



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:19 PM UTC

PDB ID : 6HG4 / pdb_00006hg4
Title : Crystal Structure of the human IL-17RC ECD in complex with human IL-17F
Authors : Rondeau, J.M.; Goepfert, A.
Deposited on : 2018-08-22
Resolution : 3.32 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

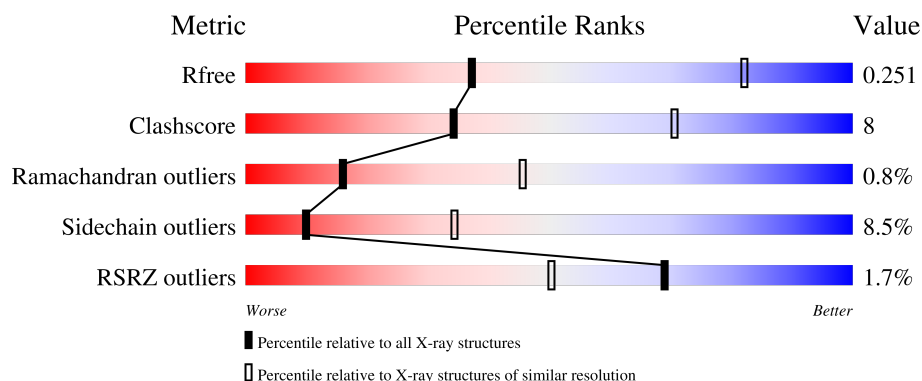
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.32 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	1153 (3.34-3.30)
Clashscore	190562	1193 (3.34-3.30)
Ramachandran outliers	187476	1172 (3.34-3.30)
Sidechain outliers	187428	1171 (3.34-3.30)
RSRZ outliers	180081	1153 (3.34-3.30)

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	139	60% 14% • 25%
2	B	451	2% 68% 23% • 5%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4176 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 Interleukin-17F.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	104	Total	C	N	O	S	0	0	0
			824	509	152	156	7			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	25	GLU	-	expression tag	UNP Q96PD4
A	26	PHE	-	expression tag	UNP Q96PD4
A	27	ARG	-	expression tag	UNP Q96PD4
A	28	HIS	-	expression tag	UNP Q96PD4
A	29	ASP	-	expression tag	UNP Q96PD4
A	30	SER	-	expression tag	UNP Q96PD4

- Molecule 2 is a protein called Interleukin-17 receptor C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	427	Total	C	N	O	S	0	0	0
			3352	2129	583	617	23			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	186	GLN	ASN	engineered mutation	UNP Q8NAC3
B	226	GLN	ASN	engineered mutation	UNP Q8NAC3
B	253	GLN	ASN	engineered mutation	UNP Q8NAC3
B	263	GLN	ASN	engineered mutation	UNP Q8NAC3
B	307	ARG	GLN	engineered mutation	UNP Q8NAC3
B	349	GLN	ASN	engineered mutation	UNP Q8NAC3
B	372	GLN	ASN	engineered mutation	UNP Q8NAC3
B	406	GLN	ASN	engineered mutation	UNP Q8NAC3
B	468	GLU	-	expression tag	UNP Q8NAC3
B	469	PHE	-	expression tag	UNP Q8NAC3

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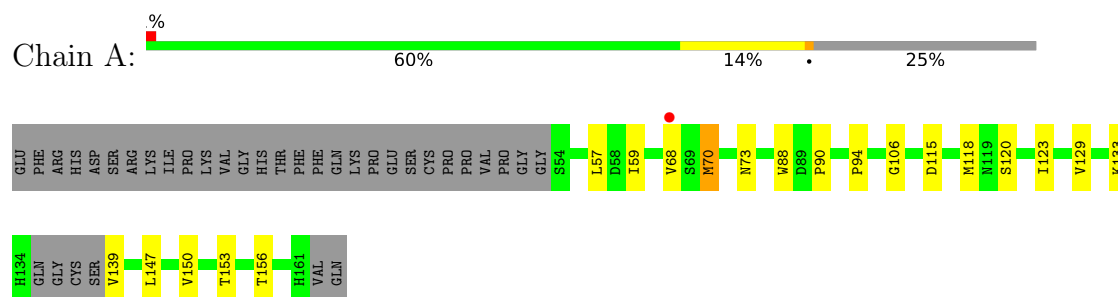
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Chain	Residue	Modelled	Actual	Comment	Reference
B	470	ARG	-	expression tag	UNP Q8NAC3
B	471	HIS	-	expression tag	UNP Q8NAC3

