



PROGETTI LIFE PER LO SVILUPPO DI METODI INNOVATIVI NELL'AMBITO DEL REGOLAMENTO REACH

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www.vegahub.eu

Do you need assistance for a property prediction?
CONTACT US

Welcome to the VEGA HUB

Offering a family of tools to evaluate chemical hazard: VEGA, ToxRead, ToxWeight, ToxDelta, and JANUS.

VEGA is the QSAR software with tens of models for individual properties.

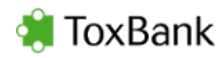
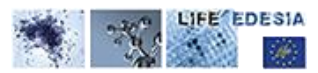
VEGA HUB - QSAR - Download -

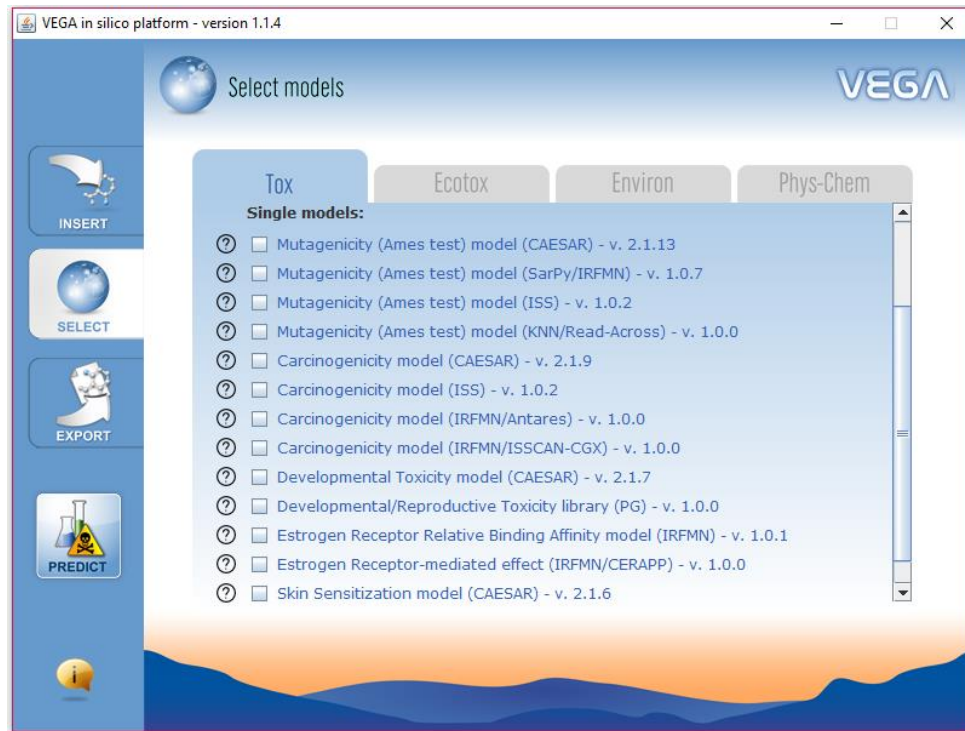
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The *in silico* methods can be very useful, if correctly applied

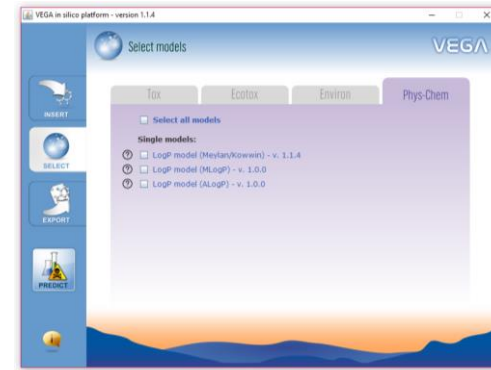
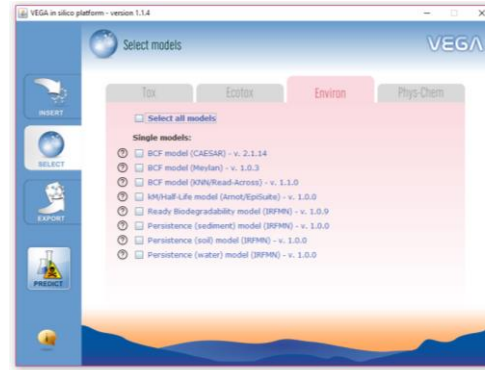
- We want to enhance**
We want to enhance the capability of the *in silico* methods to assess the properties of chemical substances.
- We sustain**
We sustain the correct use of *in silico* methods.
- We provide**
We provide computer tools to support the human experts.
- We support**
We support the evaluation of chemicals safety.

- ✓ VEGA
- ✓ ToxRead
- ✓ ToxWeight
- ✓ ToxDelta





- ✓ 15 modelli per endpoint riferiti a tossicità umana
- ✓ 7 modelli per ecotossicologia
- ✓ 8 modelli per proprietà ambientali
- ✓ 3 modelli per proprietà chimico fisiche

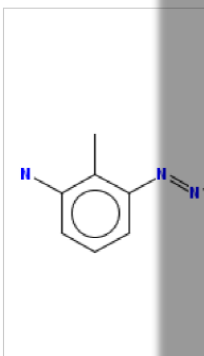


VEGA informazioni nell'output

VEGA

1. Prediction

Prediction for compound



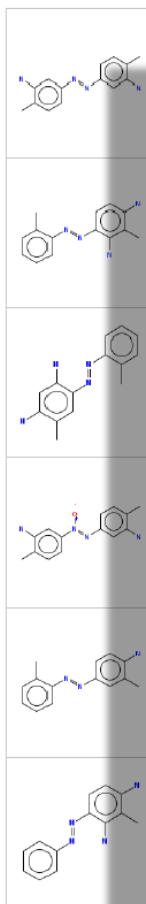
Compound: Molecule 0
 Compound SMILES: Cc1cccc(N=N)c1
 Experimental value: -
 Predicted Mutagen activity: -
 Structural alerts: SA29
 Reliability: the prediction is reliable
 Remarks:
 none

