



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 22, 2024 – 11:00 PM EDT

PDB ID : 6S3I  
Title : Crystal structure of helicase Pif1 from *Thermus oshimai* in complex with ss-DNA (dT)18 and ADP-MgF4  
Authors : Dai, Y.X.; Chen, W.F.; Teng, F.Y.; Liu, N.N.; Hou, X.M.; Dou, S.X.; Rety, S.; Xi, X.G.  
Deposited on : 2019-06-25  
Resolution : 2.46 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

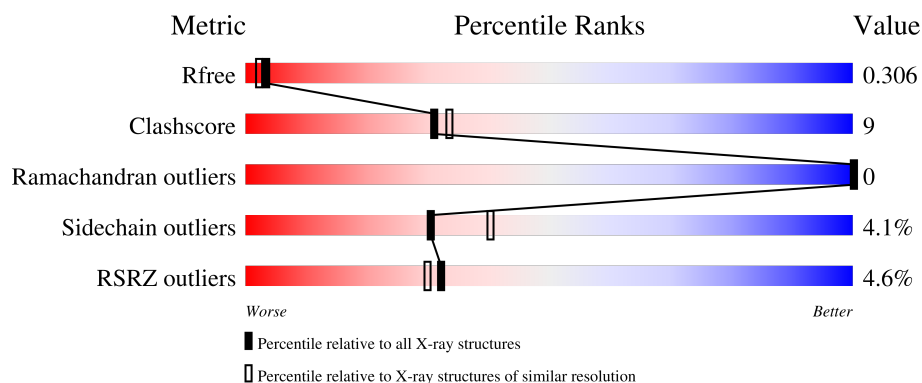
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.46 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	1544 (2.48-2.44)
Clashscore	141614	1613 (2.48-2.44)
Ramachandran outliers	138981	1598 (2.48-2.44)
Sidechain outliers	138945	1598 (2.48-2.44)
RSRZ outliers	127900	1523 (2.48-2.44)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	444	<div> <div>5%</div> <div>75%</div> <div>22%</div> <div>..</div> </div>
2	C	14	<div> <div>21%</div> <div>14%</div> <div>64%</div> </div>

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7253 atoms, of which 3619 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called PIF1 helicase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	434	Total	C	H	N	O	S	0	0	0
			7033	2231	3547	636	616	3			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	64	THR	ALA	conflict	UNP K7RJ88
A	162	ILE	MET	conflict	UNP K7RJ88
A	456	LEU	PRO	conflict	UNP K7RJ88

- Molecule 2 is a DNA chain called DNA (5'-D(P\*TP\*TP\*TP\*TP\*T)-3').

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	C	5	Total	C	H	N	O	P	0	0	0
			160	50	60	10	35	5			

- Molecule 3 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C<sub>10</sub>H<sub>15</sub>N<sub>5</sub>O<sub>10</sub>P<sub>2</sub>) (labeled as "Ligand of Interest" by depositor).

