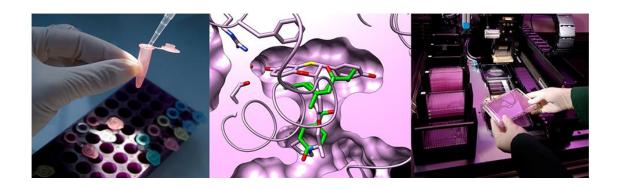


ChemBisFrance Small molecules for health and research











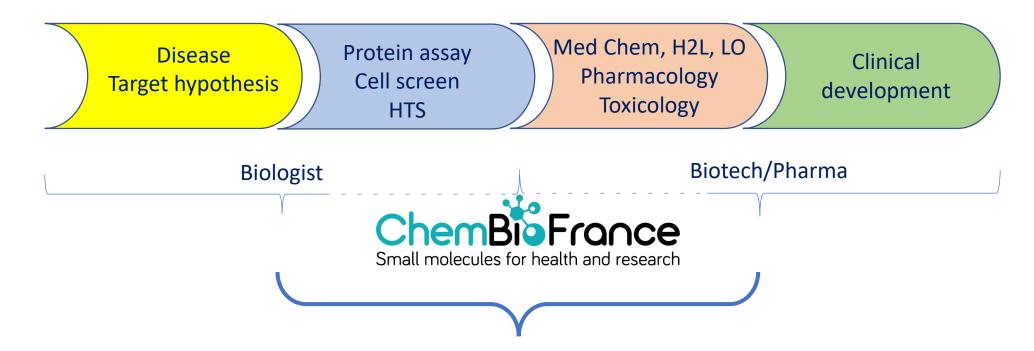












ChemBioFrance is a strategic partner in small molecule drug discovery.

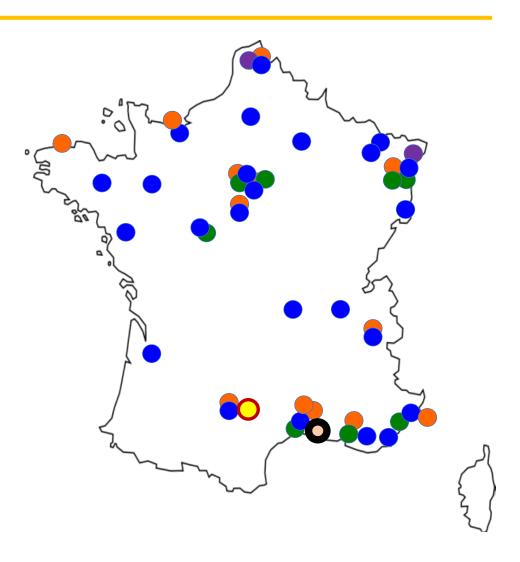
All you need in a place



A massively distributed infrastructure



- 12 Screening centers
- 8 Chemoinformatics sites
- 2 ADME-Tox platforms
- 49 contributing laboratories of the chemical library
- 1 Management unit = CNRS UAR 3035
- Chimiothèque Nationale: 1 Repository of compounds





The National Chemical Library





Chimiothèque historique du laboratoire de chimie du MNHN – A. Latzoura

Features

A compounds library enriched in drug-like.

A high Scaffold-to-compound ratio

Flexible IP management

Originality, no commercial compounds

A team of chemists open for research collaborations (Hit-to-Lead project)

A network of synthesis platforms



The National Chemical Library



Libraries ready to be screened

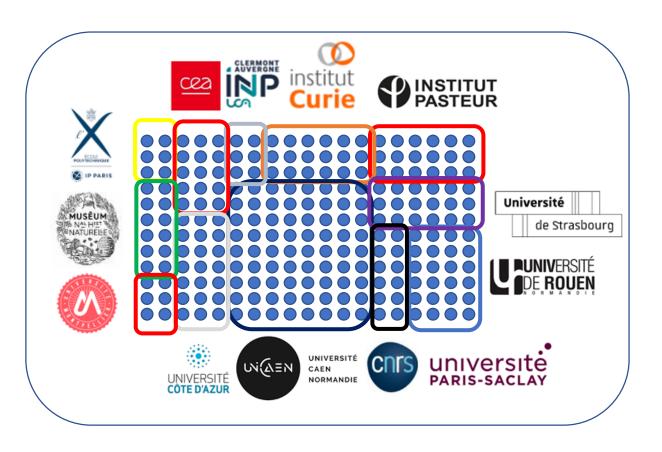
CN > 70 000 cpds at 5mM DMSO in 96 or 384 well/plates

