Data Set Citation

When using this data, please cite the data package

Epinoux C , Barbarin M , Castrec J , Churlaud C , Dabrowski M , DAVID R , Fontaine Q , Fontanaud A , Fullgrabe L , Gillet C , Gobert S , Huet V , Lejeune P , Le Vern L , Madon B , Marengo M , Pignon-Mussaud C , Pillet M , Receveur J , and Rideau P. 2022.

Water Interdisciplinary Biology and Ecology database "WIBE": Towards FAIR, open and interdisciplinary data on biomarkers to monitor the ecological status of coastal waters.

Water Interdisciplinary Biology and Ecology database "WIBE" (https://pndb.fr/metacat/metacat/Water Interdisciplinary Biology and Ecology database "WIBE"/default)

General Information

Title:	**Water Interdisciplinary Biology and Ecology database "WIBE": Towards FAIR, open and interdisciplinary data on bior waters.**
Identifier:	autogen.2022081304461235329.1
Abstract:	The Water Interdisciplinary Biology and Ecology database "WIBE" database presents data from different scientific projects prop multi-biomarker study. This database gathers biological data of selected bioindicator species and environmental contextual dat concentrations of trace elements and organic pollutants were collected in the waters as well as biomarkers of effect and exposu after various analyses, were cleaned and reworked to meet FAIR and open data principles. We focused on developing a data c standards in order to make it as reusable as possible for the ecotoxicology research community. All datasets are available on tl The current dataset can be used by port and coastal water managers but also by marine ecotoxicology researchers who will be database on marine biomarkers allowing the monitoring of coastal water contamination and thus the proposal of remediation m
Keywords:	○ biomarker ○ marine ○ aquatic ○ water ○ port ○ pollution ○ Corsica ○ La Rochelle ○ ecotoxicology ○ interdisciplinary ○ WIBE ○ ecology ○ biology ○ FAIR principles ○ open science ○ database
Publication Date:	2022-06-15

Data Table, Image, and Other Data Details:

Metadata download	Ecological Metadata Language (EM

Data Table:

Name:	**Field_organisms_samples.csv**
Description:	Content of Field_organisms_samples.csv

Physical Structure Description:

Object Name:	Field_organisms_samples.csv
Size:	95882 bytes
Authentication:	4ddea7e077afd7088ff4ffd65699ad40 Caculated By MD5
Text Format:	Number of Header Lines: 1 Record Delimiter: \r\n Attribute Orientation: column Simple Delimited: Field Delim

Number Of Records: 927

Attribute(s) Info:

Name	Column Label	Definition	Type of Value	Measurement Type	Measurement Domain
Field_organism_sample_ID		Unique identifier of the collected organism or pool of organisms.	string	nominal	Def Unique identifier of the c organism or pool of orga
Fieldwork_ID		Unique identifier of the fieldwork carried out on a specific date and located at a specific station.	string	nominal	Def Unique identifier of the f carried out on a specific and located at a specific
Field_organism_sample_species_name		Species name of the sampled organism if known. If the species is not known, the genus name followed by sp. is used.	string	nominal	Domain Info
TAXREF_ID		Identifier of the sampled organism species if known according to the taxonomic repository TAXREF. If the species is not known, the genus given by the TAXREF is used, followed by sp..	string	nominal	Domain Info
Lab_pool_ID		Unique identifier of the pool of organism(s).	string	nominal	Def Unique identifier of the p organism(s).

Data Table:

Name:	**Fieldwork.csv**
Description:	Content of Fieldwork.csv

Physical Structure Description:

Object Name:	Fieldwork.csv
Size:	13728 bytes
Authentication:	65b2516d631e355053d250b3bd049232 Caculated By MD5
Text Format:	Number of Header Lines: 1 Record Delimiter: \r\n Attribute Orientation: column Simple Delimited: Field Delime

Number Of Records: 151

Attribute(s) Info:

Name	Column Label	Definition	Type of Value	Measurement Type	Measurement Domain

Fieldwork_ID		Unique identifier of the fieldwork carried out on a specific date and located at a specific station.	string	nominal	**Def** Unique identifier of the fieldwork carried out on a specific date and located at a specific station.
Sampling_station_ID		Accurate sampling location in a port.	string	nominal	**Def** Accurate sampling location in a port.
Sampling_date		Date of the sampling.	date	dateTime	
Project_ID		Acronym of the monetary fund used by the project linked to the identifying number of the project given by the monetary fund.	string	nominal	**Def** Acronym of the monetary fund used by the projec linked to the identifying number of the project give the monetary fund.
Sampling_month_value		Month of the year in which the sampling took place.	string	nominal	**Def** Month of the year in which the sampling took pla
Sampling_year_value		Year in which the sampling took place.	date	dateTime	

Data Table:

Name:	**Project.csv**
Description:	Content of Project.csv

Physical Structure Description:

Object Name:	Project.csv
Size:	785 bytes
Authentication:	029068073a20f3e5544e86d381fdd5cb Caculated By MD5
Text Format:	Number of Header Lines: 1 Record Delimiter: \r\n Attribute Orientation: column Simple Delimited: Field Delim
Number Of Records:	2

Attribute(s) Info:

Name	Column Label	Definition	Type of Value	Measurement Type	Measurement Domain
Project_ID		Acronym of the monetary fund used by the project linked to the identifying number of the project given by the monetary fund.	string	nominal	**Def** Acronym of the monetary fund used by t project linked to the identifying number project given by the monetary fund.
Project_acronym		Acronym of the scientific project from which the data are drawn.	string	nominal	**Def** Acronym of the scientific project from wh data are drawn.
Project_name		Full name of the scientific project from which the data are drawn.	string	nominal	**Def** Full name of the scientific project from w data are drawn.
Project_number_value		Identifying number of the scientific project from which the data are drawn given by the monetary fund.	string	nominal	**Def** Identifying number of the scientific proje which the data are drawn given by the monetary fund.
Project_funding_name		Origin of the monetary fund used by the project.	string	nominal	**Def** Origin of the monetary fund used by the
Project_coordinator_name		Name of the project coordinator.	string	nominal	**Domain Info**
Project_coordinator_activity_keywords		Field of work of the project coordinator.	string	nominal	**Def** Field of work of the project coordinator.
Project_beginning_date		Start date of the project.	date	dateTime	
Project_end_date		Date of the project completion.	date	dateTime	

Data Table:

Name:	**Sampling_stations.csv**
Description:	Sampling_stations.csv

Physical Structure Description:

Object Name:	Sampling_stations.csv
Size:	1962 bytes
Authentication:	fced076171efb098b02f0e3227433891 Caculated By MD5
Text Format:	Number of Header Lines: 1 Record Delimiter: \r\n Attribute Orientation: column Simple Delimited: Field Delim

Number Of Records:	15

Attribute(s) Info:

Name	Column Label	Definition	Type of Value	Measurement Type	Measurement Domain	Missi
Sampling_station_ID		Accurate sampling location in a site.	string	nominal	**Def** Accurate sampling location in a site.	
Sampling_station_latitude_value		Latitude of the sampling station.	float	ratio	**Unit** degree **Type** real **Min** 42.07478 **Max** 46.20382	
Sampling_station_longitude_value		Longitude of the sampling station.	float	ratio	**Unit** degree **Type** real **Min** -1.19539 **Max** 9.30593	
Sampling_station_geonames_name		Longitude of the sampling station.	string	nominal	**Def** Longitude of the sampling station.	**Code** **Expl**
Sampling_station_common_name		Name of the location of the sampling used by the researchers in their publications.	string	nominal	**Def** Name of the location of the sampling used by the researchers in their publications.	
Sampling_station_administrator_name		First name and surname of the location administrator.	string	nominal	**Def** First name and surname of the location administrator.	
Sampling_station_label_name		If applicable, the name of the environmental label(s) obtained by the location.	string	nominal	**Def** If applicable, the name of the environmental label(s) obtained by the location.	**Code** **Expl**
Sampling_station_berth_number		Number of berths in the port in 2018 according to the port administrator.	string	nominal	**Def** Number of berths in the port in 2018 according to the port administrator.	
Sampling_station_municipality_name		Municipality in which the location where the sampling took place is located.	string	nominal	**Def** Municipality in which the location where the sampling took place is located.	

Data Table:

Name:	**TAXREF.csv**
Description:	Content of TAXREF.csv

Physical Structure Description:

Object Name:	TAXREF.csv
Size:	89 bytes
Authentication:	2ae58fe83b9b0a8ba61932947be75d0d Caculated By MD5
Text Format:	Number of Header Lines: 1 Record Delimiter: \r\n Attribute Orientation: column Simple Delimited: Field Delime
Number Of Records:	4

Attribute(s) Info:

Name	Column Label	Definition	Type of Value	Measurement Type	Measurement Domain
TAXREF_ID		Identifier of the sampled organism species if known according to the taxonomic repository TAXREF. If the species is not known, the genus given by the TAXREF is used, followed by sp..	string	nominal	**Def** Identifier of the sampled organism species if known accordi taxonomic repository TAXREF. If the species is not known, genus given by the TAXREF is used, followed by sp..

Data Table:

Name:	**Water_samples_measurements.csv**
Description:	Water_samples_measurements.csv

Physical Structure Description:

Object Name:	Water_samples_measurements.csv
Size:	55049 bytes
Authentication:	eeafa7536d922594288eb8c210e3ed52 Caculated By MD5
Text Format:	Number of Header Lines: 1 Record Delimiter: \r\n Attribute Orientation: column Simple Delimited: Field Delime
Number Of Records:	133

Attribute(s) Info:

Name	Column Label	Definition	Type of Value	Measurement Type	Measurement Domain
Water_sample_ID		Unique identifier of a water sample.	string	nominal	**Def** Unique identifier of a water sample.
Fieldwork_ID		Unique identifier of the fieldwork carried out on a specific date and located at a specific station.	string	nominal	**Def** Unique identifier of the fieldwork carried out on a specific date and located at a specific station.
Water_subsurface_temperature_celsius		Temperature of the subsurface water at the sampling station	float	ratio	**Unit** celsius **Type** real

		measured in celsius degrees.			**Min** 5.28 **Max** 24.71
Water_subsurface_salinity_psu		Salinity of the subsurface water at the sampling station measured in practical salinity unit.	float	ratio	**Unit** practicalSalinityUnit **Type** real **Min** 22.87 **Max** 38.45
Water_subsurface_dissolved_oxygen_mg_per_L		The concentration of oxygen in the subsurface water in mg per liter of water measured at the sampling station.	float	ratio	**Unit** milligramsPerLiter **Type** real **Min** 3.73 **Max** 31.46
Water_subsurface_pH		The chemical activity of the protons in the subsurface water at the sampling station.	float	ratio	**Unit** dimensionless **Type** real **Min** 7.68 **Max** 8.99
Water_subsurface_turbidity_fnu		Material content of the subsurface water at the sampling station measured in Formazin Nephelometric Unit.	float	ratio	**Unit** formazineNephelometricUnit **Type** real **Min** 1 **Max** 194
Water_subsurface_chlorophyll_a_concentration_µg_per_L		Concentration of chlorophyll a measured in microgram per liter of subsurface water taken at the sampling station.	float	ratio	**Unit** microgramsPerLiter **Type** real **Min** 0.18 **Max** 2.94
Water_subsurface_nitrite_concentration_µM		Concentration of nitrite measured in micromol per liter of subsurface water taken at the sampling station.	float	ratio	**Unit** micromolePerLiter **Type** real **Min** 0.03 **Max** 0.47

Water_subsurface_nitrate_concentration_µM		Concentration of nitrate measured in micromol per liter of subsurface water taken at the sampling station.	float	ratio	**Unit** micromolePerLiter **Type** real **Min** 0.16 **Max** 14.42
Water_subsurface_silica_concentration_µM		Concentration of silica measured in micromol per liter of subsurface water taken at the sampling station.	float	ratio	**Unit** micromolePerLiter **Type** real **Min** 0.29 **Max** 75.76
Water_subsurface_phosphate_concentration_µM		Concentration of phosphate measured in micromol per liter of subsurface water.	float	ratio	**Unit** micromolePerLiter **Type** real **Min** 0.02 **Max** 0.63