3.1 Applicability Domain:

Similar Compounds, with Predicted and Experimental Values



ge 1



3.2 Applicability Domain:

Measured Applicability Domain Scores



✓	Global AD AD index = 1.0 Explanation: All features are present in the training set.
✓	Similar molecules Similarity in training set: 0.977 Explanation: Similar compounds are found in the training set.
✓	Accuracy Accuracy in training set: 1.0 Explanation: All training set compounds are correctly classified.
✓	Concordance Concordance in training set: 1.0 Explanation: All training set compounds have the same predicted value.
✓	Model's descriptor Descriptor: SA29 Explanation: The descriptor is present in the training set.
✓	Atom Centric ACF index: 1.0 Explanation: All atoms are present in the training set.

Symbols explanation:

- ✓ The feature is present in the training set.
- ⚠ The feature is not present in the training set.
- ✗ The feature is not present in the training set.

4.1 Reasoning:

Relevant Chemical Fragments and Moieties



(Molecule 0) Reasoning on fragments/structural alerts:

Fragment found: SA29 Aromatic diazo

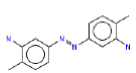


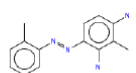
Ar = Any aromatic/heteroaromatic ring

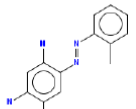
- Chemicals with a sulfonic acid group (-SO₃H) on both rings linked to the diazo group are excluded.

Aromatic diazo. If a sulfonic acid group (-SO₃H) is present on each of the rings that contain the diazo group, the substance should be not classified.

Following, the most similar compounds from the model's dataset having the same fragment.

 CAS: 119631-00-4
 Dataset id: 3482 (Training set)
 SMILES: N(=Nc1ccc(c(N)c1)C)c2ccc(c(N)c2)C
 Similarity: 0.977
 Experimental value: Mutagenic
 Predicted value: Mutagenic

 CAS: 84434-45-7
 Dataset id: 2696 (Test set)
 SMILES: N(=Nc1ccccc1C)c2ccc(N)c(c2(N))C
 Similarity: 0.951
 Experimental value: Mutagenic
 Predicted value: Mutagenic

 CAS: 7467-29-0
 Dataset id: 3021 (Training set)
 SMILES: N(=Nc1cc(c(N)cc1(N))C)c2cccc2C
 Similarity: 0.949
 Experimental value: Mutagenic
 Predicted value: Mutagenic

Uso di VEGA

ECHA

- Come software per esempio nella guida pratica “How to use and report (Q)SARs 3.1” (insieme al software dell’US-EPA EPISuite)
- Usato per “Preparation of the Annex III inventory - Technical Documentation” (insieme al Danish QSAR Database e all’OECD Toolbox)

EFSA

- Per sviluppare nuovi modelli usando i dati disponibili nel database OpenFoodTox
- Per colmare i data gap nel database OpenFoodTox usando i valori predetti da VEGA

CEFIC

- Incluso nel software AMBIT sponsorizzato dal CEFIC

Agenzia per l’ambiente tedesca

- Usato per la prioritizzazione delle sostanze (in PROMETHEUS e JANUS)

EPA danese

- VEGA incluso nel Danish QSAR database

https://echa.europa.eu/documents/10162/13655/pg_report_qsars_en.pdf

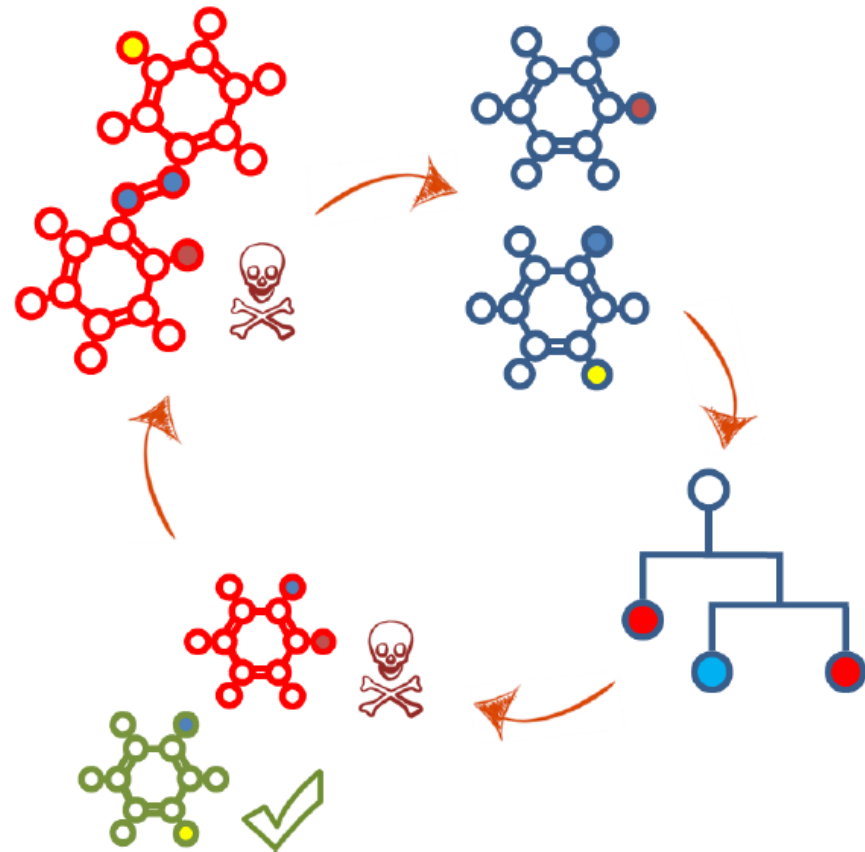
https://echa.europa.eu/documents/10162/22332820/annex_iii_preparation_inventory_en.pdf

