



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:03 AM UTC

PDB ID : 6E9Q / pdb_00006e9q
Title : The crystal structure of bovine ultralong antibody BOV-6
Authors : Dong, J.; Crowe, J.E.
Deposited on : 2018-08-01
Resolution : 3.40 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:

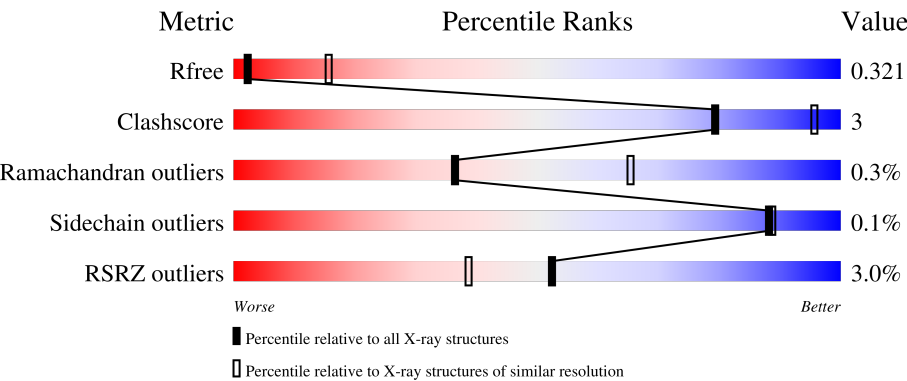
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)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

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 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	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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	263	3% 83% 5% 13%
1	C	263	2% 85% 8% 7%
1	E	263	5% 81% • 15%
1	G	263	5% 77% 6% 17%
2	B	216	94% 5%

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Mol	Chain	Length	Quality of chain
2	D	216	% 88% 11%
2	F	216	% 94% 5%
2	H	216	4% 76% 21%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11743 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 Bovine ultralong antibody BOV-6 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	229	Total	C	N	O	S	0	0	0
			1537	957	272	298	10			
1	C	245	Total	C	N	O	S	0	0	0
			1689	1058	290	328	13			
1	E	224	Total	C	N	O	S	0	0	0
			1552	969	263	310	10			
1	G	219	Total	C	N	O	S	0	0	0
			1406	875	251	271	9			

- Molecule 2 is a protein called Bovine ultralong antibody BOV-6 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	215	Total	C	N	O	S	0	0	0
			1511	931	257	318	5			
2	D	214	Total	C	N	O	S	0	0	0
			1494	918	251	320	5			
2	F	215	Total	C	N	O	S	0	0	0
			1541	947	262	327	5			
2	H	170	Total	C	N	O	S	0	0	0
			1013	610	181	218	4			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1	GLU	GLN	conflict	UNP Q3T101
B	5	ASN	THR	conflict	UNP Q3T101
B	82	ALA	PRO	conflict	UNP Q3T101
D	1	GLU	GLN	conflict	UNP Q3T101
D	5	ASN	THR	conflict	UNP Q3T101
D	82	ALA	PRO	conflict	UNP Q3T101
F	1	GLU	GLN	conflict	UNP Q3T101
F	5	ASN	THR	conflict	UNP Q3T101

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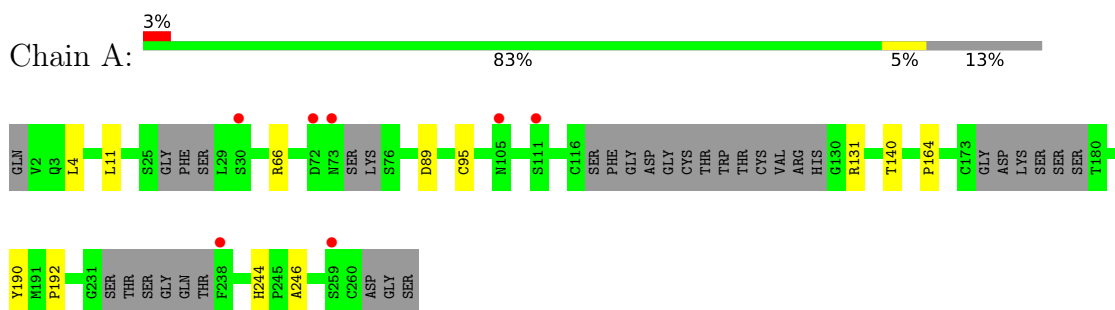
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Chain	Residue	Modelled	Actual	Comment	Reference
F	82	ALA	PRO	conflict	UNP Q3T101
H	1	GLU	GLN	conflict	UNP Q3T101
H	5	ASN	THR	conflict	UNP Q3T101
H	82	ALA	PRO	conflict	UNP Q3T101

