



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 12:33 PM UTC

PDB ID : 3AXL / pdb_00003axl
Title : Murine Valpha 10 Vbeta 8.1 T-cell receptor
Authors : Patel, O.; Rossjohn, J.
Deposited on : 2011-04-11
Resolution : 2.90 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

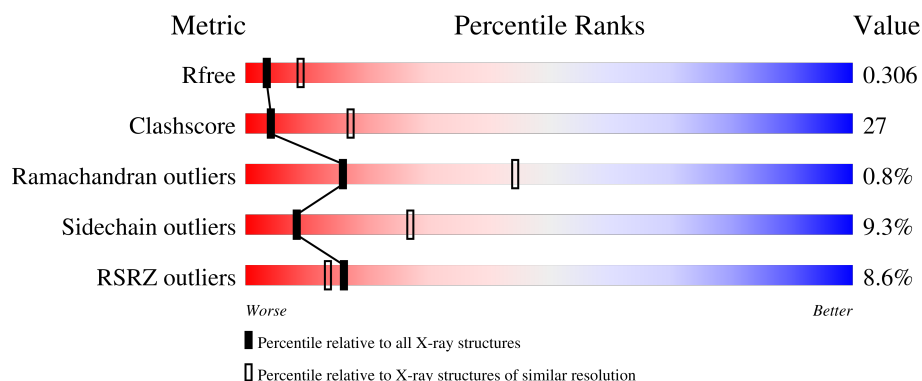
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	198	
1	G	198	
2	B	238	
2	H	238	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6728 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 Valpha 10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	183	Total	C	N	O	S	0	0	0
			1393	866	232	287	8			
1	G	190	Total	C	N	O	S	0	0	0
			1437	890	240	299	8			

- Molecule 2 is a protein called Vbeta 8.1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	238	Total	C	N	O	S	0	0	0
			1887	1193	330	358	6			
2	H	238	Total	C	N	O	S	0	0	0
			1883	1189	330	358	6			

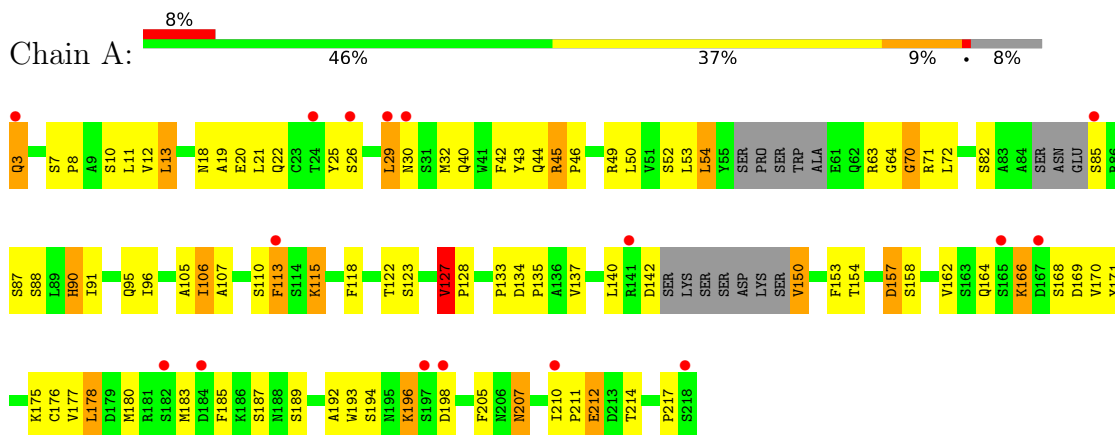
- Molecule 3 is water.

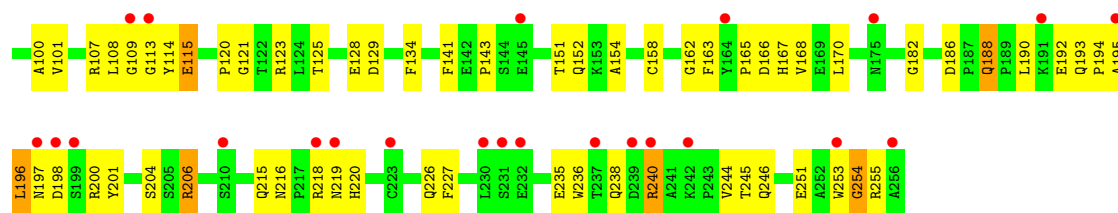
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	28	Total	O	0	0
			28	28		
3	B	43	Total	O	0	0
			43	43		
3	G	27	Total	O	0	0
			27	27		
3	H	30	Total	O	0	0
			30	30		

3 Residue-property plots [i](#)

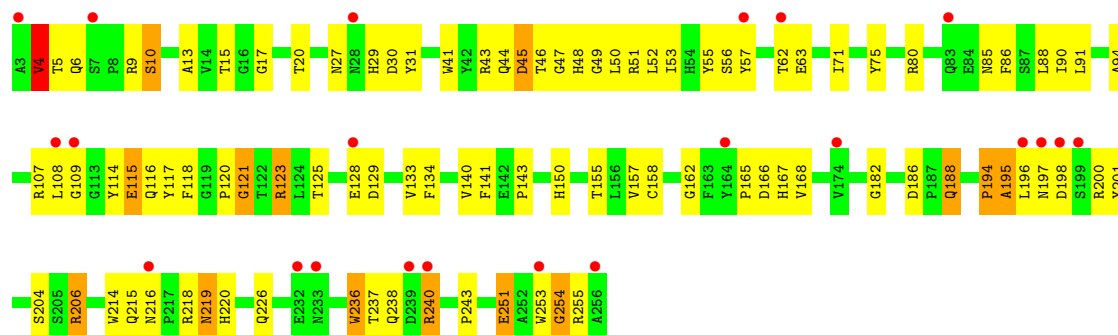
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Valpha 10





