



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:12 AM UTC

PDB ID : 7ADI / pdb_00007adi
Title : KirBac3.1 W46R: role of a highly conserved tryptophan at the membrane-water interface of Kir channel
Authors : Venien-Bryan, C.; Fagnen, C.; De Zorzi, R.; Bannwarth, L.; Oubella, I.; Haouz, A.
Deposited on : 2020-09-15
Resolution : 2.80 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

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>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

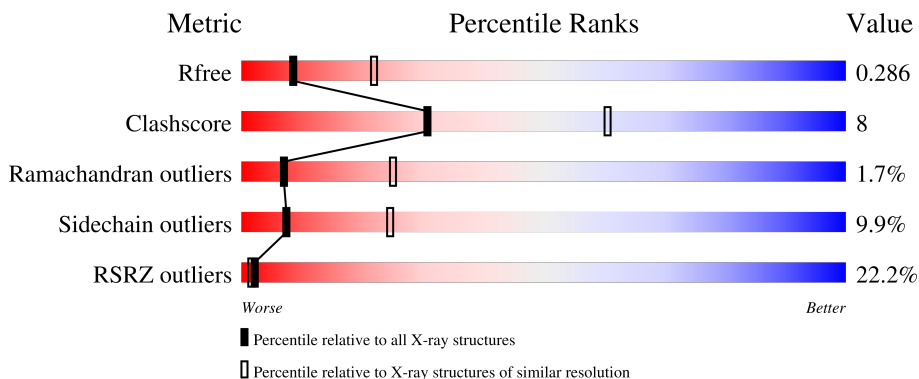
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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 ≥ 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	301	17% 70% 19% 8%
1	B	301	23% 61% 24% 12%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4303 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 Inward rectifier potassium channel Kirbac3.1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	276	Total	C	N	O	S	0	3	0
			2194	1403	390	391	10			
1	B	265	Total	C	N	O	S	0	0	0
			2082	1337	368	367	10			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	46	ARG	TRP	engineered mutation	UNP D9N164
A	296	HIS	-	expression tag	UNP D9N164
A	297	HIS	-	expression tag	UNP D9N164
A	298	HIS	-	expression tag	UNP D9N164
A	299	HIS	-	expression tag	UNP D9N164
A	300	HIS	-	expression tag	UNP D9N164
A	301	HIS	-	expression tag	UNP D9N164
B	46	ARG	TRP	engineered mutation	UNP D9N164
B	296	HIS	-	expression tag	UNP D9N164
B	297	HIS	-	expression tag	UNP D9N164
B	298	HIS	-	expression tag	UNP D9N164
B	299	HIS	-	expression tag	UNP D9N164
B	300	HIS	-	expression tag	UNP D9N164
B	301	HIS	-	expression tag	UNP D9N164

- Molecule 2 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	K	0	0
			3	3		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

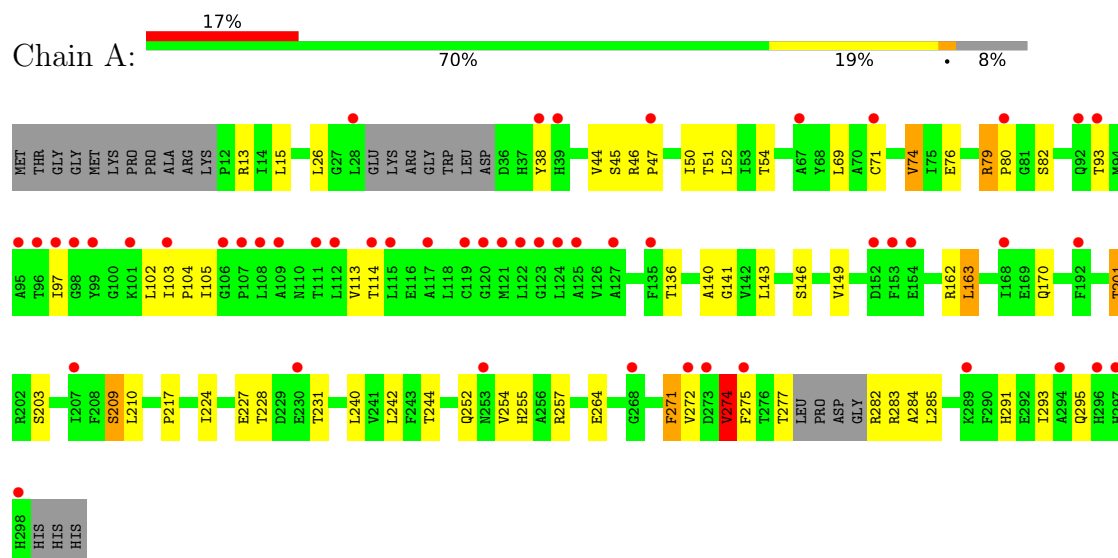
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	8	Total 8	O 8	0	0
4	B	15	Total 15	O 15	0	0

