checkCIF/PLATON report

No syntax errors found.   CIF dictionary   Interpreting this report

Datablock: mrmnd

Bond precision: C-C = 0.0095 A   Wavelength=0.71073

Cell: a=17.7892(3)   b=15.8129(3)   c=28.4025(5)
                   alpha=90   beta=90   gamma=90

Temperature: 120 K

	Calculated	Reported
Volume	7989.6(2)	7989.6(2)
Space group	P b c a	P b c a
Hall group	-P 2ac 2ab	-P 2ac 2ab
Moiety formula	C30 H46 O3 P2 Pt S, 2(C6 D6)	C30 H46 O3 P2 Pt S, 2(C6 D6)
Sum formula	C42 H46 D12 O3 P2 Pt S	C42 H46 D12 O3 P2 Pt S
Mr	912.00	912.04
Dx,g cm-3	1.516	1.516
Z	8	8
Mu (mm-1)	3.682	3.682
F000	3664.0	3664.0
F000'	3654.37	
h,k,lmax	27,24,43	27,24,43
Nref	14996	14395
Tmin,Tmax	0.498,0.576	0.767,1.000
Tmin'	0.457	

Correction method= MULTI-SCAN

Data completeness= 0.960   Theta(max)= 32.941

R(reflections)= 0.0638(11302)   wR2(reflections)= 0.1718(14395)

S = 1.101   Npar= Npar = 456

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

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PLAT094_ALERT_2_C Ratio of Maximum / Minimum Residual Density ....     2.49 Why ?
PLAT097_ALERT_2_C Large Reported Max.  (Positive) Residual Density     7.61 eA-3
PLAT250_ALERT_2_C Large U3/U1 Ratio for Average U(i,j) Tensor ....      2.1 Note
PLAT331_ALERT_2_C Small Average Phenyl   C-C Dist. C62    -C266        1.36 Ang.
PLAT342_ALERT_3_C Low Bond Precision on  C-C Bonds ..............     0.0095 Ang.
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Alert level G

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PLAT005_ALERT_5_G No _iucr_refine_instructions_details  in the CIF     Please Do !
PLAT083_ALERT_2_G SHELXL Second Parameter in WGHT Unusually Large.     82.64 Why ?
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  0 ALERT level A = Most likely a serious problem - resolve or explain
  0 ALERT level B = A potentially serious problem, consider carefully
  5 ALERT level C = Check. Ensure it is not caused by an omission or oversight
  2 ALERT level G = General information/check it is not something unexpected

  0 ALERT type 1 CIF construction/syntax error, inconsistent or missing data
  5 ALERT type 2 Indicator that the structure model may be wrong or deficient
  1 ALERT type 3 Indicator that the structure quality may be low
  0 ALERT type 4 Improvement, methodology, query or suggestion
  1 ALERT type 5 Informative message, check
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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*, 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 05/02/2014; check.def file version of 05/02/2014

NOMOVE FORCED

Prob = 50
Temp = 120

46 Y

PLATON-May 7 04:14:09 2014 - (100912)

Z 71   mrmnd   P b c a   R = 0.06   RES=   0   56 X