

# STATISTICAL ECO(-TOXICO)LOGY

IMPROVING THE UTILISATION OF DATA FOR  
ECOLOGICAL RISK ASSESSMENT

by

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Submitted Dissertation thesis for the partial fulfillment of the requirements for a

Doctor of Natural Sciences

Fachbereich 7: Natur- und Umweltwissenschaften

Universität Koblenz-Landau

17. November 2016



## THREATS TO FRESHWATER ECOSYSTEMS FROM CHEMICAL POLLUTION

Freshwater ecosystems, like streams, lakes and wetlands, make up only 0.01% of the World's water and cover only 0.8% of Earth's surface (Dudgeon et al., 2006), yet they host an important component of global biodiversity. Freshwaters are a habitat for more than 125,000 species, which represents 10% of global biodiversity and  $\frac{1}{3}$  of all vertebrate species (Balian et al., 2007; Strayer and Dudgeon, 2010) and provide essential services for human well-being (Aylward et al., 2005). Small water bodies are of particular importance, because of their high abundance (Downing et al., 2012), the high biodiversity they host (Davies et al., 2008) and the ecosystem services they provide (Biggs et al., 2016).

The earth is currently experiencing a functional change driven by human activities which are so far-reaching, that a new geological epoch "Anthropocene" has been proposed (Steffen et al., 2011; Waters et al., 2016). Consequently, these changes are also associated with biotic changes: 65% of rivers are currently at threat (Vörösmarty et al., 2010) and freshwaters are experiencing the greatest losses of biodiversity (WWF, 2016). A multitude of stressors contribute to this deterioration of freshwater biodiversity including habitat loss and degradation, overexploitation, invasive species and pollution (Dudgeon et al., 2006; Vörösmarty et al., 2010; WWF, 2016). Studies investigating water pollution have mainly focused on nutrient loading, acidification and pollution by organic loading (Schäfer et al., 2016). However, chemicals have become ubiquitous throughout humankind. Currently, more than 100,000 chemicals are registered and in daily use (Schwarzenbach et al., 2010; Schwarzman and Wilson, 2009). These substances will ultimately end somewhere in the environment.

Despite their potential negative effects on biota and humans and their intentional release, pesticides have been neglected in the past by ecological studies investigating threats to freshwaters (Schäfer et al., 2016) and it is unknown how much they contribute to biodiversity loss (Persson et al., 2013; Rockström et al., 2009). However, recent studies indicated that pollution by pesticides may be a

frequent threat to freshwaters that might have been neglected by ecological studies in the past. Malaj et al. (2014) showed that almost half of European water bodies are at risk from pesticides. In the United States, Stone et al. (2014) showed that 61% of assessed agricultural streams exceed aquatic-life benchmarks. On a global scale, Stehle and Schulz (2015) found that 52.4% of detected insecticide concentrations ( $n = 11,300$ ) exceeded risk thresholds. The high contact with adjacent land and low water volume of small streams make them particularly vulnerable to pesticide pollution (Biggs et al., 2016), however, there is currently a lack of data on pesticide pollution of small streams (Lorenz et al., 2016).

As a reaction to the degradation of freshwaters, several legal frameworks have been established to safeguard and improve the quality of freshwater ecosystems. In the European Union (EU), the Water Framework Directive (WFD) (European Union, 2000) regulates the protection of aquatic ecosystems and commits the member states to achieve a 'good' status of all water bodies. Knowing of the toxicity of pesticides and their intentional release into the environment, also the introduction and use of new pesticides are highly regulated. Sophisticated environmental risk assessment procedures have been developed and are requested by the EU (European Union, 2009) to ensure that the use of pesticides does not cause unacceptable effects to non-target organism, soil, air and water.

## ENVIRONMENTAL RISK ASSESSMENT

Environmental risk assessment (ERA) tries to estimate risks to animals, populations or ecosystems. It investigates if a chemical can be used as intended without causing detrimental impacts to the environment. Moreover, ERA is used as a tool to support decision making under uncertainty (Newman, 2015). Environmental risk is defined as a combination of the severity and the probability of occurrence of a potential adverse effect on the environment (Suter, 2007). Therefore, ERA is based on two components: Effect- and exposure assessment. A combination of both is needed to characterise environmental risks.

