

Ratiometric Luminescent Sensing of Biomarker for Sugar Consumption in Aqueous Medium by Cu(II)Coordination Polymer

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Table S1 Crystallographic and structural refinement parameters for [Cu(Hadp)₂(Bimb)]_n (KA@CP-S3).

Empirical formula	C ₂₆ H ₃₂ CuN ₄ O ₈
Formula weight	592.11
Temperature/K	293(2)
Crystal system	Triclinic
Space group	P $\bar{1}$
a/Å	8.4173(5)
b/Å	8.4366(5)
c/Å	10.5472(7)
α /°	92.125(5)
β /°	93.821(6)
γ /°	118.349(6)
Volume/Å ³	655.69(8)
Z	1
ρ_{calc} /g/cm ³	1.4994
μ /mm ⁻¹	0.890
F(000)	309.5
Crystal size/mm ³	0.47 × 0.27 × 0.14
Radiation	Mo K α (λ = 0.71073)
2 θ range for data collection/°	3.88 to 51
Index ranges	-10 ≤ h ≤ 10, -10 ≤ k ≤ 10, -13 ≤ l ≤ 13
Reflections collected	15059

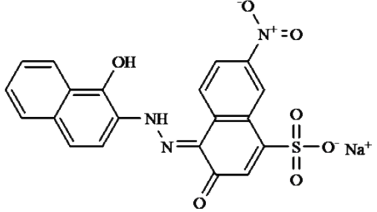
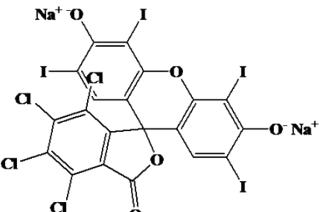
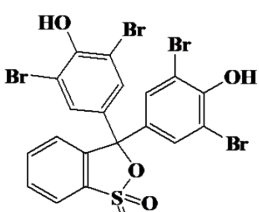
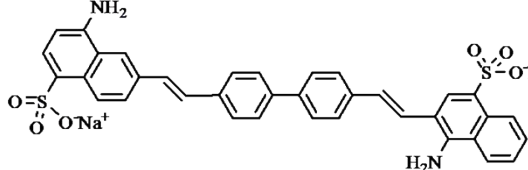
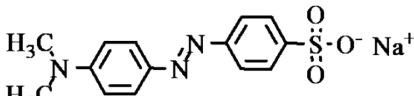
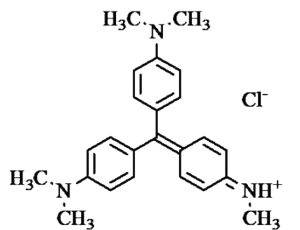
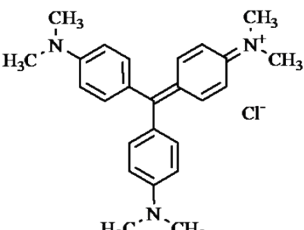
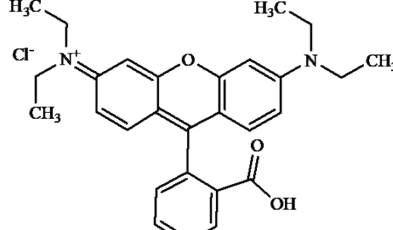
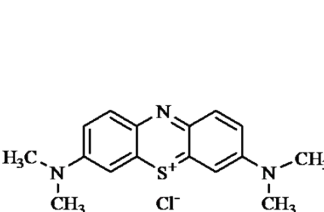
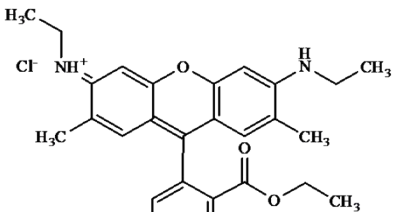
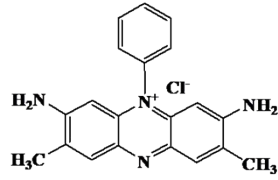
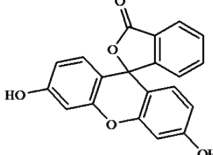
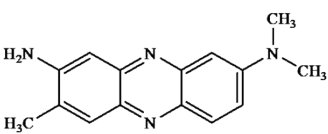
Independent reflections	2447 [$R_{\text{int}} = 0.0922$, $R_{\text{sigma}} = 0.0611$]
Data/restraints/parameters	2447/0/179
Goodness-of-fit on F^2	1.042
Final R indexes [$I \geq 2\sigma(I)$]	$R_1 = 0.1362$, $wR_2 = 0.4106$
Final R indexes [all data]	$R_1 = 0.1501$, $wR_2 = 0.4186$
Largest diff. peak/hole / $e \text{ \AA}^{-3}$	2.95/-1.13
CCDC No.	2233819

Table S2 Selected bond lengths [\AA] and bond angles [$^\circ$] for $[\text{Cu}(\text{Hadp})_2(\text{Bimb})]_n$ (**KA@CP-S3**).

<i>$[\text{Cu}(\text{Hadp})_2(\text{Bimb})]_n$ (KA@CP-S3)</i>					
<i>Bond lengths [\AA]</i>					
Cu1–O1	1.968(9)	Cu1–O3	2.735	Cu1–N1	1.981(11)
Cu1–O1 ¹	1.968(9)	Cu1–O3 ¹	2.735	Cu1–N1 ¹	1.981(11)
<i>Bond angles [$^\circ$]</i>					
O1 ¹ –Cu1–O1	180.0	N1 ¹ –Cu1–O3	89.8	N1–Cu1–O1 ¹	89.2(4)
N1 ¹ –Cu1–O1 ¹	90.8(4)	N1–Cu1–O3	90.2	N1–Cu1–O1	90.8(4)
N1 ¹ –Cu1–O1	89.2(4)	O3–Cu1–O3 ¹	180.0	N1 ¹ –Cu1–N1	180.0
O1 ¹ –Cu1–O3	94.9	O3 ¹ –Cu1–N1 ¹	90.2	O1 ¹ –Cu1–O3 ¹	85.1
O1–Cu1–O3	85.1	O3 ¹ –Cu1–N1	89.8	O1–Cu1–O3 ¹	94.9

Symmetry codes for **KA@CP-S3**: (i) 1-X,2-Y,1-Z; (ii) 1-X,1-Y,2-Z

Table S3 Chemical structures dyes included in present work.

<p>Anionic Dyes</p>	<div style="display: flex; justify-content: space-around; align-items: flex-start;"> <div style="text-align: center;">  <p>(EBT)</p> </div> <div style="text-align: center;">  <p>(RB)</p> </div> <div style="text-align: center;">  <p>(BPB)</p> </div> </div> <div style="display: flex; justify-content: space-around; align-items: flex-start; margin-top: 20px;"> <div style="text-align: center;">  <p>(CR)</p> </div> <div style="text-align: center;">  <p>(MO)</p> </div> </div>
<p>Cationic Dyes</p>	<div style="display: flex; justify-content: space-around; align-items: flex-start;"> <div style="text-align: center;">  <p>(MV)</p> </div> <div style="text-align: center;">  <p>(CV)</p> </div> <div style="text-align: center;">  <p>(RhB)</p> </div> </div> <div style="display: flex; justify-content: space-around; align-items: flex-start; margin-top: 20px;"> <div style="text-align: center;">  <p>(MB)</p> </div> <div style="text-align: center;">  <p>(Rh6G)</p> </div> <div style="text-align: center;">  <p>(SO)</p> </div> </div>
<p>Neutral Dyes</p>	<div style="display: flex; justify-content: space-around; align-items: flex-start;"> <div style="text-align: center;">  <p>(FS)</p> </div> <div style="text-align: center;">  <p>(NR)</p> </div> </div>

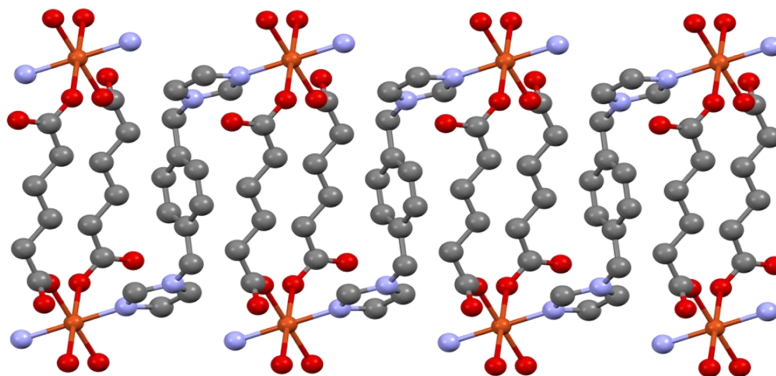


Fig. S1 View of 1D coordination polymer of **KA@CP-S3** formed through both Hadp^- and Bimb ligands along crystallographic “a” axis.

