

## Supporting Information

# Perovskite templating via bathophenanthroline additive for efficient light-emitting devices

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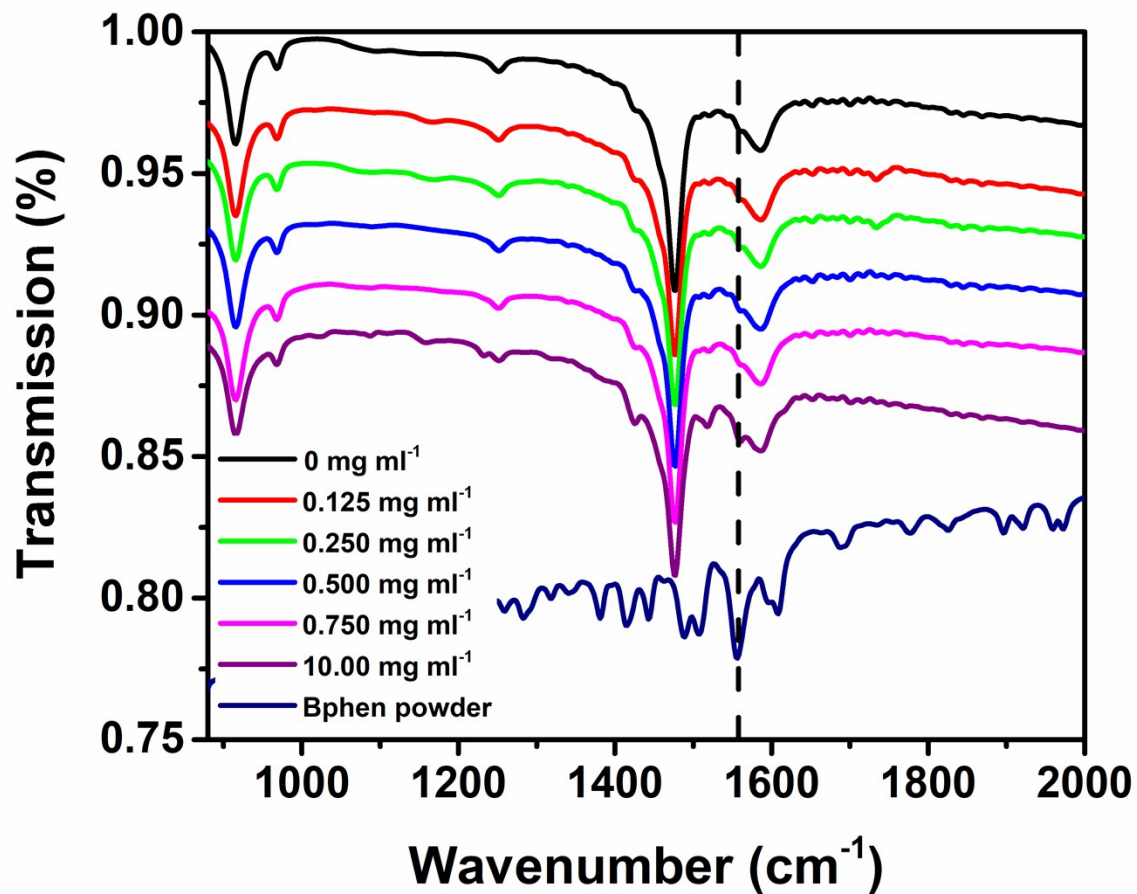
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## Additional Results



**Figure S1.** FTIR spectra of reference and BPhen-incorporated  $\text{CH}_3\text{NH}_3\text{PbBr}_3$  films as well as pure BPhen powder. Due to instrument limitation, the stretching due to the C=N bond ( $\sim 1550 \text{ cm}^{-1}$ ) was observable only in the sample prepared with  $10 \text{ mg mL}^{-1}$  BPhen concentration. The spectra are offset to provide a clearer view of the expected contribution due to C=N bond stretching.