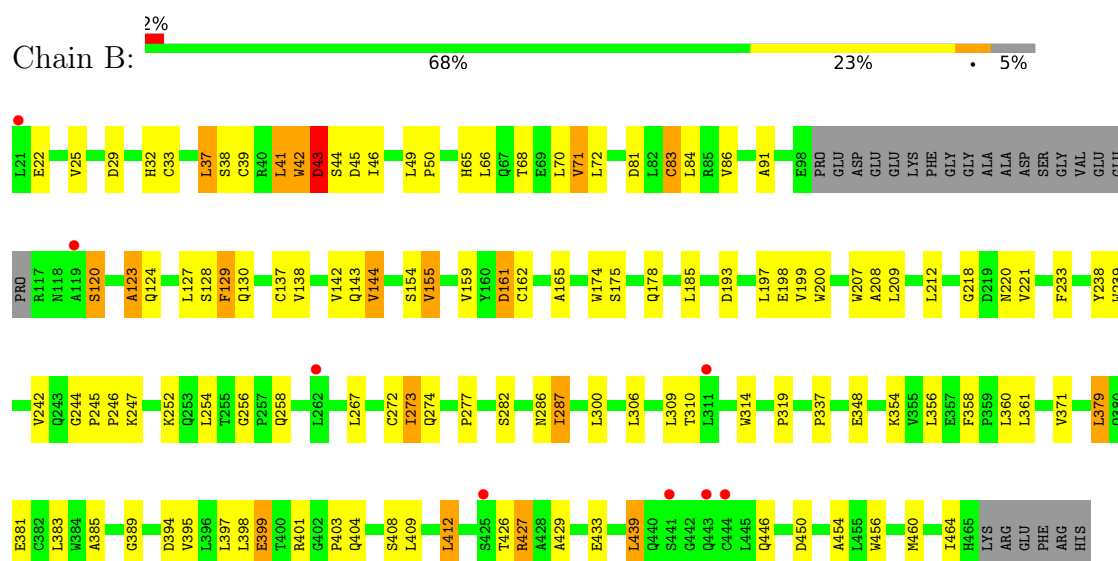
3 Residue-property plots

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: Interleukin-17F



• Molecule 2: Interleukin-17 receptor C



4 Data and refinement statistics

Property	Value	Source
Space group	P 3 2 1	Depositor
Cell constants a, b, c, α , β , γ	154.32Å 154.32Å 78.79Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	78.79 – 3.32 78.79 – 3.32	Depositor EDS
% Data completeness (in resolution range)	99.9 (78.79-3.32) 100.0 (78.79-3.32)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.23 (at 3.33Å)	Xtriage
Refinement program	BUSTER	Depositor
R, R_{free}	0.198 , 0.239 0.215 , 0.251	Depositor DCC
R_{free} test set	815 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	130.3	Xtriage
Anisotropy	0.164	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 145.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.039 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4176	wwPDB-VP
Average B, all atoms (Å ²)	164.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.98% 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

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	0.85	0/837	1.30	8/1135 (0.7%)
2	B	0.84	0/3444	1.29	25/4720 (0.5%)
All	All	0.84	0/4281	1.30	33/5855 (0.6%)

There are no bond length outliers.

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	155	VAL	N-CA-C	-9.04	104.67	113.53
2	B	245	PRO	N-CA-C	7.98	120.44	110.70
2	B	44	SER	N-CA-C	-7.25	101.32	112.99
2	B	43	ASP	CA-C-N	6.67	131.37	120.23
2	B	43	ASP	C-N-CA	6.67	131.37	120.23
2	B	161	ASP	CA-CB-CG	6.44	119.04	112.60
1	A	115	ASP	CA-C-N	6.24	130.33	120.47
1	A	115	ASP	C-N-CA	6.24	130.33	120.47
2	B	200	TRP	CA-C-N	6.16	129.38	120.38
2	B	200	TRP	C-N-CA	6.16	129.38	120.38
2	B	394	ASP	CA-CB-CG	5.85	118.45	112.60
2	B	450	ASP	CA-CB-CG	5.80	118.41	112.60
1	A	120	SER	N-CA-C	-5.56	100.74	109.25
2	B	45	ASP	CA-CB-CG	5.55	118.15	112.60
1	A	73	ASN	CA-C-N	5.49	128.27	120.53
1	A	73	ASN	C-N-CA	5.49	128.27	120.53
2	B	429	ALA	CA-C-N	5.44	127.57	120.28
2	B	429	ALA	C-N-CA	5.44	127.57	120.28
2	B	29	ASP	CA-CB-CG	5.43	118.03	112.60
2	B	22	GLU	N-CA-C	5.36	117.45	110.43
2	B	129	PHE	CA-CB-CG	5.33	119.13	113.80
1	A	156	THR	N-CA-C	-5.30	102.97	109.65
2	B	198	GLU	CA-C-N	5.27	131.45	121.97
2	B	198	GLU	C-N-CA	5.27	131.45	121.97
2	B	124	GLN	CA-C-N	5.19	129.54	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	124	GLN	C-N-CA	5.19	129.54	122.90
1	A	70	MET	CA-C-N	5.18	128.21	120.90
1	A	70	MET	C-N-CA	5.18	128.21	120.90
2	B	309	LEU	N-CA-C	-5.16	105.81	112.68
2	B	426	THR	CA-C-N	5.07	128.43	120.82
2	B	426	THR	C-N-CA	5.07	128.43	120.82
2	B	123	ALA	CA-C-N	-5.05	115.98	123.05
2	B	123	ALA	C-N-CA	-5.05	115.98	123.05