PROSIL: coloranti azoici

✓ Sviluppo di modelli specifici per i coloranti

✓ Mutagenicità di coloranti azoici studiata valutando la capacità di generare amine aromatiche

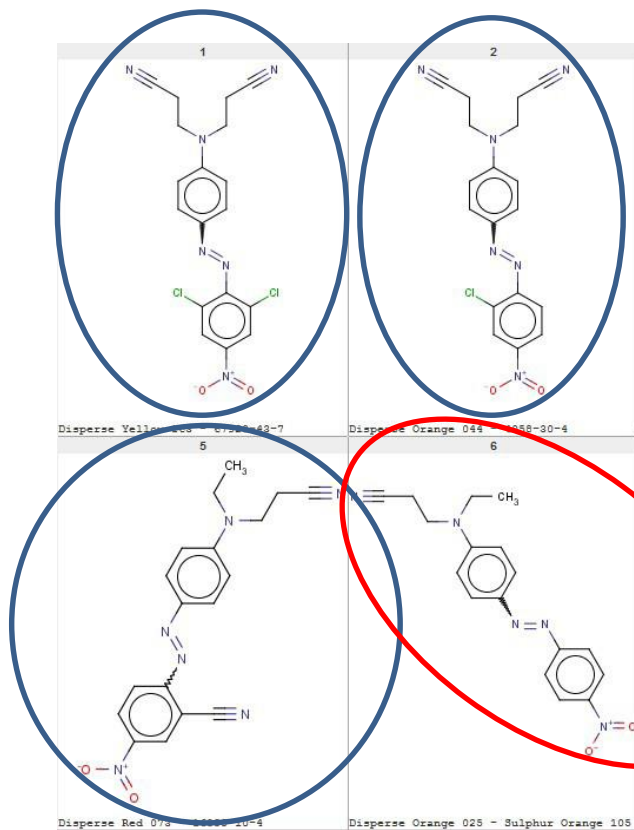
✓ Combinazione finale di 5 diversi approcci per valutare la tossicità di questa famiglia di composti



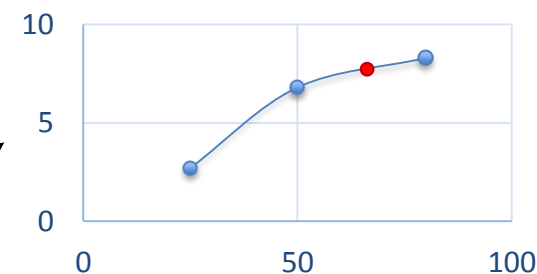
PROSIL approccio a famiglie



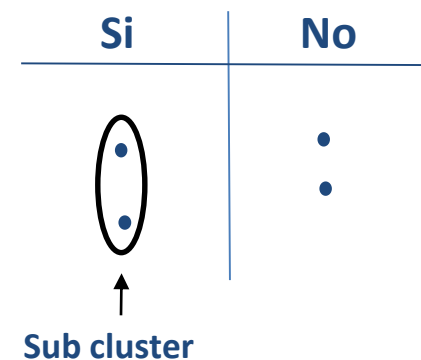
Esempio di andamento delle proprietà tossicologiche all'interno di un cluster



Dati in continuo



Dati in classificazione



Computational tool for the assessment and substitution of biocidal active substances of ecotoxicological concern

Identificazione e sostituzione di biocidi pericolosi dal punto di vista ecotossicologico, considerando l'intero ciclo vitale, dalla produzione allo smaltimento:

- **metaboliti**
- **prodotti di degradazione**
- **prodotti di trasformazione**

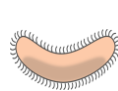


informazioni di supporto
basate sull'evidenza
(EBDSS)

