



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 09:13 PM GMT

PDB ID : 4Z4C  
Title : Human Argonaute2 Bound to t1-C Target RNA  
Authors : Schirle, N.T.; MacRae, I.J.  
Deposited on : 2015-04-02  
Resolution : 2.30 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 : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026982  
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) : rb-20026982

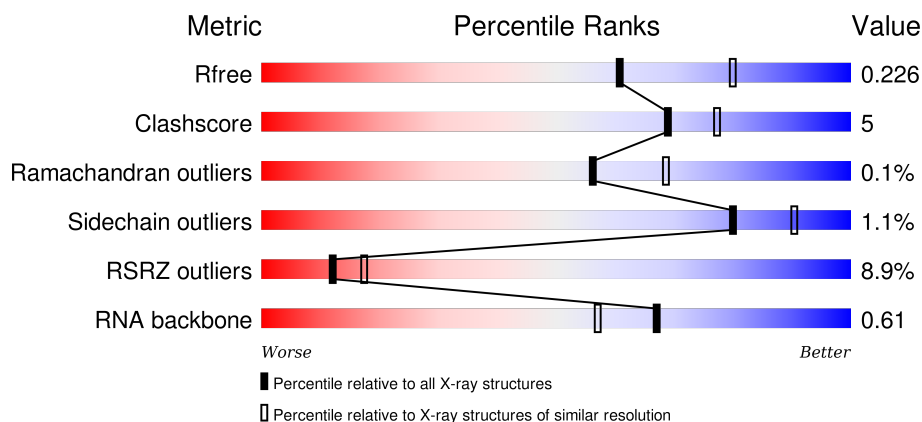
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)
RNA backbone	2183	1011 (2.84-1.76)

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	859	<div> <div>9%</div> <div>84%</div> <div>9%</div> <div>7%</div> </div>
2	B	21	<div> <div>52%</div> <div>38%</div> <div>10%</div> </div>
3	D	11	<div> <div>64%</div> <div>18%</div> <div>18%</div> </div>

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
5	IPA	A	902	-	-	-	X
6	IPH	A	903	-	-	-	X

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 7321 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 Protein argonaute-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	803	Total	C	N	O	S	0	0	0
			6429	4095	1156	1138	40			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	387	ASP	SER	engineered mutation	UNP Q9UKV8

- Molecule 2 is a RNA chain called RNA (5'-R(P\*UP\*UP\*CP\*AP\*CP\*AP\*UP\*UP\*GP\*CP\*CP\*CP\*AP\*AP\*GP\*UP\*CP\*UP\*U)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	19	Total	C	N	O	P	0	0	0
			379	168	60	132	19			

- Molecule 3 is a RNA chain called RNA (5'-R(\*CP\*AP\*AP\*UP\*GP\*UP\*GP\*AP\*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	9	Total	C	N	O	P	0	0	0
			189	86	35	60	8			

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

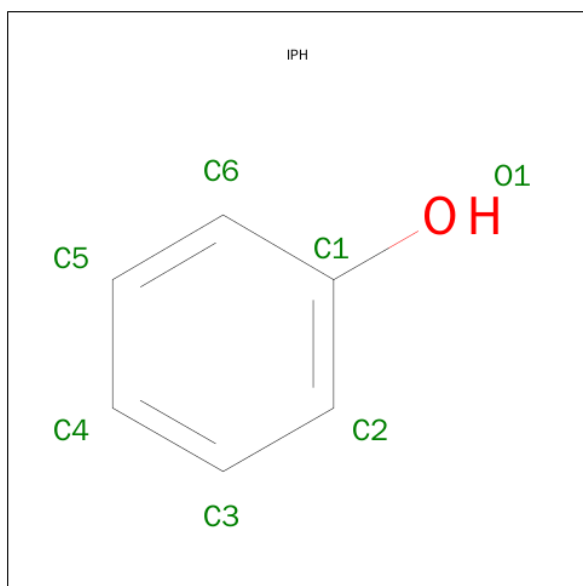
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Mg	0	0
			1	1		
4	A	1	Total	Mg	0	0
			1	1		
4	D	1	Total	Mg	0	0
			1	1		

- Molecule 5 is ISOPROPYL ALCOHOL (three-letter code: IPA) (formula: C<sub>3</sub>H<sub>8</sub>O).



