

Calaverite (AuTe_2 , *C*34) Structure:

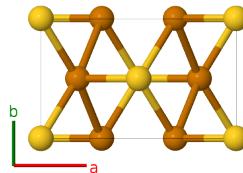
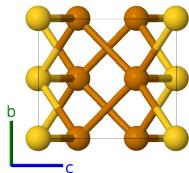
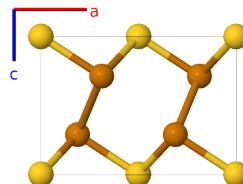
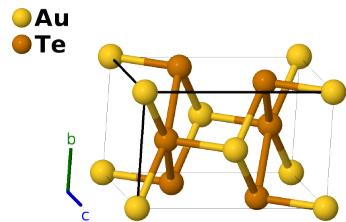
AB₂_mC₆_12_a_i-001

This structure originally had the label AB₂_mC₆_12_a_i. Calls to that address will be redirected here.

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<https://aflow.org/p/1QX2>

https://aflow.org/p/AB2_mC6_12_a_i-001



Prototype AuTe_2

AFLOW prototype label AB₂_mC₆_12_a_i-001

Strukturbericht designation *C*34

Mineral name calaverite

ICSD 72434

Pearson symbol mC6

Space group number 12

Space group symbol $C2/m$

AFLOW prototype command

```
aflow --proto=AB2_mC6_12_a_i-001  
--params=a, b/a, c/a, beta, x2, z2
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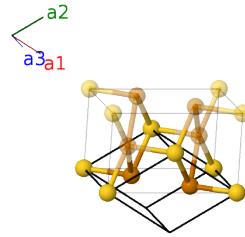
Other compounds with this structure

$\text{Au}_{10}\text{Se}_3\text{Te}_{17}$

- This structure was given the *C*34 designation by (Gottfried, 1937). (Pertlik, 1984) put calverite in space group *Pc* #7.
- AuTe_2 can also be found as krennerite (*C*46).
- *C*34 Calverite (this structure), $\alpha\text{-HgO}_2$, and NiO_2 have the same AFLOW prototype label, AB₂_mC₆_12_a_i. The structures are generated by the same symmetry operations with different sets of parameters (`--params`) specified in their corresponding CIF files.

Base-centered Monoclinic primitive vectors

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



Basis vectors

| | Lattice coordinates | = | Cartesian coordinates | Wyckoff position | Atom type |
|------------------|---|---|---|------------------|-----------|
| \mathbf{B}_1 = | 0 | = | 0 | (2a) | Au I |
| \mathbf{B}_2 = | $x_2 \mathbf{a}_1 + x_2 \mathbf{a}_2 + z_2 \mathbf{a}_3$ | = | $(ax_2 + cz_2 \cos\beta)\hat{\mathbf{x}} + cz_2 \sin\beta\hat{\mathbf{z}}$ | (4i) | Te I |
| \mathbf{B}_3 = | $-x_2 \mathbf{a}_1 - x_2 \mathbf{a}_2 - z_2 \mathbf{a}_3$ | = | $-(ax_2 + cz_2 \cos\beta)\hat{\mathbf{x}} - cz_2 \sin\beta\hat{\mathbf{z}}$ | (4i) | Te I |

References

- [1] K. Reithmayer, W. Steurer, H. Schulz, and J. L. de Boer, *High-pressure single-crystal structure study on calaverite, AuTe₂*, Acta Crystallogr. Sect. B **49**, 6–11 (1993), doi:10.1107/S0108768192007286.
- [2] C. Gottfried and F. Schossberger, eds., *Strukturbericht Band III 1933-1935* (Akademische Verlagsgesellschaft M. B. H., Leipzig, 1937).
- [3] F. Pertlik, *Kristallchemie natürlicher Telluride III: Die Kristallstruktur des Minerals Calaverit, AuTe₂*, Z. Kristallogr. **169**, 227–236 (1984), doi:10.1524/zkri.1984.169.1-4.227.