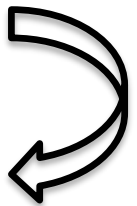


Metodi *In silico*

implementazione

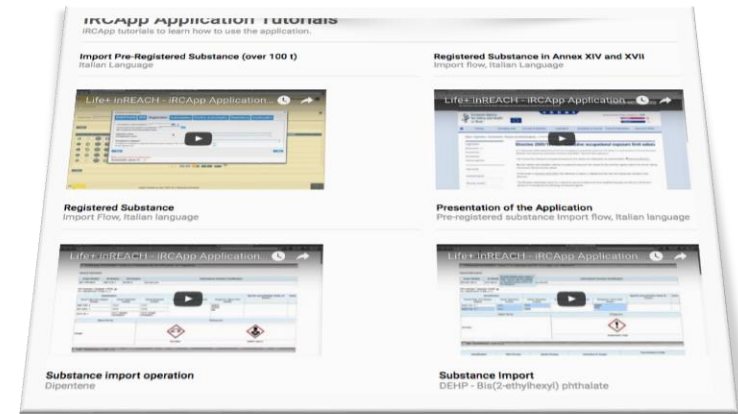


VEGA
toxRead

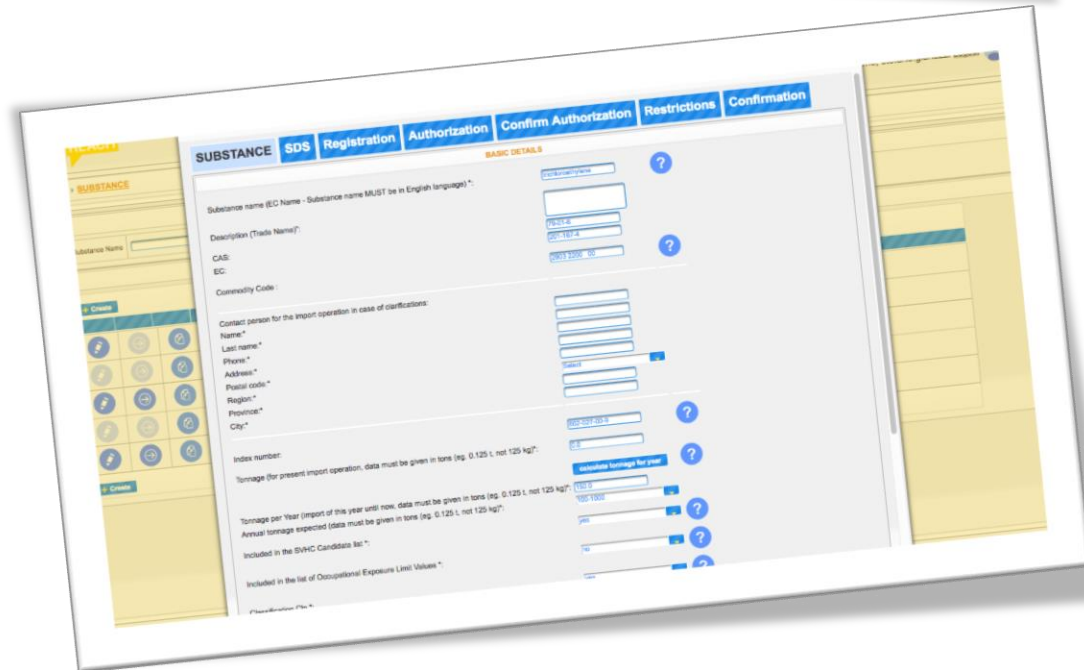


Sviluppo di un applicazione disponibile in rete

Strumento per aiutare gli stakeholder a soddisfare gli adempimenti di REACH e CLP a livello delle importazioni



Applicazione IRCApp



SUBSTANCE SDS Registration Authorization Confirm Authorization Restrictions Confirmation

BASIC DETAILS

Substance name (EC Name - Substance name MUST be in English language): *

Description (Trade Name): *

CAS: *

EC: *

Community Code: *

Contact person for the import operation in case of clarifications:

Name: *

Last name: *

Phone: *

Address: *

Postal code: *

Region: *

Province: *

City: *

Index number: *

Tonnage (for present import operation, data must be given in tons (eg. 0.125 L, not 125 kg)): *

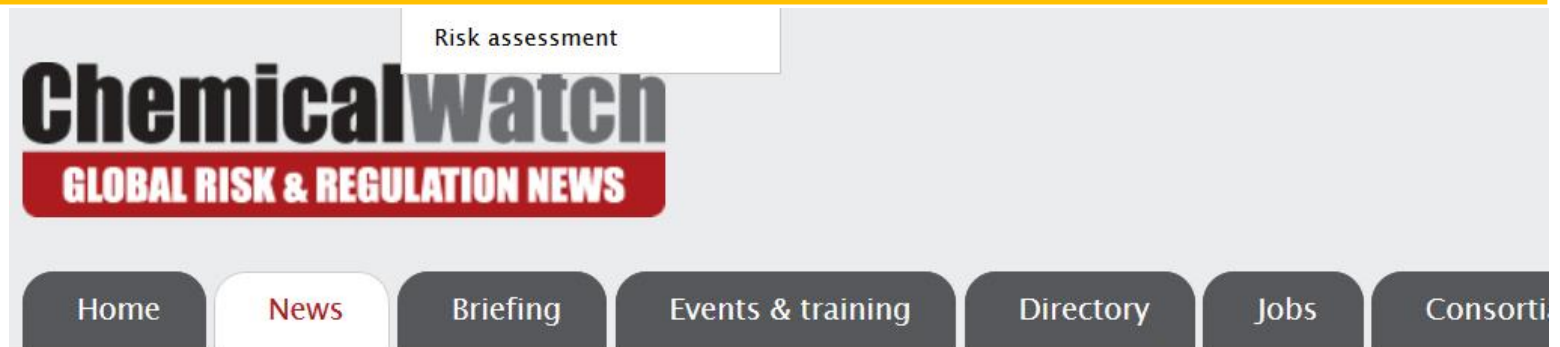
Tonnage per Year (import of this year until now, data must be given in tons (eg. 0.125 L, not 125 kg)): *

Annual tonnage expected (data must be given in tons (eg. 0.125 L, not 125 kg)): *

