

checkCIF/PLATON report

You have not supplied any structure factors. As a result the full set of tests cannot be run.

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.

No syntax errors found. CIF dictionary Interpreting this report

Datablock: I

Bond precision: C-C = 0.0153 A

Wavelength=1.54178

Cell: a=16.7510(5) b=17.2703(7) c=18.8355(6)
 alpha=109.136(3) beta=97.750(3) gamma=108.799(3)
Temperature: 123 K

	Calculated	Reported
Volume	4693.8(3)	4693.8(3)
Space group	P -1	P -1
Hall group	-P 1	-P 1
Moiety formula	2(C16 Al F36 O4), 2(C40 H60 Cu Mo4 P12), 5(C7 H8)	2(C40 H60 Cu Mo4 P12), 2(Al C16 F36 O4), 5(C7 H8)
Sum formula	C147 H160 Al2 Cu2 F72 Mo8 O8 P24	C147 H160 Al2 Cu2 F72 Mo8 O8 P24
Mr	5114.62	5114.58
Dx, g cm ⁻³	1.809	1.809
Z	1	1
Mu (mm ⁻¹)	7.604	7.604
F000	2534.0	2534.0
F000'	2545.21	
h,k,lmax	21,21,23	20,21,23
Nref	19458	18870
Tmin,Tmax	0.453,0.566	0.759,1.000
Tmin'	0.283	

Correction method= MULTI-SCAN

Data completeness= 0.970

Theta(max)= 75.433

R(reflections)= 0.0665(16507)

wR2(reflections)= 0.1867(18870)

S = 1.053

Npar= 1261

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 B

Crystal system given = triclinic

PLAT220_ALERT_2_B Large Non-Solvent F Ueq(max)/Ueq(min) Range 10.0 Ratio

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. The terminal fluorine atoms are strongly affected by this.

PLAT230_ALERT_2_B Hirshfeld Test Diff for F8 -- C44 .. 9.0 su

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. The terminal fluorine atoms are strongly affected by this.

PLAT230_ALERT_2_B Hirshfeld Test Diff for F18 -- C48 .. 18.5 su

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. The terminal fluorine atoms are strongly affected by this.

PLAT230_ALERT_2_B Hirshfeld Test Diff for C41 -- C42 .. 9.2 su

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. The terminal fluorine atoms are strongly affected by this.

PLAT230_ALERT_2_B Hirshfeld Test Diff for C45 -- C47 .. 8.6 su

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. The terminal fluorine atoms are strongly affected by this.

PLAT241_ALERT_2_B High Ueq as Compared to Neighbors for C44 Check

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. C44 is part of a CF3 group.

PLAT242_ALERT_2_B Low Ueq as Compared to Neighbors for C41 Check

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. C41 is a tertiary atom of a tert-butyl group.

PLAT250_ALERT_2_B Large U3/U1 Ratio for Average U(i,j) Tensor 5.1 Note
PLAT332_ALERT_2_B Large Phenyl C-C Range C64 -C69 0.28 Ang.

Author Response: This is a free toluene molecule in the unit cell.

PLAT362_ALERT_2_B Short C(sp3)-C(sp2) Bond C57 - C63 ... 1.31 Ang.

● **Alert level C**

PLAT213_ALERT_2_C Atom F11 has ADP max/min Ratio 3.2 prolat
PLAT213_ALERT_2_C Atom F50A has ADP max/min Ratio 3.2 prolat
PLAT213_ALERT_2_C Atom F51A has ADP max/min Ratio 3.1 prolat
PLAT213_ALERT_2_C Atom F51C has ADP max/min Ratio 3.1 prolat
PLAT213_ALERT_2_C Atom F54C has ADP max/min Ratio 4.0 prolat
PLAT213_ALERT_2_C Atom F55A has ADP max/min Ratio 4.0 prolat
PLAT213_ALERT_2_C Atom F55C has ADP max/min Ratio 3.1 oblate
PLAT213_ALERT_2_C Atom F56A has ADP max/min Ratio 3.3 prolat
PLAT213_ALERT_2_C Atom F56C has ADP max/min Ratio 3.3 prolat
PLAT213_ALERT_2_C Atom C44 has ADP max/min Ratio 3.7 prolat
PLAT213_ALERT_2_C Atom C14 has ADP max/min Ratio 3.1 prolat
PLAT213_ALERT_2_C Atom C18 has ADP max/min Ratio 3.7 prolat
PLAT213_ALERT_2_C Atom C19 has ADP max/min Ratio 3.3 prolat
PLAT220_ALERT_2_C Large Non-Solvent C Ueq(max)/Ueq(min) Range 4.8 Ratio

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. The terminal fluorine atoms are strongly affected by this.

