

# Insights on impact estimation of unmeasured micropollutants from WWTP release

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## Insights on impact estimation of unmeasured micropollutants from WWTP release

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Abstract: During wastewater treatment, incomplete elimination of micropollutants occurs. These micropollutants represent a potential threat to our health and aquatic environment. The potential impacts of these substances are underestimated due to the lack of available concentrations below the limit of quantifications (LOQ) for an important set of micropollutants. Here, we studied the potential impacts on Human health and aquatic environment of the release of 261 organic micropollutants at the scale of France and we showed that, even with a concentration below the LOQ, some micropollutants could have a significant impact. For the unmeasured concentrations, a global concentration distribution build on several datasets with different LOQ is used.

The usual convention of LOQ/2 for the unmeasured concentrations is shown to be an overestimation. Ignoring the unmeasured micropollutants led to an underestimation of the impacts of more than 300% for both human health and aquatic environment. Some substances, such as hydrazine, endrin, or 2,3,7,8-TetraCDD, could lead to a very high impact even with an unmeasured concentration. This work can be adapted to any other compartment and any other geographical context. **Keywords:** characterization factors; limit of quantification; (eco)-toxicity; continental freshwater; human health; Life Cycle Assessment. **Highlights:** Impacts for micropollutants with concentrations under LOQ are derived. Some micropollutants have a significant impact with an unmeasured concentration. Ignoring these unmeasured substances lead to a strong underestimation of the impacts. This could help to a prioritization of unmeasured micropollutants. 

## **Graphical Abstract**:

### 1. Introduction

Thanks to the implementation of the Water Framework Directive (Directive 2008/105/CE, 2008), many European countries have launched monitoring campaigns for organic micropollutants either in the effluents of urban and industrial wastewater treatment plants (WWTP) or in the receiving rivers. Indeed, many papers refer to the quantification of thousands of compounds including pharmaceuticals, hormones, pesticides, hydrocarbons, and solvents in treated wastewaters, underlying incomplete removals of the conventional treatments (Addamo et al., 2005; Bisognin et al, 2021; Yao et al., 2021).

Moreover, with increasingly accurate and efficient analytical technologies, more and more compounds are detected at low concentrations (ng.L-1 to  $\mu$ g.L-1) in rivers,

groundwater, surface water, and drinking water (Bayer et al., 2014). The incomplete removal and release of micropollutants into the aquatic environment represent a potential danger to human health and the aquatic environment of continental freshwater. The identification and quantification of the potential impact of toxic substances induce the development of new sustainable process technologies for targeted substances and represent a great challenge to the safety of aquatic ecosystems (Rosenbaum et al., 2007; Zhu et al., 2015). Recently, Aemig et al. (2021) identified 261 organic micropollutants in the effluents of French WWTP. The toxicological and ecotoxicological impacts of 94 and 88 micropollutants, respectively, were quantified by multiplying the emitted mass of the compound in the total volume of effluents from French WWTPs by their characterization factors (CFs) (Lindim et al., 2019) as usually done in the Life Cycle Assessment (LCA) framework. Some compounds cannot be assessed because they have no known concentration or their CFs were not available. Then, Servien et al. (2022a) developed a modeling method based on machine learning approaches and 40 easy-to-obtain molecular descriptors to predict missing CFs. This model allowed the prediction of toxicological and ecotoxicological CFs in continental freshwater with an acceptable margin of error. This allowed a completed assessment of the overall potential impacts (on human health and aquatic environment) of 153 organic micropollutants in continental freshwater at the scale of France. Results showed that a molecule can be highly hazardous for the aquatic environment without necessarily being an issue for human health, and vice versa. Moreover, a high impact could be driven by a high CFs, a high mass, or a combination of both. As the machine learning models allow a CFs prediction for any new compound, the missing concentrations are the only burden that prevents a complete impact estimation. Nevertheless, building

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models based on easy-to-obtain inputs seem not easily attainable for missing concentrations and the lack of data for some micropollutants is nowadays a known problem (Leenhardt et al., 2022). Cantoni et al. (2020) tackle this problem by comparing several statistical methods to estimate concentrations below the LOQ. Their work is promising but is limited to a few number of micropollutants and cannot be applied to compounds with no value above the LOQ. Not considering the micropollutants with a high number of concentrations below the LOQ in the impact estimations is equivalent to estimating that they have null impacts, which obviously leads to a strong underestimation of the potential impacts. So, the objective of this paper was to study the potential impacts of the micropollutants for which we do not have measured concentrations or for which more than 90% of the measurements are below the limit of quantification (LOQ) threshold, including those with no measurement above the LOQ. For this, we develop an estimation procedure detailed in the following Section. This allowed a complete assessment of the individual potential impacts (on human health and aquatic environment) of each of the organic micropollutants of interest in continental freshwater at the scale of France.