● Molecule 2: Vbeta 8.1



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	148.84Å 65.80Å 97.38Å 90.00° 97.14° 90.00°	Depositor
Resolution (Å)	42.95 – 2.90 42.95 – 2.90	Depositor EDS
% Data completeness (in resolution range)	96.4 (42.95-2.90) 99.6 (42.95-2.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.64 (at 2.77Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, R_{free}	0.254 , 0.327 (Not available) , 0.306	Depositor DCC
R_{free} test set	1185 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	39.7	Xtriage
Anisotropy	0.464	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 62.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	6728	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.50% 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	0.64	0/1417	1.01	6/1921 (0.3%)
1	G	0.68	0/1463	1.01	6/1986 (0.3%)
2	B	0.64	0/1940	1.08	12/2644 (0.5%)
2	H	0.75	0/1936	1.36	25/2638 (0.9%)
All	All	0.68	0/6756	1.14	49/9189 (0.5%)

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
2	B	0	1

There are no bond length outliers.

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	195	ALA	N-CA-C	8.52	123.73	113.16
2	H	240	ARG	N-CA-C	7.61	119.71	108.14
2	B	4	VAL	N-CA-C	-6.92	98.15	108.46
2	H	4	VAL	N-CA-C	-6.90	98.18	108.12
2	H	186	ASP	CA-C-N	6.79	126.59	119.05
2	H	186	ASP	C-N-CA	6.79	126.59	119.05
2	H	157	VAL	N-CA-C	6.68	117.52	108.17
2	H	155	THR	N-CA-C	6.42	118.99	108.52
2	B	186	ASP	CA-C-N	6.03	125.75	119.05
2	B	186	ASP	C-N-CA	6.03	125.75	119.05
1	G	127	VAL	CA-C-N	5.99	126.00	119.89
1	G	127	VAL	C-N-CA	5.99	126.00	119.89
2	H	194	PRO	CB-CA-C	-5.78	103.65	111.85
2	H	115	GLU	N-CA-C	5.76	118.68	110.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	240	ARG	N-CA-C	5.75	115.98	107.88
1	A	45	ARG	CA-C-N	-5.59	113.79	119.83
1	A	45	ARG	C-N-CA	-5.59	113.79	119.83
2	B	188	GLN	CA-C-N	5.59	125.54	119.78
2	B	188	GLN	C-N-CA	5.59	125.54	119.78
1	A	127	VAL	CA-C-N	5.58	125.59	119.89
1	A	127	VAL	C-N-CA	5.58	125.59	119.89
2	H	236	TRP	N-CA-C	5.50	117.38	108.26
2	H	133	VAL	N-CA-C	5.49	116.09	108.84
1	A	70	GLY	N-CA-C	5.48	121.28	114.37
1	A	162	VAL	N-CA-C	5.48	115.84	108.17
2	H	123	ARG	CA-CB-CG	5.46	125.02	114.10
2	H	150	HIS	N-CA-C	5.46	117.31	111.36
1	G	45	ARG	CA-C-N	-5.45	113.94	119.83
1	G	45	ARG	C-N-CA	-5.45	113.94	119.83
1	G	57	PRO	N-CA-C	-5.44	103.61	111.22
2	B	85	ASN	N-CA-C	5.43	118.25	109.40
2	H	196	LEU	N-CA-C	5.42	118.52	109.96
2	H	49	GLY	N-CA-C	-5.35	100.51	113.18
2	H	107	ARG	N-CA-CB	-5.32	102.09	111.39
2	B	8	PRO	N-CA-C	-5.30	101.55	112.47
2	B	7	SER	CA-C-N	5.29	126.45	119.84
2	B	7	SER	C-N-CA	5.29	126.45	119.84
2	H	188	GLN	N-CA-C	5.23	116.70	108.23
2	H	5	THR	N-CA-C	5.19	118.18	110.14
2	H	219	ASN	N-CA-C	5.16	118.03	109.46
2	H	237	THR	CB-CA-C	-5.14	100.92	109.55
2	H	140	VAL	N-CA-C	5.13	115.97	108.53
2	B	48	HIS	N-CA-C	-5.09	103.88	110.19
2	H	117	TYR	N-CA-C	5.07	117.75	109.59
2	H	116	GLN	N-CA-C	5.06	117.87	109.46
1	G	70	GLY	N-CA-C	5.05	118.99	113.58
2	B	115	GLU	N-CA-CB	-5.04	102.45	110.06
2	H	251	GLU	N-CA-C	5.01	116.81	109.24
2	H	121	GLY	N-CA-C	5.01	118.89	112.18

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	7	SER	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	1393	0	1312	88	0
1	G	1437	0	1348	94	0
2	B	1887	0	1788	119	0
2	H	1883	0	1778	91	0
3	A	28	0	0	2	0
3	B	43	0	0	18	0
3	G	27	0	0	6	0
3	H	30	0	0	5	0
All	All	6728	0	6226	352	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 27.