Effect assessment characterises the strength of effects using laboratory and semi-field experiments. It establishes relationships between the concentration of a compound and the observed effects. In the European Union a tiered approach with increasing complexity and realism. Lower tier assessment is based on highly standardised single species laboratory experiments, whereas higher tier assessment is refined by testing additional species, extended laboratory experiments or model ecosystem experiments (Brock et al., 2006). To address the

various uncertainties in effect assessment (e.g. experimental variation, variation between species, variation in environmental conditions etc.) the retrieved toxicity values are divided by an assessment factor (AF) between 100 (lower tier assessment) and 2 (higher tier assessment) depending on data quality, which yields to a regulatory acceptable concentration (RAC) (Brock et al., 2006; EFSA, 2013).

Exposure Assessment for freshwaters aims to characterise the probability of an adverse effect by deriving a predicted environmental concentration (PEC) in surface waters and sediments (Newman, 2015). It is mainly based on modelling the fate of chemicals in the environment using computer simulations. In the European Union, the FOCUS models are used (EFSA, 2013; FOCUS, 2001). To calculate PECs these models need many compound specific input parameters like the molecular weight, water solubility, partitioning coefficients and dissipation time. Additionally, information on the application regime and crop type is needed. FOCUS models the concentration within edge-of-field streams of 1 meter width (corresponding a catchment size of approx. 7km<sup>2</sup>, see Figure ??) and 30 cm depth (Erlacher and Wang, 2011). Nevertheless, recent research showed that FOCUS models fail to predict measured field concentrations of pesticides (Knäbel et al., 2014; Knäbel et al., 2012).

The final step in ERA is risk characterisation. It puts together the information gained from effect and exposure assessment. Risk can be expressed in several ways, a quantitative way being the risk quotient approach: If the ratio PEC / RAC exceeds a value of one potential risks cannot be rebutted (EFSA, 2013; Solomon et al., 2000; Suter, 2007). Consequently, pesticides can be authorised only if the risk quotient is below one indicating that harmful effects are unlikely.

## ENVIRONMENTAL MONITORING

Widespread anthropogenic activities and the induced environmental changes have resulted in concerns about the state of the environment and have led to the development of environmental monitoring programs worldwide (Nichols and Williams, 2006). After authorization, pesticides applied on agricultural fields may enter aquatic ecosystems via diffuse sources like spray-drift, surface run-off or drainage (Carter, 2000; Liess et al., 1999; Schulz, 2004; Stehle et al., 2013). These entered pesticides may have ecological effects and worsen the chemical status, acting contrary to the goal of the WFD. For monitoring the progress towards the goal of a 'good' status and for assessment of the chemical status of sur-

face waters the EU WFD established monitoring requirements for all European river basins (European Union, 2000). For chemical monitoring the WFD requires grab sampling and chemical analysis of 21 priority substances (of which 7 are pesticides) every third month and of 24 other pollutants (of which 12 are used as pesticides) every month (European Union, 2013). Additionally, 14 substances (of which 8 are used as pesticides, including all Neonicotinoids) that may pose a significant risk, have an insufficient data basis and are candidates for future priority substances are currently monitored until 2019 (European Union, 2015). Nevertheless, monitoring programs on a national scale might monitor a broader spectrum of chemical substances, e.g. for investigative monitoring. Recent studies indicate that the current sampling and chemical analyses strategy greatly underestimate the pesticide exposure (Moschet et al., 2014; Stehle et al., 2013; Xing et al., 2013).

Environmental monitoring produces humongous amounts of data containing information on pesticide concentrations in the field on a large under many conditions. Therefore, it can be complementary to ERA (Suter, 2007). Moreover, data from long-term monitoring programs can be used to study hypotheses about spatial and temporal dynamics and interactions, that are not evident from short term and short scale studies (Gitzen, 2012) and provide insights modelling approaches. If the environmental risk assessment process captured all relevant sources of risk, no concentrations above the derived RAC should be observable in European rivers. Therefore, monitoring data could be used to provide feedback for ERA after approval (Knauer, 2016). The WFD has its main focus on large water bodies  $>10 \text{ km}^2$  catchmentsize, whereas ERA has its focus on small water bodies  $<10 \text{ km}^2$  (Brock et al., 2006). However, at present little is known on pesticide concentrations in small streams comparable to those assessed in ERA (Biggs et al., 2016; Lorenz et al., 2016). Additionally, monitoring under the WFD is also performed for biological components of freshwaters and a combination with pesticide exposure data might provide valuable insights into large-scale field effects of chemical substances (Schipper et al., 2014).

