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Supporting information for article:

Tangible symmetry elements and space-group models to guide from molecular to solid-state composition

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S1. Models of axes of rotation

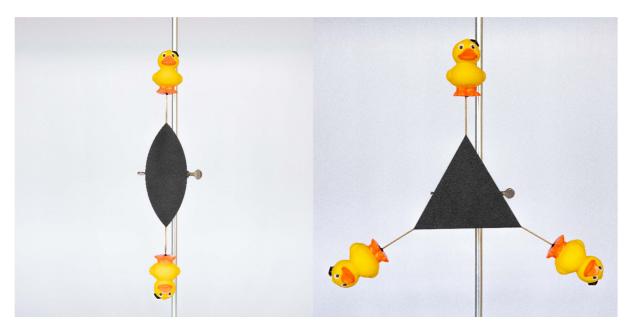


Figure S1 Models of rotational axes. Left: two-fold rotational axis. Right: three-fold rotational axis.

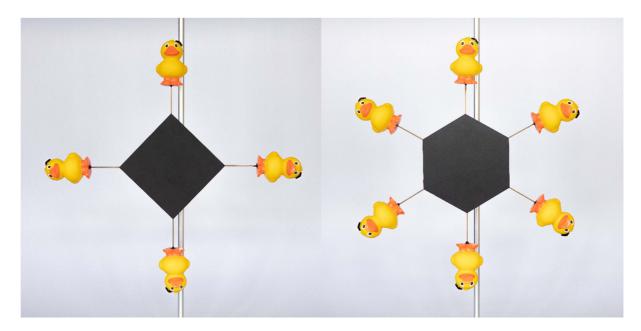


Figure S2 Models of rotational axes. Left: four-fold rotational axis. Right: six-fold rotational axis.

S2. Models of screw axes

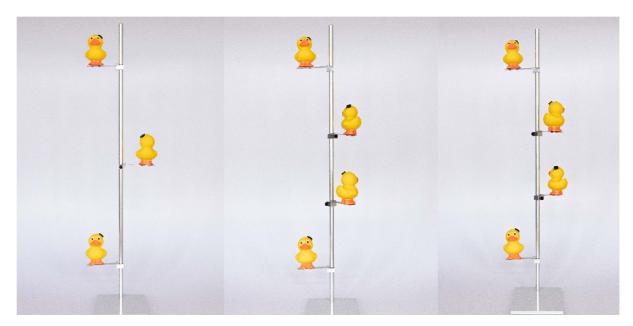


Figure S3 Models of screw axes. Left: 2_1 axis. Middle: 3_1 axis. Right: 3_2 axis.

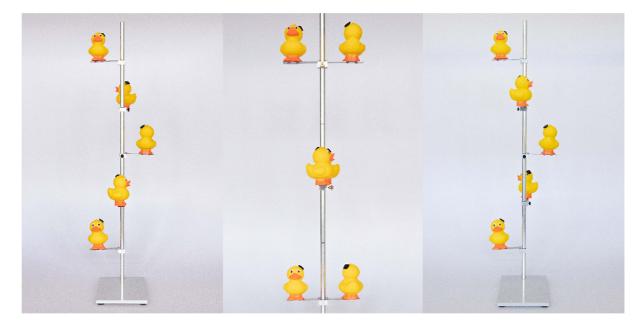


Figure S4 Models of screw axes. Left: 4_1 axis. Middle: 4_2 axis. Right: 4_3 axis.

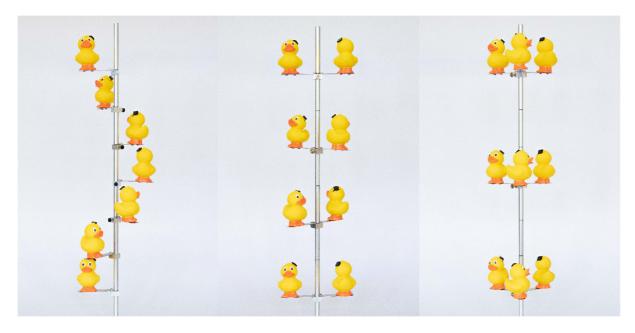


Figure S5 Models of screw axes. Left: 6_1 axis. Middle: 6_2 axis. Right: 6_3 axis.