Design of focused library by cherry picking

CNE - Essential Chemical Library

- * 1040 compounds
- * Mirror the diversity of the CN
- * Easy to screen

Natural extracts > 15 000 at 10mg/mL DMSO



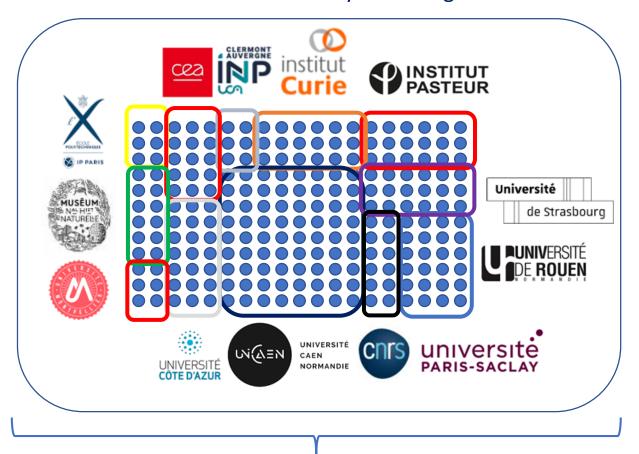
+ 36 members of ChemBioFrance



Scientific collaboration (1/2)



Chemical library screening

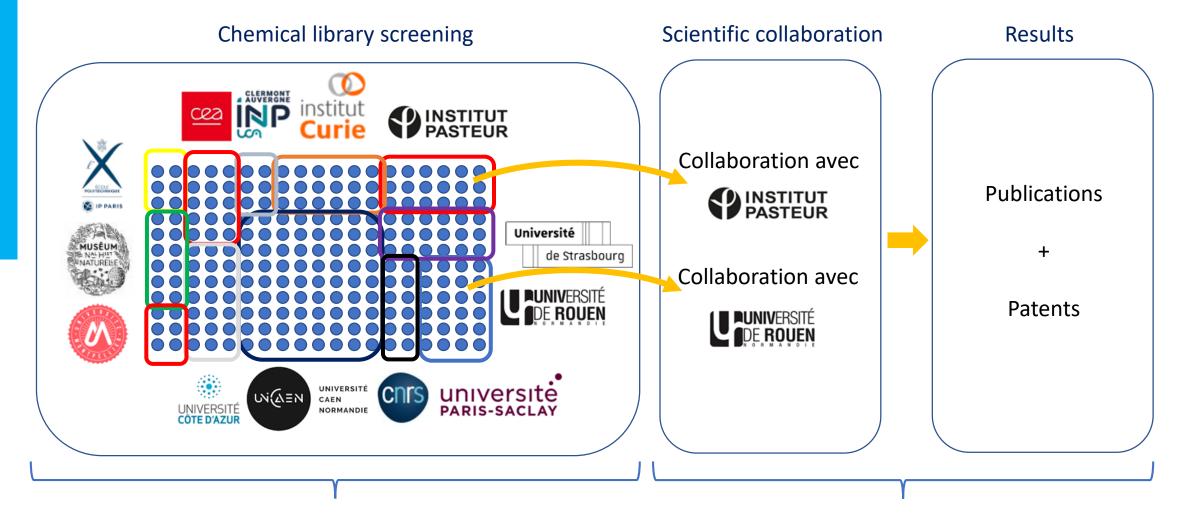


Access to the chemical library with a M.T.A.



Scientific collaboration (2/2)





Access to the chemical library with a M.T.A.

Collaboration between the biologist and the chemist



Screening



Hit identification technologies

Knowledge-based screening (kinase, GPCR)

Diversity screening

Fragment-based screening

Phenotypic screening

Structure-based drug design





A network of specialized screening centers



Functional assays

✓ Proteins (kinase, GPCR)