Included in the SVHC Candidate list: *

Included in the list of Occupational Exposure Limit Substances: *

PROMETHEUS



Germany's UBA publishes screening methodology for prioritising PBTs

Italian institute successfully trials platform

7 April 2016 / Germany, Italy, Risk assessment



Dr Emma
Davies
Reporter

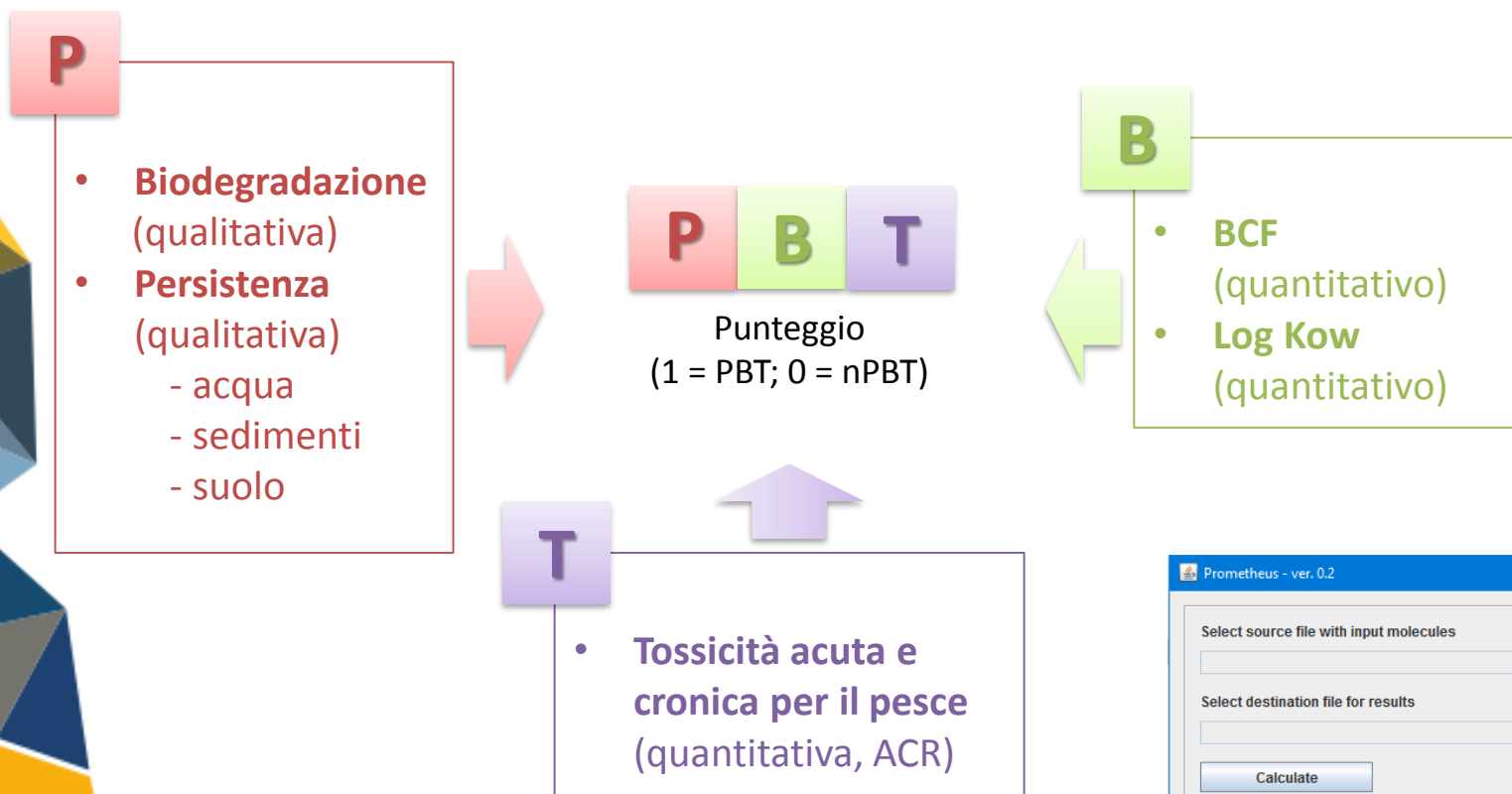
Italian researchers have created a platform for prioritising possible persistent, bioaccumulative and toxic (PBT) chemicals for further assessment.

Working on behalf of Germany's Federal Environment Agency (UBA), a team from the Mario Negri Institute for Pharmacological Research in Milan has brought together a battery of *in silico* tests to screen for chemicals that may be "of concern for the environment and human health".

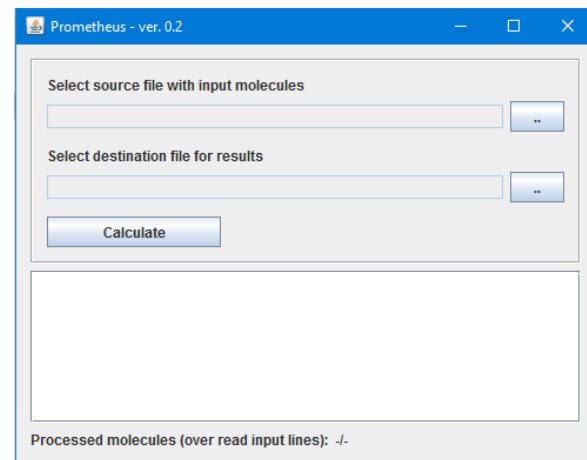


PROMETHEUS

PRioritization Of chemicals: a METHodology
Embracing PBT parameters into a Unified Strategy

