

*C53 (SrBr₂) Structure (*Obsolete*):*

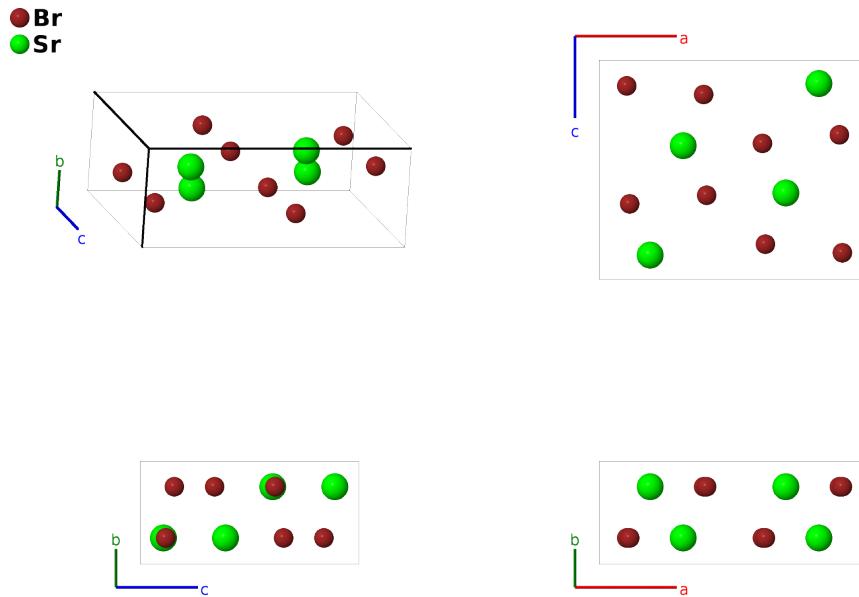
A2B_oP12_62_2c_c-010

This structure originally had the label A2B_oP12_62_2c_c.SrBr2. Calls to that address will be redirected here.

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<https://aflow.org/p/NJ0X>

https://aflow.org/p/A2B_oP12_62_2c_c-010

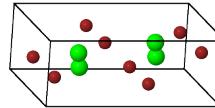


Prototype	Br ₂ Sr
AFLOW prototype label	A2B_oP12_62_2c_c-010
Strukturbericht designation	C53
ICSD	15972
Pearson symbol	oP12
Space group number	62
Space group symbol	<i>Pnma</i>
AFLOW prototype command	aflow --proto=A2B_oP12_62_2c_c-010 --params=a,b/a,c/a,x ₁ ,z ₁ ,x ₂ ,z ₂ ,x ₃ ,z ₃

- (Kamermans, 1939) determined that SrBr₂ had a distorted PbCl₂ structure and assigned it to space group *Pnma* #62. (Hermann, 1939) gave this the *Strukturbericht* designation *C53*.
- Some years later (Sass, 1963) showed that this structure was incorrect, and that α -SrBr₂ was actually in space group *P4/n* #85, rendering the *C53* symbol obsolete. We include it here for historical interest.

- (Kamermans, 1939) gave the atomic positions in terms of the *Pbnm* setting of space group #62. We have used FINDSYM to translate this to the standard *Pnma* setting.
- PbCl₂ (*C23*), HgCl₂ (*C25*), SrH₂ (*C29*), Co₂Si (*C37*), and SrBr₂ (*C53*) all share the same AFLOW label, A2B_oP12_62_2c_c. The structures are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files.