All (352) 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:H:17:GLY:HA3	3:H:275:HOH:O	1.69	0.93
2:B:58:VAL:HA	3:B:271:HOH:O	1.72	0.88
2:B:216:ASN:HB3	2:B:219:ASN:HD22	1.35	0.88
2:B:216:ASN:HB3	2:B:219:ASN:ND2	1.89	0.86
2:H:6:GLN:HG2	2:H:120:PRO:HD2	1.57	0.84
2:H:48:HIS:HD2	2:H:51:ARG:CZ	1.90	0.84
2:B:6:GLN:HG2	2:B:120:PRO:HD2	1.61	0.83
2:H:216:ASN:HB3	2:H:219:ASN:ND2	1.95	0.81
1:A:153:PHE:HB2	1:A:205:PHE:CZ	2.14	0.81
1:A:166:LYS:O	1:G:166:LYS:O	1.99	0.81
2:B:162:GLY:O	2:B:200:ARG:HD3	1.81	0.80
2:H:43:ARG:HG2	2:H:53:ILE:HD11	1.65	0.78
2:B:236:TRP:CH2	2:B:238:GLN:HG3	2.17	0.78
1:G:153:PHE:HB2	1:G:205:PHE:CZ	2.18	0.78
2:H:236:TRP:CH2	2:H:238:GLN:HG3	2.19	0.78
2:B:194:PRO:HG2	1:G:166:LYS:HG3	1.64	0.77
2:B:43:ARG:HG2	2:B:53:ILE:HD11	1.66	0.77
2:H:216:ASN:HB3	2:H:219:ASN:HD22	1.49	0.77
2:H:13:ALA:HA	3:H:270:HOH:O	1.84	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:PHE:HE2	1:A:187:SER:HB2	1.48	0.76
1:G:185:PHE:HE2	1:G:187:SER:HB2	1.51	0.75
2:B:6:GLN:HB2	3:B:279:HOH:O	1.86	0.74
2:B:162:GLY:C	2:B:200:ARG:HD3	2.13	0.74
1:G:84(A):SER:C	1:G:86:ARG:NH2	2.45	0.73
2:B:31:TYR:CE1	2:B:57:TYR:CE1	2.77	0.72
2:H:167:HIS:HE1	3:H:69:HOH:O	1.71	0.72
2:H:134:PHE:HZ	2:H:240:ARG:HH22	1.34	0.72
2:H:31:TYR:CE1	2:H:57:TYR:CE1	2.78	0.71
2:B:123:ARG:NH1	3:B:289:HOH:O	2.23	0.71
2:H:162:GLY:O	2:H:200:ARG:HD3	1.90	0.71
2:B:134:PHE:HZ	2:B:240:ARG:HH22	1.35	0.70
2:H:162:GLY:C	2:H:200:ARG:HD3	2.16	0.70
1:A:185:PHE:CE2	1:A:187:SER:HB2	2.28	0.68
1:A:110:SER:HB2	2:B:109:GLY:O	1.93	0.68
1:G:185:PHE:CE2	1:G:187:SER:HB2	2.28	0.68
1:A:113:PHE:HB2	2:B:31:TYR:OH	1.95	0.67
2:B:196:LEU:HD13	2:B:196:LEU:H	1.57	0.67
1:G:169:ASP:OD2	1:G:196:LYS:HD3	1.95	0.67
1:G:128:PRO:HG3	1:G:177:VAL:HG21	1.77	0.66
1:A:169:ASP:O	1:A:194:SER:HB2	1.95	0.66
2:B:193:GLN:HB3	2:B:196:LEU:HD11	1.77	0.66
2:B:196:LEU:HD13	2:B:196:LEU:N	2.11	0.66
2:H:15:THR:HA	2:H:94:ALA:O	1.96	0.66
2:H:134:PHE:CZ	2:H:240:ARG:NH2	2.60	0.66
1:G:110:SER:HB3	3:G:239:HOH:O	1.95	0.66
1:G:55:TYR:HE2	1:G:57:PRO:HG3	1.61	0.66
2:H:215:GLN:HA	2:H:255:ARG:O	1.95	0.66
2:B:31:TYR:HE1	2:B:57:TYR:CE1	2.15	0.65
1:G:55:TYR:CE2	1:G:57:PRO:HG3	2.31	0.65
1:A:29:LEU:HB3	1:A:106:ILE:HD13	1.78	0.65
2:B:220:HIS:HE1	2:B:251:GLU:HB2	1.62	0.65
1:A:44:GLN:HB2	1:A:50:LEU:HD23	1.78	0.65
1:A:20:GLU:OE2	1:A:90:HIS:CG	2.50	0.64
1:A:153:PHE:HB2	1:A:205:PHE:CE2	2.32	0.64
1:A:169:ASP:OD2	1:A:196:LYS:HD3	1.98	0.64
2:B:193:GLN:CB	2:B:196:LEU:HD11	2.26	0.64
1:A:18:ASN:HB3	3:A:73:HOH:O	1.98	0.64
2:B:196:LEU:HD23	2:B:197:ASN:O	1.97	0.64
2:B:48:HIS:HD2	2:B:51:ARG:CZ	2.11	0.64
2:B:215:GLN:HA	2:B:255:ARG:O	1.98	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:PHE:HB2	2:B:31:TYR:CZ	2.33	0.64
2:H:31:TYR:HE1	2:H:57:TYR:CE1	2.16	0.64
1:A:25:TYR:CD1	1:A:26:SER:N	2.66	0.63
1:A:40:GLN:NE2	2:B:115:GLU:HG3	2.13	0.63
1:G:169:ASP:O	1:G:194:SER:HB2	1.97	0.63
1:G:44:GLN:HB2	1:G:50:LEU:HD23	1.80	0.63
1:G:110:SER:HB2	2:H:109:GLY:O	1.98	0.63
1:G:140:LEU:HB3	2:H:141:PHE:HB3	1.81	0.63
1:A:32:MET:HE1	1:A:87:SER:OG	1.99	0.62
1:A:207:ASN:ND2	2:H:195:ALA:HB2	2.14	0.62
2:H:91:LEU:N	2:H:91:LEU:HD12	2.14	0.62
2:B:59:ALA:N	3:B:271:HOH:O	2.31	0.62
1:A:166:LYS:HG3	2:H:194:PRO:HG2	1.81	0.62
2:H:48:HIS:CD2	2:H:51:ARG:CZ	2.79	0.62
2:H:10:SER:HB2	2:H:167:HIS:HD2	1.65	0.61
2:B:158:CYS:O	2:B:204:SER:HA	2.00	0.61
1:G:19:ALA:HB3	1:G:91:ILE:HB	1.83	0.61
1:A:212:GLU:HG2	3:A:233:HOH:O	2.01	0.61
2:B:15:THR:HA	2:B:94:ALA:O	1.99	0.61
1:G:32:MET:HE1	1:G:87:SER:OG	2.00	0.61
1:A:135:PRO:HB2	1:A:214:THR:HA	1.82	0.60
1:G:170:VAL:HA	1:G:194:SER:HB3	1.83	0.60
1:G:128:PRO:HG3	1:G:177:VAL:CG2	2.31	0.60
1:A:113:PHE:H	1:A:113:PHE:HD2	1.48	0.60
1:G:63:ARG:C	1:G:70:GLY:H	2.10	0.60
2:H:134:PHE:CE1	2:H:240:ARG:NH2	2.70	0.60
2:B:100:ALA:C	3:B:289:HOH:O	2.43	0.60
2:H:31:TYR:HA	2:H:56:SER:O	2.02	0.60
