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Erratum

Erratum to "Inorganic perchlorato complexes" [Coordination Chemistry Reviews 178–180 (1998) 865–902]^{⋄,⋄,⋄}

Jean-Louis Pascal *. Frédéric Favier

Laboratoire des Agrégats Moléculaires et Matériaux Inorganiques, ESA 5072, Université Montpellier II, Place E. Bataillon, 34095 Montpellier, Cedex 5, France

The publisher regrets that the following errors occurred when this article was originally printed.

The footnote to the title, dedicating this paper to Jacqueline Potier, was omitted.

In Table 1, on page 869, the third compound should be NO₂ClO₄

In Table 2, on page 873, footnote (b) should read

"(b) NO_2ClO_4 and N_2O_5 were prepared according to literature methods [40,41]."

In Table 4, on page 875, footnote (e) should read

"(e) the reaction was complex and the products difficult to isolate as pure compounds [31]."

E-mail address: pasfav@univ-montp2.fr (J.-L. Pascal)

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[★] PII of original article S 0 0 1 0 - 8 5 4 5 (9 8) 0 0 1 0 2 - 7.

This paper is dedicated to Jacqueline Potier.

^{*} Corresponding author. Tel.: +33-4-67-14-33-32; fax: +33-4-67-14-3304.

In Table 5, on page 877, the headings " PO_4^2 " and " CO_3^- " should read " PO_4^3 " and " CO_3^2 ", respectively.

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In Table 5, on page 877, footnotes a-e should read. "a Ref. [165]; b Ref. [19]; Ref. [166]; d Ref. [167]; Ref. [23];
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In Table 5, on page 877, the last two sentences under the table should read

"Phosphates and sulphates polymerize [103]. However, triflates $[SO_3CF_3]$ are easily replaced [by CIO_4] [61]."

In Table 6, on page 878, the references should read.

References: H: [43,52,53]; Be: [71,72]; Mg: [73]; Sc: [74]; Y: [75]; La [75]; Ce: [76–78]; Pr: [79], [80](α); Nd: [75], [81](β); Sm: [32,75]; Eu: [32, 75]; Gd: [75]; Tb: (J.L. Pascal and F. Favier, unpublished result); Dy: [75]; Ho: [32,75]; Er: [32,75]; Tm: [32,75]; Yb: [82,158]; Lu: [32]; Th: (J.L. Pascal and F. Favier, unpublished results); U: [83,84]; Ti: [16,62,65,85,86], [87](χ), 88; Zr: [89–91]; Hf: [92,93,175]; V: [6,31]; Nb: [94]; Ta: [95]; Cr: [16,58,96,97]; Mo; [46,98]; W: [99,100]; Mn: [31,101]; Fe: [61,97,102,103]; Co: [23,31,37,104–107]; Rh: [108]; Ni: [23,26,31,107,109]; Pd: [110]; Cu: [2,23,31,42,111,112]; Zn: [31,113,114]; Cd: [62,115–117]; Hg: [62,115,118]; B: [38,59,119–123], [124](ε); Al: [60,119,122,125–134]; Ga: [46,135–137]; In: [137,138]; Tl: [138]; Si: [6,141], [142](φ)]; Ge: [62,159]; Sn: [62,97,143,144], [145](γ); As: (R.J. Gillespie et al., unpublished results) (η); Sb: [146](ι), [147] Bi: [148,149]; Te: [61]; F: [48,150]; Cl: [45–47], (G. Schrobilgen et al., unpublished results) (φ), [56], [150–157](κ); Br: [47,154,155]; I: [156]; Xe: [157](λ); C: [139,140,176](μ).

On page 880, the first sentence in the legend to Fig. 2 should read

"Molecular structure of $Sn(ClO_4)_6^{2-}$ anion from Ref. [144]"

On page 880, the first sentence in the legend to Fig. 3 should read

"Molecular structure of Sb₂Cl₆(O)(OH)ClO₄ from Ref. [146]"

On page 881, the first sentence in the legend to Fig. 4 should read

"Molecular structure of dimeric Sn₃O₂Cl₄(ClO₄)₄ from Ref. [143]"

On page 882, the legend to Fig. 5 should read

"Molecular structure of $Ti(ClO_4)_4$ from Ref. [65]. Mean distances and angles: $Ti-O_b=2.07$ Å; $Cl-O_t=1.39$ Å; $Cl-O_b=1.51$ Å; $O_b-Cl-O_b=97^{\circ}$ $O_t-Cl-O_t=115^{\circ}$ "

On page 883, the first sentence in the legend to Fig. 6 should read

"Coordination shell around Ni and Co in $Ni(ClO_4)_2$ and $Co(ClO_4)_2$ from Ref. [107]."