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

- Molecule 6 is PHENOL (three-letter code: IPH) (formula:  $C_6H_6O$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			7	6	1		
6	A	1	Total	C	O	0	0
			7	6	1		
6	A	1	Total	C	O	0	0
			7	6	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			7	6	1		

- Molecule 7 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	O	P	0	0
			5	4	1		
7	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	237	Total	O	0	0
			237	237		
8	B	25	Total	O	0	0
			25	25		
8	D	17	Total	O	0	0
			17	17		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	55.69Å 116.56Å 69.61Å 90.00° 92.43° 90.00°	Depositor
Resolution (Å)	35.30 – 2.30 35.31 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.3 (35.30-2.30) 98.3 (35.31-2.30)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.54 (at 2.31Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, $R_{free}$	0.169 , 0.219 0.183 , 0.226	Depositor DCC
$R_{free}$ test set	1881 reflections (5.12%)	DCC
Wilson B-factor (Å <sup>2</sup> )	32.4	Xtriage
Anisotropy	0.522	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 51.3	EDS
Estimated twinning fraction	0.042 for h,-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 38590 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7321	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.17% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, MG, IPA, IPH

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	0.68	0/6581	0.71	1/8906 (0.0%)
2	B	1.20	2/419 (0.5%)	1.29	5/645 (0.8%)
3	D	1.06	1/211 (0.5%)	1.47	3/327 (0.9%)
All	All	0.73	3/7211 (0.0%)	0.80	9/9878 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	U	OP3-P	-13.22	1.45	1.61
2	B	8	U	C2-N3	-5.36	1.33	1.37
3	D	7	G	C8-N7	5.24	1.34	1.30

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	375	ARG	NE-CZ-NH2	-6.42	117.09	120.30
3	D	7	G	C5-C6-O6	-6.29	124.82	128.60
2	B	8	U	N3-C4-O4	-6.24	115.03	119.40
3	D	8	A	N1-C6-N6	5.87	122.12	118.60
2	B	8	U	C5-C4-O4	5.46	129.17	125.90
2	B	7	U	C5-C4-O4	-5.35	122.69	125.90
2	B	8	U	N1-C2-N3	5.34	118.10	114.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	7	G	C5-C6-N1	5.24	114.12	111.50
2	B	1	U	N3-C2-O2	-5.14	118.60	122.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	695	GLU	Peptide

## 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	6429	0	6492	63	2
2	B	379	0	193	5	0
3	D	189	0	99	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	D	1	0	0	0	0
5	A	4	0	8	0	0
6	A	28	0	21	2	0
7	A	5	0	0	0	0
7	B	5	0	0	0	0
8	A	237	0	0	2	2
8	B	25	0	0	0	0
8	D	17	0	0	0	0
All	All	7321	0	6813	64	2

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 5.

All (64) 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:695:GLU:HG2	1:A:696:LYS:HB2	1.31	1.13

