

## Electronic Supplementary Information

# Coordination framework materials fabricated by the self-assembly of Sn(IV) porphyrins with Ag(I) ions for the photocatalytic degradation of organic dyes in wastewater

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**Fig. S20** ESI-MS spectrum (negative ion mode) of the reaction mixture of AM in the presence of **2** after 60 min of visible light irradiation.

**Table S1.** Crystallographic data and structure refinements for **1** and **2**.

	<b>1</b>	<b>2</b>
Empirical formula	C <sub>40</sub> H <sub>24</sub> Ag <sub>2</sub> N <sub>8</sub> O <sub>2</sub> Sn	C <sub>58</sub> H <sub>38</sub> Ag <sub>2</sub> F <sub>6</sub> N <sub>12</sub> O <sub>10</sub> S <sub>2</sub> Sn
Formula weight	983.1	1575.55
Crystal size (mm <sup>3</sup> )	0.20 × 0.15 × 0.15	0.2 × 0.3 × 0.5
<i>T</i> (K)	173 (2)	223(2)
Crystal system, Space group	Orthorhombic, <i>Pmna</i>	Monoclinic, <i>C2/c</i>
<i>Cell dimensions</i>		
<i>a</i> (Å)	19.8288 (15)	17.616(3)
<i>b</i> (Å)	17.6770 (14)	24.301 (4)
<i>c</i> (Å)	8.5182 (7)	19.874 (4)
$\alpha$ (deg)	90.00	90.00
$\beta$ (deg)	90.00	108.578 (3)
$\gamma$ (deg)	90.00	90.00
<i>V</i> (Å <sup>3</sup> )	2985.7 (4)	8064 (3)
<i>Z</i> , <i>D<sub>c</sub></i> (g cm <sup>-3</sup> )	2, 1.088	4, 1.298
$\mu$ (mm <sup>-1</sup> )	1.11	0.91
<i>F</i> (000)	1096	3120
$\theta$ range (°)	1.15 to 28.29	2.05 to 28.29
reflections collected	17461	23423
independent reflections ( <i>R</i> <sub>int</sub> )	3710 (0.0635)	9377 (0.041)
absorption correction	None	None
data / restraints / parameters	3710 / 0 / 123	9377 / 3 / 410
GOF on <i>F</i> <sup>2</sup>	0.9239	1.2774
<i>R</i> <sub>1</sub> , <sup>a</sup> <i>wR</i> <sub>2</sub> <sup>b</sup> [ <i>I</i> > 2 $\sigma$ ( <i>I</i> )]	0.066, 0.1892	0.091, 0.2826
<i>R</i> <sub>1</sub> , <sup>a</sup> <i>wR</i> <sub>2</sub> <sup>b</sup> (all data)	0.1233, 0.2111	0.1261, 0.3422
Largest peak/hole (e Å <sup>-3</sup> )	1.4586/ -1.6532	8.1412/-7.7329

<sup>a</sup>  $R_1 = \frac{\sum \|F_o\| - |F_c|}{\sum |F_o|}$ ; <sup>b</sup>  $wR_2 = \frac{[\sum [w(F_o^2 - F_c^2)^2]]}{[\sum [w(F_o^2)^2]]}^{1/2}$ .

**Table S2.** Selected bond lengths [Å] and angles [°] for **1**.

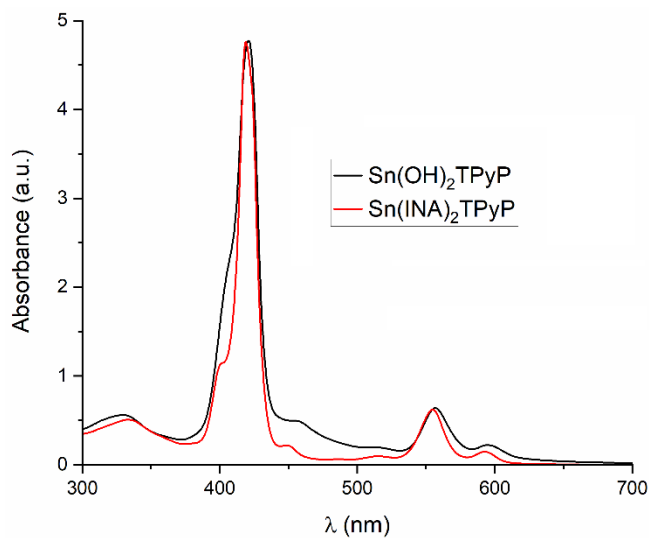
Sn1-O1	2.069(7)	Sn1-N1	2.074(4)
Ag2-N2	2.052(14)	Ag3-N3	2.166(6)
O1-Sn1-O1	180.0(4)	O1-Sn1-N1	88.7(2)
O1-Sn1-N1	91.3(2)	O1-Sn1-N1	88.7(2)
O1-Sn1-N1	91.3(2)	N1-Sn1-N1	90.3(2)
O1-Sn1-N1	91.3(2)	O1-Sn1-N1	88.7(2)
N1-Sn1-N1	180.0(8)	N1-Sn1-N1	89.7(2)
O1-Sn1-N1	91.3(2)	O1-Sn1-N1	88.7(2)
N1-Sn1-N1	89.7(2)	N1-Sn1-N1	180.0(1)
N1-Sn1-N1	90.3(2)	N2-Ag2-N2	180.0(10)
N3-Ag3-N3	180.0(5)	C2-N1-Sn1	125.8(3)
C5-N1-Sn1	125.5(4)	C9-N3-Ag3	119.5(4)
C13-N2-Ag2	122(2)	C12-N2-Ag2	127.0(16)

