



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:58 PM GMT

PDB ID : 1I6I  
Title : CRYSTAL STRUCTURE OF THE KIF1A MOTOR DOMAIN COM-  
PLEXED WITH MG-AMPPCP  
Authors : Kikkawa, M.; Sablin, E.P.; Okada, Y.; Yajima, H.; Fletterick, R.J.; Hirokawa,  
N.  
Deposited on : 2001-03-02  
Resolution : 2.00 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  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

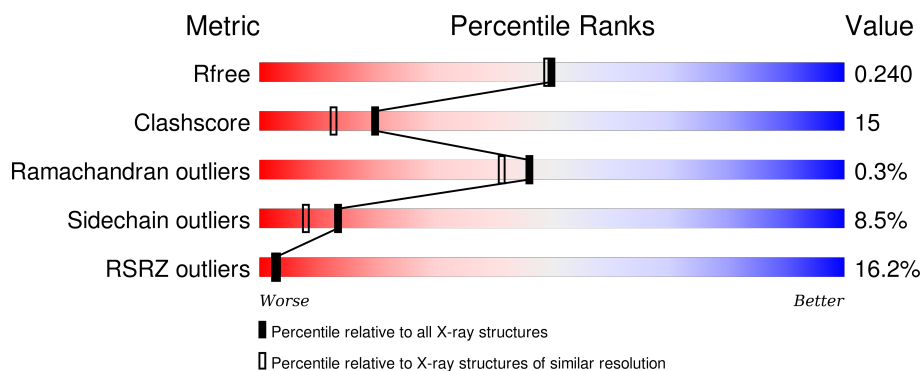
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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))
$R_{free}$	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	366	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	TRS	A	1000	-	X	-	X

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2979 atoms, of which 0 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 KINESIN-LIKE PROTEIN KIF1A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	333	Total	C	N	O	S	0	10	0
			2708	1675	480	537	16			

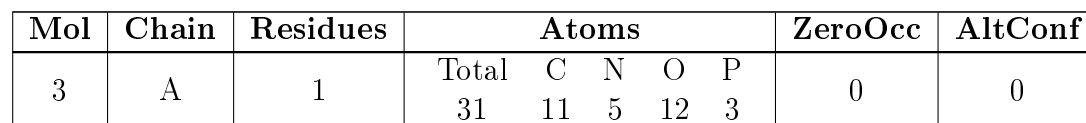
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	202	ALA	PRO	ENGINEERED	UNP P33173
A	356	ASN	-	SEE REMARK 999	UNP P33173
A	357	THR	-	SEE REMARK 999	UNP P33173
A	358	VAL	-	SEE REMARK 999	UNP P33173
A	359	SER	-	SEE REMARK 999	UNP P33173
A	360	VAL	-	SEE REMARK 999	UNP P33173
A	361	ASN	-	SEE REMARK 999	UNP P33173
A	362	HIS	-	SEE REMARK 999	UNP P33173
A	363	HIS	-	SEE REMARK 999	UNP P33173
A	364	HIS	-	SEE REMARK 999	UNP P33173
A	365	HIS	-	SEE REMARK 999	UNP P33173
A	366	HIS	-	SEE REMARK 999	UNP P33173

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 3 is PHOSPHOMETHYLPHOSPHONIC ACID ADENYLATE ESTER (three-letter code: ACP) (formula: C<sub>11</sub>H<sub>18</sub>N<sub>5</sub>O<sub>12</sub>P<sub>3</sub>).



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- Chemical structure of 2,3,4-trihydroxybutan-1-amine (THS). The structure shows a four-carbon chain. The first carbon (C1) is bonded to an amino group (NH<sub>3</sub><sup>+</sup>). The second carbon (C2) is bonded to a hydroxyl group (OH). The third carbon (C3) is bonded to a hydroxyl group (OH). The fourth carbon (C4) is bonded to a hydroxyl group (OH). The atoms are labeled as follows: C1, C2, C3, C4, NH<sub>3</sub><sup>+</sup>, OH, OH, OH.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			8	4	1	3		

