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## **Supporting Information**

## One-pot green synthesis of water-soluble carbon nanodots with multicolor photoluminescence from polyethylene glycol

Moyun Chen, Weizhi Wang\*a and Xiaoping Wu\*b



**Fig. S1** The products originated from (A) PEG, (B) PPG and (C) PTMG at 160 °C for 2 h, respectively. (Top: under daylight, down: under 365 nm UV light).

**PAHS:** We utilized an efficient reaction route for preparing PAHs via a mild intramolecular oxidative cyclodehydrogenation (Scheme S1).





Fig. S2 High resolution mass spectrum (MOLDI-TOF) of PAHs.



Fig. S3 <sup>1</sup>H NMR and <sup>13</sup>C NMR spectra of PAHs in CD<sub>2</sub>Cl<sub>2</sub>.



**Fig. S4** FTIR spectra of (a) CNDs originated from PPG by heating for 1 h, denoted as p-CNDs@1h; 2 h, denoted as p-CNDs@2h; 6 h, denoted as p-CNDs@6h, (b). CNDs originated from PTMG by heating for 3 h, denoted as t-CNDs@3h; 5 h, denoted as t-CNDs@5h; 10 h, denoted as t-CNDs@10h.



Fig. S5 <sup>1</sup>H NMR spectra of CNDs from (a) PPG and (b) PTMG.



**Fig. S6** TEM images of (a) e-CNDs@0.5h and (b) e-CNDs@6h; Diameter distribution of (c) e-CNDs@0.5h and (d) e-CNDs@6h, the red line is the Gaussian fitting curve.



**Fig. S7** TEM images of e-PAHs-CNDs@2h, inset image: diameter distribution of e-PAHs-CNDs@2h (the red line is the Gaussian fitting curve).



Fig. S8 UV-vis spectra of (a) e-PAHs-CNDs@2h and e-CNDs@0.5h at different pH varied from 1 to 11.



Fig. S9 Fluorescence excitation spectra of CNDs (red: e-CNDs@0.5h, blue: e-CNDs@6h, black: e-PAHs-CNDs@2h).



**Fig. S10** 2D-fluorescence topographical map of CNDs. (a) e-CNDs@0.5h and dispersed in toluene, (b) e-PAHs-CNDs@2h and dispersed in water.



Fig. S11 (a) Effect of pH on the fluorescence intensity at 450 nm (at 360 nm excitation).



Fig. S12 PL spectra of CNDs. 1: e-CNDs@0.5h, 2: e-CNDs@6h, 3: e-PAHs-CNDs@2h, corresponding 1\*, 2\* and 3\*: samples storaged in the ambient environment for 1 month. (e-CNDs@0.5h, e-CNDs@6h, e-PAHs-CNDs@2h were excitated at 360 nm, 390 nm and 470 nm, respectively.)



Fig. S13 Integrated fluorescence intensity versus absorbance plot CNDs and references. (a) e-CNDs@0.5h and quinine sulfate solution; (b) e-CNDs@6h and fluorescein solution.

Table S1 The value of quantum yields of CNDs.			
	М	R	Φ
Quinine sulfate	1.34E+08	0.997	55
e-CNDs@0.5h	6.15E+06	0.998	2.51
Fluorescein	1.50E+08	0.999	93
e-CNDs@6h	5.83E+06	0.996	3.58