



# Full wwPDB X-ray Structure Validation Report ⓘ

May 18, 2020 – 08:03 pm BST

PDB ID : 5JS1  
Title : Human Argonaute2 Bound to an siRNA  
Authors : Schirle, N.T.; MacRae, I.J.  
Deposited on : 2016-05-07  
Resolution : 2.50 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
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) : 2.11

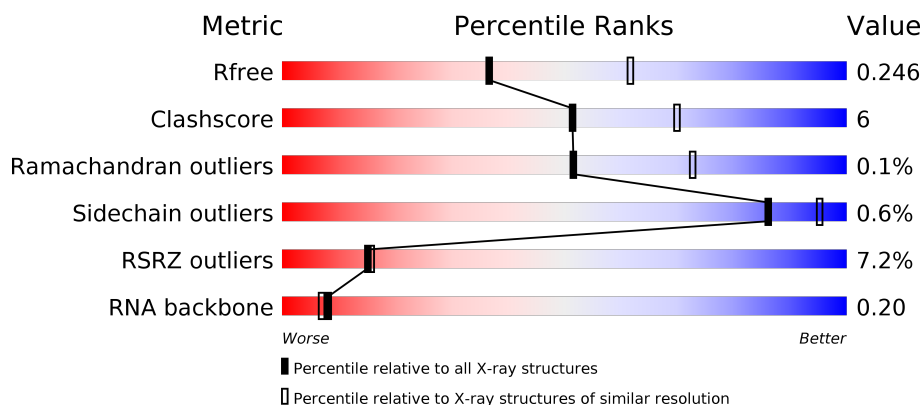
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)
RNA backbone	3102	1008 (2.84-2.16)

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 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	859	
2	B	21	

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6715 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	801	Total	C	N	O	S	0	0	0
			6439	4099	1160	1141	39			

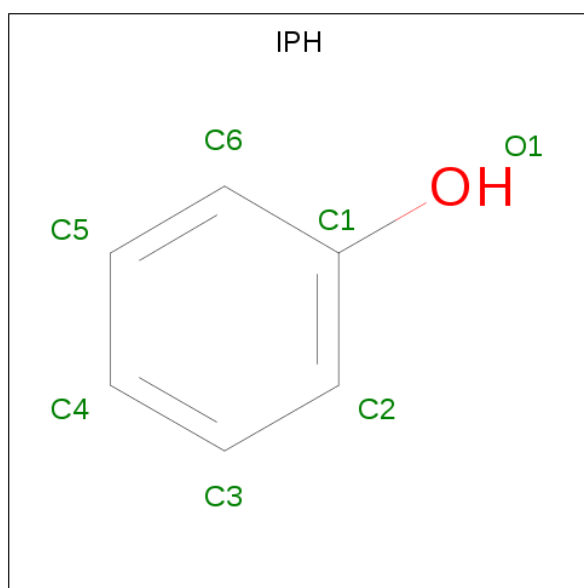
There is a discrepancy between the modelled and reference sequences:

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

- Molecule 2 is a RNA chain called siRNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	10	Total	C	N	O	P	0	0	0
			209	94	33	72	10			

- Molecule 3 is PHENOL (three-letter code: IPH) (formula: C<sub>6</sub>H<sub>6</sub>O).



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

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

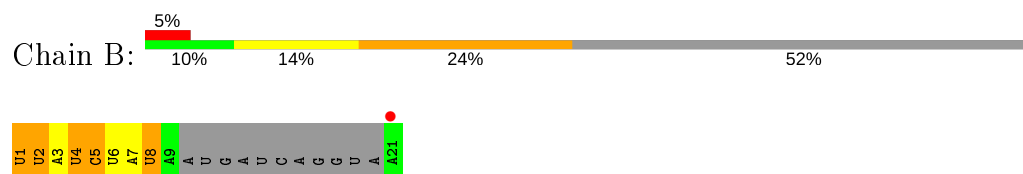
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Mg	0	0
			1	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	46	Total	O	0	0
			46	46		
5	B	6	Total	O	0	0
			6	6		



