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.

Datablock: I

```
Bond precision: Ag-S = 0.0017 A
                                         Wavelength=0.71073
                   a=9.637(2)
                                 b=12.946(3)
                                                      c=6.850(2)
Cell:
                                  beta=99.51(3)
                   alpha=90
                                                     gamma=90
                   297 K
Temperature:
                Calculated
                                          Reported
                842.9(4)
Volume
                                          842.9(4)
Space group
               Рс
                                          P 1 c 1
Hall group
              P -2yc
                                          P -2yc
Moiety formula Ag16 As7.78 S16 Sb0.22
                                          ?
Sum formula Ag16 As7.78 S16 Sb0.22
                                          Ag2 As0.973 S2 Sb0.027
               2848.36
                                          355.80
               5.611
                                          5.611
Dx,g cm-3
               1
                                          8
Mu (mm-1)
               17.806
                                          17.806
F000
               1275.9
                                          1276.0
F000'
               1264.14
h,k,lmax
               12,17,9
                                          12,17,9
Nref
                4095[ 2056]
                                          20050
Tmin, Tmax
                0.448,0.779
                                          0.467,1.000
Tmin'
                0.101
Correction method= # Reported T Limits: Tmin=0.467 Tmax=1.000
AbsCorr = MULTI-SCAN
Data completeness= 9.75/4.90 Theta(max) = 28.070
                                                    wR2 (reflections) =
R(reflections) = 0.0341(18767)
                                                    0.0870( 20050)
S = 1.758
                         Npar= 187
```

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
PLAT881_ALERT_1_A No Datum for _diffrn_reflns_av_R_equivalents ...
                                                                     Please Do !
🍭 Alert level B
                                                                      9.752
PLAT021_ALERT_4_B Ratio Unique / Expected Reflections too High ...
              # Unique Refl = 20050 / # Expected Refl =
                                                            2056
   Alert level C
PLAT041_ALERT_1_C Calc. and Reported SumFormula Strings Differ
             Calc: Ag16 As7.78 S16 Sb0.22
             Rep.: Ag2 As0.973 S2 Sb0.027
{\tt PLAT077\_ALERT\_4\_C~Unit~Cell~Contains~Non-integer~Number~of~Atoms~.}
                                                                     Please Check
PLAT241_ALERT_2_C High 'MainMol' Ueq as Compared to Neighbors of
                                                                        Ag3 Check
PLAT241_ALERT_2_C High
                        'MainMol' Ueq as Compared to Neighbors of
                                                                        Aq6 Check
                        'MainMol' Ueq as Compared to Neighbors of
PLAT241_ALERT_2_C High
                                                                        Ag7 Check
                        'MainMol' Ueq as Compared to Neighbors of
PLAT241_ALERT_2_C High
                                                                        Ag8 Check
                        'MainMol' Ueq as Compared to Neighbors of
PLAT242_ALERT_2_C Low
                                                                         S2 Check
PLAT242_ALERT_2_C Low
                        'MainMol' Ueq as Compared to Neighbors of
                                                                         S8 Check
Alert level G
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 \dots
                                                                     0.125 Check
PLAT068_ALERT_1_G Reported F000 Differs from Calcd (or Missing)...
                                                                     Please Check
PLAT301_ALERT_3_G Main Residue Disorder ......(Resd 1)
                                                                       20% Note
PLAT794_ALERT_5_G Tentative Bond Valency for Ag4
                                                                       1.07 Info
                                                    (I)
                                                               .
PLAT794_ALERT_5_G Tentative Bond Valency for Ag5
                                                                       0.98 Info
PLAT883_ALERT_1_G Absent Datum for _atom_sites_solution_primary ..
                                                                     Please Do !
PLAT966_ALERT_5_G Note: Non-Standard (i.e. 2.0) OMIT Threshold of
                                                                        3.0 Sig(I)
```

- 1 **ALERT level A** = Most likely a serious problem resolve or explain
- 1 ALERT level B = A potentially serious problem, consider carefully
- 8 **ALERT level C** = Check. Ensure it is not caused by an omission or oversight
- 9 ALERT level G = General information/check it is not something unexpected
- 5 ALERT type 1 CIF construction/syntax error, inconsistent or missing data
- 6 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
- 2 ALERT type 4 Improvement, methodology, query or suggestion
- 5 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.

Validation response form

Please find below a validation response form (VRF) that can be filled in and pasted into your CIF.

```
# start Validation Reply Form
_vrf_PLAT881_I
PROBLEM: No Datum for _diffrn_reflns_av_R_equivalents ... Please Do !
RESPONSE: ...
_vrf_PLAT021_I
PROBLEM: Ratio Unique / Expected Reflections too High ...
                                                           9.752
RESPONSE: ...
_vrf_PLAT041_I
PROBLEM: Calc. and Reported SumFormula Strings Differ
                                                           Please Check
RESPONSE: ...
_vrf_PLAT077_I
PROBLEM: Unit Cell Contains Non-integer Number of Atoms . Please Check
RESPONSE: ...
_vrf_PLAT241_I
```

```
;
PROBLEM: High 'MainMol' Ueq as Compared to Neighbors of Ag3 Check
RESPONSE: ...
;
_vrf_PLAT242_I
;
PROBLEM: Low 'MainMol' Ueq as Compared to Neighbors of S2 Check
RESPONSE: ...
;
# end Validation Reply Form
```

PLATON version of 11/11/2024; check.def file version of 11/11/2024

Datablock I - ellipsoid plot