Water_subsurface_ammonium_concentration_µM		Concentration of ammonium measured in micromol per liter of subsurface water taken at the sampling station.	float	ratio	**Unit** micromolePerLiter **Type** real **Min** 0.11 **Max** 5.21
Water_surface_Ag_concentration_µg_per_L		Concentration of silver in microgram per liter of surface water.	float	ratio	**Unit** microgramsPerLiter **Type** real
Water_surface_Al_concentration_µg_per_L		Concentration of aluminium in microgram per liter of surface water.	float	ratio	**Unit** microgramsPerLiter **Type** real **Min** 223.5 **Max** 223.5
Water_surface_As_concentration_µg_per_L		Concentration of arsenic in microgram per liter of surface	float	ratio	**Unit** microgramsPerLiter **Type** real

		water.			**Min** 14.84 **Max** 14.84
Water_surface_Ba_concentration_µg_per_L		Concentration of barium in microgram per liter of surface water.	float	ratio	**Unit** microgramsPerLiter **Type** real
Water_surface_Cd_concentration_µg_per_L		Concentration of cadmium in microgram per liter of surface water.	float	ratio	**Unit** microgramsPerLiter **Type** real
Water_surface_Co_concentration_µg_per_L		Concentration of cobalt in microgram per liter of surface water.	float	ratio	**Unit** microgramsPerLiter **Type** real **Min** 1.02 **Max** 1.66
Water_surface_Cr_concentration_µg_per_L		Concentration of chromium in microgram per liter of	float	ratio	**Unit** microgramsPerLiter **Type** real

		surface water.			
Water_surface_Cu_concentration_µg_per_L		Concentration of copper in microgram per liter of surface water.	float	ratio	**Unit** micrograms**P**er**L**iter **Type** real **Min** 2.56 **Max** 18.00
Water_surface_Fe_concentration_µg_per_L		Concentration of iron in microgram per liter of surface water.	float	ratio	**Unit** micrograms**P**er**L**iter **Type** real
Water_surface_In_concentration_µg_per_L		Concentration of indium in microgram per liter of surface water.	float	ratio	**Unit** micrograms**P**er**L**iter **Type** real

Water_surface_Mn_concentration_µg_per_L		Concentration of manganese in microgram per liter of surface water.	float	ratio	**Unit** micrograms**P**er**L**iter **Type** real **Min** 1.10 **Max** 18.94
Water_surface_Mo_concentration_µg_per_L		Concentration of molybdenum in microgram per liter of surface water.	float	ratio	**Unit** micrograms**P**er**L**iter **Type** real **Min** 11.02 **Max** 12.72
Water_surface_Ni_concentration_µg_per_L		Concentration of nickel in microgram per liter of surface water.	float	ratio	**Unit** micrograms**P**er**L**iter **Type** real
Water_surface_Pb_concentration_µg_per_L		Concentration of lead in microgram per liter of surface water.	float	ratio	**Unit** micrograms**P**er**L**iter **Type** real **Min** 1.14 **Max** 8.06

Water_surface_Sb_concentration_µg_per_L		Concentration of tin in microgram per liter of surface water.	float	ratio	**Unit** micrograms_per_liter **Type** real **Min** 2.14 **Max** 2.86
Water_surface_Se_concentration_µg_per_L		Concentration of selenium in microgram per liter of surface water.	float	ratio	**Unit** micrograms_per_liter **Type** real **Min** 47.82 **Max** 85.16
Water_surface_Sn_concentration_µg_per_L		Concentration of antimony in microgram per liter of surface water.	float	ratio	**Unit** micrograms_per_liter **Type** real
Water_surface_V_concentration_µg_per_L		Concentration of vanadium in microgram per liter of surface water.	float	ratio	**Unit** micrograms_per_liter **Type** real

Water_surface_Zn_concentration_µg_per_L		Concentration of zinc in microgram per liter of surface water.	float	ratio	**Unit** micrograms**P**er**L**iter **Type** real
Water_surface_naphthalene_concentration_ng_per_L		Concentration of naphthalene in nanogram per liter of surface water.	float	ratio	**Unit** nanograms**P**er**L**iter **Type** real **Min** 6.8 **Max** 65.4
Water_surface_C1_naphthalene_concentration_ng_per_L		Concentration in nanogram per liter of surface water of a group of molecules	float	ratio	**Unit** nanograms**P**er**L**iter **Type** real **Min** 11.4 **Max** 13.2

		with a common empirical formula, i.e. naphthalene with the addition of a methyl (CH3), but with a different structural formula, as the methyl may be bonded to different atoms of the napthalene.			
Water_surface_C2_naphthalene_concentration_ng_per_L		Concentration in nanogram per liter of surface water of a group of chemical molecules with a common empirical formula, i.e. naphthalene with the addition of two methyl (CH3) or one ethyl (C2H5), but with a different structural formula, as the methyls or the ethyl may be bonded to different atoms of the napthalene.	float	ratio	**Unit** nanogramsPerLiter **Type** real **Min** 5.3 **Max** 26.2
Water_surface_C3_naphthalene_concentration_ng_per_L		Concentration in nanogram per liter of surface water of a group of chemical molecules with a common empirical formula, i.e. naphthalene plus three methyls (CH3), one ethyl (C2H5) and one methyl, or one propyl (C3H7), but with a different structural formula, as the methyls, the ethyl and the methyl, or the propyl may be bonded to different atoms of the napthalene.	float	ratio	**Unit** nanogramsPerLiter **Type** real **Min** 5.2 **Max** 23.2
Water_surface_benzothiophene_concentration_ng_per_L		Concentration of benzothiophene in nanogram per liter of surface water.	float	ratio	**Unit** nanogramsPerLiter **Type** real

Water_surface_C1_benzothiophene_concentration_ng_per_L		Concentration in nanogram per liter of surface water of a group of molecules with a common empirical formula, i.e. benzothiophene with the addition of a methyl (CH3), but with a different structural formula, as the methyl may be bonded to different atoms of the benzothiophene.	float	ratio	**Unit** nanogramsPerLiter **Type** real
Water_surface_C2_benzothiophene_concentration_ng_per_L		Concentration in nanogram per liter of surface water of a group of chemical molecules with a common empirical formula, i.e. benzothiophene with the addition of two methyl (CH3) or one ethyl (C2H5), but with a different structural formula, as the methyls or the ethyl may be bonded to different atoms of the benzothiophene.	float	ratio	**Unit** nanogramsPerLiter **Type** real
Water_surface_C3_benzothiophene_concentration_ng_per_L		Concentration in nanogram per liter of surface water of a group of chemical molecules with a common empirical formula, i.e. benzothiophene plus three methyls (CH3), one ethyl (C2H5) and one methyl, or one propyl (C3H7), but with a different structural formula, as the methyls, the ethyl and the methyl, or the propyl may be bonded to different atoms of the benzothiophene.	float	ratio	**Unit** nanogramsPerLiter **Type** real
Water_surface_biphenyl_concentration_ng_per_L		Concentration of biphenyl in nanogram per liter of surface water.	float	ratio	**Unit** nanogramsPerLiter **Type** real