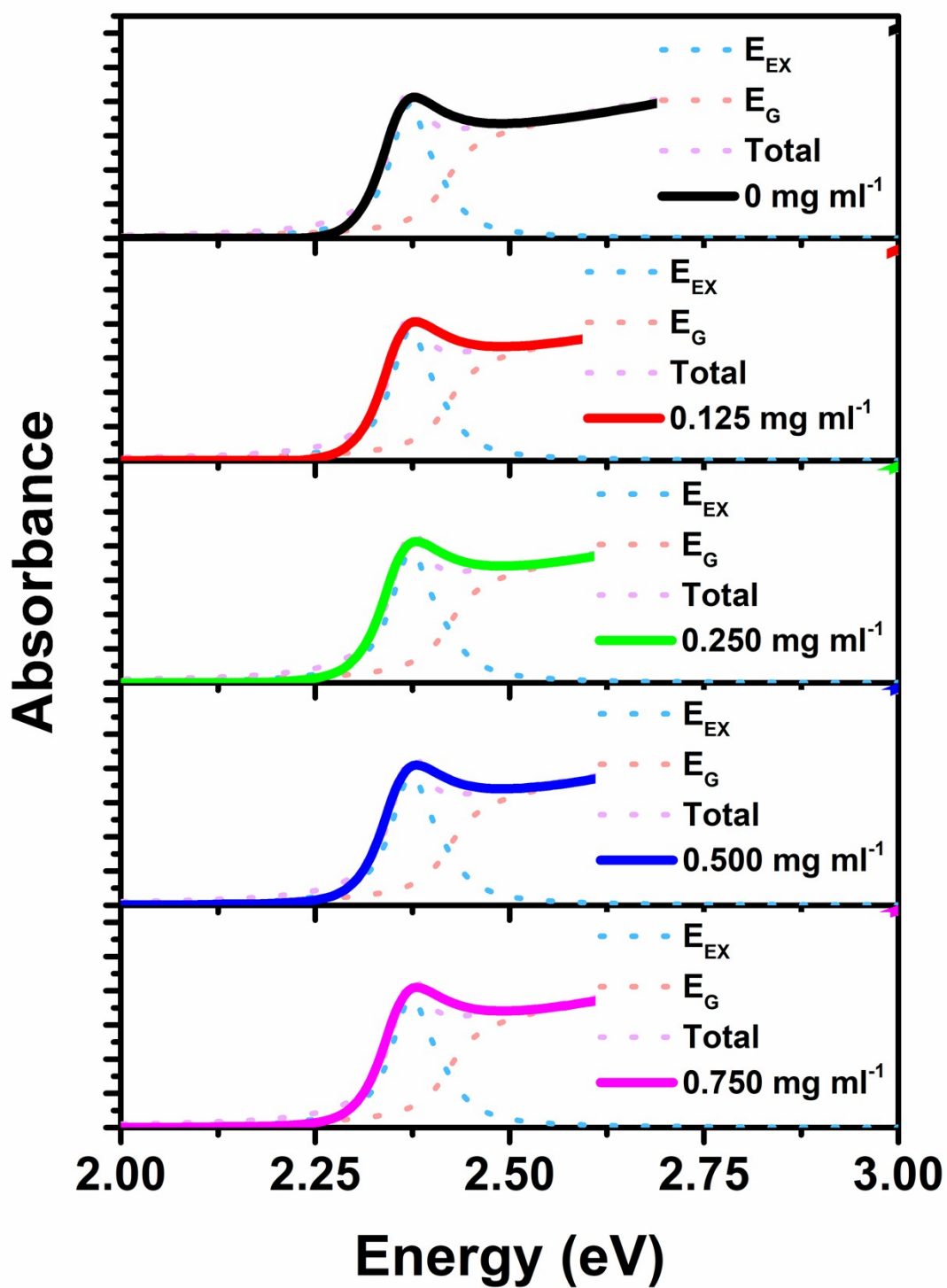


Figure S2. Elliot fitting of absorbance curves for reference and BPhen-incorporated  $\text{CH}_3\text{NH}_3\text{PbBr}_3$  films

**Table S1.** Exciton binding energy and optical band gap energy of reference and BPhen-incorporated  $\text{CH}_3\text{NH}_3\text{PbBr}_3$  films obtained by Elliot's fitting of the absorbance spectra in Figure S2.

Sample	Exciton Binding Energy ( $E_{\text{EX}}$ )	Band Gap ( $E_{\text{G}}$ )
0 mg ml <sup>-1</sup>	50 meV	2.4 eV
0.125 mg ml <sup>-1</sup>	50 meV	2.4 eV
0.250 mg ml <sup>-1</sup>	50 meV	2.4 eV
0.500 mg ml <sup>-1</sup>	50 meV	2.4 eV
0.750 mg ml <sup>-1</sup>	50 meV	2.4 eV