PLAT220_ALERT_2_C Large Non-Solvent C Ueq(max)/Ueq(min) Range 5.4 Ratio

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. The terminal fluorine atoms are strongly affected by this.

PLAT222_ALERT_3_C Large Non-Solvent H Uiso(max)/Uiso(min) .. 4.2 Ratio
PLAT230_ALERT_2_C Hirshfeld Test Diff for F1 -- C42 .. 6.8 su

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. The terminal fluorine atoms are strongly affected by this.

PLAT230_ALERT_2_C Hirshfeld Test Diff for F4 -- C43 .. 6.8 su

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. The terminal fluorine atoms are strongly affected by this.

PLAT230_ALERT_2_C Hirshfeld Test Diff for F11 -- C46 .. 5.1 su

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. The terminal fluorine atoms are strongly affected by this.

PLAT234_ALERT_4_C	Large	Hirshfeld	Difference	F3	--	C42	..	0.18	Ang.
PLAT234_ALERT_4_C	Large	Hirshfeld	Difference	F6	--	C43	..	0.18	Ang.
PLAT234_ALERT_4_C	Large	Hirshfeld	Difference	F13	--	C47	..	0.18	Ang.
PLAT234_ALERT_4_C	Large	Hirshfeld	Difference	F14	--	C47	..	0.17	Ang.
PLAT234_ALERT_4_C	Large	Hirshfeld	Difference	F16	--	C48	..	0.17	Ang.
PLAT234_ALERT_4_C	Large	Hirshfeld	Difference	F50C	--	C50	..	0.23	Ang.
PLAT234_ALERT_4_C	Large	Hirshfeld	Difference	F51A	--	C51	..	0.18	Ang.
PLAT234_ALERT_4_C	Large	Hirshfeld	Difference	F51B	--	C51	..	0.20	Ang.
PLAT234_ALERT_4_C	Large	Hirshfeld	Difference	F51C	--	C51	..	0.22	Ang.
PLAT234_ALERT_4_C	Large	Hirshfeld	Difference	F52C	--	C52	..	0.18	Ang.
PLAT234_ALERT_4_C	Large	Hirshfeld	Difference	F54A	--	C54	..	0.22	Ang.
PLAT234_ALERT_4_C	Large	Hirshfeld	Difference	F55B	--	C55	..	0.24	Ang.
PLAT234_ALERT_4_C	Large	Hirshfeld	Difference	F55C	--	C55	..	0.16	Ang.
PLAT234_ALERT_4_C	Large	Hirshfeld	Difference	F56B	--	C56	..	0.25	Ang.
PLAT234_ALERT_4_C	Large	Hirshfeld	Difference	C41	--	C44	..	0.24	Ang.
PLAT234_ALERT_4_C	Large	Hirshfeld	Difference	C45	--	C48	..	0.20	Ang.
PLAT234_ALERT_4_C	Large	Hirshfeld	Difference	C53	--	C56	..	0.19	Ang.
PLAT234_ALERT_4_C	Large	Hirshfeld	Difference	C13	--	C14	..	0.19	Ang.
PLAT241_ALERT_2_C	High		Ueq as Compared to Neighbors for				01	Check

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. C44 is part of a CF3 group.

PLAT241_ALERT_2_C	High		Ueq as Compared to Neighbors for				02	Check
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Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. C44 is part of a CF3 group.

PLAT241_ALERT_2_C	High		Ueq as Compared to Neighbors for				03	Check
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Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. C44 is part of a CF3 group.

PLAT241_ALERT_2_C	High		Ueq as Compared to Neighbors for				C48	Check
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Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. C44 is part of a CF3 group.

PLAT242_ALERT_2_C	Low		Ueq as Compared to Neighbors for				All	Check
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Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. C41 is a tertiary atom of a tert-butyl group.

PLAT242_ALERT_2_C	Low		Ueq as Compared to Neighbors for				C45	Check
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Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. C41 is a tertiary atom of a tert-butyl group.

PLAT242_ALERT_2_C Low Ueq as Compared to Neighbors for C49 Check

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. C41 is a tertiary atom of a tert-butyl group.

PLAT242_ALERT_2_C Low Ueq as Compared to Neighbors for C53 Check

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. C41 is a tertiary atom of a tert-butyl group.

