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This text is an atoms-first adaptation of OpenStax Chemistry 2e. The intention of “atoms-first” involves a few basic principles: first, it introduces atomic and molecular structure much

earlier than the traditional approach, and it threads these themes through subsequent chapters. This approach may be chosen as a way to delay the ...

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Molecular dynamics (MD) is a

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computer simulation method for analyzing the physical movements of atoms and molecules. The atoms and molecules are allowed to interact for a fixed period of time, giving a view of the

dynamic "evolution" of the system. In the most common version, the trajectories of atoms and molecules are determined by numerically solving Newton's equations of motion for a system ...

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rocksalt oxyfluorides are receiving significant attention due to their high capacity and lower voltage hysteresis compared with ordered Li-rich layered compounds. However, a deep understanding of these

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mol}^{-1} \label{5}\] The positive sign shows that you have to put in energy to perform this change. The second electron affinity of oxygen is particularly high because the electron is being

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**forced into a small, very
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Initiation, Propagation, and Termination In Free Radical Reactions. In the previous post on free radical substitution reactions we talked about why heat or light is required in free-radical

reactions. In this post we're going to go through the mechanism of a free-radical substitution reaction, which has three key types of steps: initiation, propagation, and termination.

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