

In the **Atomic Interactions** simulation, students investigate how the relationship between attractive and repulsive forces govern the interaction between atoms.

The screenshot shows the PhET Atomic Interactions simulation interface. It features a central graph of Potential Energy vs. Distance Between Atoms, a 3D model of two atoms, and a control panel on the right. Callout boxes provide instructions:

- SCALE** the graph: Points to the zoom-in (+) and zoom-out (-) icons.
- VIEW** the interaction between atoms: Points to the 3D model of two atoms.
- DRAG** the unpinned atom or the dot on the graph: Points to the unpinned atom in the 3D model and the corresponding dot on the graph.
- CHOOSE** pairs of atoms to investigate or create a custom set: Points to the atom selection menu on the right, which lists Pinned and Moving atoms (Neon, Argon, Oxygen) and a Custom Attraction option.
- HIDE or SHOW** forces between atoms: Points to the force control panel on the right, which includes options for Hide Forces, Total Force, Attractive (van der Waals), and Repulsive (electron overlap), each with a corresponding arrow icon.

At the bottom of the simulation, there are playback controls for Slow Motion and Normal, and a PhET logo with a menu icon.

Complex Controls

- The background of the simulation can be changed for easier projection by going to the PhET menu bar, selecting Options, and checking Projector mode.



Insights into Student Use

- The force arrows are off by default as students initially found them overwhelming in interviews.
- Student interviews also indicated that students found the force arrows helpful for making sense of the potential energy graph.

Model Simplifications

- The interaction between the two atoms is modeled using the Lennard-Jones potential.
- The atomic radius, which corresponds to σ (sigma), is the Van der Waals radius.
- For the oxygen-oxygen, epsilon (ϵ) roughly corresponds to the average bond energy. Sigma (σ) was calculated based on the average bond length as the bottom of the potential energy well is located at the equilibrium bond distance.
- The bonding behavior for oxygen-oxygen is not explicitly shown in this simulation, as it would require a third atom to carry off excess energy. However, the well depth and forces between the oxygen atoms are much larger, consistent with a bonding pair.

Suggestions for Use

Sample Challenge Prompts

- Describe how attractive and repulsive forces influence the attraction between two atoms.
- Explain the relationship between the attractive forces between atoms and the potential energy graph for the atom pair.
- Compare and contrast the behavior and potential energy graph for all atom pairs.
- Describe what would need to happen for oxygen to form a diatomic bond.
- Define the values of σ and ϵ .

See all published activities for Atomic Interactions [here](#).

For more tips on using PhET sims with your students, see [Tips for Using PhET](#).