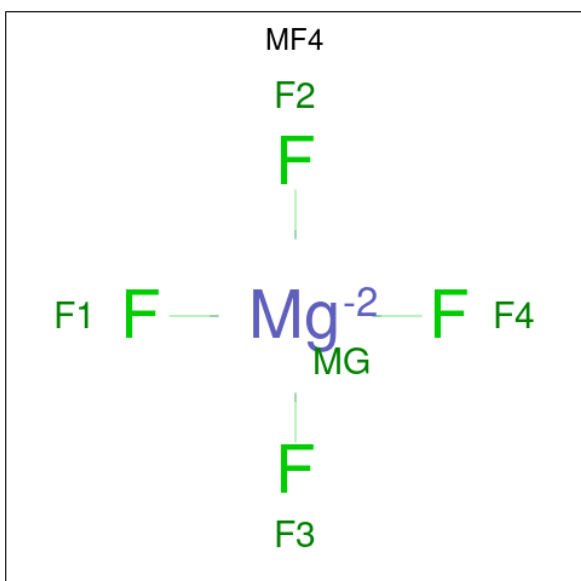


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	P	
			39	10	12	5	10	2	

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg		
			1	1	0	0

- Molecule 5 is TETRAFLUOROMAGNESATE(2-) (three-letter code: MF4) (formula: F<sub>4</sub>Mg) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	F	Mg	0	0
			5	4	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	13	Total	O	0	0
			13	13		
6	C	2	Total	O	0	0
			2	2		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	61.20Å 61.20Å 260.36Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	59.57 – 2.46 59.57 – 2.45	Depositor EDS
% Data completeness (in resolution range)	92.3 (59.57-2.46) 92.3 (59.57-2.45)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.05 (at 2.45Å)	Xtriage
Refinement program	PHENIX (1.14rc1_3177: ???)	Depositor
R, $R_{free}$	0.244 , 0.301 0.253 , 0.306	Depositor DCC
$R_{free}$ test set	913 reflections (5.18%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	63.7	Xtriage
Anisotropy	0.144	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 52.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	7253	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.62% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ADP, MG, MF4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.30	0/3568	0.52	0/4843
2	C	0.75	0/109	1.40	0/166
All	All	0.32	0/3677	0.57	0/5009

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3486	3547	3547	64	0
2	C	100	60	61	2	0
3	A	27	12	12	0	0
4	A	1	0	0	0	0
5	A	5	0	0	0	0
6	A	13	0	0	7	0
6	C	2	0	0	1	0
All	All	3634	3619	3620	66	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.



