A modular approach to the synthesis of 2,3,4-trisubstituted tetrahydrofurans

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In Table 3 the structures of the [1,3] products are shown incorrectly in the print and online pdf versions of the article as a result of a typesetting error. The corrected Table 3 is given below.

	R ₂ I Pd(OAc) ₂ , PPI K ₂ CO ₃ , <i>n</i> -Bu ₄ N MeCN/H ₂ O R ₁ 50 °C, 12-36I	MeCN -40 °C, 1h	R ₂ CHO
Entry	R ₁ , R ₂ , Yield (%), dr ^a	[1,3] Product	Yield (%), dr
1	8; R ₁ = Ph R ₂ = Ph 75, (87:13)	Ph. CHO O 9a Ph	70, (94:5:1:<1)
2	10 ; R ₁ = 2-Furyl R ₂ = Ph 65, (87:13)	Ph. CHO	94, (>95:5:<1:<1)
3	12; $R_1 = CHCHPh$ $R_2 = Ph$ 71, (83:17)	Ph. CHO Ph	88, (96:3:1:<1)
4	14; R ₁ = CH ₂ CH ₂ Ph R ₂ = Ph 65, (85:15)	Ph, CHO Ph	85, (96:3:1:<1)
5	p-O 16 ; R ₁ = CH ₂ CH ₂ Ph R ₂ = p -OMePh 59, (85:15)	OMePh, CHO Ph	84, (91:6:2:1)
6	18 ; R ₁ = Et R ₂ = Ph 64, (85:15)	Ph, CHO Et	97, (90:7:2:<1)
7	20 ; R ₁ = <i>i</i> -Pr R ₂ = Ph 67, (85:15)	Ph. CHO	83, (85:10:5:<1)
8	22 ; R ₁ = <i>t</i> -Bu R ₂ = Ph 68, (83:17)	Ph. CHO 23 t-Bu	55, (70:18:12:<1)
9	24 ; $R_1 = CH_2CH_2Ph$ $R_2 = CHCHPh$ 42 , (79:21)	CHO Ph	71, (93:6:1:<1)
10	26; R ₁ = CH ₂ CH ₂ Ph R ₂ = CHC(CH ₃)Ph 71, (78:22)	Ph Me CHO Pr	79, (83:17:<1:<1)
11	28; R ₁ = CH ₂ CH ₂ SPh R ₂ = Ph 59, (83:17)	Ph, CHO SPh	68, (83:13:4:<1)

The tuneable complexation of gold nanoparticles

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The fourth sentence of paragraph five has been revised to read as follows:

Square wave voltammetry indicated that the first reduction wave of 6 was shifted by ~ -20 to -30 mV upon addition of 5, presumably due to donor–acceptor interactions (resulting from complex formation between the naphthalene and cyclophane moieties) destabilizing the diradical dication state of the cyclophane (Fig. 2).

Porous 3-D honeycomb architecture by self-assembly of helical H-bonded molecular tapes

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The caption for Scheme 1 is incorrect. The correct caption is:

Scheme 1 Synthesis of 1,3,5-triazepan-2,6-diones 1.

The text above the arrow in Scheme 1 should read: refs. 14,15

In the second sentence of paragraph three of the main text, which begins 'Preliminary X-ray structure analysis of representative members $(R^2 \neq H)$ revealed...', R^2 should read R^3 .

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

Additions and corrections can be viewed online by accessing the original article to which they apply.