









USE ProtoPRED NOW!

- Save \$\$\$
- Very easy to use
- Pay-per-use or annual fee
- Ready-to-file regulatory results (e.g. REACH Annexes)

DON'T MISS OUT!
BEGIN USING ProtoPRED NOW AND
ENJOY OUR RELEASE PROMOTION!



Go to protoQSAR.com/protoPRED



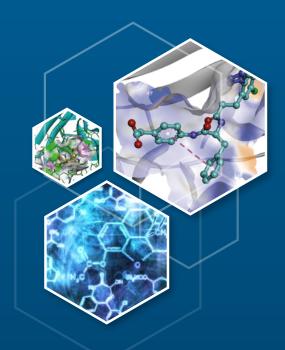
Fill out a quick form



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COMPUTATIONAL TOXICOLOGY:

FAST, ECONOMICAL & ETHICAL

ABOUT US



Company founded 2012



12 employees



100+ peer-reviewed scientific publications

WHY COMPUTATIONAL TOXICOLOGY?

QSAR and REACH

Computational toxicology allows **FAST**, **ECONOMICAL** and **ETHICAL TOXICOLOGICAL EVALUATIONS**, complying with all regulations of registry and authorization of chemicals (e.g. EU's REACH).

Computational toxicology allows you to avoid most animal tests, thus **SAVING TIME & MONEY** and complying with the **ETHICAL MANDATE** to limit animal testing (3 Rs).



FAST RESULTS

Easy and immediate application to thousands of chemicals





ETHICAL

Animal testing is replaced by our models (3 Rs). *In vivo* research to be performed only as a last resort, as recommended by regulators.

VALIDITY

OECD, ECHA, EMA, EFSA, FDA, ICH guidance on pharmaceutical impurities, European directives for registration and labeling of compounds (REACH, CLP, cosmetics, biocides, etc.)...

ProtoPRED

ProtoPRED is a a virtual lab replacing most chemical experiments and biological tests (in vitro / in vivo).

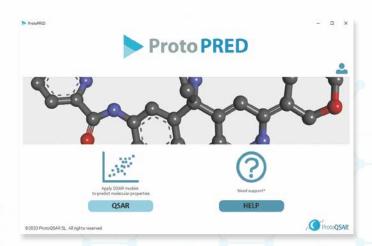
WHY ProtoPRED?

ProtoPRED goes beyond its computational toxicology competitors by providing:

- ♣ LARGER LIST OF PARAMETERS (physico-chemical, toxicological and ecotoxicological) of interest. (e.g. we cover all REACH Annexes)
- LARGE DOMAIN OF APPLICABILITY (DA):
 most chemicals can be evaluated
- REGULATORY APPLICABILITY guaranteed, models developed following the OECD & ECHA rules
- Ready for IMMEDIATE, USER-FRIENDLY USE of the tool

READY-TO-FILE RESULTS

QMRF & QPRF dossiers required by regulators provided required by regulators)



Protopred Can be used in Different areas



pharmacy



agrochemical



veterinary



functional food



cosmetic



many more!

OTHER WAYS TO COLLABORATE



CONSULTING CHEMOINFORMATICS & MOLECULAR MODELING

- Drug discovery
- Agrochemicals with lower environmental impact
- Nanomaterials risk assessment And many more!



PREDICTIVE MODELS FOR REGULATORY PURPOSES.

- Physico-chemical parameters
- Toxicity in humans
- Ecotoxicity



COLLABORATIVE PROJECTS

ProtoQSAR participates in different international research initiatives and is always open to listen to project proposals!