

# **Facile preparation of fluorescent water-soluble none-conjugated polymer dots and fabricating acetylcholinesterase biosensor**

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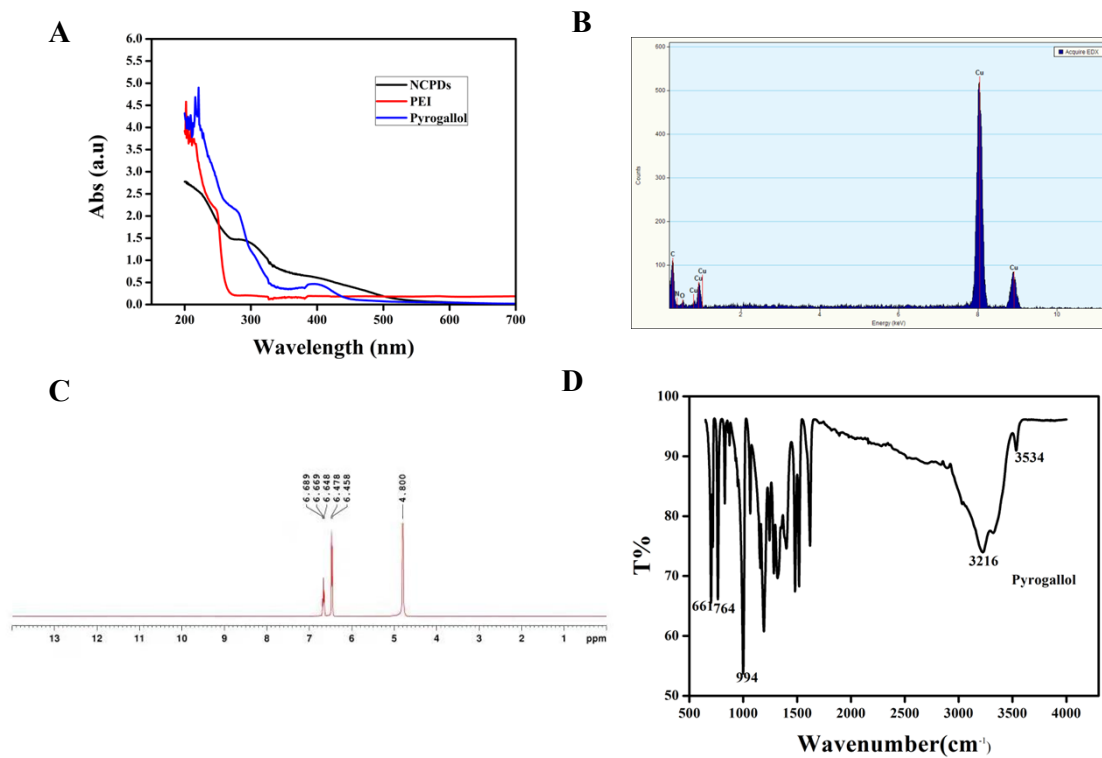
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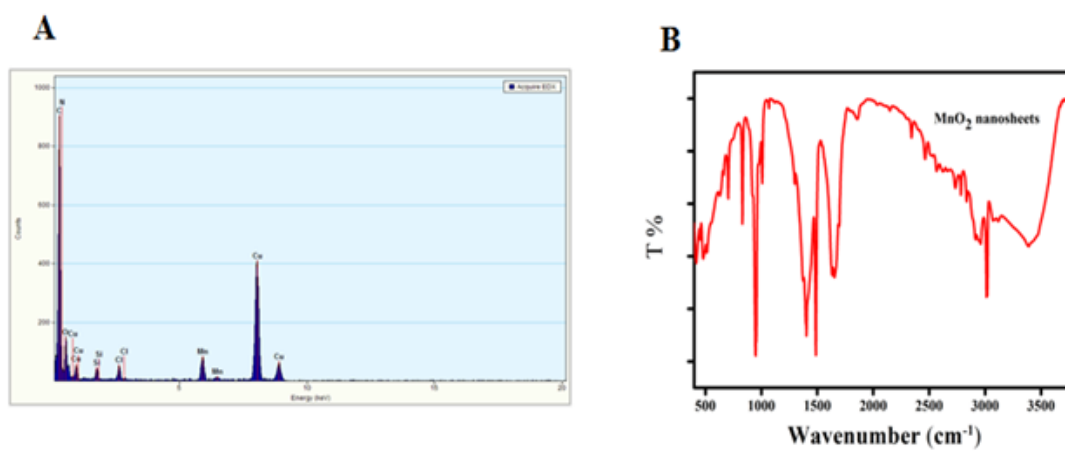
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**Fig. S1**



