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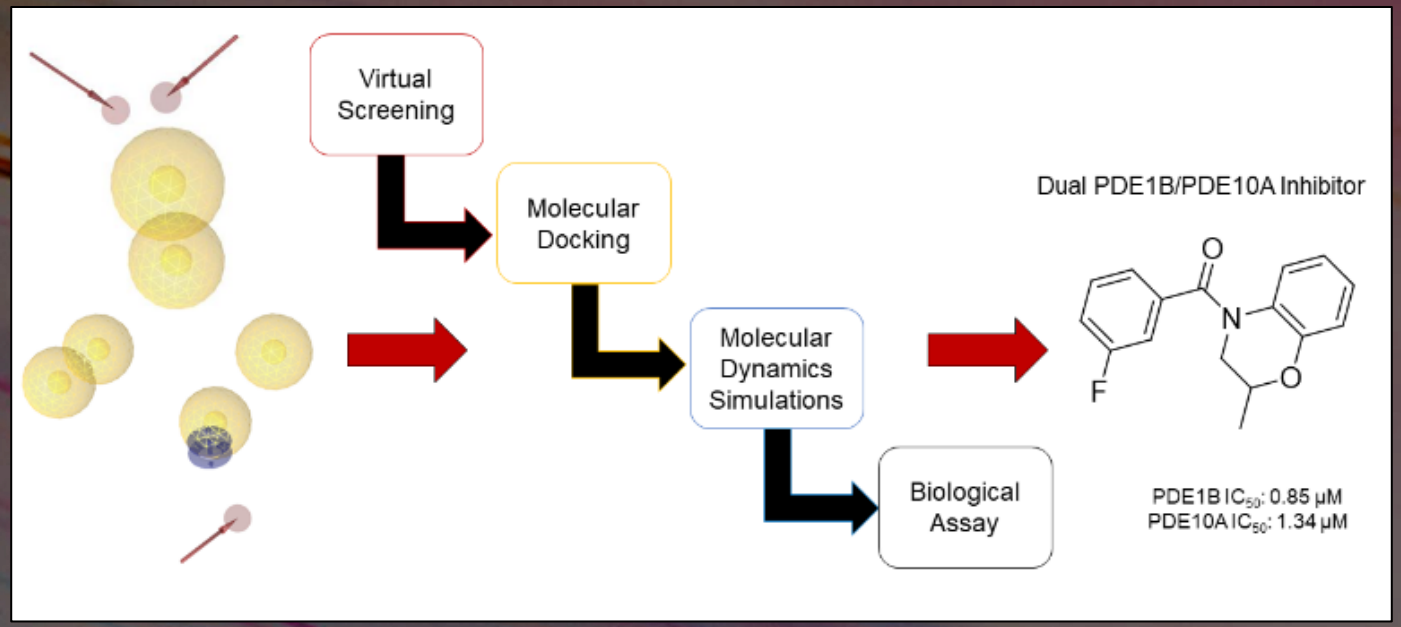

Identification of dual inhibitor of phosphodiesterase 1B/10A using structure-based drug design approach
Journal of Molecular Liquids, 2021(IF 6.165).

Areas of Industrial Collaboration

- QSAR, Computational Drug Design & Discovery, Virtual Screening, Molecular Dynamics Simulations, Drug Synthesis, Drug Evaluation

Community Service Projects

- Blood Donation Camps, COVID-19 Vaccination

Consultancy Services

- Solving pharmaceutical problems related to drug discovery & development by application of computational technology

Awards and Recognitions

- Editorial Board Member of PLOS ONE
- Visiting Lecturer Anton De Kom University of Suriname
- Recipient of RM 234100 external research grants and RM 137000 internal research grants
- Silver Award for Best Oral Presentation at 13th MI Conferenc

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