

# **Supporting Information**

for

Catalyst-free assembly of giant tris(heteroaryl)methanes: synthesis of novel pharmacophoric triads and model sterically crowded tris(heteroaryl/aryl)methyl cation salts

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**CIF** report for 9{4,7,1}

## checkCIF/PLATON report

Structure factors have been supplied for datablock(s) lg182\_2\_final\_a

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: lg182\_2\_final\_a

```
Bond precision: C-C = 0.0042 A
                                       Wavelength=0.71073
Cell:
             a=9.6968(10) b=12.2099(16) c=13.6876(14)
             alpha=90.092(4) beta=99.906(2) gamma=112.945(2)
Temperature: 200 K
               Calculated
                                        Reported
               1465.9(3)
Volume
                                        1465.9(3)
Space group
              P -1
                                        P -1
Hall group
               -P 1
                                        -P 1
Moiety formula C32 H27 Cl N2 O4, C2 H3 N C32 H27 Cl N2 O4, C2 H3 N
Sum formula
             C34 H30 Cl N3 O4
                                        C34 H30 Cl N3 O4
Mr
               580.06
                                        580.06
               1.314
                                        1.314
Dx,g cm-3
               2
Ζ
                                        2
Mu (mm-1)
               0.174
                                        0.174
F000
               608.0
                                        608.0
F000′
               608.55
h,k,lmax
               11,14,16
                                        11,14,16
                                        5207
Nref
               5208
               0.990,0.991
                                        0.521,0.562
Tmin,Tmax
Tmin'
               0.983
Correction method= # Reported T Limits: Tmin=0.521 Tmax=0.562
AbsCorr = MULTI-SCAN
Data completeness= 1.000
                                Theta(max) = 25.048
R(reflections) = 0.0490( 3206) wR2(reflections) = 0.1100( 5207)
S = 1.017
                         Npar= 383
```

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
                     'Solvent' Ueq as Compared to Neighbors of
PLAT244_ALERT_4_C Low
                                                                  C34 Check
PLAT340 ALERT 3 C Low Bond Precision on C-C Bonds ...... 0.00415 Ang.
  Alert level G
PLAT007_ALERT_5_G Number of Unrefined Donor-H Atoms .....
                                                                        1 Report
PLAT793_ALERT_4_G Model has Chirality at C1 (Centro SPGR)
                                                                        R Verify
PLAT909_ALERT_3_G Percentage of I>2sig(I) Data at Theta(Max) Still
                                                                      36% Note
PLAT910_ALERT_3_G Missing # of FCF Reflection(s) Below Theta(Min).
                                                                        2 Note
PLAT933 ALERT 2 G Number of OMIT Records in Embedded .res File ...
                                                                        2 Note
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
  2 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
  O 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
  3 ALERT type 3 Indicator that the structure quality may be low
  2 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.

