5 Info

## Datablock: 3

```
Bond precision: C-C = 0.0053 A
                                                     Wavelength=0.71073
Cell: a=7.2915(3) b=28.1816(13) c=8.9402(5)
             alpha=90
                             beta=112.196(6) gamma=90
Temperature 150 K
                    Calculated
                                                       Reported
                    1700.95(16)
                                                       1700.95(16)
Volume
Space group P 21/n P
Hall group -P 2yn -I
Moiety formula C24 H36 C12 Cu2 N8 O2, 2(B F4) ?
Sum formula C24 H36 B2 C12 Cu2 F8 N8 O2 C2
Mr 840 23
                                                       P 21/n
                                                       -P 2yn
                                                      C24 H36 B2 C12 Cu2 F8 N8 O2
                   840.23
                                                       840.21
                   1.640
Dx,g cm-3
                                                       1.640
Mu (mm-1)
               1.489
                                                       1.489
F000
                   852.0
                                                       852.0
                  854.27
9,36,11
F000'
h,k,lmax
                                                      9,36,11
Nref
                   3912
                                                      3910
                0.494,0.742
Tmin, Tmax
                                                      0.563,1.000
Tmin'
                   0.301
Correction method= # Reported T Limits: Tmin=0.563
Tmax=1.000 AbsCorr = MULTI-SCAN
Data completeness= 0.999 Theta(max)= 27.485 R(reflections)= 0.0490(3365) wR2(reflections)= 0.1256(3910)
S = 1.107
                      Npar= 221
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
PLAT094 ALERT 2 C Ratio of Maximum / Minimum Residual Density ....
                                                                               2.15 Report
<u>PLAT906 ALERT 3 C</u> Large K Value in the Analysis of Variance .....
                                                                               3.253 Check
Alert level G
PLAT063 ALERT 4 G Crystal Size Likely too Large for Beam Size ....
                                                                               0.80 mm
PLAT244 ALERT 4 G Low 'Solvent' Ueq as Compared to Neighbors of
                                                                                 B1 Check
PLAT790 ALERT 4 G Centre of Gravity not Within Unit Cell: Resd. #
                                                                                   2 Note
               B F4
PLAT978 ALERT 2 G Number C C Parks with a point value of Cul (II) . 2.22 Info PLAT978 ALERT 2 G Number C C Parks with a point (II) . 2.22 Info Please Do !
PLAT794 ALERT 5 G Tentative Bond Valency for Cul
                                                                               2.22 Info
```

```
0 ALERT level A = Most likely a serious problem - resolve or explain
0 ALERT level B = A potentially serious problem, consider carefully
2 ALERT level C = Check. Ensure it is not caused by an omission or oversight
7 ALERT level G = General information/check it is not something unexpected
1 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
2 ALERT type 3 Indicator that the structure quality may be low
3 ALERT type 4 Improvement, methodology, query or suggestion
1 ALERT type 5 Informative message, check
```

PLAT978 ALERT 2 G Number C-C Bonds with Positive Residual Density.

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 <u>full publication checks</u> 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 03/05/2019; check.def file version of 29/04/2019 **Datablock 3** - ellipsoid plot

