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Addition/Correction

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Correction to "Charge Shift Bonding Concept in Radical π-Dimers"

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The M05-2X/6-31+G(d,p) Laplacian values were reported incorrectly on pp 13942−13943. The correct calculated vales are the following:

	D (Å)	Laplacian[a]
H_2	0.74	−1.387
F_2	1.37	0.322
TCNE	2.89	0.0346

[a]Keith, T. A. AIMALL- version 10.02.09(46), http://aim.tkgristmill.com/.

The reported corrections leave the conclusions in our paper unchanged.

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