## STATISTICAL ECOTOXICOLOGY

Environmental effect assessment generates data on ecological effects using experiments. The produced datasets range from small univariate datasets (lower tier assessment) to medium sized multivariate datasets (higher tier assessment). In order to extract usable information for assessment, these datasets are anal-

ysed using statistical techniques and therefore, statistics are crucial for effect assessment (Newman, 2012). Statistical ecotoxicology combines statistics with the specific needs and constraints of ecotoxicology. Ecotoxicologists deal generally with low replicated experiments, making statistical inference difficult (Van Der Hoeven, 1998). For example, a recent analysis of eleven mesocosm studies revealed that the sample sizes for these kind of experiments range between two and five. Statistical ecotoxicology aims to provide solutions to statistical challenges in ecotoxicology (Fox and Landis, 2016a), guidance on experimental designs (Johnson et al., 2015) and tools to integrate big data (Van den Brink et al., 2016). The ultimate goal is to improve the accuracy of ERA.

The relationships between the concentration of a compound and the observed effects are usually analysed using dose-response models, which can be used to derive an effective concentration for  $x\%$  effect ( $EC_x$ ) (Ritz, 2010). Nevertheless, such relationships cannot always be established from experimental data. For example, mesocosm experiments are conducted to characterise effects on whole biological communities. However, because of multivariate responses and potential indirect effects, there is no clear dose-response relationship and no models for this kind of data available. There are also examples where fitting dose-response models is problematic (Green, 2016). In such cases, there is usually a no-observed-effect concentration (NOEC) computed.

The NOEC is the highest tested concentration that does not lead to significant deviation from the control response and therefore relies on null hypothesis significance testing (NHST). However, the use of NOEC as a toxicity measure in environmental effect assessment has been heavily criticised in the past (Chapman et al., 1996; Fox et al., 2012; Fox and Landis, 2016b; Jager, 2012; Laskowski, 1995; Warne and van Dam, 2008). One such critic is the low statistical power for NHST in common ecotoxicological experiments (Van Der Hoeven, 1998). *A priori* power calculations can provide useful guidance for choosing experimental designs (Johnson et al., 2015), but are rarely used by ecotoxicologists (Newman, 2008).

Instead of conducting experiments, toxicity could be also predicted from molecular structures using quantitative structure-activity relationships (QSAR), which are usually calculated using machine-learning techniques (Cortes-Ciriano, 2016; Murrell et al., 2015). Nevertheless, in order to improve and validate these models to give sufficient prediction accuracy more data from experiments is needed (Kühne et al., 2013). Indeed, a large amount of data is available that could be used for effect and exposure assessment. For example, the US EPA

ECOTOX database (U.S. EPA, 2016), the Pesticides Properties Database (Lewis et al., 2016) and ETOX (Umweltbundesamt, 2016) provide toxicity data that could be used for effect assessment. Databases like Physprop (Howard and Meylan, 2016) and PubChem (Kim et al., 2016) provide chemical properties that are needed as input for exposure models. Monitoring data provides information on realised concentrations, could be used for validation of models and retrospective risk assessment. This "big data" can provide new information and opportunities for ERA (Dafforn et al., 2015). However, it needs to be harmonised, linked and easily accessible in order to be used effectively in ERA.



## OBJECTIVES AND OUTLINE OF THE THESIS

The overall goal of this thesis was to contribute to the emerging field of statistical ecotoxicology, environmental risk assessment and environmental monitoring. The main objectives were (i) to scrutinise new methods in statistical ecotoxicology, (ii) explore available monitoring data and (iii) provide tools to deal with big data. Figure 1.1 provides a conceptual overview on ERA and environmental monitoring as outlined in the previous sections, as well as the parts considered in this thesis and the relations between them.

