Chemical Science

CORRECTION

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a)

2.0

1.5

0.2

0.0

Correction: Tuning radical interactions in trisradical tricationic complexes by varying host-cavity sizes

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Correction for 'Tuning radical interactions in trisradical tricationic complexes by varying host-cavity sizes' by Kang Cai et al., Chem. Sci., 2020, DOI: 10.1039/c9sc04860j.

0.6

0.4

Abs 0.2 K, =

 $0.6) \times 10^{3} M$

1200

K_ =

2 3

1200

1000

Ż 3

The authors regret that the K_a values in Fig. 1c and d are incorrect in the original manuscript. The correct figure is displayed below.

b) 2.0

1.5

0.5

0.0

600

800

 λ / nm



K_ =

0.6

0.4

SQN 0.2

Fig. 1 Vis-NIR Spectra (MeCN, 2 mm cuvette) on titrating MV⁺ into (a) PyBB²⁽⁺⁺⁾ (0.25 mM); (b) mpBB²⁽⁺⁺⁾ (0.25 mM); (c) DThBB²⁽⁺⁺⁾ (0.20 mM); (d) ThBB²⁽⁺⁺⁾ (0.25 mM). The inset shows the change in absorption at 1080 nm with titration. The curve fitting is highlighted in red.

The Royal Society of Chemistry apologises for these errors and any consequent inconvenience to authors and readers.

1200

1000

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600

800

 λ/nm



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