



QUASIMEME

Quality assurance of information
for marine environmental monitoring

Certificate of Analysis



Sediment

REFERENCE MATERIAL

Sediment sample 60



Certificate of Analysis Sediment 60

General Information

In this report an overview is given of analytical data for this sample collected in our proficiency testing program. The consensus values are calculated using a robust statistical model. With this NDA model, the mean and standard deviation are calculated using all reported data when at least 4 results are left after removal of reported 'lower than' (<) and 0 (= zero) values. No outliers are removed.

This report is divided into two sections: Consensus Values and Indicative Values. The division is made on the reliability of the data. Consensus Values are based on at least 8 results and a maximum relative uncertainty of 6.25%. Indicative Values are based on a maximum relative uncertainty of 35% and a minimum of 4 and maximum of 7 results, or a relative uncertainty greater than 6.25% when there are at least 8 results.

For each determinand the following parameters are given: mean, standard deviation, coefficient of variation, number of results, median, MAD (Median of Absolute Deviation), the uncertainty of the mean (consensus or indicative) value and the relative uncertainty.

The results of each determinand is expressed on dried sediment.

Sample information

QUASIMEME reference materials cover a range of natural Marine sediment species from contaminated waters from the North Sea and/or Mediterranean. There is no spiking, mixing or other alterations of the samples. For sample preparation the sediment samples are dried at 40 °C and milled to pass a 0.5 mm sieve.

This Sediment sample 60 of Estuary sediment spiked from Westerscheldt is prepared for the QUASIMEME proficiency programs. The results on which the values in this report are based were taken from the periods given in the following table.

Year.Round	Program	Sample Round Id
2024.1	MS6	QSP089MS
2022.2	MS6	QSP083MS
2021.2	MS6	QSP080MS
2022.1	MS7	QBC072MS
2021.2	MS8	QPF015MS
2020.2	MS8	QPF011MS



Consensus Values MS6

Method: Organometals - MS6

Element	Unit	Mean	Std.Dev.	CV %	N	Median	MAD	Uncertainty	Rel.Uncert. %
Tributyltin (TBT)	µg Sn/kg	8.99	2.22	24.7	58	9.10	1.22	0.364	4.05
Dibutyltin (DBT)	µg Sn/kg	9.16	2.45	26.8	51	9.12	1.65	0.429	4.69



Indicative Values MS6

Method: Organometals - MS6

Element	Unit	Mean	Std.Dev.	CV %	N	Median	MAD	Uncertainty	Rel.Uncert. %
Monobutyltin (MBT)	µg Sn/kg	36.7	22.2	60.6	44	37.6	14.4	4.19	11.4
Triphenyltin (TPhT)	µg Sn/kg	1.87	0.489	26.1	25	2.00	0.300	0.122	6.53
Diphenyltin (DPhT)	µg Sn/kg	2.27	0.855	37.7	11	2.32	0.569	0.322	14.2
Monophenyltin (MPhT)	µg Sn/kg	3.44	1.53	44.5	13	3.64	0.637	0.530	15.4



Consensus Values MS7

Method: Brominated Flame Retardants - MS7

Element	Unit	Mean	Std.Dev.	CV %	N	Median	MAD	Uncertainty	Rel.Uncert. %
BDE209	µg/kg	15.3	1.82	11.9	9	15.7	0.660	0.757	4.96



Indicative Values MS7

Method: Brominated Flame Retardants - MS7

Element	Unit	Mean	Std.Dev.	CV %	N	Median	MAD	Uncertainty	Rel.Uncert. %
BDE028	µg/kg	0.024	0.017	70.2	9	0.031	0.013	0.007	29.3
BDE047	µg/kg	1.30	0.227	17.4	12	1.26	0.169	0.082	6.27
BDE099	µg/kg	0.077	0.027	35.0	10	0.082	0.015	0.011	13.8



Indicative Values MS8

Method: Perfluorinated alkyl substances - MS8

Element	Unit	Mean	Std.Dev.	CV %	N	Median	MAD	Uncertainty	Rel.Uncert. %
PFHxA	µg/kg	0.017	0.006	38.4	4	0.017	0.004	0.004	24.0
PFHpA	µg/kg	0.008	0.002	28.7	4	0.008	0.001	0.001	18.0
PFOA	µg/kg	0.011	0.005	43.8	4	0.012	0.003	0.003	27.3
PFDA	µg/kg	0.012	0.003	25.0	4	0.013	0.002	0.002	15.6
PFUnDA	µg/kg	0.032	0.013	40.2	5	0.039	0.007	0.007	22.4
PFDoA	µg/kg	0.057	0.012	22.1	5	0.061	0.007	0.007	12.3
PFTTrDA	µg/kg	0.185	0.057	31.0	8	0.191	0.022	0.025	13.7
PFTeDA	µg/kg	0.133	0.022	16.8	8	0.131	0.008	0.010	7.42
n-PFHxS	µg/kg	0.029	0.007	22.8	4	0.030	0.003	0.004	14.3
total-PFOS	µg/kg	0.085	0.029	34.8	8	0.091	0.021	0.013	15.4