All (66) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:447:GLY:N	6:A:1102:HOH:O	1.91	1.03
1:A:285:ALA:HB2	1:A:466:LEU:HD21	1.47	0.96
1:A:260:ASP:OD1	6:A:1101:HOH:O	1.90	0.87
1:A:68:LEU:N	6:A:1104:HOH:O	2.08	0.86
2:C:7:DT:O3'	6:C:101:HOH:O	2.01	0.77
1:A:361:GLU:OE1	6:A:1103:HOH:O	2.05	0.74
1:A:450:LEU:HD23	1:A:479:GLU:HB3	1.69	0.72
1:A:211:ARG:NH1	1:A:487:GLU:OE2	2.24	0.70
1:A:162:ILE:O	1:A:192:ARG:NH1	2.24	0.70
1:A:466:LEU:HD23	1:A:466:LEU:O	1.94	0.67
1:A:378:LEU:HD21	1:A:391:ILE:HD11	1.80	0.62
1:A:307:ASP:OD1	1:A:424:ARG:NH2	2.34	0.61
1:A:262:LEU:H	1:A:262:LEU:HD12	1.65	0.61
1:A:314:LEU:HD22	1:A:425:LEU:HD11	1.83	0.59
1:A:241:TRP:HB2	1:A:500:LYS:HG3	1.86	0.57
1:A:366:ASP:OD1	1:A:384:ARG:HD3	2.06	0.56
1:A:111:ARG:NH2	1:A:166:GLU:OE1	2.38	0.56
1:A:166:GLU:N	1:A:166:GLU:OE2	2.39	0.56
1:A:84:HIS:O	1:A:87:HIS:NE2	2.40	0.55
1:A:92:PRO:HD2	1:A:95:THR:HG21	1.87	0.55
1:A:403:TYR:HA	1:A:410:ILE:HG22	1.91	0.53
1:A:481:LEU:HD12	1:A:481:LEU:H	1.75	0.51
1:A:225:TYR:O	1:A:229:THR:OG1	2.24	0.50
1:A:84:HIS:ND1	1:A:247:ARG:HD2	2.26	0.49
1:A:446:LEU:HD22	1:A:473:ARG:NH2	2.26	0.49
1:A:393:PRO:HA	1:A:421:VAL:HA	1.94	0.48
1:A:274:ASP:HB3	1:A:277:ALA:HB3	1.95	0.48
1:A:450:LEU:HD23	1:A:479:GLU:CB	2.41	0.48
1:A:303:ARG:NH1	1:A:337:PHE:O	2.48	0.47
1:A:446:LEU:C	6:A:1102:HOH:O	2.40	0.47
1:A:328:VAL:HG22	1:A:418:PHE:CD2	2.50	0.47
1:A:395:VAL:O	1:A:395:VAL:HG13	2.15	0.46
1:A:449:GLY:N	6:A:1102:HOH:O	2.48	0.46
1:A:304:LYS:HE2	1:A:335:THR:HA	1.98	0.45
1:A:443:HIS:ND1	1:A:470:SER:HB2	2.31	0.45
1:A:135:ARG:O	1:A:135:ARG:CG	2.65	0.45
1:A:430:THR:HG23	1:A:433:LYS:H	1.82	0.45
2:C:6:DT:C6	2:C:7:DT:H72	2.53	0.44
1:A:357:ASP:OD2	1:A:361:GLU:HG3	2.17	0.44
1:A:113:VAL:CG2	1:A:168:LEU:HD13	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:463:VAL:HG21	1:A:469:LEU:HB2	2.01	0.43
1:A:401:TYR:CD1	1:A:410:ILE:HD13	2.54	0.43
1:A:351:VAL:HB	1:A:423:VAL:HG13	2.01	0.43
1:A:338:PRO:HG2	1:A:396:TRP:CD2	2.54	0.42
1:A:285:ALA:O	1:A:471:LEU:HB2	2.19	0.42
1:A:385:ASN:HB2	1:A:387:ARG:HH11	1.84	0.42
1:A:263:PHE:CZ	1:A:267:LEU:HD11	2.54	0.42
1:A:267:LEU:HD13	1:A:461:THR:HG22	2.02	0.42
1:A:135:ARG:HD2	1:A:135:ARG:C	2.40	0.42
1:A:168:LEU:HD21	1:A:184:MET:HE3	2.01	0.42
1:A:109:LYS:C	1:A:111:ARG:H	2.23	0.42
1:A:410:ILE:H	1:A:410:ILE:HG13	1.66	0.42
1:A:79:VAL:HG21	1:A:89:ILE:HD11	2.02	0.42
1:A:116:ALA:O	1:A:129:THR:HA	2.19	0.42
1:A:262:LEU:HD12	1:A:262:LEU:N	2.33	0.42
1:A:271:ARG:NH2	6:A:1107:HOH:O	2.34	0.42
1:A:347:LYS:HE3	1:A:371:GLU:O	2.20	0.42
1:A:353:LEU:HD23	1:A:423:VAL:HG22	2.01	0.41
1:A:88:LEU:HD23	1:A:248:VAL:HG22	2.03	0.41
1:A:222:GLU:O	1:A:226:ILE:HG13	2.21	0.41
1:A:394:PHE:CD1	1:A:422:PRO:HG3	2.55	0.41
1:A:405:SER:OG	1:A:406:GLU:N	2.54	0.41
1:A:144:ARG:HD3	1:A:186:TRP:CD2	2.56	0.41
1:A:482:TRP:CD1	1:A:482:TRP:N	2.89	0.40
1:A:69:SER:HB3	1:A:72:GLN:OE1	2.21	0.40
1:A:483:ARG:HD3	1:A:483:ARG:HA	1.96	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	432/444 (97%)	425 (98%)	7 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	362/371 (98%)	347 (96%)	15 (4%)	30	40

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	97	LYS
1	A	132	SER
1	A	155	HIS
1	A	258	ARG
1	A	271	ARG
1	A	276	GLN
1	A	302	ARG
1	A	317	LEU
1	A	322	LEU
1	A	366	ASP
1	A	378	LEU
1	A	450	LEU
1	A	464	ARG
1	A	470	SER
1	A	482	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	MF4	A	1003	4,3	0,4,4	-	-	-		
3	ADP	A	1001	5,4	24,29,29	0.92	1 (4%)	29,45,45	1.40	3 (10%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ADP	A	1001	5,4	-	1/12/32/32	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1001	ADP	PA-O3A	2.21	1.61	1.59

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1001	ADP	N3-C2-N1	-4.72	122.26	128.67
3	A	1001	ADP	C1'-N9-C4	-2.40	122.42	126.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1001	ADP	C4-C5-N7	-2.31	106.90	109.34

There are no chirality outliers.

