

Electronic Supplementary Information

CoBr₂-doping-induced efficiency improvement of planar CsPbBr₃ perovskite solar cells

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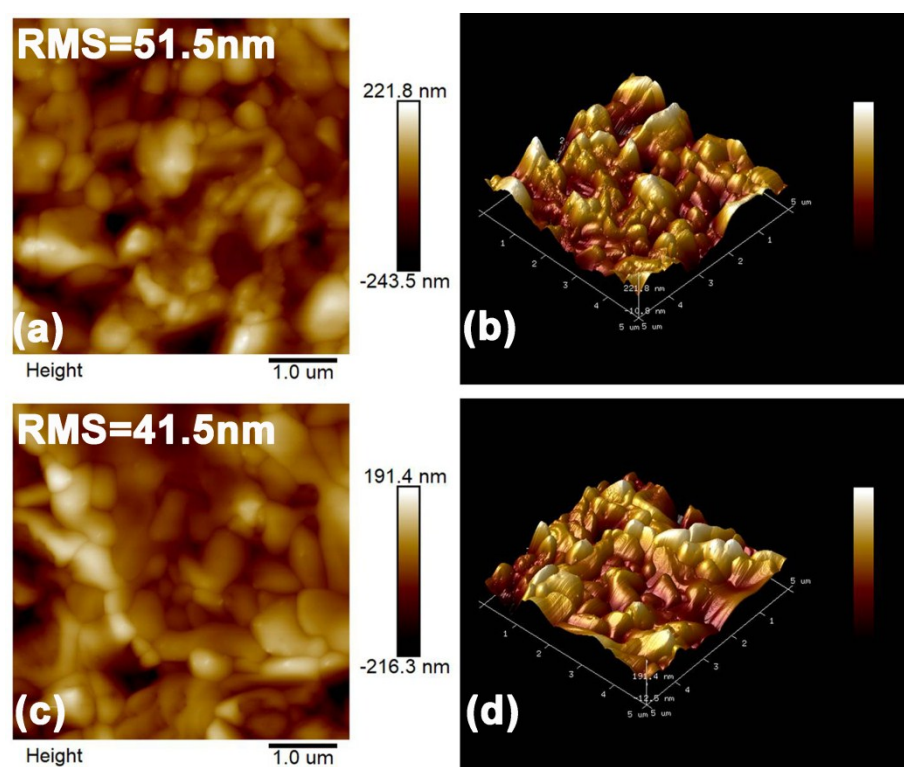


Figure S1. AFM images of CsPbBr₃ (a and b) and CsPb_{0.998}Co_{0.002}Br₃ (c and d) films.