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#### 2. Materials and Methods

### 2.1 Molecules

Aemig et al. (2021) identified 261 organic micropollutants representing a potential danger to human health and aquatic environment in continental freshwater. These molecules came from (i) the Water Framework Directive (WFD, Directive 2008/105/CE, 2008), (ii) the RSDE national action for the survey and reduction of hazardous substances in water (INERIS, 2016), and (iii) the AMPERES French project in which

micropollutants (registered in the WFD and pharmaceuticals) were analyzed in influents and effluents of 15 WWTP (Martin Ruel et al., 2016). Among these 261 organic micropollutants, the emitted mass of 153 was estimated with at least 10 sites of measures and 90% of the measured data above the limit of quantification. The impacts of these 153 compounds were assessed by Aemig et al. (2021) for those present in the USEtox® database (Rosenbaum et al., 2007) and by Servien et al. (2022a) for those who are not in this database.

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## 2.2 Characterization factors

Among the remaining 108 compounds, 58 have a CF<sub>ET</sub> and 43 a CF<sub>HT</sub> in the USEtox® database version 2.12 with a default landscape (to remain consistent with Servien et al. (2022a,b)). USEtox® is an international consensual for characterizing human and ecotoxicological impacts of chemicals (UNEP-SETAC, 2019). It was developed by the life cycle initiative under the United Nations Environmental Programme (UNEP) and the Society for Environmental Toxicology and Chemistry (SETAC) (Henderson et al., 2011) to produce a transparent and consensus characterization model. This model gathers in one single characterization factor the chemical fate, the exposure, and the effect for each of the several thousands of organic and inorganic compounds. If the structure of the USEtox® multimedia model is always the same, to determine the CF of a molecule, numerous physicochemical parameters (such as solubility, hydrophobicity, degradability) and detailed toxicological and ecotoxicological data must be provided. The remaining CFs were computed using the machine learning models developed by Servien et al. (2022a). These models predict ecotoxicological or toxicological characterization factors using 40 selected easy-to-obtain molecular descriptors (see

the list in Table S1). For more details on the choice and the calculation of these descriptors, the interested reader is referred to Servien et al. (2014). These models are based on a comparison between different linear or non-linear (machine learning) models and could be global or local using a cluster-then-predict approach. The modeled CFs were compared to USEtox® database, the differences between the modeled CFs and the USEtox® ones were small, and the models were therefore considered to be applicable. The best approaches were then selected for each cluster and each characterization factor. To apply the cluster-then-predict approaches, the cluster of each new compound needs to be estimated using a supervised clustering approach based on the clusters already obtained in Servien et al. (2022a) which was performed using the well-known k-nearest neighbor approach through the *knn* function of the R package class (Ripley, 1996). Then, the models selected in Servien et al. (2022a) were applied without any modification, assuming they had been previously tested on a large diversity of molecules covering the diversity of the compounds assessed in the present study. These previous tests have shown the reliability of the models to predict CFs that are missing in USEtox®. Note that the tributylstannanylium (CAS 36643-28-4) is made of tin and, as consequence, its molecular descriptors cannot be computed by the above-mentioned methods and, thus, its CFs. Also, chloroalkanes C10-13 (CAS 85535-84-8) is a mix of different molecular structures that can lead to different molecular descriptors and CFs and was also excluded. The 106 corresponding characterization factors are gathered in Table S2. The scheme

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of the whole study is summarized in the following Figure 1.

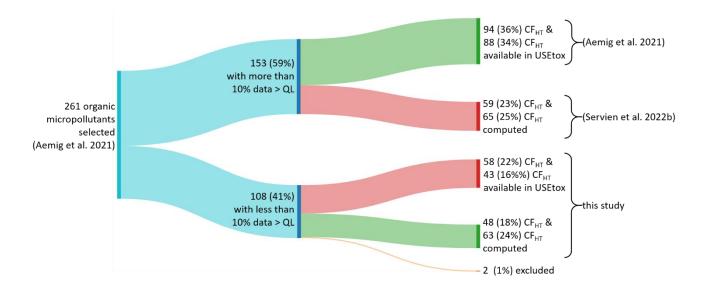


Figure 1. Position of this study in the general context.

## 2.3 Quantification of significative potential impacts

The annual volume of water effluent released into the environment by WWTP was estimated to be 5.10<sup>9</sup> m<sup>3</sup>, in France (Aemig et al., 2021). This volume was determined by multiplying the daily flow arriving at the WWTP for 365 days, assuming that it was equal to the effluent. The emitted mass of micropollutants was estimated by multiplying each concentration of the micropollutants, when available, by the volume of the effluent. The concentrations and masses are those of the Supplemental Material of Aemig et al. (2021) and are gathered in Table S3. Total impacts on human health and aquatic environment in continental freshwater were quantified by summing the impacts of all the compounds, as is usually done in LCA (Heijungs and Suh, 2002).

As explained above, the remaining 106 micropollutants are those for which not enough data were available to correctly estimate the global concentration. So, for these compounds, we chose to calculate the minimal concentration to reach 1% of the total sum of impacts calculated previously for the 153 micropollutants presented in Servien et al. (2022b). When the LOQ of a micropollutant was available, this minimal

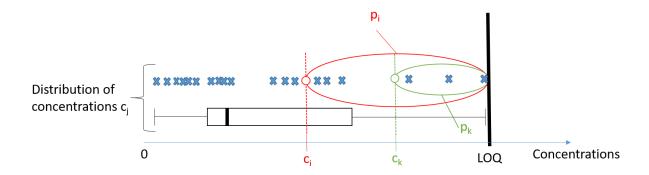
concentration was also expressed in percentage of the LOQ to be comparable among different micropollutants. For 42 compounds, no data (measurements and LOQ) were available so this minimal concentration was only expressed in  $\mu$ g.L<sup>-1</sup>. Obviously, as the potential impacts are linearly related to the estimated concentration, if a measured concentration is the minimal concentration multiplied by a constant k, the corresponding impacts will be k% of the total sum of impacts calculated previously.