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	824	0	820	7	0
2	B	3352	0	3275	57	0
All	All	4176	0	4095	63	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 (63) 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)
2:B:33:CYS:HB3	2:B:37:LEU:HD12	1.59	0.84
2:B:42:TRP:HA	2:B:42:TRP:CE3	2.27	0.69
2:B:144:VAL:HG21	2:B:155:VAL:HB	1.76	0.68
2:B:403:PRO:HB2	2:B:404:GLN:NE2	2.12	0.64
2:B:42:TRP:HA	2:B:42:TRP:HE3	1.62	0.63
1:A:88:TRP:HE1	1:A:90:PRO:HG3	1.64	0.63
2:B:403:PRO:HB2	2:B:404:GLN:HE22	1.65	0.62
2:B:427:ARG:HD3	2:B:460:MET:HE2	1.81	0.61
2:B:32:HIS:HB2	2:B:159:VAL:HB	1.82	0.61
2:B:300:LEU:HD12	2:B:319:PRO:HD2	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:306:LEU:HD12	2:B:381:GLU:HB2	1.83	0.60
2:B:66:LEU:HD21	2:B:175:SER:HB2	1.83	0.59
2:B:385:ALA:O	2:B:389:GLY:HA3	2.04	0.58
2:B:130:GLN:HB2	2:B:174:TRP:HZ3	1.69	0.57
2:B:433:GLU:H	2:B:433:GLU:CD	2.11	0.57
1:A:88:TRP:NE1	1:A:90:PRO:HG3	2.19	0.57
2:B:123:ALA:HB3	2:B:142:VAL:HG23	1.88	0.56
2:B:129:PHE:HZ	2:B:165:ALA:HB2	1.72	0.55
2:B:274:GLN:HG3	2:B:286:ASN:OD1	2.05	0.55
2:B:371:VAL:HB	2:B:379:LEU:HB2	1.88	0.55
2:B:409:LEU:HD11	2:B:439:LEU:HD22	1.89	0.54
2:B:446:GLN:HE21	2:B:454:ALA:HB1	1.73	0.52
2:B:43:ASP:HB2	2:B:138:VAL:HA	1.92	0.52
2:B:212:LEU:HD21	2:B:273:ILE:HD13	1.92	0.52
2:B:91:ALA:HA	2:B:154:SER:HA	1.93	0.51
2:B:397:LEU:HB2	2:B:456:TRP:HB2	1.93	0.51
2:B:174:TRP:HA	2:B:185:LEU:O	2.11	0.51
2:B:71:VAL:HG12	2:B:83:CYS:HB3	1.93	0.50
2:B:218:GLY:C	2:B:220:ASN:H	2.19	0.50
1:A:90:PRO:HB3	2:B:50:PRO:HB2	1.94	0.49
2:B:161:ASP:HA	2:B:162:CYS:HB3	1.95	0.49
2:B:25:VAL:HG21	2:B:161:ASP:HB3	1.93	0.49
2:B:314:TRP:HB3	2:B:358:PHE:CE2	2.48	0.49
2:B:212:LEU:HD23	2:B:287:ILE:HG21	1.96	0.47
2:B:238:TYR:CZ	2:B:272:CYS:HB2	2.49	0.47
2:B:120:SER:HB3	2:B:143:GLN:HE21	1.79	0.47
2:B:314:TRP:CZ2	2:B:356:LEU:HD22	2.51	0.46
2:B:306:LEU:HD22	2:B:314:TRP:CZ3	2.50	0.46
2:B:399:GLU:HG2	2:B:454:ALA:HB3	1.98	0.46
2:B:306:LEU:HD21	2:B:379:LEU:HD13	1.99	0.45
2:B:239:TRP:O	2:B:247:LYS:HB2	2.16	0.45
2:B:127:LEU:HD23	2:B:127:LEU:HA	1.83	0.45
2:B:128:SER:HB3	2:B:174:TRP:HB2	1.98	0.45
2:B:207:TRP:C	2:B:209:LEU:H	2.24	0.45
2:B:37:LEU:HB3	2:B:144:VAL:HG23	1.98	0.45
2:B:244:GLY:C	2:B:246:PRO:HD2	2.42	0.44
2:B:212:LEU:CD2	2:B:287:ILE:HG21	2.48	0.44
2:B:460:MET:HE3	2:B:464:ILE:HD11	2.00	0.43
2:B:277:PRO:HG2	2:B:282:SER:HB2	2.00	0.43
2:B:252:LYS:HD3	2:B:254:LEU:HD21	2.01	0.43
2:B:239:TRP:HB2	2:B:247:LYS:HB3	1.99	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:123:ILE:HD11	1:A:153:THR:HB	2.01	0.42
1:A:106:GLY:HA2	1:A:118:MET:O	2.20	0.42
2:B:233:PHE:HE2	2:B:256:GLY:HA3	1.83	0.42
2:B:398:LEU:HD11	2:B:412:LEU:HB2	2.01	0.42
1:A:94:PRO:HD2	1:A:129:VAL:HG12	2.02	0.41
2:B:68:THR:HG22	2:B:86:VAL:HG22	2.02	0.41
2:B:221:VAL:HG21	2:B:267:LEU:CD1	2.51	0.41
1:A:133:LYS:O	1:A:139:VAL:HA	2.21	0.41
2:B:49:LEU:H	2:B:178:GLN:NE2	2.19	0.41
2:B:306:LEU:HB2	2:B:383:LEU:HD21	2.03	0.41
2:B:399:GLU:HA	2:B:408:SER:O	2.22	0.41
2:B:41:LEU:HD11	2:B:43:ASP:HB3	2.01	0.40