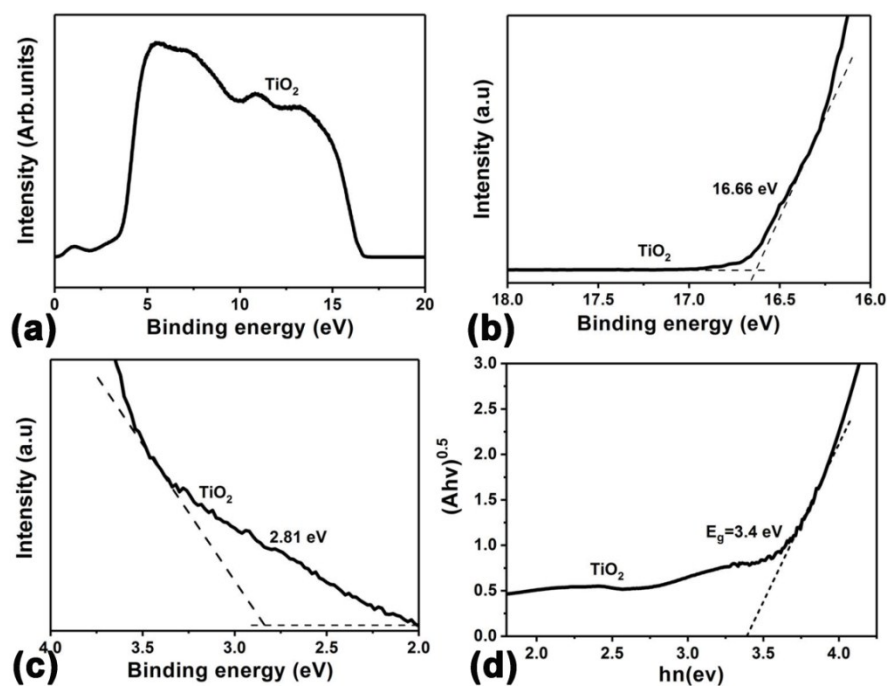


Figure S2. UPS spectrum (a), the cut-off energy ($E_{\text{cut-off}}$) region (b), the Fermi edge ($E_{\text{F, edge}}$) region (c), and Tauc plots (d) of TiO_2 ETL.

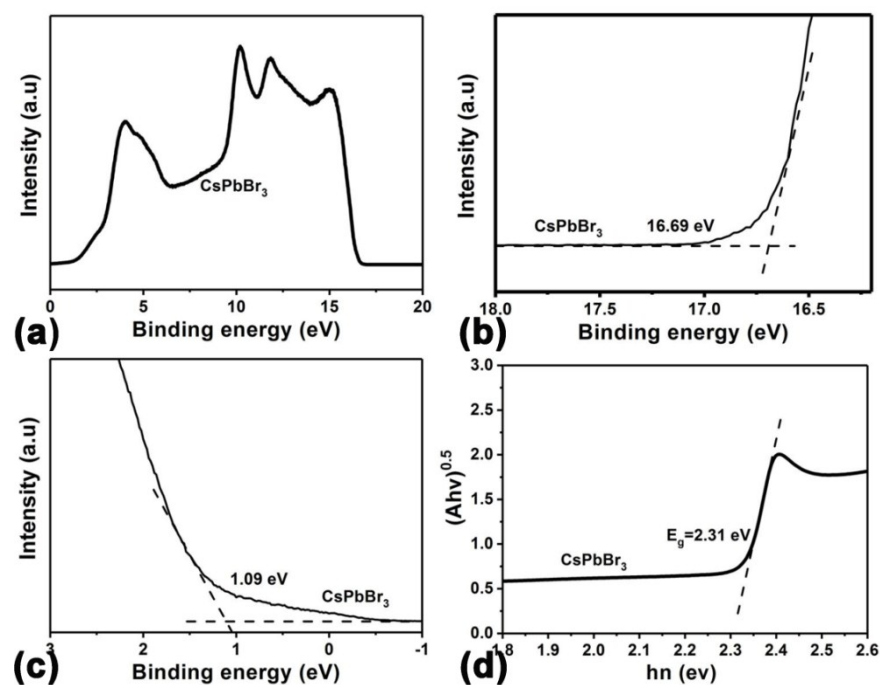


Figure S3. UPS spectrum (a), the cut-off energy ($E_{\text{cut-off}}$) region (b), the Fermi edge ($E_{\text{F, edge}}$) region (c), and Tauc plots (d) of CsPbBr_3 film.