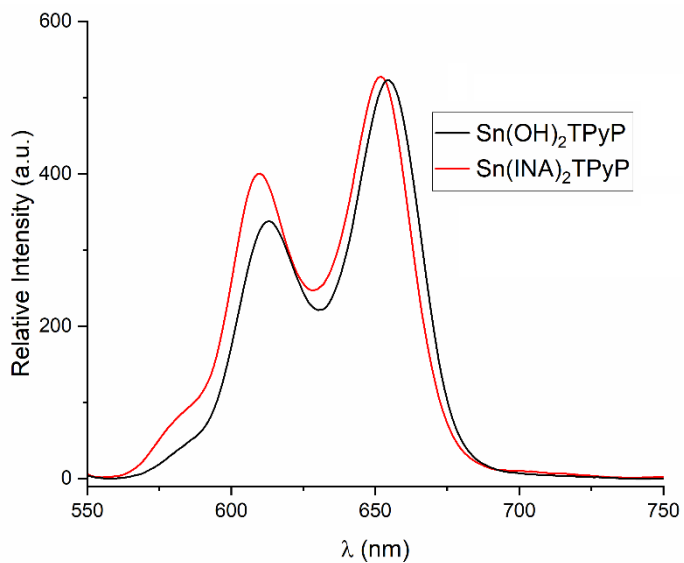
Symmetry codes: (i)  $-x+1, -y+1, -z+1$ ; (ii)  $-x+1, y, z$ ; (iii)  $x, -y+1, -z+1$ ; (iv)  $-x+1, -y+2, -z$ ; (v)  $-x, -y+1, -z+1$ .

**Table S3.** Selected bond lengths [Å] and angles [°] for **2**.

Sn1-O1	2.066(5)	Sn1-N2	2.081(5)
Sn1-N1	2.087(5)	Ag1-N6	2.198(6)
Ag1-N4	2.373(9)	Ag2-N3	2.216(6)
Ag2-N5	2.443(9)		
O1-Sn1-O1	178.9(3)	O1-Sn1-N2	94.7(2)
O1-Sn1-N2	86.1(2)	O1-Sn1-N1	94.7(2)
N2-Sn1-N2	89.6(3)	O1-Sn1-N1	85.0(2)
O1-Sn1-N1	94.2(2)	N2-Sn1-N1	179.7(2)
N2-Sn1-N1	90.3(2)	O1-Sn1-N1	94.2(2)
O1-Sn1-N1	85.0(2)	N2-Sn1-N1	90.3(2)
N2-Sn1-N1	179.7(2)	N1-Sn1-N1	89.8(3)
N6-Ag1-N6	149.7(4)	N6-Ag1-N4	105.16(18)
N3-Ag2-N3	165.0(3)	N3-Ag2-N5	97.48(16)
C23-O1-Sn1	129.2(5)	C1-N1-Sn1	126.2(4)
C4-N1-Sn1	125.2(4)	C14-N2-Sn1	125.8(4)
C11-N2-Sn1	125.2(4)	C9-N3-Ag2	124.5(6)
C8-N3-Ag2	118.1(5)	C18-N4-Ag1	121.9(5)
C22-N5-Ag2	121.3(5)	C27-N6-Ag1	118.2(6)
C26-N6-Ag1	124.2(5)		

