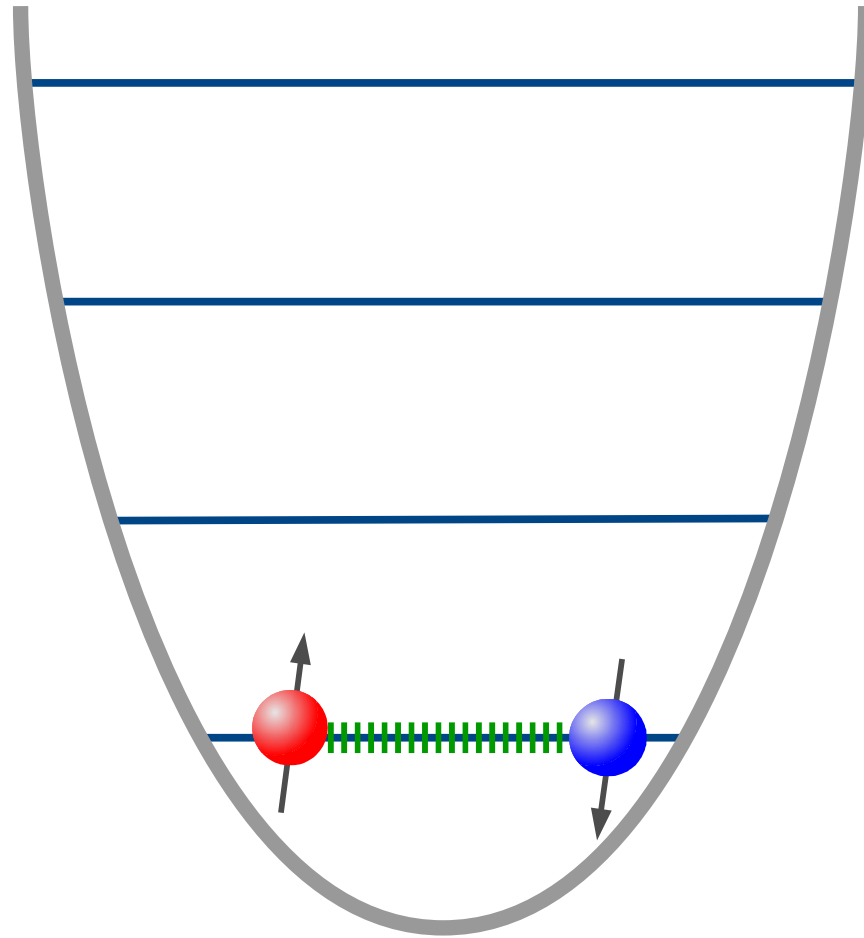
$$V_{\text{PP}}(r) = \begin{cases} \left(1 - \frac{r}{c}\right)^2 \left[ v_1 \left(\frac{1}{2} + \frac{r}{c}\right) + \sum_{i=2}^{N_v} v_i \left(\frac{r}{c}\right)^i \right] & r < c \\ 0 & r > c \end{cases}$$

$$\sum_{l=0}^{l_{\max}} \int_0^{k_F} [\delta_{0,\text{PP}} - \delta_{0,\text{cont}}]^2 dk$$

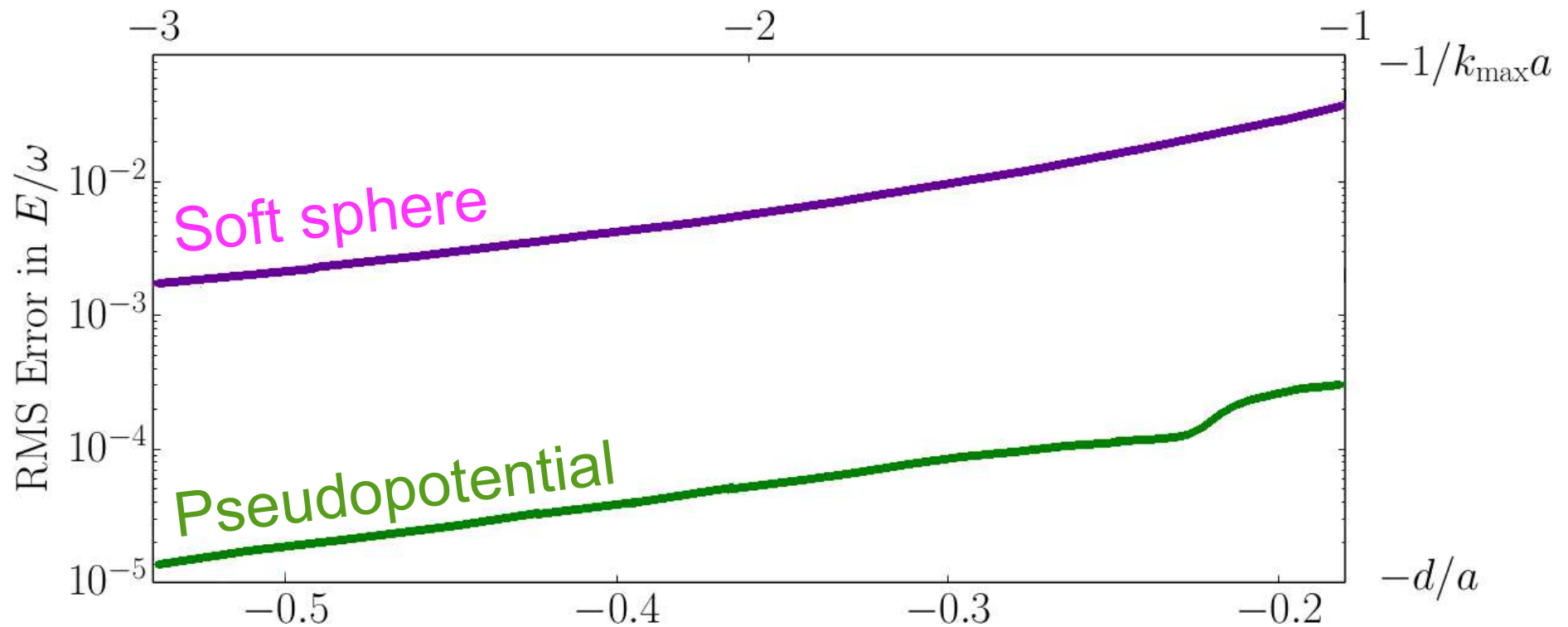
# Pseudopotential: scattering properties



