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: Zn- O = 0.0139 A Wavelength=0.71070

Cell: a=4.1307(13) b=17.9042(2) c=10.3021(10)
alpha=90 beta=90.331(6) gamma=90
Temperature: 293 K

Calculated Reported
Volume 761.9(3) 761.9(3)
Space group ? ?
Hall group ? ?
Moiety formula Br4 O7.69 Sb4 Zn3 ?
Sum formula Br4 O7.69 Sb4 Zn3 Br1 O2 Sb1 Zn0.902
Mr 1125.85 292.60
Dx,g cm-3 2.454 5.101
Z 1 8
Mu (mm-1) 11.066 23.069
F000 495.5 1028.0
F000’ 493.28
h,k,lmax 0,0,0
Nref 26084
Tmin,Tmax
Tmin’

Correction method= Not given

Data completeness= Theta(max)= 0.000
R(reflections)= 0.0982( 6011) wR2(reflections)= wR= 0.1444( 26084)
S = 1.400 Npar= Npar = 361

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

SYM004_ALERT_1_A _symmetry_equiv_pos_as_xyz loop is missing.
The symmetry equivalent positions in xyz.
The following tests will not be performed.
CELLZ_01, CHEMW_03, REF LT_03, SYMMG_01, SYMMG_02

GEOM001_ALERT_1_A _geom_bond_atom_site_label_1 is missing
Label identifying the atom site 1.

GEOM003_ALERT_1_A _geom_bond_distance is missing
Distance between atom sites 1 and 2.

GEOM006_ALERT_1_A _geom_angle_atom_site_label_2 is missing
Label identifying the atom site 2.

GEOM007_ALERT_1_A _geom_angle_atom_site_label_3 is missing
Label identifying the atom site 3.

Alert level G

PLAT814_ALERT_5_G No Validation of (In)commensurate Structure CIFs

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.