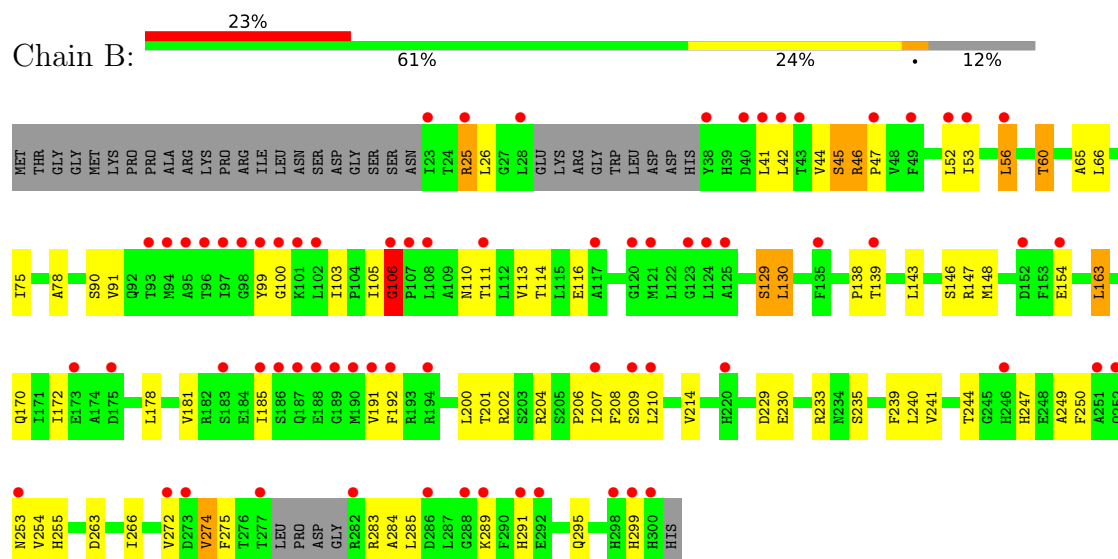
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: Inward rectifier potassium channel Kirbac3.1



- Molecule 1: Inward rectifier potassium channel Kirbac3.1



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	106.77Å 113.98Å 89.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.02 – 2.80 48.02 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.7 (48.02-2.80) 99.8 (48.02-2.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.56 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.8.0266, BUSTER	Depositor
R, R_{free}	0.222 , 0.287 0.224 , 0.286	Depositor DCC
R_{free} test set	1372 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	89.3	Xtriage
Anisotropy	0.095	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 67.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4303	wwPDB-VP
Average B, all atoms (Å ²)	105.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.22% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: K, 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	1.04	0/2249	1.48	5/3054 (0.2%)
1	B	1.08	0/2134	1.46	2/2897 (0.1%)
All	All	1.06	0/4383	1.47	7/5951 (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
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	79	ARG	CB-CA-C	5.57	117.38	109.08
1	A	271	PHE	CA-C-O	-5.56	114.88	121.10
1	B	274	VAL	N-CA-C	-5.53	107.59	112.90
1	A	44	VAL	CA-C-O	-5.47	116.27	121.64
1	B	201	THR	CA-CB-OG1	-5.46	101.41	109.60
1	A	97	ILE	N-CA-C	-5.10	106.10	110.74
1	A	274	VAL	N-CA-C	-5.07	107.90	113.43

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	141	GLY	Peptide
1	B	106	GLY	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	2194	0	2148	41	0
1	B	2082	0	2044	39	0
2	A	3	0	0	0	0
3	A	1	0	0	0	0
4	A	8	0	0	0	0
4	B	15	0	0	0	0
All	All	4303	0	4192	70	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 8.