Water_surface_acenapthtylene_concentration_ng_per_L		Concentration of acenapthtylene in nanogram per liter of surface water.	float	ratio	**Unit** nanogramsPerLiter **Type** real **Min** 2 **Max** 2
Water_surface_acenaphthene_concentration_ng_per_L		Concentration of acenaphtene in nanogram per liter of surface water.	float	ratio	**Unit** nanogramsPerLiter **Type** real
Water_surface_fluorene_concentration_ng_per_L		Concentration of fluorene in nanogram per liter of surface water.	float	ratio	**Unit** nanogramsPerLiter **Type** real **Min** 7.2 **Max** 7.2

Water_surface_C1_fluorene_concentration_ng_per_L		Concentration in nanogram per liter of surface water of a group of molecules with a common empirical formula, i.e. fluorene with the addition of a methyl (CH3), but with a different structural formula, as the methyl may be bonded to different atoms of the fluorene.	float	ratio	**Unit** nanogramsPerLiter **Type** real
Water_surface_C2_fluorene_concentration_ng_per_L		Concentration in nanogram per liter of surface water of a group of chemical molecules with a common empirical formula, i.e. fluorene with the addition of two methyl (CH3) or one ethyl (C2H5), but with a different structural formula, as the methyls or the ethyl may be bonded to different atoms of the fluorene.	float	ratio	**Unit** nanogramsPerLiter **Type** real **Min** 5.4 **Max** 9.2
Water_surface_C3_fluorene_concentration_ng_per_L		Concentration in nanogram per liter of surface water of a group of chemical molecules with a common empirical formula, i.e. fluorene plus three methyls (CH3), one ethyl (C2H5) and one methyl, or one propyl (C3H7), but with a	float	ratio	**Unit** nanogramsPerLiter **Type** real **Min** 5.4 **Max** 5.4

		different structural formula, as the methyls, the ethyl and the methyl, or the propyl may be bonded to different atoms of the fluorene.			
Water_surface_phenanthrene_concentration_ng_per_L		Concentration of phenanthrene in nanogram per liter of surface water.	float	ratio	**Unit** nanogramsPerLiter **Type** real
Water_surface_anthracene_concentration_ng_per_L		Concentration of anthracene in nanogram per liter of surface water.	float	ratio	**Unit** nanogramsPerLiter **Type** real
Water_surface_C1_phenanthrene_and_anthracene_concentration_ng_per_L		Concentration in nanogram per liter of surface water of a group of molecules with a common empirical formula, i.e. phenanthrene or anthracene with the addition of a methyl (CH3), but with a different structural formula, as the methyl may be bonded to different atoms of the phenanthrene or the anthracene.	float	ratio	**Unit** nanogramsPerLiter **Type** real

Water_surface_C2_phenanthrene_and_anthracene_concentration_ng_per_L		Concentration in nanogram per liter of surface water of a group of chemical molecules with a common empirical formula, i.e. phenanthrene or anthracene with the addition of two methyl (CH3) or one ethyl (C2H5), but with a different structural formula, as the methyls or the ethyl may be bonded to different atoms of the phenanthrene or the anthracene.	float	ratio	**Unit** nanogramsPerLiter **Type** real
Water_surface_C3_phenanthrene_and_anthracene_concentration_ng_per_L		Concentration in nanogram per liter of surface water of a group of chemical molecules with a common empirical formula, i.e. phenanthrene or anthracene plus three methyls (CH3), one ethyl (C2H5) and one methyl, or one propyl (C3H7), but with a different structural formula, as the methyls, the ethyl and the methyl, or the propyl may be bonded to different atoms of the phenanthrene or the anthracene.	float	ratio	**Unit** nanogramsPerLiter **Type** real
Water_surface_dibenzothiophene_concentration_ng_per_L		Concentration of dibenzothiophene in nanogram per liter of surface water.	float	ratio	**Unit** nanogramsPerLiter **Type** real

Water_surface_C1_dibenzothiophene_concentration_ng_per_L		Concentration in nanogram per liter of surface water of a group of molecules with a common empirical formula, i.e. dibenzothiophene with the addition of a methyl (CH3), but with a different structural formula, as the methyl may be bonded to different atoms of the dibenzothiophene.	float	ratio	**Unit** nanogramsPerLiter **Type** real
Water_surface_C2_dibenzothiophene_concentration_ng_per_L		Concentration in nanogram per liter of surface water of a group of chemical molecules with a common empirical formula, i.e. dibenzothiophene with the addition of two methyl (CH3) or one ethyl (C2H5), but with a different structural formula, as the methyls or the ethyl may be bonded to different atoms of the dibenzothiophene.	float	ratio	**Unit** nanogramsPerLiter **Type** real
Water_surface_C3_dibenzothiophene_concentration_ng_per_L		Concentration in nanogram per liter of surface water of a group of chemical molecules with a common empirical formula, i.e. dibenzothiophene plus three methyls (CH3), one ethyl (C2H5) and one methyl, or one propyl (C3H7), but with a different structural formula, as the methyls, the ethyl and the methyl, or the propyl may be bonded to different atoms of the dibenzothiophene.	float	ratio	**Unit** nanogramsPerLiter **Type** real
Water_surface_fluoranthene_concentration_ng_per_L		Concentration of fluoranthene in nanogram per liter of	float	ratio	**Unit** nanogramsPerLiter **Type** real

		surface water.			
Water_surface_pyrene_concentration_ng_per_L		Concentration of pyrene in nanogram per liter of surface water.	float	ratio	**Unit** nanogramsPerLiter **Type** real
Water_surface_C1_fluoranthene_and_pyrene_concentration_ng_per_L		Concentration in nanogram per liter of surface water of a group of molecules with a common empirical formula, i.e. fluoranthene or pyrene with the addition of a methyl (CH3), but with a different structural formula, as the methyl may be bonded to different atoms of the fluoranthene or the pyrene.	float	ratio	**Unit** nanogramsPerLiter **Type** real
Water_surface_C2_fluoranthene_and_pyrene_concentration_ng_per_L		Concentration in	float	ratio	**Unit** nanogramsPerLiter