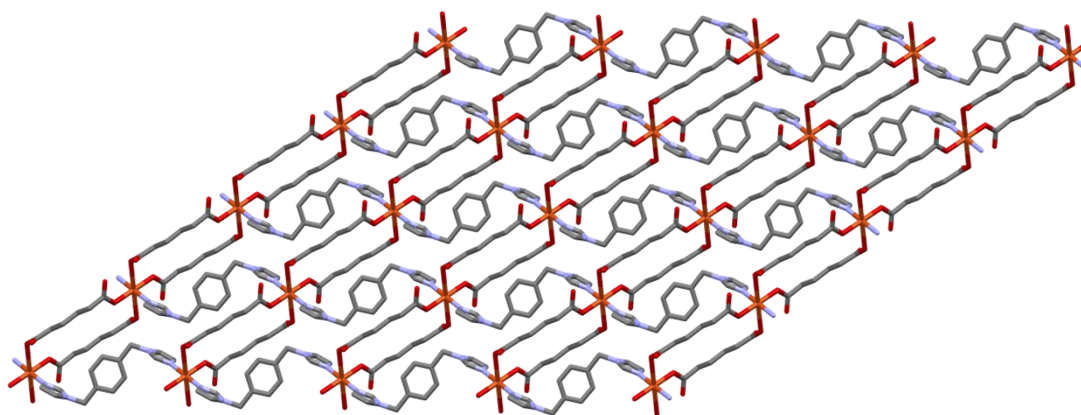


Fig. S2 View of 2D coordination network of **KA@CP-S3** along crystallographic “a” axis.

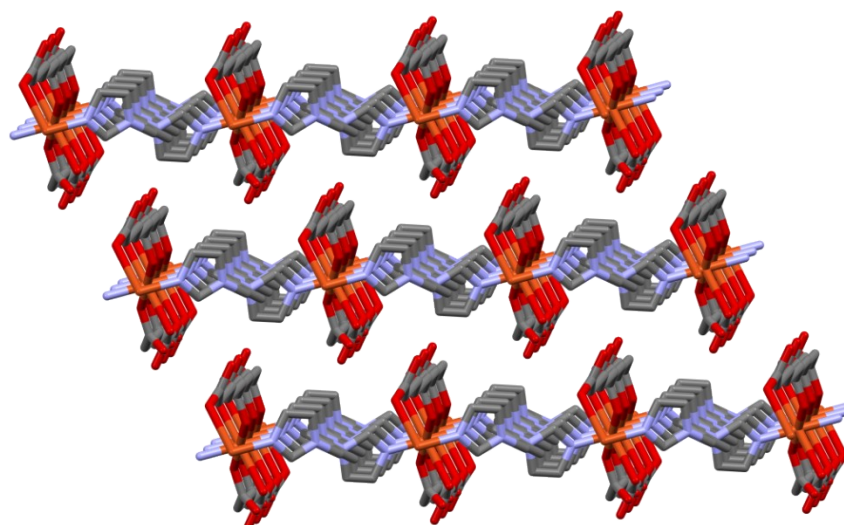


Fig. S3 View of 2D Parallel Layers of **KA@CP-S3** along crystallographic “c” axis.

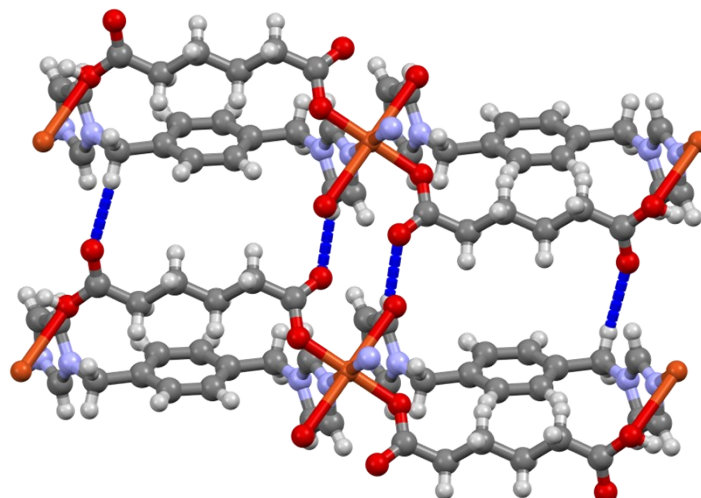


Fig. S4 View of C–H···O supramolecular hydrogen bonded 3D network of **KA@CP-S3** along crystallographic “b” axis.

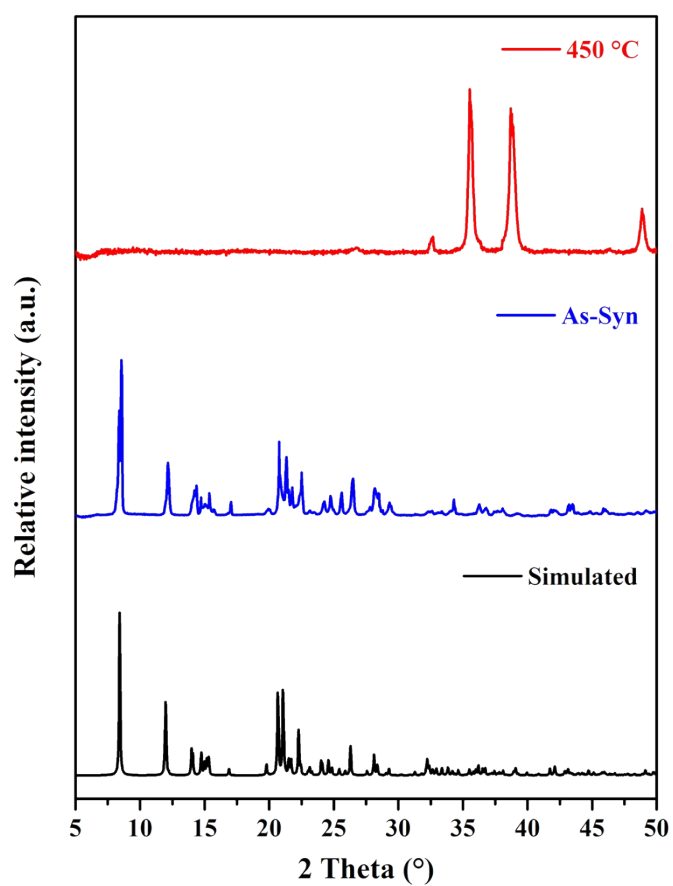


Fig. S5 Powder X-ray Diffraction spectra for **KA@CP-S3**.

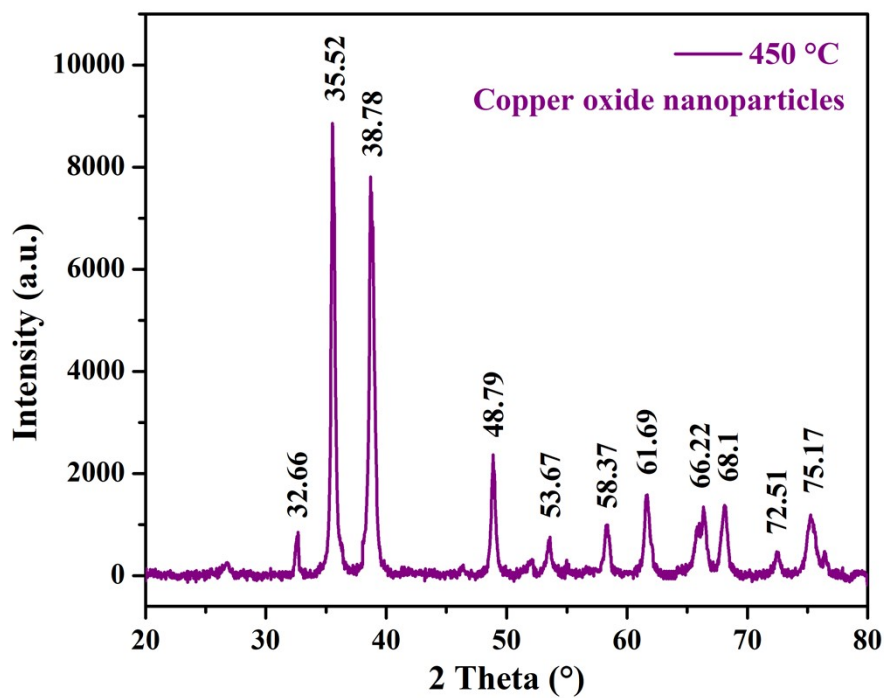


Fig. S6 PXRD patterns for CuO nanoparticles.

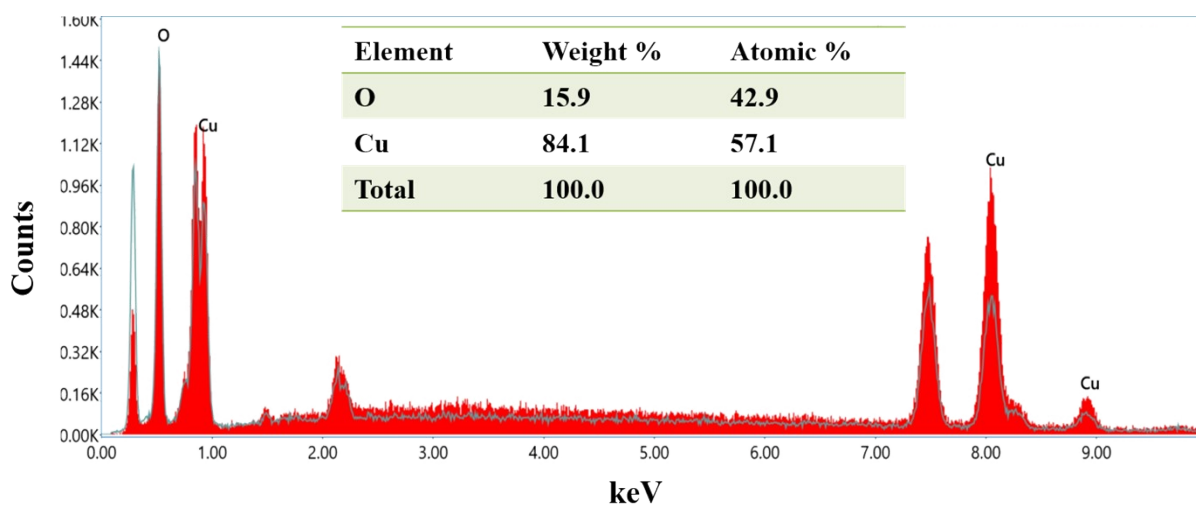


Fig. S7 EDAX diagram of CuO nanoparticles.

