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Erratum

Erratum to: "Synthesis and structural characterization of $[H(OEt_2)_2]^+[(C_3H_3N_2)\{B(C_6F_5)_3\}_2]^-$—a Brønsted acid with an imidazole-derived 'non-coordinating' anion" [Journal of Organometallic Chemistry, 641 (2002) 148–155]☆

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The publisher apologises for the error in the above-mentioned article. Section 3.6.1 of the published article should read:

3.6.1. X-ray crystal structure analysis of ***9***

Formula $C_{39}H_3N_2B_2F_{30}{*}H(C_4H_{10}O)_2$, $M = 1240.30$, colourless $0.30 \times 0.20 \times 0.15$ mm, $a = 12.848(3)$, $b = 13.603(3)$, $c = 14.144(4)$ Å, $\alpha = 81.74(1)$, $\beta = 89.79(1)$, $\gamma = 76.98(1)°$, $V = 2382.4(10)$ Å^3, $\rho_{calc} = 1.729$ g cm^{-3}, $\mu = 1.86$ cm^{-1}, empirical absorption correction via SORTAV ($0.946 \leq T \leq 0.973$), $Z = 2$, triclinic, space group $P\bar{1}$bar (No. 2), $\lambda = 0.71073$ Å, $T = 198$ K, ω and φ scans, 9927 reflections collected ($\pm h$, $\pm k$, $\pm l$), [(sin θ)/λ] = 0.62 Å^{-1}, 6203 independent ($R_{int} = 0.034$) and 3820 observed reflections [$I \geq 2\sigma(I)$], 756 refined parameters, $R = 0.071$, $wR^2 = 0.126$, max. residual electron density 0.22 (−0.23) e Å^{-3}, thermal parameters of the ether group O71–C75 indicate some disorder, not refined due to small amount of observed data from the poorly diffracting crystal, H1 between the two ether molecules from difference Fourier calculations, refined isotropically, other hydrogens calculated and refined as riding atoms.

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