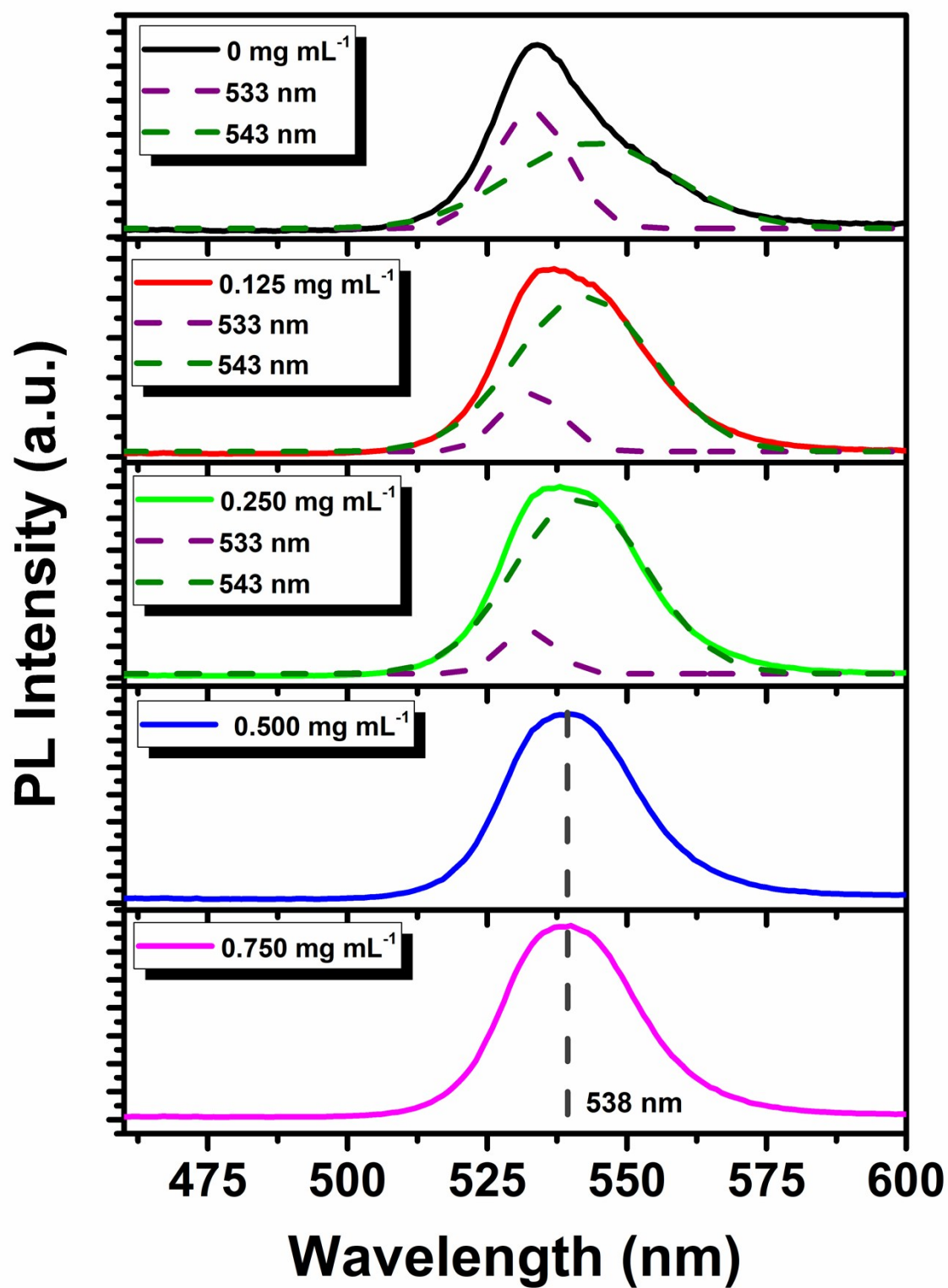
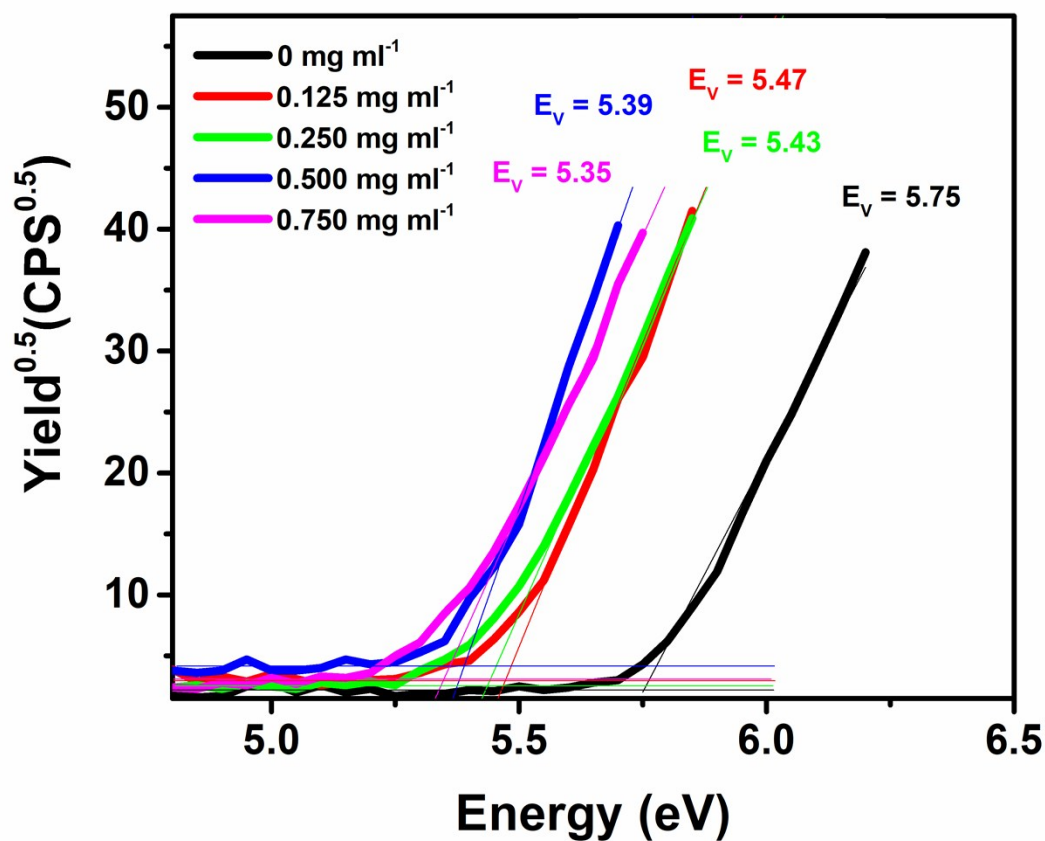