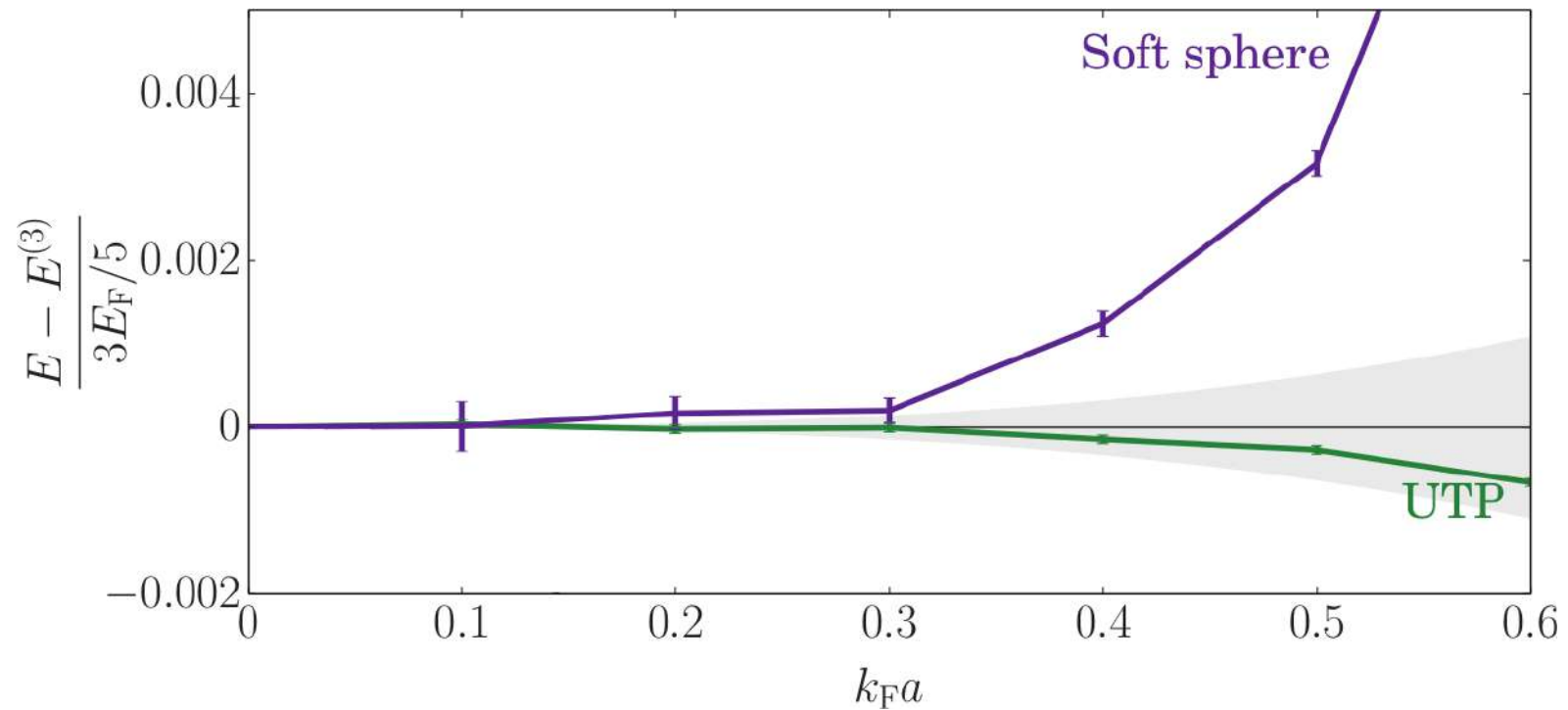
# Pseudopotential: two atoms in a trap



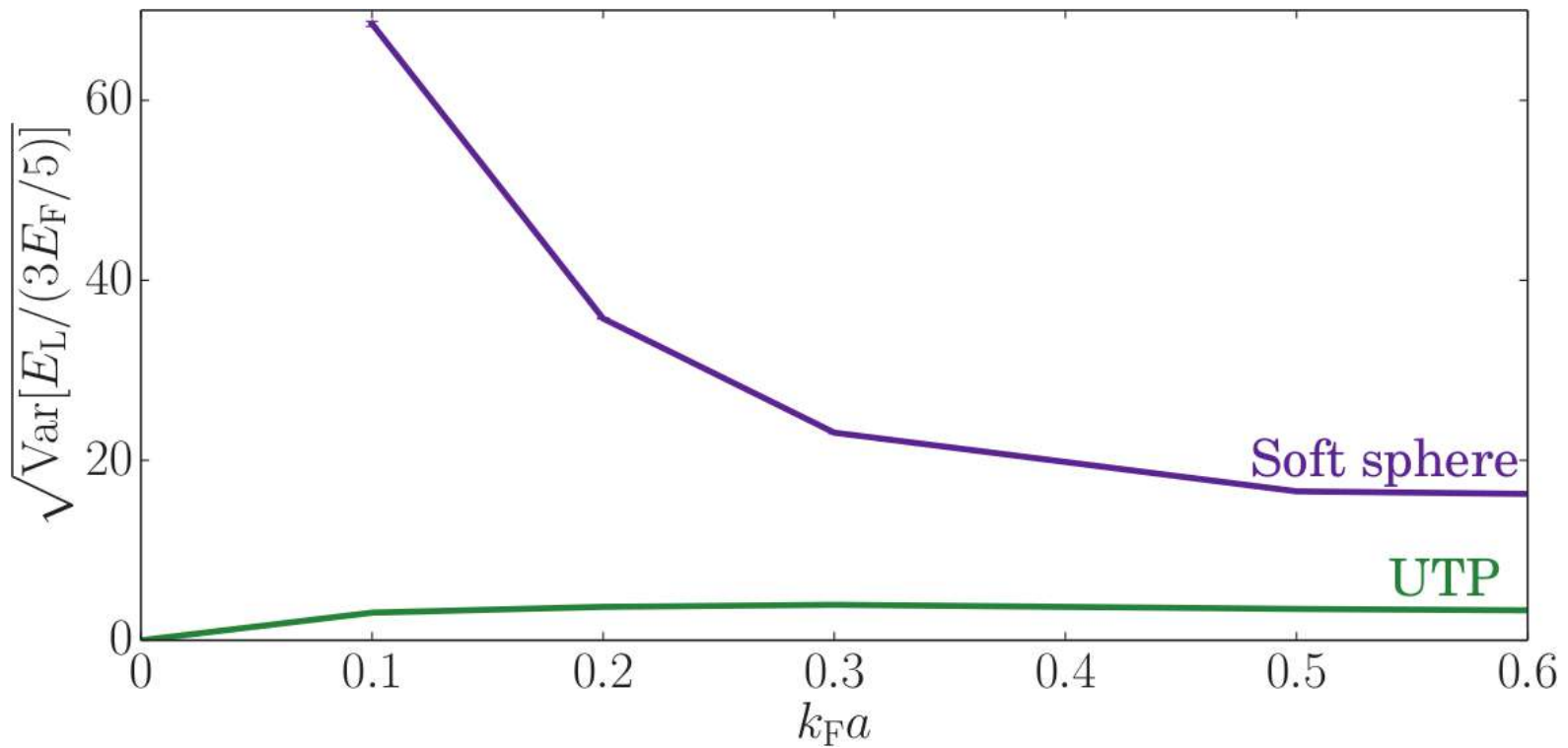
# Pseudopotential: two atoms in a trap



# Pseudopotential: homogeneous system

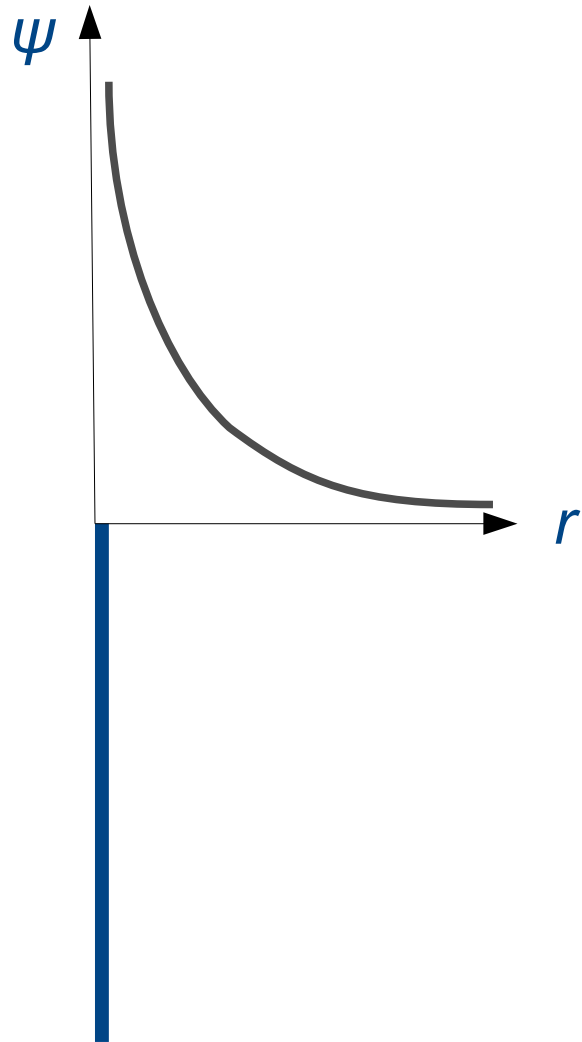


# Pseudopotential: homogeneous system

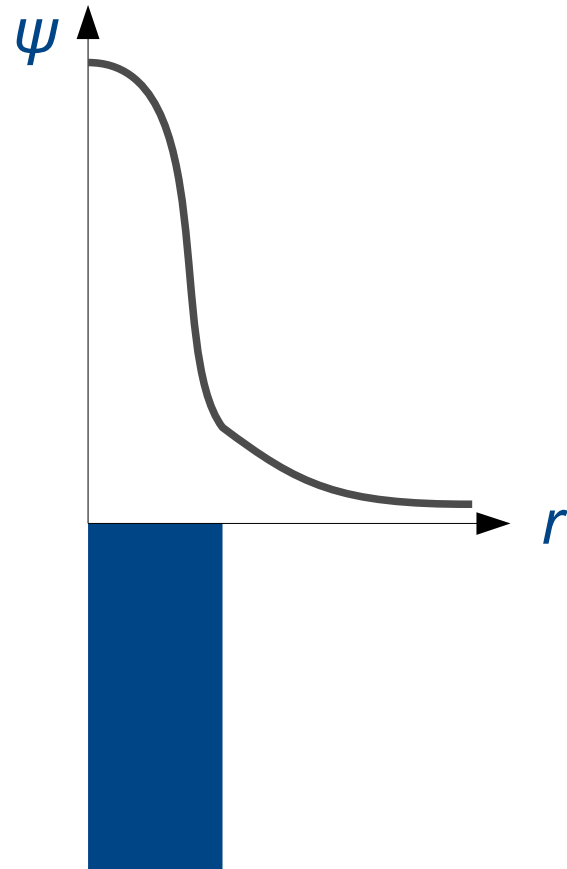


# Pseudopotential: bound state

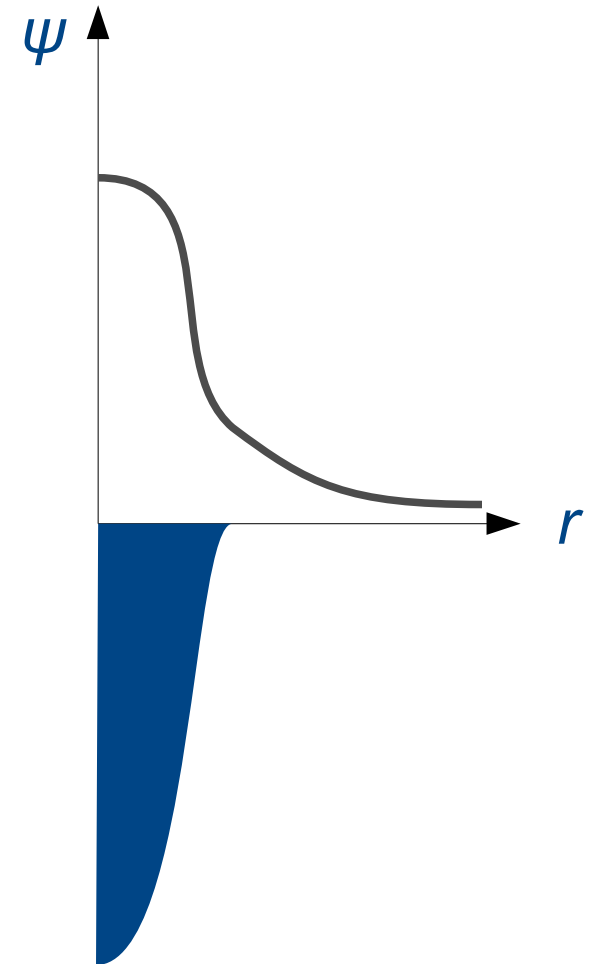
Contact



Square well

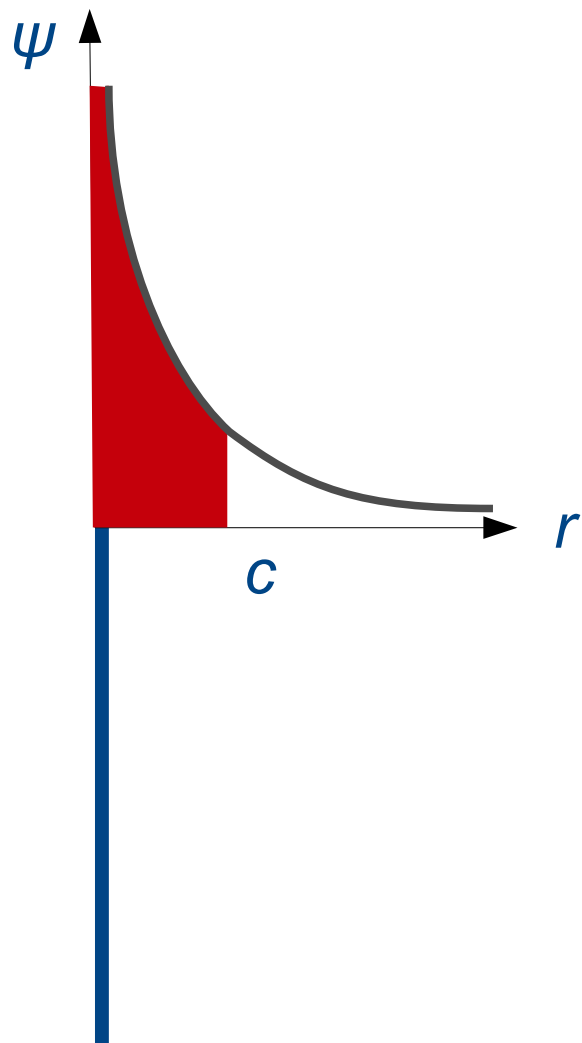


Pseudopotential

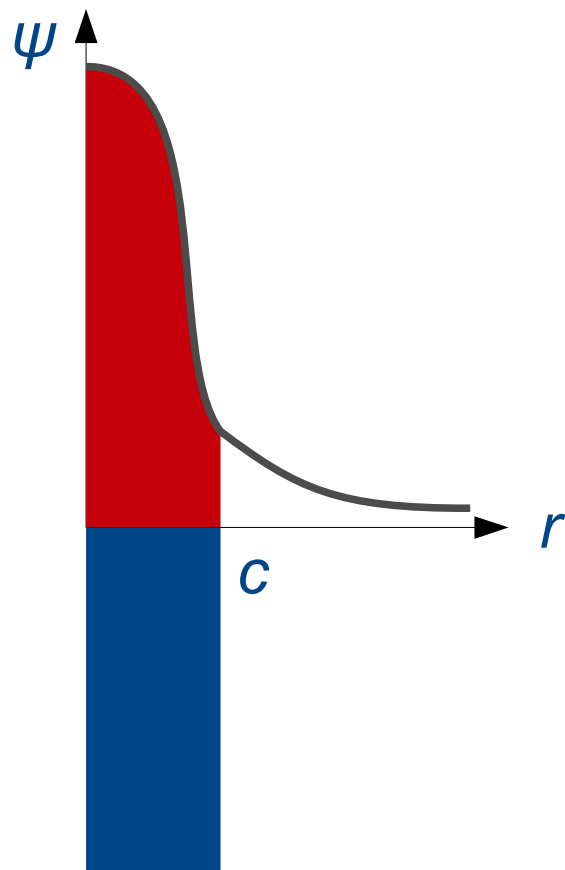


# Pseudopotential: bound state

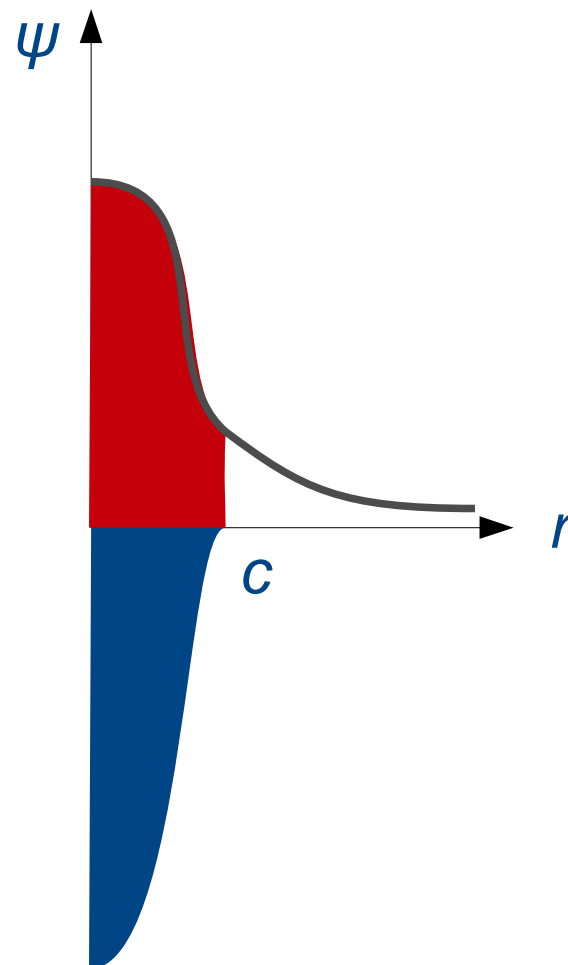
Contact



Square well



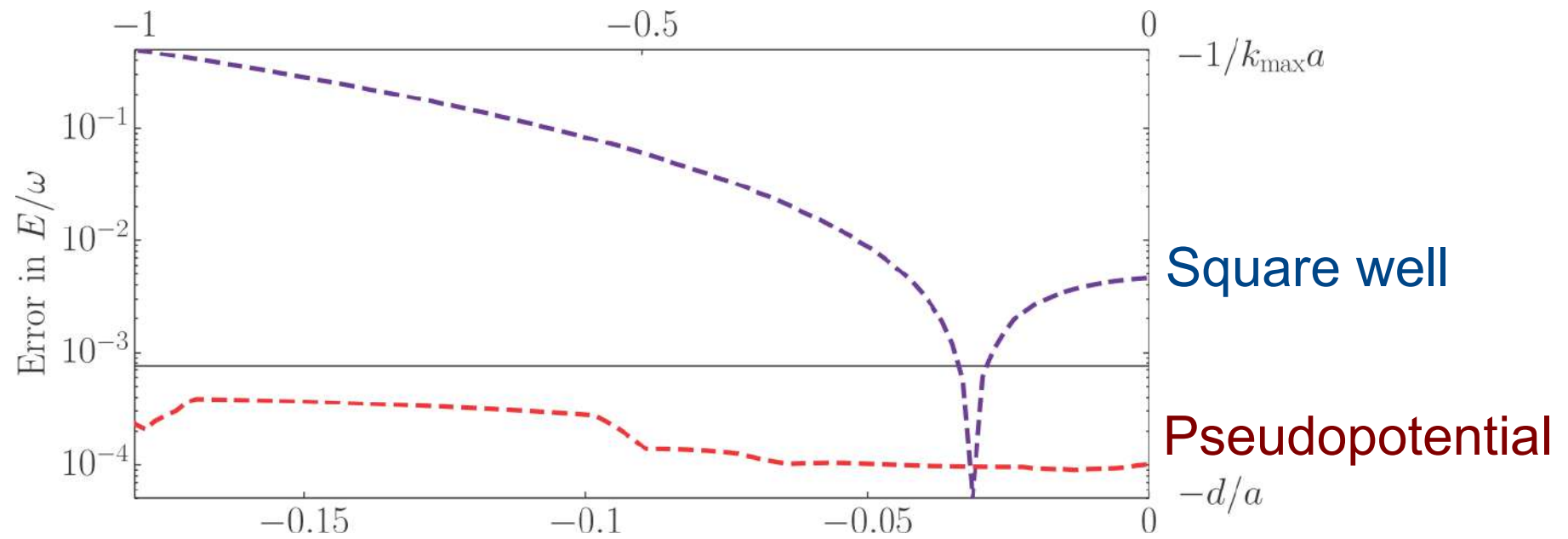
Pseudopotential