There are no symmetry-related clashes.

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	100/139 (72%)	97 (97%)	3 (3%)	0	100	100
2	B	423/451 (94%)	398 (94%)	21 (5%)	4 (1%)	14	43
All	All	523/590 (89%)	495 (95%)	24 (5%)	4 (1%)	16	45

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	197	LEU
2	B	242	VAL
2	B	208	ALA
2	B	337	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	98/129 (76%)	92 (94%)	6 (6%)	17	45
2	B	375/394 (95%)	341 (91%)	34 (9%)	9	32
All	All	473/523 (90%)	433 (92%)	40 (8%)	10	34

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	LEU
1	A	59	ILE
1	A	68	VAL
1	A	70	MET
1	A	147	LEU
1	A	150	VAL
2	B	37	LEU
2	B	38	SER
2	B	39	CYS
2	B	41	LEU
2	B	42	TRP
2	B	43	ASP
2	B	46	ILE
2	B	65	HIS
2	B	70	LEU
2	B	71	VAL
2	B	72	LEU
2	B	81	ASP
2	B	83	CYS
2	B	84	LEU
2	B	120	SER
2	B	137	CYS
2	B	144	VAL
2	B	193	ASP
2	B	199	VAL
2	B	258	GLN
2	B	273	ILE

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Mol	Chain	Res	Type
2	B	287	ILE
2	B	310	THR
2	B	348	GLU
2	B	354	LYS
2	B	360	LEU
2	B	361	LEU
2	B	379	LEU
2	B	395	VAL
2	B	399	GLU
2	B	401	ARG
2	B	412	LEU
2	B	427	ARG
2	B	439	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	ASN
1	A	73	ASN
1	A	91	ASN
1	A	104	ASN
1	A	125	GLN
1	A	134	HIS
1	A	160	HIS
2	B	143	GLN
2	B	150	GLN
2	B	153	GLN
2	B	189	GLN
2	B	339	GLN
2	B	404	GLN
2	B	406	GLN
2	B	446	GLN

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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	104/139 (74%)	-0.01	1 (0%)	79 63	94, 144, 205, 215	0
2	B	427/451 (94%)	0.01	8 (1%)	66 47	83, 164, 236, 259	0
All	All	531/590 (90%)	0.01	9 (1%)	69 50	83, 157, 234, 259	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	311	LEU	4.0
2	B	444	CYS	3.5
2	B	21	LEU	2.6
2	B	119	ALA	2.5
2	B	441	SER	2.5
2	B	443	GLN	2.4
2	B	262	LEU	2.4
1	A	68	VAL	2.2
2	B	425	SER	2.2

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.