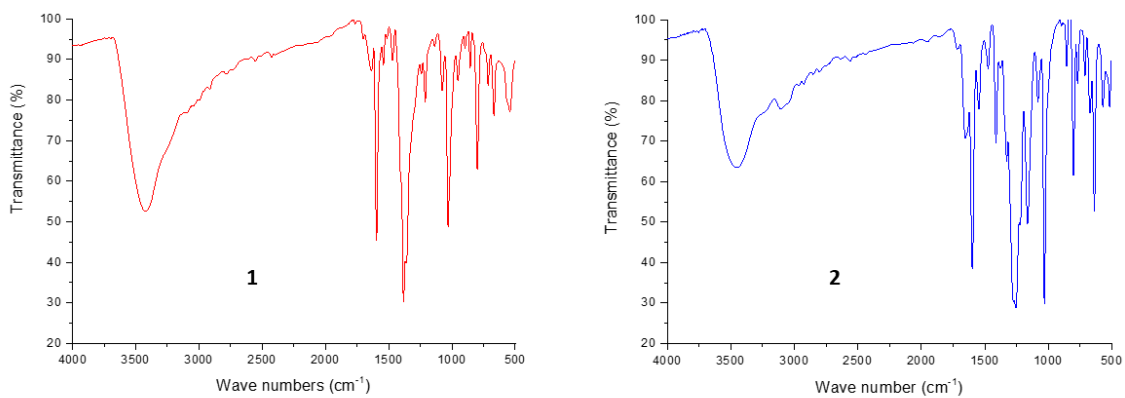
Symmetry codes: (i)  $-x+2, y, -z+3/2$ ; (ii)  $-x+2, y, -z+1/2$ ; (iii)  $-x+2, -y, -z+1$ ; (iv)  $x+1/2, y+1/2, z$ ; (v)  $x-1/2, y-1/2, z$ .

**Fig. S1** UV-vis absorption spectra of  $\text{Sn}(\text{OH})_2\text{TPyP}$  and  $\text{Sn}(\text{INA})_2\text{TPyP}$  in chloroform.



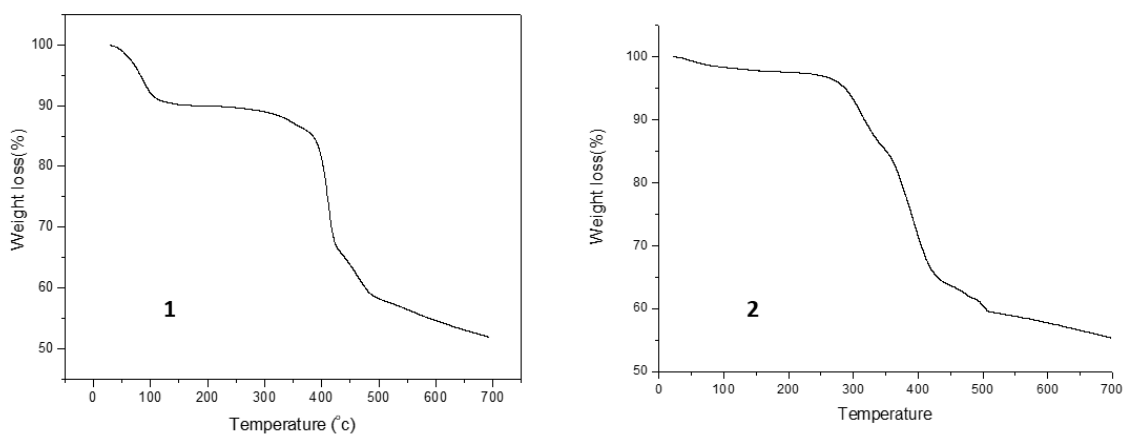
**Fig. S2** Fluorescence spectra of  $\text{Sn(OH)}_2\text{TPyP}$  and  $\text{Sn(INA)}_2\text{TPyP}$  (excited at 560 nm) in chloroform.