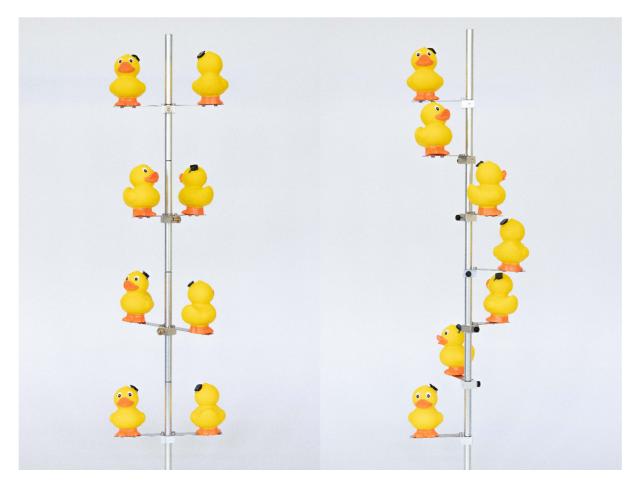


Figure S6 Models of screw axes. Left: 6_4 axis. Right: 6_5 axis.

S3. Models of space groups $P2_1$ and $P2_1/c$



Figure S7 Prototype model of space group $P2_1$ made from scrap material. Left: View along crystallographic b direction. Right: View along crystallographic a direction.



Figure S8 Small scale model of space group $P2_1$.

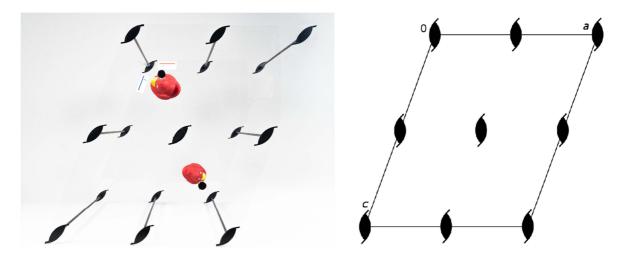


Figure S9 Left: Model of space group $P2_1$. View along crystallographic b direction. Right: Schematic depiction of space group $P2_1$. View along crystallographic b direction.

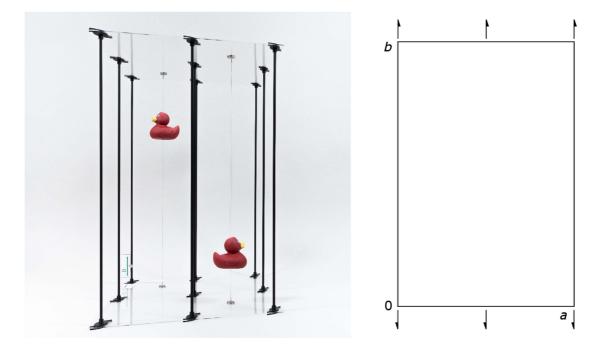


Figure S10 Left: Model of space group $P2_1$. View along crystallographic c direction. Right: Schematic depiction of space group $P2_1$. View along crystallographic c direction.

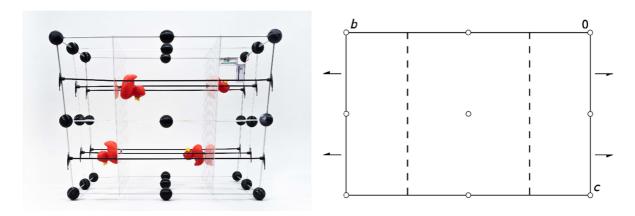


Figure S11 Left: Model of space group $P2_1/c$. View along crystallographic a direction. Right: Schematic depiction of space group $P2_1/c$. View along crystallographic a direction.

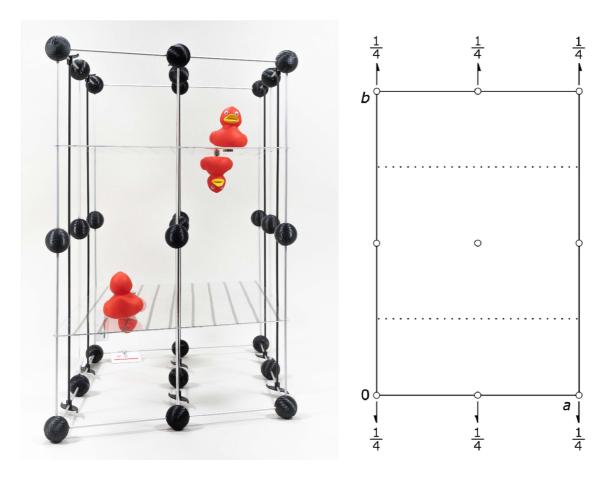


Figure S12 Left: Model of space group $P2_1/c$. View along crystallographic c direction. Right: Schematic depiction of space group $P2_1/c$. View along crystallographic c direction.