$$E = \int_0^c |\psi|^2 dr \quad \text{Troullier \& Martins PRB 43 1993 (1991)}$$



# Pseudopotential: bound state



# Contact interaction pseudopotential summary

Pseudopotential for repulsive, attractive, and bound state: 100 times more accurate, 1000 times faster

Formalism is systematically improvable, trading accuracy for speedup

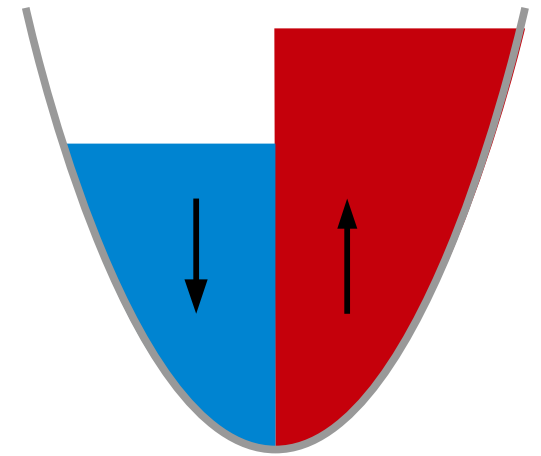
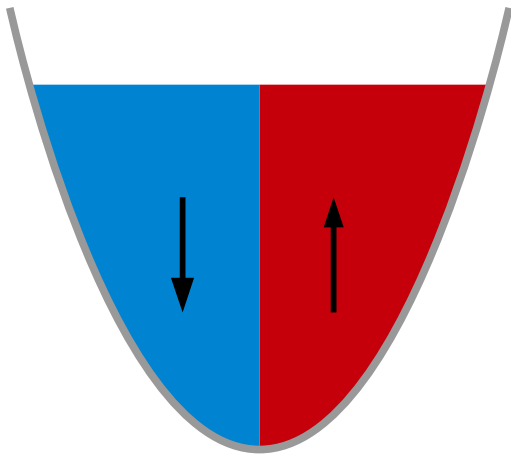
Python tool: <https://pypi.python.org/pypi/contactpp>

# Stoner Hamiltonian

$$H = -\frac{\nabla^2}{2} + 4\pi a \delta(\mathbf{r}_\uparrow - \mathbf{r}_\downarrow)$$

# Stoner Hamiltonian

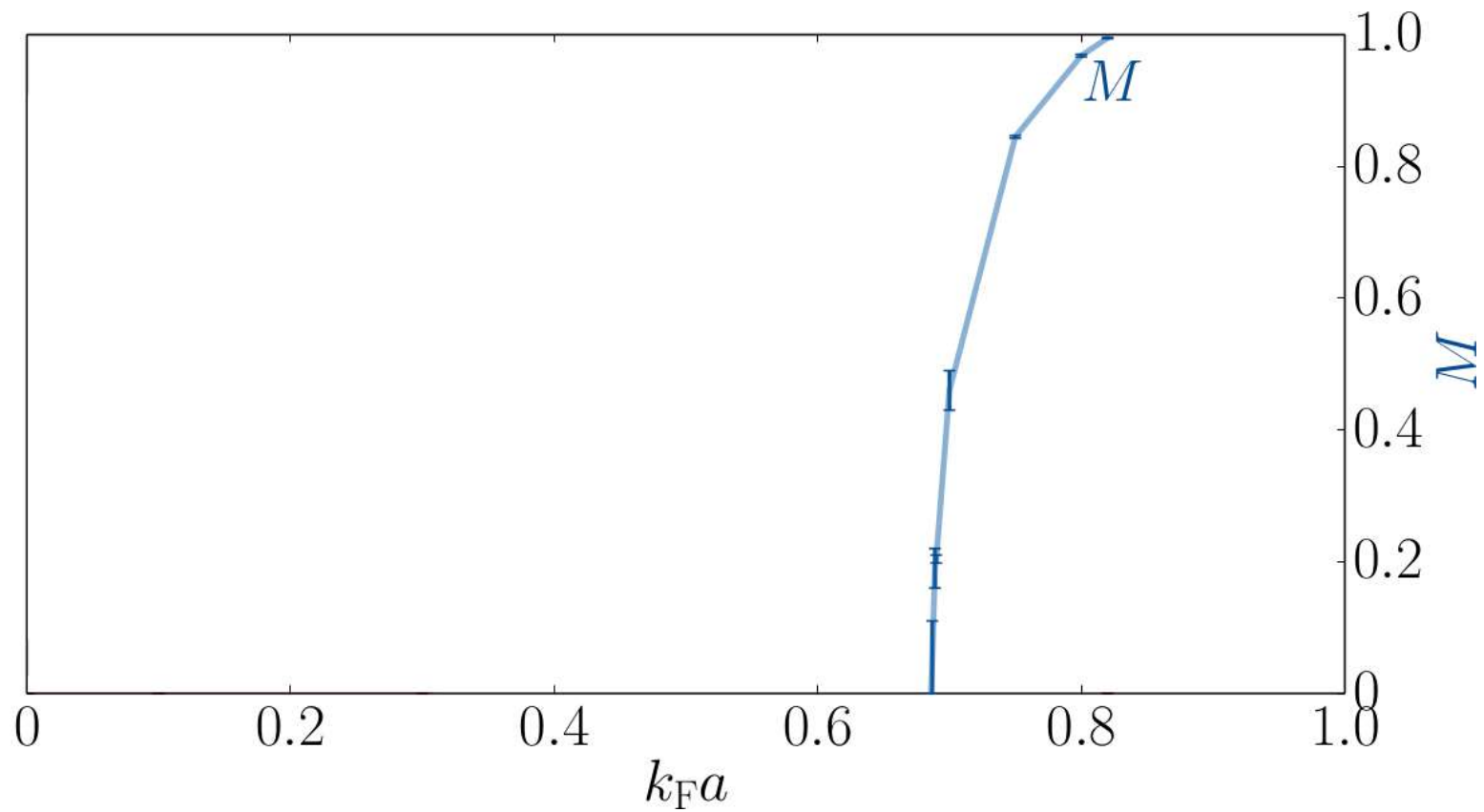
$$E = KE + gn_{\uparrow}n_{\downarrow}$$