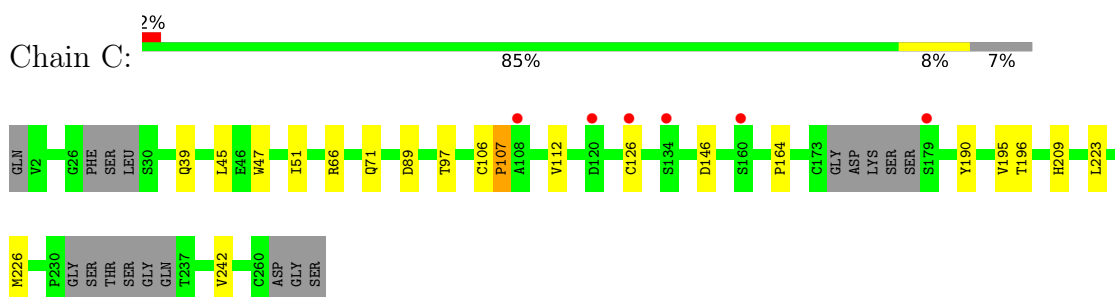
3 Residue-property plots [i](#)

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.

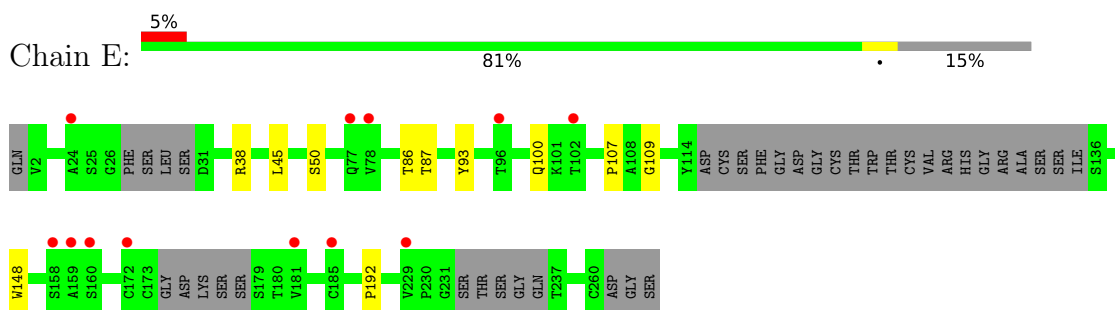
- Molecule 1: Bovine ultralong antibody BOV-6 heavy chain



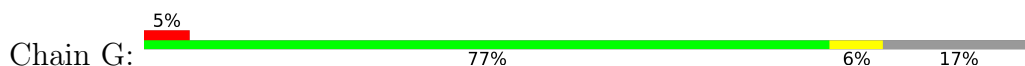
- Molecule 1: Bovine ultralong antibody BOV-6 heavy chain