1:G:25:TYR:CD1	1:G:26:SER:N	2.70	0.60
1:G:113:PHE:H	1:G:113:PHE:HD2	1.50	0.59
1:G:135:PRO:HB2	1:G:214:THR:HA	1.85	0.59
2:B:236:TRP:N	3:B:273:HOH:O	2.36	0.59
1:A:3:GLN:HB3	1:A:25:TYR:HE1	1.66	0.59
1:A:110:SER:O	2:B:109:GLY:C	2.46	0.58
1:G:29:LEU:HB3	1:G:106:ILE:HD13	1.83	0.58
1:A:170:VAL:HA	1:A:194:SER:HB3	1.85	0.58
1:G:178:LEU:HD23	3:G:240:HOH:O	2.03	0.58
2:B:91:LEU:HD12	2:B:91:LEU:N	2.18	0.58
1:G:40:GLN:NE2	2:H:115:GLU:HG3	2.18	0.58
1:G:153:PHE:HB2	1:G:205:PHE:CE2	2.39	0.58
1:A:63:ARG:C	1:A:70:GLY:H	2.12	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:220:HIS:HE1	2:H:251:GLU:HB2	1.67	0.58
2:B:128:GLU:OE1	2:B:128:GLU:HA	2.03	0.57
2:B:236:TRP:CZ2	2:B:238:GLN:HB2	2.39	0.57
2:B:134:PHE:CZ	2:B:240:ARG:NH2	2.64	0.57
2:B:216:ASN:HB2	3:B:112:HOH:O	2.04	0.57
2:B:48:HIS:CD2	2:B:51:ARG:CZ	2.88	0.57
1:A:19:ALA:HB3	1:A:91:ILE:HB	1.87	0.56
2:B:10:SER:HB2	2:B:167:HIS:HD2	1.69	0.56
2:B:31:TYR:CE1	2:B:57:TYR:CD1	2.93	0.56
1:G:20:GLU:OE2	1:G:90:HIS:CG	2.59	0.56
1:A:175:LYS:HA	1:A:189:SER:O	2.05	0.56
2:B:134:PHE:CE1	2:B:240:ARG:NH2	2.73	0.56
2:H:158:CYS:O	2:H:204:SER:HA	2.05	0.56
1:G:178:LEU:HD13	2:H:182:GLY:HA2	1.86	0.55
2:H:31:TYR:CE1	2:H:57:TYR:CD1	2.93	0.55
2:B:108:LEU:HB3	2:B:114:TYR:HB2	1.89	0.55
2:B:143:PRO:HB3	2:B:154:ALA:HB1	1.88	0.55
2:B:27:ASN:HB3	2:B:29:HIS:CD2	2.42	0.55
1:G:22:GLN:OE1	1:G:88:SER:CB	2.53	0.55
1:G:53:LEU:O	1:G:54:LEU:CD1	2.54	0.55
2:H:30:ASP:OD2	2:H:31:TYR:CE2	2.60	0.55
1:A:178:LEU:HD12	1:A:178:LEU:C	2.32	0.55
1:G:84(A):SER:C	1:G:86:ARG:HH22	2.15	0.54
1:G:3:GLN:HB3	1:G:25:TYR:HE1	1.73	0.54
1:A:128:PRO:HG3	1:A:177:VAL:HG21	1.90	0.54
1:G:127:VAL:CG1	1:G:158:SER:HB3	2.39	0.53
2:B:41:TRP:HH2	3:B:279:HOH:O	1.91	0.53
2:B:216:ASN:CB	2:B:219:ASN:ND2	2.66	0.53
1:G:175:LYS:HA	1:G:189:SER:O	2.08	0.53
2:H:243:PRO:HA	3:H:282:HOH:O	2.06	0.53
2:B:129:ASP:C	2:B:129:ASP:OD1	2.52	0.53
1:G:113:PHE:HB2	2:H:31:TYR:OH	2.09	0.53
2:H:55:TYR:CE1	2:H:63:GLU:HG3	2.43	0.53
2:B:30:ASP:OD2	2:B:31:TYR:CE2	2.61	0.53
1:A:210:ILE:O	1:A:211:PRO:C	2.51	0.53
1:G:49:ARG:NH1	2:H:118:PHE:O	2.42	0.53
2:B:48:HIS:HD2	2:B:51:ARG:NE	2.07	0.52
2:B:88:LEU:C	2:B:90:ILE:HD13	2.34	0.52
2:H:13:ALA:CA	3:H:270:HOH:O	2.51	0.52
1:A:110:SER:HB2	2:B:113:GLY:CA	2.40	0.52
1:A:128:PRO:HG3	1:A:177:VAL:CG2	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:55:TYR:CE1	2:B:63:GLU:HG3	2.43	0.52
1:G:53:LEU:O	1:G:54:LEU:HD13	2.09	0.52
2:H:20:THR:HG23	2:H:90:ILE:CD1	2.40	0.52
2:B:254:GLY:O	2:B:255:ARG:HG3	2.10	0.52
1:G:188:ASN:N	3:G:220:HOH:O	2.20	0.52
1:G:210:ILE:O	1:G:211:PRO:C	2.53	0.52
2:B:41:TRP:HD1	2:B:86:PHE:CE2	2.28	0.52
2:B:166:ASP:HB3	2:B:201:TYR:CG	2.45	0.52
1:G:189:SER:OG	2:H:206:ARG:HD2	2.10	0.52
2:B:27:ASN:CB	2:B:29:HIS:CD2	2.94	0.51
2:H:236:TRP:CZ2	2:H:238:GLN:HB2	2.46	0.51
2:H:128:GLU:HA	2:H:128:GLU:OE1	2.10	0.51
2:B:162:GLY:HA2	2:B:200:ARG:CD	2.41	0.51
1:G:40:GLN:HE22	2:H:115:GLU:HG3	1.76	0.51
1:A:166:LYS:CE	1:G:168:SER:HA	2.41	0.51
2:B:27:ASN:HB3	2:B:29:HIS:NE2	2.26	0.51
1:A:22:GLN:OE1	1:A:88:SER:CB	2.60	0.50
1:G:30:ASN:HB3	1:G:107:ALA:O	2.12	0.50
1:G:61:GLU:N	3:G:222:HOH:O	2.43	0.50
1:A:110:SER:HA	2:B:113:GLY:HA2	1.92	0.50
1:G:11:LEU:N	1:G:11:LEU:HD23	2.26	0.50
2:B:20:THR:HG23	2:B:90:ILE:CD1	2.42	0.50
1:G:8:PRO:O	1:G:122:THR:HG23	2.12	0.50
1:G:22:GLN:OE1	1:G:88:SER:HB2	2.12	0.50
2:H:88:LEU:C	2:H:90:ILE:HD13	2.37	0.49
1:A:110:SER:O	2:B:113:GLY:N	2.45	0.49
2:B:31:TYR:HA	2:B:56:SER:O	2.13	0.49
1:A:176:CYS:SG	2:B:206:ARG:NH1	2.86	0.49
1:G:29:LEU:H	1:G:29:LEU:HD13	1.78	0.49
1:G:84(A):SER:C	1:G:86:ARG:HH21	2.18	0.49
1:G:171:TYR:O	1:G:192:ALA:HA	2.13	0.49
2:H:44:GLN:HB2	2:H:50:LEU:HD23	1.93	0.49
1:A:11:LEU:N	1:A:11:LEU:HD23	2.28	0.49
1:A:150:VAL:HG22	1:A:193:TRP:HB3	1.95	0.49
1:G:113:PHE:HB2	2:H:31:TYR:CZ	2.48	0.49
1:A:113:PHE:CB	2:B:31:TYR:OH	2.59	0.49
1:A:115:LYS:HA	2:B:107:ARG:HH12	1.77	0.49
