# checkCIF/PLATON report

Structure factors have been supplied for datablock(s) exp\_3472\_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: exp\_3472\_sq

```
Bond precision: C-C = 0.0020 A
                                           Wavelength=1.54184
Cell:
                a=12.2357(4) b=13.4613(5)
                                                     c=14.9649(5)
                alpha=78.284(3) beta=77.397(3)
                                                     gamma = 76.902(3)
Temperature:
                100 K
                Calculated
                                            Reported
Volume
                2312.08(14)
                                            2312.08(15)
Space group
                P -1
                                            P -1
                -P 1
                                            -P 1
Hall group
Moiety formula C55 H43 B N2 [+ solvent]
                                            C55 H43 B N2
Sum formula
                C55 H43 B N2 [+ solvent]
                                            C55 H43 B N2
                742.72
                                            742.72
Mr
                                            1.067
Dx,g cm-3
                1.067
                2
                                            2
Mu (mm-1)
                0.464
                                            0.464
                784.0
F000
                                            784.0
F000'
                786.02
h, k, lmax
                14,16,17
                                            14, 16, 17
Nref
                8160
                                            8162
Tmin, Tmax
                0.911,0.911
                                            0.894,1.000
Tmin'
                0.911
Correction method= # Reported T Limits: Tmin=0.894 Tmax=1.000
AbsCorr = MULTI-SCAN
Data completeness= 1.000
                                    Theta (max) = 66.600
                                                      wR2 (reflections) =
R(reflections) = 0.0406(7198)
                                                       0.1038(8162)
S = 1.037
                          Npar= 529
```

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

DIFMX02\_ALERT\_1\_C The maximum difference density is > 0.1\*ZMAX\*0.75The relevant atom site should be identified.

PLAT094_ALERT_2_C Ratio of Maximum / Minimum Residual Density	2.71 Report
PLAT097_ALERT_2_C Large Reported Max. (Positive) Residual Density	0.69 eA-3
PLAT906 ALERT 3 C Large K Value in the Analysis of Variance	2.436 Check

## Alert level G

PLAT154_ALERT_1_G The s.u.'s on the Cell Angles are Equal(Note)	0.003 Degree
PLAT606_ALERT_4_G Solvent Accessible VOID(S) in Structure	! Info
PLAT869_ALERT_4_G ALERTS Related to the Use of SQUEEZE Suppressed	! Info
PLAT909_ALERT_3_G Percentage of I>2sig(I) Data at Theta(Max) Still	84% Note
PLAT941_ALERT_3_G Average HKL Measurement Multiplicity	1.8 Low
PLAT978_ALERT_2_G Number C-C Bonds with Positive Residual Density.	16 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 G = General information/check it is not something unexpected
- 2 ALERT type 1 CIF construction/syntax error, inconsistent or missing data
- 3 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
- 2 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 06/07/2023; check.def file version of 30/06/2023

