Electronic Supplementary Material (ESI) for Journal of Materials Chemistry C. This journal is © The Royal Society of Chemistry 2014

Electronic Supporting Information

Non-classical donor-acceptor-donor chromophores. A strategy for high two-photon brightness

Adina I. Ciuciu,^a Dikhi Firmansyah,^b Vincent Hugues,^c Mireille Blanchard-Desce,*^c Daniel T.

Gryko,*^{*b*, *d*} Lucia Flamigni*^{*a*}

^a Istituto per la Sintesi Organica e Fotoreattivita' (ISOF), CNR, Via P. Gobetti 101, 40129 Bologna. Italy. Fax: +39 (0)51 639 98 44; Tel: +39 (0)51 639 98 12; E mail: <u>lucia.flamigni@isof.cnr.it</u> ^b Warsay, University of Technology, Neghowskiege 2, 00 664, Warsay, Boland, Faw, +48 22 628 27

^b Warsaw University of Technology, Noakowskiego 3, 00-664, Warsaw, Poland; Fax: +48 22 628 27 41; Tel: +48 22 234 58 01

^c Université de Bordeaux, ISM (UMR5255 CNRS),F 33400, Bordeaux, France, E-mail: <u>mireille.blanchard-desce@iu-bordeaux.fr</u>

^d Institute of Organic Chemistry Polish Academy of Sciences, Kasprzaka 44/52 01-224, Warsaw, Poland. Fax: +48 22 632 66 81; Tel: +48 22 343 30 63, E-mail: <u>dtgryko@icho.edu.pl</u>

Table of Contents

1.	Linear optical absorption and emission spectroscopy	page	2
2.	2PA properties	page	10
3.	NMR Spectral data	page	11



Fig. S1 Absorption spectra of the studied compounds in DCM.



Fig. S2 Luminescence spectra of optically matched solutions of reference monomers in DCM at room temperature. Excitation wavelength was 325 nm and $A_{325 \text{ nm}} = 0.07$.



Fig. S3 Luminescence spectra of optically matched solutions of the dimers **8b-c** and **10a-c** in DCM at room temperature. Excitation wavelength was 424 nm and $A_{424 \text{ nm}} = 0.07$ for **10a-c** and $A_{424 \text{ nm}} = 0.03$ for **8b-c**.































Fig. S4 Excitation spectra (black) overlaid with the corresponding absorption spectra (red) in TOL and DCM.

Compound	Solvent	λ_{abs}^{max}/nm	$\lambda_{\mathrm{em}}^{\mathrm{max}}/\mathrm{nm}$	Stokes shift/ cm ⁻¹	$\phi_{\mathrm{fl}}{}^{\mathrm{a}}$
90	DMSO	480	567	3200	0.54
ðC	H ₂ O	488	655	5220	0.04
10.0	DMSO	451	518	2870	0.43
100	H ₂ O	494	556	2260	0.13

Table S1 Absorption and fluorescence properties of compounds 8c and 10c in water and dimethyl sulfoxide.

^a ϕ_{fl} = fluorescence quantum yield, standard Rhodamine 6G in EtOH (ϕ_{fl} =0.94)



Fig. S5 Arbitrary scaled RT and 77K total and delayed luminescence of monomers in TOL, upon excitation at 325 nm.



Fig. S6 Lifetime measured for **3c** at 77 K in TOL rigid matrixes in the phosphorescence region (full circles), $\lambda = 507$ nm, and in the delayed fluorescence region (empty circles), $\lambda = 406$ nm.



Fig. S7 Arbitrary scaled RT luminescence in TOL and 77 K total and delayed luminescence of 8c and 10a, 10c in TOL with 50% EtI, upon excitation at 424 nm.



Fig. S8 Comparison of one photon and 2PA absorption spectra for the bis-imidazo[1,2-*a*]pyridines (recorded in DCM).































