

## 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: new\_sq

---

Bond precision:    C-C = 0.0126 Å                      Wavelength=0.70000

Cell:                      a=29.915(4)              b=29.915(4)              c=34.378(7)  
                                    alpha=90              beta=90              gamma=90  
Temperature:              193 K

	Calculated	Reported
Volume	30765(10)	30765(11)
Space group	P 4/m n c	P 4/m n c
Hall group	-P 4 2n	-P 4 2n
Moiety formula	2(C79 H41.60 N7.40 O17.90 Zn5.50) [+ solvent]	?
Sum formula	C158 H83.20 N14.80 O35.80 Zn11 [+ solvent]	C158 H83.20 N14.80 O35.80 Zn11
Mr	3480.89	3480.66
Dx, g cm <sup>-3</sup>	0.752	0.751
Z	4	4
Mu (mm <sup>-1</sup> )	0.845	0.845
F000	7004.8	7005.0
F000'	7020.42	
h,k,lmax	34,34,39	34,34,39
Nref	12144	12091
Tmin,Tmax	0.816,0.859	0.707,0.863
Tmin'	0.689	

Correction method= # Reported T Limits: Tmin=0.707 Tmax=0.863  
AbsCorr = MULTI-SCANS

Data completeness= 0.996                      Theta(max)= 23.498

R(reflections)= 0.1291( 10477)              wR2(reflections)= 0.4435( 12091)

S = 2.113                      Npar= 511

---

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

THETM01\_ALERT\_3\_B The value of  $\sin(\theta_{\max})/\lambda$  is less than 0.575  
Calculated  $\sin(\theta_{\max})/\lambda = 0.5696$   
PLAT084\_ALERT\_3\_B High wR2 Value (i.e. > 0.25) ..... 0.44 Report  
PLAT213\_ALERT\_2\_B Atom C8P has ADP max/min Ratio ..... 4.1 prolat  
PLAT241\_ALERT\_2\_B High 'MainMol' Ueq as Compared to Neighbors of 02 Check  
PLAT241\_ALERT\_2\_B High 'MainMol' Ueq as Compared to Neighbors of 02W Check  
PLAT241\_ALERT\_2\_B High 'MainMol' Ueq as Compared to Neighbors of 04 Check  
PLAT241\_ALERT\_2\_B High 'MainMol' Ueq as Compared to Neighbors of 06 Check  
PLAT241\_ALERT\_2\_B High 'MainMol' Ueq as Compared to Neighbors of N2 Check  
PLAT241\_ALERT\_2\_B High 'MainMol' Ueq as Compared to Neighbors of C32 Check  
PLAT242\_ALERT\_2\_B Low 'MainMol' Ueq as Compared to Neighbors of Zn5 Check  
PLAT242\_ALERT\_2\_B Low 'MainMol' Ueq as Compared to Neighbors of C26 Check

---

### Alert level C

ABSMU01\_ALERT\_1\_C The ratio of given/expected absorption coefficient lies  
outside the range 0.99 <> 1.01  
Calculated value of  $\mu = 0.881$   
Value of  $\mu$  given = 0.845  
GOODF01\_ALERT\_2\_C The least squares goodness of fit parameter lies  
outside the range 0.80 <> 2.00  
Goodness of fit given = 2.113  
RADNW01\_ALERT\_1\_C The radiation wavelength lies outside the expected range  
for the supplied radiation type. Expected range 0.71065-0.71075  
Wavelength given = 0.70000  
PLAT082\_ALERT\_2\_C High R1 Value ..... 0.13 Report  
PLAT087\_ALERT\_2\_C Unsatisfactory S value (Too High) ..... 2.11 Check  
PLAT094\_ALERT\_2\_C Ratio of Maximum / Minimum Residual Density .... 2.07 Report  
PLAT213\_ALERT\_2\_C Atom O1W has ADP max/min Ratio ..... 3.3 prolat  
PLAT213\_ALERT\_2\_C Atom C26 has ADP max/min Ratio ..... 3.1 prolat  
PLAT213\_ALERT\_2\_C Atom C6P has ADP max/min Ratio ..... 3.5 prolat  
PLAT220\_ALERT\_2\_C Non-Solvent Resd 1 C Ueq(max)/Ueq(min) Range 3.2 Ratio  
PLAT220\_ALERT\_2\_C Non-Solvent Resd 1 O Ueq(max)/Ueq(min) Range 4.4 Ratio  
PLAT241\_ALERT\_2\_C High 'MainMol' Ueq as Compared to Neighbors of 01 Check  
PLAT241\_ALERT\_2\_C High 'MainMol' Ueq as Compared to Neighbors of 03 Check  
PLAT241\_ALERT\_2\_C High 'MainMol' Ueq as Compared to Neighbors of 05 Check  
PLAT241\_ALERT\_2\_C High 'MainMol' Ueq as Compared to Neighbors of N3 Check  
PLAT241\_ALERT\_2\_C High 'MainMol' Ueq as Compared to Neighbors of N4 Check  
PLAT241\_ALERT\_2\_C High 'MainMol' Ueq as Compared to Neighbors of C17 Check  
PLAT241\_ALERT\_2\_C High 'MainMol' Ueq as Compared to Neighbors of C22 Check  
PLAT241\_ALERT\_2\_C High 'MainMol' Ueq as Compared to Neighbors of C25 Check  
PLAT241\_ALERT\_2\_C High 'MainMol' Ueq as Compared to Neighbors of C28 Check  
PLAT241\_ALERT\_2\_C High 'MainMol' Ueq as Compared to Neighbors of C34 Check  
PLAT242\_ALERT\_2\_C Low 'MainMol' Ueq as Compared to Neighbors of Zn2 Check  
PLAT242\_ALERT\_2\_C Low 'MainMol' Ueq as Compared to Neighbors of N1P Check  
PLAT242\_ALERT\_2\_C Low 'MainMol' Ueq as Compared to Neighbors of C13 Check  
PLAT242\_ALERT\_2\_C Low 'MainMol' Ueq as Compared to Neighbors of C18 Check  
PLAT242\_ALERT\_2\_C Low 'MainMol' Ueq as Compared to Neighbors of C20 Check  
PLAT242\_ALERT\_2\_C Low 'MainMol' Ueq as Compared to Neighbors of C29 Check  
PLAT242\_ALERT\_2\_C Low 'MainMol' Ueq as Compared to Neighbors of C31 Check  
PLAT242\_ALERT\_2\_C Low 'MainMol' Ueq as Compared to Neighbors of C33 Check  
PLAT250\_ALERT\_2\_C Large U3/U1 Ratio for Average U(i,j) Tensor .... 2.4 Note  
PLAT341\_ALERT\_3\_C Low Bond Precision on C-C Bonds ..... 0.01256 Ang.  
PLAT731\_ALERT\_1\_C Bond Calc 1.483(13), Rep 1.48(7) ..... 5 su-Rat  
C34 -C18 1.555 7.665 ..... # 77 Check