1:A:171:TYR:O	1:A:192:ALA:HA	2.13	0.49
2:B:4:VAL:O	2:B:120:PRO:HD3	2.12	0.48
1:G:44:GLN:HA	3:G:233:HOH:O	2.13	0.48
2:H:141:PHE:N	2:H:141:PHE:CD2	2.81	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:129:ASP:OD1	2:H:129:ASP:C	2.56	0.48
1:A:29:LEU:HB3	1:A:106:ILE:CD1	2.42	0.48
1:G:3:GLN:NE2	1:G:26:SER:HB2	2.28	0.48
1:A:207:ASN:HD22	2:H:195:ALA:HB2	1.77	0.48
2:B:52:LEU:O	2:B:71:ILE:HG12	2.13	0.48
1:A:127:VAL:CG1	1:A:158:SER:HB3	2.42	0.48
2:B:193:GLN:HB2	2:B:196:LEU:HD11	1.95	0.48
1:G:84(B):ASN:N	1:G:86:ARG:HH21	2.12	0.48
1:A:7:SER:HB3	1:A:22:GLN:HB2	1.95	0.48
1:A:13:LEU:HD23	1:G:8:PRO:HB3	1.95	0.48
2:H:41:TRP:CD1	2:H:86:PHE:CD2	3.02	0.48
1:G:178:LEU:C	1:G:178:LEU:HD12	2.38	0.48
2:H:41:TRP:HD1	2:H:86:PHE:CE2	2.32	0.48
1:A:40:GLN:HE22	2:B:115:GLU:HG3	1.77	0.48
2:B:11:LYS:NZ	3:B:288:HOH:O	2.39	0.48
2:H:27:ASN:HB3	2:H:29:HIS:CD2	2.49	0.48
2:H:75:TYR:HB3	2:H:88:LEU:HD11	1.96	0.48
2:H:114:TYR:N	2:H:114:TYR:CD2	2.81	0.48
1:A:29:LEU:HD13	1:A:29:LEU:H	1.79	0.47
1:A:54:LEU:HD12	1:A:54:LEU:HA	1.59	0.47
1:G:84(B):ASN:O	1:G:84(C):GLU:C	2.56	0.47
2:H:45:ASP:OD1	2:H:45:ASP:N	2.47	0.47
2:H:48:HIS:HD2	2:H:51:ARG:NE	2.12	0.47
1:A:46:PRO:HG3	3:B:285:HOH:O	2.15	0.47
1:A:178:LEU:HD13	2:B:182:GLY:HA2	1.97	0.47
1:G:40:GLN:HG3	1:G:54:LEU:O	2.14	0.47
1:A:30:ASN:HB3	1:A:107:ALA:O	2.15	0.47
2:H:30:ASP:OD2	2:H:31:TYR:CD2	2.68	0.47
1:A:3:GLN:NE2	1:A:26:SER:HB2	2.29	0.47
2:B:41:TRP:CD1	2:B:86:PHE:CD2	3.01	0.47
1:G:22:GLN:OE1	1:G:88:SER:HB3	2.13	0.47
1:G:29:LEU:H	1:G:29:LEU:CD1	2.27	0.47
2:H:168:VAL:HA	2:H:226:GLN:O	2.15	0.47
1:A:3:GLN:HB3	1:A:25:TYR:CE1	2.49	0.47
2:H:4:VAL:O	2:H:120:PRO:HD3	2.14	0.47
2:B:43:ARG:NH1	3:B:268:HOH:O	2.47	0.47
1:G:57:PRO:O	1:G:58:SER:O	2.33	0.46
1:A:29:LEU:H	1:A:29:LEU:CD1	2.28	0.46
1:G:3:GLN:HB3	1:G:25:TYR:CE1	2.50	0.46
2:H:125:THR:HG21	2:H:165:PRO:CG	2.45	0.46
1:A:22:GLN:OE1	1:A:88:SER:HB2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:168:SER:HA	1:G:166:LYS:CE	2.46	0.46
1:A:71:ARG:C	1:A:72:LEU:HD12	2.40	0.46
1:G:63:ARG:O	1:G:70:GLY:N	2.48	0.46
1:G:84(B):ASN:N	1:G:86:ARG:NH2	2.63	0.46
2:H:46:THR:HA	2:H:47:GLY:HA2	1.73	0.46
2:H:162:GLY:HA2	2:H:200:ARG:CD	2.46	0.46
1:A:8:PRO:O	1:A:122:THR:HG23	2.15	0.46
1:G:113:PHE:CB	2:H:31:TYR:OH	2.64	0.46
2:B:53:ILE:HD12	2:B:53:ILE:N	2.31	0.46
1:G:29:LEU:HB3	1:G:106:ILE:CD1	2.43	0.46
2:H:108:LEU:HB3	2:H:114:TYR:HB2	1.98	0.46
2:B:101:VAL:HG23	3:B:289:HOH:O	2.15	0.46
2:B:236:TRP:CZ2	2:B:238:GLN:HG3	2.51	0.46
1:G:196:LYS:N	3:G:228:HOH:O	2.48	0.46
1:A:133:PRO:HB3	1:A:157:ASP:HB3	1.98	0.45
1:G:54:LEU:HD12	1:G:54:LEU:HA	1.67	0.45
2:B:28:ASN:ND2	3:B:278:HOH:O	2.49	0.45
1:A:134:ASP:O	1:A:135:PRO:C	2.58	0.45
2:B:235:GLU:HA	3:B:273:HOH:O	2.16	0.45
1:A:63:ARG:O	1:A:70:GLY:N	2.49	0.45
2:H:51:ARG:HB3	2:H:71:ILE:CD1	2.46	0.45
2:H:134:PHE:CE1	2:H:240:ARG:NH1	2.85	0.45
2:H:254:GLY:O	2:H:255:ARG:HG3	2.16	0.45
2:B:195:ALA:HB2	1:G:207:ASN:ND2	2.32	0.45
1:G:178:LEU:CD1	2:H:182:GLY:HA2	2.47	0.45
1:G:127:VAL:HG11	1:G:158:SER:HB3	1.99	0.45
1:A:43:TYR:CD2	1:A:43:TYR:C	2.94	0.45
2:B:226:GLN:NE2	3:B:266:HOH:O	2.46	0.45
1:G:84(C):GLU:O	1:G:86:ARG:N	2.49	0.45
1:A:82:SER:O	1:A:87:SER:HA	2.17	0.45
1:G:176:CYS:SG	2:H:206:ARG:NH1	2.89	0.45
2:H:6:GLN:CD	2:H:121:GLY:H	2.25	0.45
2:B:192:GLU:HG3	2:B:200:ARG:O	2.17	0.45
2:B:216:ASN:ND2	2:B:219:ASN:ND2	2.65	0.45
2:H:218:ARG:O	2:H:253:TRP:HZ3	2.00	0.44
2:B:162:GLY:C	2:B:200:ARG:CD	2.88	0.44
2:H:10:SER:HA	2:H:123:ARG:O	2.17	0.44
2:H:236:TRP:CZ2	2:H:238:GLN:HG3	2.51	0.44
2:B:75:TYR:HB3	2:B:88:LEU:HD11	1.99	0.44
2:B:151:THR:O	2:B:152:GLN:HB2	2.17	0.44
2:B:162:GLY:CA	2:B:200:ARG:HD3	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:216:ASN:O	2:B:254:GLY:HA3	2.16	0.44
1:G:63:ARG:C	1:G:70:GLY:N	2.75	0.44
1:G:113:PHE:HB2	2:H:31:TYR:CE1	2.53	0.44
1:A:8:PRO:HB2	1:G:12:VAL:O	2.16	0.44
