

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: MgDy

Bond precision:	C-C = 0.0076 A	Wavelength=0.71073
Cell:	a=15.4531(3)	b=19.5753(4) c=14.0907(3)
	alpha=90	beta=109.082(2) gamma=90
Temperature:	150 K	
	Calculated	Reported
Volume	4028.21(15)	4028.19(15)
Space group	P 21/c	P 1 21/c 1
Hall group	-P 2ybc	-P 2ybc
Moiety formula	C50 H90 Dy2 Mg2 O22, 2(C6 H15 N)	C50 H90 Dy2 Mg2 O22, 2(C6 H15 N)
Sum formula	C62 H120 Dy2 Mg2 N2 O22	C62 H120 Dy2 Mg2 N2 O22
Mr	1619.22	1619.21
Dx,g cm-3	1.335	1.335
Z	2	2
Mu (mm-1)	1.920	1.920
F000	1676.0	1676.0
F000'	1676.10	
h,k,lmax	19,24,17	19,24,17
Nref	8241	8225
Tmin,Tmax	0.506,0.681	0.773,1.000
Tmin'	0.459	

Correction method= # Reported T Limits: Tmin=0.773 Tmax=1.000
AbsCorr = MULTI-SCAN

Data completeness= 0.998 Theta(max)= 26.372

R(reflections)= 0.0385(7274) wR2(reflections)= 0.0860(8225)

S = 1.093 Npar= 400

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

PLAT213_ALERT_2_C	Atom O3	has ADP max/min Ratio	3.2	prolat
PLAT213_ALERT_2_C	Atom C8	has ADP max/min Ratio	3.7	prolat
PLAT213_ALERT_2_C	Atom C23	has ADP max/min Ratio	3.7	prolat
PLAT213_ALERT_2_C	Atom C25	has ADP max/min Ratio	3.1	prolat
PLAT220_ALERT_2_C	Large Non-Solvent C	Ueq(max)/Ueq(min) Range	4.7	Ratio
PLAT220_ALERT_2_C	Large Non-Solvent O	Ueq(max)/Ueq(min) Range	4.0	Ratio
PLAT241_ALERT_2_C	High	Ueq as Compared to Neighbors for	03	Check
PLAT241_ALERT_2_C	High	Ueq as Compared to Neighbors for	09	Check
PLAT242_ALERT_2_C	Low	Ueq as Compared to Neighbors for	C2	Check
PLAT242_ALERT_2_C	Low	Ueq as Compared to Neighbors for	C12	Check
PLAT242_ALERT_2_C	Low	Ueq as Compared to Neighbors for	C22	Check
PLAT244_ALERT_4_C	Low	'Solvent' Ueq as Compared to Neighbors of	N1	Check



Alert level G

PLAT002_ALERT_2_G	Number of Distance or Angle Restraints on AtSite		4	Note
PLAT003_ALERT_2_G	Number of Uiso or Uij Restrained non-H Atoms ...		1	Report
PLAT007_ALERT_5_G	Number of Unrefined Donor-H Atoms		2	Report
PLAT083_ALERT_2_G	SHELXL Second Parameter in WGHT Unusually Large.		8.26	Why ?
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		3	Report
PLAT231_ALERT_4_G	Hirshfeld Test (Solvent) N1 -- C26 ..		7.3	su
PLAT231_ALERT_4_G	Hirshfeld Test (Solvent) N1 -- C29 ..		8.7	su
PLAT300_ALERT_4_G	Atom Site Occupancy of >C30 is Constrained at		0.623	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of <C32 is Constrained at		0.377	Check
PLAT302_ALERT_4_G	Anion/Solvent Disorder	Percentage =	14	Note
PLAT432_ALERT_2_G	Short Inter X...Y Contact O3 .. C30 ..		2.94	Ang.
PLAT720_ALERT_4_G	Number of Unusual/Non-Standard Labels		1	Note
PLAT764_ALERT_4_G	Overcomplete CIF Bond List Detected (Rep/Expd) .		1.18	Ratio
PLAT860_ALERT_3_G	Number of Least-Squares Restraints		9	Note

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

- 0 ALERT type 1 CIF construction/syntax error, inconsistent or missing data
15 ALERT type 2 Indicator that the structure model may be wrong or deficient
1 ALERT type 3 Indicator that the structure quality may be low
10 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*, 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.