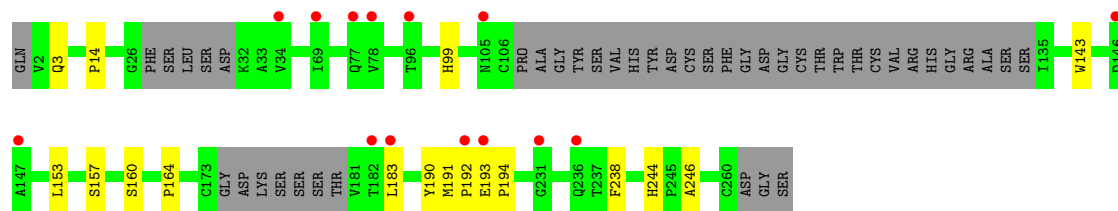


- Molecule 1: Bovine ultralong antibody BOV-6 heavy chain



- Molecule 1: Bovine ultralong antibody BOV-6 heavy chain





- Molecule 2: Bovine ultralong antibody BOV-6 light chain

Chain B: 94% 5%



- Molecule 2: Bovine ultralong antibody BOV-6 light chain

Chain D: 88% 11%



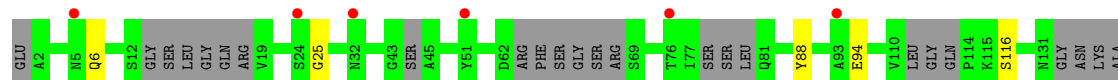
- Molecule 2: Bovine ultralong antibody BOV-6 light chain

Chain F: 94% 5%



- Molecule 2: Bovine ultralong antibody BOV-6 light chain

Chain H: 76% 4% 21%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	64.04Å 72.69Å 452.86Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	23.69 – 3.40 23.69 – 3.40	Depositor EDS
% Data completeness (in resolution range)	100.0 (23.69-3.40) 99.6 (23.69-3.40)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.16 (at 3.40Å)	Xtriage
Refinement program	PHENIX (1.14_3211: ???)	Depositor
R, R_{free}	0.254 , 0.311 0.264 , 0.321	Depositor DCC
R_{free} test set	1464 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	90.7	Xtriage
Anisotropy	0.405	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 82.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	11743	wwPDB-VP
Average B, all atoms (Å ²)	89.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.93% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.08	0/1568	0.27	0/2148
1	C	0.10	0/1728	0.30	0/2366
1	E	0.09	0/1583	0.28	0/2169
1	G	0.11	0/1437	0.36	0/1977
2	B	0.08	0/1541	0.26	0/2108
2	D	0.09	0/1524	0.26	0/2085
2	F	0.08	0/1571	0.26	0/2144
2	H	0.07	0/1022	0.22	0/1398
All	All	0.09	0/11974	0.28	0/16395

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	G	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	G	191	MET	Peptide

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	1537	0	1277	8	0
1	C	1689	0	1442	15	0
1	E	1552	0	1351	6	0
1	G	1406	0	1037	9	0
2	B	1511	0	1360	8	0
2	D	1494	0	1327	16	0
2	F	1541	0	1431	7	0
2	H	1013	0	667	4	0
All	All	11743	0	9892	59	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 3.