Figure S3: Peak fitting for PL spectra of reference and BPhen-incorporated  $\text{CH}_3\text{NH}_3\text{PbBr}_3$  film

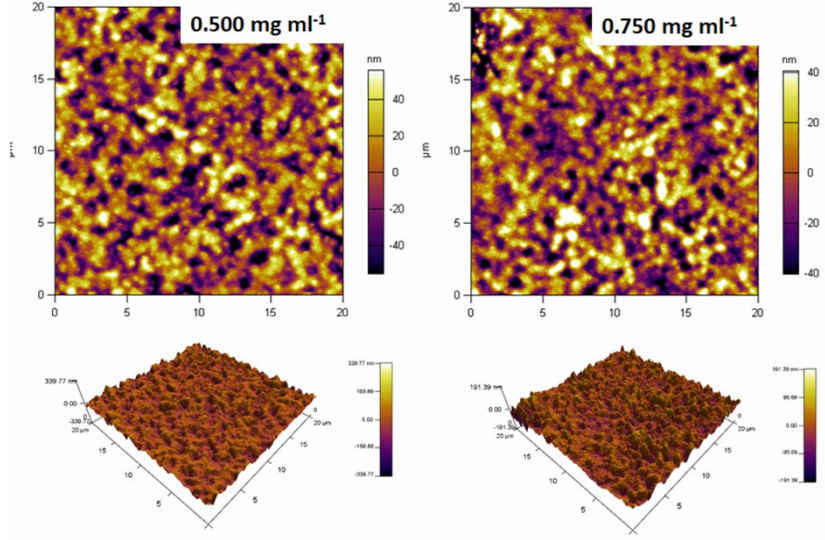
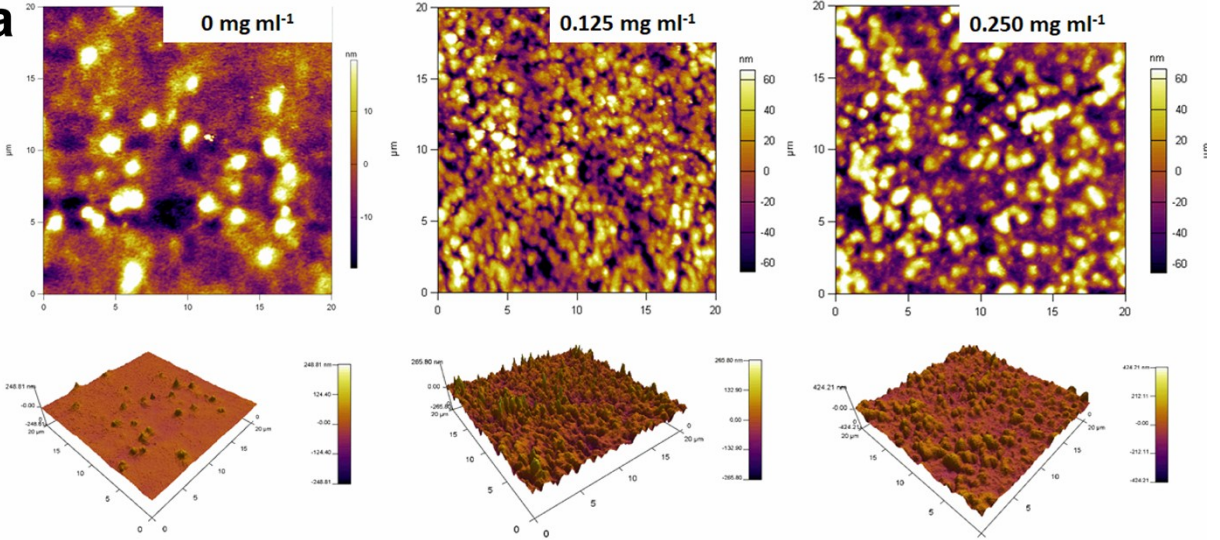


**Figure S4.** Valence band maximum estimation obtained using PESA measurements for reference and BPhen-incorporated  $\text{CH}_3\text{NH}_3\text{PbBr}_3$  films.