		nanogram per liter of surface water of a group of chemical molecules with a common empirical formula, i.e. fluoranthene or pyrene with the addition of two methyl (CH3) or one ethyl (C2H5), but with a different structural formula, as the methyls or the ethyl may be bonded to different atoms of the fluoranthene or the pyrene.			**Type** real
Water_surface_C3_fluoranthene_and_pyrene_concentration_ng_per_L		Concentration in nanogram per liter of surface water of a group of chemical molecules with a common empirical formula, i.e. fluoranthene or pyrene plus three methyls (CH3), one ethyl (C2H5) and one methyl, or one propyl (C3H7), but with a different structural formula, as the methyls, the ethyl and the methyl, or the propyl may be bonded to different atoms of the fluoranthene or the pyrene.	float	ratio	**Unit** nanogramsPerLiter **Type** real
Water_surface_benz_a_anthracene_concentration_ng_per_L		Concentration of benz[a]anthracene in nanogram per liter of surface water.	float	ratio	**Unit** nanogramsPerLiter **Type** real
Water_surface_chrysene_concentration_ng_per_L		Concentration of chrysene in nanogram per liter of surface water.	float	ratio	**Unit** nanogramsPerLiter **Type** real **Min** 1.7

					Max 1.7
Water_surface_C1_chrysene_concentration_ng_per_L		Concentration in nanogram per liter of surface water of a group of molecules with a common empirical formula, i.e. chrysene with the addition of a methyl (CH3), but with a different structural formula, as the methyl may be bonded to different atoms of the chrysene.	float	ratio	**Unit** nanogramsPerLiter **Type** real
Water_surface_C2_chrysene_concentration_ng_per_L		Concentration in nanogram per liter of surface water of a group of chemical molecules with a common empirical formula, i.e. chrysene with the addition of two methyl (CH3) or one ethyl (C2H5), but with a different structural formula, as the methyls or the ethyl may be bonded to different atoms of the chrysene.	float	ratio	**Unit** nanogramsPerLiter **Type** real
Water_surface_C3_chrysene_concentration_ng_per_L		Concentration in nanogram per liter of surface water of a group of chemical molecules with a common empirical formula, i.e. chrysene plus three methyls (CH3), one ethyl (C2H5) and one methyl, or one propyl (C3H7), but with a different structural formula, as the	float	ratio	**Unit** nanogramsPerLiter **Type** real

		methyls, the ethyl and the methyl, or the propyl may be bonded to different atoms of the chrysene.			
Water_surface_benzo_b_k_fluranthene_concentration_ng_per_L		Concentration of benzo[b+k]fluranthene in nanogram per liter of surface water.	float	ratio	**Unit** nanogramsPerLiter **Type** real **Min** 1.9 **Max** 4.1
Water_surface_benzo_e_pyrene_concentration_ng_per_L		Concentration of benzo[e]pyrene in nanogram per liter of surface water.	float	ratio	**Unit** nanogramsPerLiter **Type** real **Min** 2.3 **Max** 2.3
Water_surface_benzo_a_pyrene_concentration_ng_per_L		Concentration of benzo[a]pyrene in nanogram per liter of surface water.	float	ratio	**Unit** nanogramsPerLiter **Type** real **Min** 1 **Max** 1
Water_surface_perylene_concentration_ng_per_L		Concentration of perylene in nanogram per liter of surface water.	float	ratio	**Unit** nanogramsPerLiter **Type** real **Min** 1.1 **Max** 1.1

Water_surface_indeno_1_2_3_cd_pyrene_concentration_ng_per_L		Concentration of indeno[1,2,3-cd]pyrene in nanogram per liter of surface water.	float	ratio	**Unit** nanogramsPerLiter **Type** real
Water_surface_dibenz_a_h_anthracene_concentration_ng_per_L		Concentration of dibenz[a,h]pyrene in nanogram per liter of surface water.	float	ratio	**Unit** nanogramsPerLiter **Type** real
Water_surface_benzo_ghi_perylene_concentration_ng_per_L		Concentration of benzo[ghi]perylene in nanogram per liter of surface water.	float	ratio	**Unit** nanogramsPerLiter **Type** real

Water_measurements_protocol_DOI		Digital object identifier of the publication where the protocol followed for the water sampling may be found.	string	nominal	**Def** Digital object identifier of the publication where the protocol followed for the water sampling may be found.

Other Entity:

Name:	**ConceptualDataModel.jpg**
Data Object Type:	image/png

Other Entity:

Name:	**WIBE_Data_Dictionary.csv**
Data Object Type:	text/csv

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Data Set Characteristics

Geographic Region:

Geographic Description:	Calvi_port
Bounding Coordinates:	West: 8.76163 degrees East: 8.76163 degrees North: 42.56726 degrees South: 42.56726 degrees

Geographic Region:

Geographic Description:	Calvi_fairing
Bounding Coordinates:	West: 8.75694 degrees

	East: 8.75694 degrees
	North: 42.56359 degrees
	South: 42.56359 degrees

Geographic Region:

Geographic Description:	Calvi_oil
Bounding Coordinates:	West: 8.75768 degrees
	East: 8.75768 degrees
	North: 42.56439 degrees
	South: 42.56439 degrees

Geographic Region:

Geographic Description:	Ile_rousse_port
Bounding Coordinates:	West: 8.93639 degrees
	East: 8.93639 degrees
	North: 42.64151 degrees
	South: 42.64151 degrees

Geographic Region:

Geographic Description:	Ile_rousse_fairing
Bounding Coordinates:	West: 8.93554 degrees
	East: 8.93554 degrees
	North: 42.64023 degrees
	South: 42.64023 degrees

Geographic Region:

Geographic Description:	Ile_rousse_oil
Bounding Coordinates:	West: 8.9359129 degrees
	East: 8.9359129 degrees
	North: 42.6406397 degrees
	South: 42.6406397 degrees

Geographic Region:

Geographic Description:	St_florent_port
Bounding Coordinates:	West: 9.30031 degrees
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	North: 42.67993 degrees
	South: 42.67993 degrees

Geographic Region:

Geographic Description:	St_florent_fairing
Bounding Coordinates:	West: 9.3009414 degrees
	East: 9.3009414 degrees

North:	42.6798025 degrees
South:	42.6798025 degrees

Geographic Region:

Geographic Description:	St_florent_oil
Bounding Coordinates:	
West:	9.29801 degrees
East:	9.29801 degrees
North:	42.68093 degrees
South:	42.68093 degrees

