Supplementary information

Preparation of highly porous carbon from sustainable α -cellulose for superior performance removal of tetracycline and sulfamethazine from water

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| Compounds | Chemical formula | Molecular weight | S _W ^a (mmol L ⁻¹) | log <i>K_{OW}b</i> (L L ⁻¹) | p <i>K</i> a |
|-----------|-----------------------|------------------|--|--|--------------|
| ТС | $C_{22}H_{24}N_2O_8$ | 444.43 | 3.81 | -1.19 ¹ | 3.3(1) |
| | | | | | 7.7(2) |
| | | | | | 9.7(3)1 |
| SMZ | $C_{12}H_{14}N_4O_2S$ | 278.33 | 1.46 ² | 0.89 ² | 2.3(1) |
| | | | | | $7.5(2)^2$ |

Table S1 Summary of Sorbate Properties.

^a S_W - Water Solubility (25 °C);

^b *Kow* - octanol–water partition coefficient.

¹ J Tolls. Environ. Sci. Technol.35 (2001) 3397-3406.

² L Warisara, O Say Kee, T Moorman. Chemosphere 76 (2009) 558-564.

| Samples | Elemental composition ^a (%) | | | Yield | S _{BET} ^b | V_P^{c} | $S_{micro}{}^{\mathrm{d}}$ | V _{micro} e | d _{aver} ^f |
|----------|--|-------|------|------------|-----------------------------------|------------------------------------|----------------------------|------------------------------------|--------------------------------|
| | N | С | Н | - (wt%) | (m ² g ⁻¹) | (cm ³ g ⁻¹) | $(m^2 g^{-1})$ | (cm ³ g ⁻¹) | (nm) |
| Са | 0.87 | 60.03 | 1.21 | 47.22 | 26.14 | 0.042 | 1.92 | 0.002 | 2.769 |
| Са-850-0 | 0.51 | 89.12 | 0.75 | 59.61 | 56.34 | 0.061 | 12.01 | 0.013 | 1.688 |
| Cα-850-4 | 0.41 | 89.77 | 0.76 | 11.41 | 3187.91 | 1.781 | 2655.35 | 1.407 | 1.628 |

Table S2 Characteristics of α -cellulose, C α , C α - β - γ .

^a: Determined by Vario Micro cube element analyzer.

^b: BET specific surface area calculated in the relative pressure region $P/P_o = 0.05 - 0.35$.

^c: Total pore volume which is determined at $P/P_o = 1.0000$.

^d: Micropore surface area calculated by the t-plot method.

^e: Micropore volume calculated by the t-plot method.

^f: Average pore diameter obtained from DFT equation using N₂ isotherms.



Fig. S1 Chemical structures of tetracycline and sulfamethazine. The regions framed by dashed lines represent the structural moieties associated with the acidic dissociation constants (pKa).



Fig. S2 High-resolution XPS spectra of (a) O1s and (b) C1s peaks of $C\alpha$ -850-4.