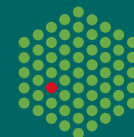
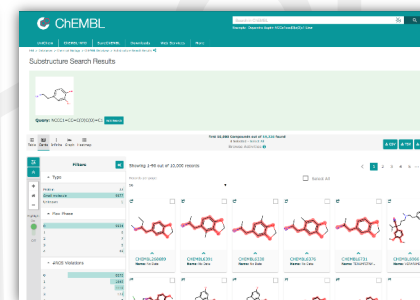
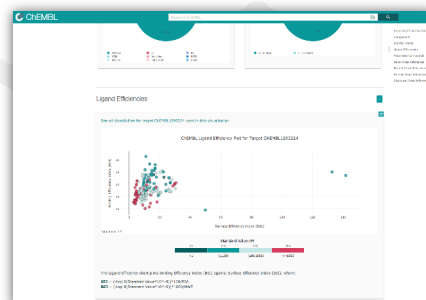
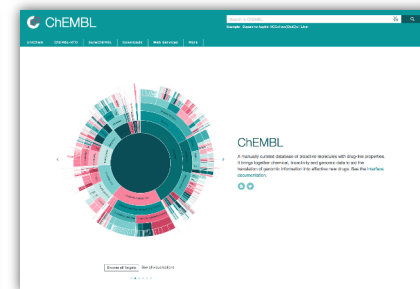
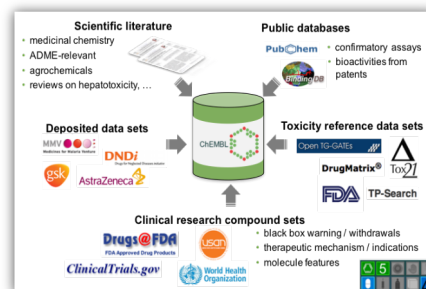
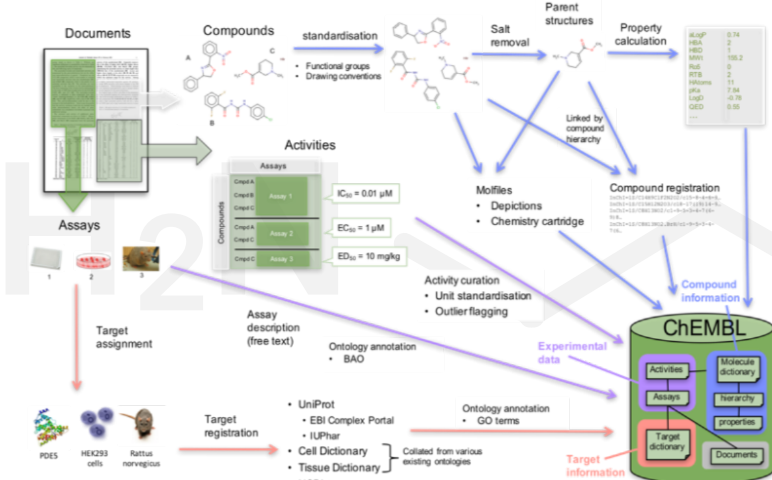


A FAIRer ChEMBL website

EMBL-EBI



CURATION & STANDARDISATION



Why & How?

Free, publically accessible database

High quality, extensive, curated data:

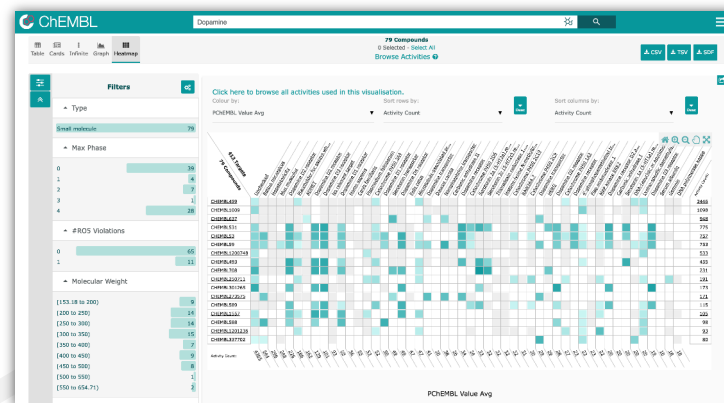
- 12,091 biological targets (including proteins and organisms).
- 1.8 million distinct compounds (small molecules or peptides).
- 15 million bioactivities (binding, functional assays and ADMET).
- From 69,000 scientific papers.

<https://www.ebi.ac.uk/chembl/beta>

Scan me!



BIOACTIVITY DATA ANALYSIS ON THE GO



Going FAIR

Findable

- Unique and persistent identifiers.
- Machine readable metadata available.
- Data is discoverable through search indexes.

Accessible

- Our entire database is downloadable.
- The interface code is open source and universally implementable.

Interoperable

- The data is exposed in RDF format.
- ChEMBL uses well-known ontologies such as BAO, Uberon, EFO among others.

Reusable

- Data is available under CC BY-SA 3.0.
- Source datasets are linked and retrievable.
- Broadly used within the community.