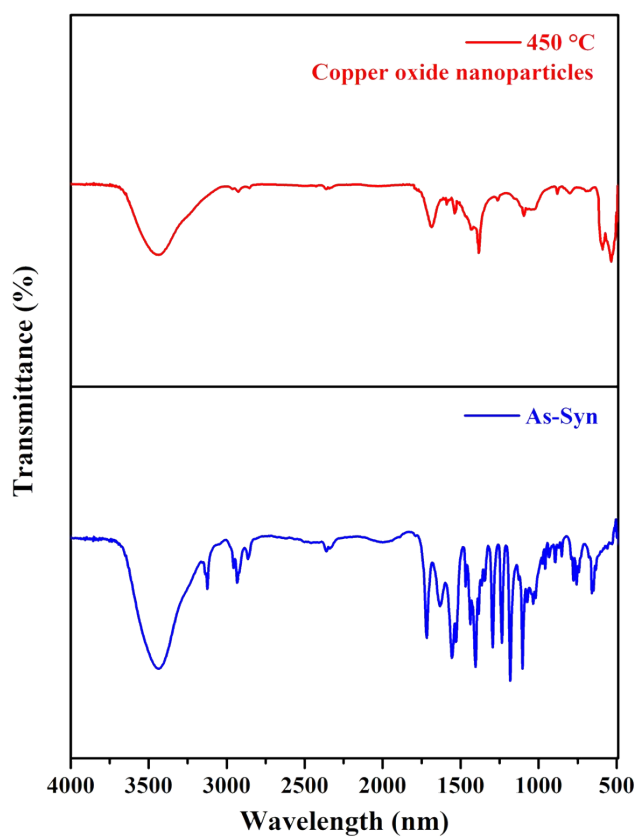


Fig. S8 FT-IR spectra for KA@CP-S3.

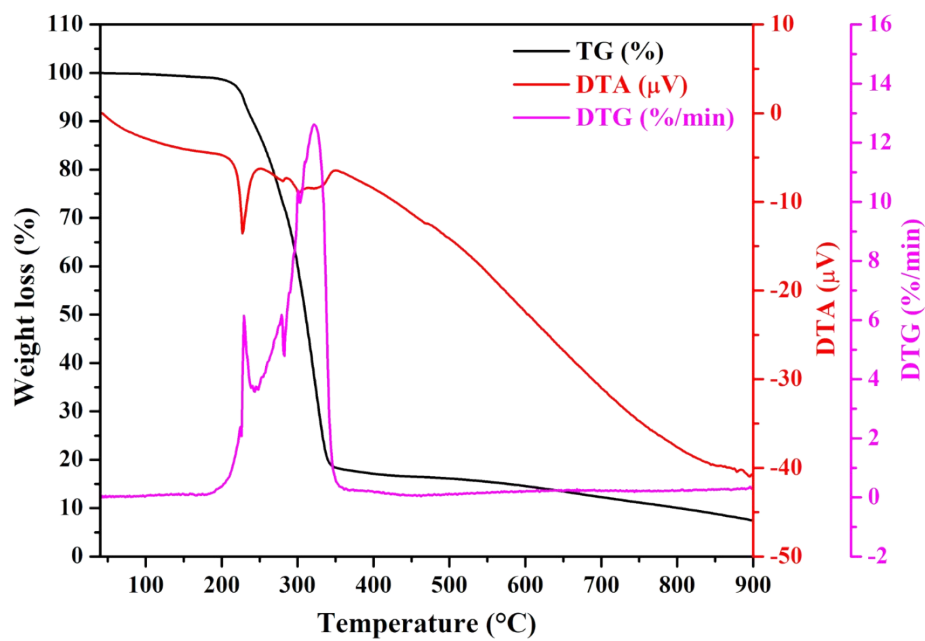


Fig. S9 TG-DTA-DTG curves for KA@CP-S3.

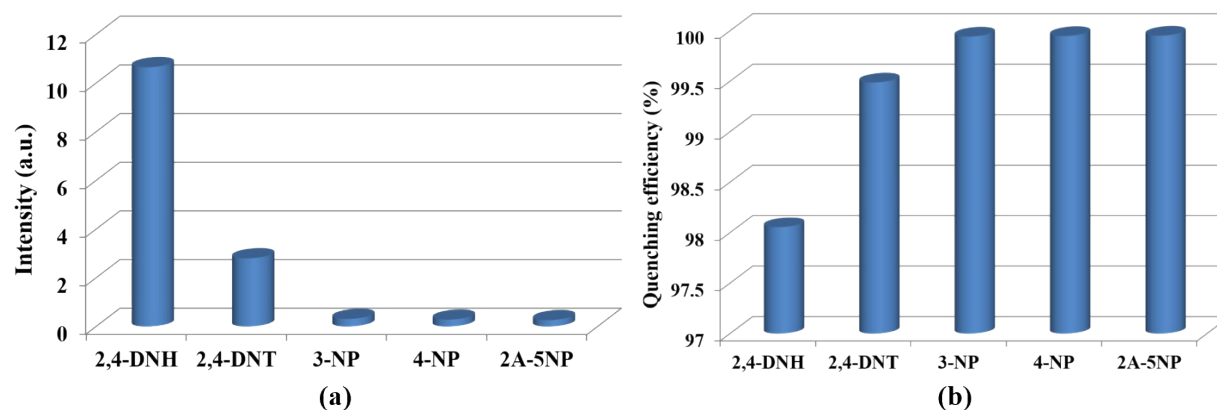


Fig. S10 (a) Fluorescence emission intensities and (b) Relative fluorescence quenching efficiencies of KA@CP-S3 with different nitrobenzene derivatives.

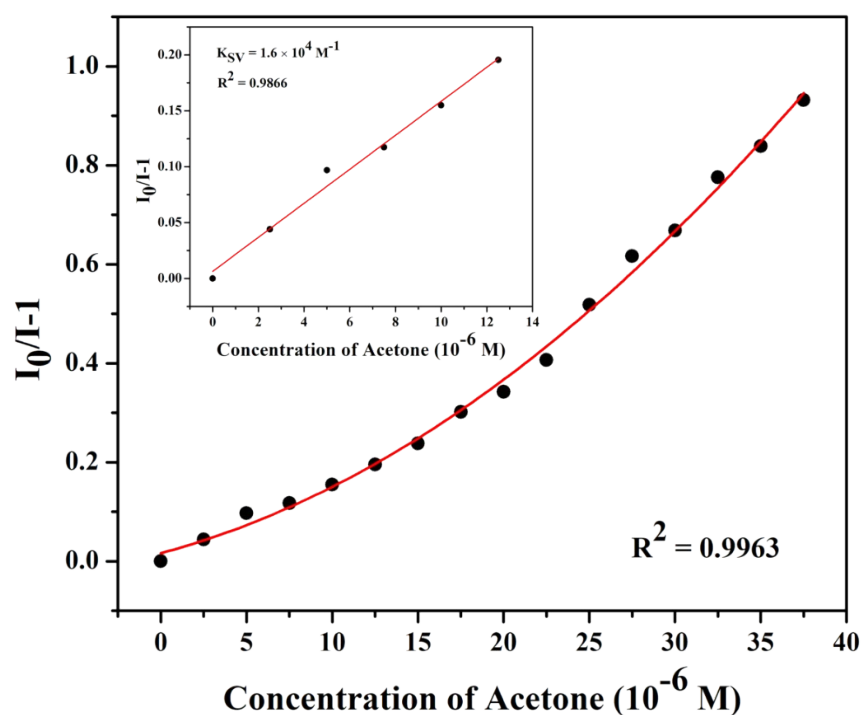


Fig. S11 S-V plot of KA@CP-S3 dispersed in ethanolic solution after gradual addition of Acetone (1 mM), the linear S-V curve of KA@CP-S3 at low concentration of Acetone (inset).

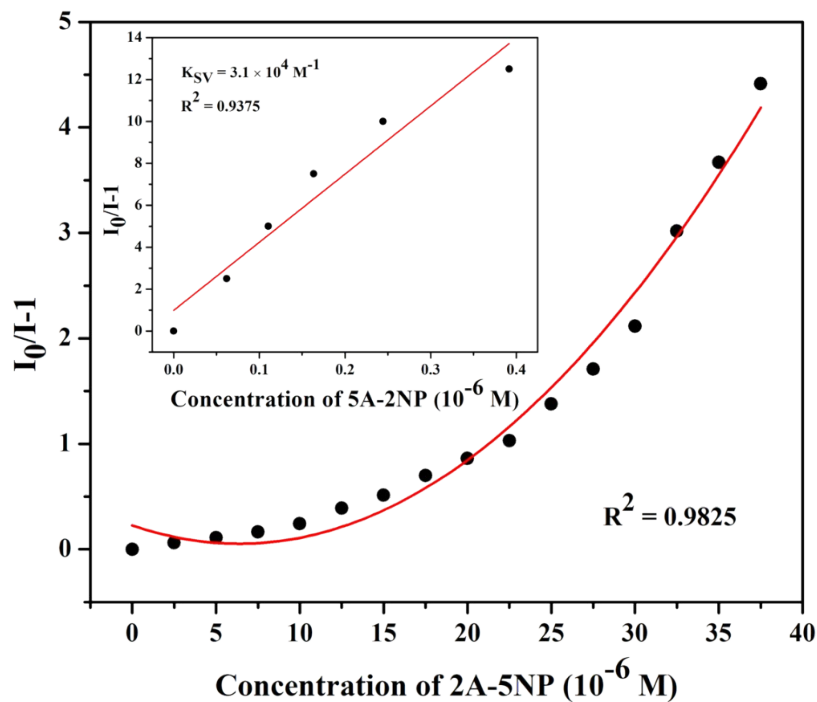


Fig. S12 S - V plot of **KA@CP-S3** dispersed in ethanolic solution after gradual addition of 2A-5NP (1 mM), the linear S - V curve of **KA@CP-S3** at low concentration of 2A-5NP (inset).