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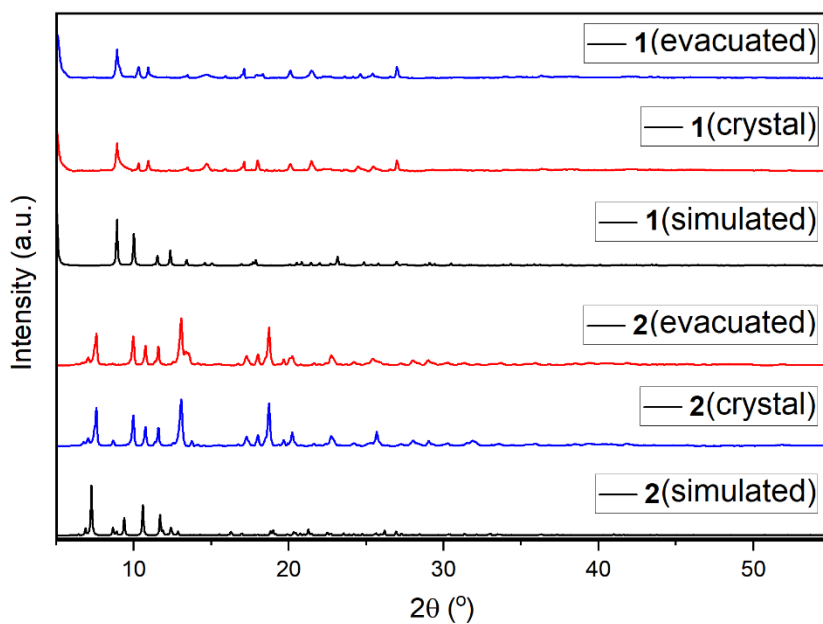
**Fig. S3** FT-IR spectra of **1** and **2**.

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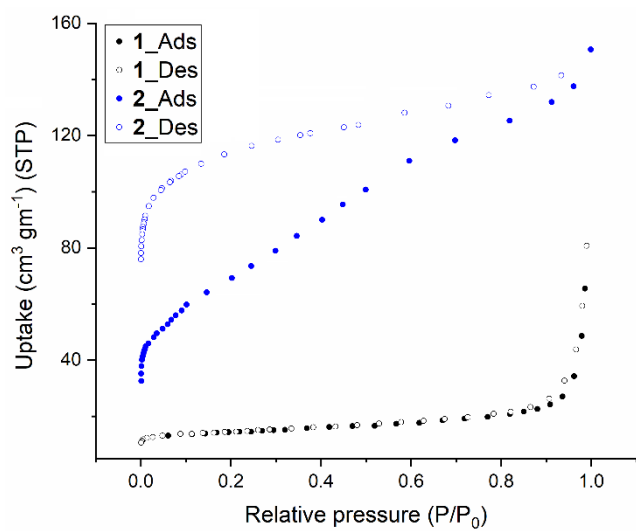
**Fig. S4** TGA curves of **1** and **2**.

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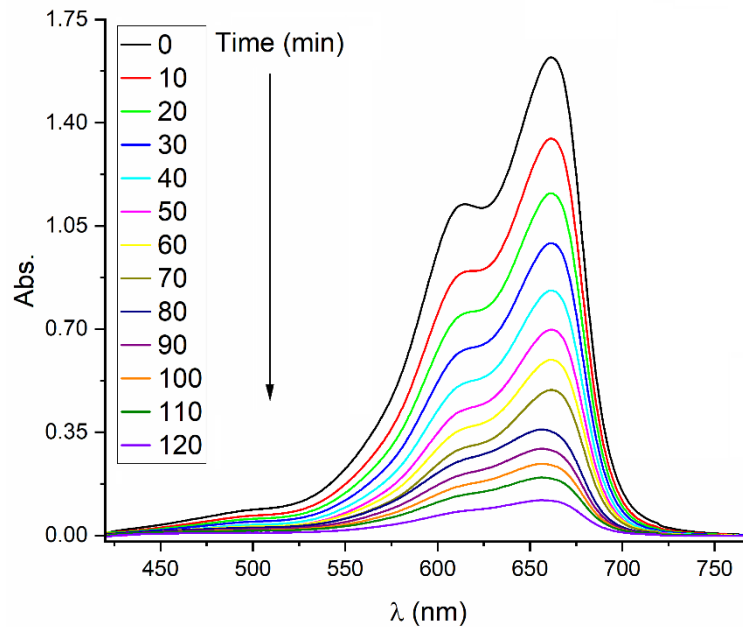
**Fig. S5** Powder X-ray diffraction (PXRD) patterns of **1** and **2**.

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**Fig. S6** Adsorption and desorption isotherms of N<sub>2</sub> for **1** and **2** at 77 K.

