

# The Prediction of the binding Affinity of BACE-1 using DeepLearning and Quantum Chemical Descriptors

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## Abstract

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## Descriptors

QSAR descriptors<sup>1</sup>

## References

- (1) Wang, L.; Ding, J.; Pan, L.; Cao, D.; Jiang, H.; Ding, X. Quantum chemical descriptors in quantitative structure–activity relationship models and their applications. *Chemometrics and Intelligent Laboratory Systems* **2021**, *217*, 104384.