2:B:80:ARG:HD2	2:B:85:ASN:O	2.17	0.44
2:B:143:PRO:CB	2:B:154:ALA:HB1	2.48	0.44
2:B:163:PHE:CE2	2:B:201:TYR:HB2	2.51	0.44
1:G:134:ASP:O	1:G:135:PRO:C	2.60	0.44
2:B:134:PHE:CE1	2:B:240:ARG:NH1	2.86	0.44
1:G:169:ASP:OD1	1:G:169:ASP:N	2.49	0.44
1:A:63:ARG:C	1:A:70:GLY:N	2.76	0.43
1:A:185:PHE:C	1:A:185:PHE:CD2	2.96	0.43
2:B:141:PHE:CD2	2:B:141:PHE:N	2.85	0.43
1:A:53:LEU:O	1:A:54:LEU:CD1	2.66	0.43
1:A:140:LEU:HB3	2:B:141:PHE:HB3	2.00	0.43
2:H:10:SER:HB2	2:H:167:HIS:CD2	2.49	0.43
2:B:41:TRP:CD1	2:B:86:PHE:CE2	3.06	0.43
2:B:125:THR:HG21	2:B:165:PRO:CG	2.49	0.43
2:H:166:ASP:HB3	2:H:201:TYR:CG	2.54	0.43
2:H:216:ASN:CB	2:H:219:ASN:ND2	2.76	0.43
1:A:12:VAL:O	1:G:8:PRO:HB2	2.18	0.43
1:A:113:PHE:HB2	2:B:31:TYR:CE1	2.53	0.43
2:B:96:LEU:HD11	2:B:128:GLU:OE2	2.19	0.43
2:H:53:ILE:N	2:H:53:ILE:HD12	2.33	0.43
2:H:166:ASP:HB3	2:H:201:TYR:CD2	2.54	0.43
2:H:219:ASN:O	2:H:253:TRP:HA	2.19	0.43
2:H:52:LEU:O	2:H:71:ILE:HG12	2.19	0.43
2:H:197:ASN:O	2:H:198:ASP:C	2.61	0.43
2:H:27:ASN:CB	2:H:29:HIS:CD2	3.02	0.43
2:B:101:VAL:CG2	3:B:289:HOH:O	2.66	0.43
1:A:72:LEU:HD12	1:A:72:LEU:N	2.34	0.42
2:H:41:TRP:CD1	2:H:86:PHE:CE2	3.07	0.42
2:H:91:LEU:N	2:H:91:LEU:CD1	2.82	0.42
2:B:218:ARG:O	2:B:253:TRP:HZ3	2.02	0.42
1:A:29:LEU:CB	1:A:106:ILE:HD13	2.47	0.42
1:G:53:LEU:O	1:G:54:LEU:HD12	2.18	0.42
1:A:53:LEU:O	1:A:54:LEU:HD13	2.18	0.42
1:A:212:GLU:HG2	1:A:212:GLU:O	2.19	0.42
2:B:190:LEU:C	2:B:190:LEU:HD12	2.44	0.42
1:G:71:ARG:C	1:G:72:LEU:HD12	2.44	0.42
1:G:6:GLN:HA	1:G:22:GLN:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:ALA:HB2	1:A:118:PHE:CD1	2.55	0.42
2:B:168:VAL:HA	2:B:226:GLN:O	2.20	0.42
2:H:27:ASN:HB3	2:H:29:HIS:NE2	2.35	0.42
1:A:21:LEU:HD22	1:A:122:THR:HG21	2.02	0.42
2:B:10:SER:HB2	2:B:167:HIS:CD2	2.51	0.42
2:B:216:ASN:HD22	2:B:219:ASN:HD21	1.68	0.42
1:A:45:ARG:HG3	1:A:46:PRO:O	2.19	0.41
2:B:43:ARG:HH21	2:B:100:ALA:HB2	1.85	0.41
2:B:45:ASP:OD1	2:B:45:ASP:N	2.52	0.41
2:B:196:LEU:N	2:B:196:LEU:CD1	2.81	0.41
1:A:113:PHE:N	1:A:113:PHE:CD2	2.88	0.41
2:B:30:ASP:OD2	2:B:31:TYR:CD2	2.73	0.41
1:G:210:ILE:O	1:G:210:ILE:HG13	2.18	0.41
2:B:43:ARG:NH2	2:B:100:ALA:HB2	2.36	0.41
2:B:255:ARG:HB3	3:B:290:HOH:O	2.20	0.41
2:H:236:TRP:CZ2	2:H:243:PRO:HD3	2.56	0.41
2:B:6:GLN:CD	2:B:121:GLY:H	2.28	0.41
2:B:227:PHE:O	2:B:245:THR:HG23	2.20	0.41
2:B:255:ARG:NH1	3:B:270:HOH:O	2.53	0.41
1:A:180:MET:HE2	1:A:183:MET:CG	2.51	0.41
1:G:82:SER:O	1:G:87:SER:HA	2.20	0.41
1:G:127:VAL:HG13	1:G:158:SER:OG	2.20	0.41
1:G:11:LEU:HD21	1:G:124:LEU:HD13	2.03	0.41
1:G:29:LEU:CD1	1:G:29:LEU:N	2.84	0.41
2:B:114:TYR:HB3	2:B:115:GLU:H	1.61	0.41
2:B:197:ASN:O	2:B:198:ASP:C	2.63	0.41
1:G:12:VAL:HA	1:G:125:SER:O	2.21	0.41
1:G:131:GLN:O	1:G:133:PRO:HD3	2.21	0.41
1:A:142:ASP:C	1:A:142:ASP:OD1	2.62	0.41
2:B:244:VAL:O	2:B:246:GLN:HG2	2.20	0.41
1:G:133:PRO:HB3	1:G:157:ASP:HB3	2.03	0.41
1:G:193:TRP:CD1	1:G:193:TRP:C	2.95	0.41
2:H:143:PRO:HD2	2:H:214:TRP:CE2	2.56	0.41
2:H:162:GLY:CA	2:H:200:ARG:HD3	2.51	0.41
1:A:42:PHE:CD1	1:A:52:SER:HA	2.56	0.41
2:B:46:THR:HA	2:B:47:GLY:HA2	1.74	0.41
2:H:80:ARG:HD2	2:H:85:ASN:O	2.21	0.41
1:A:166:LYS:HE2	1:G:168:SER:HA	2.02	0.40
1:A:3:GLN:HB3	1:A:3:GLN:HE21	1.69	0.40
1:A:169:ASP:OD1	1:A:169:ASP:N	2.53	0.40
2:B:44:GLN:HB2	2:B:50:LEU:HD23	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:170:LEU:C	2:B:170:LEU:HD23	2.46	0.40
2:B:219:ASN:O	2:B:253:TRP:HA	2.22	0.40
2:H:51:ARG:CB	2:H:71:ILE:HD11	2.52	0.40
2:H:62:THR:HG22	2:H:63:GLU:N	2.36	0.40
1:A:137:VAL:O	1:A:217:PRO:HD2	2.21	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	157/198 (79%)	145 (92%)	11 (7%)	1 (1%)	21	51
1	G	184/198 (93%)	170 (92%)	12 (6%)	2 (1%)	11	36
2	B	192/238 (81%)	180 (94%)	10 (5%)	2 (1%)	12	39
2	H	236/238 (99%)	225 (95%)	10 (4%)	1 (0%)	30	59
All	All	769/872 (88%)	720 (94%)	43 (6%)	6 (1%)	16	44