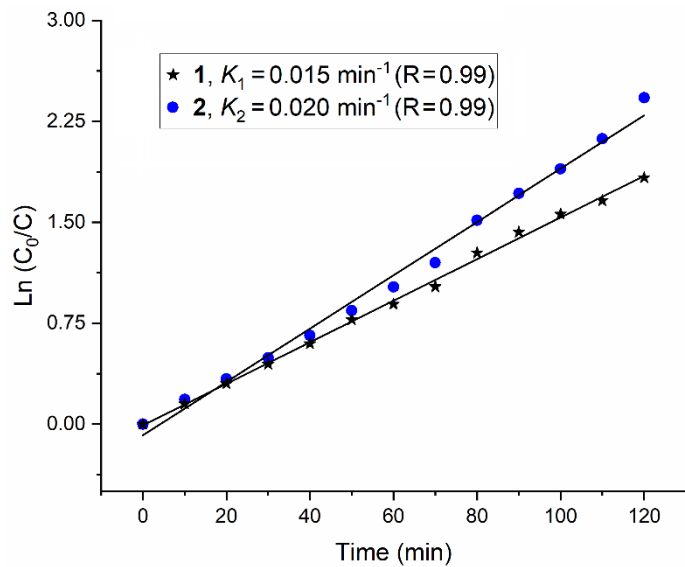
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**Fig. S7** Time-dependent absorption spectra of MB in the presence of **2** under visible light irradiation.

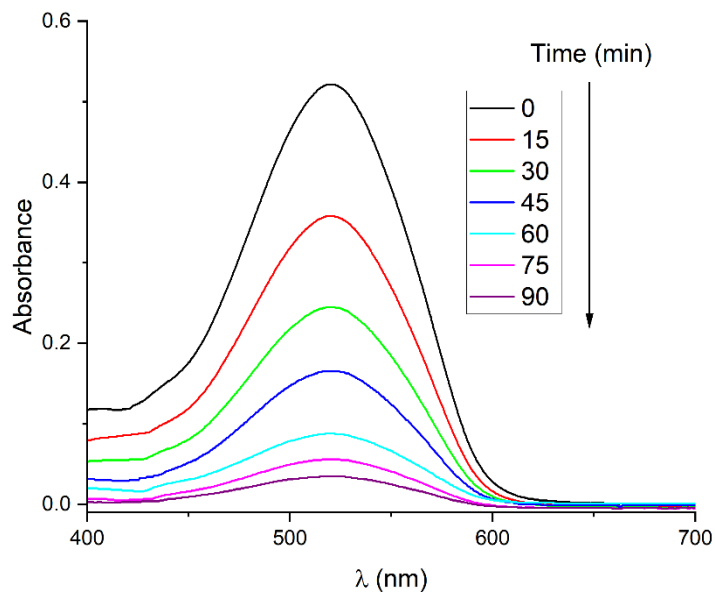
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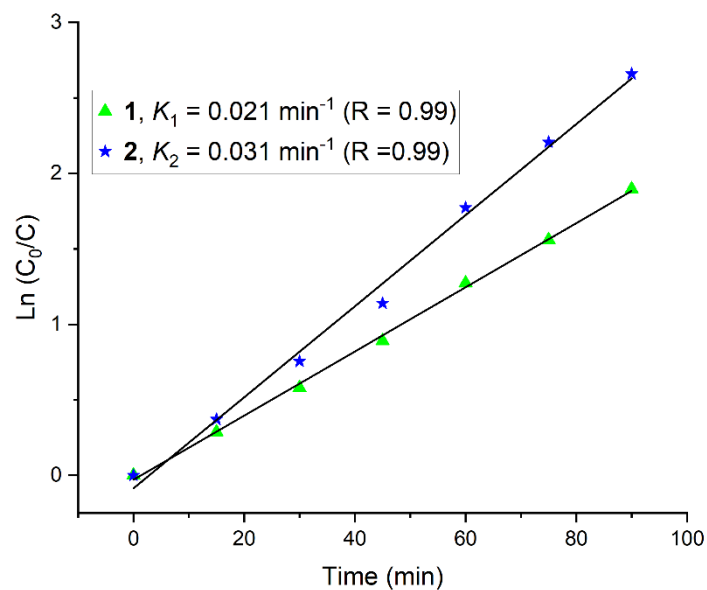
**Fig. S8** Kinetics for the photocatalytic degradation of MB under visible light irradiation by photocatalysts **1** and **2**.