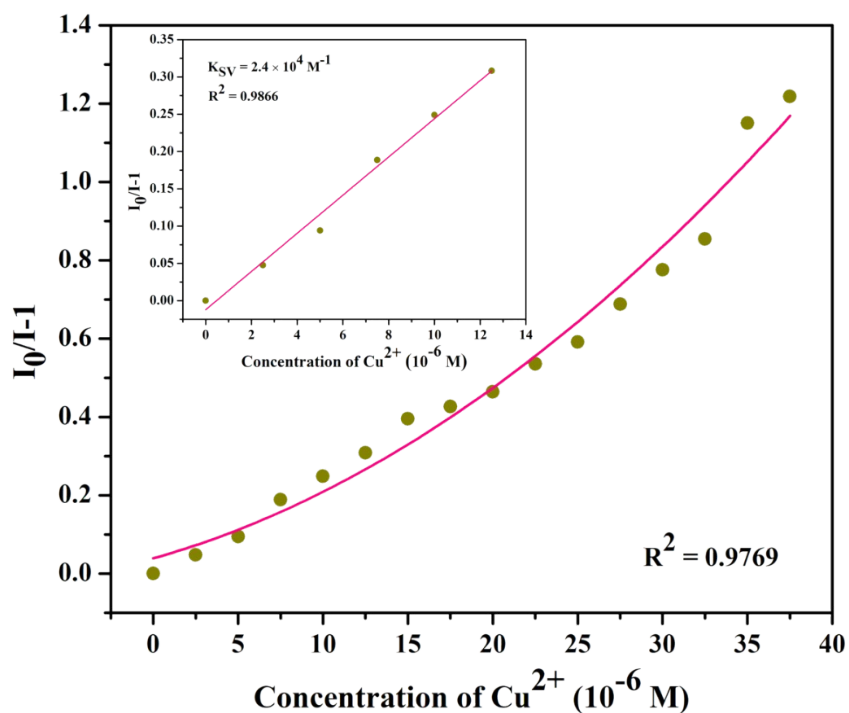


Fig. S13 S - V plot of **KA@CP-S3** dispersed in ethanolic solution after gradual addition of Cu^{2+} (1 mM), the linear S - V curve of **KA@CP-S3** at low concentration of Cu^{2+} (inset).

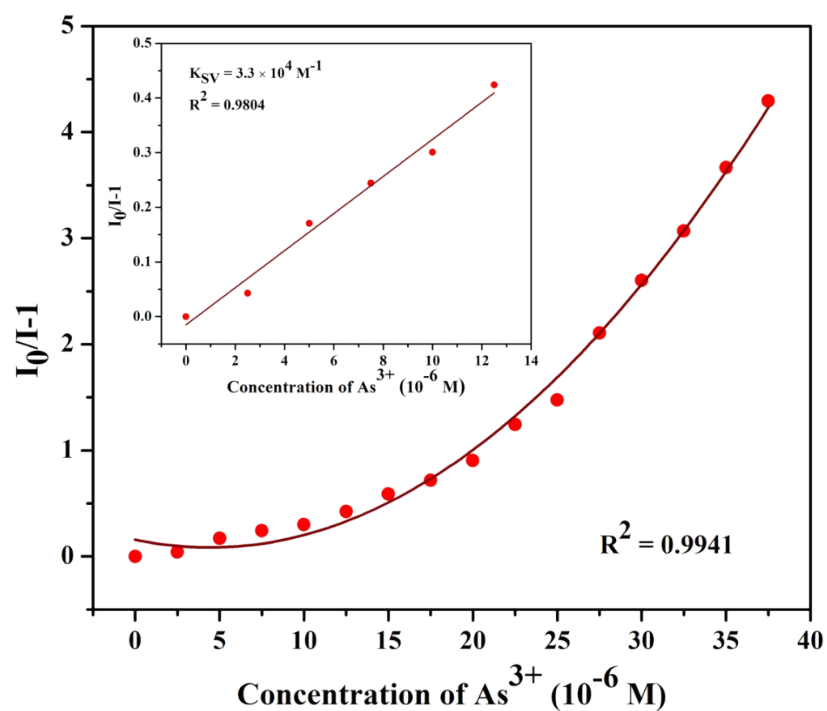


Fig. S14 S - V plot of **KA@CP-S3** dispersed in ethanolic solution after gradual addition of As^{3+} (1 mM), the linear S - V curve of **KA@CP-S3** at low concentration of As^{3+} (inset).

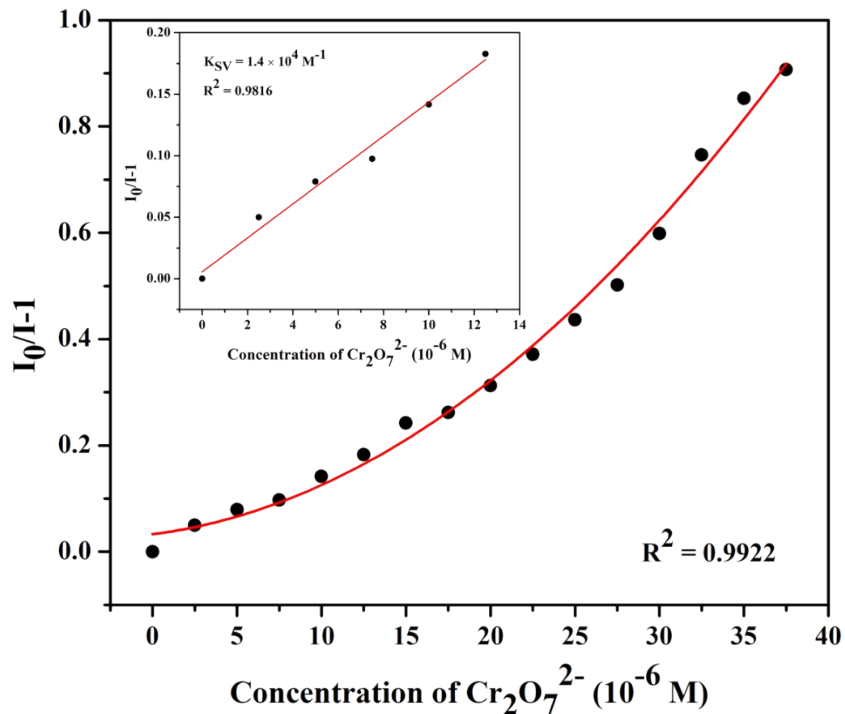


Fig. S15 S - V plot of **KA@CP-S3** dispersed in ethanolic solution after gradual addition of $\text{Cr}_2\text{O}_7^{2-}$ (1 mM), the linear S - V curve of **KA@CP-S3** at low concentration of $\text{Cr}_2\text{O}_7^{2-}$ (inset).

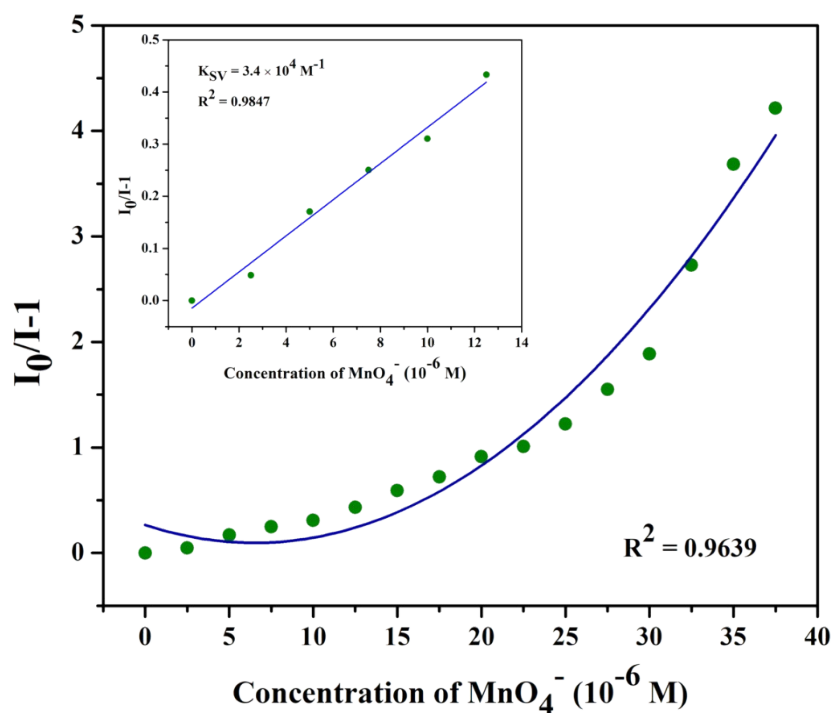


Fig. S16 S - V plot of **KA@CP-S3** dispersed in ethanolic solution after gradual addition of MnO_4^- (1 mM), the linear S - V curve of **KA@CP-S3** at low concentration of MnO_4^- (inset).

