Review of Deep Learning Algorithms in Computational biochemistry

Biochemistry actually is a fertile pool of resources, including modeling, DL and optimization techniques for chemical data and associated phenomena. Computational Biochemistry is a very effective entity. These include rapid literature research, physical and quantum chemical propaganda analyzes, transition states, chemical structures, chemical reactions as well as new catalysts and candidates for drugs. DL algorithms can integrate raw input into intermediary layers of features and successfully fit the desired compound into optimum combinations. This review study introduces a series of biochemical applications with the most exciting discoveries involving the use of DL in wide range of applications such as Modeling bio-processes, algorithms and methods to promote the design and synthesis of bio compounds, material design, binding forecasting and molecular activity are given special importance.