- Molecule 1: Protein argonaute-2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	62.93 Å 107.12 Å 68.20 Å 90.00° 106.98° 90.00°	Depositor
Resolution (Å)	31.53 – 2.50 38.93 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.0 (31.53-2.50) 98.0 (38.93-2.50)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	0.04	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.63 (at 2.51 Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, $R_{free}$	0.216 , 0.247 0.215 , 0.246	Depositor DCC
$R_{free}$ test set	1491 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	49.0	Xtriage
Anisotropy	0.189	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 50.4	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	0.93	EDS
Total number of atoms	6715	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.25% 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 [i](#)

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

Bond lengths and bond angles in the following residue types are not validated in this section: IPH, 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	0.46	2/6589 (0.0%)	0.51	4/8915 (0.0%)
2	B	1.67	4/231 (1.7%)	1.95	16/352 (4.5%)
All	All	0.55	6/6820 (0.1%)	0.63	20/9267 (0.2%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	U	OP3-P	-14.04	1.44	1.61
1	A	700	PRO	N-CD	13.81	1.67	1.47
1	A	674	GLY	C-O	8.06	1.36	1.23
2	B	5	C	O3'-P	-7.19	1.52	1.61
2	B	4	U	O3'-P	-6.03	1.53	1.61
2	B	3	A	O3'-P	-5.08	1.55	1.61

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	8	U	N1-C1'-C2'	-8.69	102.44	112.00
2	B	2	U	N1-C1'-C2'	-7.98	103.22	112.00
2	B	2	U	C1'-C2'-O2'	-7.71	87.48	110.60
2	B	7	A	N9-C1'-C2'	-7.70	103.53	112.00
2	B	2	U	C4'-C3'-O3'	7.35	127.69	113.00
1	A	674	GLY	CA-C-O	6.93	133.07	120.60
2	B	5	C	N1-C1'-C2'	-6.84	104.48	112.00
2	B	7	A	C4'-C3'-O3'	6.25	125.51	113.00
2	B	5	C	C1'-C2'-O2'	-6.21	91.98	110.60
2	B	3	A	C1'-C2'-O2'	-6.13	92.20	110.60
2	B	3	A	N9-C1'-C2'	-6.12	105.27	112.00
1	A	673	GLU	O-C-N	-6.10	112.83	123.20
2	B	6	U	C1'-C2'-O2'	-5.88	92.97	110.60

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	6	U	N1-C1'-C2'	-5.75	105.68	112.00
2	B	3	A	C4'-C3'-O3'	5.68	124.35	113.00
1	A	248	LYS	C-N-CD	5.61	140.18	128.40
2	B	1	U	C4'-C3'-O3'	-5.59	97.65	109.40
2	B	4	U	C4'-C3'-O3'	5.56	124.12	113.00
1	A	742	HIS	C-N-CD	5.48	139.90	128.40
2	B	4	U	C1'-C2'-O2'	-5.07	95.39	110.60

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	6439	0	6499	73	0
2	B	209	0	107	1	0
3	A	14	0	11	1	0
4	A	1	0	0	0	0
5	A	46	0	0	3	0
5	B	6	0	0	0	0
All	All	6715	0	6617	74	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 6.

All (74) 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:211:TRP:O	1:A:212:LYS:HG2	1.28	1.26
1:A:269:ILE:HG22	1:A:271:HIS:HB2	1.39	1.04
1:A:210:LEU:HG	1:A:211:TRP:HD1	1.29	0.97
1:A:211:TRP:O	1:A:212:LYS:CG	2.12	0.97
1:A:676:PHE:CZ	1:A:768:LEU:HD21	2.06	0.91
1:A:210:LEU:HG	1:A:211:TRP:CD1	2.06	0.89
1:A:269:ILE:CG2	1:A:271:HIS:HB2	2.04	0.87