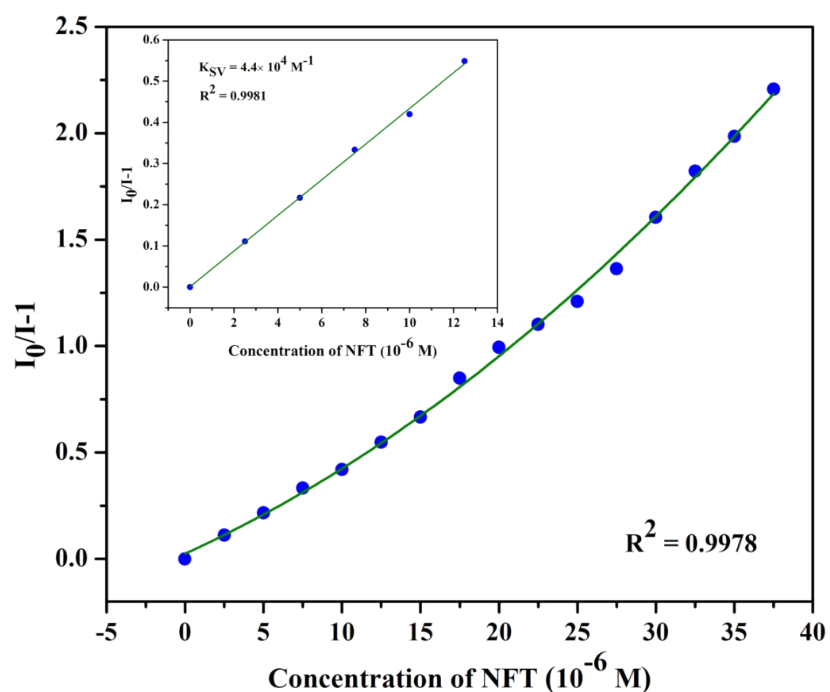


Fig. S17 S - V plot of **KA@CP-S3** dispersed in aqueous solution after gradual addition of **NFT** (1 mM), the linear S - V curve of **KA@CP-S3** at low concentration of **NFT** (inset).

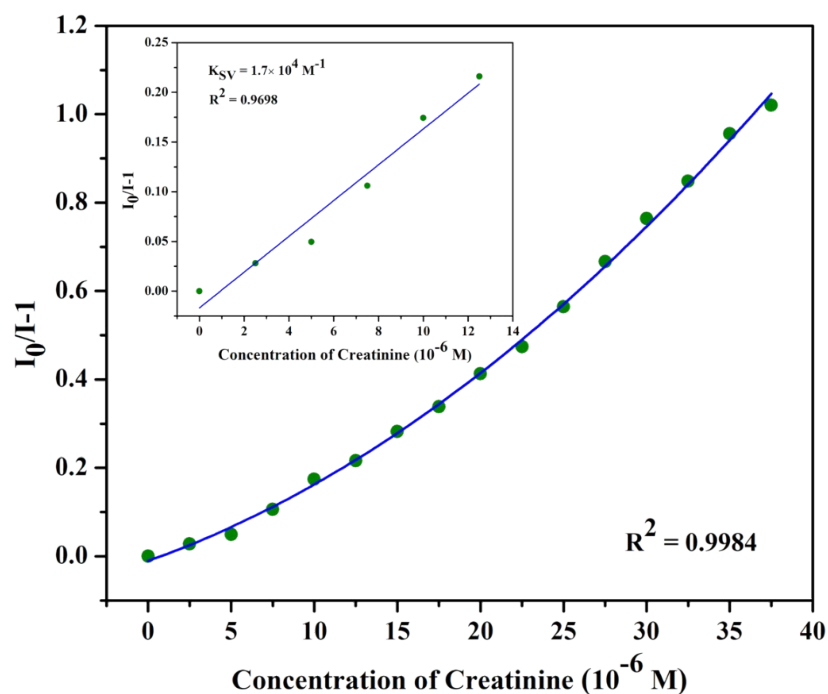


Fig. S18 S – V plot of **KA@CP-S3** dispersed in aqueous solution after gradual addition of Creatinine (1 mM), the linear S – V curve of **KA@CP-S3** at low concentration of Creatinine (inset).

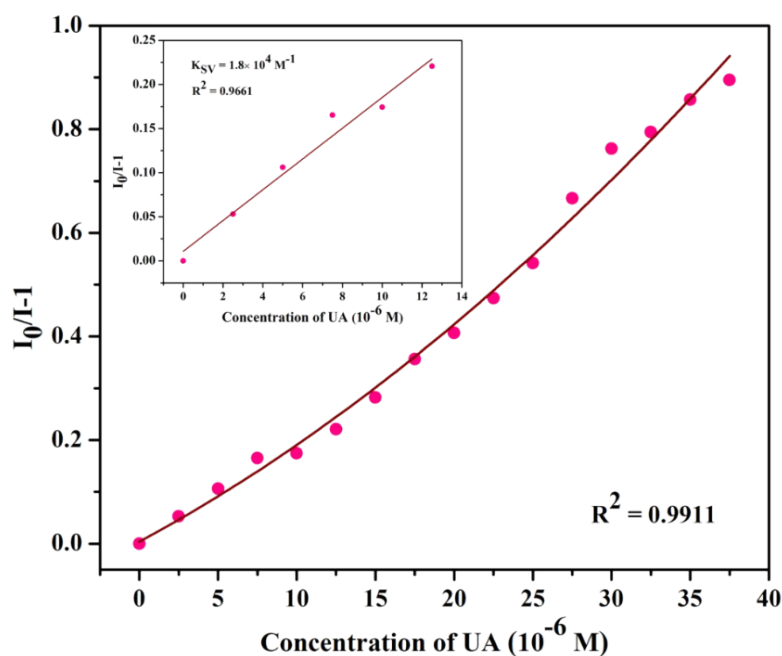


Fig. S19 S – V plot of **KA@CP-S3** dispersed in aqueous solution after gradual addition of UA (1 mM), the linear S – V curve of **KA@CP-S3** at low concentration of UA (inset).

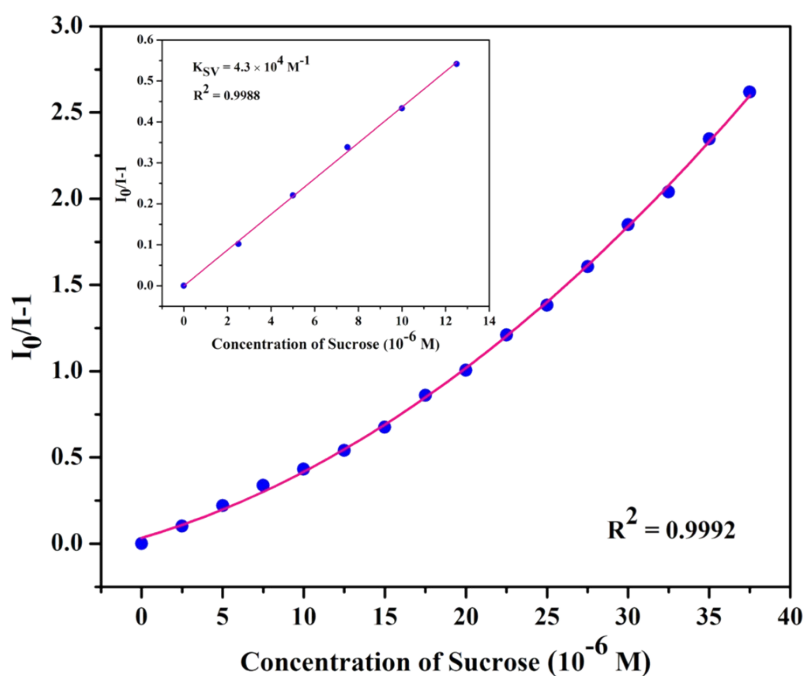


Fig. S20 $S-V$ plot of **KA@CP-S3** dispersed in aqueous solution after gradual addition of Sucrose (1 mM), the linear $S-V$ curve of **KA@CP-S3** at low concentration of Sucrose (inset).

