

Supplementary Table 6. Data collection and refinement statistics for structural model of co-crystallized MLL1-SET and Borealin K143 peptide

	MLL1-SET <sub>3785-3969</sub> + Borealin peptide <sub>137-145</sub>
<b>Data collection</b>	
Wavelength (Å)	1.12705
Space group	P31 2 1
Cell dimensions	
<i>a</i> , <i>b</i> , <i>c</i> (Å)	54.48, 54.48, 105.63
α, β, γ (°)	90, 90, 120
Resolution (Å)	47.18 - 2.59 (2.63 - 2.59)
<i>R</i> <sub>sym</sub> or <i>R</i> <sub>merge</sub> (%)	4.6 (78.9)
<i>I</i> / <i>σ</i> <i>I</i>	11.2 (0.8)
Completeness (%)	97.96 (79.46)
Redundancy	5.7
<b>Refinement</b>	
Resolution (Å)	43.08 - 2.59 (2.68 - 2.59)
No. reflections	5878
<i>R</i> <sub>work</sub> / <i>R</i> <sub>free</sub>	24.77 / 30.36
No. atoms	
Protein	1000
Ligand/ion	1 (Zn <sup>2+</sup> ), 26 (AdoHcy),
<i>B</i> -factors	
Protein	61.12
Ligand/ion	142.86 (Zn <sup>2+</sup> ), 61.92 (AdoHcy)
R.m.s. deviations	
Bond lengths (Å)	0.01
Bond angles (°)	1.16
Ramachandran plot (%)	
Favored	93
Outliers	0
<b>Protein Data Bank code</b>	7U5V