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:211:TRP:C	1:A:212:LYS:HG2	2.01	0.79
1:A:196:ARG:NH1	1:A:352:CYS:SG	2.57	0.77
1:A:329:GLN:HE22	1:A:333:GLU:HG3	1.50	0.77
1:A:478:SER:OG	1:A:483:MET:O	2.03	0.75
1:A:676:PHE:CE2	1:A:768:LEU:HD21	2.21	0.75
1:A:638:ILE:HD13	1:A:679:VAL:HG22	1.72	0.72
1:A:660:LYS:NZ	1:A:695:GLU:OE1	2.23	0.71
1:A:638:ILE:CD1	1:A:679:VAL:HG22	2.22	0.70
1:A:676:PHE:CE2	1:A:768:LEU:CD2	2.75	0.70
1:A:110:ARG:O	1:A:110:ARG:HG2	1.90	0.69
1:A:493:LYS:HG3	1:A:494:TYR:O	1.93	0.69
1:A:370:ARG:NH1	1:A:378:GLU:OE2	2.28	0.66
1:A:329:GLN:NE2	1:A:333:GLU:HG3	2.11	0.65
1:A:210:LEU:CG	1:A:211:TRP:HD1	2.07	0.63
1:A:62:LYS:HG3	1:A:129:LYS:HB2	1.82	0.62
1:A:583:ARG:NH1	1:A:620:ALA:O	2.32	0.62
1:A:755:GLY:N	5:A:1003:HOH:O	2.34	0.59
1:A:269:ILE:HG22	1:A:271:HIS:CB	2.24	0.59
1:A:384:ARG:NH2	5:A:1001:HOH:O	2.28	0.59
1:A:629:VAL:HG22	1:A:846:VAL:HG11	1.85	0.58
1:A:245:GLU:O	1:A:246:GLN:HB3	2.06	0.56
1:A:436:ASP:OD2	1:A:438:ARG:NE	2.35	0.55
1:A:817:LEU:HD21	1:A:846:VAL:HG23	1.88	0.54
1:A:475:ARG:NH2	1:A:487:GLY:O	2.41	0.54
1:A:112:LYS:CA	1:A:133:LYS:HG3	2.38	0.53
1:A:531:GLU:OE1	1:A:534:ARG:NH2	2.41	0.53
2:B:4:U:H2'	2:B:5:C:C6	2.44	0.52
1:A:817:LEU:HD21	1:A:846:VAL:CG2	2.39	0.52
1:A:216:ASN:ND2	5:A:1009:HOH:O	2.42	0.51
1:A:62:LYS:HA	1:A:63:PRO:C	2.30	0.51
1:A:676:PHE:HE2	1:A:768:LEU:HD22	1.76	0.51
1:A:240:PHE:CZ	1:A:246:GLN:HG2	2.46	0.50
1:A:631:VAL:HG13	1:A:846:VAL:HG21	1.93	0.50
1:A:90:ARG:HH12	1:A:107:PRO:HD3	1.77	0.49
1:A:676:PHE:CE2	1:A:768:LEU:HD22	2.48	0.49
1:A:113:VAL:O	1:A:131:SER:HA	2.13	0.48
1:A:110:ARG:CG	1:A:110:ARG:O	2.58	0.48
1:A:169:LEU:HB3	1:A:170:PRO:HD3	1.96	0.47
1:A:475:ARG:HH22	1:A:488:GLN:HA	1.79	0.47
1:A:676:PHE:HE2	1:A:768:LEU:CD2	2.28	0.47
1:A:243:ILE:HG13	1:A:243:ILE:O	2.13	0.47

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:419:LEU:HD12	1:A:420:TYR:N	2.30	0.46
1:A:419:LEU:HD12	1:A:420:TYR:O	2.15	0.46
1:A:422:GLY:HA3	1:A:441:GLN:HB2	1.96	0.46
1:A:112:LYS:H	1:A:133:LYS:HE3	1.80	0.46
1:A:607:LYS:O	1:A:607:LYS:HG2	2.16	0.46
1:A:676:PHE:HZ	1:A:768:LEU:HD21	1.73	0.46
1:A:739:LYS:HE3	1:A:739:LYS:HB3	1.72	0.46
1:A:112:LYS:HA	1:A:133:LYS:HG3	1.99	0.45
1:A:424:ASN:O	1:A:425:LYS:HB2	2.17	0.45
1:A:94:PHE:CE2	1:A:96:GLY:HA2	2.52	0.45
1:A:675:GLN:O	1:A:679:VAL:HG23	2.16	0.44
1:A:455:CYS:HA	1:A:521:ILE:HB	2.00	0.44
1:A:511:THR:HG23	1:A:512:TYR:CD2	2.52	0.44
1:A:688:ARG:NH2	1:A:769:TRP:NE1	2.66	0.44
1:A:269:ILE:CG2	1:A:271:HIS:CB	2.86	0.44
1:A:651:ILE:HD11	1:A:694:LEU:HD11	1.99	0.44
1:A:210:LEU:CG	1:A:211:TRP:CD1	2.90	0.43
1:A:698:TYR:CZ	3:A:901:IPH:H4	2.54	0.42
1:A:269:ILE:HD13	1:A:269:ILE:HA	1.83	0.42
1:A:49:ILE:HD12	1:A:213:MET:HG3	2.01	0.41
1:A:384:ARG:HD2	1:A:384:ARG:HH11	1.74	0.41
1:A:104:MET:HB3	1:A:105:PRO:HD2	2.02	0.41
1:A:167:ARG:HB2	1:A:181:PHE:HZ	1.85	0.41
1:A:524:GLY:HA2	1:A:548:GLN:HE21	1.86	0.41
1:A:282:CYS:SG	1:A:329:GLN:HG2	2.61	0.41
1:A:177:VAL:HG11	1:A:351:ARG:HD3	2.03	0.41