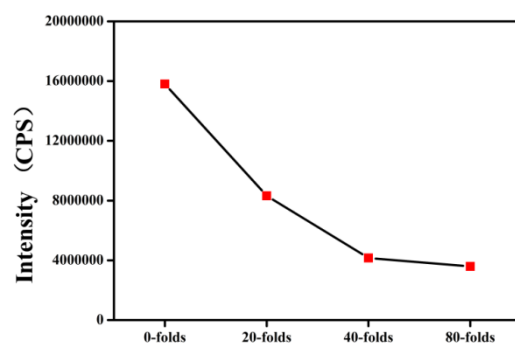
(A) UV-vis absorption spectra of NCPDs, PEI and Pyrogallol, respectively (B) The EDS of NCPDs (C) NMR spectra of Pyrogallol (D) FT-IR spectroscopy of Pyrogallol

**Fig.S2**



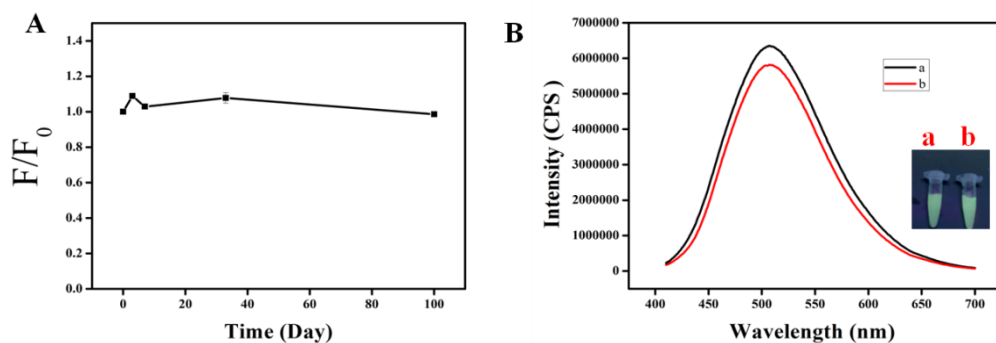
(A) The EDS of MnO<sub>2</sub> nanosheets (B) The FT-IR spectra of MnO<sub>2</sub> nanosheets

**Fig. S3**



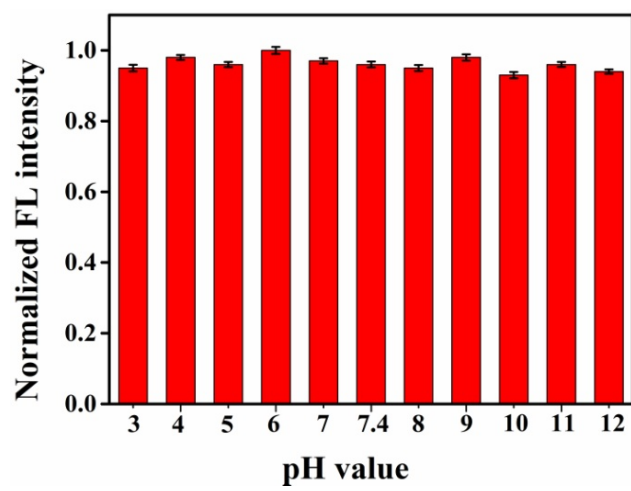
Relationship between fluorescent intensity and diluted folds of NCPDs

**Fig. S4**



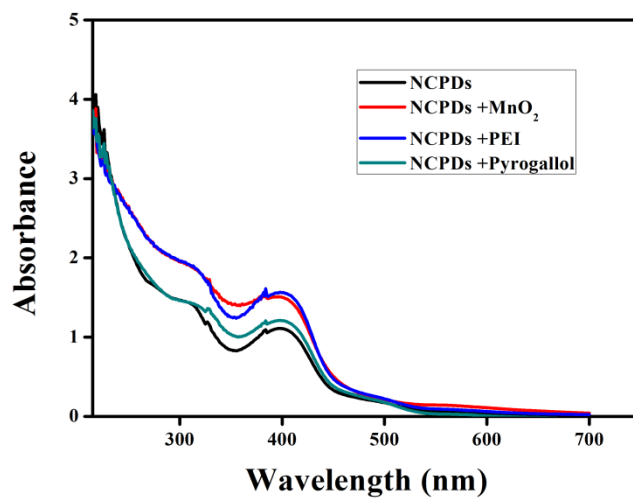
(A) The relationship between relatively fluorescent intensity ( $F/F_0$ ) of NCPDs and storage time (B) Fluorescent emission spectra of NCPDs fresh prepared (black line) and after preserved 6 months (red line) The inset photographs showed new and old NCPDs under UV light