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:127:ILE:HG22	1:A:127:ILE:O	1.59	0.98
1:A:695:GLU:HG2	1:A:696:LYS:CB	1.99	0.92
1:A:61:ILE:HD13	1:A:74:ILE:HD11	1.71	0.71
1:A:61:ILE:CD1	1:A:74:ILE:HD11	2.21	0.71
1:A:127:ILE:O	1:A:127:ILE:CG2	2.35	0.68
1:A:86:ILE:HG22	1:A:87:PHE:N	2.09	0.67
1:A:86:ILE:HG22	1:A:87:PHE:CD2	2.32	0.64
1:A:78:MET:HE1	1:A:130:VAL:HB	1.81	0.63
1:A:689:GLU:OE1	1:A:693:LYS:NZ	2.32	0.58
1:A:696:LYS:HG3	1:A:696:LYS:O	2.04	0.57
1:A:86:ILE:HG22	1:A:87:PHE:CG	2.40	0.57
1:A:71:ASN:O	1:A:74:ILE:HG13	2.05	0.57
1:A:395:ARG:HA	1:A:395:ARG:NE	2.20	0.56
1:A:243:ILE:HG12	1:A:244:GLU:H	1.72	0.55
1:A:60:ASP:OD1	1:A:131:SER:OG	2.26	0.54
1:A:238:LEU:HD11	1:A:250:LEU:HD21	1.90	0.53
1:A:692:ILE:HG12	6:A:903:IPH:H4	1.90	0.53
1:A:801:ALA:N	1:A:802:PRO:HD2	2.24	0.52
1:A:311:TYR:CZ	1:A:315:ARG:HD3	2.46	0.51
1:A:86:ILE:CG2	1:A:87:PHE:N	2.72	0.51
1:A:279:TYR:HB3	1:A:330:VAL:HB	1.92	0.51
1:A:250:LEU:CD1	1:A:325:LEU:HD22	2.41	0.51
2:B:3:C:H2'	2:B:4:A:C8	2.47	0.50
1:A:282:CYS:SG	1:A:333:GLU:HG3	2.52	0.50
1:A:185:SER:HG	1:A:188:CYS:HG	1.59	0.49
1:A:250:LEU:CD1	1:A:325:LEU:CD2	2.91	0.49
1:A:182:PHE:CE1	1:A:222:THR:HG22	2.48	0.49
1:A:521:ILE:HD12	1:A:552:VAL:HG21	1.94	0.49
1:A:57:TYR:HE1	1:A:108:ILE:HG21	1.79	0.48
1:A:143:LEU:HB2	1:A:158:THR:HG21	1.97	0.47
1:A:311:TYR:OH	1:A:315:ARG:NH1	2.48	0.47
1:A:59:LEU:O	1:A:98:LYS:HG2	2.16	0.46
1:A:85:GLN:HG2	1:A:85:GLN:O	2.15	0.46
1:A:78:MET:HE1	1:A:130:VAL:CB	2.46	0.46
1:A:231:ILE:HG23	1:A:243:ILE:HD12	1.96	0.46
1:A:52:ILE:HG22	1:A:53:ASP:N	2.31	0.46
1:A:68:ARG:NH1	2:B:14:A:OP1	2.45	0.46
1:A:86:ILE:HG22	1:A:87:PHE:H	1.81	0.45
1:A:468:LYS:NZ	8:A:1012:HOH:O	2.41	0.45
1:A:699:GLN:HG2	6:A:903:IPH:C4	2.46	0.45
1:A:102:THR:HG21	1:A:106:LEU:HD21	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:696:LYS:CG	1:A:696:LYS:O	2.66	0.44
1:A:145:ASP:HB3	1:A:150:ARG:HB2	1.99	0.44
1:A:336:HIS:ND1	2:B:21:U:O2	2.51	0.44
1:A:696:LYS:O	1:A:697:ASP:CB	2.65	0.44
1:A:395:ARG:HE	1:A:395:ARG:HA	1.83	0.44
1:A:559:THR:HG23	2:B:2:U:C6	2.54	0.43
1:A:86:ILE:CG2	1:A:87:PHE:CD2	3.01	0.42
1:A:565:LEU:O	1:A:569:VAL:HG23	2.19	0.42
1:A:372:ALA:HB3	1:A:373:PRO:HD3	2.02	0.42
1:A:169:LEU:HD21	1:A:173:ARG:NH2	2.34	0.42
1:A:113:VAL:O	1:A:131:SER:HA	2.20	0.41
1:A:64:GLU:OE1	1:A:64:GLU:N	2.54	0.41
1:A:57:TYR:CE1	1:A:108:ILE:HG21	2.56	0.41
1:A:337:THR:HA	2:B:21:U:H1'	2.03	0.41
1:A:243:ILE:HG23	1:A:244:GLU:HG3	2.03	0.41
1:A:555:THR:HA	8:A:1180:HOH:O	2.19	0.41
1:A:199:TRP:HB2	1:A:224:PHE:CD1	2.56	0.41
1:A:250:LEU:HD11	1:A:325:LEU:HD22	2.03	0.40
1:A:45:PHE:CZ	1:A:383:MET:HG3	2.56	0.40
1:A:86:ILE:CG2	1:A:87:PHE:H	2.34	0.40
1:A:56:HIS:NE2	1:A:99:ASN:OD1	2.45	0.40
1:A:695:GLU:CG	1:A:696:LYS:CB	2.87	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:468:LYS:NZ	8:A:1222:HOH:O[2_746]	0.49	1.71
1:A:468:LYS:CE	8:A:1222:HOH:O[2_746]	1.06	1.14

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	793/859 (92%)	758 (96%)	34 (4%)	1 (0%)	56 68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	149	GLY

### 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	708/752 (94%)	700 (99%)	8 (1%)	80 90

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	112	LYS
1	A	134	TRP
1	A	345	ASN
1	A	548	GLN
1	A	658	ARG
1	A	746	PHE
1	A	804	TYR
1	A	815	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	345	ASN
1	A	461	GLN

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	16/21 (76%)	0	0
3	D	8/11 (72%)	0	0
All	All	24/32 (75%)	0	0

There are no RNA backbone outliers to report.

There are no RNA pucker outliers to report.