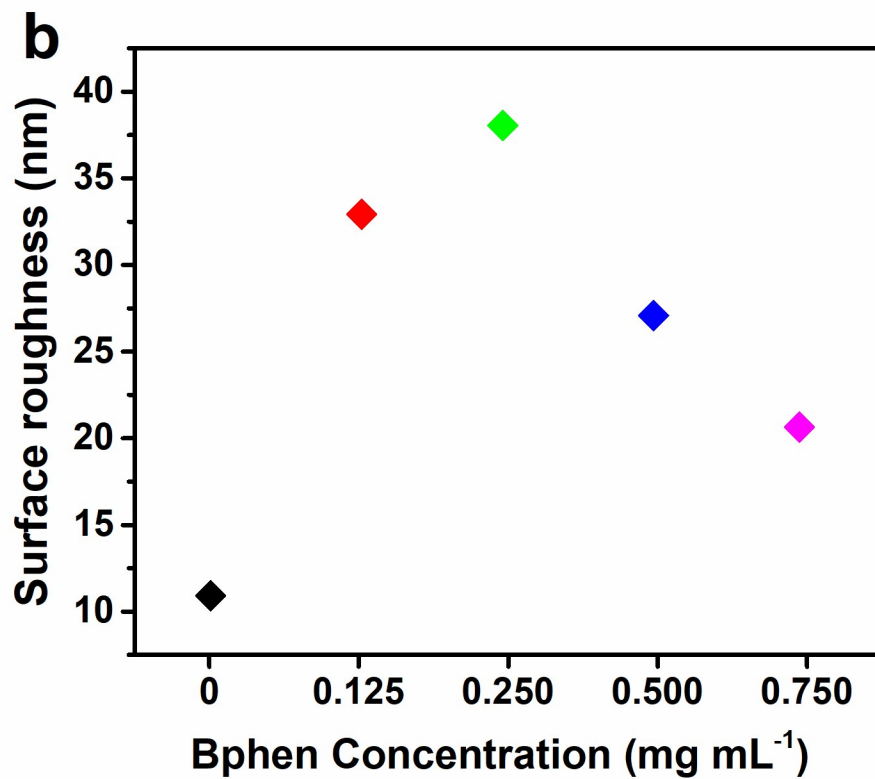
**Table S2.** Summary of decay lifetimes of reference and BPhen-incorporated  $\text{CH}_3\text{NH}_3\text{PbBr}_3$  films.

Sample	$a_1$	$A_1$	$T_1$	$a_2$	$A_2$	$T_2$	$T_{\text{ave}}$
<b>0 mg ml<sup>-1</sup></b>	0.258	0.2716	2.026	0.692	0.72842	18.426	13.9721
<b>0.125 mg ml<sup>-1</sup></b>	0.342	0.3548	3.112	0.622	0.64523	20.413	14.2751
<b>0.250 mg ml<sup>-1</sup></b>	0.389	0.3994	2.421	0.585	0.60062	14.693	9.79176
<b>0.500 mg ml<sup>-1</sup></b>	0.401	0.4199	2.016	0.554	0.5801	12.541	8.1216
<b>0.750 mg ml<sup>-1</sup></b>	0.456	0.4423	2.303	0.575	0.55771	13.812	8.7217

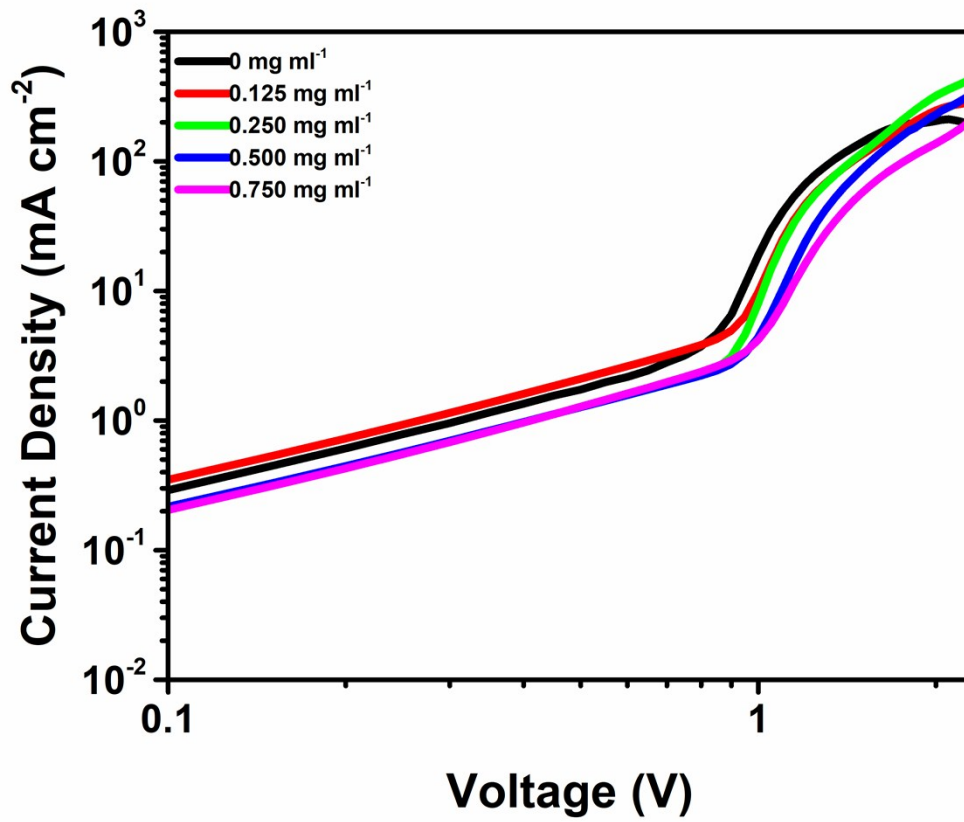
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**Figure S5.** (a) Surface topology taken using AFM and (b) estimated surface roughness of reference and BPhen-incorporated  $\text{CH}_3\text{NH}_3\text{PbBr}_3$  films.