All (70) 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:114:THR:HG22	1:B:91:VAL:HG11	1.58	0.84
1:B:244:THR:HG22	1:B:255:HIS:CD2	2.18	0.78
1:A:257:ARG:NH2	1:B:170:GLN:OE1	2.18	0.77
1:B:56:LEU:O	1:B:60:THR:OG1	2.03	0.76
1:A:170:GLN:HE21	1:A:209:SER:HB3	1.55	0.71
1:A:244:THR:HG22	1:A:255[A]:HIS:CD2	2.26	0.69
1:B:44:VAL:HG22	1:B:45:SER:H	1.58	0.69
1:A:71:CYS:O	1:A:74:VAL:CG2	2.42	0.67
1:A:13:ARG:NH2	1:A:264:GLU:OE1	2.29	0.66
1:A:103:ILE:HG23	1:A:104:PRO:HD2	1.78	0.65
1:B:244:THR:CG2	1:B:255:HIS:CD2	2.80	0.64
1:A:224:ILE:HA	1:A:227:GLU:HG3	1.79	0.64
1:B:163:LEU:HD12	1:B:208:PHE:CZ	2.35	0.61
1:B:46:ARG:HB3	1:B:47:PRO:HD3	1.82	0.61
1:A:146:SER:H	1:A:291:HIS:HD2	1.48	0.61
1:B:146:SER:H	1:B:291:HIS:HD2	1.51	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:275:PHE:HA	1:A:284:ALA:O	2.03	0.57
1:B:138:PRO:HB2	1:B:250:PHE:CD2	2.40	0.56
1:A:103:ILE:CG2	1:A:104:PRO:HD2	2.35	0.55
1:A:79:ARG:O	1:A:82:SER:HB3	2.07	0.55
1:A:46:ARG:HB3	1:A:47:PRO:HD3	1.89	0.54
1:A:93:THR:HG21	1:A:113:VAL:HG22	1.91	0.53
1:A:255[A]:HIS:CD2	1:B:172:ILE:CD1	2.92	0.53
1:A:257:ARG:HD2	1:B:207:ILE:HD11	1.91	0.53
1:B:200:LEU:HD13	1:B:214:VAL:HG13	1.91	0.52
1:B:100:GLY:O	1:B:103:ILE:CG2	2.58	0.52
1:B:274:VAL:HG23	1:B:285:LEU:HD12	1.92	0.51
1:B:148:MET:HE1	1:B:239:PHE:CG	2.46	0.51
1:B:138:PRO:HB2	1:B:250:PHE:CE2	2.46	0.51
1:A:140:ALA:HB2	1:A:254:VAL:HB	1.93	0.50
1:A:274:VAL:HG13	1:A:285:LEU:HD12	1.94	0.50
1:B:75:ILE:O	1:B:78:ALA:HB3	2.13	0.49
1:A:51:THR:O	1:A:54:THR:OG1	2.30	0.48
1:A:162:ARG:CZ	1:A:271:PHE:HB3	2.43	0.48
1:A:224:ILE:HA	1:A:227:GLU:CG	2.43	0.48
1:A:242:LEU:HD11	1:B:206:PRO:HB2	1.98	0.45
1:A:15:LEU:HD13	1:A:240:LEU:HD13	1.98	0.45
1:B:90:SER:OG	1:B:116:GLU:OE2	2.35	0.45
1:B:129:SER:O	1:B:130:LEU:C	2.60	0.44
1:A:114:THR:HG22	1:B:91:VAL:HG21	2.00	0.44
1:A:136:THR:HB	1:A:252:GLN:HE22	1.83	0.44
1:A:210:LEU:HD13	1:A:275:PHE:HE1	1.83	0.44
1:A:143:LEU:O	1:A:163:LEU:HA	2.18	0.44
1:A:50:ILE:O	1:A:54:THR:HG23	2.19	0.43
1:A:146:SER:N	1:A:291:HIS:HD2	2.12	0.43
1:B:110:ASN:O	1:B:114:THR:HG23	2.17	0.43
1:A:74:VAL:HG12	1:A:105:ILE:O	2.18	0.43
1:A:102:LEU:HD12	1:A:102:LEU:HA	1.88	0.43
1:A:201:THR:HG22	1:A:217:PRO:CG	2.49	0.42
1:A:114:THR:CG2	1:B:91:VAL:HG21	2.49	0.42
1:B:178:LEU:HD13	1:B:241:VAL:HG22	2.02	0.42
1:B:65:ALA:O	1:B:66:LEU:C	2.61	0.42
1:B:229:ASP:OD1	1:B:233:ARG:NE	2.42	0.42
1:B:191:VAL:O	1:B:191:VAL:HG13	2.20	0.41
1:B:25:ARG:HG3	1:B:26:LEU:O	2.19	0.41
1:A:69:LEU:HD21	1:A:82:SER:O	2.20	0.41
1:B:105:ILE:O	1:B:106:GLY:O	2.38	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:240:LEU:HD21	1:B:207:ILE:HD12	2.01	0.41
1:B:247:HIS:ND1	1:B:249:ALA:HB3	2.35	0.41
1:B:146:SER:N	1:B:291:HIS:HD2	2.16	0.41
1:A:79:ARG:HA	1:A:80:PRO:HD3	1.92	0.41
1:A:228:THR:OG1	1:A:231:THR:HG23	2.21	0.41
1:B:275:PHE:HA	1:B:284:ALA:O	2.21	0.41
1:B:148:MET:HE1	1:B:239:PHE:CD1	2.56	0.41
1:A:242:LEU:HD21	1:B:207:ILE:HG13	2.03	0.41
1:A:170:GLN:NE2	1:A:209:SER:HB3	2.31	0.40
1:A:255[A]:HIS:CD2	1:B:172:ILE:HD11	2.55	0.40
1:A:149:VAL:CG1	1:A:293:ILE:HD13	2.51	0.40
1:B:143:LEU:O	1:B:163:LEU:HA	2.21	0.40
1:B:274:VAL:HG23	1:B:285:LEU:CD1	2.51	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	273/301 (91%)	246 (90%)	25 (9%)	2 (1%)	18	47
1	B	259/301 (86%)	228 (88%)	24 (9%)	7 (3%)	4	15
All	All	532/602 (88%)	474 (89%)	49 (9%)	9 (2%)	7	25