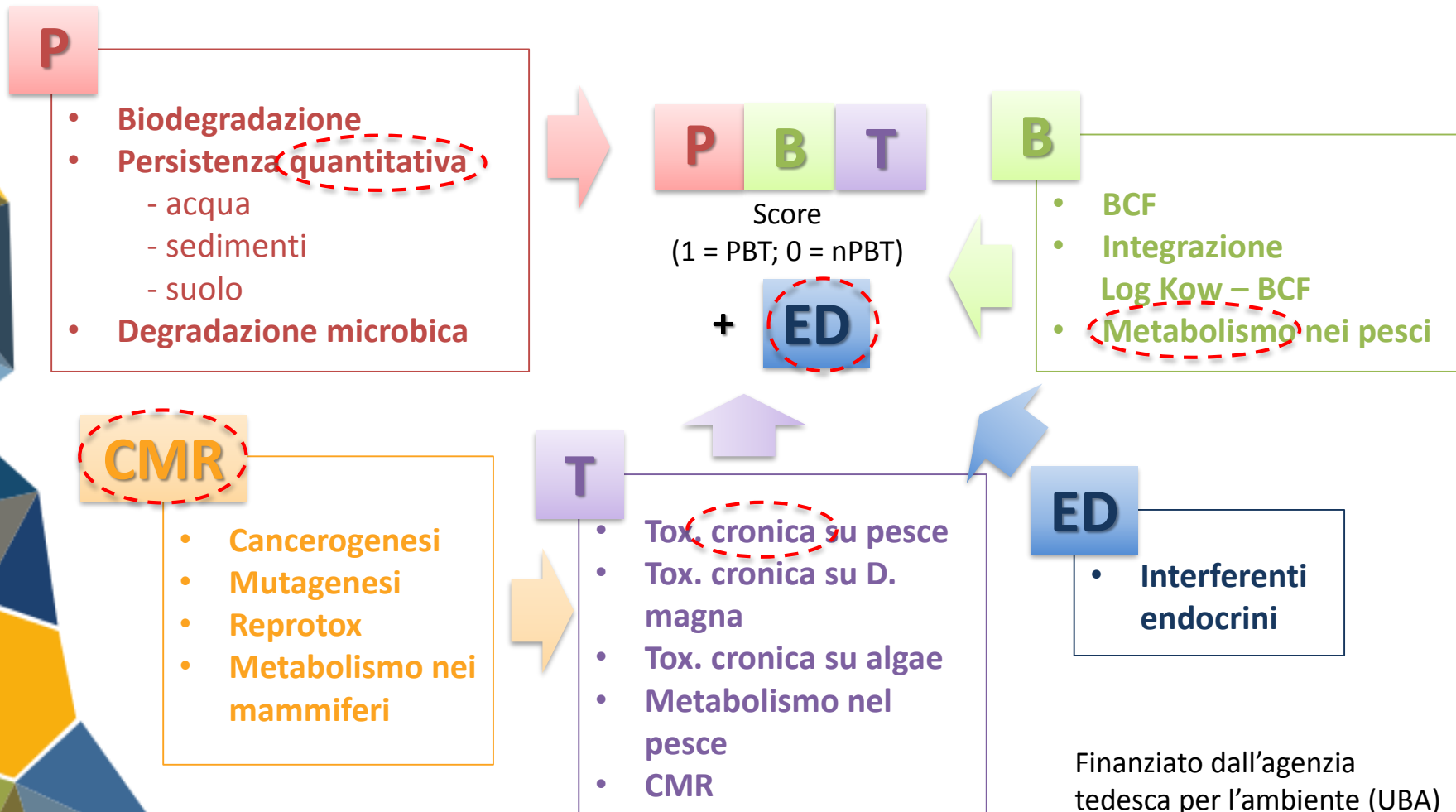


Finanziato dall'agenzia
tedesca per l'ambiente (UBA)



JANUS

Joining environmental, ecotoxicological and toxicological Assessment of chemical substances with Non-testing methods within a Unified Screening

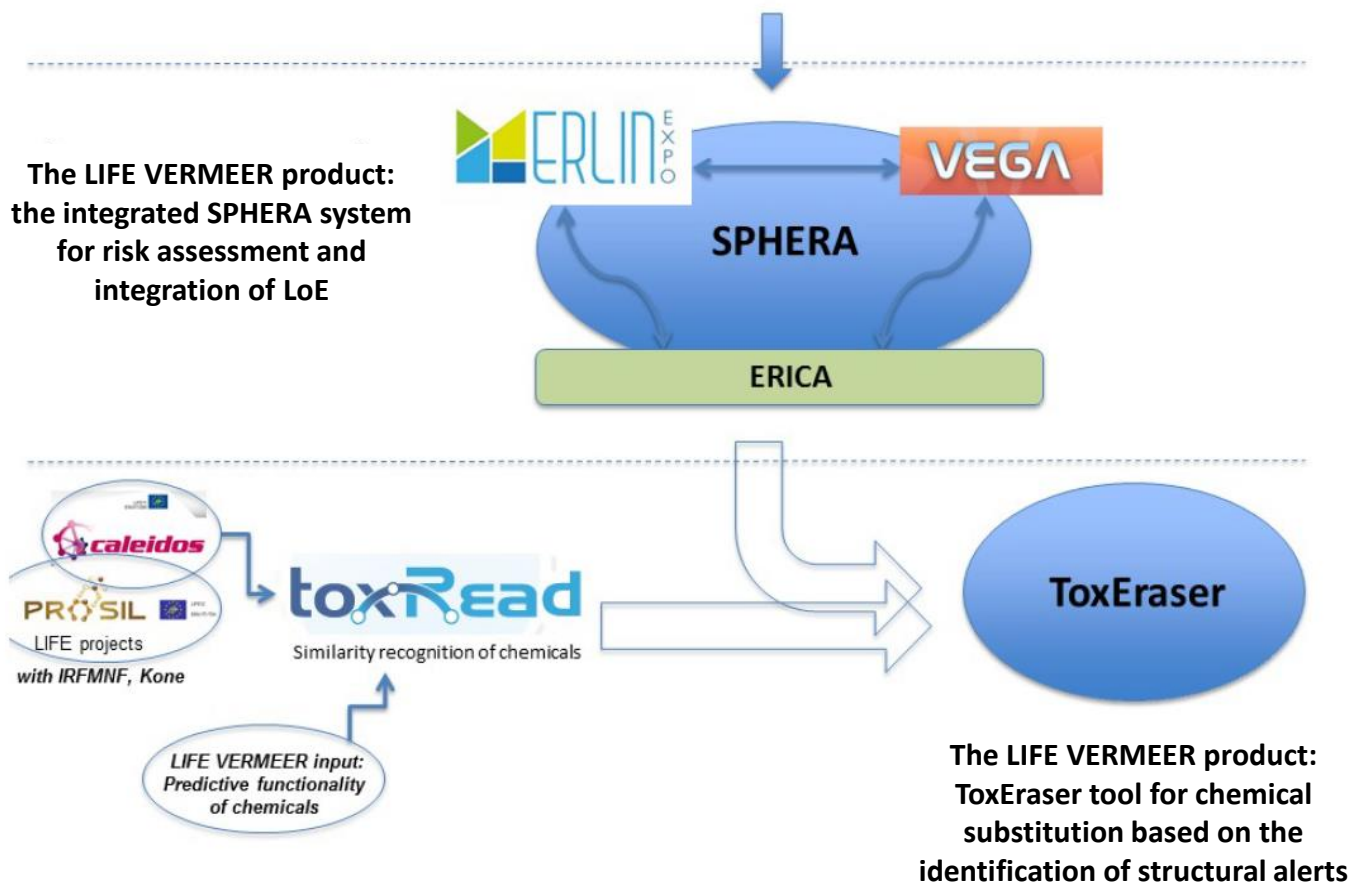


Finanziato dall'agenzia tedesca per l'ambiente (UBA)