PLAT243_ALERT_4_C High 'Solvent' Ueq as Compared to Neighbors of C57 Check
PLAT243_ALERT_4_C High 'Solvent' Ueq as Compared to Neighbors of C59 Check
PLAT243_ALERT_4_C High 'Solvent' Ueq as Compared to Neighbors of C65 Check
PLAT243_ALERT_4_C High 'Solvent' Ueq as Compared to Neighbors of C69 Check
PLAT244_ALERT_4_C Low 'Solvent' Ueq as Compared to Neighbors of C58 Check
PLAT244_ALERT_4_C Low 'Solvent' Ueq as Compared to Neighbors of C61 Check
PLAT244_ALERT_4_C Low 'Solvent' Ueq as Compared to Neighbors of C64 Check
PLAT244_ALERT_4_C Low 'Solvent' Ueq as Compared to Neighbors of C67 Check
PLAT342_ALERT_3_C Low Bond Precision on C-C Bonds 0.0153 Ang.
PLAT368_ALERT_2_C Short C(sp2)-C(sp2) Bond C66 - C67 ... 1.20 Ang.
PLAT368_ALERT_2_C Short C(sp2)-C(sp2) Bond C67 - C68 ... 1.22 Ang.
PLAT411_ALERT_2_C Short Inter H...H Contact H62 .. H93 .. 2.02 Ang.

● **Alert level G**

PLAT002_ALERT_2_G Number of Distance or Angle Restraints on AtSite 54 Note
PLAT003_ALERT_2_G Number of Uiso or Uij Restrained non-H Atoms ... 13 Report
PLAT042_ALERT_1_G Calc. and Reported MoietyFormula Strings Differ Please Check
PLAT072_ALERT_2_G SHELXL First Parameter in WGHT Unusually Large. 0.11 Report
PLAT083_ALERT_2_G SHELXL Second Parameter in WGHT Unusually Large. 17.28 Why ?
PLAT154_ALERT_1_G The su's on the Cell Angles are Equal 0.00300 Degree
PLAT171_ALERT_4_G The CIF-Embedded .res File Contains EADP Records 1 Report
PLAT172_ALERT_4_G The CIF-Embedded .res File Contains DFIX Records 14 Report
PLAT175_ALERT_4_G The CIF-Embedded .res File Contains SAME Records 2 Report
PLAT176_ALERT_4_G The CIF-Embedded .res File Contains SADI Records 7 Report
PLAT177_ALERT_4_G The CIF-Embedded .res File Contains DELU Records 2 Report
PLAT178_ALERT_4_G The CIF-Embedded .res File Contains SIMU Records 5 Report
PLAT230_ALERT_2_G Hirshfeld Test Diff for F50A -- C50 .. 6.0 su

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. The terminal fluorine atoms are strongly affected by this.

PLAT230_ALERT_2_G Hirshfeld Test Diff for F52B -- C52 .. 8.9 su

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. The terminal fluorine atoms are strongly affected by this.

PLAT230_ALERT_2_G Hirshfeld Test Diff for F54B -- C54 .. 6.7 su

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. The terminal fluorine atoms are strongly affected by this.

PLAT230_ALERT_2_G Hirshfeld Test Diff for F55A -- C55 .. 5.9 su

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. The terminal fluorine atoms are strongly affected by this.

PLAT230_ALERT_2_G Hirshfeld Test Diff for F56A -- C56 .. 8.8 su

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. The terminal fluorine atoms are strongly affected by this.

PLAT230_ALERT_2_G Hirshfeld Test Diff for C49 -- C51 .. 5.9 su

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. The terminal fluorine atoms are strongly affected by this.

PLAT230_ALERT_2_G Hirshfeld Test Diff for C49 -- C52 .. 7.2 su

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. The terminal fluorine atoms are strongly affected by this.

PLAT232_ALERT_2_G Hirshfeld Test Diff (M-X) Mo2 -- P6 .. 5.5 su
PLAT232_ALERT_2_G Hirshfeld Test Diff (M-X) Mo3 -- P9 .. 5.6 su
PLAT232_ALERT_2_G Hirshfeld Test Diff (M-X) Mo3 -- P12 .. 6.0 su
PLAT232_ALERT_2_G Hirshfeld Test Diff (M-X) Mo4 -- P8 .. 6.3 su
PLAT232_ALERT_2_G Hirshfeld Test Diff (M-X) Mo4 -- P10 .. 5.3 su
PLAT232_ALERT_2_G Hirshfeld Test Diff (M-X) Mo4 -- P11 .. 5.7 su
PLAT242_ALERT_2_G Low Ueq as Compared to Neighbors for C42 Check

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. C41 is a tertiary atom of a tert-butyl group.

