JUN 30, 2020

OPEN ACCESS

DOI:
dx.doi.org/10.17504/protocols.io.bh3mj8k6

Protocol Citation: Maria .
 2020. Virtual Screening for
 insilico drug designing.
protocols.io
<https://dx.doi.org/10.17504/protocols.io.bh3mj8k6>

License: This is an open
 access protocol distributed
 under the terms of
 the [Creative Commons
 Attribution License](#), which
 permits unrestricted use,
 distribution, and reproduction
 in any medium, provided the
 original author and source are
 credited

Protocol status: Working
 We use this protocol and it's
 working

Created: Jun 30, 2020

Last Modified: Jun 30, 2020

PROTOCOL integer ID:
 38733

Virtual Screening for insilico drug designing

Maria .¹

¹Institute of Chemical Sciences, University of Peshawar, Pakistan

Maria .

DISCLAIMER

DISCLAIMER – FOR INFORMATIONAL PURPOSES ONLY; USE AT YOUR OWN RISK

The protocol content here is for informational purposes only and does not constitute legal, medical, clinical, or safety advice, or otherwise; content added to [protocols.io](#) is not peer reviewed and may not have undergone a formal approval of any kind. Information presented in this protocol should not substitute for independent professional judgment, advice, diagnosis, or treatment. Any action you take or refrain from taking using or relying upon the information presented here is strictly at your own risk. You agree that neither the Company nor any of the authors, contributors, administrators, or anyone else associated with [protocols.io](#), can be held responsible for your use of the information contained in or linked to this protocol or any of our Sites/Apps and Services.

1 1Retrieval and Refinement of Crystal Structure

2 2Combinatorial Library Designing

3 3Docking and Virtual Screening

4 4Docking Protocol applied by MOE

5 5Docking Analysis