

checkCIF/PLATON report

Structure factors have been supplied for datablock(s) c01a_1a_sq

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: c01a_1a_sq

Bond precision:	C-C = 0.0039 A	Wavelength=0.68890
Cell:	a=32.77108(17)	b=30.00687(16) c=40.3365(2)
	alpha=90	beta=96.1279(5) gamma=90
Temperature:	100 K	
	Calculated	Reported
Volume	39438.6(4)	39438.6(3)
Space group	C 2/c	C 1 2/c 1
Hall group	-C 2yc	-C 2yc
Moiety formula	C336 H262 Co8 N72, 2.84(C9 H6 O2), B2 F5, 7(B F4) [+ solvent]	C336 H262 Co8 N72, 3.5(B2 F8), B2 F5, 1.434(C18 H12 O4), 7[BF4]
Sum formula	C361.56 H279.04 B9 Co8 F33 N72 O5.68 [+ solvent]	C398.15 H435.65 B16 Co8 F64 N74.85 O39.35
Mr	6918.96	8759.54
Dx, g cm-3	1.165	1.475
Z	4	4
Mu (mm-1)	0.369	0.405
F000	14223.4	18141.0
F000'	14239.45	
h,k,lmax	47,43,58	47,43,58
Nref	62896	62798
Tmin,Tmax		0.959,1.000
Tmin'		

Correction method= # Reported T Limits: Tmin=0.959 Tmax=1.000
AbsCorr = EMPIRICAL

Data completeness= 0.998 Theta(max)= 29.947

R(reflections)= 0.0651(29550) wR2(reflections)= 0.2292(62798)

S = 0.957 Npar= 2330

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

PLAT026_ALERT_3_C	Ratio Observed / Unique Reflections (too) Low ..	47%	Check
PLAT077_ALERT_4_C	Unitcell Contains Non-integer Number of Atoms ..		Please Check
PLAT241_ALERT_2_C	High 'MainMol' Ueq as Compared to Neighbors of		C44B Check
PLAT260_ALERT_2_C	Large Average Ueq of Residue Including O14H	0.113	Check
PLAT260_ALERT_2_C	Large Average Ueq of Residue Including F92	0.157	Check
PLAT260_ALERT_2_C	Large Average Ueq of Residue Including F22	0.106	Check
PLAT260_ALERT_2_C	Large Average Ueq of Residue Including F32	0.110	Check
PLAT260_ALERT_2_C	Large Average Ueq of Residue Including F42	0.198	Check
PLAT260_ALERT_2_C	Large Average Ueq of Residue Including F32A	0.135	Check
PLAT905_ALERT_3_C	Negative K value in the Analysis of Variance ...	-6.951	Report
PLAT911_ALERT_3_C	Missing FCF Refl Between Thmin & STh/L= 0.600	47	Report
PLAT918_ALERT_3_C	Reflection(s) with I(obs) much Smaller I(calc) .	7	Check
PLAT978_ALERT_2_C	Number C-C Bonds with Positive Residual Density.	0	Info

● Alert level G

FORMU01_ALERT_1_G There is a discrepancy between the atom counts in the
_chemical_formula_sum and _chemical_formula_moiety. This is
usually due to the moiety formula being in the wrong format.
Atom count from _chemical_formula_sum: C398.1499 H435.6499 B16 Co8 F
Atom count from _chemical_formula_moiety:C361.8120 H279.2079 B16 Co8 F

FORMU01_ALERT_1_G There is a discrepancy between the atom counts in the
_chemical_formula_sum and the formula from the _atom_type* data.
Atom count from _chemical_formula_sum:C398.1499 H435.6499 B16 Co8 F64
Atom count from the _atom_type data:

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:C398.1499 H435.6499 B16 Co8 F64
Atom count from the _atom_site data: C361.5599 H279.0399 B9 Co8 F33 N

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 4
From the CIF: _chemical_formula_sum C398.15 H435.65 B16 Co8 F64 N74.85
TEST: Compare cell contents of formula and atom_site data

atom	Z*formula	cif sites	diff
C	1592.60	1446.24	146.36
H	1742.60	1116.16	626.44
B	64.00	36.00	28.00
Co	32.00	32.00	0.00
F	256.00	132.00	124.00
N	299.40	288.00	11.40
O	157.40	22.72	134.68

PLAT002_ALERT_2_G Number of Distance or Angle Restraints on AtSite 238 Note

PLAT003_ALERT_2_G Number of Uiso or Uij Restrained non-H Atoms ... 255 Report

PLAT014_ALERT_1_G N.O.K. _shelx_fab_checksum Found in CIF Please Check

PLAT041_ALERT_1_G Calc. and Reported SumFormula Strings Differ Please Check

PLAT042_ALERT_1_G Calc. and Reported MoietyFormula Strings Differ Please Check

PLAT072_ALERT_2_G SHELXL First Parameter in WGHT Unusually Large 0.14 Report

PLAT092_ALERT_4_G Check: Wavelength Given is not Cu,Ga,Mo,Ag,In Ka 0.68890 Ang.

