FOR IMMEDIATE RELEASE

# CHEMICAL COMPUTING GROUP RELEASES MOE 2016.08

**MONTREAL, Canada – October 20, 2016** – Chemical Computing Group is pleased to announce the 2016 release of the Molecular Operating Environment (MOE). The 2016.08 version of MOE contains a unified web-based framework for ligand analytics as well as enhanced spectral analysis tools for structure determination. The key features for MOE 2016 include:

- **MOEsaic: Web-Application for Ligand Analytics**
- **Spectral Analysis for Structure Determination**
- **Enhanced Protein Patch Analyzer**
- **Integrated Antibody Project Database and Antibody Homology Modeler**
- **Small Footprint MOE to Facilitate Large Scale Deployments**
- **Physical and Virtual Rendering of Structures**

Additional features in MOE 2016.08 include 2D Protein Patch maps, Protein Patch Descriptors for QSPR modeling and updates to the Protein-Protein Docking and Protein Design applications. For more information on MOE 2016.08, please contact: info@chemcomp.com

## About Chemical Computing Group

CCG (Chemical Computing Group) is a leading supplier of software solutions for life sciences. With a proven track record in scientific innovation, CCG continues to provide state-of-the-art applications in drug discovery to pharmaceutical, biotechnology and academic researchers. CCG's software platform is the Molecular Operating Environment (MOE) which is used by computational chemists, medicinal chemists and biologists in the major pharmaceutical and biotechnology companies throughout the world. CCG has a very strong reputation for collaborative scientific support, maintaining support offices in both Europe and North America. Founded in 1994, CCG is headquartered in Montreal, Canada. For more information, visit: www.chemcomp.com

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