---

## Alert level G

ABSTY01\_ALERT\_1\_G Extra text has been found in the \_exptl\_absorpt\_correction\_type field, which should be only a single keyword. A literature citation should be included in the \_exptl\_absorpt\_process\_details field.

PLAT002_ALERT_2_G	Number of Distance or Angle Restraints on AtSite	11	Note
PLAT003_ALERT_2_G	Number of Uiso or Uij Restrained non-H Atoms ...	5	Report
PLAT004_ALERT_5_G	Polymeric Structure Found with Maximum Dimension	3	Info
PLAT068_ALERT_1_G	Reported F000 Differs from Calcd (or Missing)...		Please Check
PLAT072_ALERT_2_G	SHELXL First Parameter in WGHT Unusually Large	0.20	Report
PLAT092_ALERT_4_G	Check: Wavelength given is not Cu,Ga,Mo,Ag,In Ka	0.70000	Ang.
PLAT172_ALERT_4_G	The CIF-Embedded .res File Contains DFIX Records	8	Report
PLAT173_ALERT_4_G	The CIF-Embedded .res File Contains DANG Records	4	Report
PLAT174_ALERT_4_G	The CIF-Embedded .res File Contains FLAT Records	3	Report
PLAT177_ALERT_4_G	The CIF-Embedded .res File Contains DELU Records	2	Report
PLAT178_ALERT_4_G	The CIF-Embedded .res File Contains SIMU Records	2	Report
PLAT186_ALERT_4_G	The CIF-Embedded .res File Contains ISOR Records	3	Report
PLAT300_ALERT_4_G	Atom Site Occupancy of Zn4 is Constrained at	0.55	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of Zn3 is Constrained at	0.45	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of O1W is Constrained at	0.45	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of C1P is Constrained at	0.5	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of C2P is Constrained at	0.5	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of N2P is Constrained at	0.45	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of C4P is Constrained at	0.45	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of C5P is Constrained at	0.45	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of C6P is Constrained at	0.45	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of C7P is Constrained at	0.45	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of C8P is Constrained at	0.45	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of H1P is Constrained at	0.5	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of H2P is Constrained at	0.5	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of H4P is Constrained at	0.45	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of H5P is Constrained at	0.45	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of H7P is Constrained at	0.45	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of H8P is Constrained at	0.45	Check
PLAT301_ALERT_3_G	Main Residue Disorder .....(Resd 1)..	9%	Note
PLAT606_ALERT_4_G	VERY LARGE Solvent Accessible VOID(S) in Structure	!	Info
PLAT779_ALERT_4_G	Suspect or Irrelevant (Bond) Angle in CIF .... #	150	Check
	ZN4 -O2 -ZN3 1.555 1.555 1.555	27.30	Deg.
PLAT779_ALERT_4_G	Suspect or Irrelevant (Bond) Angle in CIF .... #	154	Check
	ZN4 -O4 -ZN3 1.555 1.555 1.555	22.06	Deg.
PLAT779_ALERT_4_G	Suspect or Irrelevant (Bond) Angle in CIF .... #	158	Check
	ZN3 -O6 -ZN4 1.555 1.555 1.555	19.73	Deg.
PLAT779_ALERT_4_G	Suspect or Irrelevant (Bond) Angle in CIF .... #	192	Check
	ZN4 -O2W -ZN3 1.555 1.555 1.555	28.51	Deg.
PLAT860_ALERT_3_G	Number of Least-Squares Restraints .....	54	Note
PLAT869_ALERT_4_G	ALERTS Related to the use of SQUEEZE Suppressed	!	Info

---

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

5 **ALERT type 1** CIF construction/syntax error, inconsistent or missing data  
40 **ALERT type 2** Indicator that the structure model may be wrong or deficient  
5 **ALERT type 3** Indicator that the structure quality may be low  
30 **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 13/08/2017; check.def file version of 27/07/2017**

