

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: MgHDTMP_180R_publ

Bond precision: C-C = 0.0178 A Wavelength=1.54060

Cell: a=12.36315(16) b=9.07202(13) c=9.55250(13)
 alpha=90.6651(8) beta=106.7290(9) gamma=111.8975(8)
Temperature: 298 K

	Calculated	Reported
Volume	943.46(2)	943.45(2)
Space group	P -1	P -1
Hall group	-P 1	-P 1
Moiety formula	C10 N2 O12 P4	?
Sum formula	C10 N2 O12 P4	C10 Mg N2 O12 P4
Mr	464.00	488.31
Dx,g cm-3	1.633	0.000
Z	2	2
Mu (mm-1)	4.235	0.000
F000	460.0	0.0
F000'	464.01	
h,k,lmax	12,9,9	12,9,9
Nref	1948	1947
Tmin,Tmax		1.000,1.000
Tmin'		

Correction method= # Reported T Limits: Tmin=1.000 Tmax=1.000
AbsCorr = ?

Data completeness= 0.999 Theta(max)= 49.992

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 B

PLAT043_ALERT_1_B Calculated and Reported Mol. Weight Differ by .. 24.31 Check
PLAT340_ALERT_3_B Low Bond Precision on C-C Bonds 0.0178 Ang.

The atomic mass of the Mg atom has not been taking into account for the molecular weight. The checkcif program is not interpreting correctly the cif file creates by the GSAS program

Alert level C

PLAT041_ALERT_1_C Calc. and Reported SumFormula Strings Differ Please Check
PLAT074_ALERT_1_C Occupancy Parameter = 0.0 for MG1 Check
PLAT074_ALERT_1_C Occupancy Parameter = 0.0 for MG2 Check
PLAT155_ALERT_4_C The Triclinic Unitcell is NOT Reduced Please Do !

Alert level G

FORMU01_ALERT_2_G There is a discrepancy between the atom counts in the
_chemical_formula_sum and the formula from the _atom_site* data.

Atom count from _chemical_formula_sum: C10 Mg1 N2 O12 P4

Atom count from the _atom_site data: C10 N2 O12 P4

CELLZ01_ALERT_1_G Difference between formula and atom_site contents detected.

CELLZ01_ALERT_1_G ALERT: Large difference may be due to a

symmetry error - see SYMMG tests

From the CIF: _cell_formula_units_Z 2

From the CIF: _chemical_formula_sum C10 Mg N2 O12 P4

TEST: Compare cell contents of formula and atom_site data

atom	Z*formula	cif sites	diff
P	8.00	8.00	0.00
O	24.00	24.00	0.00
C	20.00	20.00	0.00
N	4.00	4.00	0.00
Mg	2.00	0.00	2.00

PLAT004_ALERT_5_G Polymeric Structure Found with Maximum Dimension 3 Info
PLAT040_ALERT_1_G No H-atoms in this Carbon Containing Compound .. Please Check
PLAT092_ALERT_4_G Check: Wavelength Given is not Cu,Ga,Mo,Ag,In Ka 1.54060 Ang.
PLAT112_ALERT_2_G ADDSYM Detects New (Pseudo) Symm. Elem n 92 %Fit
PLAT113_ALERT_2_G ADDSYM Suggests Possible Pseudo/New Space Group I2/c Check
PLAT300_ALERT_4_G Atom Site Occupancy of Mg1 Constrained at 0.0 Check
PLAT300_ALERT_4_G Atom Site Occupancy of Mg2 Constrained at 0.0 Check
PLAT860_ALERT_3_G Number of Least-Squares Restraints 61 Note
PLAT981_ALERT_1_G No non-zero f" Anomalous Scattering Values Found Please Check
PLAT986_ALERT_1_G No non-zero f' Anomalous Scattering Values Found Please Check

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

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

PLATON version of 14/07/2018; check.def file version of 05/06/2018

