Erratum: Assessing the accuracy of the van der Waals density functionals for rare-gas and small molecular systems [Phys. Rev. B **91, 195103 (2015)]**

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The binding energies of the Kr crystal obtained using different the van der Waals density functionals presented in the original article are erroneous and Table III should be corrected as follows.

Method	Ne		Ar		Kr	
	a_0	E_b	a_0	E_b	a_0	E_b
vdW-DF2	4.18	56	5.26	142	5.70	164
optB88-vdW	4.25	45	5.22	143	5.61	180
optB86b-vdW	4.34	60	5.30	131	5.67	183
rev-vdW-DF2	4.43	30	5.35	90	5.71	120
CCSD(T) ^a	4.314	26.4	5.284	82.8	5.670	114
RPA@LDA ^b	4.7	11	5.4	59	5.8	88
RPA@PBE ^b	4.5	17	5.3	83	5.7	112
Expt. ^a	4.35	27.2	5.23	88.9	5.61	122.5

^aReference [58].^bReference [60].

The corrected binding energies of the Kr crystal are now consistent with the interaction energy curves shown in Fig. 3(c). The conclusion of the paper is unaffected by this change.

I.H. thanks Fabien Tran for pointing out the error.