**Figure S6.** Current density against voltage curves for reference and BPhen-incorporated CH<sub>3</sub>NH<sub>3</sub>PbBr<sub>3</sub> hole-only devices. The device architecture for the hole-only device consists of ITO/PEDOT: PSS/perovskite/Au.

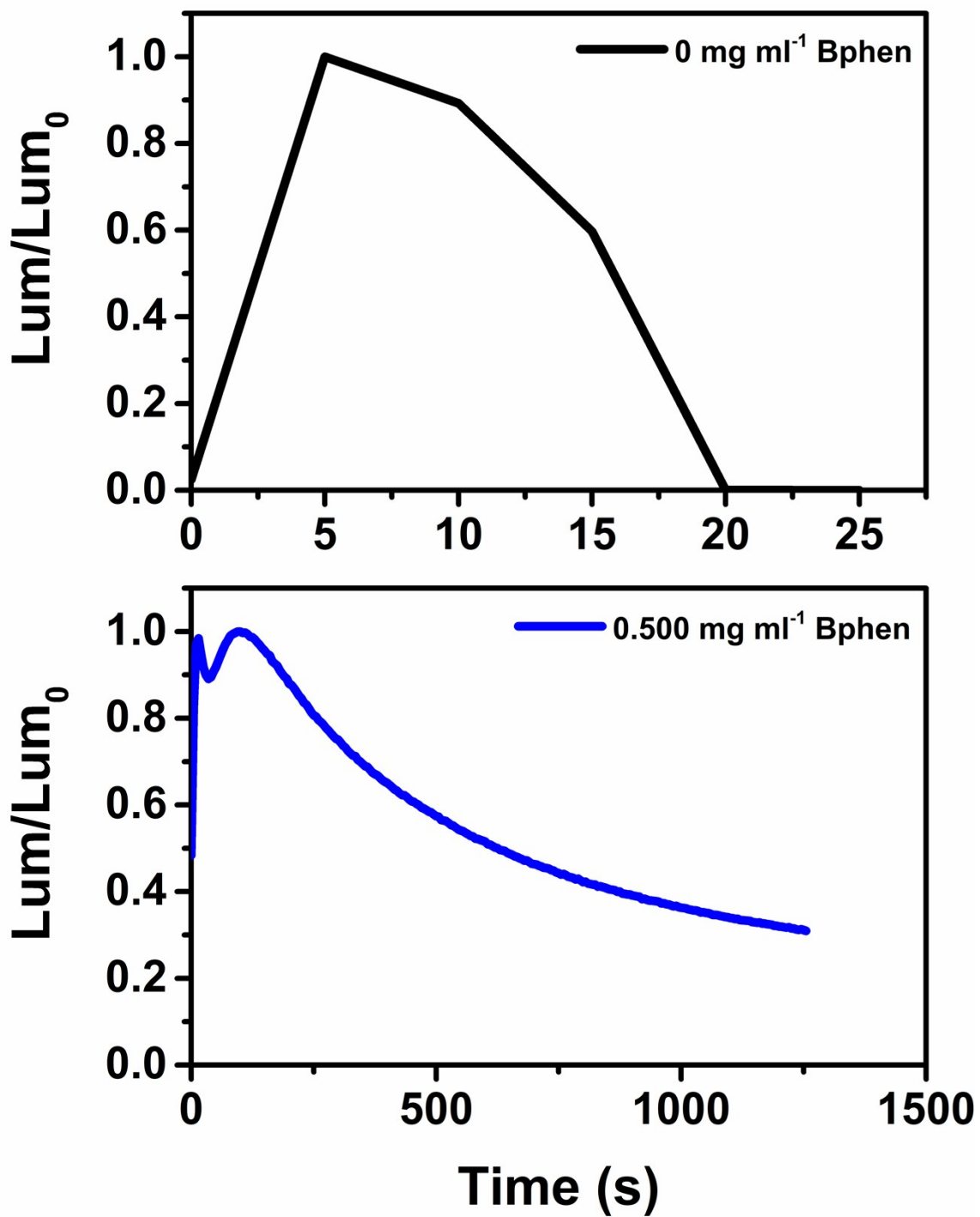
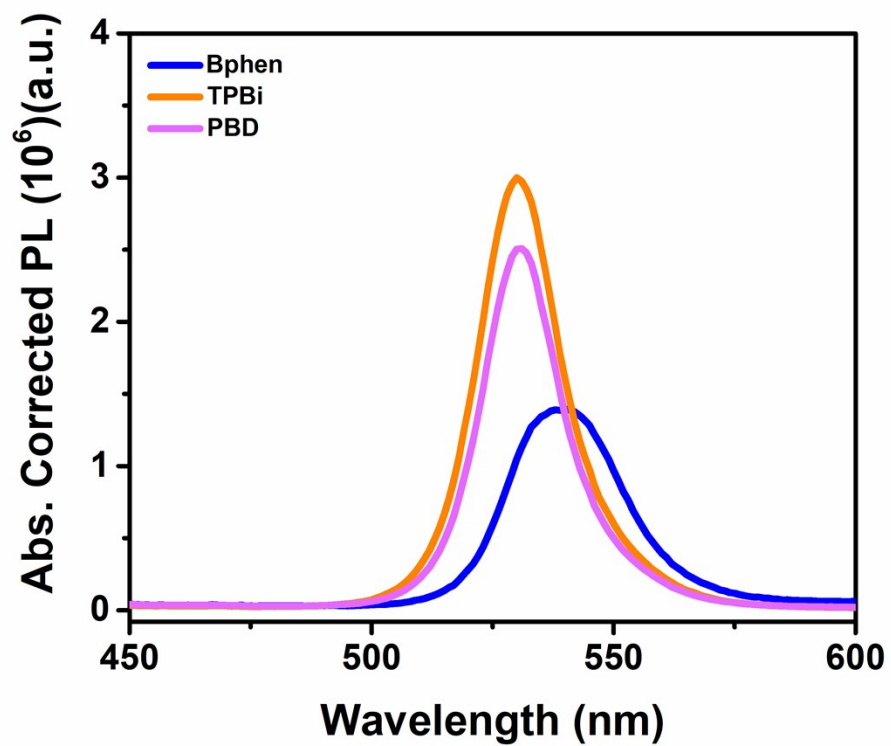
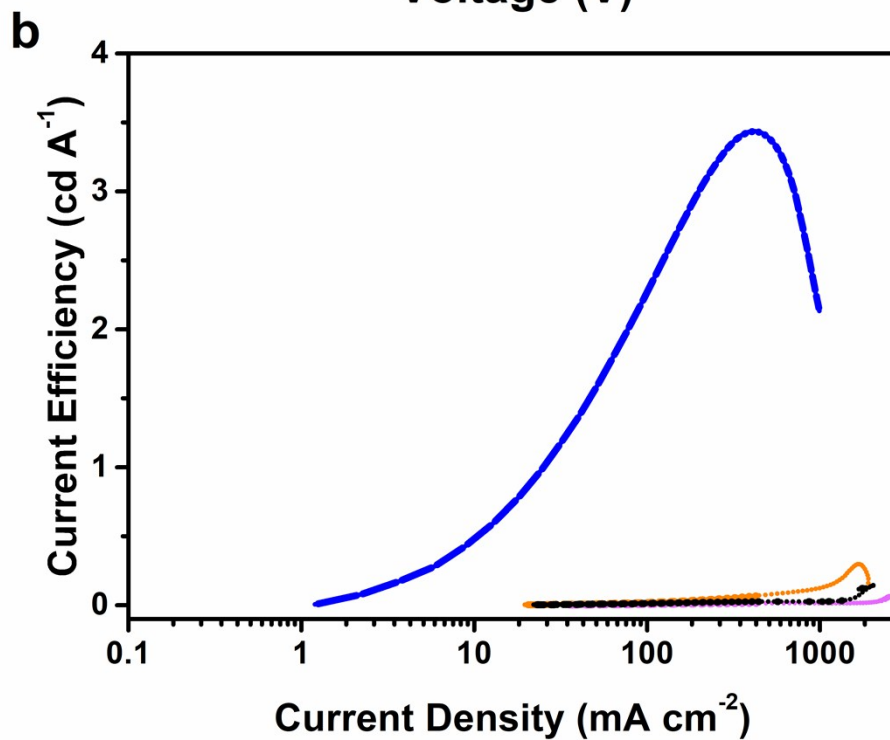
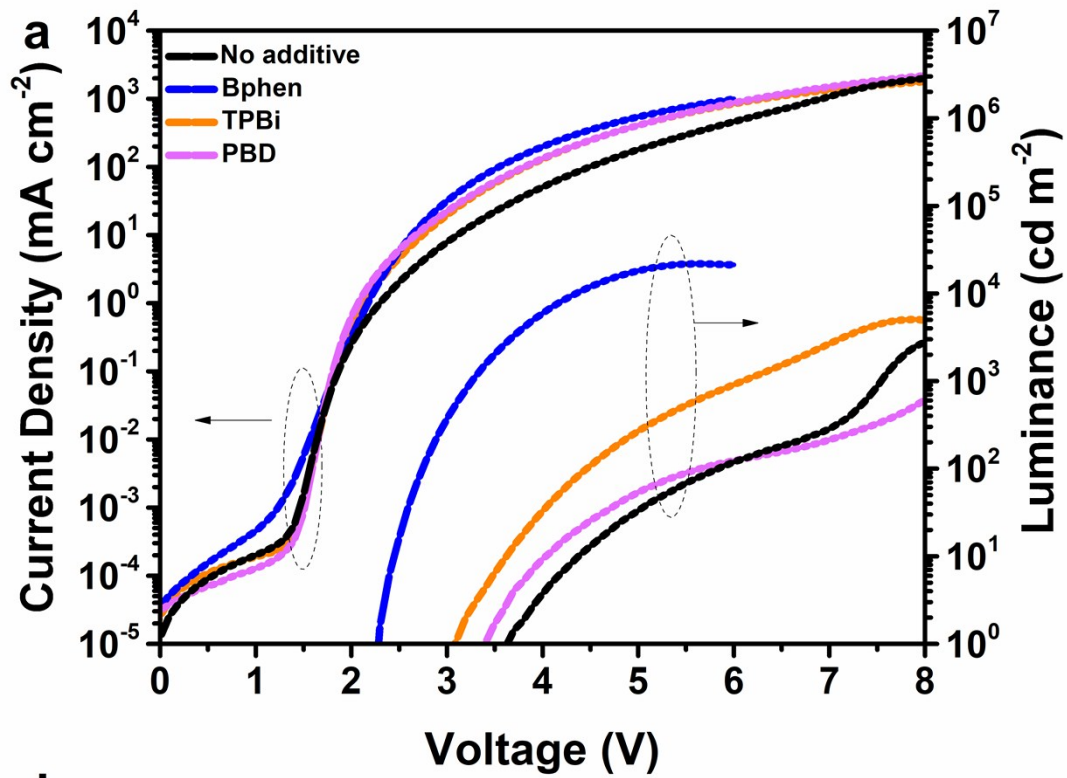


