



Efficient Chemical
Drawing



Chemical & Biological
Intelligence



Eye Catching
Visual Chemistry



Integrations with Key
Research Applications
& Customizations



Quick & Painless
Chemistry Reporting



Modern Cloud
Architecture

ChemDraw Prime



- Comprehensive Chemical Drawing Tools
- Hotkeys & Shortcuts
- Generic Structures
- Structure & Reaction Clean-up
- Smart Copy/Paste (SMILES to Structure, auto-positioning)



- Enhanced stereo perception
- Enhanced Stereo Perception
- M/P Stereochemistry Labeling of Allenes and Atropisomers
- Physico-chemical Property Predictions: pKa, logP, logS and more (on Windows)

ChemDraw Professional



- Comprehensive Chemical Drawing Tools
- Hotkeys & Shortcuts
- Generic Structures
- Structure & Reaction Clean-up
- Smart Copy/Paste (SMILES to Structure, auto-positioning)



- Name-to-Structure & Structure-to-Name
- Enhanced Stereo Perception
- M/P Stereochemistry Labeling of Allenes and Atropisomers
- Physico-chemical Property Predictions: pKa, logP, logS and more (on Windows)
- ¹H & ¹³C NMR Prediction
- Advanced HELM Biopolymer Representation
- Support for Atom-mapped SMILES in HELM Strings, HELM In-line Annotations & FASTA



- Biopolymer Clean-up & Enhanced Display
- Biopolymer Colored Document Settings
- Atom / Bond / Ring-Fill Color Highlighting
- 3D Clean-up
- Structure Perspective Depth-fading
- Structure from CAS-RN with ChemACX



- CAS SciFinder™ Integration
- Elsevier Reaxys® Integration

Signals ChemDraw



- Comprehensive Chemical Drawing Tools
- Hotkeys (includes HELM Hotkeys) & Shortcuts
- Generic Structures
- Structure & Reaction Clean-up
- Smart Copy/Paste (SMILES to Structure, auto-positioning)



- Name-to-Structure & Structure-to-Name
- Enhanced Stereo Perception
- M/P Stereochemistry Labeling of Allenes and Atropisomers
- Properties, Query & Guided Stereo panels
- Physico-chemical property predictions: pKa, logP, logS and more (on Windows)
- ¹H & ¹³C NMR Prediction
- Advanced HELM Biopolymer Representation (includes Watson-Crick base pairing tools)
- Support for Atom-mapped SMILES in HELM Strings, HELM in-line annotations & FASTA
- Centralized Biopolymer Library Management in Signals (HELM Monomer Curation)



- Biopolymer Clean-up & Enhanced Display
- Biopolymer Colored Document Settings
- Atom / Bond / Ring-Fill Color Highlighting
- 3D Clean-up
- Structure Perspective Depth-fading
- Save and Copy as 3D Printable Format (3MF)



- CAS SciFinder™ Integration
- Elsevier Reaxys® Integration
- Centralized HELM Monomer Libraries
- ChemACX Explorer Add-in
- Google Scholar/Patents Add-in
- PubChem Safety Add-in
- Custom, Dynamic Add-in SDK



- Cloud Anywhere Access and Automatic Updates
- Create & Share Drawings and Notebooks
- Create Collections of Molecules/Reactions
- Add/Edit Properties in Collections
- Export Collections to .sdf/.pptx Files
- Create Reports of Signals Notebook Experiments
- Extract ChemDraw/BIOVIA Draw Drawings from Documents
- Perform Structure Searches of Drawings



- Centralized User Management (SaaS)
- ChemDraw Desktop Automatic Updates
- Create, Manage & Curate Chemical Drawings from your Web Browser

Revvity Signals Software, Inc
77 4th Avenue | Waltham, MA 02451 USA
P: (800) 762-4000 or (+1) 203-925-4602
revvitysignals.com/company/contact

Copyright ©, Revvity, Inc. All rights reserved. Revvity™ is a registered trademark of Revvity, Inc. All other trademarks are the property of their respective owners.

40 Years of ChemDraw by
revvity signals

revvity
signals

JUNE 2025