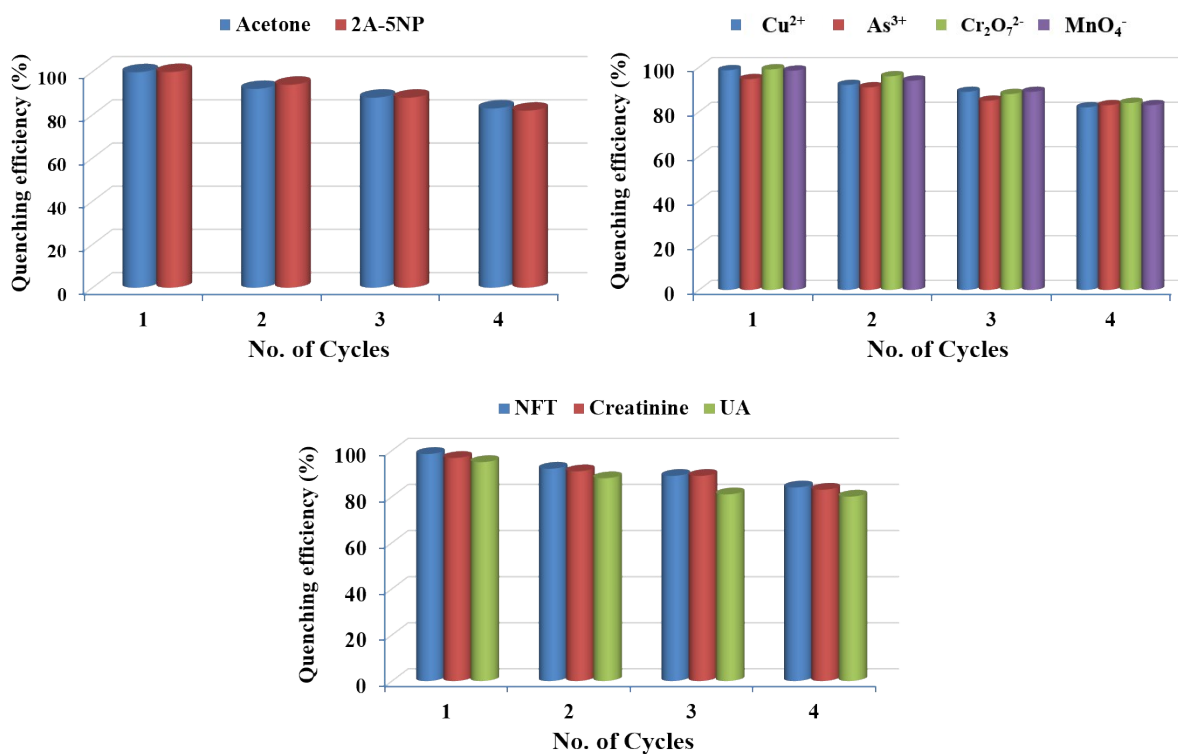


Fig. S21 Catalysts reusability of **KA@CP-S3** for each cycle.

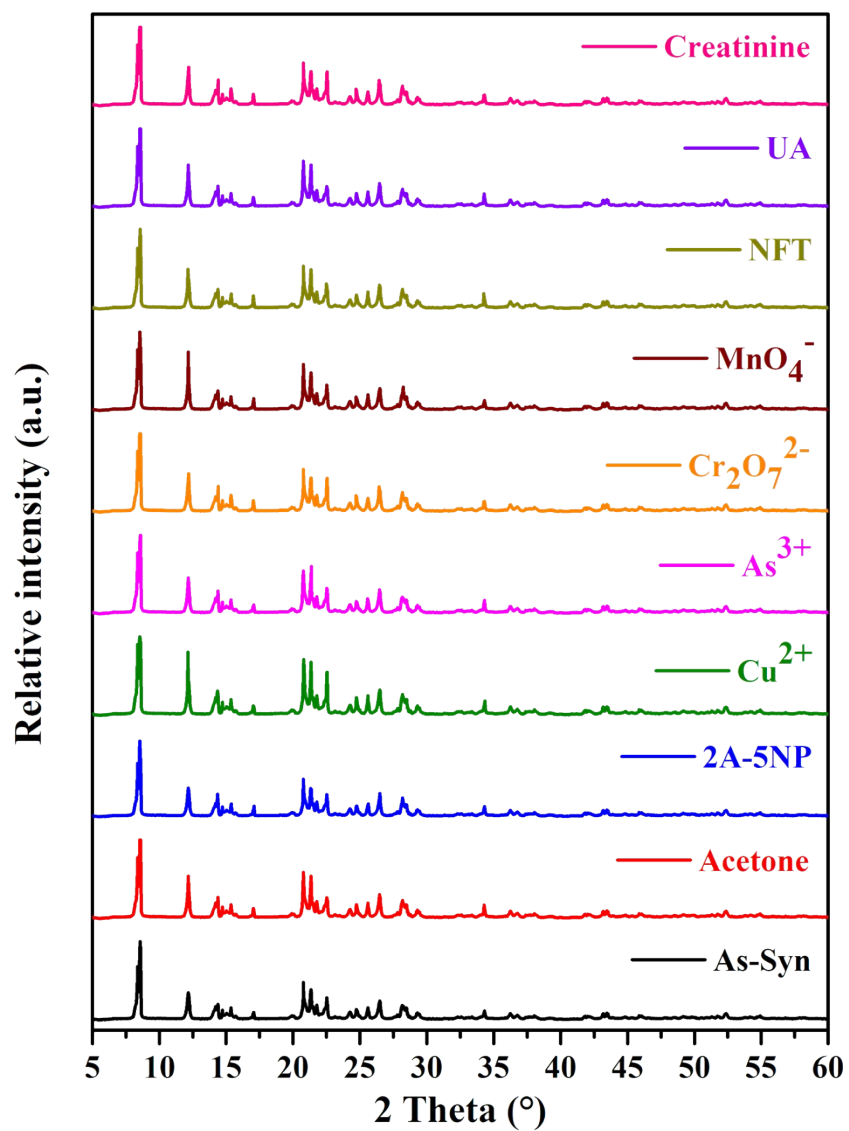


Fig. S22 PXRD patterns of KA@CP-S3 before and after fluorescence sensing of various analytes.

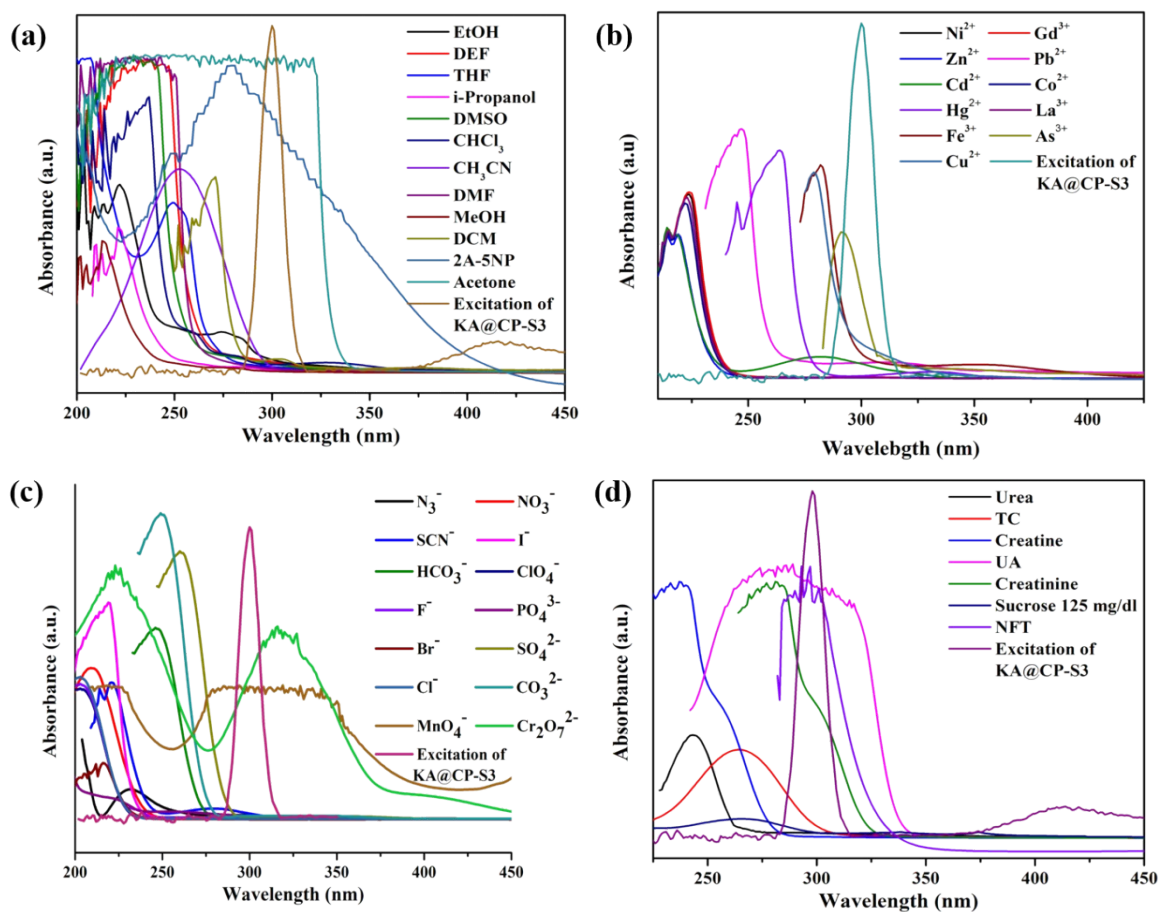


Fig. S23 Spectral overlap between the UV-Vis absorption spectra of various analytes and the excitation spectra of KA@CP-S3.