## 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 10 ligands modelled in this entry, 3 are monoatomic - leaving 7 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	IPA	A	902	-	3,3,3	0.58	0	3,3,3	0.20	0
6	IPH	A	903	-	7,7,7	0.44	0	8,8,8	0.61	0
6	IPH	A	904	-	7,7,7	0.37	0	8,8,8	0.69	0
6	IPH	A	905	-	7,7,7	0.33	0	8,8,8	0.48	0
6	IPH	A	906	-	7,7,7	0.54	0	8,8,8	0.39	0
7	PO4	A	907	-	4,4,4	0.57	0	6,6,6	0.25	0
7	PO4	B	102	-	4,4,4	0.67	0	6,6,6	0.23	0

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
5	IPA	A	902	-	-	0/0/0/0	0/0/0/0
6	IPH	A	903	-	-	0/0/0/0	0/1/1/1
6	IPH	A	904	-	-	0/0/0/0	0/1/1/1
6	IPH	A	905	-	-	0/0/0/0	0/1/1/1
6	IPH	A	906	-	-	0/0/0/0	0/1/1/1
7	PO4	A	907	-	-	0/0/0/0	0/0/0/0
7	PO4	B	102	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

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
6	A	903	IPH	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	803/859 (93%)	0.33	74 (9%) <span>11</span> <span>16</span>	17, 38, 88, 136	0
2	B	19/21 (90%)	0.03	0 <span>100</span> <span>100</span>	22, 37, 106, 116	0
3	D	9/11 (81%)	-0.49	0 <span>100</span> <span>100</span>	31, 33, 46, 75	0
All	All	831/891 (93%)	0.32	74 (8%) <span>12</span> <span>17</span>	17, 37, 89, 136	0

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	311	TYR	6.2
1	A	423	ARG	5.8
1	A	248	LYS	5.5
1	A	86	ILE	4.9
1	A	65	LYS	4.8
1	A	85	GLN	4.6
1	A	149	GLY	4.5
1	A	150	ARG	4.4
1	A	134	TRP	4.4
1	A	243	ILE	4.3
1	A	246	GLN	4.3
1	A	152	PRO	4.2
1	A	312	PHE	4.1
1	A	108	ILE	3.9
1	A	331	GLY	3.9
1	A	84	THR	3.8
1	A	110	ARG	3.7
1	A	424	ASN	3.7
1	A	89	ASP	3.7
1	A	87	PHE	3.7
1	A	133	LYS	3.6
1	A	315	ARG	3.6
1	A	64	GLU	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	313	LYS	3.5
1	A	23	PHE	3.4
1	A	355	LYS	3.4
1	A	88	GLY	3.4
1	A	142	ALA	3.4
1	A	244	GLU	3.3
1	A	151	LEU	3.3
1	A	55	TYR	3.1
1	A	247	GLN	3.0
1	A	338	TYR	3.0
1	A	321	ARG	3.0
1	A	116	GLU	2.9
1	A	119	LEU	2.9
1	A	52	ILE	2.9
1	A	104	MET	2.9
1	A	140	LEU	2.9
1	A	61	ILE	2.8
1	A	250	LEU	2.8
1	A	296	LEU	2.8
1	A	241	LYS	2.8
1	A	295	PRO	2.8
1	A	803	ALA	2.7
1	A	335	LYS	2.7
1	A	318	LEU	2.6
1	A	154	VAL	2.6
1	A	120	PRO	2.6
1	A	319	VAL	2.6
1	A	336	HIS	2.5
1	A	808	LEU	2.5
1	A	83	LYS	2.5
1	A	342	GLU	2.5
1	A	153	SER	2.5
1	A	239	ASP	2.5
1	A	293	THR	2.4
1	A	307	THR	2.4
1	A	132	ILE	2.4
1	A	107	PRO	2.4
1	A	294	PHE	2.4
1	A	118	THR	2.3
1	A	316	HIS	2.3
1	A	115	LEU	2.3
1	A	279	TYR	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	337	THR	2.2
1	A	282	CYS	2.2
1	A	114	GLU	2.2
1	A	106	LEU	2.2
1	A	117	VAL	2.1
1	A	63	PRO	2.1
1	A	334	GLN	2.1
1	A	288	PRO	2.0
1	A	130	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 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
6	IPH	A	903	7/7	0.90	0.18	4.29	66,67,67,68	0
5	IPA	A	902	4/4	0.92	0.14	2.05	49,52,52,56	0
6	IPH	A	904	7/7	0.98	0.16	1.88	36,36,39,43	0
6	IPH	A	906	7/7	0.97	0.14	0.08	32,33,35,36	0
6	IPH	A	905	7/7	0.96	0.10	-0.12	35,37,39,39	0
4	MG	A	901	1/1	0.91	0.10	-1.46	34,34,34,34	0
7	PO4	B	102	5/5	0.86	0.20	-	142,143,143,143	0
7	PO4	A	907	5/5	0.71	0.27	-	112,113,113,114	0
4	MG	B	101	1/1	0.95	0.15	-	52,52,52,52	0
4	MG	D	101	1/1	0.93	0.05	-	35,35,35,35	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.