Electronic Supplementary Information

Review on the Raman spectroscopy of different types of layered

materials

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Table S1 Character table for D_{6h} point group

	E	2C ₆	2C ₃	C ₂	3C'2	3C"2	i	2S ₃	2S ₆	σ_{h}	$3\sigma_d$	3 σ _v	Linear, rotations	Quadratic
A _{1g}	1	1	1	1	1	1	1	1	1	1	1	1		x^2+y^2 , z^2
A _{2g}	1	1	1	1	-1	-1	1	1	1	1	-1	-1	Rz	
B _{1g}	1	-1	1	-1	1	-1	1	-1	1	-1	1	-1		
B _{2g}	1	-1	1	-1	-1	1	1	-1	1	-1	-1	1		
E _{1g}	2	1	-1	-2	0	0	2	1	-1	-2	0	0	(R_x, R_y)	(xz, yz)
E _{2g}	2	-1	-1	2	0	0	2	-1	-1	2	0	0		(x^2-y^2, xy)
A_{1u}	1	1	1	1	1	1	-1	-1	-1	-1	-1	-1		
A _{2u}	1	1	1	1	-1	-1	-1	-1	-1	-1	1	1	Z	
B _{1u}	1	-1	1	-1	1	-1	-1	1	-1	1	-1	1		
B _{2u}	1	-1	1	-1	-1	1	-1	1	-1	1	1	-1		
E _{1u}	2	1	-1	-2	0	0	-2	-1	1	2	0	0	(x, y)	
E_{2u}	2	-1	-1	2	0	0	-2	1	1	-2	0	0		



Fig. S1 Raman spectra of WSe₂, GaSe, CuS, GeSe and Bi_2Te_2S under 633 nm. The modes (251 cm⁻¹ (A_{1g}) for WSe₂, 209.3 cm⁻¹ (E¹_{1g}) for GaSe, 65.0 cm⁻¹ (E²_{2g}) for CuS, 175 cm⁻¹ (A²_g) for GeSe and 34.9 cm⁻¹ (E¹g) for Bi₂Te₂S) which are absent under 532-nm excitation had been detected under 633-nm excitation.



Fig. S2 Raman spectra of bulk SnSe2 under 532-nm and 633-nm excitations. These weak modes are marked in green.