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	787/859 (92%)	768 (98%)	18 (2%)	1 (0%)	51 73

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	523	PRO

### 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	712/752 (95%)	708 (99%)	4 (1%)	86 95

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	271	HIS
1	A	709	LYS
1	A	795	ARG
1	A	804	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	292	GLN
1	A	329	GLN

### 5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	9/21 (42%)	2 (22%)	1 (11%)

All (2) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	2	U
2	B	8	U

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	1	U

## 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 carbohydrates 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	$\# Z  > 2$
3	IPH	A	901	-	7,7,7	0.32	0	8,8,8	0.28	0
3	IPH	A	902	-	7,7,7	0.71	0	8,8,8	0.49	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
3	IPH	A	901	-	-	-	0/1/1/1
3	IPH	A	902	-	-	-	0/1/1/1

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 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	901	IPH	1	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	801/859 (93%)	0.50	57 (7%) 16 16	26, 55, 89, 123	0
2	B	10/21 (47%)	0.41	1 (10%) 7 6	43, 55, 87, 104	0
All	All	811/880 (92%)	0.49	58 (7%) 15 16	26, 55, 90, 123	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	423	ARG	6.8
1	A	151	LEU	6.7
1	A	271	HIS	4.6
1	A	110	ARG	4.5
2	B	21	A	4.3
1	A	486	GLN	4.3
1	A	451	TRP	4.1
1	A	490	CYS	3.9
1	A	334	GLN	3.9
1	A	247	GLN	3.8
1	A	211	TRP	3.7
1	A	449	LYS	3.6
1	A	332	GLN	3.6
1	A	452	ALA	3.6
1	A	485	ILE	3.5
1	A	787	CYS	3.5
1	A	448	ILE	3.4
1	A	819	ASP	3.4
1	A	85	GLN	3.3
1	A	838	ASP	3.3
1	A	104	MET	3.3
1	A	514	GLY	3.2
1	A	513	ALA	3.1
1	A	120	PRO	3.1

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	A	512	TYR	3.1
1	A	488	GLN	3.0
1	A	150	ARG	3.0
1	A	113	VAL	2.9
1	A	607	LYS	2.9
1	A	318	LEU	2.9
1	A	439	ASN	2.8
1	A	468	LYS	2.8
1	A	438	ARG	2.7
1	A	460	ARG	2.6
1	A	462	CYS	2.6
1	A	127	ILE	2.6
1	A	156	PHE	2.6
1	A	447	GLU	2.6
1	A	475	ARG	2.6
1	A	128	PHE	2.5
1	A	511	THR	2.5
1	A	493	LYS	2.5
1	A	494	TYR	2.3
1	A	425	LYS	2.3
1	A	149	GLY	2.3
1	A	487	GLY	2.2
1	A	115	LEU	2.2
1	A	277	ARG	2.2
1	A	111	ASP	2.2
1	A	496	GLN	2.1
1	A	839	HIS	2.1
1	A	55	TYR	2.1
1	A	757	GLN	2.1
1	A	841	ALA	2.1
1	A	353	ILE	2.0
1	A	677	GLN	2.0
1	A	505	PHE	2.0
1	A	484	PRO	2.0

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

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. 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
3	IPH	A	902	7/7	0.89	0.43	110,111,112,112	0
3	IPH	A	901	7/7	0.92	0.27	51,52,57,64	0
4	MG	A	903	1/1	0.99	0.12	45,45,45,45	0

### 6.5 Other polymers [i](#)

There are no such residues in this entry.