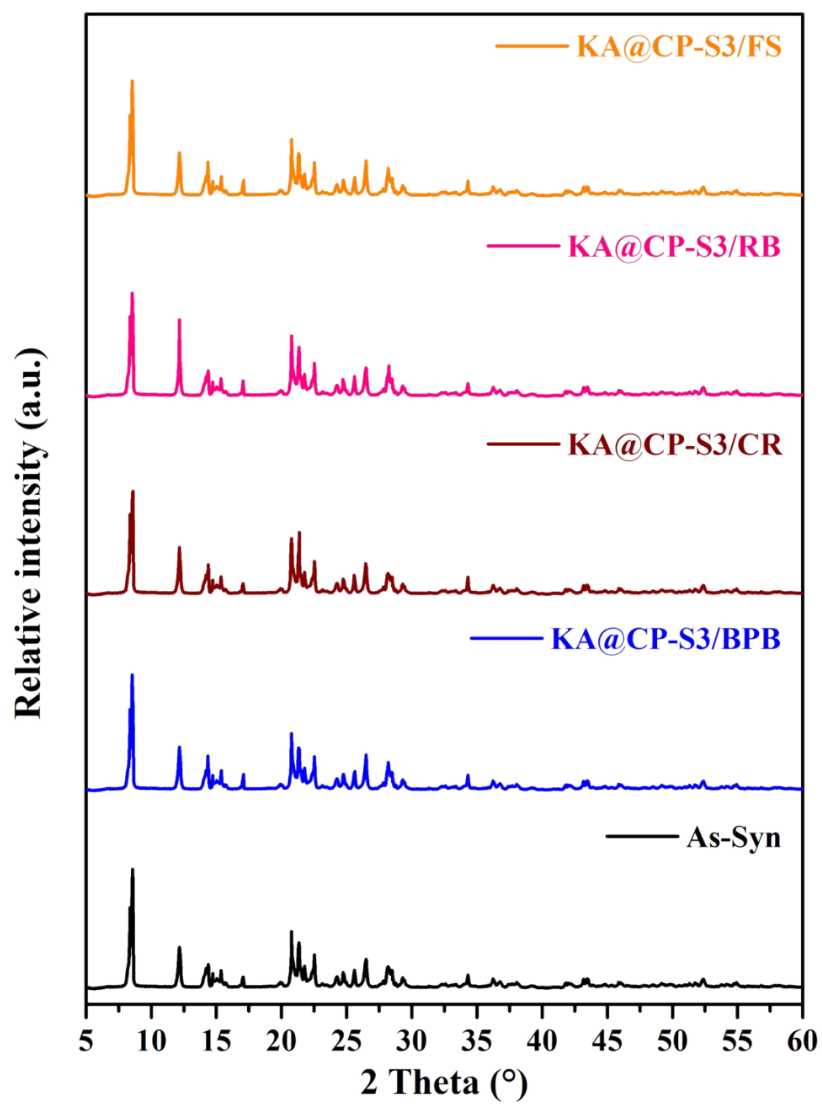


Fig. S24 PXRD patterns of **KA@CP-S3** before and after photocatalytic experiments with dyes (BPB, CR, RB and FS).

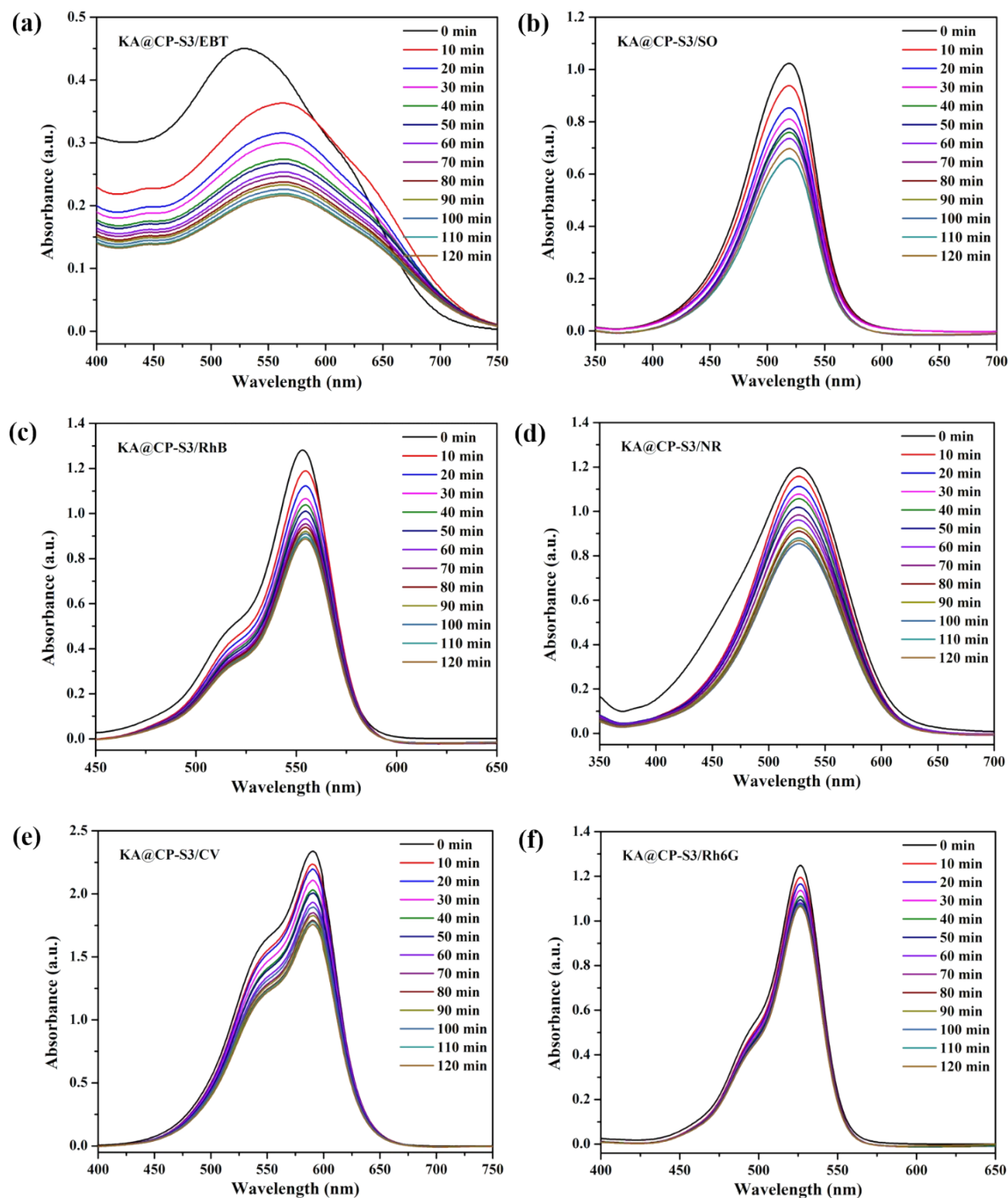


Fig. S25 Time-dependent UV-Vis absorption spectra of dyes (a) EBT, (b) SO, (c) RhB, (d) NR, (e) CV and (f) Rh6G with catalysts **KA@CP-S3**.

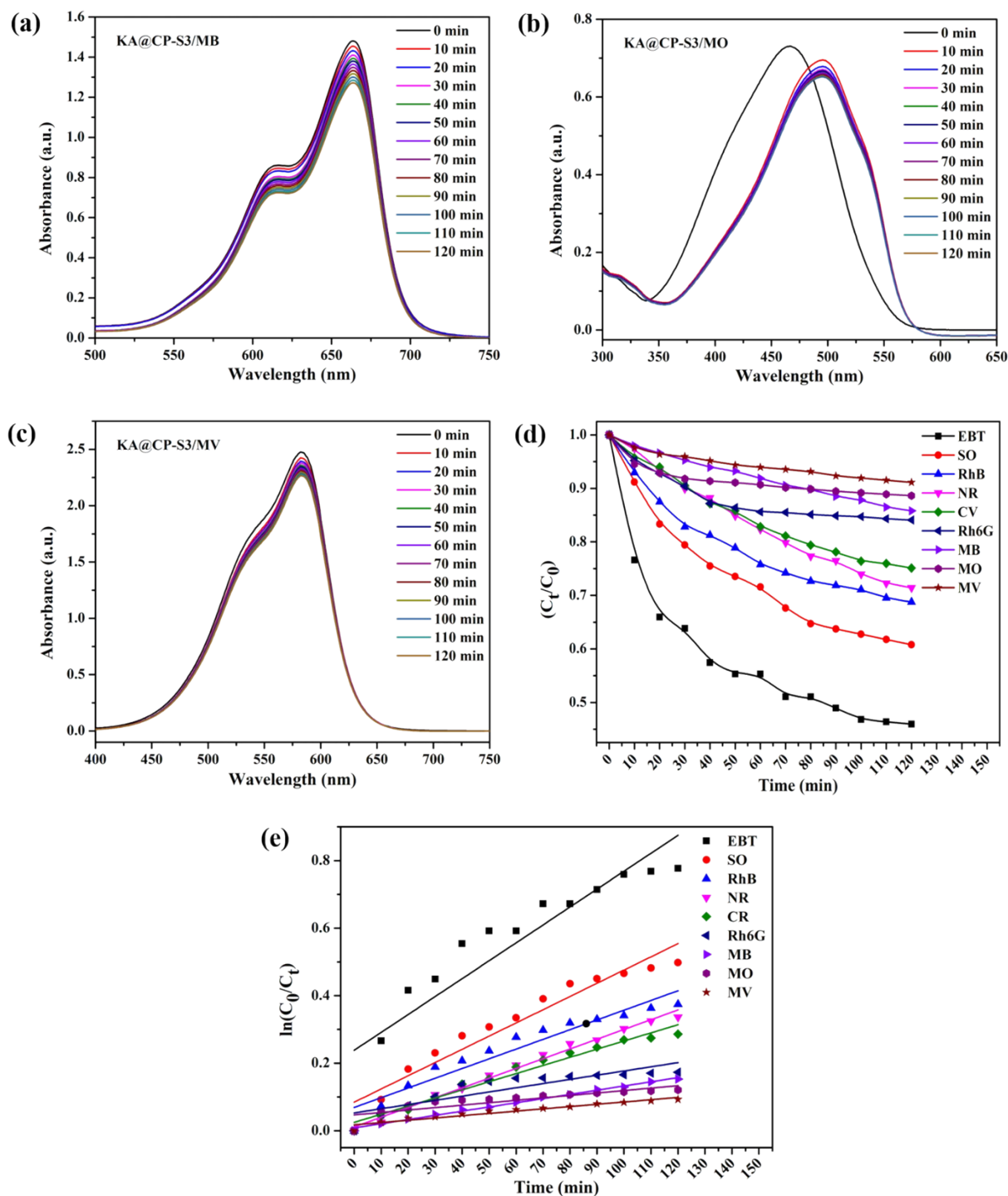


Fig. S26 Time-dependent UV-Vis absorption spectra of dyes (a) MB, (b) MO and (c) MV with catalysts **KA@CP-S3** (d) The Photocatalytic degradation kinetic plot of such dyes C_t/C_0 versus time interval plot and (e) Plot of $\ln(C_0/C_t)$ versus time interval in the presence of catalysts.