PLAT242_ALERT_2_G Low Ueq as Compared to Neighbors for C46 Check

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. C41 is a tertiary atom of a tert-butyl group.

PLAT242_ALERT_2_G Low Ueq as Compared to Neighbors for C47 Check

Author Response: The weakly coordinating aluminate anion shows heavy disorder and thermal motion. C41 is a tertiary atom of a tert-butyl group.

PLAT301_ALERT_3_G	Main Residue Disorder	Percentage =	21	Note
PLAT302_ALERT_4_G	Anion/Solvent Disorder	Percentage =	20	Note
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C6	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C7	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C8	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C9	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C10	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C16	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C17	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C18	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C19	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C20	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C26	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C27	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C28	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C29	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C30	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C36	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C37	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C38	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C39	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C40	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C63	Check
PLAT380_ALERT_4_G	Incorrectly? Oriented X(sp2)-Methyl Moiety		C70	Check
PLAT434_ALERT_2_G	Short Inter HL..HL Contact F2 .. F56C .		2.68	Ang.
PLAT434_ALERT_2_G	Short Inter HL..HL Contact F4 .. F56C .		2.79	Ang.
PLAT434_ALERT_2_G	Short Inter HL..HL Contact F11 .. F50C .		2.72	Ang.
PLAT790_ALERT_4_G	Centre of Gravity not Within Unit Cell: Resd. #		2	Note
	C40 H60 Cu Mo4 P12			
PLAT860_ALERT_3_G	Number of Least-Squares Restraints		241	Note

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

2 ALERT type 1 CIF construction/syntax error, inconsistent or missing data
62 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
56 ALERT type 4 Improvement, methodology, query or suggestion
0 ALERT type 5 Informative message, check

checkCIF publication errors

Alert level A

PUBL002_ALERT_1_A The contact author's address is missing,
_publ_contact_author_address.

PUBL005_ALERT_1_A _publ_contact_author_email, _publ_contact_author_fax and
_publ_contact_author_phone are all missing.
At least one of these should be present.

PUBL006_ALERT_1_A _publ_requested_journal is missing
e.g. 'Acta Crystallographica Section C'

PUBL008_ALERT_1_A _publ_section_title is missing. Title of paper.

PUBL009_ALERT_1_A _publ_author_name is missing. List of author(s) name(s).

PUBL010_ALERT_1_A _publ_author_address is missing. Author(s) address(es).

PUBL012_ALERT_1_A _publ_section_abstract is missing.
Abstract of paper in English.

7 **ALERT level A** = Data missing that is essential or data in wrong format
0 **ALERT level G** = General alerts. Data that may be required is missing

Publication of your CIF

You should 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 nature of your study may justify the reported deviations from journal submission requirements and the more serious of these should 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.

If level A alerts remain, which you believe to be justified deviations, and you intend to submit this CIF for publication in a journal, you should additionally insert an explanation in your CIF using the Validation Reply Form (VRF) below. This will allow your explanation to be considered as part of the review process.

Validation response form

Please find below a validation response form (VRF) that can be filled in and pasted into your CIF.

```
# start Validation Reply Form
_vrf_PUBL002_GLOBAL
;
PROBLEM: The contact author's address is missing,
RESPONSE: ...
;
_vrf_PUBL005_GLOBAL
;
PROBLEM: _publ_contact_author_email, _publ_contact_author_fax and
RESPONSE: ...
;
_vrf_PUBL006_GLOBAL
;
PROBLEM: _publ_requested_journal is missing
RESPONSE: ...
;
_vrf_PUBL008_GLOBAL
;
PROBLEM: _publ_section_title is missing. Title of paper.
RESPONSE: ...
;
_vrf_PUBL009_GLOBAL
;
PROBLEM: _publ_author_name is missing. List of author(s) name(s).
RESPONSE: ...
;
```

```
_vrf_PUBL010_GLOBAL
;
PROBLEM: _publ_author_address is missing. Author(s) address(es).
RESPONSE: ...
;
_vrf_PUBL012_GLOBAL
;
PROBLEM: _publ_section_abstract is missing.
RESPONSE: ...
;
# end Validation Reply Form
```

If you wish to submit your CIF for publication in Acta Crystallographica Section C or E, you should upload your CIF via the web. If your CIF is to form part of a submission to another IUCr journal, you will be asked, either during electronic submission or by the Co-editor handling your paper, to upload your CIF via our web site.

PLATON version of 20/08/2014; check.def file version of 18/08/2014

Datablock I - ellipsoid plot