VERMEER



Integrazione di VEGA, ToxRead, MERLIN-Expo e ERICA in una piattaforma per la valutazione del rischio e la **sostituzione** di sostanze problematiche



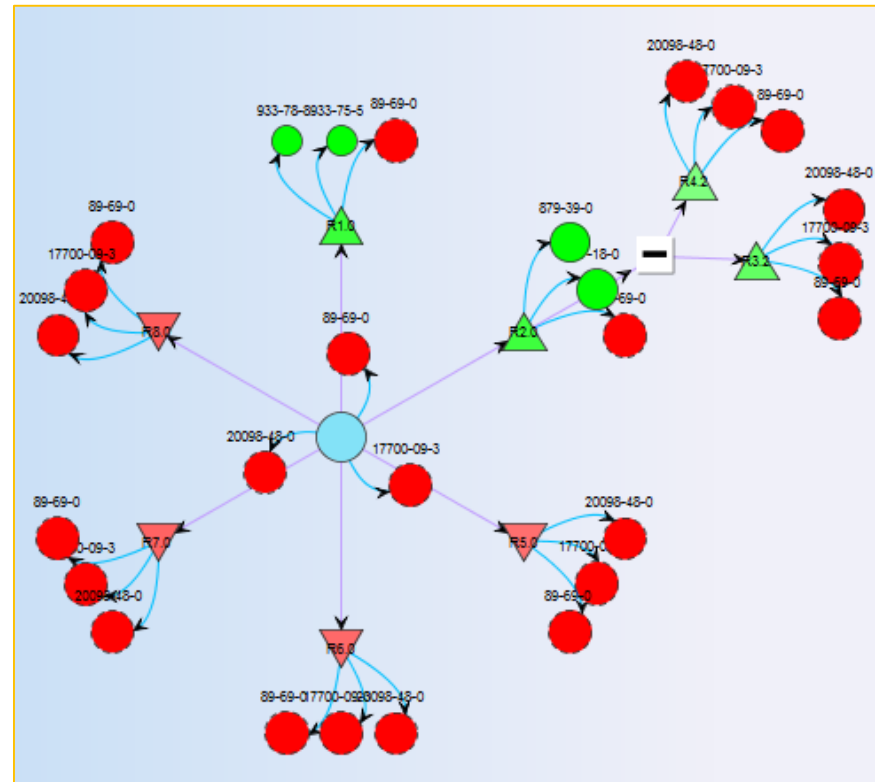


GRAZIE!

toxRead user interface (mutagenicity)

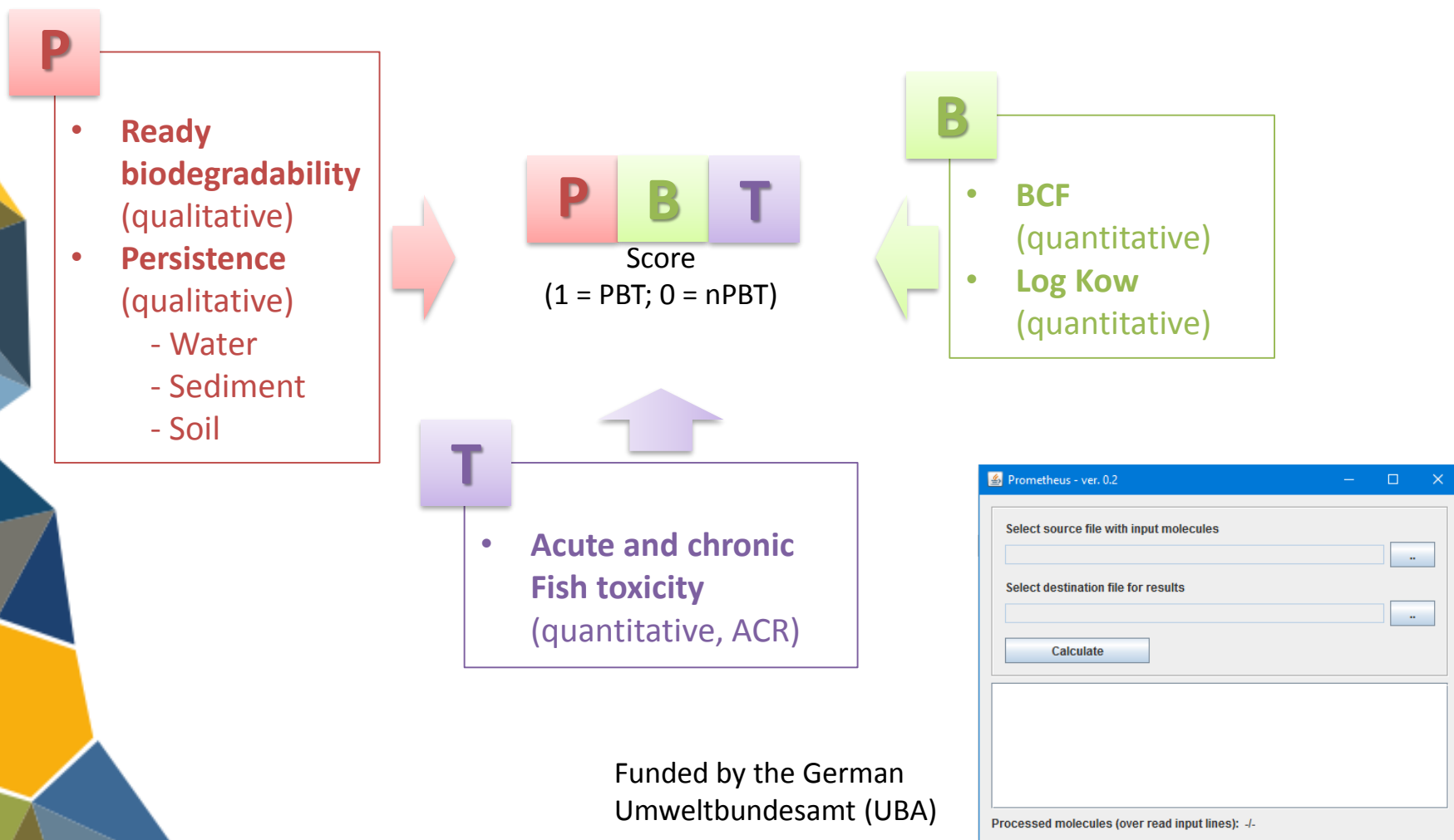
- Shape: circles are molecules, triangles are structural alerts
 - Target molecule in the center of the visualization panel
 - Target compound directly connected to most similar molecules (in inner circle)

