

checkCIF/PLATON report

Structure factors have been supplied for datablock(s) muf-91-6h-50pc

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: muf-91-6h-50pc

Bond precision: = 0.0000 A

Wavelength=0.95372

Cell: a=17.101 (9) b=17.101 (9) c=17.101 (9)

alpha=90

beta=90

gamma=90

Temperature: 100 K

	Calculated	Reported
Volume	5001 (8)	5001 (8)
Space group	P -4 3 m	P -4 3 m
Hall group	P -4 2 3	P -4 2 3
Moiety formula	C13.15 H7.51 N0.75 O2.44 Zn0.75	C84 H48 N6 O13 Zn4, C21.139 H12.079 O6.543 Zn2.013
Sum formula	C13.15 H7.51 N0.75 O2.44 Zn0.75	C105.14 H60.08 N6 O19.54 Zn6.01
Mr	264.25	2112.86
Dx, g cm ⁻³	0.702	0.702
Z	8	1
Mu (mm ⁻¹)	1.632	1.609
F000	1070.0	1070.0
F000'	1070.18	
h, k, lmax	14, 14, 14	14, 13, 14
Nref	624 [358]	623
Tmin, Tmax		0.468, 0.746
Tmin'		

Correction method= # Reported T Limits: Tmin=0.468 Tmax=0.746

AbsCorr = MULTII-SCAN

Data completeness= 1.74/1.00

Theta(max)= 23.409

R(reflections)= 0.1808(370)

wR2(reflections)=
0.4170(623)

S = 1.543

Npar= 33

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 A

THETM01_ALERT_3_A The value of sine(theta_max)/wavelength is less than 0.550

Calculated sin(theta_max)/wavelength = 0.4166

PLAT601_ALERT_2_A Unit Cell Contains Solvent Accessible VOIDS of . 1608 Ang**3

Alert level B

PLAT049_ALERT_1_B Calculated Density Less Than 1.0 gcm-3 0.7019 Check

PLAT082_ALERT_2_B High R1 Value 0.18 Report

PLAT084_ALERT_3_B High wR2 Value (i.e. > 0.25) 0.42 Report

Alert level C

RINTA01_ALERT_3_C The value of Rint is greater than 0.12

Rint given 0.131

STRVA01_ALERT_4_C Flack test results are ambiguous.

From the CIF: _refine_ls_abs_structure_Flack 0.360

From the CIF: _refine_ls_abs_structure_Flack_su 0.030

PLAT020_ALERT_3_C The Value of Rint is Greater Than 0.12 0.131 Report

PLAT041_ALERT_1_C Calc. and Reported SumFormula Strings Differ Please Check

PLAT053_ALERT_1_C Minimum Crystal Dimension Missing (or Error) ... Please Check

PLAT054_ALERT_1_C Medium Crystal Dimension Missing (or Error) ... Please Check

PLAT055_ALERT_1_C Maximum Crystal Dimension Missing (or Error) ... Please Check

PLAT077_ALERT_4_C Unitcell Contains Non-integer Number of Atoms .. Please Check

PLAT250_ALERT_2_C Large U3/U1 Ratio for Average U(i,j) Tensor 2.9 Note

PLAT260_ALERT_2_C Large Average Ueq of Residue Including Zn1 0.296 Check

PLAT905_ALERT_3_C Negative K value in the Analysis of Variance ... -0.004 Report

PLAT918_ALERT_3_C Reflection(s) with I(obs) much Smaller I(calc) . 2 Check

PLAT934_ALERT_3_C Number of (Iobs-Icalc)/Sigma(W) > 10 Outliers .. 1 Check

PLAT939_ALERT_3_C Large Value of Not (SHELXL) Weight Optimized S . 17.97 Check

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: C105.1399 H60.08 N6 O19.54 Zn6.0

Atom count from the _atom_site data: C78.81600 H54.04799 N6 O9.086700

ABSMU01_ALERT_1_G Calculation of _exptl_absorpt_correction_mu

not performed for this radiation type.

