

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

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

Cell: a=5.22237(8) b=5.22237(8) c=11.4581(2)
 alpha=90 beta=90 gamma=90
Temperature: 300 K

	Calculated	Reported
Volume	312.499(11)	312.497(11)
Space group	I 41/a	I41/a:2
Hall group	-I 4ad	-I 4ad
Moiety formula	Ag2 Gd1.20 Ho0.10 Mo4 O16 Yb0.70	Ag2 Gd1.20 Ho0.10 Mo4 O16 Yb0.70
Sum formula	Ag2 Gd1.20 Ho0.10 Mo4 O16 Yb0.70	Ag2 Gd1.20 Ho0.10 Mo4 O16 Yb0.70
Mr	1181.82	1181.82
Dx, g cm-3	6.280	6.280
Z	1	1
Mu (mm-1)	108.107	108.109
F000	522.5	0.0
F000'	507.36	
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=

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.

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 0.175 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 0.3 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) . 5.97 Info

PLAT811_ALERT_5_G No ADDSYM Analysis: Too Many Excluded Atoms ! Info

0 **ALERT level A** = Most likely a serious problem - resolve or explain

0 **ALERT level B** = A potentially serious problem, consider carefully

3 **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

2 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

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.