### 2.4 Estimation of values under the LOQ

For these 106 micropollutants, according to Table S3, a very high number (more than 90%) of the values are below the LOQ of the corresponding study. Data inferior to LOQ are usually estimated at half of the quantification limit as usually applied in environmental studies (INERIS, 2016). This assumption is a very strong one that could have a major impact on the estimated concentration (and thus the potential impact) when there are a lot of concentrations below LOQ. Testing this strong assumption could be of great help.

In the RSDE datasets, the compounds are measured using routine methods so most of the LOQ are pretty high. This is not always the case for the other sources of quantification data. So, we have some quantifications that came from other sources and that are below the LOQ of the main dataset of the study, the RSDE one. These values are then scaled relative to the RSDE LOQ and are gathered for all compounds to have enough data. This gives a dataset of 141 samples with concentrations denoted  $(c_j)_{1 \le j \le 141}$  coming from 25 different compounds. The distribution of these samples is then used to compute the probability that the minimal concentration  $c_i$  of compound i has been reached in case of quantification below the LOQ. This probability

 $p_i$  is computed as the number of measurements above the corresponding minimal concentration  $c_i$  divided by the total number of measurements (i.e. 141):  $p_i = \frac{\sum_{j=1}^{141} \mathbb{I}_{\left(c_i < c_j\right)}}{141}$  and is detailed in the following Figure 2. In other terms, when the minimal concentration  $c_i$  is near the LOQ, the probability to get a value above  $c_i$  is low.



**Figure 2**: The distribution of concentrations  $(c_j)_{1 \le j \le 141}$  is represented here by some blue crosses and its corresponding boxplot. A new minimal concentration  $c_i$  is estimated for a compound i, this results in the corresponding  $p_i$ , calculated by counting all the concentrations  $c_j > c_i$ . Another minimal concentration  $c_k$  is estimated for another compound k with  $c_k > c_i$ . Obviously, the probability  $p_k$  that the minimal concentration for the compound k will be attained is below  $p_i$ .

Note that if a value is below the two LOQ (the RSDE one and the one of the other dataset) it won't be taken into account, so the final distribution will probably be a little overestimated. Indeed, taking into account these unknown measurements relying on different relative LOQ values would bring a lot of uncertainties in the below LOQ distribution calculations.

## 2.5 Calculation of the total impacts

As usually done in LCA, the total impacts on human health and aquatic environment in continental freshwater were quantified by summing the impacts of all the compounds. Human impact is expressed in DALY representing the number of negatively impacted human years, and the ecotoxicological impact is expressed in PDF representing the potential fraction of disappeared species. To compute these impacts, we impute values under the LOQ by the median values under the LOQ for all the compounds (see Figure 3), namely LOQ/3. For the compounds without a LOQ, the median value of the LOQ, namely 0.05 µg.L-1, was used.

### 3. Results and discussion

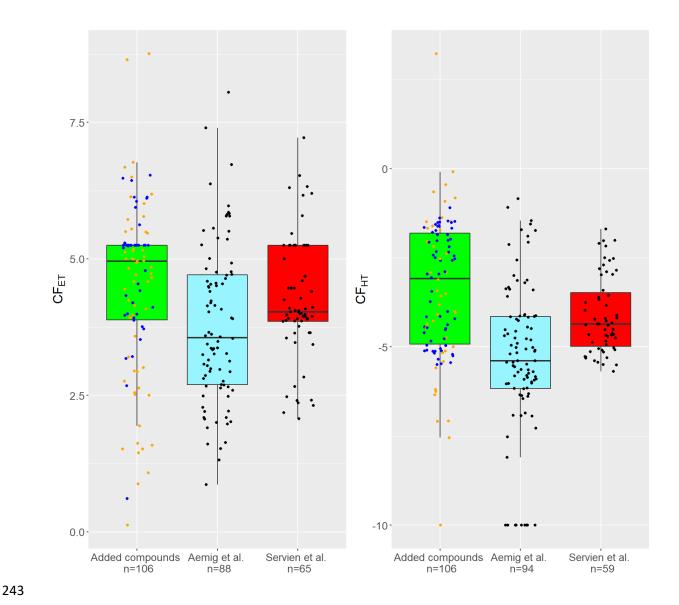
## 3.1 Global analysis of characterization factors

Regarding the CFs, the different values are reported in the boxplots of Figure 3 below.

As usual, the CFs are log-transformed with the addition of 1e-10 for the CF<sub>HT</sub> to avoid

a computational problem with log<sub>10</sub>(0). 1e-10 has been chosen to be below the

minimum of the USEtox® database (5e-9).



**Figure 3**. Comparison of the CF<sub>HT</sub> and the CF<sub>ET</sub> (both in log10 scale) for the compounds taken into account in previous studies and the added compounds. For the added compounds, the CF in orange came from USEtox® whereas those in blue are computed using Servien et al. (2022a).

