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Addition/Correction

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Correction to "Through-Space Charge Transfer in Rod-Like Molecules: Lessons from Theory"

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We realized that the data and graph listed for oligomeric series **O3** are incorrect in our contribution. Below are the corrected values and graphs. Consequently, the paragraph starting from "With that respect, phenylacetylene..." (bottom of page 11948) up to "..transition in **O3** than in **O4**" should be removed from the original article.

Table 1. Computed Transition Wavelengths (nm) and Oscillator Strength for O3[a]

N	λ	f	q^{CT}	d^{CT}	H
1	302	0.47	0.60	2.23	0.78
2	356	1.18	0.70	4.17	4.10
3	374	2.07	0.65	5.51	5.87
4	381	3.14	0.57	5.55	7.47
5	387	4.28	0.53	4.73	8.80
6	391	5.38	0.51	3.75	10.08
7	395	6.45	0.50	2.99	11.34
8	397	7.49	0.50	2.38	12.60
9	400	8.52	0.49	1.86	13.89
10	401	9.57	0.49	1.60	15.18

[a] q^{CT} is in e, d^{CT} and H are in Å. All values have been obtained at the CAM-B3LYP/6-311++G(d,p)//PBE0/6-311G(d,p) level.

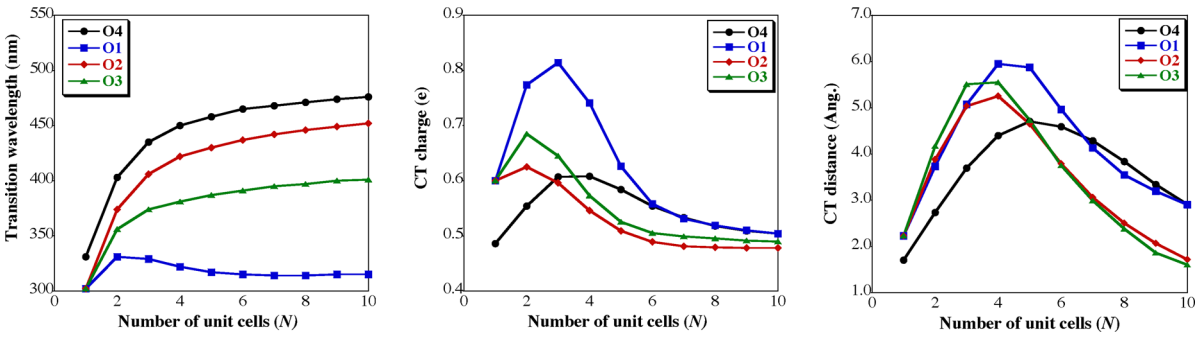

Figure 1. Evolution of the main parameters for increasingly long chains. Left: the vertical transition wavelength (nm). Center: amount of transferred charge (e^-). Right: charge-transfer distance (Å).

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