See discussions, stats, and author profiles for this publication at: https://www.researchgate.net/publication/51752924

Erratum to: Pivotal role of the strictly conserved aromatic residue F15 in the cytochrome c 7 family

ARTICLE in EUROPEAN JOURNAL OF BIOCHEMISTRY · OCTOBER 2011

Impact Factor: 2.54 · DOI: 10.1007/s00775-011-0854-z · Source: PubMed

READS

17

8 AUTHORS, INCLUDING:



Yuri Y Londer

New England Biolabs

38 PUBLICATIONS 745 CITATIONS

SEE PROFILE



Phani Raj Pokkuluri

Argonne National Laboratory

75 PUBLICATIONS 1,049 CITATIONS

SEE PROFILE



Ricardo Louro

New University of Lisbon

87 PUBLICATIONS 1,093 CITATIONS

SEE PROFILE



Carlos A Salgueiro

New University of Lisbon

80 PUBLICATIONS 1,195 CITATIONS

SEE PROFILE

ERRATUM

Erratum to: Pivotal role of the strictly conserved aromatic residue F15 in the cytochrome c_7 family

Joana M. Dantas · Leonor Morgado · Yuri Y. Londer · Ana P. Fernandes · Ricardo O. Louro · P. Raj Pokkuluri · Marianne Schiffer · Carlos A. Salgueiro

Published online: 29 October 2011

© SBIC 2011

Erratum to: J Biol Inorg Chem DOI 10.1007/s00775-011-0821-8

Unfortunately the original version has been published with errors in one figure.

The corrected figure is given here (Fig. 9).

The online version of the original article can be found under doi:10.1007/s00775-011-0821-8.

J. M. Dantas · L. Morgado · A. P. Fernandes · C. A. Salgueiro (☒)
Requimte-CQFB, Departamento de Química,
Faculdade de Ciências e Tecnologia,
Universidade Nova de Lisboa,
Campus Caparica, 2829-516 Caparica, Portugal
e-mail: csalgueiro@dq.fct.unl.pt

Y. Y. Londer · P. R. Pokkuluri · M. Schiffer Biosciences Division, Argonne National Laboratory, Argonne, IL 60439, USA

Present Address:
Y. Y. Londer
New England Biolabs,
240 County Road, Ipswich, MA 01938, USA

R. O. Louro Instituto de Tecnologia Química e Biológica, Universidade Nova de Lisboa, Av. da Republica (EAN), 2780-157 Oeiras, Portugal



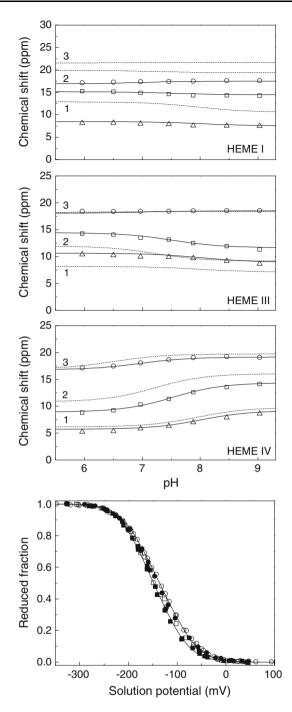


Fig. 9 Fitting of the thermodynamic model to the experimental data for PpcAF15L. The *solid lines* are the result of the simultaneous fitting of the NMR and visible data. The *three upper panels* show the pH dependence of heme methyl chemical shifts at oxidation stages 1 (*triangles*), 2 (*squares*), and 3 (*circles*). The chemical shifts of the heme methyls in the fully reduced stage (stage 0) are not plotted since they are unaffected by the pH. The *dashed lines* in each panel represent the best fit for the wild-type protein and the nearest label (1, 2, 3) indicates the oxidation stage represented by the *curve*. The *bottom panel* corresponds to the reduced fractions of PpcAF15L determined by visible spectroscopy at pH 7.0 (*circles*) and pH 8.0 (*squares*). The *open symbols* and the *filled symbols* represent the data points in the reductive and oxidative titrations, respectively