# Theories of ferromagnetism

Stoner mean-field theory	Second order	$k_{Fa}=1.57$
Fluctuations beyond Hertz-Millis	First order	-
Polaron theory	First order	-
Field theory	First order	$k_{Fa}=1.054$
Tan relations	No magnetism	-
DMC hard sphere	First order	$k_{Fa}=0.81(2)$
QMC square well	First / second order	$k_{Fa}=0.83(2)$

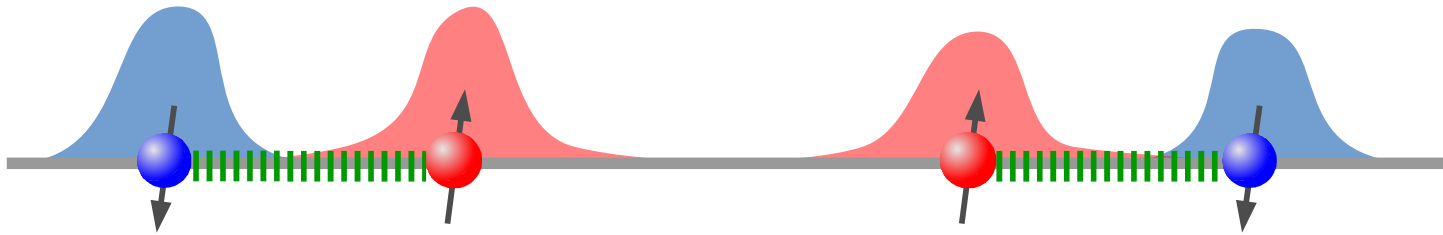
# Stoner Hamiltonian



# Theories of ferromagnetism

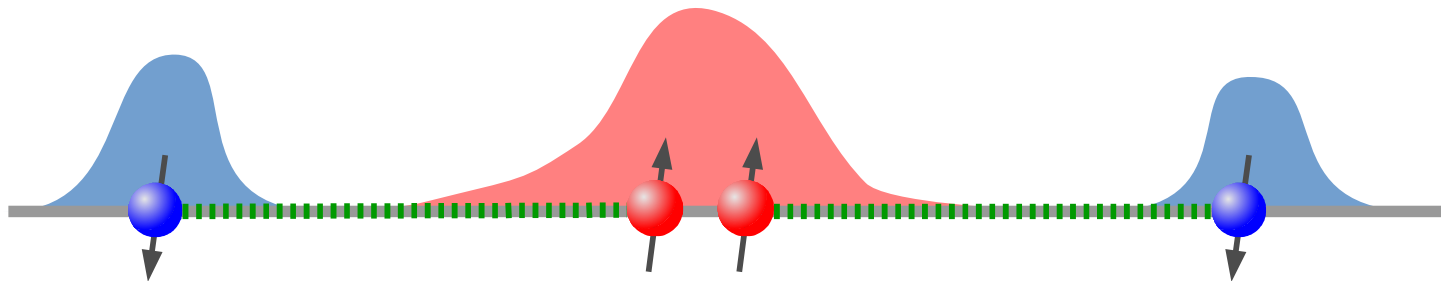
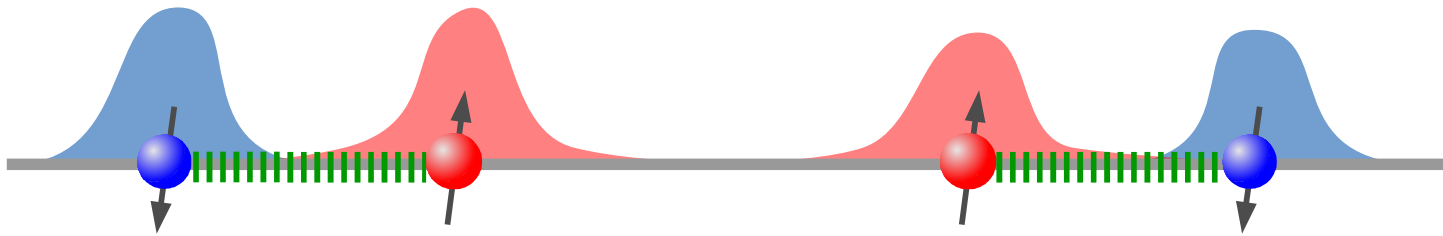
Stoner mean-field theory	Second order	$k_{Fa}=1.57$
Fluctuations beyond Hertz-Millis	First order	-
Polaron theory	First order	-
Field theory	First order	$k_{Fa}=1.054$
Tan relations	No magnetism	-
DMC hard sphere	First order	$k_{Fa}=0.81(2)$
QMC square well	First / second order	$k_{Fa}=0.83(2)$
DMC pseudopotential	Second order	$k_{Fa}=0.683(1)$

