checkCIF/PLATON report

Structure factors have been supplied for datablock(s) skm312a_0m_a_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.

Datablock: skm312a_0m_a_sq

```
Wavelength=0.71073
Bond precision: C-C = 0.0018 A
Cell:
                  a=11.1145(3)
                                   b=13.7766(3)
                                                        c=12.9339(3)
                                   beta=94.722(1)
                  alpha=90
                                                        gamma=90
Temperature:
                  170 K
                 Calculated
                                             Reported
Volume
                 1973.72(8)
                                             1973.72(8)
Space group
                P 21/n
                                             P 21/n
                                             −P 2yn
Hall group
                −P 2yn
Moiety formula C21 H20 N2 O S [+ solvent] C21 H20 N2 O S
Sum formula
                C21 H20 N2 O S [+ solvent] C21 H20 N2 O S
                 348.45
                                             348.45
Mr
                                             1.173
Dx,g cm-3
                 1.173
                 0.174
                                             0.174
Mu (mm-1)
F000
                 736.0
                                             736.0
F000'
                 736.75
h, k, lmax
                 15, 19, 18
                                             15, 19, 18
Nref
                 6085
                                             5994
Tmin, Tmax
                 0.941,0.967
                                             0.920,0.970
Tmin'
                 0.937
Correction method= # Reported T Limits: Tmin=0.920 Tmax=0.970
AbsCorr = NUMERICAL
Data completeness= 0.985
                                     Theta (max) = 30.626
                                                        wR2 (reflections) =
R(reflections) = 0.0439(5057)
                                                        0.1449(5994)
S = 1.070
                           Npar= 229
```

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

PLAT910_ALERT_3_C Missing # of FCF Reflection(s) Below Theta(Min).	5 Note
PLAT911_ALERT_3_C Missing FCF Refl Between Thmin & STh/L= 0.600	67 Report
PLAT913_ALERT_3_C Missing # of Very Strong Reflections in FCF	43 Note
PLAT918 ALERT 3 C Reflection(s) with I(obs) much Smaller I(calc) .	2 Check

Alert level G

PLAT605_ALERT_4_G Largest Solvent Accessible VOID in the Structure	116 A**3
PLAT793_ALERT_4_G Model has Chirality at C7 (Centro SPGR)	R Verify
PLAT869_ALERT_4_G ALERTS Related to the Use of SQUEEZE Suppressed	! Info
PLAT912_ALERT_4_G Missing # of FCF Reflections Above STh/L= 0.600	20 Note
PLAT933_ALERT_2_G Number of HKL-OMIT Records in Embedded .res File	14 Note
PLAT978_ALERT_2_G Number C-C Bonds with Positive Residual Density.	19 Info

- 0 **ALERT level A** = Most likely a serious problem resolve or explain
- 0 **ALERT level B** = A potentially serious problem, consider carefully
- 4 **ALERT level C** = Check. Ensure it is not caused by an omission or oversight
- 6 ALERT level ${\bf G}$ = General information/check it is not something unexpected
- 0 ALERT type 1 CIF construction/syntax error, inconsistent or missing data
- 2 ALERT type 2 Indicator that the structure model may be wrong or deficient
- 4 ALERT type 3 Indicator that the structure quality may be low
- 4 ALERT type 4 Improvement, methodology, query or suggestion
- O 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.

PLATON version of 28/11/2022; check.def file version of 28/11/2022

