MOE is a leading drug discovery software platform that integrates visualization, modeling and simulations, as well as methodology development, in one package. MOE scientific applications are used by biologists, medicinal chemists and computational chemists in pharmaceutical, biotechnology and academic research. MOE runs on Windows, Linux, Unix, and MAC OS X.

**MAIN APPLICATION AREAS**

- Structure-Based Design
- Fragment-Based Design
- Pharmacophore Discovery
- Medicinal Chemistry Applications
- Biologics Applications
- Protein and Antibody Modeling
- Molecular Modeling and Simulations
- Cheminformatics & QSAR

**About MOE**

To learn more about MOE, please visit: chemcomp.com