User manual

Toolbox 4.4 Release Notes

# **Document history**

Version	Comment
Version 1.0	February 2020: Toolbox 4.4 Release Notes

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If you have questions or comments that relate to this document, please send them to <a href="mailto:ehscont@oecd.org">ehscont@oecd.org</a> or visit the QSAR Toolbox discussion forum at <a href="https://community.oecd.org/community/toolbox">https://community.oecd.org/community/toolbox</a> forum

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## 1 Overview

The Toolbox 4.4 installation is a major update of Toolbox 4.3. It can be installed as a separate product alongside previous major releases of Toolbox (4.3, 4.2, 4.1, 4.0, 3.4, 3.3, etc.)

# 2 System Requirements

HDD: 20 GB free hard drive space

Microsoft .NET Framework 4.5.1

File system: NTFS

# 3 Change log

#### I. General

- New Simplified Toolbox interface
- Development of a Toolbox web repository where different modules which could be docked to Toolbox are stored
- Development of Toolbox repository client within the QSAR Toolbox which manages the docking of external modules via direct connection to web repository or uploading from the local computer
- Possibility to "unlock" the ECHA REACH database export in Toolbox which allows exporting of data after accepting the terms and condition for usage.
- One working mode only 2.5D mode available

## II. Input

- Define target endpoint:
  - ECHA REACH endpoints could be defined
  - New options for easier definition of the target endpoint "Define here",
    Copy-Paste defined target endpoint.
- Loading of list with SMILES (or CAS) shows in a window all substances corresponding to the entered identifier

# III. Profiling

**71 profilers** (Predefined: 5; General Mechanistic: 29; Endpoint Specific: 25; Empiric: 9; Toxicological: 1; Custom: 2) and 16 metabolisms (5 observed and 11 simulated) are available.

• New "QSAR domain query" in the profiling editor.

#### A. New profilers

1. Custom profile: Skin sensitization for DASS

### B. Updated profilers

1. Acute Oral Toxicity

- 2. Biodeg Probability (Biowin 5)
- 3. Biodeg Probability (Biowin 6)
- 4. Carcinogenicity (genotox and nongenotox) alerts by ISS
- 5. DART scheme
- 6. DNA binding by OASIS
- 7. DNA alerts for AMES, CA and MNT by OASIS
- 8. Hydrolysis Half life (pH 6.5 7.0)
- 9. in vitro mutagenicity (Ames test) alerts by ISS
- 10. in vivo mutagenicity (Micronucleus) alerts by ISS
- 11. Organic Functional groups
- 12. Organic Functional groups (nested)
- 13. Organic functional groups (US EPA)
- 14. Protein binding alerts for Chromosomal aberration by OASIS
- 15. Protein binding alerts for skin sensitization by OASIS
- 16. Protein binding by OASIS
- 17. Protein binding potency by GHS
- 18. Substance type

#### C. Updated simulators

- 1. Autoxidation simulator
- 2. Hydrolysis simulator (neutral)
- 3. in vivo Rat metabolism
- 4. Microbial metabolism simulator
- 5. Rat liver S9 metabolism simulator
- 6. Skin metabolism simulator

#### IV. Data

#### 57 databases with 92 134 substances and 2 634 458 data points are available

• IUCLID – new options when import/export data

#### A. Updated databases

- 1. Aquatic Japan MoE
  - Available: 664 substances and 4 577 data points
  - New: 5 substances and 14 data points.
- 2. ECHA REACH (formerly ECHA CHEM)

- Available: 13 305 substances and 802 230 data points
- New: 1 562 substances and 133 926 data points

#### 3. ECOTOX

- Available: 11 822 substances and 969 352 data points
- New: 167 substances and 52 306 data points
- 4. Genotoxicity OASIS
  - Available: 8 031 substances and 30 943 data points
  - New: 46 substances and 496 data points
- 5. pKa OASIS
  - Available: 2 898 substances and 3 492 data points
  - New: 968 substances and 1 183 data points
- 6. Repeated Dose Toxicity HESS
  - Available: 745 substances and 485 842 data points
  - New: 45 substances and 45 446 data points
- 7. Toxicity Japan MHLW
  - Available: 390 substances and 3 971 data points
  - New: 138 substances and 1 057 data points

#### **B.** Updated inventories

- 1. NICNAS
  - Available: 40 180 substances
  - New: 486 substances

## V. Category definition

- New clustering options
- Option to sort chemicals when defined a category based on their structural similarity to the target

## VI. Data Gap Filling

- New AW for Skin sensitization for Defined approaches purposes
- Implementation of the latest EPISUITE models versions

## VII. Reporting/Exporting

- New section with all options for export in the Report module
- New template of exporting profiling result when applying export of the data matrix
- New RAAF templates for environmental and ecotoxicity endpoints
- Addition of QSAR predictions in the Data matrix report.

### VIII. IT improvements

- Data matrix responsiveness improvement
- Database deployment consistency checking
- Reduced memory use on server
- Acceleration of:
  - Starting the program
  - Clustering
  - Automated workflows
  - Profiling

### IX. Additional new features

- Providing numbers of chemicals, available data and predictions for each list in the document tree
- Option for removing the chemicals without experimental data on the data matrix
- Option for searching information in the data matrix
- Freezing target chemical column
- Possibility to edit an entered substance with composition
- Highlights the row of the document tree where a prediction is accepted; sections in the report containing RAAF assessment elements.
- Improved biological taxonomy
- Caching of generated metabolites in Toolbox database (new combination of metabolic simulators and databases are cached)
- Other small improvements

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