## Datablock: moc40

```
Bond precision:
                                             Wavelength=0.71073
                  C-C = 0.0037 A
Cell:
          a=7.7916(6) b=8.9848(6) c=11.9955(7)
           alpha=104.394(6)beta=106.396(6) gamma=104.368(6)
Temperature 150 K
                 Calculated
                                              Reported
                 733.10(10)
Volume
                                              733.10(9)
Space group
                 P -1
                                              P -1
-P 1
                                             C26 H22 C14 Cu2 N6 O8
                                             C26 H22 C14 Cu2 N6 O8
                 815.40
                                              815.37
                1.847
                                              1.847
Dx,g cm-3
Mu (mm-1)
                 1.877
                                              1.877
F000
                 410.0
                                              410.0
               411.37
10,11,15
F000'
h,k,lmax
                                              10,11,15
Nref
                3357
                                              3355
               0.893,0.910
Tmin, Tmax
                                              0.831,1.000
Tmin'
                0.687
Correction method= # Reported T Limits: Tmin=0.831
Tmax=1.000 AbsCorr = MULTI-SCAN
Data completeness= 0.999 Theta(max) = 27.476
                                        wR2(reflections)=
R(reflections) = 0.0339(2769)
                                        0.0748 (3355)
S = 1.070
                    Npar= 210
```

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 G
```

```
PLAT154 ALERT 1 G The s.u.'s on the Cell Angles are Equal .. (Note) 0.006 Degree
PLAT794 ALERT 5 G Tentative Bond Valency for Cu1 (II) . 2.12 Info
PLAT910 ALERT 3 G Missing # of FCF Reflection(s) Below Theta(Min). 1 Note
PLAT912 ALERT 4 G Missing # of FCF Reflections Above STh/L= 0.600 1 Note
PLAT941 ALERT 3 G Average HKL Measurement Multiplicity ....... 1.9 Low
PLAT978 ALERT 2 G Number C-C Bonds with Positive Residual Density. 4 Info
```

```
O ALERT level A = Most likely a serious problem - resolve or explain
O ALERT level B = A potentially serious problem, consider carefully
O ALERT level C = Check. Ensure it is not caused by an omission or oversight
ALERT level G = General information/check it is not something unexpected

1 ALERT type 1 CIF construction/syntax error, inconsistent or missing data
1 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
1 ALERT type 4 Improvement, methodology, query or suggestion
1 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.

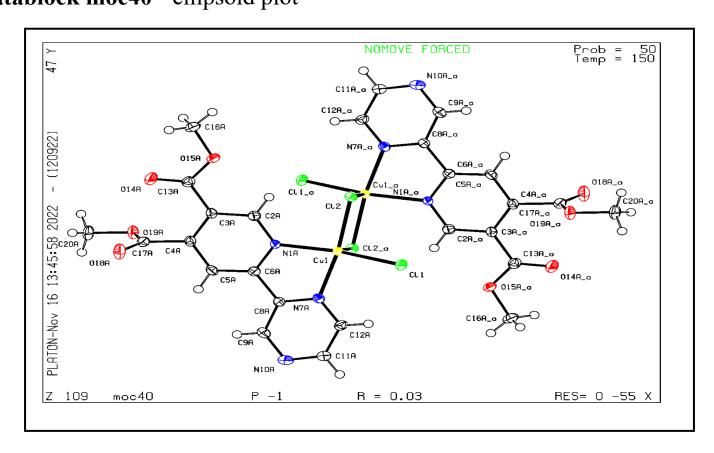
## Publication of your CIF in IUCr journals

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 <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 12/09/2022; check.def file version of 09/08/2022 **Datablock moc40** - ellipsoid plot



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