

Correction to “Interlayer Exciton–Phonon Coupling in MoSe₂/WSe₂ Heterostructures”

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Supporting Information

Corrections to [Table 2](#) in the original article are as follows.

Table 2. Deformation Potentials for the A_{1g}^M and A_{1g}^W Heterostructure Eigenvectors on the MoSe₂ and WSe₂ Bands at K (° = M, W)^a

	$M_{A_{1g}^M/XA^\circ}$ (eV/Å)	$M_{A_{1g}^W/XA^W}$ (eV/Å)	ratio
H-type Stacking			
$M_{A_{1g}^M/XA^\circ}$	104	22	0.21
$M_{A_{1g}^W/XA^\circ}$	4	194	0.02
R-type Stacking			
$M_{A_{1g}^M/XA^\circ}$	160	28	0.18
$M_{A_{1g}^W/XA^\circ}$	17	199	0.09

^aThe other heterostructure stackings are reported in [Supporting Information Table I](#).

■ ASSOCIATED CONTENT

SI Supporting Information

The Supporting Information is available free of charge at <https://pubs.acs.org/doi/10.1021/acs.nanolett.4c05914>.

Experimental details and sample production as well as a description of the fitting procedure using eq 1 to fit the Raman spectra shown in Figure 3; additional figures showing the sample area, PL spectra from the monolayers, and a table of optical deformation potentials as well as how they were calculate ([PDF](#))

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