Figure 3 suggests that the added compounds have globally higher CF than already included in the previous studies. In more detail, for the  $CF_{HT}$  and the  $CF_{ET}$ , the

maximum, the 1st and 3rd quartile, the median and the mean of the added compounds are above those of the previous studies. But there are also very low CF values largely below those of the previous study. All the descriptive values are given in Table S4. We could also see that we have extremely high values for some compounds already included in USEtox®. For the CFET, these two extreme values are for the 1,2,3,7,8-PentaCDD (CAS 40321-76-4) and the 2.3,7,8-TetraCDF (CAS 51207-31-9) from the polychlorodibenzodioxin/furane family. Their very extreme values would lead to significant impacts if their concentrations are not too low. For the CF<sub>HT</sub>, the very extreme value is for the 2,3,7,8-TetraCDD (CAS 1746-01-6). This is the unique value of the whole USEtox® database that is log-positive for this variable. Nevertheless, the estimation of the CF<sub>HT</sub> available in the environmental footprint database (Fazio et al., 2018) seems to corroborate this very high value. This dioxin is a well-studied and recognized endocrine disrupter presenting the highest affinity to the aryl hydrocarbon receptor (AhR), a cytosolic protein present in most vertebrate tissues, that mediated the toxic and biological effects of such compounds (Van den Berg et al., 2006). Its very high CF<sub>ET</sub> and CF<sub>HT</sub> are linked to its high endocrine-disrupting potency. If we don't consider this compound, five others have CF<sub>HT</sub> above the maximal value already included in previous studies. These compounds are all pesticides: chlordane (CAS 57-74-9), 1,1-bis(4-Chlorophenyl)-2,2-Dichloroethane (CAS 72-54-8), dieldrin (CAS 60-57-1), heptachlor epoxide (CAS 1024-57-3) and toxaphene (CAS 8001-35-2).

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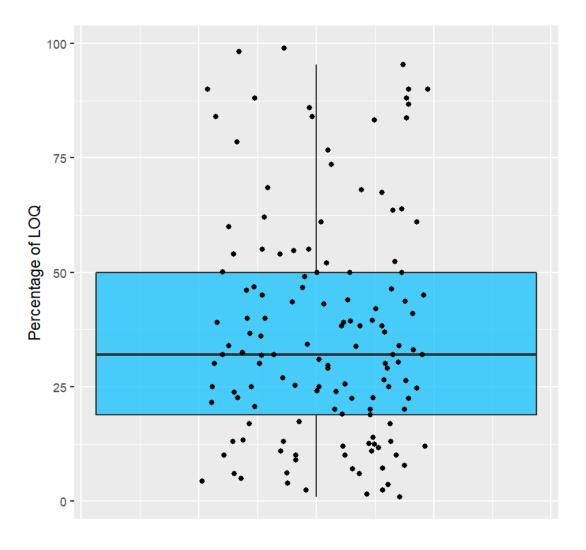
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### 3.2 Estimation of concentrations under the LOQ

In the following Figure 4, the distribution of the concentrations measured below the RSDE LOQ is represented.



**Figure 4:** Boxplot of the 141 measured values below the RSDE LOQ on another dataset with lower LOQ (expressed in % of the RSDE LOQ).

We could see that the median (and so is the mean) value of this distribution seems to be close to LOQ/3, and not LOQ/2 as usually used for calculation when data are below the LOQ. Furthermore, as the concentration values under the smaller LOQ are not taken into account, even LOQ/3 is probably an overestimation, when faced with an under LOQ value in the RSDE dataset. This is also confirmed by the conclusions of Cantoni et al. (2020) where the LOQ and LOQ/2 are shown to be overestimation.

## 3.3 Impact on the aquatic environment

For the aquatic environment, all the 64 minimal concentrations that could be expressed in percentage of LOQ are gathered in Table S6 and the smallest (i.e. that are below the LOQ) are reported in Table 1.

**Table 1** – Estimation of the concentrations needed to reach 1% of the global impacts on ecotoxicity, when the minimal concentration is below the RSDE LOQ. The probability is calculated when the percentage of LOQ is under 100%. It is computed using the distribution of subsection 3.2 and is the estimation that a measurement under the LOQ is above the minimal concentration. The results for all compounds are gathered in Table S6.

	Minimal	Minimal	
	concentration	concentration (% of	
Name	(μg.L <sup>-1</sup> )	RSDE LOQ)	Probability
Hydrazine	4,0E+00	4%	0.96
Endrin	1,0E-02	19%	0.75
Chlordecone	9,0E-02	46%	0.29
Chlorpyrifos	2,0E-02	88%	0.04
	Hydrazine Endrin Chlordecone	$\begin{array}{c} \text{concentration} \\ \text{Name} & (\mu g. L^{-1}) \\ \\ \text{Hydrazine} & 4,0E+00 \\ \\ \text{Endrin} & 1,0E-02 \\ \\ \text{Chlordecone} & 9,0E-02 \\ \end{array}$	$\begin{array}{c} \text{Concentration} & \text{concentration (\% of} \\ \text{Name} & (\mu g.L^{-1}) & \text{RSDE LOQ)} \\ \\ \text{Hydrazine} & 4,0E+00 & 4\% \\ \\ \text{Endrin} & 1,0E-02 & 19\% \\ \\ \text{Chlordecone} & 9,0E-02 & 46\% \\ \end{array}$