PLAT142_ALERT_4_G s.u. on b - Axis Small or Missing 0.00016 Ang.

PLAT143_ALERT_4_G	s.u. on c - Axis Small or Missing	0.00020	Ang.
PLAT175_ALERT_4_G	The CIF-Embedded .res File Contains SAME Records	32	Report
PLAT178_ALERT_4_G	The CIF-Embedded .res File Contains SIMU Records	1	Report
PLAT187_ALERT_4_G	The CIF-Embedded .res File Contains RIGU Records	1	Report
PLAT244_ALERT_4_G	Low 'Solvent' Ueq as Compared to Neighbors of	B91	Check
PLAT244_ALERT_4_G	Low 'Solvent' Ueq as Compared to Neighbors of	B11	Check
PLAT244_ALERT_4_G	Low 'Solvent' Ueq as Compared to Neighbors of	B21	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of F42	0.5	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of F43	0.5	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of F44	0.5	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of F45	0.5	Check
PLAT300_ALERT_4_G	Atom Site Occupancy of B41	0.5	Check
PLAT302_ALERT_4_G	Anion/Solvent/Minor-Residue Disorder (Resd 3)	100%	Note
PLAT302_ALERT_4_G	Anion/Solvent/Minor-Residue Disorder (Resd 7)	100%	Note
PLAT302_ALERT_4_G	Anion/Solvent/Minor-Residue Disorder (Resd 8)	100%	Note
PLAT302_ALERT_4_G	Anion/Solvent/Minor-Residue Disorder (Resd 9)	100%	Note
PLAT304_ALERT_4_G	Non-Integer Number of Atoms in Resd 3	7.14	Check
PLAT304_ALERT_4_G	Non-Integer Number of Atoms in Resd 7	3.13	Check
PLAT304_ALERT_4_G	Non-Integer Number of Atoms in Resd 8	2.50	Check
PLAT304_ALERT_4_G	Non-Integer Number of Atoms in Resd 9	1.88	Check
PLAT343_ALERT_2_G	Unusual sp? Angle Range in Main Residue for	C24F	Check
PLAT432_ALERT_2_G	Short Inter X...Y Contact F34A ..C43E	2.97	Ang.
	x,y,z =	1_555	Check
PLAT606_ALERT_4_G	VERY LARGE Solvent Accessible VOID(S) in Structure	!	Info
PLAT779_ALERT_4_G	Suspect or Irrelevant (Bond) Angle(s) in CIF . #	747	Check
	B91 -F93 -B91 1.555 1.555 2.656	30.20	Deg.
PLAT779_ALERT_4_G	Suspect or Irrelevant (Bond) Angle(s) in CIF . #	748	Check
	B91 -F94 -B91 2.656 1.555 1.555	29.20	Deg.
PLAT789_ALERT_4_G	Atoms with Negative _atom_site_disorder_group #	17	Check
PLAT794_ALERT_5_G	Tentative Bond Valency for Co1 (II) .	1.54	Info
PLAT794_ALERT_5_G	Tentative Bond Valency for Co2 (II) .	1.56	Info
PLAT794_ALERT_5_G	Tentative Bond Valency for Co3 (II) .	1.51	Info
PLAT794_ALERT_5_G	Tentative Bond Valency for Co4 (II) .	1.50	Info
PLAT802_ALERT_4_G	CIF Input Record(s) with more than 80 Characters	7	Info
PLAT860_ALERT_3_G	Number of Least-Squares Restraints	5962	Note
PLAT869_ALERT_4_G	ALERTS Related to the Use of SQUEEZE Suppressed	!	Info
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).	1	Note
PLAT912_ALERT_4_G	Missing # of FCF Reflections Above STh/L= 0.600	52	Note
PLAT933_ALERT_2_G	Number of OMIT Records in Embedded .res File ...	32	Note
PLAT984_ALERT_1_G	The Co-f' = 0.3538 Deviates from the B&C-Value	0.3480	Check
PLAT985_ALERT_1_G	The Co-f" = 0.9121 Deviates from the B&C-Value	0.9239	Check

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

11 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
 6 ALERT type 3 Indicator that the structure quality may be low
 30 ALERT type 4 Improvement, methodology, query or suggestion
 4 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.