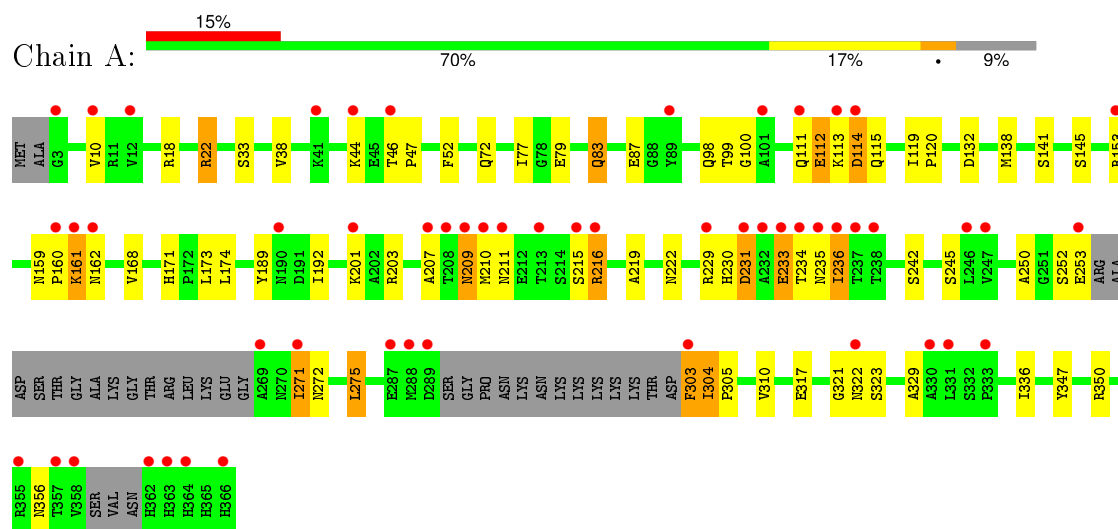
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	231	Total 231	O 231	0	0

### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

#### • Molecule 1: KINESIN-LIKE PROTEIN KIF1A



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	41.99 Å   56.40 Å   156.12 Å 90.00°   90.00°   90.00°	Depositor
Resolution (Å)	19.70 – 2.00 25.06 – 2.00	Depositor EDS
% Data completeness (in resolution range)	92.1 (19.70-2.00) 92.2 (25.06-2.00)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	7.19 (at 2.01 Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.219   ,   0.230 0.228   ,   0.240	Depositor DCC
$R_{free}$ test set	1150 reflections (4.86%)	DCC
Wilson B-factor (Å <sup>2</sup> )	28.6	Xtriage
Anisotropy	0.161	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 56.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 24025 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	2979	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.85% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.45	0/2751	0.71	1/3711 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	153	ARG	NE-CZ-NH1	-5.75	117.43	120.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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	2708	0	2638	81	0
2	A	1	0	0	0	0
3	A	31	0	14	2	0
4	A	8	0	12	0	0
5	A	231	0	0	8	0
All	All	2979	0	2664	81	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 15.



