

# checkCIF/PLATON report

Structure factors have been supplied for datablock(s) 20190115-yxg-a560

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: 20190115-yxg-a560

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Bond precision:    C-C = 0.0033 Å                      Wavelength=0.71073

Cell:                      a=9.5869(3)              b=18.6505(7)              c=23.2010(9)  
                            alpha=90              beta=90              gamma=90  
Temperature:              293 K

|                | Calculated             | Reported               |
|----------------|------------------------|------------------------|
| Volume         | 4148.4(3)              | 4148.3(3)              |
| Space group    | P b c a                | P b c a                |
| Hall group     | -P 2ac 2ab             | -P 2ac 2ab             |
| Moiety formula | C23 H17 N3 O4 Zn, H2 O | C23 H17 N3 O4 Zn, H2 O |
| Sum formula    | C23 H19 N3 O5 Zn       | C23 H19 N3 O5 Zn       |
| Mr             | 482.80                 | 482.78                 |
| Dx,g cm-3      | 1.546                  | 1.546                  |
| Z              | 8                      | 8                      |
| Mu (mm-1)      | 1.226                  | 1.226                  |
| F000           | 1984.0                 | 1984.0                 |
| F000'          | 1987.06                |                        |
| h,k,lmax       | 13,25,32               | 12,25,29               |
| Nref           | 5710                   | 4878                   |
| Tmin,Tmax      | 0.779,0.874            | 0.944,1.000            |
| Tmin'          | 0.684                  |                        |

Correction method= # Reported T Limits: Tmin=0.944 Tmax=1.000  
AbsCorr = MULTI-SCAN

Data completeness= 0.854                      Theta(max)= 29.391

R(reflections)= 0.0353( 3771)              wR2(reflections)= 0.0864( 4878)

S = 1.044                      Npar= 298

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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

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|  |       |       |
|--|-------|-------|
| PLAT906_ALERT_3_C Large K Value in the Analysis of Variance .....  | 2.421 | Check |
| PLAT910_ALERT_3_C Missing # of FCF Reflection(s) Below Theta(Min). | 8     | Note  |

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### Alert level G

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|  |      |        |
|--|------|--------|
| PLAT007_ALERT_5_G Number of Unrefined Donor-H Atoms .....          | 3    | Report |
| PLAT199_ALERT_1_G Reported _cell_measurement_temperature ..... (K) | 293  | Check  |
| PLAT200_ALERT_1_G Reported _diffrn_ambient_temperature ..... (K)   | 293  | Check  |
| PLAT232_ALERT_2_G Hirshfeld Test Diff (M-X) Zn1 --N2 .             | 5.4  | s.u.   |
| PLAT794_ALERT_5_G Tentative Bond Valency for Zn1 (II) .            | 2.07 | Info   |
| PLAT912_ALERT_4_G Missing # of FCF Reflections Above STh/L= 0.600  | 753  | Note   |
| PLAT952_ALERT_5_G Calculated (ThMax) and CIF-Reported Lmax Differ  | 3    | Units  |
| PLAT958_ALERT_1_G Calculated (ThMax) and Actual (FCF) Lmax Differ  | 3    | Units  |
| PLAT978_ALERT_2_G Number C-C Bonds with Positive Residual Density. | 13   | 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  
2 **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  
2 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  
1 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.

Datablock 20190115-yxg-a560 - ellipsoid plot