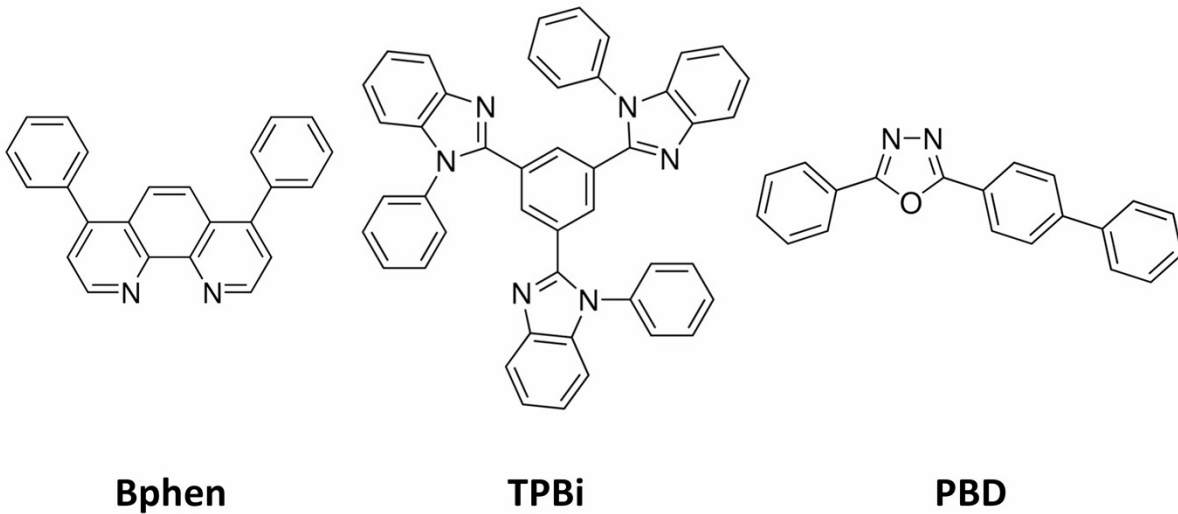
Figure S7. Plot of luminance against time for the 0 and 0.500 mg mL<sup>-1</sup> devices measured under constant current



**Figure S8.** Absorbance-corrected PL intensity of BPhen, TPBi and PBD-incorporated  $\text{CH}_3\text{NH}_3\text{PbBr}_3$  films with fixed concentration of  $0.500 \text{ mg mL}^{-1}$  in toluene.



**Figure S9.** (a) Current density and luminance against voltage and (b) current efficiency against current density of BPhen, TPBi and PBD-incorporated  $\text{CH}_3\text{NH}_3\text{PbBr}_3$  films with fixed concentration of  $0.500 \text{ mg mL}^{-1}$  in toluene.



**Figure S10.** Molecular structure of the small molecule additives used in this work: BPhen, TPBi and PBD.