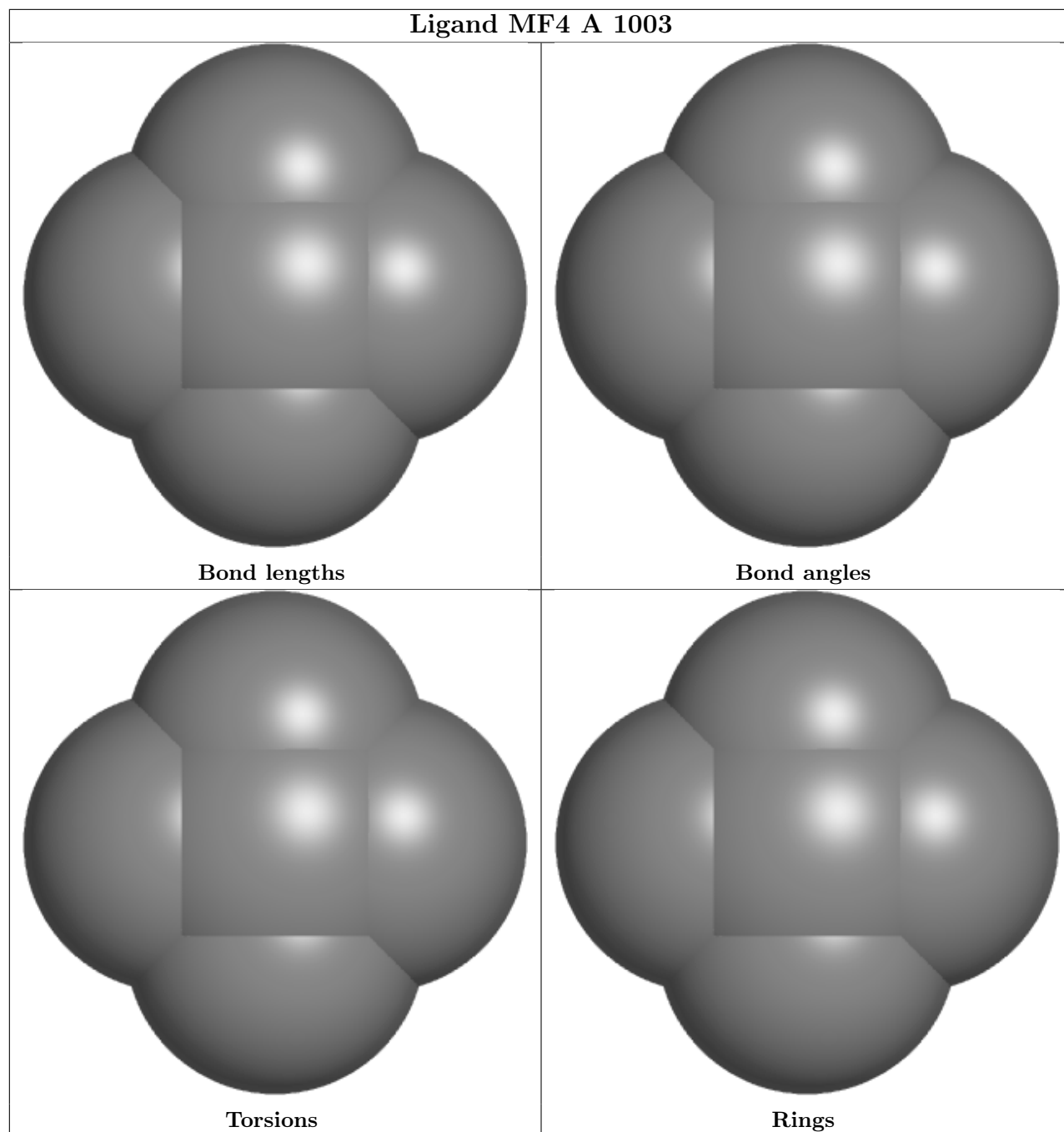
All (1) torsion outliers are listed below:

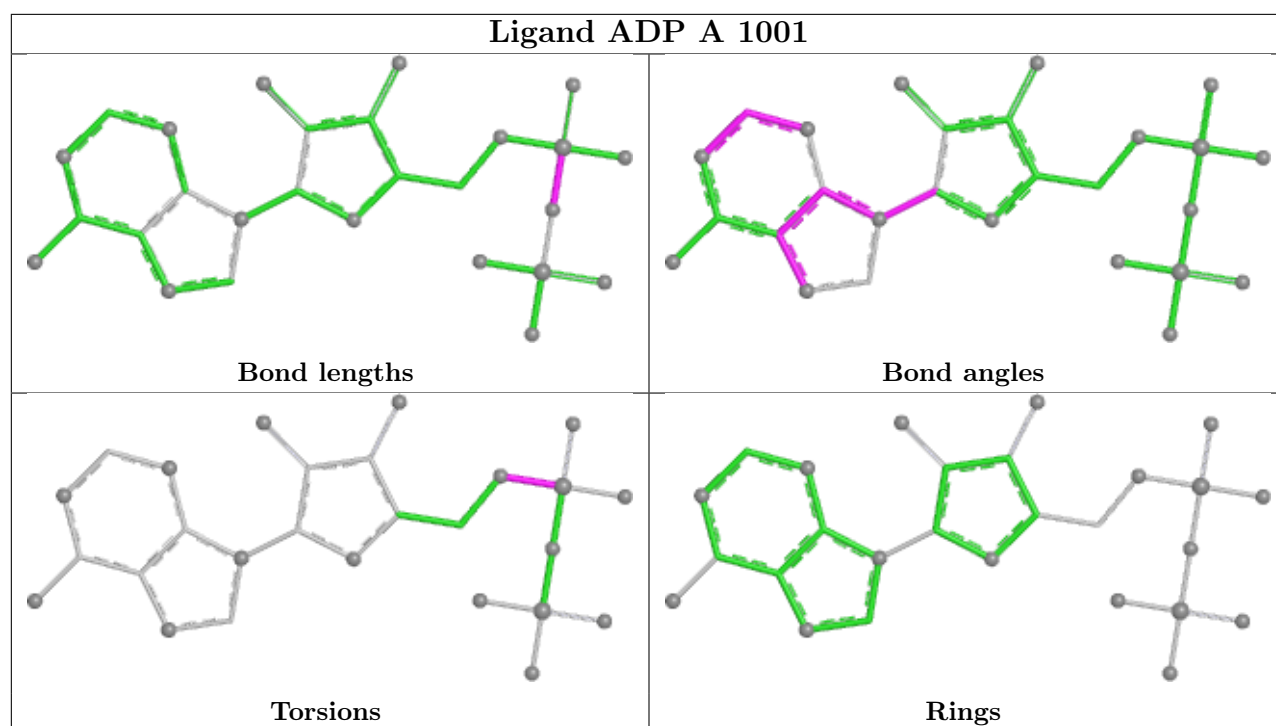
Mol	Chain	Res	Type	Atoms
3	A	1001	ADP	C5'-O5'-PA-O1A