There are only 4 compounds for which a concentration under the LOQ could still cause a significant impact on ecotoxicity and for 2 of them the minimal concentration is also below LOQ/3. For the 60 other compounds, the minimal concentration being above the LOQ and the measured one frequently below, their contribution to the environmental impact could be considered as very low. For the hydrazine, a very small concentration of only 4% of the RSDE LOQ, that have a high probability of occurrence, will lead to a significant impact. Hydrazine and its derivatives are widely used compounds in chemical industries and pharmaceuticals (propellant, blowing agent, pesticide, antituberculosis...). It is a very small molecule (32 g/mol), very soluble and volatile, no BCF is found in literature, nor DT50. It was classified by US EPA as a human carcinogen (group B2): an intermediate inhalation minimal risk level of 4 µg/L was calculated and a minimum limit for cancer was fixed at 10 µg/L (Qi et al., 2020). The

minimal concentration was found to be 4  $\mu$ g/L and is largely below the RSDE LOQ, which is very high (100  $\mu$ g/L) by comparison to the one of endrin (0,052  $\mu$ g/L) and also others compounds. However, very recent papers refer to the development of spectroscopic methods and sensors to analyze hydrazine in environmental samples including water up to pico-molar level (0.03 to 3.2  $\mu$ g/L, Arulraj et al., 2015; Qi et al., 2020). However, no data are provided in real water samples.

For the three pesticides, due to their hydrophobicity, they are usually quantified in solid matrices like soil, sediment, and biota. However, concentrations of endrin were reported in river water between 0.17 and 484 ng/L (Matsumo et al., 2009) and concentrations of chlorpyrifos between 1 and 32 ng/L (Wan et al., 2021). Their presence in such environment is mainly due to their use in agriculture. They are also targeted by the WFD. The Environmental Quality Standard (yearly average value) for the sum of the 4 cyclodiene pesticides in the WFD (aldrin, isodrin, endrin, dieldrin) is 10 ng/L (inner water surface) and 5 ng/L (other water). For chlorpyrifos, the EQS is 30 ng/L. The three calculated minimal concentrations to reach 1% of the global impact are in the same range of these environmental and regulatory values.

The compounds without any information on LOQ are gathered in the following Table 3.

**Table 3**. Minimal concentrations for each compound to reach 1% of the global impacts on ecotoxicity. The compounds in this table are those without any measure or LOQ.

CAS number	Name	Minimal concentration (μg.L <sup>-1</sup> )
40321-76-4	1,2,3,7,8-P5CDD	9,62E-05
51207-31-9	2,3,7,8-T4CDF	1,23E-04
1746-01-6	2,3,7,8-T4CDD	1,16E-02
57653-85-7	1,2,3,6,7,8-H6CDD	3,61E-02
72-54-8	DDD 44'	4,04E-02
28159-98-0	Cybutryne	5,31E-02
72-55-9	DDE 44'	1,56E-01

19408-74-3	1,2,3,7,8,9-H6CDD	3,08E-01
39227-28-6	1,2,3,4,7,8-H6CDD	3,08E-01
57117-44-9	1,2,3,6,7,8-H6CDF	3,08E-01
70648-26-9	1,2,3,4,7,8-H6CDF	3,08E-01
72918-21-9	1,2,3,7,8,9-H6CDF	3,08E-01
57117-41-6	1,2,3,7,8-P5CDF	3,09E-01
3268-87-9	1,2,3,4,6,7,8,9-O8CDD	3,10E-01
35822-46-9	1,2,3,4,6,7,8-H7CDD	3,10E-01
39001-02-0	1,2,3,4,6,7,8,9-O8CDF	3,10E-01
446255-20-5	Heptabromodiphenylethers	3,10E-01
55673-89-7	1,2,3,4,7,8,9-H7CDF	3,10E-01
57117-31-4	2,3,4,7,8-P5CDF	3,10E-01
60851-34-5	2,3,4,6,7,8-H6CDF	3,10E-01
67562-39-4	1,2,3,4,6,7,8-H7CDF	3,10E-01
789-02-6	DDT 24'	5,58E-01
2032-65-7	Methiocarb	6,08E-01
25167-83-3	Tetrachlorophenols	6,50E-01
131860-33-8	Azoxystrobin	1,43E+00
139968-49-3	Metaflumizone	1,59E+00
118-79-6	2,4,6-tribromophenol	1,98E+00
564-25-0	Doxycycline	2,11E+00
153719-23-4	Thiamethoxam	2,61E+00
111988-49-9	Thiacloprid	4,24E+00
2303-17-5	Triallate	5,88E+00
21312-10-7	Acetyl sulfamethoxazole	6,67E+00
5466-77-3	Octinoxate	7,36E+00
615-58-7	2,4-dibromophenol	9,62E+00
135410-20-7	Acetamiprid	1,05E+01
210880-92-5	Clothianidin	1,65E+01
14297-93-9	4-epi-chlortetracycline	3,37E+01
25167-80-0	Monochlorophenol	3,67E+01
25167-82-2	Trichlorophenols	5,41E+01
128-37-0	Butylhydroxytoluene	6,19E+01
57-62-5	Chlortetracycline	1,27E+02
95-56-7	2-bromophenol	1,55E+02

