SUPPLEMENTARY MATERIALS

Fluoxetine Removal from Aqueous Solutions Using a Lignocellulosic Substrate Colonized by the White-Rot Fungus *Pleurotus ostreatus*

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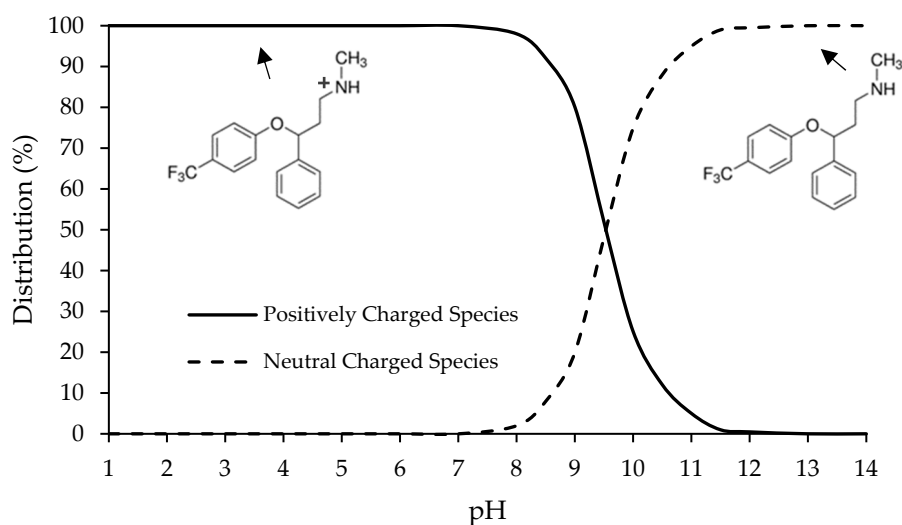
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Figure S1. Species distribution diagram of fluoxetine as a function of pH (adapted from [2] and [3]).
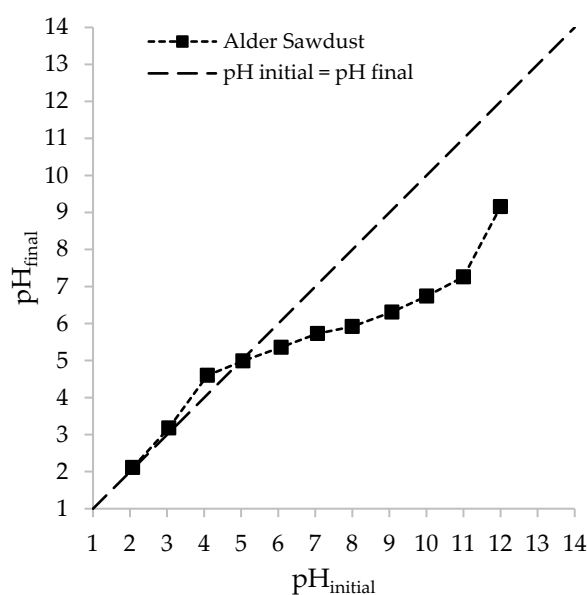
Figure S2. The point of zero charge (pH_{PZC}) determination in alder sawdust.

Table S1. Physicochemical properties of fluoxetine hydrochloride.

Molecular formula [1]	M.w. (g mol^{-1})	S.w. (mg L^{-1})	*p*Ka [1]	Log K$_{ow}$ [1]
$C_{17}H_{18}F_3NO \cdot HCl$	345.79	4000	9.8	4.17

M.w.: Molecular weight; S.w.: Solubility in water; Ka: Acid dissociation constant; K$_{ow}$: Octanol-water partition coefficient.

[1] Source [1].

Table S2. List of reagents used in the study.

Reagent	Supplier	Country
Potassium bromide (KBr, for IR spectroscopy)	PanReac	Spain
Potassium nitrate (KNO_3)	Sigma-Aldrich®	USA
Sodium hydroxide (NaOH, purity >99.0%)	Labkem	Spain
Hydrochloric acid (HCl, purity 37% wt)	Honeywell Fluka™	Austria
2.6-dimethoxyphenol (2.6-DMP, purity 99%)	Labkem	Spain
Sodium acetate anhydrous (CH_3COONa, purity ≥99%)	Sigma-Aldrich	USA
Acetic acid glacial (CH_3COOH, purity ≥99.8%)	Carlo Erba	Spain
1-aminobenzotriazole (1-ABT)	Sigma-Aldrich	USa
Ortho-boric acid (H_3BO_3, purity 100%)	VWR BDH® Chemicals	Belgium
Acetic acid glacial (CH_3COOH, purity 99.8%, density 1.0501-1.0521)	CARLO ERBA Reagents	France
Ortho-phosphoric acid (H_3PO_4, purity 85%, density 1.70 kg L^{-1})	PanReac	Spain
Acetonitrile (CH_3CN, purity 99.99%)	VWR BDH® Chemicals HPLC-super gradient	France
Formic acid (HCOOH, purity 99%)	CARLO ERBA Reagents	France

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Methanol (CH₃OH, purity >99.9%)	VWR BDH® Chemicals HPLC-super gradient	Netherlands

Table S3. Band assignments of FT-IR spectrum of alder sawdust.

Band	Wavenumber (cm⁻¹)	Band assignments [1]	References
A	3408 ± 0	ν(O-H)	[4-8]
B	2925 ± 0	$\nu_{as}(CH_2)$ and $\nu_s(CH_2)$	[4-7]
C	1650 ± 1	ν(C=O)	[4, 7-9]
D	1539 ± 3	δ(C=C)	[8]
E	1401 ± 25	$\delta(CH_2)$ and $\delta(CH_3)$	[4, 8]
F	1238 ± 1	ν(C-O)	[8]
G	1079 ± 52	ν(C-O)	[4, 8, 9]
H	668 ± 1	Bending vibration modes of aromatic compounds.	[8]
I	601 ± 1		

[1] ν: stretching; ν_s: symmetric stretching; ν_{as}: asymmetric stretching; δ: deformation.

Table S4. pH and temperature values recorded at the beginning and the end of batch assays.

Assay	pH		Temperature (°C)	
	Beginning	End	Beginning	End
CMS Assays (1st day)	6.83 ± 0.19	6.69 ± 0.01	22.0 ± 0.0	24.0 ± 0.0
CMS Assays (15th day)	7.01 ± 0.01	6.66 ± 0.03	20.0 ± 0.0	24.0 ± 0.0
Biosorption Assays	6.94 ± 0.01	6.46 ± 0.00	24.0 ± 0.0	25.0 ± 0.0
CYP450 Inhibition Assays	6.75 ± 0.02	6.74 ± 0.05	23.0 ± 0.0	24.0 ± 0.0
CEE Assays	6.68 ± 0.01	6.68 ± 0.01	24.5 ± 0.7	24.5 ± 0.7

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