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	434/444 (97%)	0.45	20 (4%) 32 30	45, 69, 92, 110	0
2	C	5/14 (35%)	0.27	0 100 100	61, 67, 74, 97	0
All	All	439/458 (95%)	0.45	20 (4%) 32 30	45, 69, 92, 110	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	501	SER	12.0
1	A	404	ASP	5.5
1	A	343	LEU	3.6
1	A	482	TRP	3.2
1	A	295	GLY	3.0
1	A	358	PRO	2.6
1	A	313	ARG	2.6
1	A	478	THR	2.6
1	A	497	ILE	2.3
1	A	396	TRP	2.3
1	A	327	GLN	2.2
1	A	318	PRO	2.2
1	A	328	VAL	2.1
1	A	299	LEU	2.1
1	A	294	PRO	2.1
1	A	479	GLU	2.1
1	A	481	LEU	2.0
1	A	227	ALA	2.0
1	A	329	LYS	2.0
1	A	423	VAL	2.0



## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 6.4 Ligands [i](#)

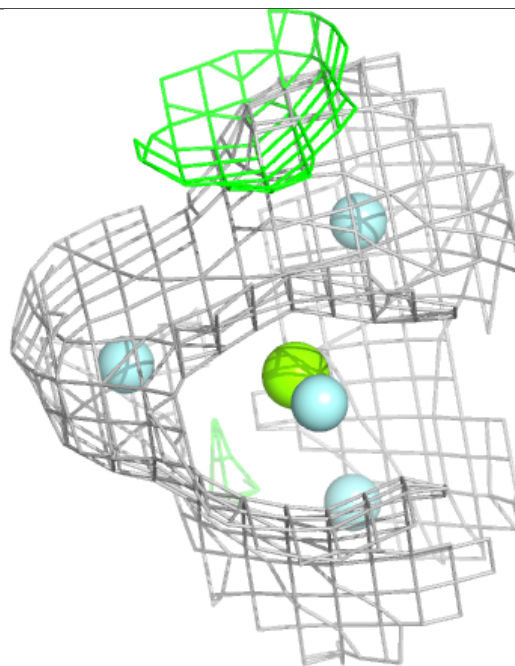
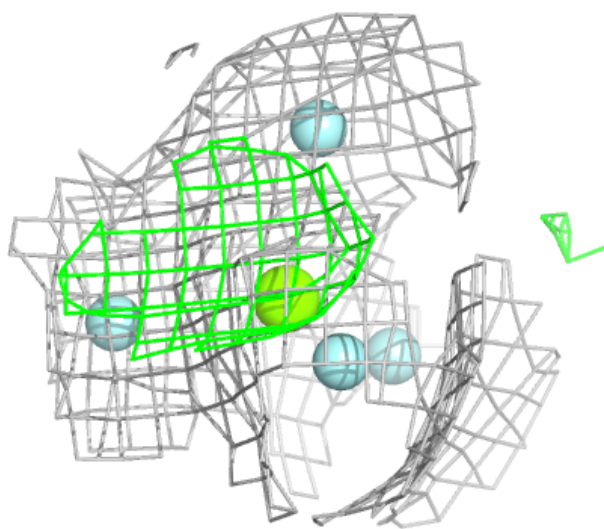
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	MF4	A	1003	5/5	0.95	0.11	44,44,61,72	0
4	MG	A	1002	1/1	0.96	0.12	62,62,62,62	0
3	ADP	A	1001	27/27	0.97	0.15	44,68,89,91	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