For these compounds, only the minimal concentration in the usual unity (µg.L-¹) is available. It leads to results that are more difficult to interpret but we could see that there are 2 compounds from the dioxin/furane family (1,2,3,7,8-PentaCDD and 2,3,7,8-TetraCDF) for which a concentration above 0,01 and 0,1 ng.L-¹ will be enough to lead to a significant impact. The 15 other congeners of this family present minimal

concentration between 11 and 310 ng/L. These compounds, produced during waste incineration, and chemical manufacturing, are classified as very hydrophobic and persistent compounds, and accumulate in environmental compartments such as soil, and sediment as well as in biota (Srogi, 2008). As an example, Villanneau et al (2011) found soil concentrations of 23,8 ng I-TEQ/kgDM for the ∑17 PCDD/F. The regulatory limit in meats for the sum of dioxin and furan is 1 to 5 ng I-TEQ / kg of fat. In such organic and fat matrices, low concentrations can be found, but the probability to find such low concentrations in water matrices may be very low. Dioxins and furans are also on the list of priority substances of the WFD but no EQS for water is provided only EQS for the biota is fixed at 6,5 ng/kg DW (I-TEQ for 7 PCDD+10 PCDF+12 PCB). Such low concentrations could probably be reached locally (see Lu et al., 2016, for PCDD/F concentrations in raw and treated drinking water and Ellis et al., 2008, for data on lake Chelan). DDD 44' and DDE 44' are transformation products of DDT that is no more used in western countries since the 90th. All congeners are hydrophobic and persistent compounds with accumulation in biota (Garrison et al., 2014). Total DDT, including several DDT isomers, DDE, and DDD, is listed on the WFD with an EQS of 25 ng/L implying its monitoring in the French RSDE program. Cybutryne is a biocide from the triazine family and is used mainly in anti-fouling paints for the protection of boat hulls. It was included on the WFD priority list for monitoring and control in 2012 with an EQS of 2.5 ng/L (yearly average) and a maximal admissible concentration of 16 ng/L. In Switzerland, Ollers et al. (2001) have quantified cybutryne in effluent at very low frequencies around 15 ng/l, which is below the calculated minimal concentration (53 ng/L).

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At the bottom of Table 3, the listed compounds present high minimal concentrations ranging from 10 to 155 µg/L. As such, chlorophenols, present on EU and US EPA priority lists, are widely used for dyestuff, pharmaceutical and pesticide synthesis and are also by-products of their degradation in the environment (Ganigar et al., 2010). These compounds were mainly monitored in industrial wastewaters where they are treated by biological and physicochemical processes before discharge in environment or domestic sewage; this could explain why no concentration of such compound was found in the domestic effluent of the French studied WWTP. Gao et al. (2008) have quantified 2,4,6-trichlorophenol in 54.4 % of their river water samples (600 sites) at a median value of 2 ng/L. These low measured values suggest also the existence of low limit of quantifications as the one given by Suliman et al. (2006) around µg/L for TCP. These data suggest that it may not be realistic to get such minimal concentration in effluent. To reinforce this idea, we can have a look on compounds like chlortetracycline and its metabolite, 4-epi-chlortetracycline, whose minimal concentrations to reach 1% of the total ecotoxicological impact are respectively 127 and 33 µg/L: these compounds were found in the effluent of many WWTP but at very lower concentrations than the minimal one. For example, Li and Zhang (2011) measured chlortetracycline levels around 24 and 33 ng/L in secondary effluents of 2 WWTP in Hong Kong. Chlortetracycline concentrations in effluents of American WWTPs (<8.1–60 ng L<sup>-1</sup>) (Yang et al., 2005; Spongberg and Witter, 2008) and Australian WWTPs (5-250 ng L-1) (Watkinson et al., 2009) were also found.

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## 3.4 Impact on Human health

For human toxicity, all the 64 minimal concentrations that could be expressed in percentage of LOQ are gathered in Table S5 and the smallest (i.e. that are below their LOQ) are reported in the following Table 2.

Table 2 – Estimation of the concentrations needed to reach 1% of the global impacts on human toxicity, when the minimal concentration is below the RSDE LOQ. The probability is calculated when the percentage of LOQ is under 100%. It is computed using the distribution of subsection 3.2 and is the estimation that a measurement under the LOQ is above the minimal concentration. When the probability is equal to one, it means that no measured value is below this percentage of LOQ. The results for all compounds are gathered in Table S5.