# Fluctuation contributions

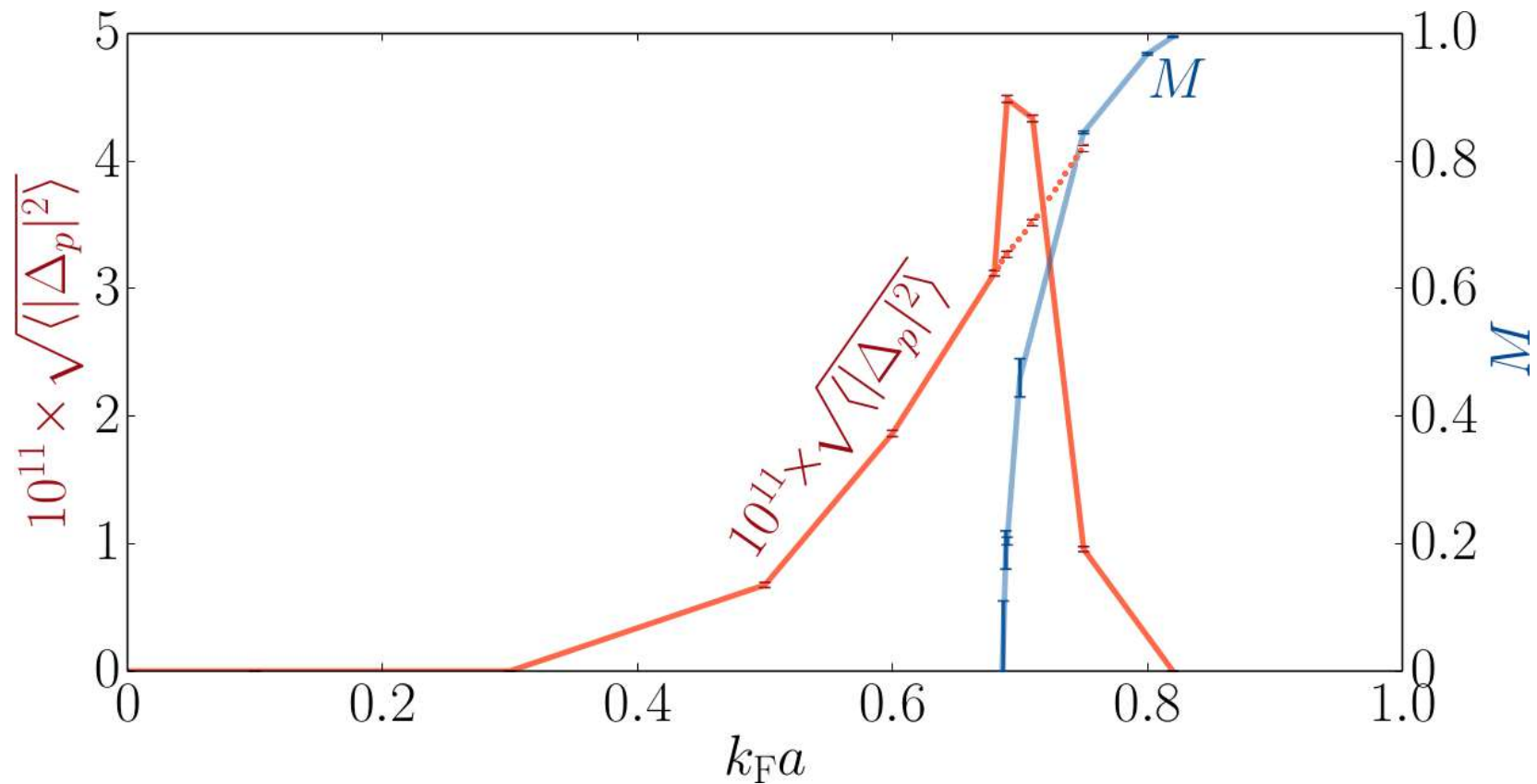




# Fluctuation contributions



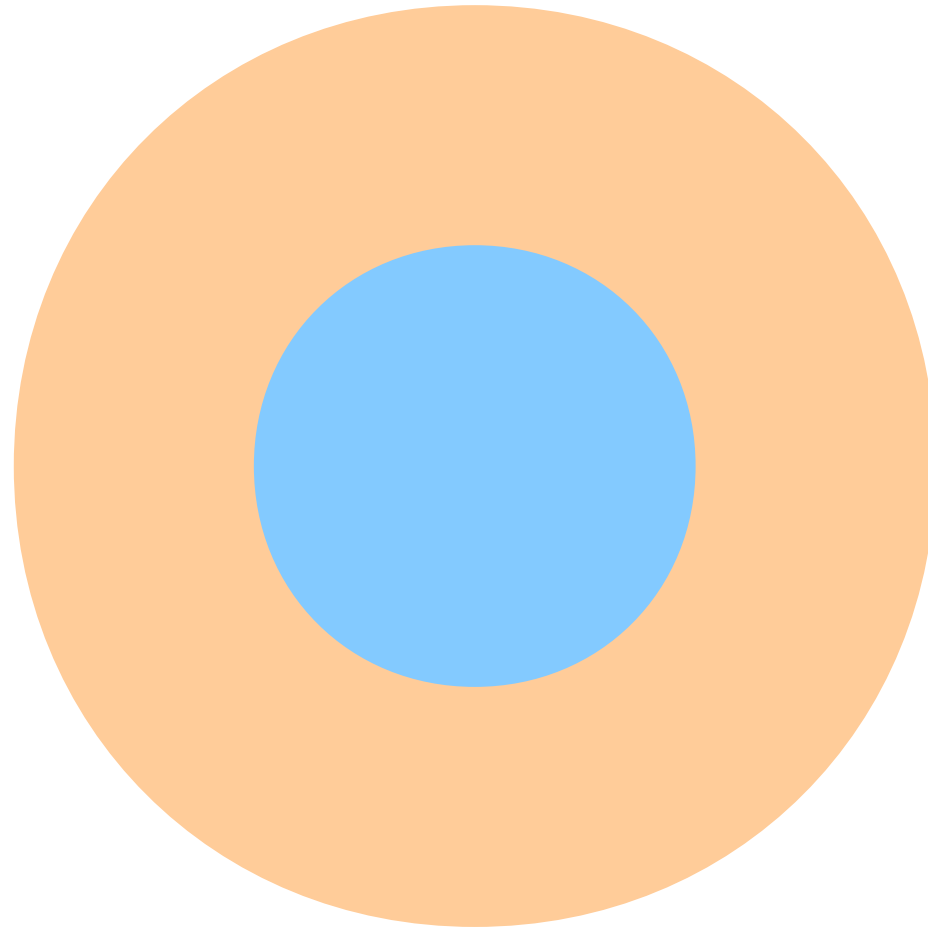
# Stoner Hamiltonian



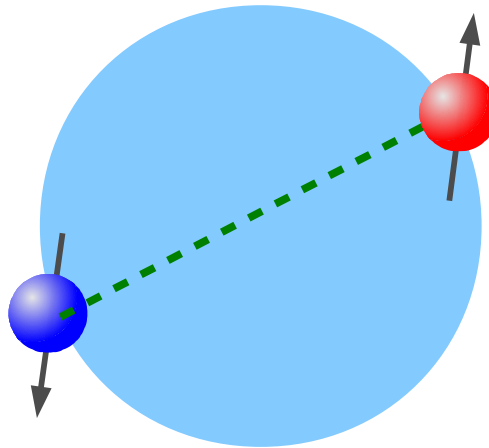
# Finite ranged interactions

$$\cot \delta_0 = -\frac{1}{ka} + \frac{1}{2} k r_{eff}$$

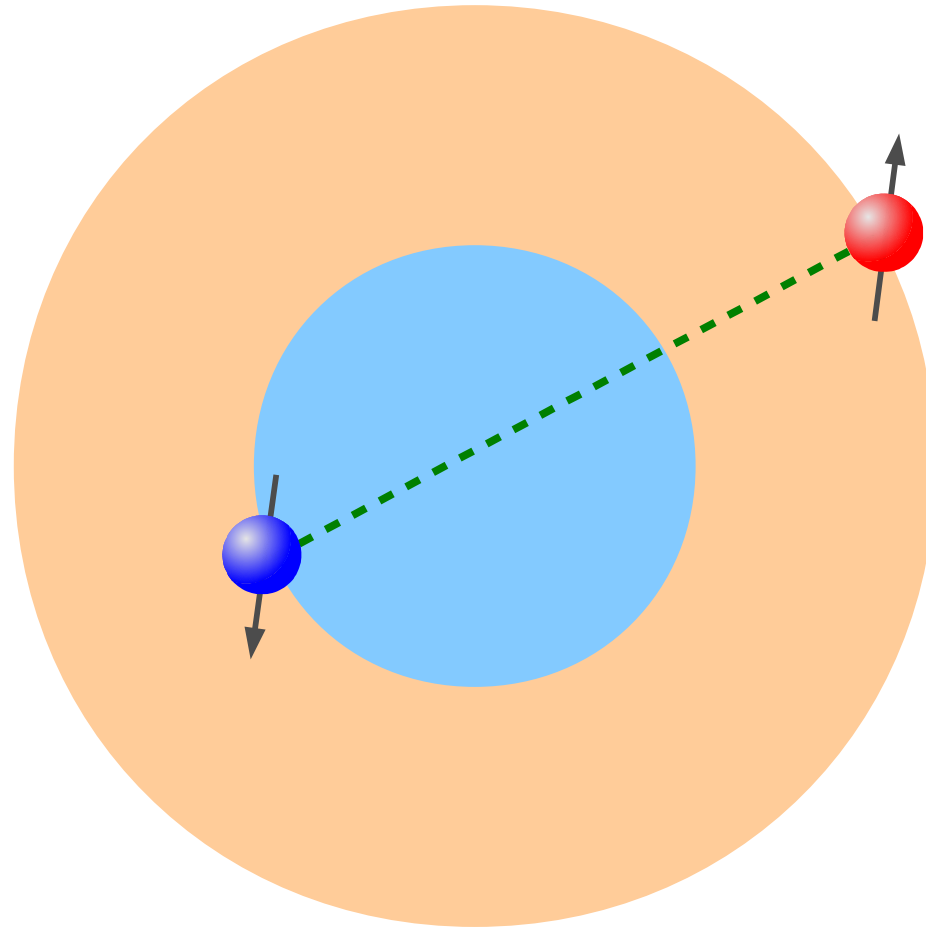
# Imbalanced superfluid



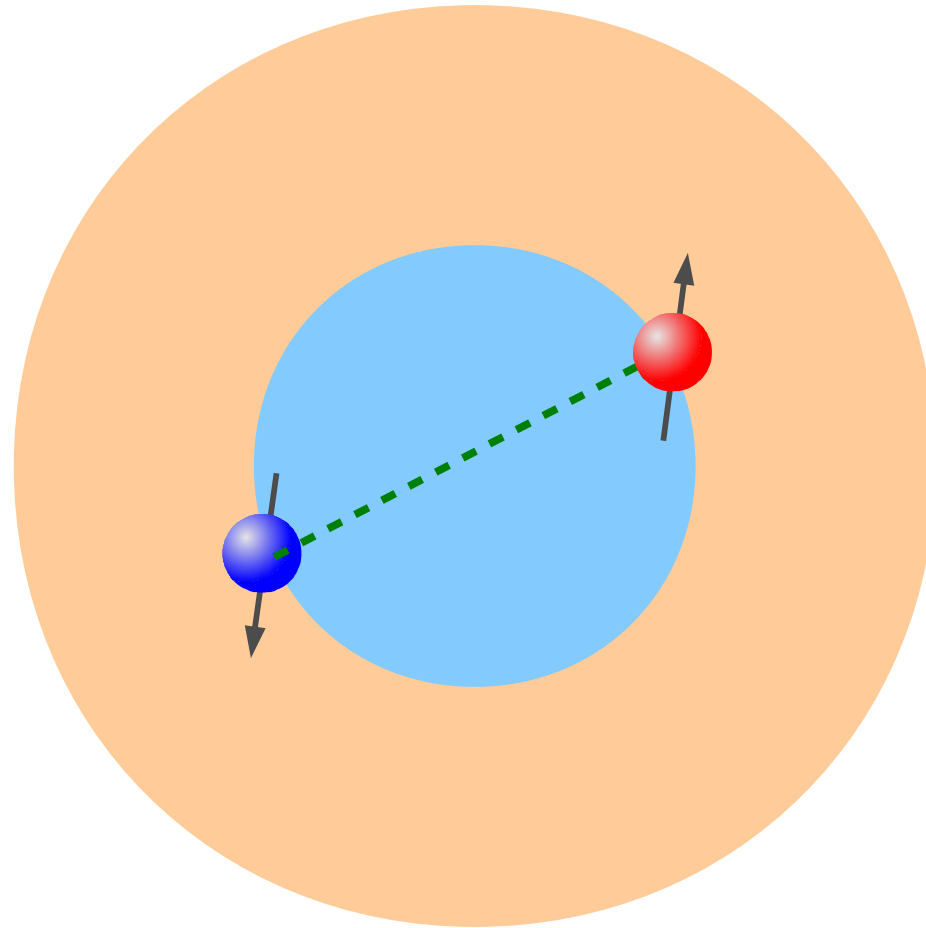