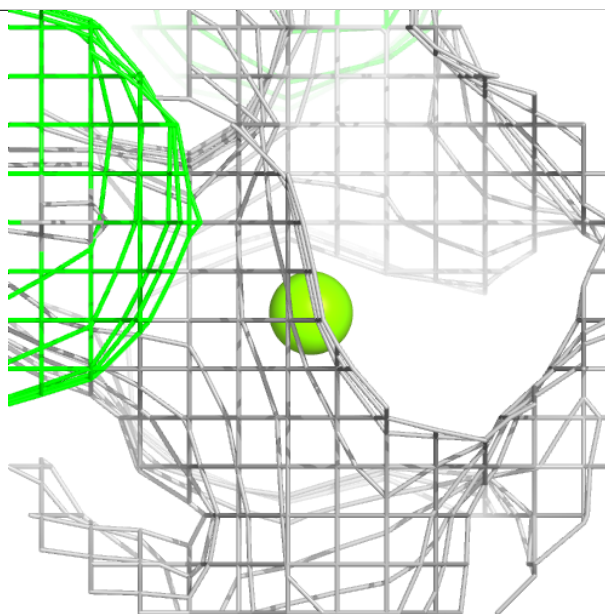
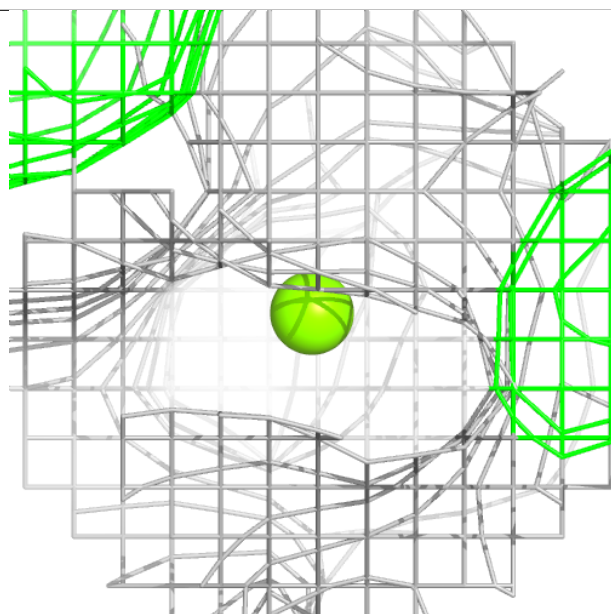
**Electron density around MF4 A 1003:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



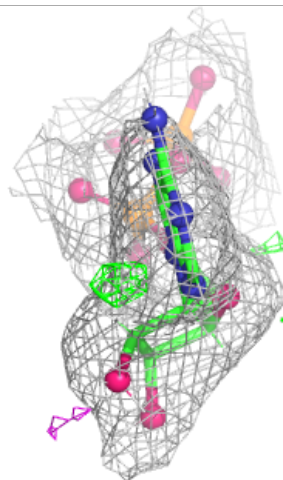
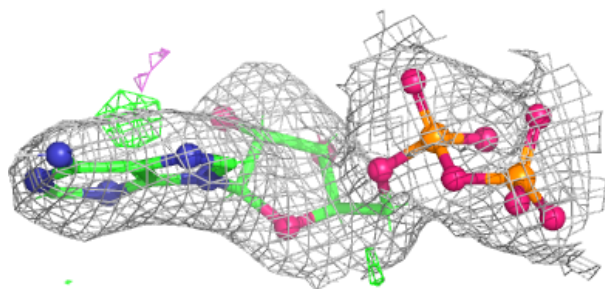
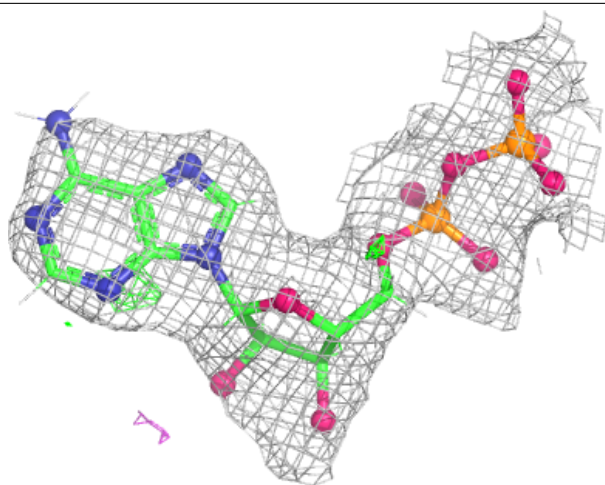
**Electron density around MG A 1002:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around ADP A 1001:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



## 6.5 Other polymers [i](#)

There are no such residues in this entry.