

https://doi.org/10.1038/s41467-019-08920-9

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## Author Correction: A shape-shifting redox foldase contributes to *Proteus mirabilis* copper resistance

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Correction to: Nature Communications; https://doi.org/10.1038/ncomms16065; published online 19 July 2017

This Article contains errors in Fig. 1, Table 1 and the Methods section. In panel c, the labels for PmScsC and EcDsbC in the upper two curves are interchanged. In Table 1 and the Methods section entitled 'Extended structure', the space group of the extended PmScsC structure is incorrectly referred to as H3<sub>2</sub> and should read H32. Correct versions of Fig. 1 and Table 1 are presented below; the errors have not been corrected in the Article.

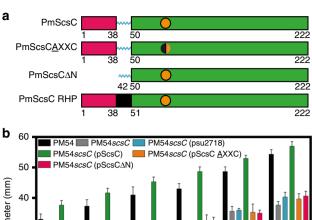
Published online: 01 March 2019

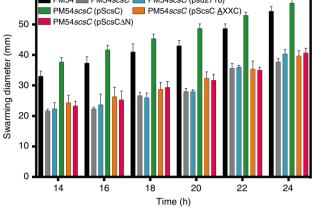
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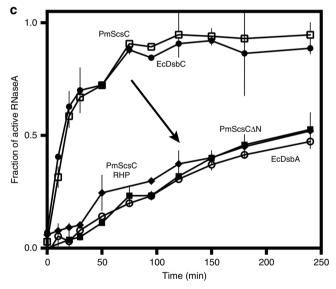


Fig. 1

Table 1 PmScsC crystal structure statistics			
	Compact (4XVW)	Transitional (5IDR)	Extended (5ID4)
Data collection			
Space group	P2 <sub>1</sub>	14	H32
Cell dimensions			
a, b, c (Å)	137.5, 163.9,	193.1, 193.1,	86.7, 86.7,
	181.9	105.8	330.9
$\alpha, \beta, \gamma$ ( $^{\circ}$ )	90, 90, 90	90, 90, 90	90, 90, 120
Resolution (Å)	91.15-2.60	136.51-2.56	110.29-2.92
	(2.74-2.60)	(2.57-2.56)	(2.93-2.92)
R <sub>merge</sub>	0.072 (0.617)		
1 /σΙ	11.0 (2.0)	14.9 (2.2)	14.2 (2.8)
Completeness	98.6 (95.4)	99.4 (100.0)	99.2 (100.0)
(%)			
Redundancy	3.8 (3.7)	4.1 (4.1)	4.1 (4.2)
Refinement	0445 0 40	40.00.054	40.04.000
Resolution (Å)	91.15-2.60	42.82-2.56	40.36-2.92
No. of reflections	-,	62,069	10,652
$R_{\text{work}}/R_{\text{free}}$ (%)	24.8/28.2	17.1/22.2	25.1/26.3
No. of atoms	40.050	10.262	1720
Protein	40,850	10,262	1720
Ligand/ion	NA	NA	NA
Water	281	82	0
B factors (Å <sup>2</sup> )	F0.7	FO (	122.2
Protein	59.7	50.6	122.2
Ligand/ion	NA 41.F	NA 43.0	NA
Water RMS deviations	41.5	43.0	NA
Bond length (Å)	0.006	0.008	0.010
Bond length (A) Bond angles (º)	1.21	1.05	1.17
bonu angles (=)	1.21	1.05	1.17

Single crystals were used to collect each dataset. Values for highest resolution shell are shown in parentheses