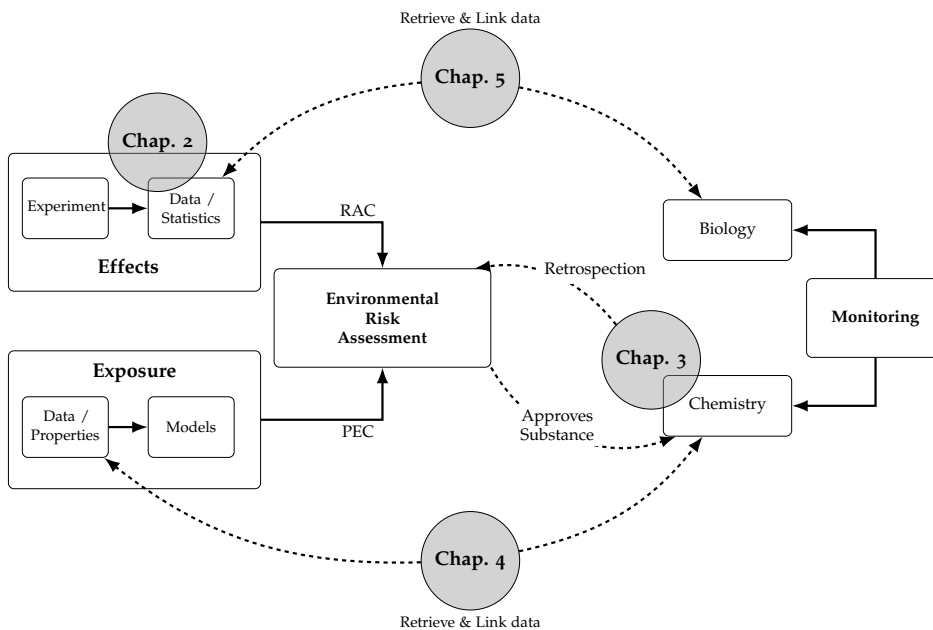


Figure 1.1: Conceptual overview on environmental risk assessment, environmental monitoring and the parts addressed by this thesis.

The thesis starts with a comparison of statistical methods to analyse ecotoxicological experiments using NHST in effect assessment (Chapter ??). Specific questions addressed were:

- Are newer statistical methods, explicitly considering the type of analysed data, more powerful than currently used methods for NHST?
- How much statistical power do current experimental designs in ecotoxicology exhibit?

Risk assessment procedures in the European Union has its main focus on small waterbodies adjacent to agricultural fields where plant protection products are applied. Therefore, chapter ?? focuses on measured large-scale environmental concentrations in small streams, the drivers thereof and comparison with RACs derived from ERA. Specific goals of this study were:

- Compile monitoring data on pesticides in small streams in Germany and check if the available data is suitable to inform ERA.
- Explore the relationship between agricultural land use and stream size and RAC exceedances.
- Scrutinise the annual dynamics of pesticide exposure, as well as the influence of precipitation on measured pesticide concentrations.
- We use RACs derived from ERA to assess the current pollution in German streams and identify pesticides exhibiting currently a risk to freshwaters.

The compilation of monitoring data from different data sources in Chapter ??, resulted in a big inhomogeneous amount of data. Moreover, Biologists, Chemists and ecotoxicologists face similar problems with the need to identify and harmonise their biological and chemical data. Chapters ?? (chemical data) and ?? (biological data) describe software solutions to simplify and accelerate the workflow of:

- validating and harmonising chemical and taxonomic data
- linking datasets from different databases
- retrieving properties and identifiers

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# 2 | GENERAL DISCUSSION AND OUTLOOK

## TOPICS IN STATISTICAL ECOTOXICOLOGY

The simulation study performed in chapter ?? clearly showed that common experimental designs exhibit unacceptably low statistical power (Szöcs and Schäfer, 2016; Van Der Hoeven, 1998). This underpins the criticism accumulated over the last 30 years towards the usage of NOEC as an endpoint for ERA (Fox and Landis, 2016). Nevertheless, the NOEC is still one of the standard endpoints of mesocosm experiments in higher tier risk assessment (EFSA, 2013). Clearly, further advances in the statistical evaluation of mesocosm experiments is needed.

Recently, *a posteriori* calculations of statistical power have been proposed to counteract these limitations and aid the interpretation treatment-related effects in model ecosystems (Brock et al., 2015). The "minimum detectable difference" (MDD) estimates the difference between two means that must exist in order to produce a statistically significant result ( $p < 0.05$  (Gelman and Stern, 2006)) and could be used to interpret NOEC. However, *a posteriori* calculations have been shown to have logical flaws when used for interpretation of non-significant results (Hoenig and Heisey, 2001; Nakagawa and Foster, 2004). However, conducting and reporting of *a priori* power calculations, as performed in chapter ??, might provide researchers important information to optimise their study designs, ensuring that their experimental designs have appropriate power and can lead to interpretable results (Johnson et al., 2015).

