Version 23.0 NEW FEATURES	Platform	ChemDraw Prime	ChemDraw Professional
Dark Mode Style Sheet	win/mac	●	●
Atropisomer perception	win/mac	●	●
Ignore Top Level Chiral flag	win/mac	●	●
Smart Paste (no overlapping on paste actions)	win/mac	●	●
Hydrogen Bonding in 3D cleanup	win/mac		●

Recent Additions	Platform	ChemDraw Prime	ChemDraw Professional
Magic Hotkeys Enhancements	Win/Mac	●	●
Shortcuts Enhancements	Win/Mac	●	●
Join function improvements	Win/Mac	●	●
Smart Copy/Paste (SMILES, InChI, HELM)	Win/Mac	●	●
Aromatic Cycle Display Toggle and Preferences	Win/Mac	●	●
Stereochemistry handling improvements	Win/Mac	●	●
Improved Polymer Brackets (Average MW)	Win/Mac	●	●
Hydrogen Bond Tool	Win/Mac	●	●
Open CIF Files	Win/Mac	●	●
Atom/Bond Color Highlighting	Win/Mac		●
Ring-Fill Coloring	Win/Mac		●
Search into SciFinder-n	Win/Mac		●
Search into Reaxys	Win/Mac		●
Improved HELM Monomer Toolbar	Win/Mac		●
HELM Monomer Versioning Support	Win/Mac		●

HELM Cartoon Representation	Win/Mac		●
Support for ambiguous FASTA/HELM Monomers	Win/Mac		●
Includes	**Platform**	**ChemDraw Prime**	**ChemDraw Professional**
Read and Save as .cdx / .cdxml Files	Win/Mac	●	●
Read and Save as .rxn Files (v2000, v3000)	Win/Mac	●	●
Read and Save as .skc Files	Win/Mac	●	●
Read and Save as .mol Files (v2000, v3000)	Win/Mac	●	●
Read and Save as .sdf Files (v2000, v3000)	Win/Mac	●	●
Read and Save as .rdf Files (v2000, v3000)	Win/Mac	●	●
Save ChemDraw Style Sheet	Win/Mac	●	●
Structure Clean-up	Win/Mac	●	●
Reaction Clean-up	Win/Mac	●	●
Magic Hotkeys	Win/Mac	●	●
Chemical Bonds Tools	Win/Mac	●	●
Text Tool	Win/Mac	●	●
3D Perspective Tool	Win/Mac	●	●
Chemical Rings Tools	Win/Mac	●	●
Arrow Tool	Win/Mac	●	●
Orbitals Tool	Win/Mac	●	●
Brackets Tool	Win/Mac	●	●
Pen Tools	Win/Mac	●	●
Shapes Tool	Win/Mac	●	●
Chemical Polymers Tools	Win/Mac	●	●

Mass Fragmentation Tools	Win/Mac	●	●
Thin Layer Chromatography Tool	Win/Mac	●	●
Gel Electrophoresis Tool	Win/Mac	●	●
Insert OLE Object in ChemDraw	Win	●	●
Copy ChemDraw Structures as OLE Object	Win	●	●
Show Stereochemistry	Win/Mac	●	●
Relative Stereochemistry (ISIS compatibility)	Win/Mac	●	●
Reaction Interpretation	Win/Mac	●	●
Reaction Mapping	Win/Mac	●	●
Calculate MW	Win/Mac	●	●
Calculate Exact Mass	Win/Mac	●	●
Calculate Chemical Formula	Win/Mac	●	●
Calculate Elemental Analysis	Win/Mac	●	●
Calculate m/z	Win/Mac	●	●
Copy/Paste as CDXML	Win/Mac	●	●
Copy/Paste as SMILES	Win/Mac	●	●
Copy/Paste as SYBYL (SLN)	Win/Mac	●	●
Copy/Paste as InChI	Win/Mac	●	●
Copy/Paste as Mol File / Mol3000	Win/Mac	●	●
pKa / Log P / Log S	Win/Mac	●	●
tPSA	Win/Mac	●	●
Atom List Generic Structures (Enumeration)	Win/Mac	●	●
Variable Attachment Generic Structures (Enumeration)	Win/Mac	●	●

Label Repeating Units Generic Structures (Enumeration)	Win/Mac	●	●
Polymer Repeating Units Generic Structures (Enumeration)	Win/Mac	●	●
Chemical Structures Templates	Win/Mac	●	●
Laboratory Equipment Templates	Win/Mac	●	●
Analyze/Check Structures	Win/Mac	●	●
Expand/Contract Labels	Win/Mac	●	●
Define/Use Nicknames	Win/Mac	●	●
Document Metadata/Tagging	Win/Mac	●	●
Multiple ChemDraw Items Folder	Win/Mac	●	●
Multicenter Attachments	Win/Mac	●	●
Save as JPEG image	Win/Mac	●	●
Save as PNG image	Win/Mac	●	●
Save as TIFF image	Win/Mac	●	●
Save as Scalable Vector Graphics (SVG)	Win/Mac	●	●
Save as Encapsulated Post Script (EPS)	Win/Mac	●	●
Name-to-Structure / Structure-to-Name	Win/Mac		●
Predict 1H NMR	Win/Mac		●
Predict 13C NMR	Win/Mac		●
Search SciFinder	Win/Mac		●
Search SciFinder-n	Win/Mac		●
Search Reaxys	Win/Mac		●
Reaction Stoichiometry Grid	Win/Mac		●
R-Group Table Generic Structures (Enumeration)	Win/Mac		●

BioDraw Toolbar	Win/Mac		●
cLogP	Win/Mac		●
HELM Toolbar	Win/Mac		●
Copy/Paste as HELM	Win/Mac		●
Copy/Paste as FASTA Peptide	Win/Mac		●
Copy/Paste as FASTA DNA/RNA	Win/Mac		●
Support for HELM notation	Win/Mac		●
CAS RN to Structure from ChemACX.com	Win/Mac		●
Enhanced Stereochemistry Support	Win/Mac		●
ChemDraw for Excel	Win		●
CombiChem for Excel	Win		●
Name-to-Structure / Structure-to-Name for ChemDraw for Excel	Win		●
Chem3D Professional	Win		●
ChemFinder Standard	Win		●
ChemScript	Win		●
ClogP/CMR for ChemDraw for Excel/Chem3D	Win		●
Molecular Networks (pKa/Log P/Log S) for ChemDraw for Excel/ Chem 3D	Win		●
Molecular Topology for Chem Draw for Excel/Chem 3D	Win		●
ChemProp Std Properties for Chem Draw for Excel/Chem 3D	Win		●