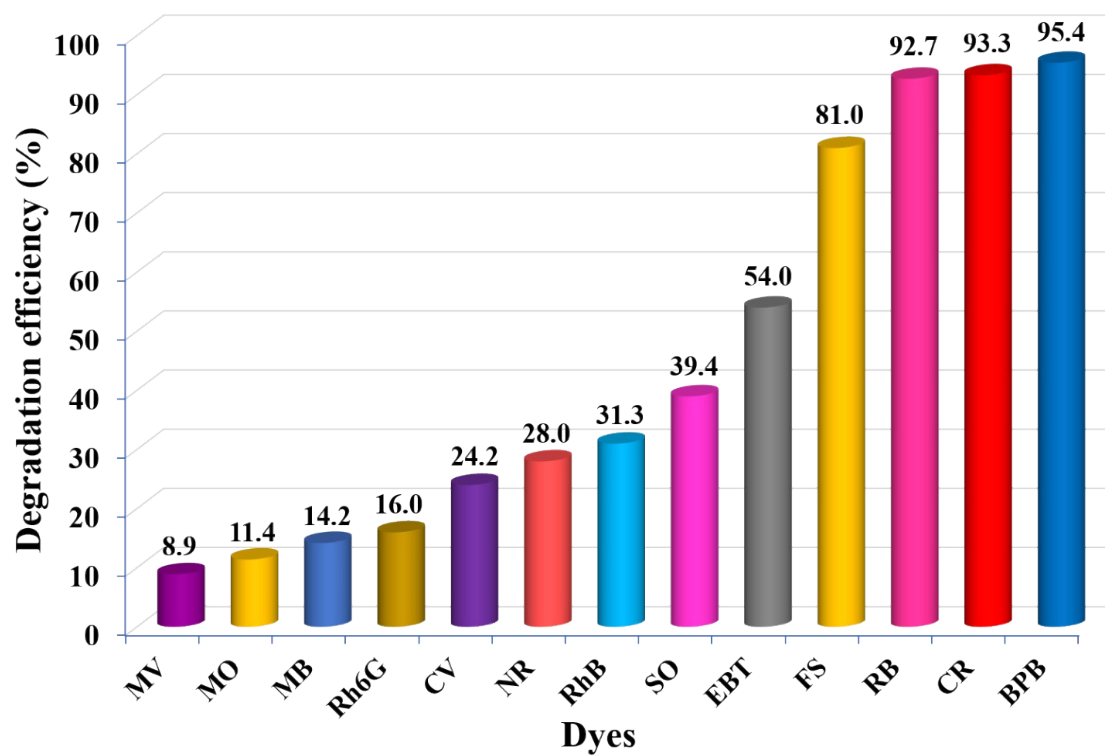


Fig. S27 Photocatalytic degradation efficiencies of KA@CP-S3 for dyes.

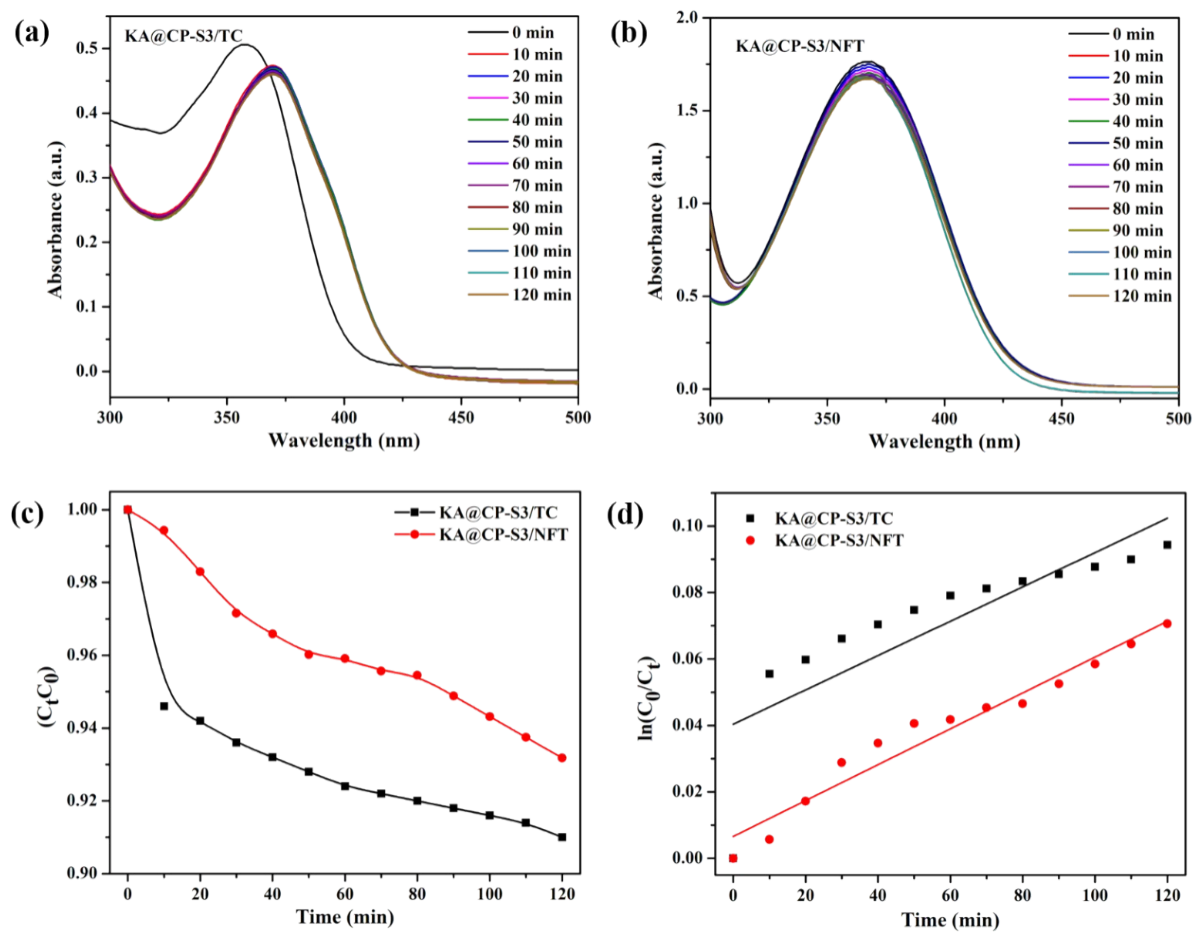


Fig. S28 Time-dependent UV-Vis absorption spectra of (a) TC and (c) NFT with catalysts KA@CP-S3 (d) The Photocatalytic degradation kinetic plot C_t/C_0 versus time interval plot and (e) Plot of $\ln(C_0/C_t)$ versus time interval in the presence of catalysts.

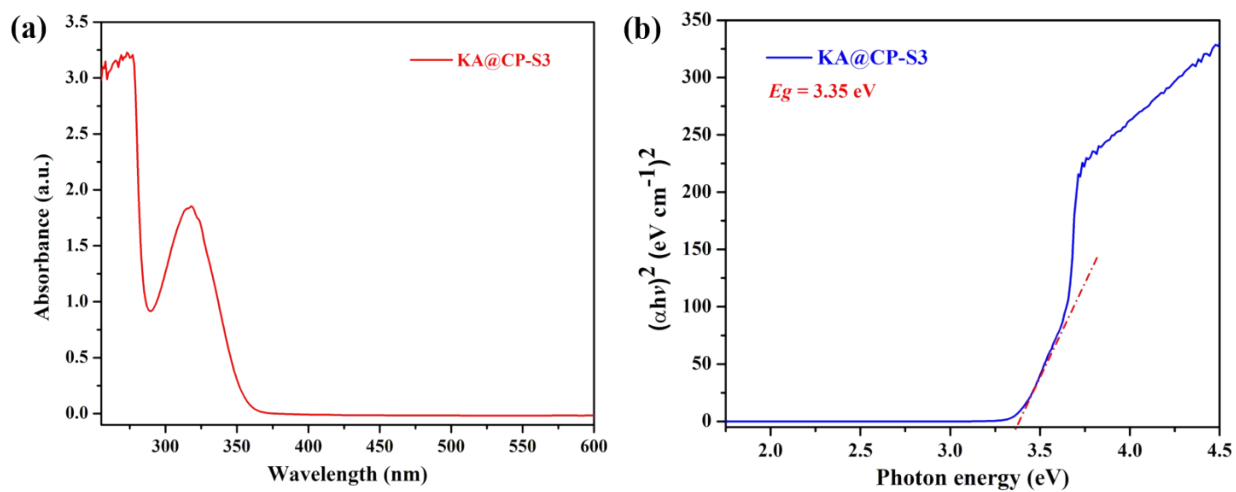


Fig. S29 (a) UV-Vis absorption spectra and (b) Tauc plot ($(\alpha h\nu)^2$ vs energy (eV)) for band gap study of KA@CP-S3.