Moreover, similar simulations could not only be used to study factorial but also regression designs. Indeed, simulations could be used to determine the optimal experimental design for dose-response models and  $EC_x$  determination, balancing precision and usage of resources. Regression designs are generally more powerful and provide more information than factorial designs (Cottingham et al., 2005). In mesocosm experiments, such designs, assigning the replicates to more tested concentrations, might also provide additional insights. Currently, statistical tools to analyse a community dose-response relationship are not well explored and no equivalent  $EC_{x,mesocosm}$  can be derived. On possibility could be to fit separate dose-response models to each species, leading

to a  $EC_x$  for each species in a mesocosm study. Subsequently, these  $EC_x$  values could be combined and summarised using Species Sensitivity Distributions (Posthuma et al., 2002), providing a hazardous concentration ( $HC_{x,mesocosm}$ ) for  $x$  % of species affected in mesocosms (Maltby et al., 2005). Another possibility would be to use a logistic type of ordination (van den Brink et al., 2003). Reduced-Rank vector generalised linear models (RR-VGLM) could be used to fit such type of models (Yee, 2015; Yee and Hastie, 2003), but they have not been applied in ecotoxicology yet.

In a similar vein, community ecology is currently experiencing a shift towards a new class of multivariate methods, incorporating statistical models for abundances across many taxa simultaneously (ter Braak and Šmilauer, 2015; Warton et al., 2015a; Warton et al., 2015b; Warton et al., 2012). However, these methods have not been applied frequently and their applicability to ecotoxicological data is currently unclear (Szöcs et al., 2015). All these models have in common, that the choice of statistical model is primarily based on the grounds of data properties. In chapter ?? we showed, that using statistical models that fit the type of data analysed, can provide higher statistical power. Simultaneously, Ives (2015) published a study reaching contradictory conclusions (*"For testing the significance of regression coefficients, go ahead and log-transform count data"*). It must be noted, that the simulation designs differed significantly between both studies: We used a low-replicated factorials design, whereas Ives (2015) simulated a well-replicated regression designs with two predictors. We both found that the negative-binomial GLMs were surprisingly prone to Type I errors, although the assumptions of this model closely matched the data. Nevertheless, as we show in chapter ??, the parametric bootstrap might provide a solution to this problem, but is computationally intensive and not widely used. The parametric bootstrap is akin to Bayesian methods (Gelman et al., 2014), which might provide an alternative inference method for inference. The main point, leading Ives (2015) to his conclusions, was that GLM showed undesirable Type I errors in case of correlated predictors, a case not commonly encountered in ecotoxicology and not studied in chapter ?. Recently, the current state-of-the-art was discussed by Warton et al. (2016): i) choose the statistical model based on the grounds of data properties; ii) fix Type I errors using parametric bootstrap or resampling; iii) take mean-variance relationship into account, which is in line with the findings of chapter ?. However, there are still open questions regarding the use of GLMs for count data (see e.g. raised by Prof. John Maindonald, <http://uni-ko-ld.de/fb>). To diagnose issues such as overdispersion

and excess of zeros in count data models new tools like the recently developed “*Rootograms*” (Kleiber and Zeileis, 2016) provide useful additions.

In chapter ?? we applied new statistical modelling techniques that explicitly consider the limit of quantification. The currently most often used methods to deal with such censored data is to omit or substitute non-detects. Censoring is very common when dealing with chemical and ecological datasets, but is rarely taken into account (Fox et al., 2015). Recent examples from ecotoxicology and environmental chemistry show that omission (Hansen et al., 2015), randomization (Goulson, 2015) or substitution by a fixed value (Helsel, 2010; Helsel, 2006) can lead to biased results. Hansen et al. (2015) used a Tobit regression (Tobin, 1958) that takes the amount of censored data into account, assuming a (log-) normal distribution of concentrations. In chapter ?? we used a slightly different approach, using a zero adjusted gamma model (ZAGA). We modelled measured concentrations as two separate processes, generating i) zero values and ii) non-zero values assuming a gamma distribution of concentrations. In ecological statistics this type models is also known as *hurdle* models (Martin et al., 2005). The log-normal Tobit model has no probability mass at zero, whereas ZAGA model has a probability at zero. Generally, the difference between Tobit and two-part models are small (Min and Agresti, 2002) and the same holds for differences between the log-normal and Gamma distribution. Indeed, a Tobit-like model could be also fitted assuming a Gamma distribution (Sigrist and Stahel, 2010).