On page 883, the first sentence in the legend to Fig. 7 should read

"Molecular structure of Eu(ClO₄)₃ displaying the Eu coordination shell, a slightly distorted tricapped trigonal prism, and bridging tridentate [ClO₄) groups [32]."

On page 884, the first sentence in the legend of Fig. 8 should read

"Molecular structure of Yb(ClO₄)₃ · H₂O from Ref. [158]."

On page 885, the first sentence in the legend of Fig. 9 should read

"Molecular structure of Lu(ClO₄)₃, from Ref. [32]."

On page 886, the first sentence in the legend of Fig. 10 should read

"One set of the disordered atoms in Cu(ClO₄)₂ from Ref. [112]."

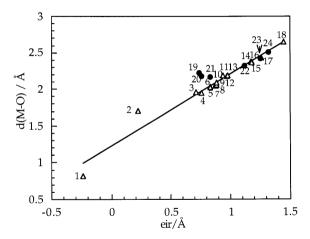


Fig. 11. Plot of M–O distances related to effective ionic radii (eir) [160]. 1—HClO₄, 2—Cl₂O₇, 3—Cu(ClO₄)₂, 4—Ga(ClO₄)₃, 5—Sn(ClO₄)₆²–, 6—Ni(ClO₄)₂, 7—Zn(ClO₄)₂, 8—Co(ClO₄)₃⁻, 9—Ti(ClO₄)₄, 10—Co(ClO₄)₂, 11—In(ClO₄)₃, 12—Hf(ClO₄)₄, 13—Zr(ClO₄)₄, 14—Lu(ClO₄)₃, 15—Yb(ClO₄)₃ (LT), 16—Yb(ClO₄)₃ (HT), 17—Eu(ClO₄)₃, 18—Bi₂(ClO₄)₁₀⁴–, 19—Sb₂Cl₆O(OH)ClO₄, 20—Sn₃O₂Cl₄(ClO₄)₄ (five-coordination around Sn), 21—Sn₃O₂Cl₄(ClO₄)₄ (six-coordination around Sn), 22—Yb(ClO₄)₃·H₂O, 23—Nd₂(OH)₃(ClO₄)₃·5H₂O, 24—Pr₂(OH)₃H₂O(ClO₄)₃. The curve fit, dM—O = 1.23 + 0.98 eir, was calculated from complexes including exclusively ClO₄ as a ligand (open triangles). Shaded circles correspond to complexes incorporating more basic ligands than ClO₄: O, OH, H₂O, and it is clearly show that with small metallic eir, [ClO₄] is scattered out the coordination sphere while with greater eir, [ClO₄] draws nearer.

On page 887, the first sentence in the legend of Fig. 11 should read

"Plot of M-O distances related to effective ionic radii (eir) [160]."

On page 887, one of the points was missing of the plot of Fig. 11. The corrected Fig. is printed above.

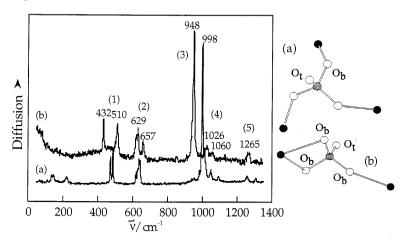
On page 889, the last part of the legend to Fig. 12 should read

"(7) $v_{as}ClO_4^-$, $HgCl^+$ was assumed to be polymeric cation $(HgCl)_n^{n+}$ [115]".

On page 891, the second sentence in the legend of Fig. 15 should read

"Bridging bidentate $[ClO_4]$ involved in a layered bidimentional polymeric unit (c.f. $Ga(ClO_4)_3$ [137])."

On page 893, the frequency spectrum part of Fig. 18 was incorrect, the correct figure is printed below.



On pages 899 and 901, the following references should read

"[32] J.L. Pascal, F. Favier, F. Cunin, A.N. Fitch, G. Vaughan, J. Solid State Chem. 139 (1998) 259."

"[36] A.C. Pavia, C.R. Acad. Sci. Paris, Serie II 261 (1965) 5118."

"[47] K.O. Christie, C.J. Shack, E.C. Curtis, Inorg. Chem. 10 (1971) 1589."

"[57] K.M. Tobias, M. Jansen, Z. Anorg. Allg. Chem. 550 (1987) 16."

"[66] R.G. Pearson, J. Am. Chem. Soc. 85 (1963) 3533."

"[68] D. Datta, Inorg. Chem. 31 (1992) 2797.

"[156] K.O. Christie, C.J. Shack, Inorg. Chem. 11 (1972) 1682."