		Minimal	Minimal concentration	
CAS number	Name	concentration (µg.L <sup>-1</sup> )	(% of RSDE LOQ)	Probability
302-01-2	Hydrazine	2,1E-01	0.2%	1
1024-57-3	Heptachlor epoxide	8,2E-05	0.4%	1
8001-35-2	Toxaphene	2,9E-04	0.6%	1
143-50-0	Chlordecone	1,6E-03	0.8%	1
60-57-1	Dieldrin	4,4E-04	0.9%	1
57-74-9	Chlordane	5,5E-04	2.8%	0.97
309-00-2	Aldrin	2,0E-03	4.0%	0.96
2385-85-5	Mirex	2,7E-03	5.4%	0.94
72-20-8	Endrin	3,5E-03	7.0%	0.91
62-53-3	Aniline	6,6E+00	13%	0.79
76-44-8	Heptachlor	3,1E-03	15%	0.77
50-29-3	DDT 44'	1,0E-02	20%	0.71
50-32-8	Benzo(a)pyrene	2,1E-03	21%	0.7
470-90-6	Chlorfenvinphos	2,3E-02	23%	0.67
67-56-1	Methanol	2,4E+03	24%	0.67
36483-60-0	Hexabromodiphenylethers	6,2E-03	31%	0.52
36355-01-8	Hexabromobiphenyl	6,9E-03	35%	0.44
75-01-4	Chloroethene	1,9E+00	38%	0.42
32536-52-0	Octabromodiphenylethers	2,5E-02	50%	0.26
120-12-7	Anthracene	2,3E-02	56%	0.18
39635-31-9	PCB 189	2,9E-03	59%	0.18
124495-18-7	Quinoxyfen	6,0E-02	60%	0.18
118-74-1	Hexachlorobenzene	7,2E-03	72%	0.12
3424-82-6	DDE 24'	7,6E-02	76%	0.11
87-68-3	Hexachlorobutadiene	4,3E-01	86%	0.06

More compounds (26) could have a significant impact here if we compare with the previous subsection on environmental impact and 16 have a minimal concentration below LOQ/3 and will let to a significant increase in the total sum of impacts if they were taken into account. Among the four previous compounds (hydrazine, endrin, chlordecone, chlorpyrifos) with a significant ET impact, only chlorpyrifos is not present here. Many of these compounds are pesticides, hydrocarbons, and fire retardants that are persistent and hydrophobic compounds usually quantified in soil, sediment, and biota for which the minimal concentration is between 10 pg.L<sup>-1</sup> (heptachlor epoxide) and 100 ng.L<sup>-1</sup> (hexachlorobutadiene). There are also smaller molecules like aniline, chloroethene, and methanol, used as solvents and intermediate chemicals for dye and pharmaceutical industries whose minimal concentrations are higher between μg.L<sup>-1</sup> and mg.L<sup>-1</sup> (methanol). These compounds are present in industrial wastewater at very high concentrations and are quite well degraded in WWTP (Chaturvedi and Katoch, 2020), suggesting low residual concentrations in urban effluents. The compounds without any information on LOQ are gathered in Table 4.

**Table 4**. Minimal concentrations for each compound to reach 1% of the global HT impacts. The compounds in this table are those without any measure or LOQ.

CAS number	Name	Minimal concentration (μg.L <sup>-1</sup> )
1746-01-6	2,3,7,8-T4CDD	3,97E-08
72-54-8	DDD 44'	1,87E-04
39001-02-0	1,2,3,4,6,7,8,9-O8CDF	8,37E-04
3268-87-9	1,2,3,4,6,7,8,9-O8CDD	1,60E-03
55673-89-7	1,2,3,4,7,8,9-H7CDF	1,93E-03
35822-46-9	1,2,3,4,6,7,8-H7CDD	2,02E-03
67562-39-4	1,2,3,4,6,7,8-H7CDF	2,10E-03
57653-85-7	1,2,3,6,7,8-H6CDD	2,14E-03
39227-28-6	1,2,3,4,7,8-H6CDD	2,18E-03

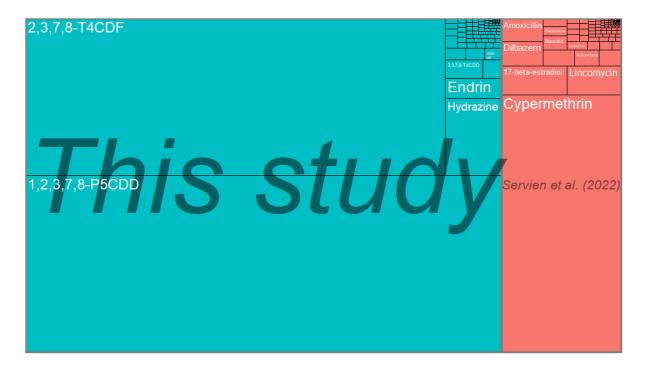
72918-21-9	1,2,3,7,8,9-H6CDF	2,41E-03
19408-74-3	1,2,3,7,8,9-H6CDD	2,65E-03
70648-26-9	1,2,3,4,7,8-H6CDF	2,73E-03
57117-44-9	1,2,3,6,7,8-H6CDF	2,78E-03
57117-41-6	1,2,3,7,8-P5CDF	3,11E-03
60851-34-5	2,3,4,6,7,8-H6CDF	3,56E-03
57117-31-4	2,3,4,7,8-P5CDF	4,12E-03
40321-76-4	1,2,3,7,8-P5CDD	4,72E-03
51207-31-9	2,3,7,8-T4CDF	1,03E-02
446255-20-5	Heptabromodiphenylethers	1,45E-02
72-55-9	DDE 44'	1,59E-02
789-02-6	DDT 24'	2,09E-02
118-79-6	2,4,6-tribromophenol	3,53E-01
25167-83-3	Tetrachlorophenols	9,29E-01
131860-33-8	Azoxystrobin	9,47E-01
139968-49-3	Metaflumizone	1,08E+00
2303-17-5	Triallate	1,56E+00
564-25-0	Doxycycline	1,59E+00
615-58-7	2,4-dibromophenol	1,82E+00
14297-93-9	4-epi-chlortetracycline	1,92E+00
57-62-5	Chlortetracycline	2,05E+00
25167-82-2	Trichlorophenols	2,66E+00
28159-98-0	Cybutryne	8,56E+00
2032-65-7	Methiocarb	8,65E+00
95-56-7	2-bromophenol	8,79E+00
111988-49-9	Thiacloprid	9,89E+00
153719-23-4	Thiamethoxam	1,07E+01
210880-92-5	Clothianidin	1,17E+01
25167-80-0	Monochlorophenol	1,46E+01
128-37-0	Butylhydroxytoluene	1,70E+01
5466-77-3	Octinoxate	1,81E+01
135410-20-7	Acetamiprid	1,95E+01
21312-10-7	Acetyl sulfamethoxazole	2,05E+01

