

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: p4mnc_sq

Bond precision: C-C = 0.0125 Å Wavelength=0.70000

Cell: a=30.326(4) b=30.326(4) c=34.476(7)
 alpha=90 beta=90 gamma=90
Temperature: 253 K

	Calculated	Reported
Volume	31706(11)	31706(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 ⁻³	0.729	0.729
Z	4	4
Mu (mm ⁻¹)	0.820	0.820
F000	7004.8	7005.0
F000'	7020.42	
h,k,lmax	34,34,39	34,34,39
Nref	12545	12486
Tmin,Tmax	0.821,0.863	0.714,0.866
Tmin'	0.697	

Correction method= # Reported T Limits: Tmin=0.714 Tmax=0.866
AbsCorr = MULTI-SCANS

Data completeness= 0.995 Theta(max)= 23.500

R(reflections)= 0.1242(10497) wR2(reflections)= 0.4354(12486)

S = 2.031 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
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 C16 Check
PLAT242_ALERT_2_B Low 'MainMol' Ueq as Compared to Neighbors of Zn5 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 $\mu = 0.855$
Value of μ given = 0.820
GOODF01_ALERT_2_C The least squares goodness of fit parameter lies
outside the range 0.80 <> 2.00
Goodness of fit given = 2.031
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 0.12 Report
PLAT087_ALERT_2_C Unsatisfactory S value (Too High) 2.03 Check
PLAT213_ALERT_2_C Atom O1W has ADP max/min Ratio 3.1 prolat
PLAT213_ALERT_2_C Atom C28 has ADP max/min Ratio 3.3 prolat
PLAT213_ALERT_2_C Atom C29 has ADP max/min Ratio 3.2 prolat
PLAT220_ALERT_2_C Non-Solvent Resd 1 C Ueq(max)/Ueq(min) Range 3.1 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 N2 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 C8 Check
PLAT241_ALERT_2_C High 'MainMol' Ueq as Compared to Neighbors of C24 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 C15 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 C26 Check
PLAT242_ALERT_2_C Low 'MainMol' Ueq as Compared to Neighbors of C34 Check
PLAT250_ALERT_2_C Large U3/U1 Ratio for Average U(i,j) Tensor ... 2.2 Note
PLAT341_ALERT_3_C Low Bond Precision on C-C Bonds 0.01249 Ang.
PLAT369_ALERT_2_C Long C(sp2)-C(sp2) Bond C13 - C14 .. 1.55 Ang.
PLAT369_ALERT_2_C Long C(sp2)-C(sp2) Bond C25 - C29 .. 1.55 Ang.
PLAT731_ALERT_1_C Bond Calc 1.510(14), Rep 1.51(7) 5 su-Rat
C34 -C18 1.555 3.755 # 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	9	Note
PLAT003_ALERT_2_G	Number of Uiso or Uij Restrained non-H Atoms ...	32	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	7	Report
PLAT173_ALERT_4_G	The CIF-Embedded .res File Contains DANG Records	5	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	5	Report
PLAT178_ALERT_4_G	The CIF-Embedded .res File Contains SIMU Records	5	Report
PLAT186_ALERT_4_G	The CIF-Embedded .res File Contains ISOR Records	6	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 ClP 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	27.53	Deg.
PLAT779_ALERT_4_G	Suspect or Irrelevant (Bond) Angle in CIF #	154	Check
	ZN4 -O4 -ZN3 1.555 1.555 1.555	21.49	Deg.
PLAT779_ALERT_4_G	Suspect or Irrelevant (Bond) Angle in CIF #	158	Check
	ZN3 -O6 -ZN4 1.555 1.555 1.555	20.30	Deg.
PLAT779_ALERT_4_G	Suspect or Irrelevant (Bond) Angle in CIF #	192	Check
	ZN4 -O2W -ZN3 1.555 1.555 1.555	29.40	Deg.
PLAT860_ALERT_3_G	Number of Least-Squares Restraints	425	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
 8 **ALERT level B** = A potentially serious problem, consider carefully
 28 **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
 33 **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

