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These Terms of Use are applicable to all users of ChemBioServer. Users should read, understand and accept these Terms when accessing ChemBioServer. If a given user has not done this, or does not agree with the contents of these Terms of Use, that user should not make use of any of the services provided via ChemBioServer.

By accepting these Terms of Use, the user enters into an agreement with ChemBioServer made up by the contents of these Terms, in order to regulate the user's access and use of the ChemBioServer. Users which accept these Terms of Use hereby represent and warrant that: They accept to be bound by these Terms of Use, either individually, or in representation of an organisation;

They are capable, under the terms of the law applicable to them, to be legally bound by and comply with these Terms of use;

Their use of the ChemBioServer will be carried out in accordance with all applicable laws and regulations;

In case users publish results produced from ChemBioServer or using ChemBioServer by any means, they are obligated to cite:

Karatzas E, Zamora JE, Athanasiadis E, Dellis D, Cournia Z, Spyrou G. "ChemBioServer2.0: An Advanced Web Server for Filtering, Clustering and Networking of Chemical Compounds Facilitating Both Drug Discovery and Repurposing"

Bioinformatics 2020, pii: btz976

<http://dx.doi.org/10.1093/bioinformatics/btz976>

In the event of doubts concerning the Terms of Use, ChemBioServer may be contacted at: zcournia@bioacademy.gr or georges@cing.ac.cy.

For information on the processing of personal data on users which is carried out via the ChemBioServer, please see the ChemBioServer privacy policy.

1. Services provided

ChemBioServer is a publicly available web-application for effectively mining and filtering chemical compounds used in drug discovery. Features: (i) browse and visualize compounds along with their properties, (ii) filter chemical compounds for a variety of properties such as steric clashes and toxicity, (iii) apply perfect match substructure search, (iv) cluster compounds according to their physicochemical properties providing representative compounds for each cluster, (v) build custom compound mining pipelines and (vi) quantify through property graphs the top ranking compounds in drug discovery procedures. ChemBioServer allows for pre-processing of compounds prior to an in silico screen, as well as for post-processing of top-ranked molecules resulting from a docking exercise with the aim to increase the efficiency and the quality of compound selection that will pass to the experimental test phase.

2. Copyright / Intellectual Property

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3. Membership or subscription fees

ChemBioServer is free for every user and no registration is required.

4. User data protection

The output of ChemBioServer belongs to the users and we have no authority to use these results for our own benefit or to share them with others. The output link and output data are kept for 1 (one) week and then they are automatically deleted.

5. Liability

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