# BCS theory



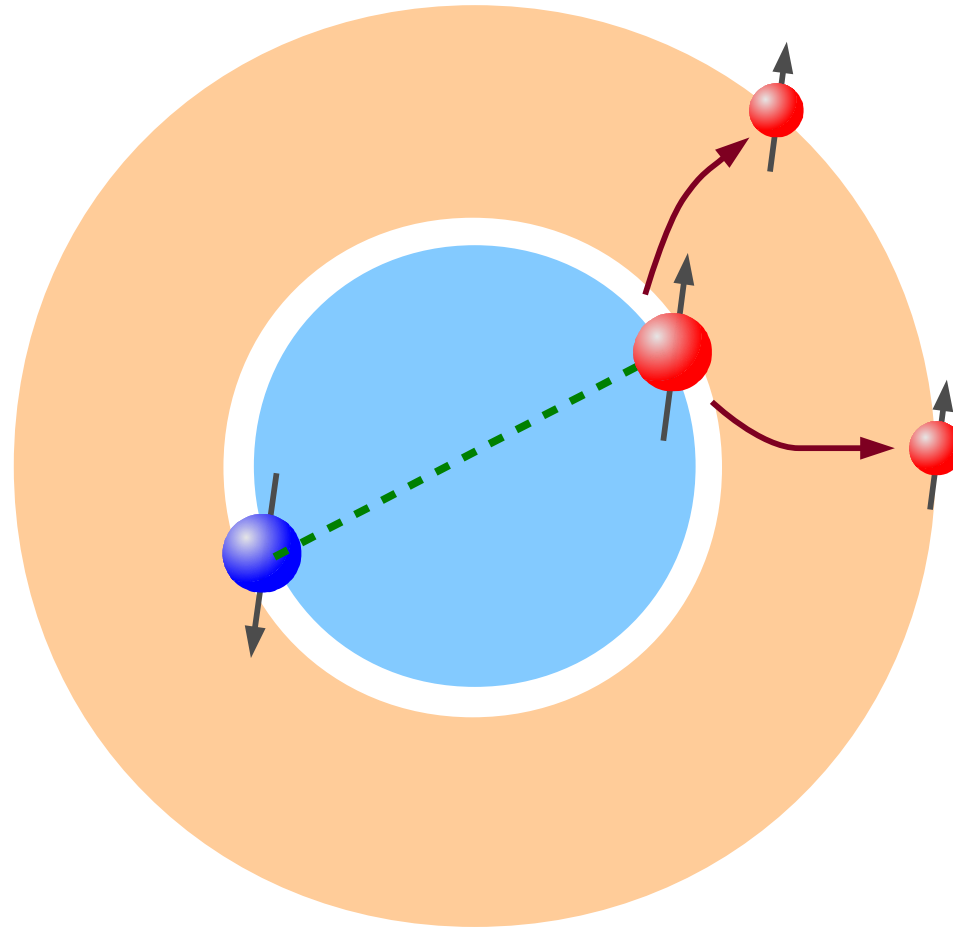
# FFLO superfluid



# Breached superfluid

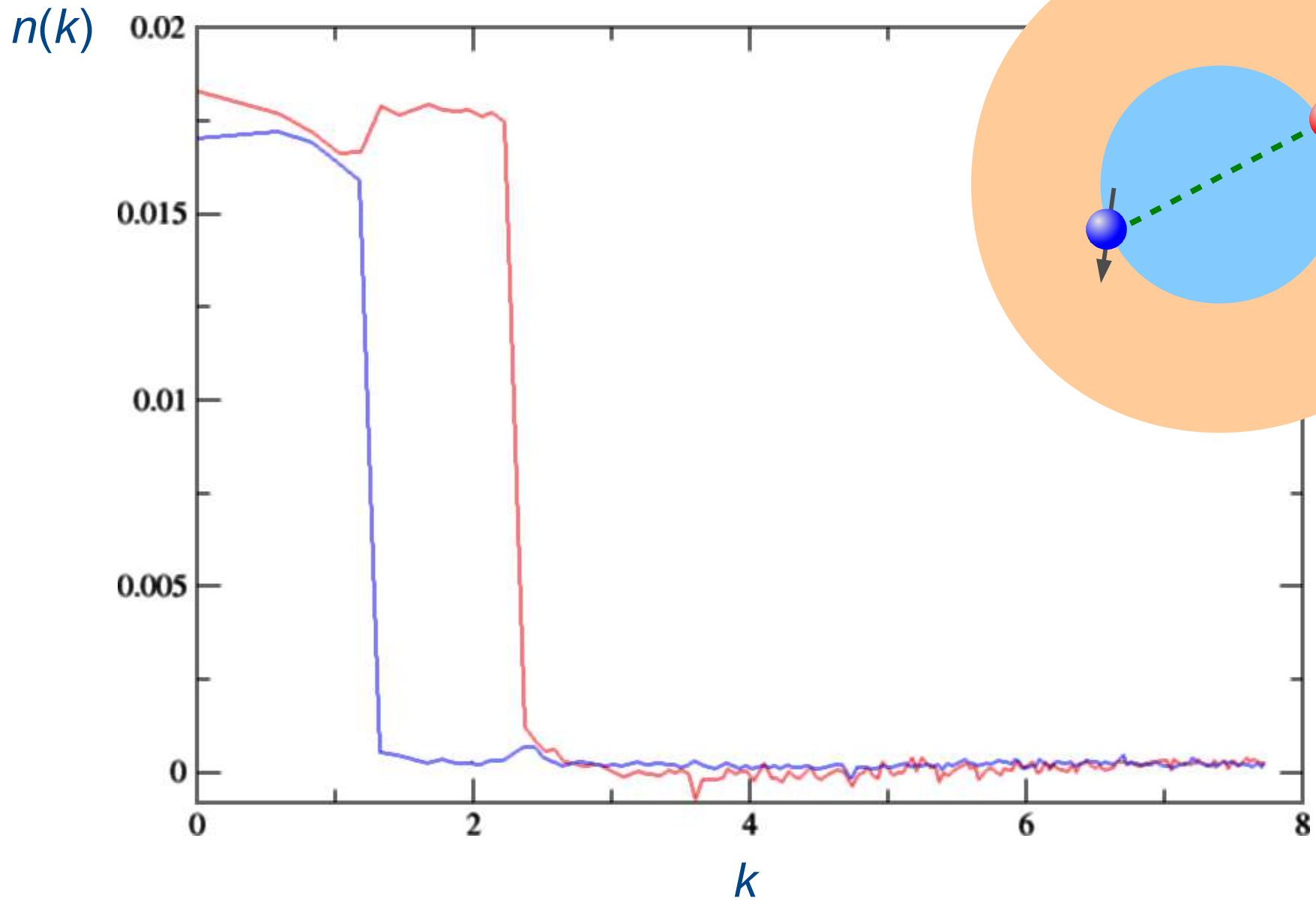


# Breached superfluid

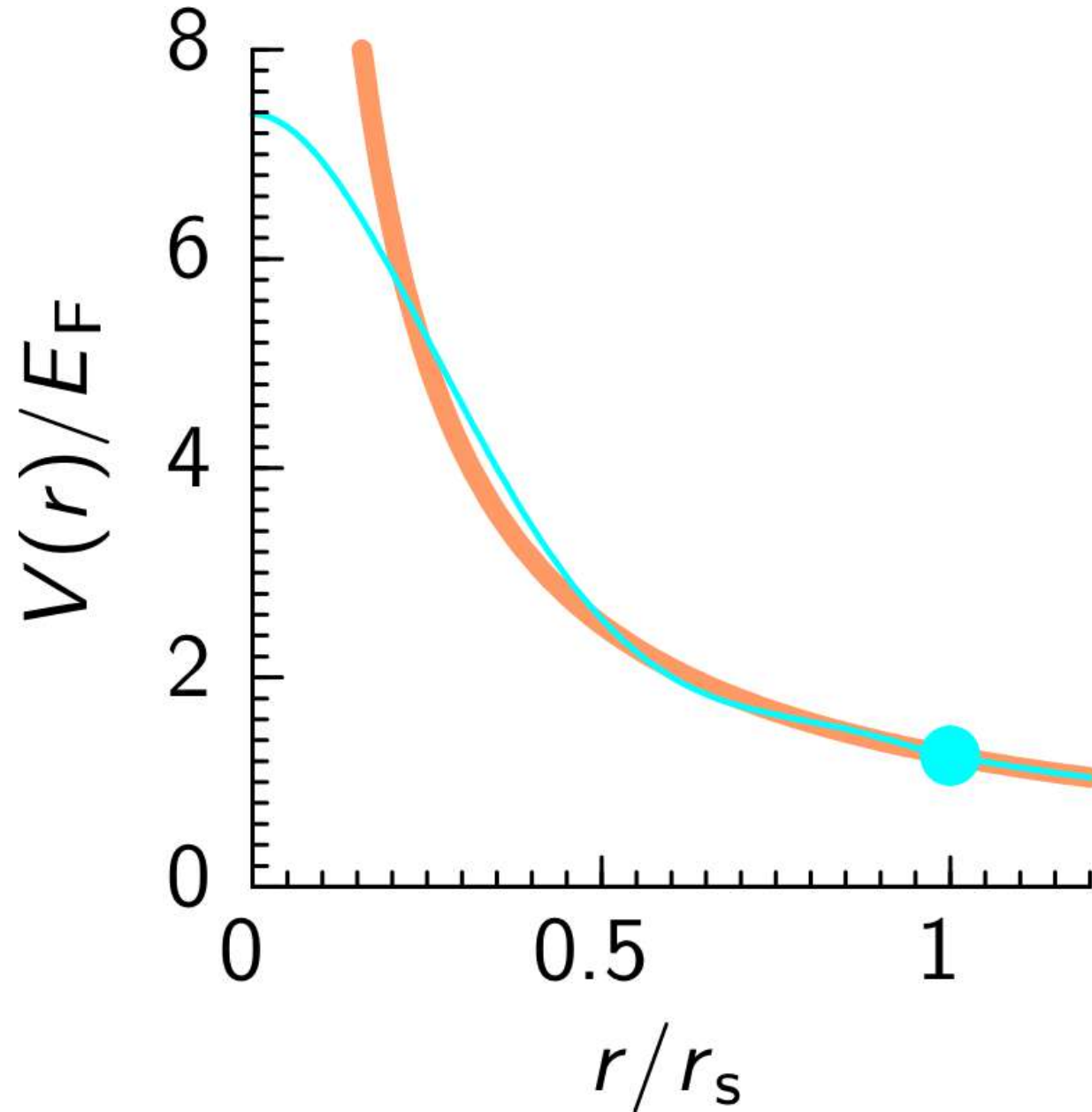




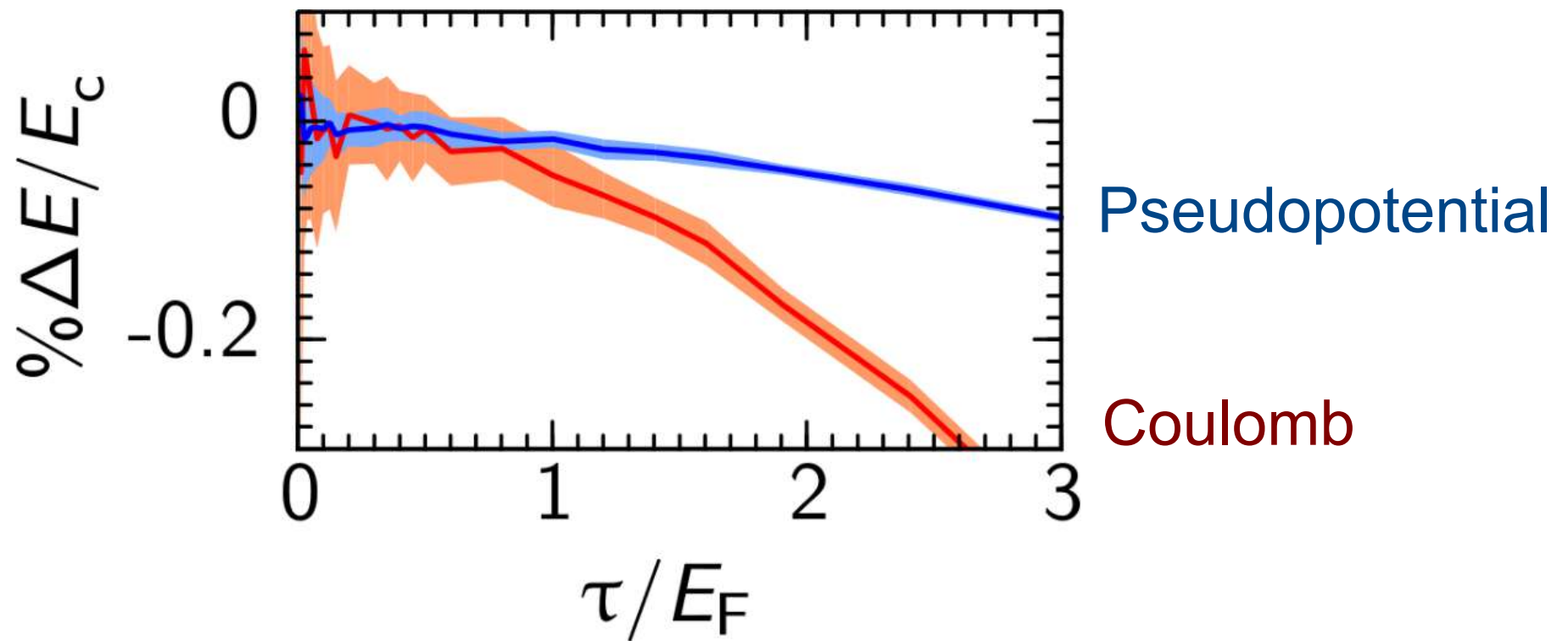
# Breached superfluid



# Coulomb pseudopotential

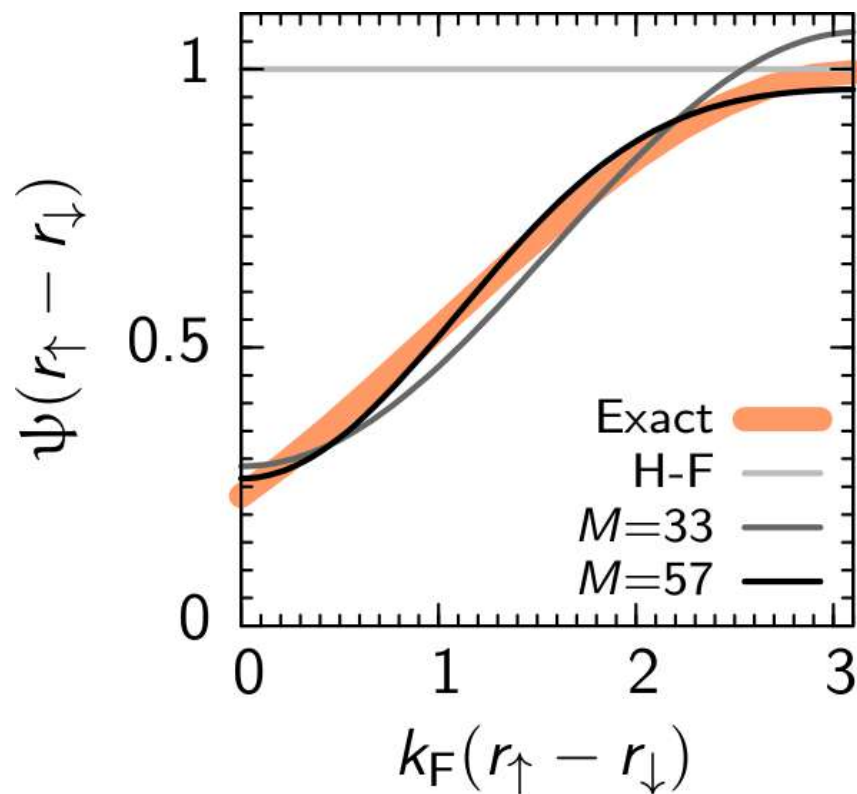


# Homogeneous electron gas

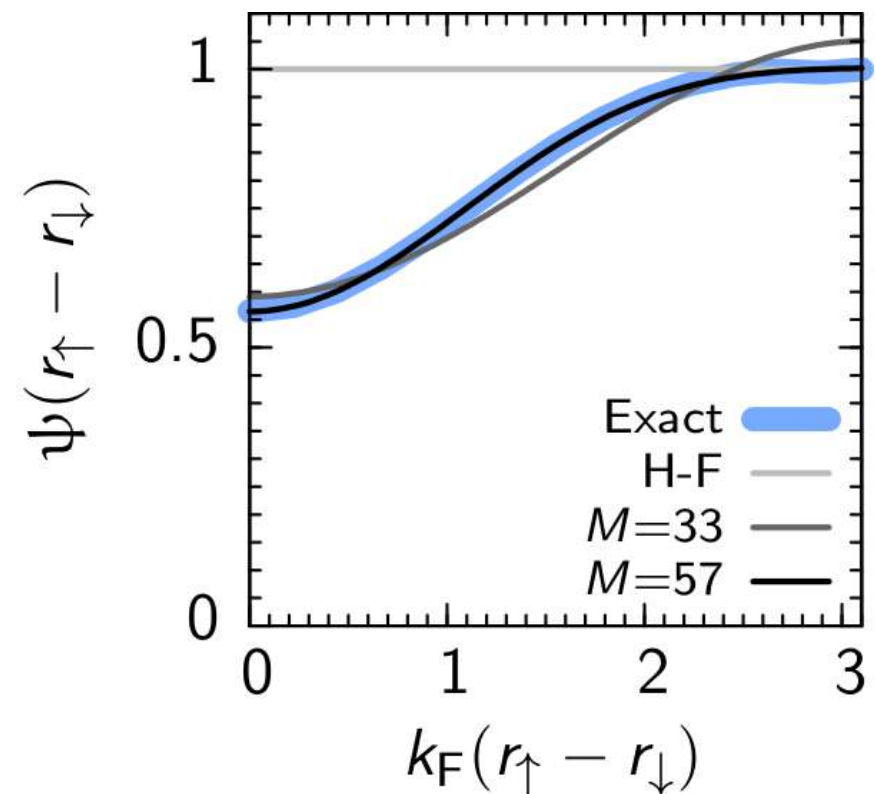