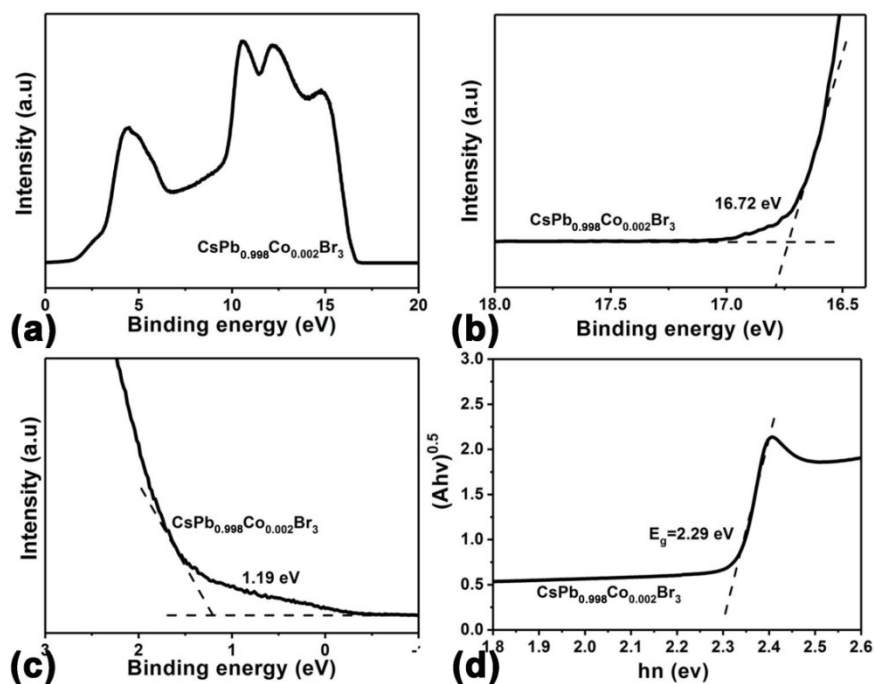


Figure S4. UPS spectrum (a), the cut-off energy ($E_{\text{cut-off}}$) region (b), the Fermi edge ($E_{\text{F, edge}}$) region (c), and Tauc plots (d) of CsPb_{0.998}Co_{0.002}Br₃ film.

Table S1. Energy levels of the TiO₂ ETL, CsPbBr₃ and CsPb_{0.998}Co_{0.002}Br₃ layers calculated from the UPS and Tauc plots data*.

Functional layer	$E_{\text{cut-off}}$ (eV)	$E_{\text{F, edge}}$ (eV)	E_{g} (eV)	E_{F} (eV)	E_{VB} (eV)	E_{CB} (eV)
TiO ₂	16.66	2.81	3.40	4.56	-7.35	-3.95
CsPbBr ₃	16.69	1.09	2.31	4.53	-5.62	-3.31
CsPb _{0.998} Co _{0.002} Br ₃	16.72	1.19	2.29	4.50	-5.69	-3.40

*The Fermi levels (E_{F}) is calculated from the equation of $E_{\text{F}} = E_{\text{cut-off}}$ (cut-off binding energy) - 21.2 eV (emission energy from He irradiation); the valence band maximum (E_{VB}) is calculated from the equation of $E_{\text{VB}} = E_{\text{F}} - E_{\text{F, edge}}$; the energy bandgap (E_{g}) is obtained from the Tauc plots; the E_{CB} is determined from E_{g} and E_{VB} [S1].

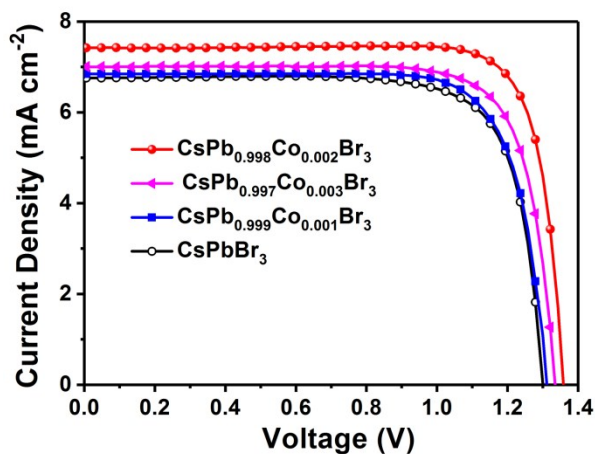


Figure S5. J-V curves of the CsPbBr₃ and CoBr₂ doped CsPbBr₃ planar PSCs.

Table S2. Photovoltaic parameters of the CsPbBr₃ and CoBr₂ doped CsPbBr₃ planar PSCs.

Perovskite composition	V _{OC} /V	J _{SC} /mA cm ⁻²	FF/%	PCE/%
CsPbBr ₃	1.301	6.79	77.08	6.81
CsPb _{0.999} Co _{0.001} Br ₃	1.311	6.85	79.99	7.18
CsPb _{0.998} Co _{0.002} Br ₃	1.357	7.45	84.84	8.57
CsPb _{0.997} Co _{0.003} Br ₃	1.332	6.92	79.92	7.36