Geographic Region:

Geographic Description:	STARESO_port
Bounding Coordinates:	
West:	8.72542 degrees
East:	8.72542 degrees
North:	42.58044 degrees
South:	42.58044 degrees

Geographic Region:

Geographic Description:	Diana_mussel_farm
Bounding Coordinates:	
West:	9.536647 degrees
East:	9.5366473 degrees
North:	42.131517 degrees
South:	42.131517 degrees

Geographic Region:

Geographic Description:	L'Houmeau_marsh_reference
Bounding Coordinates:	
West:	-1.19539 degrees
East:	-1.19539 degrees
North:	46.20382 degrees
South:	46.20382 degrees

Geographic Region:

Geographic Description:	Minimes_fairing
Bounding Coordinates:	
West:	-1.16187 degrees
East:	-1.16187 degrees
North:	46.14675 degrees
South:	46.14675 degrees

Geographic Region:

Geographic Description:	Minimes_rain_waters
Bounding Coordinates:	
West:	-1.16605 degrees
East:	-1.16605 degrees
North:	46.14143 degrees

South:	46.14143 degrees

Geographic Region:

Geographic Description:	Minimes_oil
Bounding Coordinates:	West: -1.16749 degrees East: -1.16749 degrees North: 46.14688 degrees South: 46.14688 degrees

Time Period:

Begin:	2019-06-03
End:	2022-12-31

Taxonomic Range:

Classification: 1

Rank Name:	kingdom
Rank Value:	Animalia
Classification:	50

Rank Name:	phylum
Rank Value:	Echinodermata
Classification:	222

Rank Name:	class
Rank Value:	Holothuroidea
Classification:	9505681

Rank Name:	order
Rank Value:	Holothuriida
Classification:	9548

Rank Name:	
Rank Value:	
Classification:	

Classification: 1

Rank Name:	kingdom
Rank Value:	Animalia
Classification:	52

Rank Name:	phylum
Rank Value:	Mollusca
Classification:	137

Rank Name:	class

Rank Value:	Bivalvia
Classification:	9250425
Rank Name:	order
Rank Value:	Pectinida
Classification:	3472
Rank Name:	family
Rank Value:	Pectinidae
Classification:	2285867
Rank Name:	
Rank Value:	
Classification:	

Classification:	1
Rank Name:	kingdom
Rank Value:	Animalia
Classification:	52
Rank Name:	phylum
Rank Value:	Mollusca
Classification:	137
Rank Name:	class
Rank Value:	Bivalvia
Classification:	9330464
Rank Name:	order
Rank Value:	Mytilida
Classification:	3476
Rank Name:	family
Rank Value:	Mytilidae
Classification:	2285679
Rank Name:	ge
Rank Value:	M
Classification:	22

Classification:	5
Rank Name:	kingdom
Rank Value:	Fungi
Classification:	95

		Rank Name:	phylum
		Rank Value:	Ascomycota
		Classification:	316
			Rank Name: class
			Rank Value: Pezizomycetes
			Classification: 1057
			Rank Name: order
			Rank Value: Pezizales
			Classification: 4131
			Rank Name:
			Rank Value:
			Classification:
Classification:	Rank Value:	Holothuria sp.	
Classification:	Rank Value:	Mimachlamys varia	
Classification:	Rank Value:	Mytilus galloprovincialis	
Classification:	Rank Value:	Patella sp.	

Sampling, Processing and Quality Control Methods

Step by Step Procedures

Step 1:	
Description:	## Method of the QUAMPO project: Three species (*Mytilus galloprovincialis, Patella sp. and Holothuria sp.*) w different ports of North Corsica (France). St-Florent (42.67993, 9.30031), Ile Rousse (42.64151, 8.93639) and C site while the STARESO (42.58044, 8.72542) was chosen as the control site. ## Method of the QUALIPERTUIS Minimes* (La Rochelle, Region Nouvelle-Aquitaine). These sites were chosen following previous studies carried anthropogenic contamination: careening station (CAR), fuel station (NFS), storm water outfall (PLU) (Breitwiese 2017). The reference site selected for this study is a former oyster farm, in a semi-open environment, and not su
Sampling Area And Frequency:	## QUAMPO In situ and caging experiment In January 2020, mussels (n=17) were found only in St-Florent port everywhere except in St-Florent port. Sea cucumbers (n = 4-7 per location) were collected in the four ports in Ja afterwards. In September 2020, limpets (n = 15-16 per location) were collected in the four ports. As mussels we were used for the following collections. Mussels were bought from the mussel farm of the Etang de Diana in Jur then installed in the different ports for three months. In September 2020, they (n = 16 per location) were collecte mussels survived and were used for bioaccumulation analyses. In January 2021, limpets (n = 10 in each locatio per location) were collected in the ports of Calvi, St-Florent and Île Rousse but not in STARESO. In September collected in all of the four ports. Finally, in January 2022, mussels (n = 20) were collected in each port except in in the four studied ports. ### Test on the “pool” techniques During the January 2020 sampling, additional organi “individual” sampling techniques: 13 mussels (length = 68.56 ± 8.84 mm) and 4 sea cucumber (length = 193.25 (length = 32.34 ± 4.93 mm) at Île Rousse. To test the “pool” sampling technique, the tissues of 7 individuals fron (limpet) were dissected, pooled and processed as detailed below for all the analyses (enzymes and bioaccumul biological model used for this study is the same as that used in previous publications: the black scallop Mimachl caging. For this, individuals were fished at the Loix site (Ile de Ré, less subject to port influences, (Breitwieser et