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 1

From the CIF: _chemical_formula_sum C105.14 H60.08 N6 O19.54 Zn6.01

TEST: Compare cell contents of formula and atom_site data

	atom	Z*formula	cif sites	diff	
	C	105.14	78.82	26.32	
	H	60.08	54.05	6.03	
	N	6.00	6.00	0.00	
	O	19.54	9.09	10.45	
	Zn	6.01	1.00	5.01	
PLAT002_ALERT_2_G	Number of Distance or Angle Restraints on AtSite				10 Note
PLAT003_ALERT_2_G	Number of Uiso or Uij Restrained non-H Atoms ...				18 Report
PLAT007_ALERT_5_G	Number of Unrefined Donor-H Atoms				1 Report
PLAT042_ALERT_1_G	Calc. and Reported Moiety Formula Strings Differ				Please Check
PLAT045_ALERT_1_G	Calculated and Reported Z Differ by a Factor ...				8.0000 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				0.95372 Ang.
PLAT171_ALERT_4_G	The CIF-Embedded .res File Contains EADP Records				3 Report
PLAT172_ALERT_4_G	The CIF-Embedded .res File Contains DFIX Records				11 Report
PLAT174_ALERT_4_G	The CIF-Embedded .res File Contains FLAT Records				3 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				1 Report
PLAT300_ALERT_4_G	Atom Site Occupancy of Zn0A Constrained at				0.1667 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of O2 Constrained at				0.5 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of O1 Constrained at				0.0417 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of C12 Constrained at				0.5 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of C13 Constrained at				0.5 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of N1 Constrained at				0.25 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of C10 Constrained at				0.25 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of C11 Constrained at				0.25 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of C14 Constrained at				0.25 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of C15 Constrained at				0.25 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of C16 Constrained at				0.25 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of C17 Constrained at				0.25 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of C18 Constrained at				0.25 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of C19 Constrained at				0.25 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of C20 Constrained at				0.25 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of C21 Constrained at				0.25 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of H12 Constrained at				0.5 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of H13 Constrained at				0.25 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of H17 Constrained at				0.25 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of H18 Constrained at				0.25 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of H19 Constrained at				0.25 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of H20 Constrained at				0.25 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of H21 Constrained at				0.25 Check
PLAT301_ALERT_3_G	Main Residue Disorder(Resd 1)				100% Note
PLAT304_ALERT_4_G	Non-Integer Number of Atoms in (Resd 1)				24.60 Check
PLAT315_ALERT_2_G	Singly Bonded Carbon Detected (H-atoms Missing).				C14 Check
PLAT315_ALERT_2_G	Singly Bonded Carbon Detected (H-atoms Missing).				C34 Check
PLAT432_ALERT_2_G	Short Inter X...Y Contact C14 ..C15 .				3.18 Ang.
			1-x,y,1-z =	9_656 Check	
PLAT432_ALERT_2_G	Short Inter X...Y Contact C14 ..C15 .				3.18 Ang.
			1-z,y,1-x =	23_656 Check	
PLAT720_ALERT_4_G	Number of Unusual/Non-Standard Labels				1 Note
PLAT789_ALERT_4_G	Atoms with Negative _atom_site_disorder_group #				31 Check
PLAT860_ALERT_3_G	Number of Least-Squares Restraints				75 Note
PLAT883_ALERT_1_G	No Info/Value for _atom_sites_solution_primary .				Please Do !
PLAT910_ALERT_3_G	Missing # of FCF Reflection(s) Below Theta(Min).				4 Note
PLAT913_ALERT_3_G	Missing # of Very Strong Reflections in FCF				3 Note
PLAT984_ALERT_1_G	The O-f' = 0.0218 Deviates from the B&C-Value				0.0205 Check

PLAT984_ALERT_1_G The Zn-f' = -0.1382 Deviates from the B&C-Value -0.1517 Check
 PLAT985_ALERT_1_G The Zn-f" = 2.4327 Deviates from the B&C-Value 2.3734 Check

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

14 ALERT type 1 CIF construction/syntax error, inconsistent or missing data
 12 ALERT type 2 Indicator that the structure model may be wrong or deficient
 12 ALERT type 3 Indicator that the structure quality may be low
 34 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.

Validation response form

Please find below a validation response form (VRF) that can be filled in and pasted into your CIF.

```
# start Validation Reply Form
_vrf_THETM01_muf-91-6h-50pc
;
PROBLEM: The value of sine(theta_max)/wavelength is less than 0.550
RESPONSE: ...
;
```

_vrf_PLAT601_muf-91-6h-50pc

;

PROBLEM: Unit Cell Contains Solvent Accessible VOIDS of .

1608 Ang**3

RESPONSE: ...

;

end Validation Reply Form

PLATON version of 20/01/2022; check.def file version of 19/01/2022

Datablock muf-91-6h-50pc - ellipsoid plot