**Fig. S5**



The stability of NCPDs at various PH values

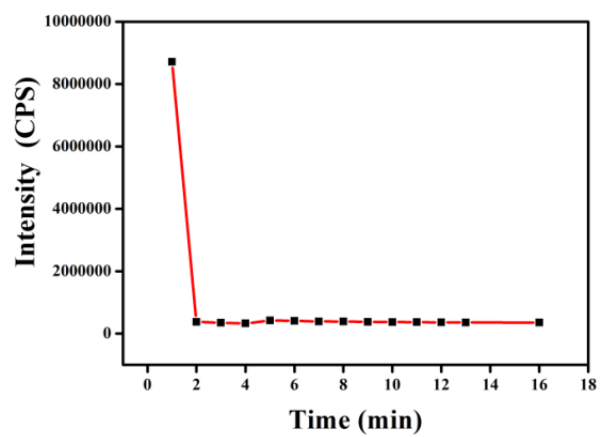
**Fig. S6**



UV-vis absorption spectra of NCPDs adding with MnO<sub>2</sub> nanosheets, PEI and Pyrogallol

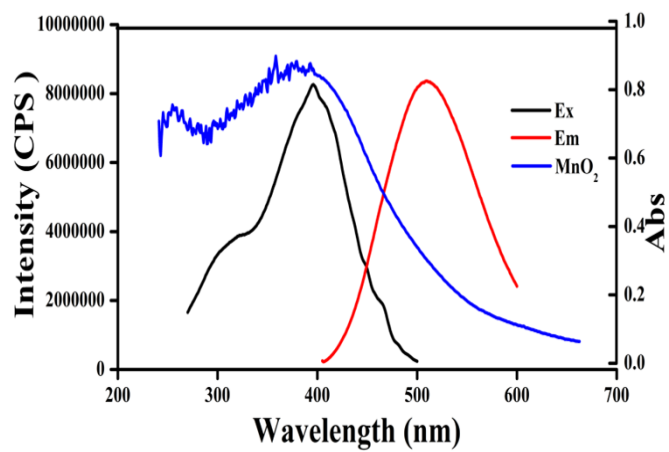


**Fig.S7**



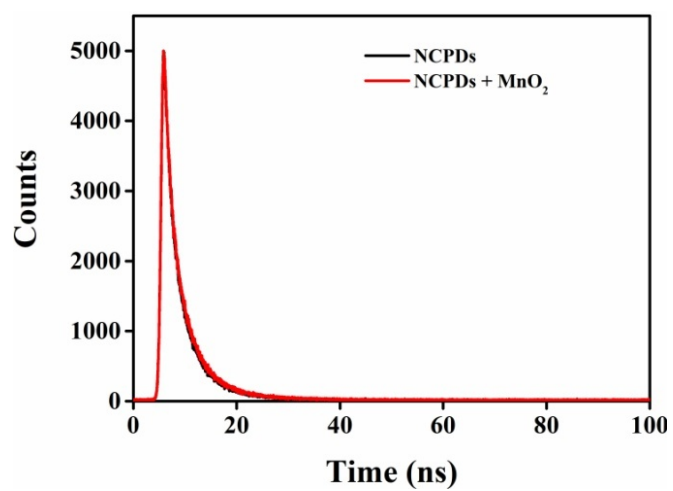
Response time of the NCPDs-MnO<sub>2</sub> biosystem (12.3 μM)

**Fig. S8**



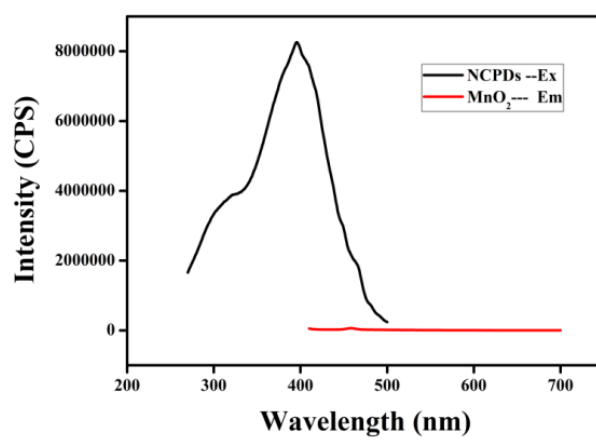
The overlap of excitation and emission spectra of NCPDs and UV-vis absorption spectra of MnO<sub>2</sub> nanosheets

**Fig. S9**



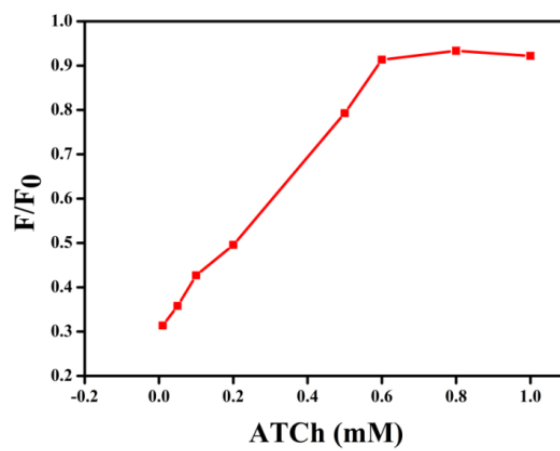
Decay curves of NCPDs in the absence (black curve) and presence (red curve) of MnO<sub>2</sub> nanosheets,  $\lambda_{\text{ex}} = 396 \text{ nm}$  and  $\lambda_{\text{em}} = 507 \text{ nm}$