✓ Cells 2D and 3D

✓ Nucleic acid

Technologies

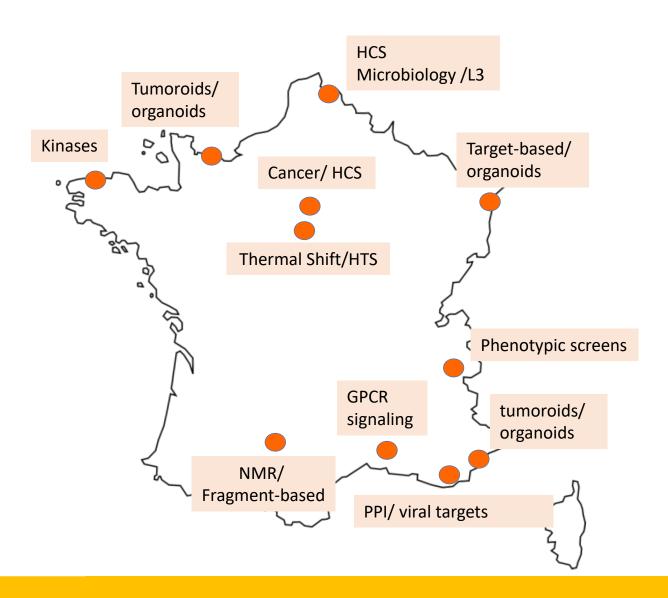
✓ Fluorescence (HTRF, FRET, AlphaScreen)

✓ Luminescence

✓ Radioactivity

Specialized platforms

- ✓ Tumoroids
- ✓ Infectious agents
- ✓ Fragment-based screening
- ✓ Inflammation







A Chemoinformatics network

Virtual Screenings

- ✓ ligand-based
- ✓ structure-based

Libraries Design

- ✓ CN subset
- ✓ protein-protein inhibitors
- ✓ interfering peptides

Raw data analysis and SAR building

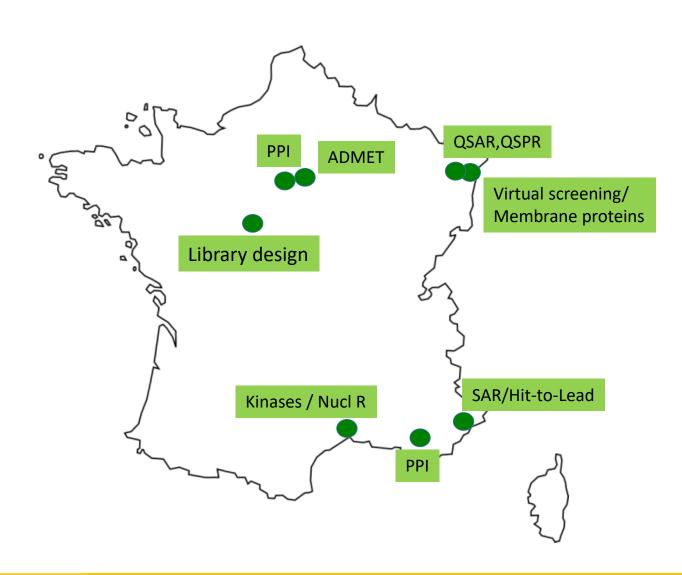
- ✓ PAINS/reactivity alert
- ✓ analogs in catalogs
- ✓ properties prediction LogP, Sol.,Kd, K_{on}/K_{off}

Hit-to-lead

- ✓ prioritizing cpds
- ✓ prioritizing reactants
- ✓ scaffold hopping

Bio-Profiling & ADME-Tox Prediction

- ✓ Metabolism (site, CYP450s...)
- ✓ Off-targets







ADME-Tox Platforms

Physicochemistry

✓ Solubility

✓ LogP/D

✓ Pka

✓ Plasma stability, binding

Absorption

✓ Permeability (Caco2, PAMPA)

✓ In vivo (P.o.; i.v. etc..)

Distribution

✓ In vivo (mice, rat)

✓ All tissues

✓ Blood-Brain barrier

Metabolism

✓ Metabolic stability (Hepatocytes, microsomes)

