Erratum: Comment on "Analytic value of the atomic three-electron correlation integral with Slater wave functions" [Phys. Rev. A 68, 016501 (2003)]

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In our Comment on Remiddi's ''Analytic value of the atomic three-electron correlation integral with Slater wave functions'' [Phys. Rev. A **68**, 016501 (2003)] the last eight digits in the Remiddi column of Table I are in error. The corrected table is given below.

TABLE I. $\mathbf{Z}(w_1,w_2,w_3)$ results for various w_1, w_2, and w_3.

w_1	w_2	w_3	Remiddi
1.50	1.00	1.00	8.783505801548186149325724126 64
1.50	1.00	1.50	2.539564605260415333027454370 04
1.50	1.00	2.00	1.053910366507817767776457486 52
1.50	1.50	1.00	2.539564605260415333027454370 04
1.50	2.00	1.50	3.511260150977273542310199408 89
2.00	1.00	1.00	2.986672982272132343817654866 58
2.00	2.00	1.00	3.243503012860346267583161790 89

