checkCIF/PLATON report

Structure factors have been supplied for datablock(s) 20250498a_auto

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: 20250498a_auto

```
Bond precision: C-C = 0.0068 A
                                         Wavelength=1.54184
                a=13.1194(4)
Cell:
                                b=12.8734(4)
                                                    c=10.0344(5)
                                beta=100.177(4)
                alpha=90
                                                   gamma=90
                293 K
Temperature:
               Calculated
                                          Reported
Volume
               1668.06(11)
                                          1668.06(11)
Space group
               Сс
                                          C 1 c 1
Hall group
             C -2yc
                                          C -2yc
Moiety formula C20 H16 N2 O4
                                         C20 H16 N2 O4
Sum formula C20 H16 N2 O4
                                          C20 H16 N2 O4
Mr
               348.35
                                          348.35
               1.387
                                          1.387
Dx,g cm-3
               4
Mu (mm-1)
              0.808
                                          0.808
F000
               728.0
                                          728.0
F000'
               730.37
h,k,lmax
              16,15,12
                                         15,15,12
Nref
               3221[ 1614]
                                          2199
Tmin, Tmax
              0.890,0.922
                                          0.895,1.000
Tmin'
               0.886
Correction method= # Reported T Limits: Tmin=0.895 Tmax=1.000
AbsCorr = MULTI-SCAN
Data completeness= 1.36/0.68 Theta(max)= 70.946
                                                    wR2 (reflections) =
R(reflections) = 0.0457(2054)
                                                    0.1343 (2199)
S = 1.057
                         Npar= 237
```

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.

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Alert level C
PLAT089 ALERT 3 C P
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PLAT089_ALERT_3_C Poor Data / Parameter Ratio (Zmax < 18) 6.72 Note
PLAT241_ALERT_2_C High 'MainMol' Ueq as Compared to Neighbors of C8 Check
PLAT241_ALERT_2_C High 'MainMol' Ueq as Compared to Neighbors of C18 Check
PLAT340_ALERT_3_C Low Bond Precision on C-C Bonds 0.00684 Ang.

Alert level G

PLAT111_ALERT_2_G ADDSYM Detects New (Pseudo) Centre of Symmetry . 92 %Fit PLAT113_ALERT_2_G ADDSYM Suggests Possible Pseudo/New Space Group C2/c Check Check Model Parameter Symmetry for Reflection Data Support PLAT199_ALERT_1_G Reported _cell_measurement_temperature (K) 293 Check PLAT200_ALERT_1_G Reported __diffrn_ambient_temperature (K) 293 Check PLAT398_ALERT_2_G Deviating C-O-C Angle From 120 for O2 109.8 Degree PLAT792_ALERT_1_G Model has Chirality at C3 R Verify (Polar SpGr) PLAT792_ALERT_1_G Model has Chirality at C12 (Polar SpGr) S Verify ${\tt PLAT883_ALERT_1_G~Absent~Datum~for~_atom_sites_solution_primary~..}$ Please Do ! PLAT912_ALERT_4_G Missing # of FCF Reflections Above STh/L= 0.600 21 Note PLAT915_ALERT_3_G No Flack x Check Done: Low Friedel Pair Coverage 38 % PLAT941_ALERT_3_G Average HKL Measurement Multiplicity 3.9 Low PLAT969_ALERT_5_G The 'Henn et al.' R-Factor-gap value 3.910 Note Predicted wR2: Based on SigI**2 3.43 or SHELX Weight 12.70 PLAT978_ALERT_2_G Number C-C Bonds with Positive Residual Density. 2 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
- 13 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
 - 4 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 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 02/02/2025; check.def file version of 02/02/2025