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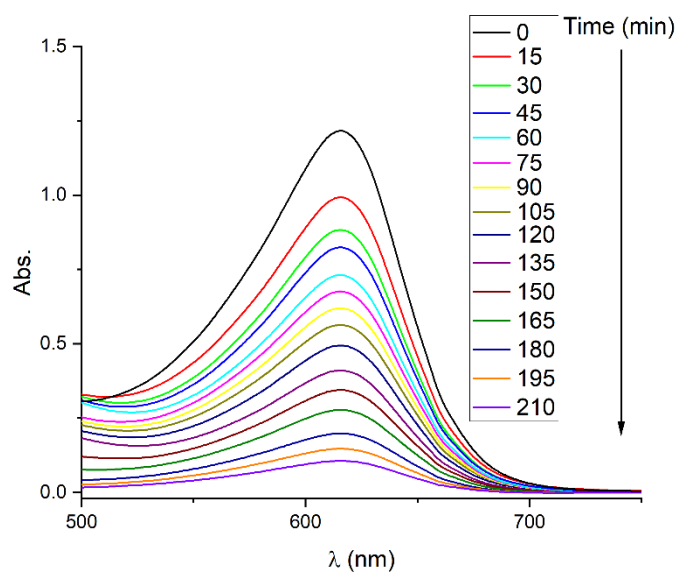
**Fig. S9** Time-dependent absorption spectra of AM in the presence of **2** under visible light irradiation.

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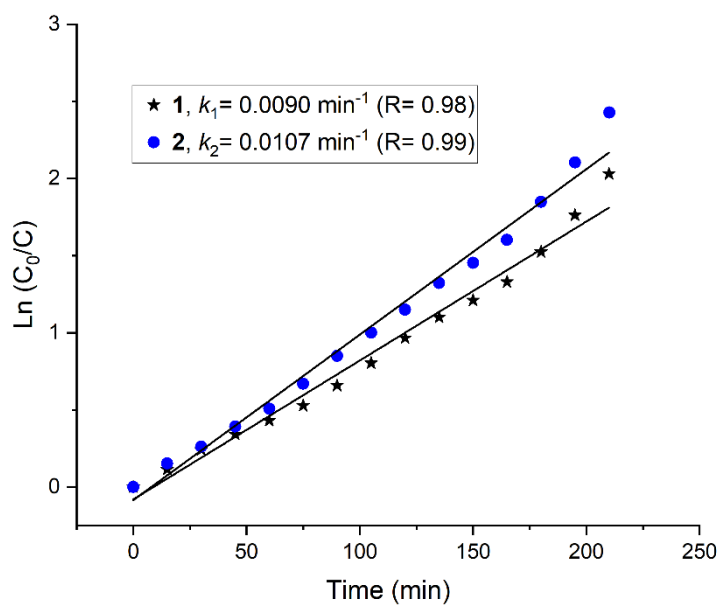
**Fig. S10** Kinetics-for the photocatalytic degradation of AM under visible light irradiation by photocatalysts **1** and **2**.

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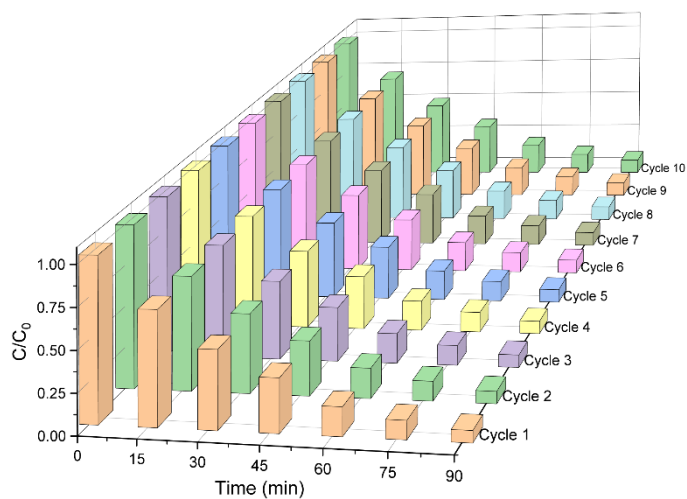
**Fig. S11** Time-dependent absorption spectra of BCG in the presence of **2** under visible light irradiation.