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	64	GLY
2	B	254	GLY
2	H	254	GLY
2	B	8	PRO
1	G	64	GLY
1	G	211	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	160/177 (90%)	135 (84%)	25 (16%)	2	9
1	G	166/177 (94%)	140 (84%)	26 (16%)	2	9
2	B	172/205 (84%)	164 (95%)	8 (5%)	23	56
2	H	201/205 (98%)	195 (97%)	6 (3%)	36	70
All	All	699/764 (92%)	634 (91%)	65 (9%)	8	27

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	10	SER
1	A	13	LEU
1	A	29	LEU
1	A	49	ARG
1	A	54	LEU
1	A	85	SER
1	A	90	HIS
1	A	95	GLN
1	A	96	ILE
1	A	106	ILE
1	A	113	PHE
1	A	115	LYS
1	A	123	SER
1	A	127	VAL
1	A	150	VAL
1	A	154	THR
1	A	157	ASP
1	A	164	GLN
1	A	166	LYS
1	A	178	LEU
1	A	196	LYS
1	A	198	ASP
1	A	207	ASN

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Mol	Chain	Res	Type
1	A	212	GLU
2	B	4	VAL
2	B	9	ARG
2	B	10	SER
2	B	45	ASP
2	B	48	HIS
2	B	188	GLN
2	B	196	LEU
2	B	206	ARG
1	G	3	GLN
1	G	10	SER
1	G	13	LEU
1	G	27	THR
1	G	29	LEU
1	G	49	ARG
1	G	54	LEU
1	G	85	SER
1	G	90	HIS
1	G	95	GLN
1	G	96	ILE
1	G	106	ILE
1	G	113	PHE
1	G	115	LYS
1	G	123	SER
1	G	127	VAL
1	G	150	VAL
1	G	154	THR
1	G	157	ASP
1	G	164	GLN
1	G	166	LYS
1	G	178	LEU
1	G	196	LYS
1	G	198	ASP
1	G	207	ASN
1	G	212	GLU
2	H	4	VAL
2	H	9	ARG
2	H	10	SER
2	H	45	ASP
2	H	188	GLN
2	H	206	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	30	ASN
1	A	132	ASN
1	A	159	GLN
1	A	161	ASN
2	B	44	GLN
2	B	48	HIS
2	B	152	GLN
2	B	167	HIS
2	B	219	ASN
2	B	220	HIS
2	B	246	GLN
1	G	3	GLN
1	G	132	ASN
1	G	159	GLN
2	H	44	GLN
2	H	48	HIS
2	H	83	GLN
2	H	132	ASN
2	H	167	HIS
2	H	219	ASN
2	H	220	HIS

