

# Li<sub>3</sub>N Structure:

## A3B\_hP4\_191\_bc\_a-001

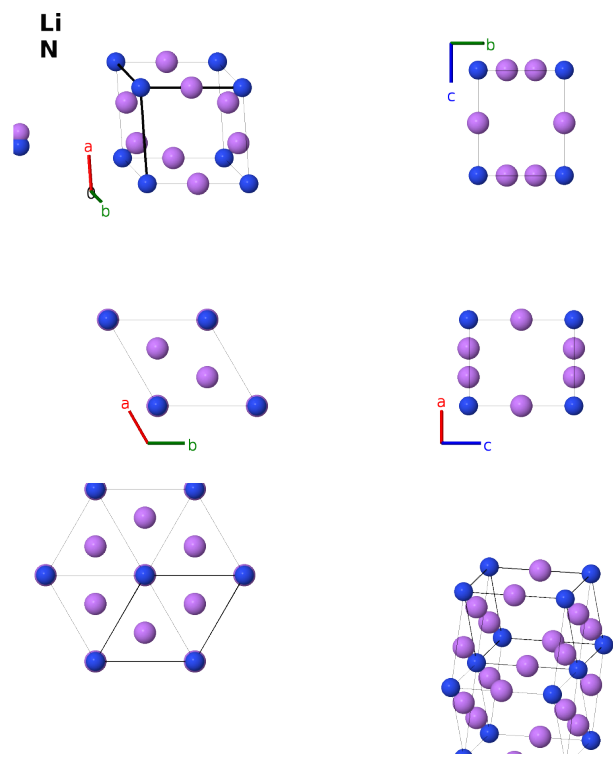
This structure previously used the label(s) **A3B\_hP4\_191\_bc\_a**. Calls to these addresses will be redirected here.

Cite this page as:

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<https://aflow.org/prototype-encyclopedia/RYZL>

[https://aflow.org/prototype-encyclopedia/A3B\\_hP4\\_191\\_bc\\_a-001](https://aflow.org/prototype-encyclopedia/A3B_hP4_191_bc_a-001)

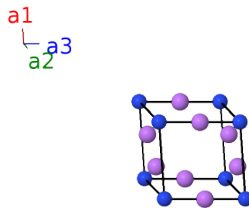


|                         |                                                                     |
|-------------------------|---------------------------------------------------------------------|
| Prototype               | Li <sub>3</sub> N                                                   |
| AFLOW prototype label   | A3B_hP4_191_bc_a-001                                                |
| ICSD                    | 95939                                                               |
| CCDC                    | 1655315                                                             |
| Pearson symbol          | hP4                                                                 |
| Space group number      | 191                                                                 |
| Space group symbol      | $P6/mmm$                                                            |
| AFLOW prototype command | <code>aflow --proto=A3B_hP4_191_bc_a-001<br/>--params=a, c/a</code> |

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## Hexagonal primitive vectors

$$\begin{aligned}\mathbf{a}_1 &= \frac{1}{2}a\hat{\mathbf{x}} - \frac{\sqrt{3}}{2}a\hat{\mathbf{y}} \\ \mathbf{a}_2 &= \frac{1}{2}a\hat{\mathbf{x}} + \frac{\sqrt{3}}{2}a\hat{\mathbf{y}} \\ \mathbf{a}_3 &= c\hat{\mathbf{z}}\end{aligned}$$




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## Basis vectors

|                | Lattice<br>coordinates |                                                     | Cartesian<br>coordinates | Wyckoff<br>position                                                  | Atom<br>type |
|----------------|------------------------|-----------------------------------------------------|--------------------------|----------------------------------------------------------------------|--------------|
| $\mathbf{B}_1$ | $=$                    | $0$                                                 | $=$                      | $0$                                                                  | N I          |
| $\mathbf{B}_2$ | $=$                    | $\frac{1}{2}\mathbf{a}_3$                           | $=$                      | $\frac{1}{2}c\hat{\mathbf{z}}$                                       | Li I         |
| $\mathbf{B}_3$ | $=$                    | $\frac{1}{3}\mathbf{a}_1 + \frac{2}{3}\mathbf{a}_2$ | $=$                      | $\frac{1}{2}a\hat{\mathbf{x}} + \frac{\sqrt{3}}{6}a\hat{\mathbf{y}}$ | Li II        |
| $\mathbf{B}_4$ | $=$                    | $\frac{2}{3}\mathbf{a}_1 + \frac{1}{3}\mathbf{a}_2$ | $=$                      | $\frac{1}{2}a\hat{\mathbf{x}} - \frac{\sqrt{3}}{6}a\hat{\mathbf{y}}$ | Li II        |

## References

- [1] D. H. Gregory, P. M. O’Meara, A. G. Gordon, J. P. Hodges, S. Short, and J. D. Jorgensen, *Structure of Lithium Nitride and Transition-Metal-Doped Derivatives,  $\text{Li}_{3-x-y}\text{M}_x\text{N}$  ( $M = \text{Ni}, \text{Cu}$ ): A Powder Neutron Diffraction Study*, Chem. Mater. **14**, 2063–2070 (2002), doi:10.1021/cm010718t.

## Found in

- [1] P. Villars, K. Cenzual, J. Daams, R. Gladyshevskii, O. Shcherban, V. Dubenskyy, N. Melnichenko-Koblyuk, O. Pavlyuk, I. Savysyuk, S. Stoyko, and L. Sysa, *Structure Types. Part 3: Space Groups (194) $P6_3/mmc$ -(190) $P-62c$ :  $\text{Li}_3\text{N}$*  (2006). Landolt-Börnstein - Group III Condensed Matter 43A6 (Springer-Verlag Berlin Heidelberg).

## First cited in

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