# Configuration interaction

## Coulomb

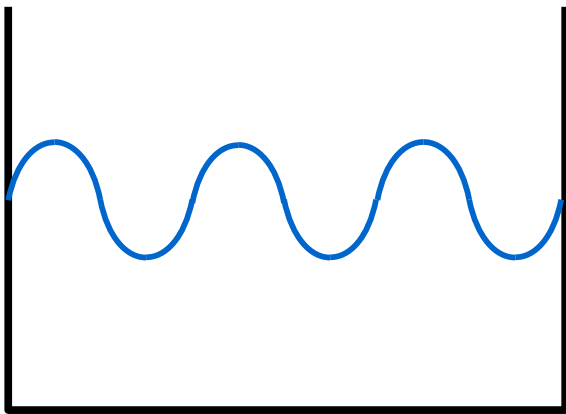


## Pseudopotential



# Kinetic energy pseudopotential

$$H = \text{KE} + V_{e-i} + V_{e-e}$$



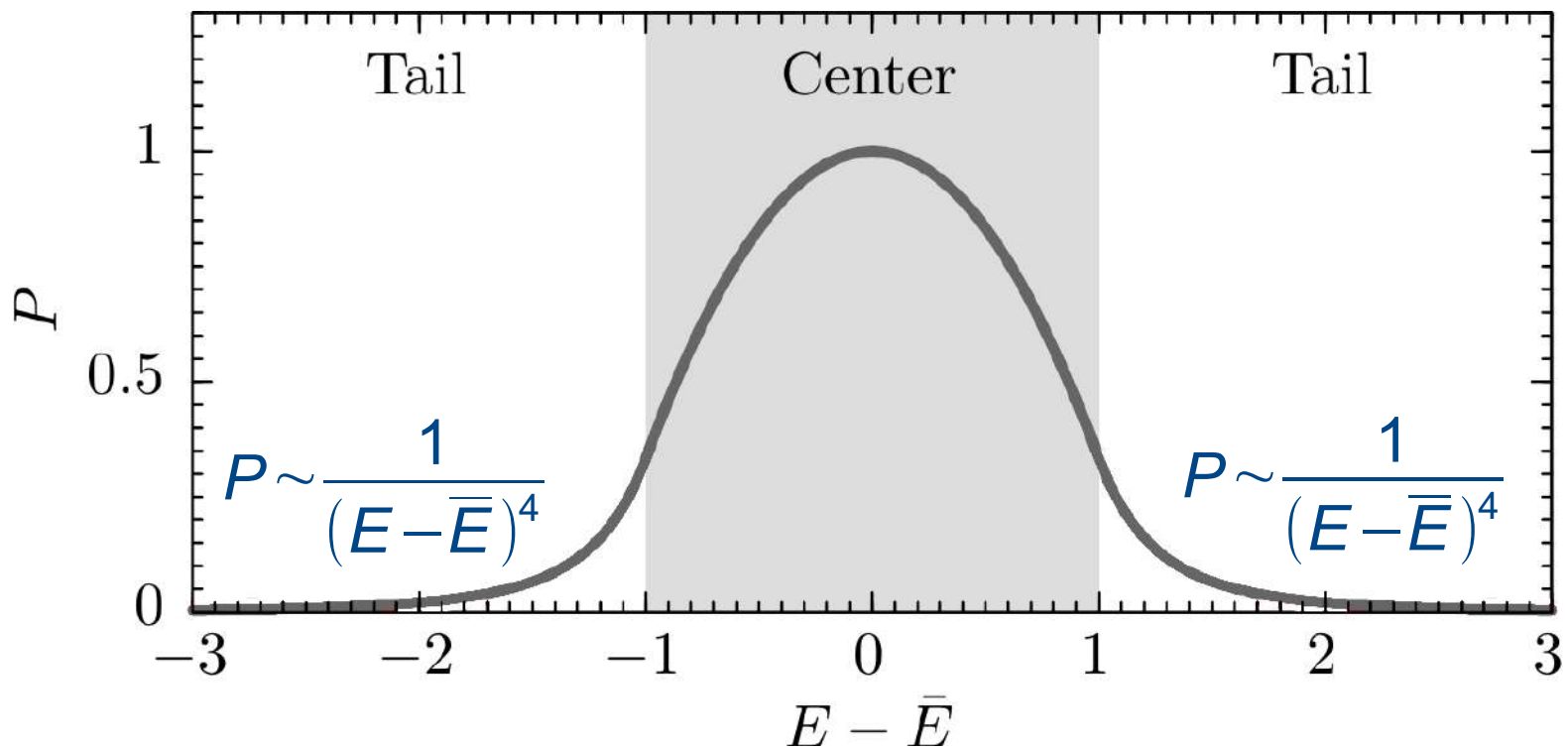
# Kinetic energy pseudopotential

$$H = KE + V_{e-i} + V_{e-e}$$

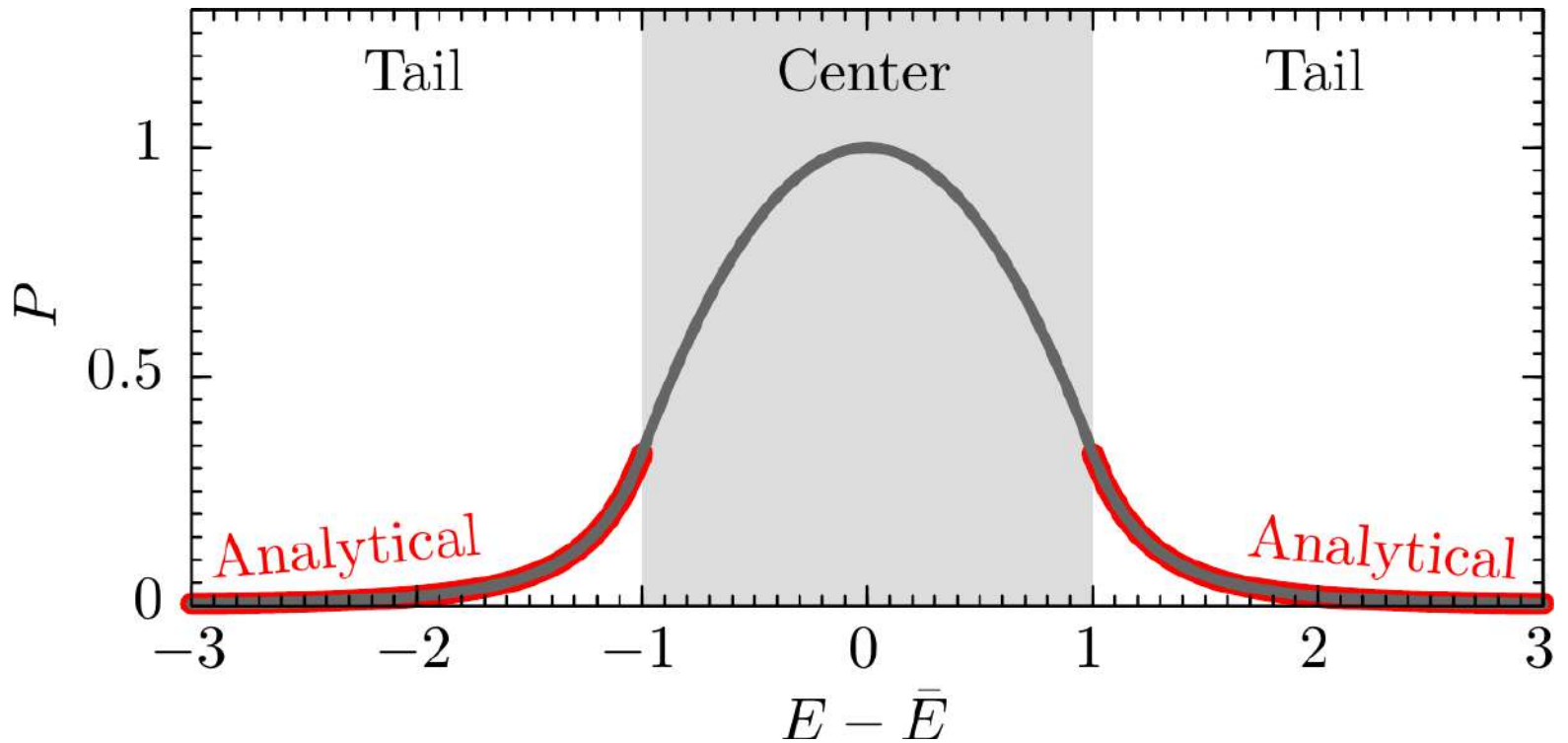
$$E = \frac{\int \bar{\psi} H \psi d\mathbf{r}}{\int \bar{\psi} \psi d\mathbf{r}} = \int_{\psi^2} \frac{H \psi}{\psi}$$