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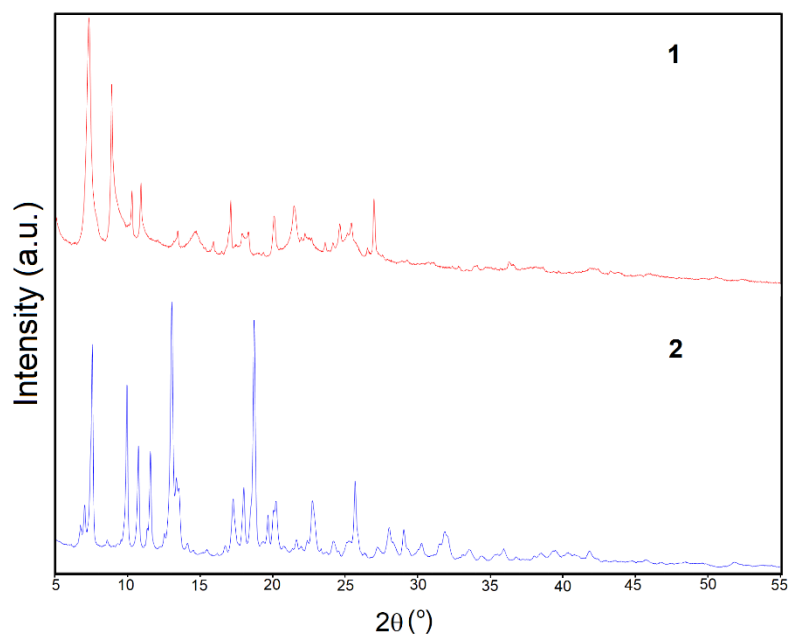
**Fig. S12** Kinetics for the photocatalytic degradation of BCG under visible light irradiation by photocatalysts **1** and **2**.

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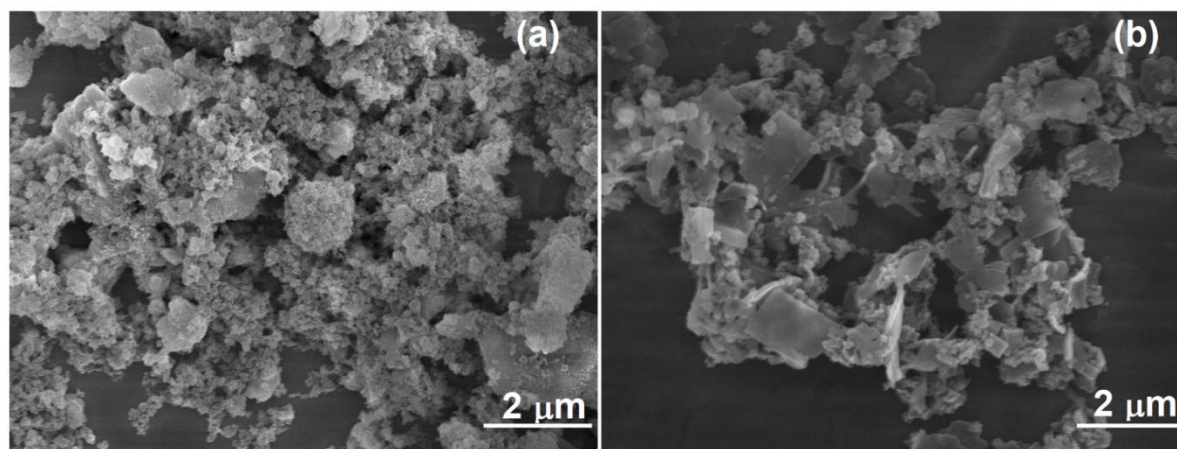
**Fig. S13** Catalytic cycles (up to 10 cycles) for photo-catalyst **2** for the degradation of AM dye.

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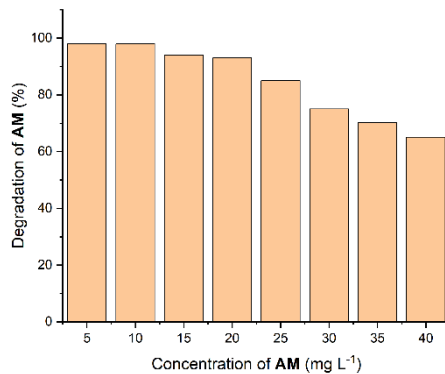
**Fig. S14** Powder X-ray diffraction (PXRD) patterns of **1** and **2** after use for photocatalytic degradation of AM dye.

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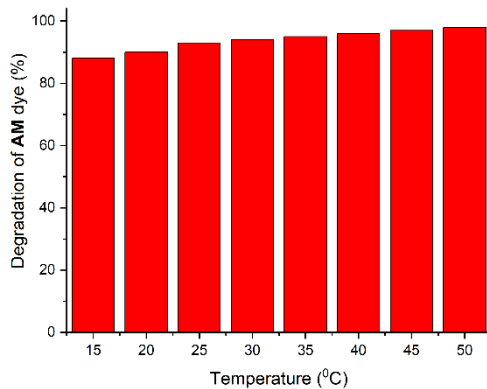
**Fig. S15** FE-SEM images of photocatalyst **1** and **2** after the degradation of AM dye.