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	38	TYR
1	B	45	SER
1	B	106	GLY
1	A	283	ARG
1	B	99	TYR

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Mol	Chain	Res	Type
1	B	41	LEU
1	B	129	SER
1	B	130	LEU
1	B	154	GLU

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	236/256 (92%)	222 (94%)	14 (6%)	18	48
1	B	221/256 (86%)	190 (86%)	31 (14%)	3	12
All	All	457/512 (89%)	412 (90%)	45 (10%)	7	25

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	LEU
1	A	45	SER
1	A	52	LEU
1	A	74	VAL
1	A	76	GLU
1	A	163	LEU
1	A	201	THR
1	A	203	SER
1	A	209	SER
1	A	272	VAL
1	A	274	VAL
1	A	277	THR
1	A	282	ARG
1	A	295	GLN
1	B	25	ARG
1	B	42	LEU
1	B	46	ARG
1	B	52	LEU
1	B	53	ILE

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Mol	Chain	Res	Type
1	B	56	LEU
1	B	60	THR
1	B	111	THR
1	B	113	VAL
1	B	139	THR
1	B	147	ARG
1	B	163	LEU
1	B	181	VAL
1	B	185	ILE
1	B	192	PHE
1	B	202	ARG
1	B	204	ARG
1	B	209	SER
1	B	210	LEU
1	B	230	GLU
1	B	235	SER
1	B	240	LEU
1	B	253	ASN
1	B	254	VAL
1	B	263	ASP
1	B	266	ILE
1	B	272	VAL
1	B	283	ARG
1	B	289	LYS
1	B	295	GLN
1	B	299	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	ASN
1	A	61	ASN
1	A	92	GLN
1	A	170	GLN
1	A	221	HIS
1	A	252	GLN
1	A	253	ASN
1	A	291	HIS
1	A	295	GLN
1	A	297	HIS
1	B	61	ASN
1	B	177	HIS

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Mol	Chain	Res	Type
1	B	221	HIS
1	B	253	ASN
1	B	255	HIS
1	B	291	HIS
1	B	295	GLN

5.3.3 RNA [i](#)

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

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.

No monomer is involved in short contacts.