Simple Orthorhombic primitive vectors



Basis vectors

	Lattice coordinates	=	Cartesian coordinates	Wyckoff position	Atom type
\mathbf{B}_1	$x_1 \mathbf{a}_1 + \frac{1}{4} \mathbf{a}_2 + z_1 \mathbf{a}_3$	=	$a x_1 \hat{\mathbf{x}} + \frac{1}{4} b \hat{\mathbf{y}} + c z_1 \hat{\mathbf{z}}$	(4c)	Br I
\mathbf{B}_2	$-(x_1 - \frac{1}{2}) \mathbf{a}_1 + \frac{3}{4} \mathbf{a}_2 + (z_1 + \frac{1}{2}) \mathbf{a}_3$	=	$-a(x_1 - \frac{1}{2}) \hat{\mathbf{x}} + \frac{3}{4} b \hat{\mathbf{y}} + c(z_1 + \frac{1}{2}) \hat{\mathbf{z}}$	(4c)	Br I
\mathbf{B}_3	$-x_1 \mathbf{a}_1 + \frac{3}{4} \mathbf{a}_2 - z_1 \mathbf{a}_3$	=	$-a x_1 \hat{\mathbf{x}} + \frac{3}{4} b \hat{\mathbf{y}} - c z_1 \hat{\mathbf{z}}$	(4c)	Br I
\mathbf{B}_4	$(x_1 + \frac{1}{2}) \mathbf{a}_1 + \frac{1}{4} \mathbf{a}_2 - (z_1 - \frac{1}{2}) \mathbf{a}_3$	=	$a(x_1 + \frac{1}{2}) \hat{\mathbf{x}} + \frac{1}{4} b \hat{\mathbf{y}} - c(z_1 - \frac{1}{2}) \hat{\mathbf{z}}$	(4c)	Br I
\mathbf{B}_5	$x_2 \mathbf{a}_1 + \frac{1}{4} \mathbf{a}_2 + z_2 \mathbf{a}_3$	=	$a x_2 \hat{\mathbf{x}} + \frac{1}{4} b \hat{\mathbf{y}} + c z_2 \hat{\mathbf{z}}$	(4c)	Br II
\mathbf{B}_6	$-(x_2 - \frac{1}{2}) \mathbf{a}_1 + \frac{3}{4} \mathbf{a}_2 + (z_2 + \frac{1}{2}) \mathbf{a}_3$	=	$-a(x_2 - \frac{1}{2}) \hat{\mathbf{x}} + \frac{3}{4} b \hat{\mathbf{y}} + c(z_2 + \frac{1}{2}) \hat{\mathbf{z}}$	(4c)	Br II
\mathbf{B}_7	$-x_2 \mathbf{a}_1 + \frac{3}{4} \mathbf{a}_2 - z_2 \mathbf{a}_3$	=	$-a x_2 \hat{\mathbf{x}} + \frac{3}{4} b \hat{\mathbf{y}} - c z_2 \hat{\mathbf{z}}$	(4c)	Br II
\mathbf{B}_8	$(x_2 + \frac{1}{2}) \mathbf{a}_1 + \frac{1}{4} \mathbf{a}_2 - (z_2 - \frac{1}{2}) \mathbf{a}_3$	=	$a(x_2 + \frac{1}{2}) \hat{\mathbf{x}} + \frac{1}{4} b \hat{\mathbf{y}} - c(z_2 - \frac{1}{2}) \hat{\mathbf{z}}$	(4c)	Br II
\mathbf{B}_9	$x_3 \mathbf{a}_1 + \frac{1}{4} \mathbf{a}_2 + z_3 \mathbf{a}_3$	=	$a x_3 \hat{\mathbf{x}} + \frac{1}{4} b \hat{\mathbf{y}} + c z_3 \hat{\mathbf{z}}$	(4c)	Sr I
\mathbf{B}_{10}	$-(x_3 - \frac{1}{2}) \mathbf{a}_1 + \frac{3}{4} \mathbf{a}_2 + (z_3 + \frac{1}{2}) \mathbf{a}_3$	=	$-a(x_3 - \frac{1}{2}) \hat{\mathbf{x}} + \frac{3}{4} b \hat{\mathbf{y}} + c(z_3 + \frac{1}{2}) \hat{\mathbf{z}}$	(4c)	Sr I
\mathbf{B}_{11}	$-x_3 \mathbf{a}_1 + \frac{3}{4} \mathbf{a}_2 - z_3 \mathbf{a}_3$	=	$-a x_3 \hat{\mathbf{x}} + \frac{3}{4} b \hat{\mathbf{y}} - c z_3 \hat{\mathbf{z}}$	(4c)	Sr I
\mathbf{B}_{12}	$(x_3 + \frac{1}{2}) \mathbf{a}_1 + \frac{1}{4} \mathbf{a}_2 - (z_3 - \frac{1}{2}) \mathbf{a}_3$	=	$a(x_3 + \frac{1}{2}) \hat{\mathbf{x}} + \frac{1}{4} b \hat{\mathbf{y}} - c(z_3 - \frac{1}{2}) \hat{\mathbf{z}}$	(4c)	Sr I

References

- [1] M. A. Kamermans, *The Crystal Structure of SrBr₂*, Z. Kristallogr. **101**, 406–411 (1939), doi:10.1524/zkri.1939.101.1.406.
- [2] R. L. Sass, T. Brackett, and E. Brackett, *The Crystal Structure Of Strontium Bromide*, J. Phys. Chem. **67**, 2862–2863 (1963), doi:10.1021/j100806a516.

Found in

- [1] K. Herrmann, ed., *Strukturbericht Band VII 1939* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1943).