All (59) 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:C:107:PRO:HB2	1:G:3:GLN:HE22	1.43	0.83
1:C:209:HIS:HD2	2:D:172:GLN:HB3	1.59	0.67
1:C:39:GLN:HB2	1:C:45:LEU:HD23	1.76	0.67
1:A:11:LEU:HD13	1:A:192:PRO:HG3	1.83	0.60
1:A:140:THR:HA	2:B:32:ASN:HD21	1.67	0.58
1:C:164:PRO:HB3	1:C:190:TYR:HB3	1.85	0.58
1:E:50:SER:OG	1:E:100:GLN:OE1	2.15	0.58
1:C:47:TRP:NE1	2:D:98:SER:O	2.39	0.54
1:A:164:PRO:HB3	1:A:190:TYR:HB3	1.90	0.53
1:C:226:MET:HE1	2:D:142:SER:HB3	1.90	0.53
1:G:164:PRO:HB3	1:G:190:TYR:HB3	1.90	0.53
1:C:209:HIS:CD2	2:D:172:GLN:HB3	2.42	0.53
1:A:4:LEU:HD21	1:A:95:CYS:HB3	1.92	0.52
2:F:27:SER:HA	2:F:31:GLY:HA3	1.94	0.50
1:G:99:HIS:O	1:G:143:TRP:HA	2.11	0.50
1:G:183:LEU:HD12	1:G:238:PHE:HD2	1.77	0.49
1:E:45:LEU:H	1:E:45:LEU:HD23	1.78	0.49
1:C:47:TRP:CD1	2:D:100:ALA:H	2.31	0.48
2:D:6:GLN:NE2	2:D:88:TYR:O	2.39	0.48
1:G:244:HIS:CE1	1:G:246:ALA:HB3	2.48	0.48
1:C:107:PRO:HB2	1:G:3:GLN:NE2	2.19	0.48
1:G:14:PRO:HD3	1:G:157:SER:O	2.14	0.48
1:C:195:VAL:HG12	1:C:223:LEU:HD21	1.95	0.48
1:C:66:ARG:NH1	1:C:89:ASP:OD2	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:97:THR:O	1:C:146:ASP:N	2.47	0.47
2:B:22:THR:HG22	2:B:74:THR:HG22	1.96	0.46
2:F:53:ASP:OD2	2:F:68:ARG:NE	2.34	0.46
1:G:193:GLU:N	1:G:194:PRO:CD	2.79	0.46
1:E:107:PRO:O	1:E:109:GLY:N	2.43	0.46
1:A:244:HIS:CE1	1:A:246:ALA:HB3	2.51	0.45
2:D:123:PHE:HE2	2:D:140:LEU:HD12	1.80	0.45
1:G:153:LEU:HD21	1:G:193:GLU:OE1	2.16	0.45
1:E:38:ARG:NH1	1:E:93:TYR:OH	2.49	0.45
1:A:131:ARG:HE	1:A:131:ARG:HB2	1.63	0.44
2:B:129:GLU:OE2	2:B:136:THR:OG1	2.31	0.44
2:B:137:LEU:HB2	2:B:183:LEU:HB3	1.99	0.44
2:D:118:PRO:HD3	2:D:202:HIS:CD2	2.53	0.44
1:A:140:THR:HA	2:B:32:ASN:ND2	2.31	0.44
2:H:6:GLN:NE2	2:H:88:TYR:O	2.47	0.44
2:D:32:ASN:OD1	2:D:33:GLY:N	2.51	0.44
2:D:118:PRO:HD3	2:D:202:HIS:HD2	1.83	0.43
2:H:116:SER:N	2:H:145:TYR:O	2.49	0.43
2:B:28:SER:N	2:D:28:SER:OG	2.51	0.43
2:B:95:ASP:OD2	2:D:27:SER:OG	2.36	0.43
2:F:2:ALA:N	2:H:25:GLY:O	2.52	0.43
2:D:68:ARG:HA	2:D:73:ALA:HA	2.01	0.42
1:A:66:ARG:NH1	1:A:89:ASP:OD2	2.51	0.42
2:F:71:ASN:ND2	2:H:94:GLU:OE2	2.52	0.42
1:C:196:THR:O	1:C:242:VAL:HA	2.19	0.42
2:D:7:PRO:O	2:D:106:THR:OG1	2.35	0.42
2:D:94:GLU:C	2:D:96:SER:H	2.27	0.42
1:E:148:TRP:CD2	2:F:46:PRO:HB2	2.54	0.41
2:F:16:GLY:HA2	2:F:79:SER:HA	2.02	0.41
1:C:112:VAL:HA	1:C:126:CYS:SG	2.60	0.41
2:F:174:ASN:C	2:F:176:LYS:H	2.29	0.41
2:B:63:ARG:HB3	2:B:78:SER:O	2.21	0.41
1:C:51:ILE:HD13	1:C:71:GLN:HB2	2.02	0.40
2:D:152:VAL:HG23	2:D:199:GLU:HB3	2.03	0.40
1:E:86:THR:OG1	1:E:87:THR:N	2.55	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	217/263 (82%)	201 (93%)	16 (7%)	0	100	100
1	C	237/263 (90%)	219 (92%)	17 (7%)	1 (0%)	30	59
1	E	214/263 (81%)	202 (94%)	11 (5%)	1 (0%)	24	54
1	G	211/263 (80%)	197 (93%)	12 (6%)	2 (1%)	14	42
2	B	213/216 (99%)	201 (94%)	12 (6%)	0	100	100
2	D	212/216 (98%)	204 (96%)	7 (3%)	1 (0%)	24	54
2	F	213/216 (99%)	197 (92%)	16 (8%)	0	100	100
2	H	148/216 (68%)	137 (93%)	11 (7%)	0	100	100
All	All	1665/1916 (87%)	1558 (94%)	102 (6%)	5 (0%)	36	65