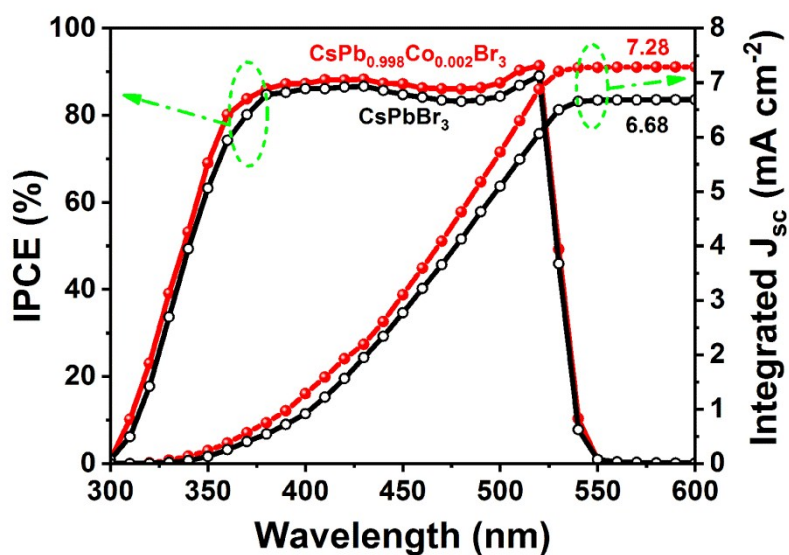


Figure S6. IPCE curves and integrated J_{sc} of the CsPbBr₃ and CsPb_{0.998}Co_{0.002}Br₃ planar PSCs.

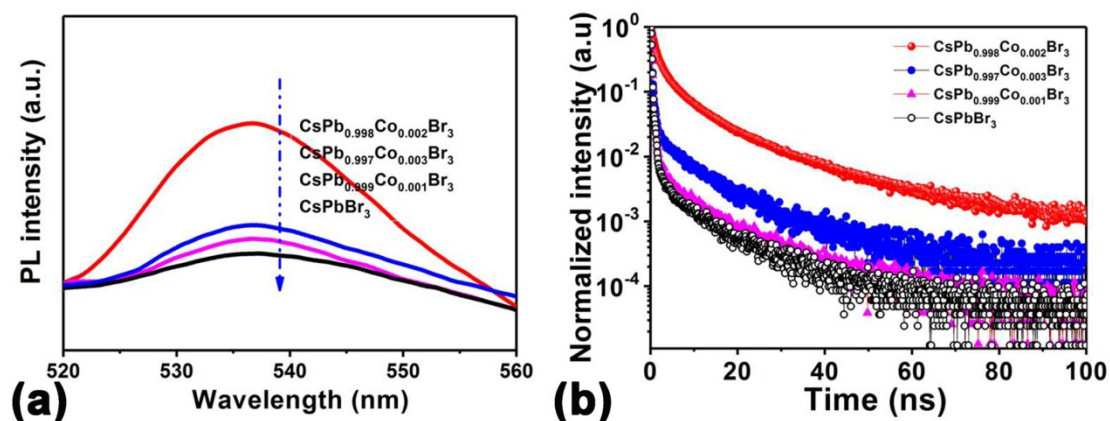


Figure S7. PL (a) and TRPL (b) spectra of the CsPbBr₃ and CoBr₂ doped CsPbBr₃ films.

Table S3. TRPL key parameters of the CsPbBr₃ and CoBr₂ doped CsPbBr₃ films.

Samples	τ_{ave} /ns	τ_1 /ns	τ_2 /ns	A_1 /%	A_2 /%
CsPbBr ₃	2.55	8.60	0.24	72.36	27.64
CsPb _{0.999} Co _{0.001} Br ₃	3.34	9.20	0.25	64.71	35.29
CsPb _{0.998} Co _{0.002} Br ₃	10.71	15.74	2.96	60.58	39.42
CsPb _{0.997} Co _{0.003} Br ₃	6.34	9.74	0.26	64.09	35.91

[S1] J. Song, W. N. Zhang, D. Wang, K. M. Deng, J. H. Wu, Z. Lan, Colloidal synthesis of Y-doped SnO₂ nanocrystals for efficient and slight hysteresis planar perovskite solar cells, *Solar Energy* 2019, 185, 508–515.