

# checkCIF/PLATON report

You have not supplied any structure factors. As a result the full set of tests cannot be run.

THIS REPORT IS FOR GUIDANCE ONLY. IF USED AS PART OF A REVIEW PROCEDURE FOR PUBLICATION, IT SHOULD NOT REPLACE THE EXPERTISE OF AN EXPERIENCED CRYSTALLOGRAPHIC REFEREE.

No syntax errors found.      CIF dictionary      Interpreting this report

## Datablock: a

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Bond precision:    Mo- O = 0.0090 Å                      Wavelength=1.54430

Cell:                      a=5.21239(8)              b=5.21239(8)              c=11.4394(3)  
                                alpha=90                      beta=90                      gamma=90  
Temperature:              300 K

|                        | Calculated                   | Reported                     |
|------------------------|------------------------------|------------------------------|
| Volume                 | 310.797(13)                  | 310.798(12)                  |
| Space group            | I 41/a                       | I41/a:2                      |
| Hall group             | -I 4ad                       | -I 4ad                       |
| Moiety formula         | Ag2 Gd0.90 Ho0.10 Mo4 O16 Yb | Ag2 Gd0.90 Ho0.10 Mo4 O16 Yb |
| Sum formula            | Ag2 Gd0.90 Ho0.10 Mo4 O16 Yb | Ag2 Gd0.90 Ho0.10 Mo4 O16 Yb |
| Mr                     | 1186.56                      | 1186.56                      |
| Dx, g cm <sup>-3</sup> | 6.340                        | 6.340                        |
| Z                      | 1                            | 1                            |
| Mu (mm <sup>-1</sup> ) | 102.522                      | 102.524                      |
| F000                   | 524.3                        | 0.0                          |
| F000'                  | 509.48                       |                              |
| h,k,lmax               | 4,4,10                       |                              |
| Nref                   | 63                           |                              |
| Tmin,Tmax              |                              |                              |
| Tmin'                  |                              |                              |

Correction method= Not given

Data completeness= 0.000                      Theta(max)=

R(reflections)=                      wR2(reflections)=

S =                      Npar=

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The following ALERTS were generated. Each ALERT has the format

**test-name\_ALERT\_alert-type\_alert-level.**

Click on the hyperlinks for more details of the test.



### Alert level C

CRYSR01\_ALERT\_1\_C \_exptl\_crystal\_size\_rad not in the CIF when expected.  
PLAT076\_ALERT\_1\_C Occupancy 0.250 Less Than 1.0 for Sp.pos . YB  
PLAT077\_ALERT\_4\_C Unitcell Contains Non-integer Number of Atoms .. Please Check  
PLAT313\_ALERT\_2\_C Oxygen with Three Covalent Bonds (rare) ..... 0 Check



### Alert level G

RADNT01\_ALERT\_1\_G Extra text has been found in the \_diffrn\_radiation\_type field.  
Radiation given as Cu K\alpha~1~2~  
Radiation identified as Cu K\alpha~1~  
PLAT004\_ALERT\_5\_G Polymeric Structure Found with Maximum Dimension 3 Info  
PLAT300\_ALERT\_4\_G Atom Site Occupancy of Yb Constrained at 0.25 Check  
PLAT300\_ALERT\_4\_G Atom Site Occupancy of Ho Constrained at 0.025 Check  
PLAT300\_ALERT\_4\_G Atom Site Occupancy of Gd Constrained at 0.225 Check  
PLAT300\_ALERT\_4\_G Atom Site Occupancy of Ag Constrained at 0.5 Check  
PLAT301\_ALERT\_3\_G Main Residue Disorder .....(Resd 1 ) 41% Note  
PLAT794\_ALERT\_5\_G Tentative Bond Valency for Mo (VI) . 5.98 Info  
PLAT811\_ALERT\_5\_G No ADDSYM Analysis: Too Many Excluded Atoms .... ! Info

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0 **ALERT level A** = Most likely a serious problem - resolve or explain  
0 **ALERT level B** = A potentially serious problem, consider carefully  
4 **ALERT level C** = Check. Ensure it is not caused by an omission or oversight  
9 **ALERT level G** = General information/check it is not something unexpected

3 ALERT type 1 CIF construction/syntax error, inconsistent or missing data  
1 ALERT type 2 Indicator that the structure model may be wrong or deficient  
1 ALERT type 3 Indicator that the structure quality may be low  
5 ALERT type 4 Improvement, methodology, query or suggestion  
3 ALERT type 5 Informative message, check

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It is advisable to attempt to resolve as many as possible of the alerts in all categories. Often the minor alerts point to easily fixed oversights, errors and omissions in your CIF or refinement strategy, so attention to these fine details can be worthwhile. In order to resolve some of the more serious problems it may be necessary to carry out additional measurements or structure refinements. However, the purpose of your study may justify the reported deviations and the more serious of these should normally be commented upon in the discussion or experimental section of a paper or in the "special\_details" fields of the CIF. checkCIF was carefully designed to identify outliers and unusual parameters, but every test has its limitations and alerts that are not important in a particular case may appear. Conversely, the absence of alerts does not guarantee there are no aspects of the results needing attention. It is up to the individual to critically assess their own results and, if necessary, seek expert advice.

### **Publication of your CIF in IUCr journals**

A basic structural check has been run on your CIF. These basic checks will be run on all CIFs submitted for publication in IUCr journals (*Acta Crystallographica*, *Journal of Applied Crystallography*, *Journal of Synchrotron Radiation*); however, if you intend to submit to *Acta Crystallographica Section C* or *E* or *IUCrData*, you should make sure that full publication checks are run on the final version of your CIF prior to submission.

### **Publication of your CIF in other journals**

Please refer to the *Notes for Authors* of the relevant journal for any special instructions relating to CIF submission.

