

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: I

Bond precision:	= 0.0000 A	Wavelength=0.41066
Cell:	a=6.9188(2)	b=6.9188(2) c=6.2124(3)
	alpha=90	beta=90 gamma=90
Temperature:	293 K	
	Calculated	Reported
Volume	297.39(2)	297.384(19)
Space group	I 41/a m d:	I 41/a m d
Hall group	I 4bw 2bw -	-I 4bd;-2
	As3.34 Dy0.29 Er0.15 Gd0.16	
Moiety formula	Ho0.14 Lu0.06 O16 P0.66 Y3 ?	
	Yb0.10	
	As3.34 Dy0.29 Er0.15 Gd0.16 O4 As0.82 P0.17 Y0.76	
Sum formula	Ho0.14 Lu0.06 O16 P0.66 Y3 Gd0.04 Er0.04 Ho0.03 Dy0.07	
	Yb0.10	Yb0.04 Lu0.02
Mr	943.43	238.00
Dx, g cm-3	5.268	5.316
Z	1	4
Mu (mm-1)	6.934	7.121
F000	426.1	426.0
F000'	427.02	
h, k, lmax	11, 11, 10	3, 11, 10
Nref	206	101
Tmin, Tmax		
Tmin'		
Correction method=	Not given	
Data completeness=	0.490	Theta(max)= 20.070

R(reflections)= 0.0207(91)

wR2(reflections)=

wR= 0.0256(101)

S = 1.570

Npar= 11

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

EXPT005_ALERT_1_A _exptl_crystal_description is missing

Crystal habit description.

The following tests will not be performed.

CRYSR_01

PLAT027_ALERT_3_A _diffrn_reflns_theta_full value (too) Low	8.16 Degree
PLAT029_ALERT_3_A _diffrn_measured_fraction_theta_full value Low .	0.790 Why?
PLAT183_ALERT_1_A Missing _cell_measurement_reflns_used Value	Please Do !
PLAT184_ALERT_1_A Missing _cell_measurement_theta_min Value	Please Do !
PLAT185_ALERT_1_A Missing _cell_measurement_theta_max Value	Please Do !
PLAT699_ALERT_1_A Missing _exptl_crystal_description Value	Please Do !

Alert level C

PLAT041_ALERT_1_C Calc. and Reported SumFormula Strings Differ Please Check

Calc: As0.83 Dy0.07 Er0.04 Gd0.04 Ho0.03 Lu0.02 O4 P0.17 Y0.75 Yb0

Rep.: O4 As0.82 P0.17 Y0.76 Gd0.04 Er0.04 Ho0.03 Dy0

.07 Yb0.04 Lu0.02

PLAT043_ALERT_1_C Calculated and Reported Mol. Weight Differ by .. 8.57 Check

PLAT051_ALERT_1_C Mu(calc) and Mu(CIF) Ratio Differs from 1.0 by . 2.62 %

PLAT052_ALERT_1_C Info on Absorption Correction Method Not Given Please Do !

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

PLAT127_ALERT_1_C Implicit Hall Symbol Inconsistent with Explicit -I 4bd;-2 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: As0.82 Dy.07 Er.04 Gd.04 Ho.03 L

Atom count from the _atom_site data: As0.8349 Dy.0735 Er.0379 Gd.0393

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: check formula stoichiometry or atom site occupancies.

From the CIF: _cell_formula_units_Z 4

From the CIF: _chemical_formula_sum O4 As0.82 P0.17 Y0.76 Gd0.04 Er0.0

TEST: Compare cell contents of formula and atom_site data

atom	Z*formula	cif sites	diff
O	16.00	16.00	0.00
As	3.28	3.34	-0.06
P	0.68	0.66	0.02
Y	3.04	3.01	0.03