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	192	PRO
1	G	192	PRO
1	C	107	PRO
1	G	160	SER
2	D	53	ASP

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	130/224 (58%)	130 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	154/224 (69%)	153 (99%)	1 (1%)	78	80
1	E	147/224 (66%)	147 (100%)	0	100	100
1	G	94/224 (42%)	94 (100%)	0	100	100
2	B	156/184 (85%)	156 (100%)	0	100	100
2	D	154/184 (84%)	154 (100%)	0	100	100
2	F	169/184 (92%)	169 (100%)	0	100	100
2	H	63/184 (34%)	63 (100%)	0	100	100
All	All	1067/1632 (65%)	1066 (100%)	1 (0%)	88	89

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	106	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	5	ASN
2	B	39	GLN
1	C	60	ASN
1	C	99	HIS
2	D	99	ASN
2	F	71	ASN
1	G	3	GLN
1	G	60	ASN
1	G	100	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 ⓘ

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	229/263 (87%)	0.35	7 (3%) 51 38	55, 85, 141, 195	0
1	C	245/263 (93%)	0.23	6 (2%) 59 45	52, 73, 149, 203	0
1	E	224/263 (85%)	0.34	12 (5%) 31 24	67, 95, 148, 168	0
1	G	219/263 (83%)	0.47	14 (6%) 25 20	63, 105, 163, 271	0
2	B	215/216 (99%)	-0.01	0 100 100	49, 64, 99, 148	0
2	D	214/216 (99%)	0.11	2 (0%) 81 68	50, 72, 107, 134	0
2	F	215/216 (99%)	0.25	2 (0%) 81 68	60, 84, 119, 204	0
2	H	170/216 (78%)	0.70	9 (5%) 32 24	66, 117, 173, 198	0
All	All	1731/1916 (90%)	0.29	52 (3%) 52 39	49, 85, 148, 271	0

All (52) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	2	ALA	6.8
1	C	126	CYS	3.4
2	H	93	ALA	3.3
1	A	105	ASN	3.2
1	C	120	ASP	3.0
2	F	189	ASP	2.9
1	G	231	GLY	2.9
1	A	30	SER	2.9
1	E	96	THR	2.9
1	C	108	ALA	2.8
1	C	134	SER	2.8
1	G	192	PRO	2.7
2	H	197	SER	2.7
1	E	172	CYS	2.6
2	H	24	SER	2.6
1	E	159	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
1	G	193	GLU	2.5
1	E	181	VAL	2.5
1	G	34	VAL	2.5
1	E	78	VAL	2.4
1	A	73	ASN	2.4
1	G	77	GLN	2.4
2	D	66	GLY	2.4
2	H	51	TYR	2.3
1	E	229	VAL	2.3
1	G	183	LEU	2.3
1	A	238	PHE	2.3
1	G	78	VAL	2.3
2	H	32	ASN	2.3
1	A	111	SER	2.2
2	H	5	ASN	2.2
2	H	198	CYS	2.2
1	G	96	THR	2.2
1	A	259	SER	2.2
1	E	102	THR	2.2
1	G	182	THR	2.2
2	H	76	THR	2.2
2	D	177	TYR	2.2
1	C	179	SER	2.2
1	E	158	SER	2.2
1	E	77	GLN	2.1
1	G	236	GLN	2.1
1	G	147	ALA	2.1
1	G	146	ASP	2.1
1	G	69	ILE	2.1
2	H	177	TYR	2.1
1	E	160	SER	2.1
1	A	72	ASP	2.1
1	C	160	SER	2.0
1	G	105	ASN	2.0
1	E	24	ALA	2.0
1	E	185	CYS	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 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.