	They are caged in the reference site (Marsh) for a minimum of 15 days to acclimate before being transplanted to the individuals were counted in order to assess mortality, then seven individuals per site were taken (the D0 sam in liquid nitrogen in order to preserve enzymatic activities. Low mortality (less than 10 %) at both sites was obse -80°C before being prepared for biochemical analysis. Once in the laboratory, the individuals are processed in a recovered from each sample become one sample). A previous study compared pooled and individual data, the r the two methods (Breitwieser et al., 2018). Statistical analysis is performed using R software (RStudio, version 1 by seasons (7), sites (4) and sampling days (4). For all variables, the homogeneity of variances is assessed usi Shapiro-Wilk test.
Sampling Description:	## QUAMPO ### Tissue sampling After sampling, individuals were immediately placed in dry ice and shipped to lab, half (n = ~7) of the mussels, limpets and sea cucumber tissues were kept at -20°C for trace element bioacc for all the other analyses. ### Biomarkers of exposure (= trace element bioaccumulation) The soft tissues of mu and pestle. The body wall of sea cucumber was cut in pieces as small as possible. The samples were freeze dri materials (DOLT-5, dogfish liver, and TORT-3, lobster hepatopancreas, National Research Council Canada) we (v/v) 67-70% HNO3/34-37% HCl mixture (Fisher, trace metal quality). The digestion process was carried out ov digestion (30 min with constantly increasing temperature up to 120 °C, then 15 min at this temperature). At the e water until 50 mL. aluminium (Al), silver (Ag), arsenic (As), barium (Ba), cadmium (Cd), cobalt (Co), chromium ((Mo), nickel (Ni), lead (Pb), antimony (Sb), selenium (Se), tin (Sn), vanadium (V) and zinc (Zn) contents were ar California, USA) and an ICP-MS XSeries 2 (ThermoFisher Scientific, Waltham, 192 Massachusetts, USA).Mear 90.5% (Cd), 89.7% (Co), 89% (Cr), 91.2% (Cu), 93.5% (Fe), 90% (Mn), 87.7% (Mo), 81% (Ni), 88% (Pb), 96.7% 19.28 (Al), 0.01 (Ag), 0.1 (As), 0.1 (Ba), 0.01 (Cd), 0.01 (Co), 0.1 (Cr), 0.96 (Cu), 3.86 (Fe), 0.01 (Mn), 0.1 (Mo), (V) and 3.86 (Zn) Î¼g.g-1 of dry weight. The trace element contamination of the water samples was also analys seawater standards, in 2 % nitric acid) and ICP-MS (dilution to 1/20 for water samples and seawater standards, Tissues samples processing Individuals were dissected to keep the digestive gland for mussel, the whole soft tis tissues were weighted grinded in liquid nitrogen using a MM400 Retsch© (GmbH, Éragny, Luxemburg) mixer m phosphate buffered saline (PBS) solution (100 mM, pH containing 0.1 % Triton X-100 and 1 mM ethylenediamin samples were centrifuged 15 min at 12500 g at 4°C (Sorvall Legeng Micro 17R, ThermoFisher Scientific, Waltha divided into aliquots and used for the subsequent biomarkers of effect and protein analyses. #### Enzymatic bio evaluated using superoxide dismutase (SOD), catalase (CAT) and glutathione peroxidase (GPx) activities. SOD superoxide-driven NADH oxidation and was measured following the rate of NADH oxidation at 340 nm (Paoletti (1984) by measuring the rate of decomposition of hydrogen peroxide at 240 nm. GPx activity was assayed acco Janssens et al. (2000) by monitoring the consumption of NADPH at 340 nm. The glutathione S-transferase (GST detoxification process, were respectively determined by measuring the production of GS-DNB at 340 nm (Habig 340 nm (Carlberg and Mannervik 1985). Acetylcholinesterase (AChE) is widely used to estimate neurotoxic imp carbamates, several heavy metals and surfactants. Its activity was assayed by recording the production of 5-thio the laccase-type phenoloxidase (LAC) is involved in the immune, antioxidant and detoxification processes (Luna measured according to (Luna-Acosta et al. 2010) y following the oxidation of p-phenylenediamine at 420 nm. Th acid protein assay kit (Sigma-Aldrich®) that is similar to the Lowry et al. (1951) procedure. All activities measure spectrophotometer (BMG labtech, Champigny-sur-Marne, France). All chemicals were obtained from Sigma-Ald standard methods adapted for a microplate reader. For all enzymatic analysis, homogenates were diluted to obt 25°C. Total and specific enzymatic activities were measured and expressed as 242 U (Î¼mol.min1).g1 of wet tis peroxidation Malondialdehyde (MDA) content, a secondary product of lipid peroxidation, was quantified using th following Heath and Packer (1968) and Bird and Draper (1984), adapted by Torres et al. (2020). This method es thiobarbituric acid. The thiobarbituric acid-malondialdehyde complex produced during the reaction is followed by nonspecific turbidity is applied by recorded the absorbance at 600 nm. #### Integrated biomarker response (IBR (IBR, (Devin et al. 2014)) was calculated using the response of biomarkers of effect. On R software, the functior are based on Beliaeff and Burgeot (2002) revisited by Devin et al. (2014). ## QUALIPERTUIS All biochemical a using the same instrument: SpectroStar Nano with specific absorbance for each type of biomarker. The assays varia) from all the study sites at D0, D07, D21 and D30 (one pool (n=7 individuals) per site and per date). Data a Data obtained for MDA are in Î¼M.mg protein. Total protein concentrations are determined using the BCA kit m BCA kit contains bovine serum albumin (BSA) as a standard and involves the reduction of alkaline Cu(II) proteir nm. Superoxide dismutase (SOD) activity plays an important role in the oxidative stress response (Paoletti et al. et al., 2017). The specific activity of glutathione S-transferase (GST) plays a major role in the detoxification of xe enzymes) is determined according to the Sigma kit method (CS0410-1 KT) (Breitwieser et al., 2017). In addition by quantifying malondialdehyde (MDA), a chemical metabolite of cellular lipid breakdown. The concentration of International) (Breitwieser et al., 2017; Milinkovitch et al., 2015). AChE is an enzyme involved in the neurotransm acetylcholine to choline and acetic acid. Its inhibition is directly related to the mechanism of toxic action of organ on the inhibition of cholinesterase activity as a diagnostic tool for organophosphorus pesticide contamination in (Galgani et al., 1992; Zinkl et al., 1987), but few studies have been carried out on marine invertebrates, which a