- Circle dimension: related to similarity
- Paths connect molecules sharing the same structural alert
- Color: red or green with different saturation indicates active or non active at different levels
 - Clicking on nodes shows structure, explanation, etc.



PROMETHEUS

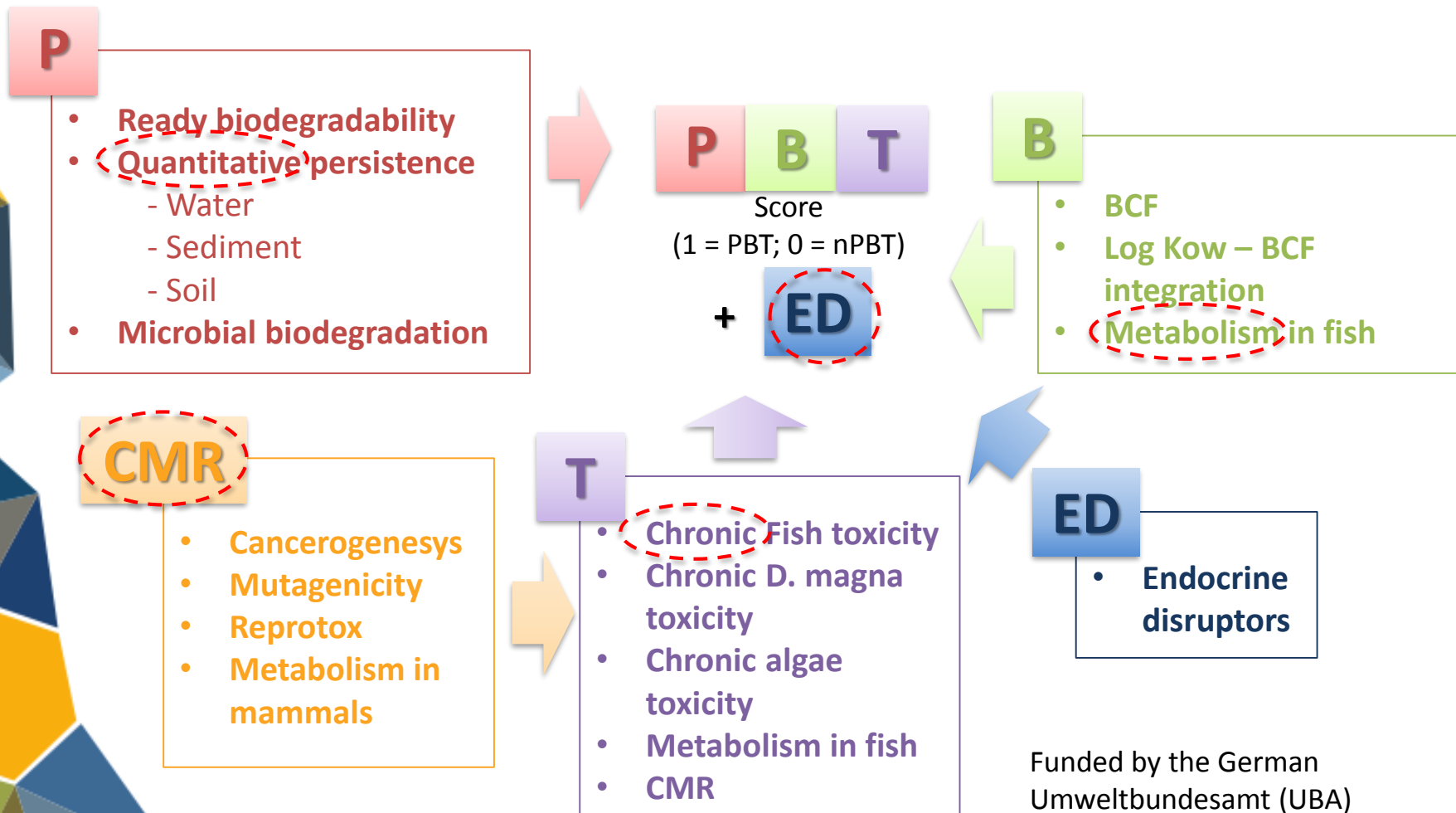
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