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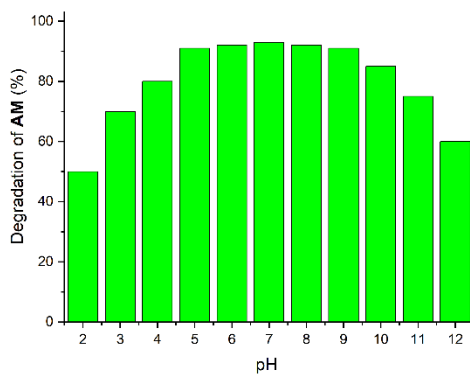
**Fig. S16** Concentration effect of AM dye on the photo-degradation by **2** (5 mg) within 90 min of visible light irradiation.

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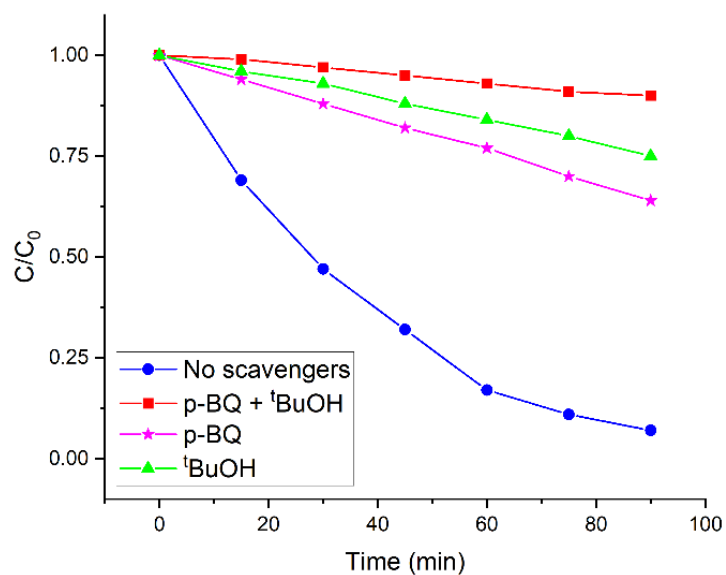
**Fig. S17** Effect of temperature on the degradation of AM dye by **2**.

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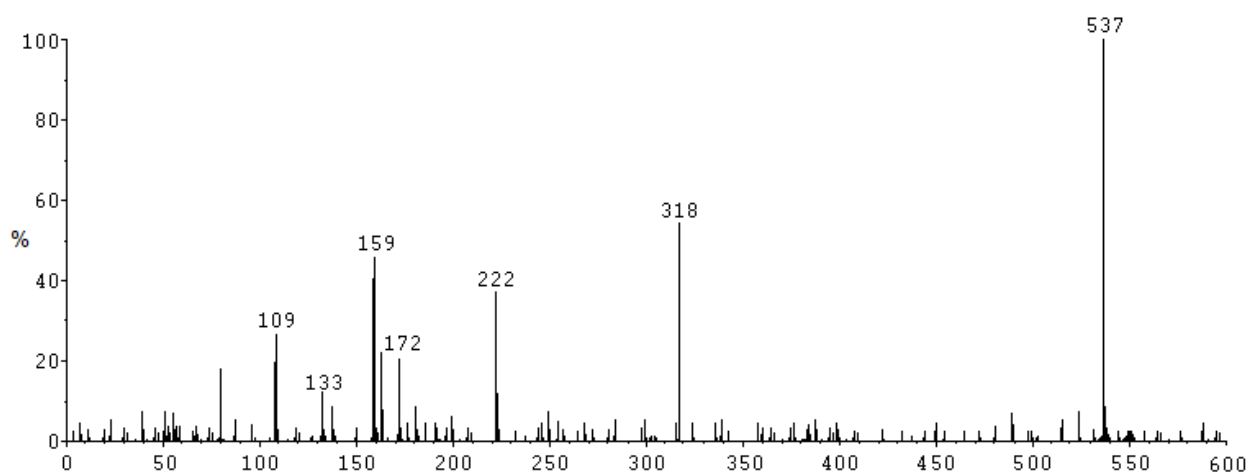
**Fig. S18** pH Effect of AM dye solution on the photo-degradation of by **2**.

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**Fig. S19** Effect of various scavengers on the degradation of AM dye in the presence of **2** under visible light irradiation.

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**Fig. S20** ESI-MS spectrum (negative ion mode) of the reaction mixture of AM in the presence of **2** after 60 min of visible light irradiation.

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