**Fig. S10**



The overlap of excitation spectra of NCPDs and emission spectrum of MnO<sub>2</sub> nanosheets

**Fig. S11**



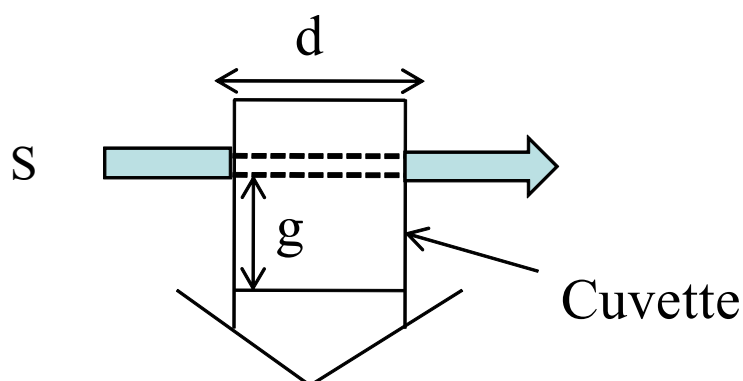
Optimization the substrate concentration

IE was analyzed by the following equation:

$$IE = \frac{F_{no-inhibitor} - F_{inhibitor}}{F_{no-inhibitor} - F_0} \quad (1)$$

$F_{inhibitor}$  and  $F_{no-inhibitor}$  represent the fluorescent intensity of AChE-NCPDs-MnO<sub>2</sub> system and AChE-NCPDs-MnO<sub>2</sub>-inhibitor system, respectively.  $F_0$  refers to the fluorescent intensity of the NCPDs-MnO<sub>2</sub> system without AChE and inhibitor. <sup>1</sup>

**Fig. S12**



Cuvette geometry and parameters used in equation

$$\frac{F_{cor}}{F_{obsd}} = \frac{2.3dA_{ex}}{1 - 10^{-dA_{ex}}} 10^{gA_{em}} \frac{2.3sA_{em}}{1 - 10^{-sA_{em}}} \quad (2)$$

**Table S1** IFE of MnO<sub>2</sub>nanosheets on the fluorescence of NCPDs

MnO <sub>2</sub> μM	A <sub>ex</sub>	A <sub>em</sub>	F <sub>obsd</sub>	F <sub>cor</sub>	CF	F <sub>cor</sub> /F <sub>coro</sub>
0.00	0.28152	0.075	5.03E+06	7.44E+06	1.48	1.00
1.27	0.34388	0.095	4.40E+06	7.09E+06	1.61	1.05
1.86	0.36919	0.105	4.10E+06	6.84E+06	1.67	1.09
2.92	0.44468	0.129	3.85E+06	7.12E+06	1.85	1.13
3.42	0.49078	0.148	3.64E+06	6.57E+06	1.98	1.03
6.81	0.71132	0.229	2.78E+06	7.34E+06	2.64	1.01

$F_{obsd}$  is the observed fluorescence intensity and  $F_{cor}$  is the corrected fluorescence intensity by removing IFE contribution from . and represent the absorbance at the excitation wavelength ( =396 nm) and maximum emission wavelength ( <sub>m</sub>=507 nm), respectively; s is the thickness of excitation beam (5 nm), g is the distance between the edge of the excitation beam and the edge of the cuvette

(0.60 cm in this case) and  $d$  is the width of the cuvette (1.00 cm) (Table S1 summarizes the parameters used in calculating the contribution of IFE to the fluorescence quenching process. Corrected factor (CF) is defined as  $CF = F_{\text{obsd}} / F_0$ . The maximum value of CF could not exceed 3; otherwise, the correction is not convincing.  $F_0$  and  $F$  are the corrected fluorescence intensities of NCPDs in the absence and presence of  $\text{MnO}_2$  nanosheets, respectively.

### Calculation of the Quenching Efficiency

The quenching efficiency (E) was calculated according to the formula

$$E = \left(1 - \frac{F}{F_0}\right) \times 100\% \quad (3)$$

$F_0$  is the fluorescence intensity of NCPDs and  $F$  represents the fluorescence intensity of NCPDs quenched by  $\text{MnO}_2$  nanosheets.

### Reference:

1. Y. Zhang, T. Hei, Y. Cai, Q. Gao and Q. Zhang, *Anal. Chem.*, 2012, **84**, 2830-2836.