As seen previously, the very high value for the CF<sub>HT</sub> for the 2,3,7,8-TetraCDD lead to a very low value for its minimal concentration. Many other compounds of the same family only need around 1 ng.L<sup>-1</sup> to have significant impacts. However, as already discussed for environmental impact, such low concentration will be measured with low probability in aqueous media. Other compounds as doxycycline, chlortetracycline, and acetyl sulfamethoxazole were already quantified in effluents of WWTP but at

concentrations largely below the calculated minimal concentrations (Verlicchi et al., 2012).

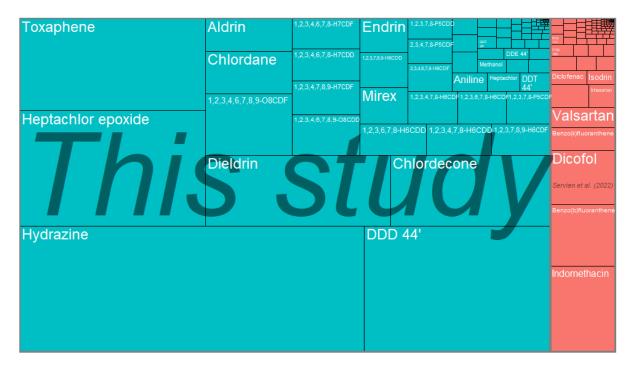
## 3.5 Influence on the sum of the impacts

According to all the assumptions made previously we were able to compare the sum of the impacts of the 106 micropollutants of this study to the 153 of Servien et al. (2022b). For the environmental impacts, the sum of the impacts of this study was nearly 4 times higher than the previous one. This could be seen in the following Figure 5 where the environmental impacts of the 259 compounds are represented.



**Figure 5**. Graphical representation of the environmental impacts of the 259 compounds studied in Servien et al. (2022b) (in red) and this study (in blue). Each compound has an estimated environmental impact proportional to its rectangular area.

For the impacts on human health it was more than 8 times higher, even without taking into account the molecule 2,3,7,8-TetraCDD. This unique molecule has an impact of more than 500 000 higher than all the other ones, if we consider it, highlighting its very high potential impact, even at a very small concentration. The other impacts are represented in the following Figure 6.



**Figure 6**. Graphical representation of the human impacts of the 258 compounds studied in Servien et al. (2022b) (in red) and this study (in blue). Note that the 2,3,7,8-T4CDD has been removed. Each compound has an estimated environmental impact proportional to its rectangular area.

The choice of the imputation value for the measured concentrations below the LOQ is also crucial. Indeed, a choice of LOQ/2 instead of LOQ/3 would have led to an overestimation of 50.0% of the overall impacts on human health and 49.7% on environmental impacts.

## 4 Conclusion

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In a previous study, 261 organic micropollutants were selected to study their potential impacts on human health and aquatic environment in continental freshwater at the scale of France. However, the lack of data allows quantifying the impacts for only 1/3 of these substances (88 for aquatic environment and 94 for human health). Using a new modeling approach, we were then able to estimate the impact of 153 organic micropollutants, i.e. for all with an estimated mass released in the environment (Servien et al., 2022). The hundred remaining compounds without an accurate concentration measure were not taken into account. It means that their impacts were considered null, even if their characterization factors were high and could be worrying. In this work, the potential impacts of all these compounds are considered. We derive individual compound results, that could help to select substances in which a special effort should be made on the tertiary treatments to implement in WWTP. It was particularly possible to obtain a rough estimation of the probability that a compound produces a non-negligible impact, even if a measured concentration is below a LOQ. Obviously, due to the lack of concentration data, no precise results on the total sum of the impacts could be obtained. Nevertheless, with some strong assumptions, it was possible to conduct such a study and to show that ignoring compounds with unmeasured concentrations could lead to strong underestimations of total impacts. A future work could be to balance the individual impacts of these compounds with a probability of occurrence. Also, it has to be underlined that this whole methodology (from the computation of missing CFs using machine learning models to the individual compound impact estimations with missing concentrations) can be adapted to any other compartment and any other geographical context, with predictive models still to develop.

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