Grab sampling likely underestimates chemical concentrations (Xing et al., 2013) because of short term peak concentrations (Wittmer et al., 2010). Although this leads to an increased variation in chemical measurements, we still can learn from the process generating values above LOQ, even if the absolute value is subject to error. This is also highlighted by the results of chapter ??, with estimated coefficients for the absolute concentration showing much larger uncertainty than coefficients for the probability of exceeding LOQ (Figure ??). Currently, models explicitly taking the censored nature of chemical monitoring data are not well explored and seldom applied. Further research on those is needed and might provide useful information for analysing monitoring data, assessing the chemical status and trends thereof.

## LEVERAGING MONITORING DATA FOR ECOLOGICAL RISK ASSESSMENT

In chapter ?? we compiled and analysed

## CHALLENGES UTILISING 'BIG DATA' IN ECOLOGICAL RISK ASSESSMENT

Effect assessment and environmental monitoring produce huge amounts of data. However, the profoundness of ecological risk assessment often determined by the available data (Van den Brink et al., 2016). Useful data for ERA is currently spread over several largely unconnected databases. E.g. ecotoxicity data is spread over database maintained by the U.S. EPA (ECOTOX, U.S. EPA (2016)), the University of Hertfordshire (PPDB, Lewis et al. (2016)), the German Environment Agency (ETOX, Umweltbundesamt (2016)) and others. Chemical information is similarly spread over several databases, like PubChem (Kim et al., 2016) or Chemspider (Pence and Williams, 2010). Additional complications arise because these databases use different identifiers for chemical substances. The U.S. EPA (U.S. EPA, 2016) uses solely the CAS-Number for identification, whereas other databases uses SMILES (Weininger, 1990) or InChI and InChIKeys (Heller et al., 2015). Integrating these databases is currently a challenge in ERA, which is complicated by ambiguous identifiers, e.g. should different salts be considered separately for aquatic risk assessment? Projects like the NORMAN EMPODAT database (Brack et al., 2012) or the STOFF-IDENT (Huckele and Track 2013, <http://uni-ko-ld.de/fc>) are first attempts for such an integration. Integration monitoring data and risk assessment data is a mandatory requirement for landscape level ecotoxicology and risk assessment (Cairns and Niederlehner, 1996; Focks, 2014) and needed for model validation (Knäbel et al., 2012). Chapter ?? is an example for such an integration, but represents only a preliminary assessment and spatial-temporal risk dynamics should be further investigated.

The webchem package, presented in chapter ??, can foster such an integration, however data must also be accessible. Unfortunately, major parts of data produce for environmental risk assessment are not available (Schäfer et al., 2013). Recently, it has been demonstrated that data from the European Registration, Evaluation, Authorisation, and Restriction of Chemicals (REACH) database can be used to improve the characterisation of ecotoxicity in life cycle assessment

(LCA) (Müller et al., 2016). Although this database hosts data used risk assessment, it is not available in a convenient way. Indeed, a systematic data collection contravenes the legal usage of the REACH database (<http://uni-ko-ld.de/fd>). This may be also the reason, why the quality of chemical property data submitted this database is currently unknown (Müller et al., 2016; Stieger et al., 2014).

The software tools described in chapters ?? and ?? assist researchers handling and cleaning their data. Aggregating taxonomic data to a higher taxonomic level is a common task when analysing data from mesocosm experiments or from field sampling. Taxize facilitates the retrieval of taxonomic classification, which is the basis also for more sophisticated aggregation methods (Cuffney et al., 2007). Today, taxize has been used in more than thirty scientific publications. Recent applications of the webchem package, have been demonstrated by Münch and Galizia (2016) and Ranke (2016): Münch and Galizia (2016) compiled a database for odorant responses of *Drosophila melanogaster* and webchem "likely saved [him] hundreds of working hours". Ranke (2016) is using webchem to compile and store chemical information for further analyses. The analyses performed in chapter ?? needed to integrate monitoring, chemical and risk assessment data which would not have been difficult without the webchem package. These examples show that researchers have been missing such tools in the past. If they can reduce their time spend on data retrieval and handling, they could focus more on the quintessence of their research.

## CONCLUSIONS



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