ChemOffice® 16 Suite of Products

SciFinder®

PerkinElmer is collaborating with Chemical Abstracts Service to combine the power of two leading chemistry solutions, ChemDraw® and SciFinder®.

New Features	Platform	ChemDraw Prime	ChemDraw Professional	ChemOffice Professional
Advanced Retrosynthesis Tool	Win/Mac		X	X
Access to ChemDraw Cloud	Win/Mac		X	X
Access and Integration to ChemDraw E-Notebook	Win			X
Includes	**Platform**	**ChemDraw Prime**	**ChemDraw Professional**	**ChemOffice Professional**
ChemDraw	Win/Mac	X	X	X
Multiple ChemDraw Items Folders	Win/Mac	X	X	X
Save and Read Graphic Files	Win/Mac	X	X	X
Save and Read Chemical Files	Win/Mac	X	X	X
Printing Options	Win/Mac	X	X	X
Chemical Templates	Win/Mac	X	X	X
Equipment Templates	Win/Mac	X	X	X
Analyze/Check Structures	Win/Mac	X	X	X
Insert OLE Object in ChemDraw	Win	X	X	X
In-place OLE Editing of ChemDraw Objects	Win	X	X	X
Show Stereochemistry	Win/Mac	X	X	X
Relative Stereochemistry (ISIS compatibility)	Win/Mac	X	X	X
Reaction Interpretation	Win/Mac	X	X	X
Reaction Mapping	Win/Mac	X	X	X
Calculate Properties	Win/Mac	X	X	X
pKa LogP LogS	Win	X	X	X
Advanced Retrosynthesis Tool	Win/Mac	X	X	X
Auto-numbering of multiple structures	Win/Mac	X	X	X
tPSA	Win/Mac	X	X	X
Hotlink Properties with Structure	Win/Mac	X	X	X
Read JCamp & Galactic Spectra Files	Win/Mac	X	X	X
Manual spectrum/structure assignments	Win/Mac	X	X	X
Chemical Polymer Tools	Win/Mac	X	X	X
Structure Clean Up	Win/Mac	X	X	X
Database Hotlink	Win/Mac	X	X	X
Expand/Contract Labels	Win/Mac	X	X	X
Create/Use Nicknames	Win/Mac	X	X	X
Expand Generic Structure	Win/Mac	X	X	X
Multicenter Attachments	Win/Mac	X	X	X

PerkinElmer®
For the Better

Includes	Platform	ChemDraw Prime	ChemDraw Professional	ChemOffice Professional
TLC/GEP Tools	Win/Mac	X	X	X
Fragmentation Tools	Win/Mac	X	X	X
ChemDraw Active X Plugin	Win/Mac	X	X	X
Special Copy/Paste				
SMILES	Win/Mac	X	X	X
CDX	Win/Mac	X	X	X
Molfile	Win/Mac	X	X	X
CDXML	Win/Mac	X	X	X
ChemDraw Cloud	Win/Mac		X	X
Search SciFinder	Win/Mac		X	X
Name = Structure/Structure = Name	Win/Mac		X	X
cLogP	Win/Mac		X	X
Biopolymer Toolbar	Win/Mac		X	X
BioDraw	Win/Mac		X	X
Reaction Stoichiometry Grid	Win/Mac		X	X
Calculate ^{1}H ^{13}C NMR Spectra	Win/Mac		X	X
Sequences	Win/Mac		X	X
Query Features	Win/Mac		X	X
Advanced Stereochemistry	Win/Mac		X	X
ChemDraw E-Notebook	Win			X
ChemFinder (Std in CD Prof; Ultra in CO)	Win		X	X
ChemDraw for Excel	Win		X	X
Name=Struct for ChemDraw for XL	Win		X	X
ChemScript + Python	Win		X	X
CombiChem for Excel	Win		X	X
3D Search	Win		X	X
Chem3D (Pro in CDPro; Ultra in CO)	Win		X	X
Chem3D Hotlink	Win		X	X
Chem3D Active X Plugin	Win		X	X
GAMESS	Win			X
Interface to MOPAC 2016	Win			X
Interface to Gaussian	Win			X
Interface to Conflex	Win			X
Interface to Autodock	Win			X
ChemFinder / Oracle	Win			X
ChemFinder for Office	Win			X
BioViz in ChemFinder Ultra	Win			X
Compound Profiles in ChemFinder Ultra	Win			X
Clustering in ChemFinder Ultra	Win			X

Visit www.cambridgesoft.com/software/overview.aspx for more information.

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