

This is a full version of Chemwindows 6.0.rar, which contains the latest edition of Chemwindows and its predecessor 5.0 If you're looking for a handy chemistry app to enhance your desktop experience, try ChemWindows! This free and open-source project includes a variety of tools: you can easily convert between units and calculate properties such as density or vapor pressure; there's also an interactive periodic table that lets you look up information about different elements from Latin to English; you can view 2D animations of molecular structures, explore 3D molecule structures in space, or make models from scratch with the Molecule Editor. ChemWindows is free to download, use, and modify. Its 3D molecular models are available under the Creative Commons Attribution-NonCommercial 4.0 International license. "ChemWindows": <http://www.chemwindows.org> "Molecule Editor": <http://www.chemwindows.org/software/moledit3dframe.html>

The main developer of ChemWindows is Nicolas Lecompte, who relaunched the project on SourceForge in October 2007 with Version 2, then released an update in May 2008 that prompted him to rename it "ChemWindows". The project's motto is "La science communicative", which roughly translates as "communicating science". Version 3 has been released, with special focus on the Molecule Editor.

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The package contains these files: 1 .tar file : main ChemWindows 6 files (19MB) - ChemWindows 6.

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