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Erratum: Reaction of (*E*)-Bis(2,4,6-tri-*tert*-butylphenyl)diphosphene with Tetrachloro-*o*-benzoquinone

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Legends for Figures 1 through 4 were incorrect as published. For clarity, the figures and corrected legends appear below:

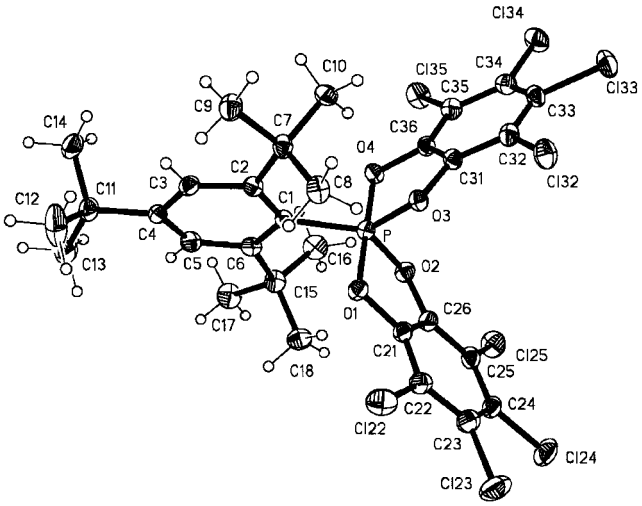

FIGURE 1 A thermal ellipsoid drawing of **2** with 50% probability levels. Selected bond lengths (pm) and angles (°): P–O1, 170.9(2); P–O2, 165.1(2); P–O3, 164.8(2); P–O4, 179.6(2); P–C1, 185.2(3); O1–P–O2, 90.81(11); O1–P–O3, 88.62(11); O1–P–O4, 179.19(10); O2–P–O3, 102.65(10); O2–P–O4, 89.24(11); O3–P–O4, 90.58(11); C1–P–O1, 89.25(11); C1–P–O2, 128.18(11); C1–P–O3, 129.15(12); C1–P–O4, 91.36(11).

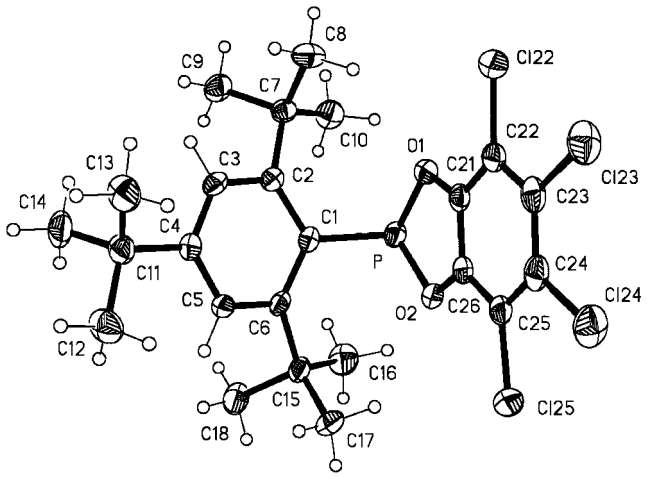

FIGURE 2 A thermal ellipsoid drawing of one of the two independent molecules of **3** with 50% probability levels. Selected bond lengths (pm) and angles (°): P–O2, 170.05 (15); P–O1, 170.20(16); P–C1, 185.0(2); O2–P–O1, 91.83(7); O2–P–C1, 99.45(8); O1–P–C1, 99.51(8).

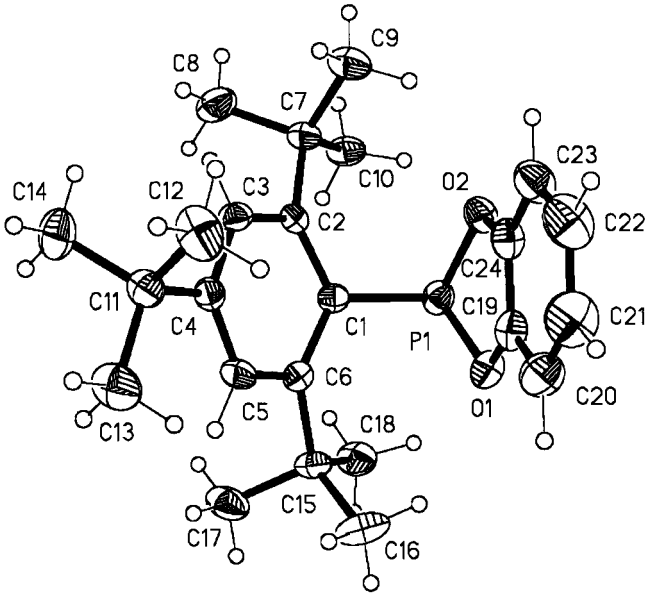

FIGURE 3 A thermal ellipsoid drawing of one of the four independent molecules of **6** with 50% probability levels. The second position of the *p-tert*-butyl group is omitted for clarity. Selected bond lengths (pm) and angles (°): P–O2, 167.81 (15); P–O1, 168.16(14); P–C1, 186.02(19); O2–P–O1, 92.03(7); O2–P–C1, 97.82(8); O1–P–C1, 97.47(8).

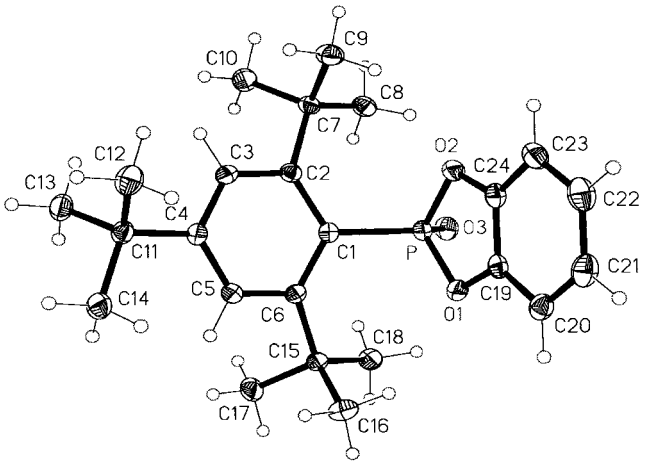

FIGURE 4 A thermal ellipsoid drawing of one of the two independent molecules of **7** with 50% probability levels. Selected bond lengths (pm) and angles (°): P–O3, 145.52 (8); P–O2, 163.72(8); P–O1, 162.66(9); P–C1, 180.71(10); O3–P–O1, 114.25(5); O3–P–O2, 114.24(4); O1–P–O2, 95.08(4); O3–P–C1, 121.25(5); O1–P–C1, 104.57(5); O2–P–C1, 103.81(4).

Additionally, reference 20 should read:

[20] Perrin, D. D.; Armarego, W. L. F. Purification of Laboratory Chemicals, 3rd ed.; Pergamon Press: Oxford, New York, 1988.