	programs (Day and Scott, 1990). AChE activity (IU.mg protein-1) was measured using the AcetylCholinEsterase phenoloxidase (Laccase) activity reveals an alteration of the bivalve immune system. Its enzymatic activity is se Luna-Acosta et al., 2010a) and plays a crucial role in the immune defense mechanism of marine invertebrates. I (Breitwieser et al., 2017). ## Environmental parameters ### QUAMPO One water sample was taken from each element contamination of the environment. The water samples were collected using a Niskin bottle and Falcon t (temperature, salinity, dissolved oxygen and pH) were measured in-situ using a YSI-Exo2 multi-parameters prob nutrient concentrations as well as chlorophyll a concentrations are described by Fullgrabe et al. (2020) while the (PCBs, PAHs) are described by Frantzen et al. (2016). The trace element contamination of the water samples w samples and seawater standards, in 2 % nitric acid) and ICP-MS (dilution to 1/20 for water samples and seawat each day of sampling (D0, D07, D21, D30), the physico-chemical parameters of the water (Temperature (°C) an (multiparameter HI9829). The probe was calibrated before each sampling period.
Quality Control Step 1:	
Description:	## QUAMPO Statistics analyses ### Test on the pool technique To test the pool technique, the mean of the ind the pooled samples. The percentage of difference between these two values was compared for effect and biom homogeneity of variances were verified by Shapiro and Levene tests and data were log transformed to avoid he case of heteroscedasticity) correlations between total and specific enzymatic activity responses were verified fo ANOVAs were used to test the effects of location and sampling period on the biomarkers of effect. When signific to compare means (Î± = 0.05). Except for LAC activity in limpet, no significant interaction between exposure and effect, analyses for simple main effects were then applied to the data (Tybout and Sternthal 2001). Student t-tes applied to test the effect of sampling period on each location. One-way ANOVAs (or non-parametric Kruskall-Wa for each sampling period, followed by respectively Tukey or Mann-Whitney Wilcoxon post-hoc test when a signi location in January, Student t-tests (or non-parametric Mann-Whitney Wilcoxon test) were applied to test the eff ANOVAs (or non-parametric Kruskall-Wallis tests), followed by respectively Tukey or Mann-Whitney Wilcoxon p performed to analyse the effect of location in September on mussel. Finally, multi-correlations were run betwee parameters of individuals to visualize the link between the response of organism and its condition. ### Biomark January samples, the September samples were processed in pool. First, the normality and homogeneity of varia effect of location in January, one-way ANOVAs (or non-parametric Kruskall-Wallis tests) were performed, follow post-hoc test when a significant effect was found. ### Environmental parameters Graphs and analysis were dor were weekly regularized in order to obtain 48 data per year (four per month) using the pastecs R package (regu (https://github.com/phgrosjean/pastecs). Average abundances over different time periods (year, month, week) w dynamics were described by integrating the time series over one year using boxplot diagrams in order to observ (described as data point standing 1.5 times outside the interquartile range above the upper quartile and below th group were further explored by computing the smoothed annual variation of the median, 10th and 90th percentil highlight both annual and interannual variability. Smoothing was performed using the lowess R function which u 1981). Further on, cumulative relative abundances were computed over an average year in order to observe the order to describe the succession of the peaks of absolute or relative abundance of each group over an average cluster analysis on zooplankton samples based on the complete linkage method using Bray-Curtis dissimilarity (grouped depending on both abundance and general composition similarities. A cut at a distance of 0.80 in the re each resulting cluster suggesting a different community state. These eight community states were labelled A to et al., 2009) was then applied. The analysis used as criteria of classification both absolute and relative abundan decision node of the resulting classification tree corresponded to a condition to which zooplankton samples migl zooplankton samples were classified in one of the eight community states. The succession of the main commun Multiple Factor Analysis (MFA) enables an integrative analysis of all parameters including factors and was perfo community states, months, zooplankton and environmental time series. Cube root transformation was applied o highly abundant groups, whilst other variables were standardized to zero mean and scaled to unit variance. The structure was explored by depicting the succession of the community states over the 13-year time series, along abundance as well as the weekly anomalies of water temperature. Finally, in order to further explore the underly particular emphasis was placed on the spring peak event, since it was found to be a major aspect characterizing (Dauby, 1980). Therefore we considered the abundance of the group responsible for the spring peak, i.e., calan March to the third week of May). The 0.80 quantile values were computed to minimize the influence of extreme spring peak. To explore potential link between each of the environmental variables with the spring peak magnitu from two to seven months were tested using Spearman's rank correlation coefficient. Relationships with the mos linear or cubic spline regressions. Similarly, the relationship between spring peak magnitude and the winter Nor //climatedataguide.ucar.edu/climate-data/hurrell-north-atlantic- oscillation-nao-index-station-based) was also inv analysis is performed using R software (RStudio, version 1.4.1103) and Excel 2013. The data exploration is org all variables (five biomarkers (SOD, GST, MDA, AChE, LAC), two physicochemical water parameters (Salinity, T using Fligner-Killeen tests. Normality is also tested using a Shapiro-Wilk test. Several tests were then carried ou

equation for each biomarker, taking into account each of the confounding factors (Press et al., 2013; Saporta 20
regression formulae were obtained for each biomarker and using these equations, predicted values and residua
measuring a deviation from the model, were classified into three groups for each biomarker using the k-means c
The resulting clusters were then used to determine, for each biomarker, the expected ranges of variation accorc

Data Set Usage Rights

Additional Metadata


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additionalMetadata
        |___text '\n    '
        |___element 'metadata'
        |     |___text '\n      '
        |     |___element 'unitList'
        |     |     |___text '\n        '
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        |     |     |        \___attribute 'name' = 'formazineNephelometricUnit'
        |     |     |        \___attribute 'unitType' = 'dimensionless'
        |     |     |        \___attribute 'id' = 'formazineNephelometricUnit'
        |     |     |        \___attribute 'parentSI' = ''
        |     |     |        \___attribute 'multiplierToSI' = '1.0'
        |     |     |        \___attribute 'constantToSI' = '0.0'
        |     |     |___text '\n        '
        |     |     |___element 'unit'
        |     |     |        \___attribute 'name' = 'micromolePerLiter'
        |     |     |        \___attribute 'unitType' = 'spatialDensity'
        |     |     |        \___attribute 'id' = 'micromolePerLiter'
        |     |     |        \___attribute 'parentSI' = 'molePerLiter'
        |     |     |        \___attribute 'multiplierToSI' = '0.000001'
        |     |     |        \___attribute 'constantToSI' = '0.0'
        |     |     |___text '\n        '
        |     |     |___element 'unit'
        |     |     |        \___attribute 'name' = 'nanogramsPerLiter'
        |     |     |        \___attribute 'unitType' = 'spatialDensity'
        |     |     |        \___attribute 'id' = 'nanogramsPerLiter'
        |     |     |        \___attribute 'parentSI' = 'gramsPerLiter'
        |     |     |        \___attribute 'multiplierToSI' = '0.000000001'
        |     |     |        \___attribute 'constantToSI' = '0.0'
        |     |     |___text '\n      '
```

```
|     |___text '\n    '
|___text '\n  '
```