✓ Blood partitioning

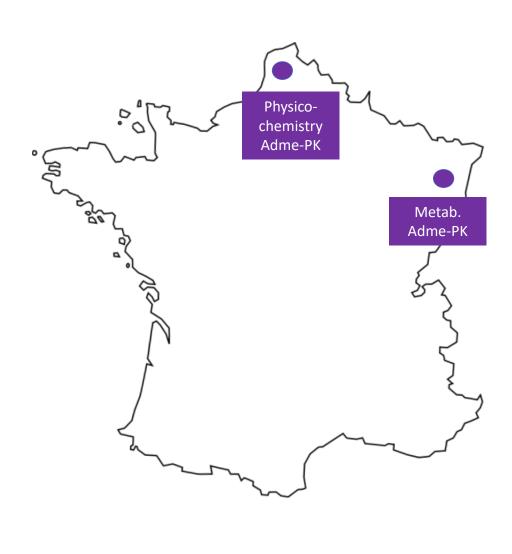
✓ Hemolysis

Pharmacokinetic

✓ In vivo

✓ Kidney clearance

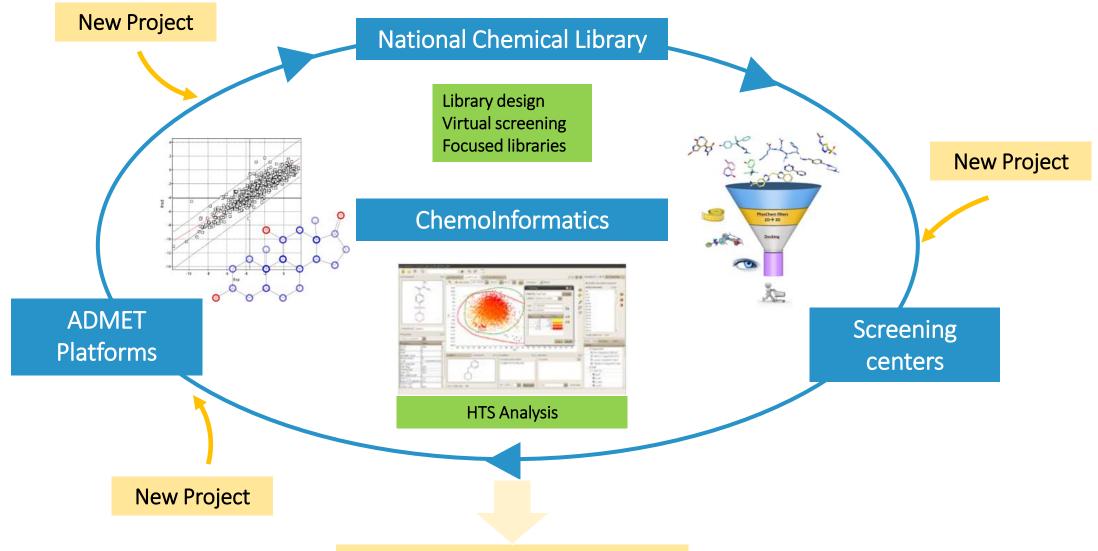
✓ Secondary metabolites











Drug Development (pre-clinical & clinical studies)

https://chembiofrance.cn.cnrs.fr/fr/projet



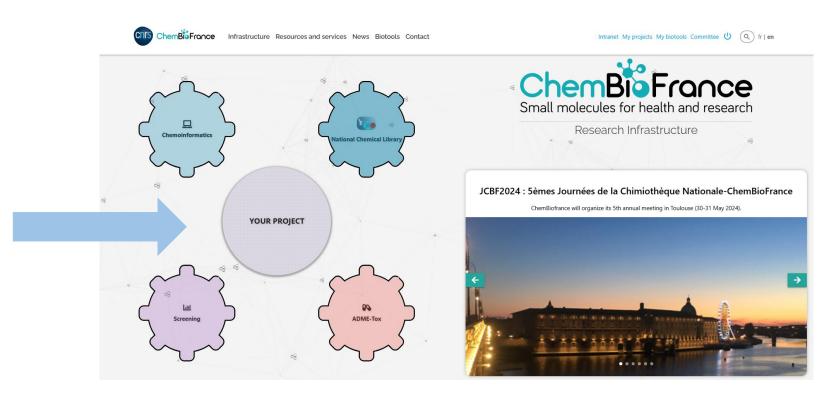
Projects processing



Kindel media

I have a project :

I need a chemical library
I need a screening platform



https://chembiofrance.cn.cnrs.fr/fr



Projects processing



I have a project:

I need a chemical library
I need a screening platform

- 1- First, you need to register
- 2- Submit your project



https://chembiofrance.cn.cnrs.fr/en

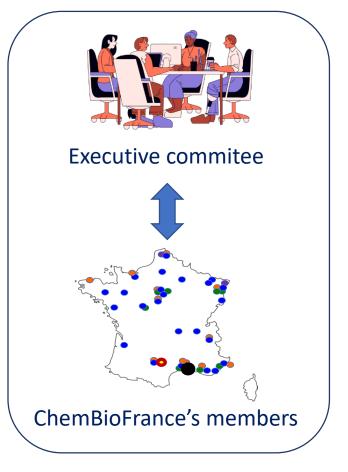


Evaluation



Projects processing

Evaluation









Kindel media

Response within 15 days





Business model



- Fee-for-Service
- Collaborative research program



Kindel media

Cost

Chemical compounds and natural extracts

 Auditable pricing (+access fee for private company)

Chemoinformatics

• 50€/hr

Screening

Project-specific quote

ADME-Tox

Project-specific quote





Thank You

Contact Us: mikael.le-clech@cnrs.fr