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](#)

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	183/198 (92%)	0.66	16 (8%) 16 13	10, 29, 55, 82	0
1	G	190/198 (95%)	0.50	6 (3%) 50 41	11, 28, 56, 83	0
2	B	238/238 (100%)	1.01	29 (12%) 8 7	20, 35, 63, 98	0
2	H	238/238 (100%)	0.87	22 (9%) 14 12	17, 33, 61, 98	0
All	All	849/872 (97%)	0.78	73 (8%) 16 13	10, 32, 58, 98	0

All (73) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	240	ARG	6.2
2	H	197	ASN	5.2
2	B	113	GLY	4.1
2	H	216	ASN	4.0
1	G	197	SER	4.0
1	A	24	THR	3.7
1	G	84(C)	GLU	3.7
1	G	182	SER	3.7
2	H	199	SER	3.6
2	H	62	THR	3.6
1	A	3	GLN	3.5
2	B	199	SER	3.5
2	B	197	ASN	3.5
2	H	7	SER	3.4
2	B	239	ASP	3.4
1	A	182	SER	3.2
2	B	195	ALA	3.1
1	G	198	ASP	3.1
2	B	57	TYR	3.1
2	B	198	ASP	3.0
1	A	218	SER	2.9

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Mol	Chain	Res	Type	RSRZ
2	B	3	ALA	2.9
2	B	45	ASP	2.9
1	A	197	SER	2.9
2	B	240	ARG	2.9
2	B	242	LYS	2.8
1	A	210	ILE	2.8
2	B	231	SER	2.8
2	H	198	ASP	2.7
1	A	198	ASP	2.7
2	H	108	LEU	2.7
2	H	256	ALA	2.7
2	H	233	ASN	2.7
2	H	196	LEU	2.7
2	B	164	TYR	2.6
2	B	237	THR	2.6
2	B	175	ASN	2.6
2	B	219	ASN	2.6
1	G	84(A)	SER	2.6
2	B	109	GLY	2.6
1	A	113	PHE	2.6
2	H	232	GLU	2.5
2	B	256	ALA	2.5
2	B	218	ARG	2.5
2	H	3	ALA	2.5
1	A	165	SER	2.5
2	H	253	TRP	2.5
2	B	210	SER	2.4
2	B	65	GLY	2.4
1	A	26	SER	2.4
2	H	57	TYR	2.4
2	B	46	THR	2.4
1	A	167	ASP	2.3
1	G	3	GLN	2.3
1	A	85	SER	2.3
2	H	164	TYR	2.2
2	H	128	GLU	2.2
2	B	191	LYS	2.2
2	B	86	PHE	2.2
2	H	83	GLN	2.1
2	B	223	CYS	2.1
2	H	239	ASP	2.1
1	A	184	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
2	B	230	LEU	2.1
2	H	109	GLY	2.1
2	B	253	TRP	2.1
2	H	174	VAL	2.1
1	A	29	LEU	2.1
1	A	30	ASN	2.0
2	H	28	ASN	2.0
1	A	141	ARG	2.0
2	B	145	GLU	2.0
2	B	232	GLU	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.