Gd	0.16	0.16	0.00	
Er	0.16	0.15	0.01	
Ho	0.12	0.14	-0.02	
Dy	0.28	0.29	-0.01	
Yb	0.16	0.10	0.06	
Lu	0.08	0.06	0.02	
PLAT004_ALERT_5_G	Polymeric Structure Found with Maximum Dimension			3 Info
PLAT005_ALERT_5_G	No Embedded Refinement Details Found in the CIF			Please Do !
PLAT045_ALERT_1_G	Calculated and Reported Z Differ by a Factor ...			0.250 Check
PLAT092_ALERT_4_G	Check: Wavelength Given is not Cu,Ga,Mo,Ag,In Ka			0.41066 Ang.
PLAT120_ALERT_1_G	Reported I41/amd Inconsistent with Explicit			I41/amd Check
PLAT199_ALERT_1_G	Reported _cell_measurement_temperature (K)			293 Check
PLAT200_ALERT_1_G	Reported _diffn_ambient_temperature (K)			293 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of Lu	Constrained at		0.0151 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of Yb	Constrained at		0.0261 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of Er	Constrained at		0.0379 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of Ho	Constrained at		0.0348 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of Dy	Constrained at		0.0735 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of Gd	Constrained at		0.0393 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of Y	Constrained at		0.7517 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of As	Constrained at		0.8349 Check
PLAT300_ALERT_4_G	Atom Site Occupancy of P	Constrained at		0.1651 Check
PLAT301_ALERT_3_G	Main Residue Disorder(Resd 1)			79% Note
PLAT808_ALERT_5_G	No Parseable SHELXL Style Weighting Scheme Found			Please Check
PLAT811_ALERT_5_G	No ADDSYM Analysis: Too Many Excluded Atoms			! Info
PLAT883_ALERT_1_G	No Info/Value for _atom_sites_solution_primary .			Please Do !
PLAT950_ALERT_5_G	Calculated (ThMax) and CIF-Reported Hmax Differ			8 Units
PLAT966_ALERT_5_G	Note: Non-Standard (i.e. 2.0) OMIT Threshold of			3.0 Sig(I)
PLAT984_ALERT_1_G	The As-f' = 0.2878	Deviates from the B&C-Value		0.2864 Check
PLAT984_ALERT_1_G	The Dy-f' = -0.6999	Deviates from the B&C-Value		-0.6916 Check
PLAT984_ALERT_1_G	The Er-f' = -0.6332	Deviates from the B&C-Value		-0.6207 Check
PLAT984_ALERT_1_G	The Gd-f' = -0.7802	Deviates from the B&C-Value		-0.7780 Check
PLAT984_ALERT_1_G	The Ho-f' = -0.6648	Deviates from the B&C-Value		-0.6545 Check
PLAT984_ALERT_1_G	The Lu-f' = -0.5571	Deviates from the B&C-Value		-0.5418 Check
PLAT984_ALERT_1_G	The Y-f' = 0.2097	Deviates from the B&C-Value		0.2116 Check
PLAT984_ALERT_1_G	The Yb-f' = -0.5804	Deviates from the B&C-Value		-0.5650 Check
PLAT985_ALERT_1_G	The As-f'' = 0.7669	Deviates from the B&C-Value		0.7641 Check
PLAT985_ALERT_1_G	The Dy-f'' = 1.7292	Deviates from the B&C-Value		1.7162 Check
PLAT985_ALERT_1_G	The Er-f'' = 1.9570	Deviates from the B&C-Value		1.9421 Check
PLAT985_ALERT_1_G	The Gd-f'' = 1.5225	Deviates from the B&C-Value		1.5097 Check
PLAT985_ALERT_1_G	The Ho-f'' = 1.8406	Deviates from the B&C-Value		1.8265 Check
PLAT985_ALERT_1_G	The Lu-f'' = 2.3386	Deviates from the B&C-Value		2.3193 Check
PLAT985_ALERT_1_G	The Yb-f'' = 2.2053	Deviates from the B&C-Value		2.1884 Check

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- 7 **ALERT level A** = Most likely a serious problem - resolve or explain
 0 **ALERT level B** = A potentially serious problem, consider carefully
 9 **ALERT level C** = Check. Ensure it is not caused by an omission or oversight
 41 **ALERT level G** = General information/check it is not something unexpected
- 36 ALERT type 1 CIF construction/syntax error, inconsistent or missing data
 1 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
 11 ALERT type 4 Improvement, methodology, query or suggestion
 6 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 1 - ellipsoid plot