All (81) 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:159:ASN:ND2	1:A:161:LYS:HG3	1.62	1.15
1:A:18[B]:ARG:HH21	1:A:336:ILE:HD12	1.15	1.08
1:A:233:GLU:HG3	1:A:234:THR:H	0.90	1.07
1:A:252:SER:O	1:A:253:GLU:HG2	1.53	1.06
1:A:233:GLU:HG3	1:A:234:THR:N	1.73	1.02
1:A:160:PRO:HA	5:A:771:HOH:O	1.61	1.00
1:A:233:GLU:CG	1:A:234:THR:H	1.76	0.96
1:A:159:ASN:HD22	1:A:161:LYS:HG3	1.26	0.95
1:A:216:ARG:HH11	1:A:216:ARG:HG3	1.33	0.92
1:A:252:SER:HA	1:A:275:LEU:HD11	1.50	0.92
1:A:252:SER:HA	1:A:275:LEU:CD1	2.02	0.90
1:A:234:THR:HB	1:A:236:ILE:HD11	1.56	0.86
1:A:317:GLU:OE1	1:A:322:ASN:HB3	1.74	0.86
1:A:234:THR:HB	1:A:236:ILE:CD1	2.10	0.80
1:A:18[B]:ARG:NH2	1:A:336:ILE:HD12	1.98	0.78
1:A:216:ARG:HH11	1:A:216:ARG:CG	1.98	0.75
1:A:159:ASN:HD22	1:A:161:LYS:CG	2.02	0.72
1:A:252:SER:O	1:A:253:GLU:CG	2.37	0.72
1:A:168:VAL:HG21	1:A:310:VAL:HG13	1.72	0.70
1:A:250:ALA:HA	5:A:751:HOH:O	1.94	0.66
1:A:231:ASP:O	1:A:233:GLU:HG2	1.96	0.66
1:A:159:ASN:ND2	1:A:161:LYS:CG	2.52	0.65
1:A:171:HIS:CD2	1:A:174:LEU:H	2.18	0.61
1:A:303:PHE:C	1:A:303:PHE:CD2	2.72	0.61
1:A:303:PHE:O	1:A:303:PHE:CD2	2.56	0.59
1:A:171:HIS:HD2	1:A:173:LEU:H	1.51	0.59
1:A:83[A]:GLN:NE2	1:A:87:GLU:OE2	2.36	0.58
1:A:252:SER:HA	1:A:275:LEU:HD12	1.83	0.57
1:A:252:SER:CA	1:A:275:LEU:HD11	2.31	0.57
1:A:159:ASN:HD21	1:A:161:LYS:HG3	1.64	0.56
1:A:100:GLY:H	3:A:500:ACP:H3B1	1.68	0.56
1:A:18[B]:ARG:HH21	1:A:336:ILE:CD1	2.04	0.56
1:A:119:ILE:HB	1:A:120:PRO:HD3	1.87	0.56
1:A:207:ALA:HA	1:A:211:ASN:OD1	2.06	0.56
1:A:77[B]:ILE:HD12	1:A:77[B]:ILE:H	1.70	0.56
1:A:18[A]:ARG:HE	1:A:22[A]:ARG:NH2	2.03	0.56
1:A:159:ASN:C	1:A:161:LYS:H	2.09	0.55
1:A:234:THR:CB	1:A:236:ILE:HD11	2.31	0.55
1:A:171:HIS:HD2	1:A:174:LEU:H	1.52	0.54
1:A:22[B]:ARG:NH1	5:A:607:HOH:O	2.40	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:321:GLY:O	1:A:356:ASN:HB3	2.08	0.53
1:A:171:HIS:CD2	1:A:173:LEU:H	2.28	0.52
1:A:216:ARG:HD2	1:A:216:ARG:N	2.24	0.52
1:A:83[A]:GLN:O	1:A:87:GLU:HG3	2.10	0.52
1:A:33:SER:HB2	1:A:52:PHE:O	2.10	0.51
1:A:83[B]:GLN:O	1:A:87:GLU:HG3	2.10	0.51
1:A:209:ASN:O	1:A:210[A]:MET:HB2	2.11	0.51
1:A:138:MET:SD	1:A:229:ARG:HB2	2.52	0.50
1:A:317:GLU:CD	1:A:322:ASN:H	2.15	0.50
1:A:216:ARG:NH1	1:A:216:ARG:CG	2.67	0.50
1:A:201:LYS:HE3	5:A:757:HOH:O	2.11	0.50
1:A:215:SER:C	1:A:216:ARG:HD2	2.33	0.48
1:A:222:ASN:HD22	1:A:245[B]:SER:HA	1.78	0.48
1:A:215:SER:OG	1:A:216:ARG:HD2	2.13	0.47
1:A:230:HIS:CG	1:A:231:ASP:N	2.83	0.47
1:A:234:THR:CB	1:A:236:ILE:CD1	2.88	0.47
1:A:216:ARG:NH1	5:A:697:HOH:O	2.48	0.46
1:A:209:ASN:O	1:A:210[B]:MET:HB2	2.15	0.46
1:A:216:ARG:NH1	1:A:216:ARG:HG3	2.15	0.46
1:A:222:ASN:HD22	1:A:245[A]:SER:HA	1.79	0.46
1:A:189:TYR:O	1:A:192:ILE:HG22	2.16	0.45
1:A:203:ARG:CZ	1:A:219:ALA:HB2	2.47	0.45
1:A:234:THR:HB	1:A:236:ILE:HD13	1.93	0.45
1:A:72[A]:GLN:NE2	5:A:612:HOH:O	2.49	0.45
1:A:216:ARG:NH1	1:A:272:ASN:OD1	2.49	0.45
1:A:100:GLY:N	3:A:500:ACP:H3B1	2.32	0.45
1:A:161:LYS:H	1:A:161:LYS:HG2	1.25	0.44
1:A:233:GLU:CG	1:A:234:THR:N	2.50	0.44
1:A:216:ARG:HA	1:A:250:ALA:HB1	2.00	0.44
1:A:38:VAL:HG22	1:A:47:PRO:HG3	1.99	0.44
1:A:112:GLU:O	1:A:114:ASP:N	2.51	0.44
1:A:271:ILE:H	1:A:271:ILE:HG12	1.52	0.43
1:A:10:VAL:HG12	1:A:329:ALA:HB3	2.00	0.42
1:A:18[A]:ARG:HE	1:A:22[A]:ARG:HH22	1.66	0.42
1:A:83[A]:GLN:NE2	1:A:87:GLU:CD	2.73	0.42
1:A:98:GLN:HG2	1:A:99:THR:N	2.35	0.41
1:A:271:ILE:HD13	1:A:271:ILE:N	2.36	0.41
1:A:347:TYR:CE2	1:A:350:ARG:NH1	2.89	0.41
1:A:79:GLU:HG2	5:A:708:HOH:O	2.20	0.41
1:A:304:ILE:HA	1:A:305:PRO:HD3	1.90	0.40
1:A:203:ARG:HD3	5:A:670:HOH:O	2.20	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	335/366 (92%)	321 (96%)	13 (4%)	1 (0%)	46 41

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	233	GLU

### 5.3.2 Protein sidechains [i](#)

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	303/321 (94%)	274 (90%)	29 (10%)	10 6