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, 95th 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 ²)	Q<0.9
1	A	276/301 (91%)	0.91	51 (18%) 3 2	39, 100, 159, 193	3 (1%)
1	B	265/301 (88%)	1.26	69 (26%) 1 1	58, 107, 167, 181	0
All	All	541/602 (89%)	1.09	120 (22%) 2 1	39, 103, 164, 193	3 (0%)

All (120) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	111	THR	9.8
1	A	124	LEU	8.5
1	B	124	LEU	8.0
1	B	292	GLU	7.9
1	A	97	ILE	7.5
1	B	121	MET	7.4
1	A	96	THR	7.3
1	B	108	LEU	7.2
1	B	188	GLU	7.1
1	B	173	GLU	6.9
1	A	121	MET	6.8
1	B	96	THR	6.8
1	A	95	ALA	6.3
1	B	190	MET	6.1
1	B	289	LYS	5.9
1	B	111	THR	5.9
1	B	253	ASN	5.9
1	A	38	TYR	5.9
1	A	114	THR	5.8
1	B	95	ALA	5.6
1	A	296	HIS	5.4
1	B	97	ILE	5.2
1	A	273	ASP	5.2
1	A	108	LEU	5.1

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Mol	Chain	Res	Type	RSRZ
1	A	98	GLY	5.1
1	A	107	PRO	5.1
1	A	154	GLU	5.0
1	A	123	GLY	5.0
1	A	152	ASP	4.9
1	B	272	VAL	4.9
1	A	268	GLY	4.7
1	B	38	TYR	4.7
1	A	153	PHE	4.7
1	B	23	ILE	4.5
1	B	42	LEU	4.3
1	B	125	ALA	4.2
1	B	102	LEU	4.1
1	B	186	SER	4.0
1	B	299	HIS	4.0
1	B	101	LYS	3.9
1	B	252	GLN	3.9
1	A	297	HIS	3.9
1	B	93	THR	3.9
1	B	41	LEU	3.9
1	A	99	TYR	3.8
1	B	154	GLU	3.7
1	B	152	ASP	3.6
1	A	294	ALA	3.6
1	B	120	GLY	3.6
1	B	246	HIS	3.5
1	B	123	GLY	3.5
1	B	43	THR	3.4
1	A	71	CYS	3.4
1	B	192	PHE	3.3
1	A	272	VAL	3.3
1	B	183	SER	3.3
1	B	187	GLN	3.2
1	A	120	GLY	3.2
1	B	107	PRO	3.1
1	A	119	CYS	3.0
1	A	207	ILE	3.0
1	A	93	THR	2.9
1	B	298	HIS	2.9
1	B	286	ASP	2.9
1	B	185	ILE	2.9
1	B	191	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	106	GLY	2.8
1	A	47	PRO	2.8
1	A	101	LYS	2.8
1	B	25	ARG	2.8
1	B	175	ASP	2.8
1	B	98	GLY	2.8
1	A	127	ALA	2.8
1	B	273	ASP	2.8
1	A	298	HIS	2.7
1	B	47	PRO	2.7
1	A	135	PHE	2.7
1	A	275	PHE	2.7
1	B	300	HIS	2.7
1	B	277	THR	2.7
1	B	251	ALA	2.7
1	B	288	GLY	2.6
1	B	28	LEU	2.6
1	B	135	PHE	2.6
1	B	291	HIS	2.6
1	B	53	ILE	2.6
1	A	103	ILE	2.6
1	B	52	LEU	2.6
1	A	125	ALA	2.6
1	A	230	GLU	2.5
1	B	94	MET	2.5
1	B	207	ILE	2.5
1	A	117	ALA	2.5
1	B	56	LEU	2.4
1	B	49	PHE	2.4
1	B	189	GLY	2.4
1	A	289	LYS	2.4
1	B	117	ALA	2.4
1	B	40	ASP	2.3
1	A	122	LEU	2.3
1	B	194	ARG	2.3
1	A	253	ASN	2.3
1	B	139	THR	2.3
1	A	28	LEU	2.3
1	A	80	PRO	2.2
1	A	112	LEU	2.2
1	A	115	LEU	2.2
1	A	109	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	220	HIS	2.1
1	B	106	GLY	2.1
1	B	209	SER	2.1
1	A	92	GLN	2.1
1	B	210	LEU	2.1
1	B	282	ARG	2.1
1	A	168	ILE	2.1
1	A	39	HIS	2.0
1	B	100	GLY	2.0
1	A	67	ALA	2.0
1	B	99	TYR	2.0
1	A	192	PHE	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 oligosaccharides 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, 95th 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(Å ²)	Q<0.9
2	K	A	1001	1/1	0.66	0.32	110,110,110,110	1
2	K	A	1003	1/1	0.68	0.29	106,106,106,106	1
3	MG	A	1004	1/1	0.95	0.09	37,37,37,37	1
2	K	A	1002	1/1	0.96	0.12	78,78,78,78	1

6.5 Other polymers [i](#)

There are no such residues in this entry.