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## **Supporting Information**

## A Highly Sensitive "ON-OFF-ON" Dual Optical Sensor for Cu(II) Ion

## and Triazole Pesticides Based on Novel BODIPY-Substituted Cavitand

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## 1. Fluorescence Quantum Yield

Fluorescence quantum yield ( $\Phi_F$ ) is determined by the comparative method according to the Equation 1:

$$\Phi_F = \Phi_{FStd} \frac{F.A_{Std}.n^2}{F_{Std}.A.n_{Std}^2}$$
(1)

where F and F<sub>Std</sub> are the areas under the fluorescence emission curves of cavitand **3** and the standard, respectively. A and A<sub>Std</sub> are the respective absorbance of the cavitand **3** and standard at the excitation wavelengths, respectively. n<sup>2</sup> and n<sup>2</sup><sub>std</sub> are the refractive indices of solvents used for the sample and standard, respectively. Rhodamine 6G (in water) was employed as a standard compound ( $\Phi_F = 0.95$ ) in this study. The absorbances of the studied cavitand **3** and the standard Rhodamine 6G were kept ca. 0.05 at the excitation wavelength.



**Fig. S1** Full <sup>11</sup>B NMR spectra of (a) ethynyl-BODIPY **2** and (b) cavitand **3**; full <sup>19</sup>F NMR spectra of (c) ethynyl-BODIPY **2** and (d) cavitand **3**.



Fig. S2 FT-IR spectra (3500-800 cm<sup>-1</sup>) of (a) cavitand 1; (b) ethynyl-BODIPY 2 and (c) cavitand 3.



Fig. S3 MALDI-TOF spectrum of BODIPY functionalized resorcin[4]arene cavitand (3) (Matrix: DIT).



**Fig. S4** MALDI-TOF spectrum of BODIPY functionalized resorcin[4]arene cavitand (**3**) after the addition of Cu<sup>2+</sup> solution (Matrix: DIT).