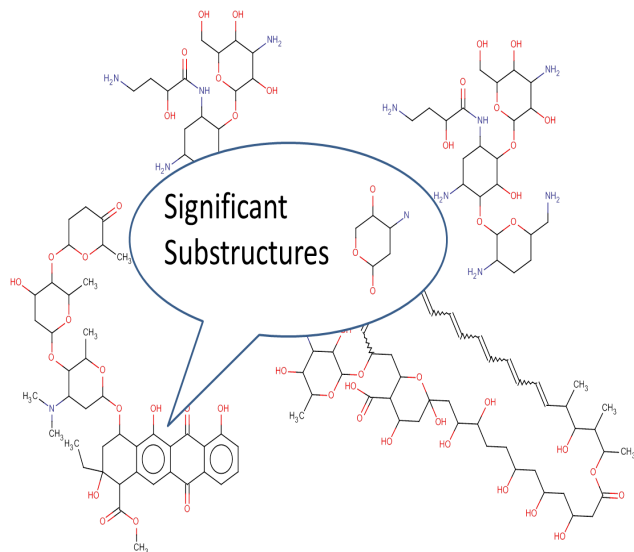
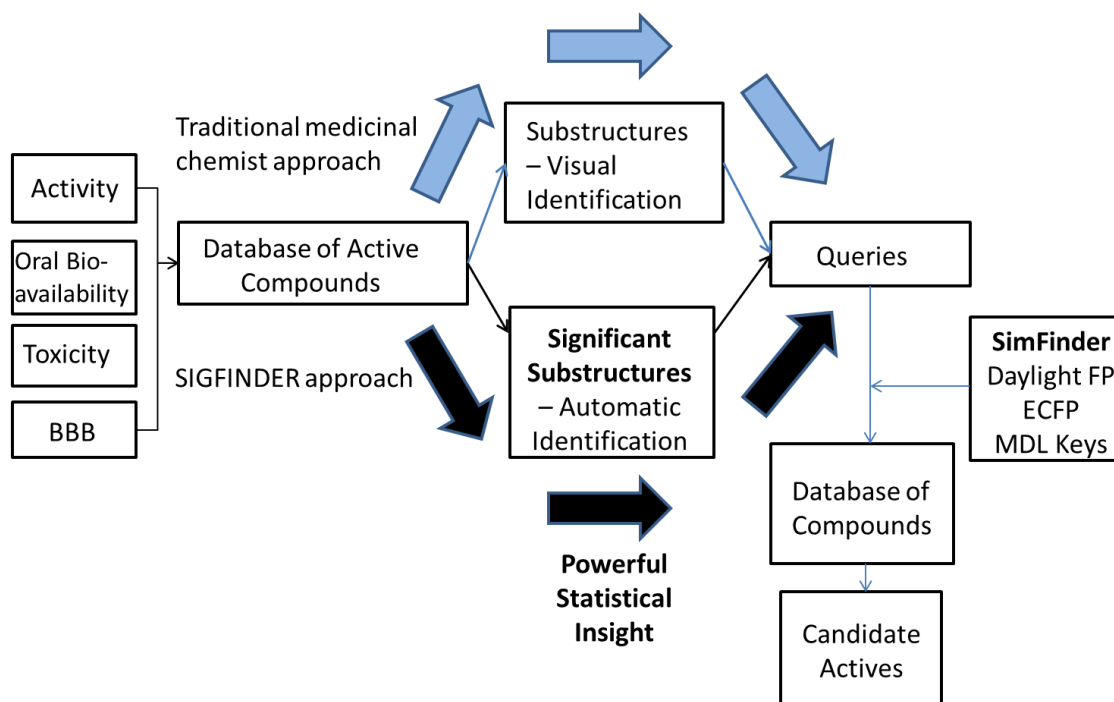




# SigFinder: Powerful Substructure Mining from Chemical Databases



SigFinder is the first tool of its kind that allows the user to automatically extract significant substructures from any chemical database. Using our powerful statistical measures, SigFinder can not only identify substructures that are frequently occurring (something which is already intuitive to a medicinal chemist) but also substructures that are absent or significant in a given database. In addition, SigFinder can classify and predict the properties of drug candidates such as binding, toxicity, and permeability. For more information please visit our website at <http://www.ancelot.com/SigFinder.html>



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