

LINKING ALGAL GROWTH INHIBITION TO CHEMICAL ACTIVITY

Stine N. Schmidt^{1*}, James M. Amitage², Jon Arnot², Kresten Ole Kusk¹ and Philipp Mayer¹ ¹Department of Environmental Engineering, Technical University of Denmark, Kgs. Lyngby, Denmark and ²ARC Arnot Research & Consulting Inc., Toronto, Canada. *e-mail: stnsch@env.dtu.dk

PART 1: Challenging the chemical activity range for baseline toxicity

Background

- Recently, high-quality data were published on the algal growth inhibition caused by 50 non-polar narcotic compounds, of which 39 were liquids with defined water solubility (Figure 1a, [1]).
- In the present study [2], the toxicity data for these liquids were applied to challenge the chemical activity range for baseline toxicity [3-6].
- Chemical activity (a, unitless) quantifies the energetic level of an organic compound relative to its pure liquid [0-1], and several studies have reported that baseline toxicity requires a chemical activity of at least 0.01-0.1 [e.g., 3-9].

Objectives

- To convert toxicity data for liquids to chemical activity.
- To challenge the proposed chemical activity range of 0.01-0.1 for baseline toxicity.
- To propose a framework for using the obtained findings.

PART 2: Extending the utilisation of the chemical activity concept

Background

- As for the above data [1], aquatic toxicity data are most often expressed on a concentration basis (e.g., EC_{50} in mg L⁻¹).
- Whereas the data are useful within regulatory risk assessment, the actual EC₅₀ values offer no direct information on whether the compounds exert baseline toxicity or excess toxicity.
- In the present study, algal growth inhibition data were plotted relative to a regression for the (sub-cooled) liquid solubility, which served as a visual reference for chemical activity of unity.
- The data came from three comprehensive and carefully conducted algal toxicity experiments [1, 10, 11].

Objectives

- To express aquatic toxicity data on a chemical activity basis.
- To identify and quantify excess toxicity and thereby compounds of concern.



[1] Aruoja et al. 2014. *Chemosphere* 96: 23-32.



Figure 1. (a) Quantitative structure-activity relationship (QSAR) relating effective concentrations (EC₅₀) and octanol to water partition ratios (K_{ow}) for 50 non-polar narcotic compounds [1]. Water miscible and solid compounds (n=11) are indicated by light circles. (b) Ratios of EC₅₀ (mg L⁻¹) and water solubility (S_{water}, mg L⁻¹) for the remaining 39 liquids plotted against K_{ow}.



[2] Schmidt and Mayer. 2015. Chemosphere 120: 305-308

Reichenberg and Mayer. 2006. Environ. Toxicol. Chem. 25: 1239-1245.

[4] Mayer and Holmstrup. 2008. Environ. Sci. Technol. 42: 7516-7521.

- 5] Mackay et al. 2009. SAR. QSAR. Environ. Res. 20: 393-414.
- 6] Mackay et al. 2014. SAR. QSAR. Environ. Res. 25: 343-355.
- [7] Smith et al. 2010. Aquat. Toxicol. 98: 15-24.
- [8] Mackay et al. 2011. Integr. Environ. Assess. Manag. 7: 248-255.

Procedure

- 1b, [2]).

Results

- limit would indicate excess toxicity.
- the CEFIC LRI-ECO30 project.

Procedure

- of 1 (Figure 2).
- in the same chart (Figure 2).

Preliminary Results

- [9] Lee et al. 2013. Ecotoxicol. Environ. Saf. 94: 116-122.
- [10] Aruoja et al. 2011. Chemosphere 84: 1310-1320
- [11] Kusk et al. To be submitted.
- [12] Mackay et al. 1980. *Chemosphere* 9: 701-711.



• Ratios of effective concentration (EC₅₀, mg L⁻¹) and water solubility (S_{water}, mg L⁻¹) were determined, which essentially equals the effective chemical activity (Ea_{50}).

• The EC_{50}/S_{water} ratios were plotted as a function of K_{ow} (Figure

• Most EC_{50}/S_{water} ratios were within the expected chemical activity range of 0.01-0.1, and none were significantly below 0.01.

• These findings suggest EC_{50} values for baseline toxicity to be at or above 1% of liquid saturation, and EC_{50} values well below this

• Methods for applying the chemical activity concept for ecological risk assessment and environmental quality guidelines for baseline (mixture) toxicity are being proposed [2] and refined in

• A regression of sub-cooled liquid solubility (S₁, mmol L⁻¹) [12] was plotted as a function of K_{ow} , representing a chemical activity

 EC_{50} values (mmol L⁻¹) [1, 10, 11] were plotted against their K_{ow}

• The data analysis confirmed baseline toxicity for 68 compounds, which were characterised as polar and non-polar narcotic compounds by Aruoja and co-workers (Figure 2a).

• The data analysis of 188 compounds reported by Kusk and coworkers revealed 34 compounds exerting toxicity well below the limit of 1% of liquid saturation (a=0.01), and thereby indicating excess toxicity to algae (red circles, Figure 2b).

• In depth interpretation and further analyses on algal toxicity data, covering a wide range of solids and liquids, several expected modes of action and several algal species will be conducted.

