

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: shelx_Sam_AH151

Bond precision: C-C = 0.0110 Å Wavelength=0.71073

Cell: a=9.6212(3) b=16.3828(4) c=15.2790(3)
 alpha=90 beta=91.989(2) gamma=90

Temperature: 120 K

	Calculated	Reported
Volume	2406.86(11)	2406.86(11)
Space group	P 21/n	P 21/n
Hall group	-P 2yn	-P 2yn
Moiety formula	C33 Cl3 N3 Nd2 O18 S6	C33 Cl3 N3 Nd2 O18 S6
Sum formula	C33 Cl3 N3 Nd2 O18 S6	C33 Cl3 N3 Nd2 O18 S6
Mr	1313.55	1313.55
Dx,g cm-3	1.813	1.812
Z	2	2
Mu (mm-1)	2.630	2.630
F000	1260.0	1260.0
F000'	1262.14	
h,k,lmax	11,19,18	11,19,18
Nref	4267	4251
Tmin,Tmax	0.854,0.949	0.946,1.000
Tmin'	0.854	

Correction method= # Reported T Limits: Tmin=0.946 Tmax=1.000
AbsCorr = MULTI-SCAN

Data completeness= 0.996 Theta(max)= 25.036

R(reflections)= 0.0553(3221) wR2(reflections)= 0.1375(4251)

S = 1.029 Npar= 343

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

PLAT220_ALERT_2_B Non-Solvent Resd 1 C Ueq(max)/Ueq(min) Range 8.8 Ratio

Author Response: It refers to the thermal ellipsoids of particular atoms in our structure.

PLAT241_ALERT_2_B High 'MainMol' Ueq as Compared to Neighbors of C4D Check

Author Response: ISOR restraints were applied to atom S22D in the disordered coordinated DMSO. The atom was still not ideally shaped, however, this does not indicate an incorrect atom-type assignment.

PLAT241_ALERT_2_B High 'MainMol' Ueq as Compared to Neighbors of C22D Check

Author Response: ISOR restraints were applied to atom S22D in the disordered coordinated DMSO. The atom was still not ideally shaped, however, this does not indicate an incorrect atom-type assignment.



Alert level C

PLAT213_ALERT_2_C Atom Old has ADP max/min Ratio 3.3 prolat

PLAT241_ALERT_2_C High 'MainMol' Ueq as Compared to Neighbors of Old Check

Author Response: ISOR restraints were applied to atom S22D in the disordered coordinated DMSO. The atom was still not ideally shaped, however, this does not indicate an incorrect atom-type assignment.

PLAT241_ALERT_2_C High 'MainMol' Ueq as Compared to Neighbors of Cl3D Check

Author Response: ISOR restraints were applied to atom S22D in the disordered coordinated DMSO. The atom was still not ideally shaped, however, this does not indicate an incorrect atom-type assignment.

PLAT242_ALERT_2_C Low 'MainMol' Ueq as Compared to Neighbors of C1 Check

PLAT342_ALERT_3_C Low Bond Precision on C-C Bonds 0.011 Ang.



Alert level G

PLAT003_ALERT_2_G Number of Uiso or Uij Restrained non-H Atoms ... 7 Report

PLAT004_ALERT_5_G Polymeric Structure Found with Maximum Dimension 1 Info

PLAT040_ALERT_1_G No H-atoms in this Carbon Containing Compound .. Please Check

PLAT083_ALERT_2_G SHELXL Second Parameter in WGHT Unusually Large 16.90 Why ?

PLAT171_ALERT_4_G The CIF-Embedded .res File Contains EADP Records 3 Report

PLAT177_ALERT_4_G The CIF-Embedded .res File Contains DELU Records 1 Report

PLAT186_ALERT_4_G The CIF-Embedded .res File Contains ISOR Records 1 Report

PLAT300_ALERT_4_G Atom Site Occupancy of Cl1 Constrained at 0.5 Check

PLAT300_ALERT_4_G Atom Site Occupancy of Cl4 Constrained at 0.5 Check

PLAT300_ALERT_4_G Atom Site Occupancy of C111	Constrained at	0.5	Check
PLAT300_ALERT_4_G Atom Site Occupancy of S1D	Constrained at	0.5	Check
PLAT300_ALERT_4_G Atom Site Occupancy of S2D	Constrained at	0.5	Check
PLAT300_ALERT_4_G Atom Site Occupancy of S11D	Constrained at	0.5	Check
PLAT300_ALERT_4_G Atom Site Occupancy of S12D	Constrained at	0.5	Check
PLAT300_ALERT_4_G Atom Site Occupancy of S21D	Constrained at	0.75	Check
PLAT300_ALERT_4_G Atom Site Occupancy of S22D	Constrained at	0.25	Check
PLAT300_ALERT_4_G Atom Site Occupancy of N100	Constrained at	0.5	Check
PLAT300_ALERT_4_G Atom Site Occupancy of N111	Constrained at	0.5	Check
PLAT300_ALERT_4_G Atom Site Occupancy of N400	Constrained at	0.5	Check
PLAT300_ALERT_4_G Atom Site Occupancy of C14D	Constrained at	0.5	Check
PLAT300_ALERT_4_G Atom Site Occupancy of C15D	Constrained at	0.5	Check
PLAT300_ALERT_4_G Atom Site Occupancy of C100	Constrained at	0.5	Check
PLAT300_ALERT_4_G Atom Site Occupancy of C111	Constrained at	0.5	Check
PLAT300_ALERT_4_G Atom Site Occupancy of C400	Constrained at	0.5	Check
PLAT301_ALERT_3_G Main Residue Disorder(Resd 1)		24%	Note
PLAT432_ALERT_2_G Short Inter X...Y Contact C14 ..C14D		3.20	Ang.
	$1/2-x, -1/2+y, 3/2-z =$	2_546	Check
PLAT432_ALERT_2_G Short Inter X...Y Contact C111 ..S2D		3.14	Ang.
	$1/2+x, 1/2-y, -1/2+z =$	4_665	Check
PLAT432_ALERT_2_G Short Inter X...Y Contact N400 ..C14D		2.94	Ang.
	$1/2-x, -1/2+y, 3/2-z =$	2_546	Check
PLAT432_ALERT_2_G Short Inter X...Y Contact C15D ..C15D		2.91	Ang.
	$-x, 1-y, 1-z =$	3_566	Check
PLAT773_ALERT_2_G Check long C-C Bond in CIF: C11 --C111		1.75	Ang.
PLAT860_ALERT_3_G Number of Least-Squares Restraints		42	Note

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

1 ALERT type 1 CIF construction/syntax error, inconsistent or missing data
 14 ALERT type 2 Indicator that the structure model may be wrong or deficient
 3 ALERT type 3 Indicator that the structure quality may be low
 20 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 19/10/2018; check.def file version of 15/10/2018