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22[A]	ARG
1	A	22[B]	ARG
1	A	44	LYS
1	A	46	THR
1	A	83[A]	GLN
1	A	83[B]	GLN
1	A	111	GLN

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Mol	Chain	Res	Type
1	A	112	GLU
1	A	113	LYS
1	A	114	ASP
1	A	115	GLN
1	A	132	ASP
1	A	141	SER
1	A	145[A]	SER
1	A	145[B]	SER
1	A	161	LYS
1	A	162	ASN
1	A	209	ASN
1	A	216	ARG
1	A	231	ASP
1	A	235	ASN
1	A	236	ILE
1	A	242[A]	SER
1	A	242[B]	SER
1	A	271	ILE
1	A	275	LEU
1	A	303	PHE
1	A	304	ILE
1	A	323	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	159	ASN
1	A	162	ASN
1	A	171	HIS
1	A	222	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 5.5 Carbohydrates

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

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	$\# Z  > 2$
4	TRS	A	1000	-	7,7,7	2.54	3 (42%)	9,9,9	3.39	7 (77%)
3	ACP	A	500	2	25,33,33	1.49	4 (16%)	31,52,52	2.51	2 (6%)

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
4	TRS	A	1000	-	-	0/9/9/9	0/0/0/0
3	ACP	A	500	2	-	0/15/38/38	0/3/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1000	TRS	C-N	-5.67	1.42	1.50
3	A	500	ACP	PG-O3G	-2.37	1.49	1.54
4	A	1000	TRS	C1-C	2.08	1.57	1.53
4	A	1000	TRS	C2-C	2.25	1.57	1.53
3	A	500	ACP	PG-O2G	2.34	1.60	1.54
3	A	500	ACP	C4-N3	2.70	1.39	1.35
3	A	500	ACP	O4'-C1'	3.91	1.46	1.41

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	500	ACP	N3-C2-N1	-12.59	119.26	128.89
4	A	1000	TRS	O1-C1-C	-4.64	101.78	111.18
4	A	1000	TRS	C2-C-C1	-4.53	100.97	110.78
3	A	500	ACP	C4-C5-N7	-3.82	105.96	109.48
4	A	1000	TRS	C1-C-N	-3.06	102.52	108.09
4	A	1000	TRS	O2-C2-C	-2.17	106.80	111.18
4	A	1000	TRS	C3-C-C2	-2.09	106.26	110.78
4	A	1000	TRS	C2-C-N	4.32	115.94	108.09
4	A	1000	TRS	C3-C-N	4.37	116.04	108.09

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	500	ACP	2	0

## 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	333/366 (90%)	0.91	54 (16%) <b>3</b> <b>3</b>	15, 30, 54, 59	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	269	ALA	11.3
1	A	234	THR	8.8
1	A	236	ILE	7.2
1	A	303	PHE	6.7
1	A	366	HIS	6.6
1	A	210[A]	MET	6.3
1	A	231	ASP	6.2
1	A	208	THR	5.8
1	A	232	ALA	5.7
1	A	289	ASP	5.3
1	A	358	VAL	4.8
1	A	211	ASN	4.5
1	A	3	GLY	4.5
1	A	238	THR	4.4
1	A	213	THR	4.3
1	A	288	MET	4.3
1	A	114	ASP	4.2
1	A	287	GLU	3.9
1	A	113	LYS	3.9
1	A	111	GLN	3.7
1	A	235	ASN	3.4
1	A	362	HIS	3.3
1	A	216	ARG	3.3
1	A	322	ASN	3.2
1	A	161	LYS	3.2
1	A	357	THR	3.1
1	A	160	PRO	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	207	ALA	2.7
1	A	237	THR	2.7
1	A	162	ASN	2.7
1	A	247	VAL	2.7
1	A	233	GLU	2.6
1	A	363	HIS	2.6
1	A	271	ILE	2.5
1	A	101	ALA	2.5
1	A	209	ASN	2.5
1	A	331	LEU	2.5
1	A	364	HIS	2.5
1	A	355	ARG	2.5
1	A	153	ARG	2.4
1	A	253	GLU	2.4
1	A	10	VAL	2.4
1	A	44	LYS	2.4
1	A	330	ALA	2.4
1	A	190	ASN	2.3
1	A	46	THR	2.3
1	A	246	LEU	2.3
1	A	89	TYR	2.2
1	A	201	LYS	2.2
1	A	229	ARG	2.1
1	A	41	LYS	2.1
1	A	12	VAL	2.1
1	A	333	PRO	2.1
1	A	215	SER	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains.



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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	TRS	A	1000	8/8	0.74	0.37	4.55	55,55,55,56	0
3	ACP	A	500	31/31	0.94	0.16	-0.41	21,27,34,37	0
2	MG	A	501	1/1	0.89	0.08	-	28,28,28,28	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.