at a node  $\psi \sim x$  so  $E_L = \frac{H \psi}{\psi} \sim \frac{1}{x}$

# Energy distribution

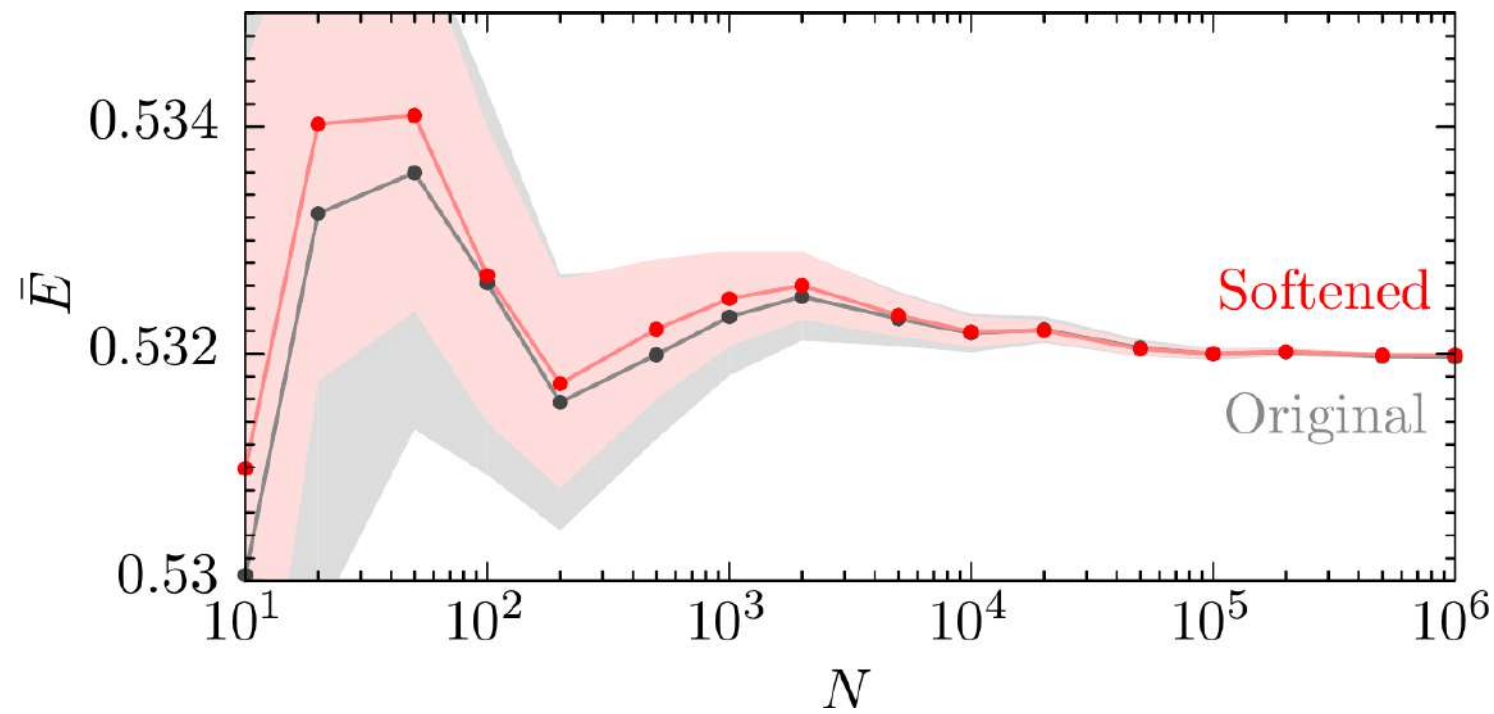


# Softening the energy distribution

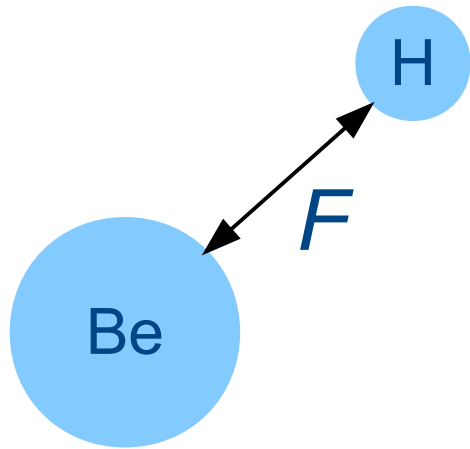




# Accuracy of the homogeneous electron gas

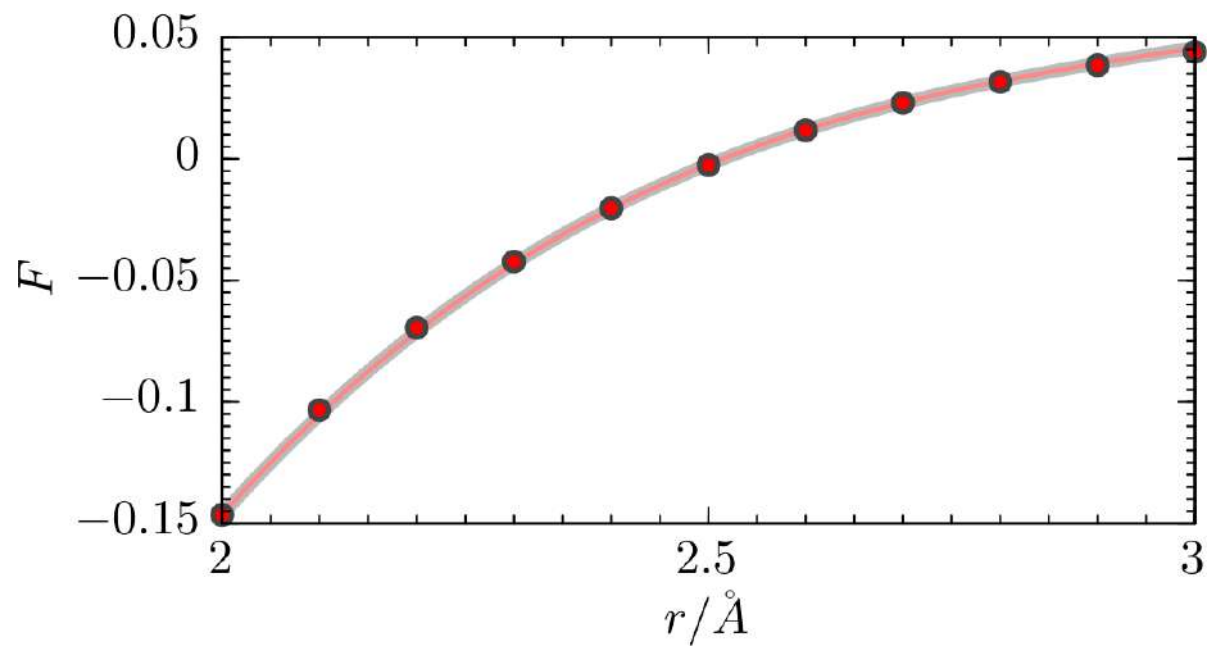


# Force distribution

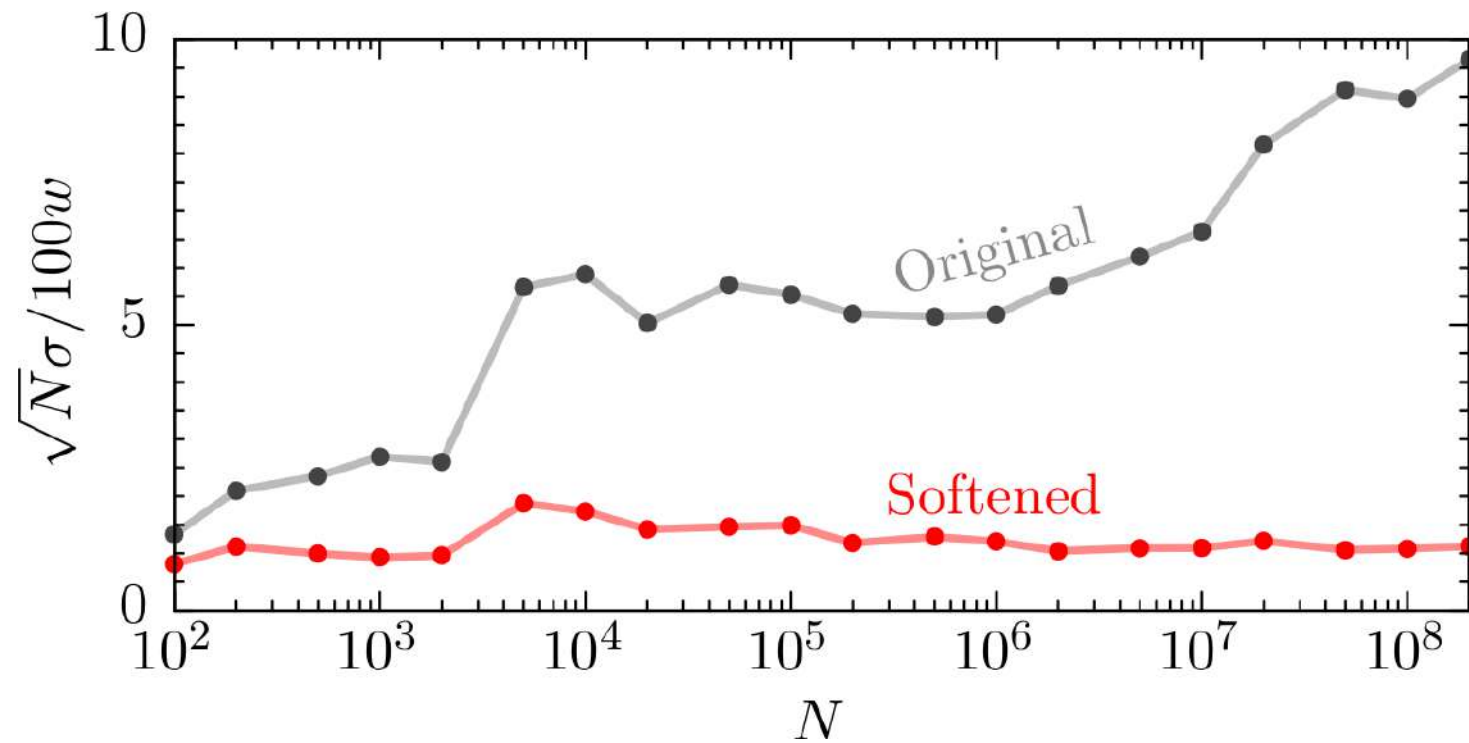


$$P \sim \frac{1}{F^{5/2}}$$

# Force distribution



# Force variance



# Forces in DMC

Van der Waals interaction

Peierls distortion

Bond breaking

# Summary

Developed a pseudopotential for the contact, Coulomb, and dipolar interactions

Studied many-body phenomena

Proposed a formalism to soften the energy and interatomic force distribution

Python tool: <https://pypi.python.org/pypi/contactpp>