

# Neural Network and Random Forest Algorithms as Catalysts in QSAR/QSAAR Modeling: Targeting Carbonic Anhydrase for Antituberculosis Drug Design

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The authors have withdrawn their manuscript owing to the inability to complete the revisions or follow up on the manuscript at this time, due to personal circumstances. Therefore, the authors do not wish this work to be cited as reference for the project. If you have any questions, please contact the corresponding author.