



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:37 PM UTC

PDB ID : 4FZ5 / pdb_00004fz5
Title : Crystal Structure of Human TIRAP TIR-domain
Authors : Woo, J.R.; Kim, S.; Shoelson, S.E.; Park, S.
Deposited on : 2012-07-06
Resolution : 3.60 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

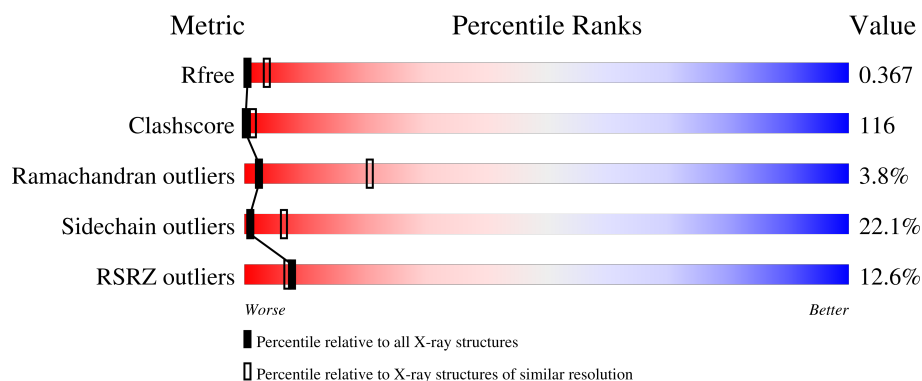
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.60 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	1747 (3.70-3.50)
Clashscore	190562	1827 (3.70-3.50)
Ramachandran outliers	187476	1773 (3.70-3.50)
Sidechain outliers	187428	1772 (3.70-3.50)
RSRZ outliers	180081	1745 (3.70-3.50)

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	154	9% 18% 38% 15% 5% 25%
1	B	154	10% 23% 32% 15% 5% 26%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1798 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 Toll/interleukin-1 receptor domain-containing adapter protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	116	Total	C	N	O	S	0	0	0
			899	567	153	171	8			
1	B	114	Total	C	N	O	S	0	0	0
			899	572	151	167	9			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	68	GLY	-	expression tag	UNP P58753
A	69	SER	-	expression tag	UNP P58753
A	70	HIS	-	expression tag	UNP P58753
A	71	MET	-	expression tag	UNP P58753
B	68	GLY	-	expression tag	UNP P58753
B	69	SER	-	expression tag	UNP P58753
B	70	HIS	-	expression tag	UNP P58753
B	71	MET	-	expression tag	UNP P58753

4 Data and refinement statistics

Property	Value	Source
Space group	P 43 2 2	Depositor
Cell constants a, b, c, α , β , γ	100.25Å 100.25Å 78.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.60 50.00 – 3.61	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-3.60) 98.8 (50.00-3.61)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.94 (at 3.56Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.324 , 0.346 0.350 , 0.367	Depositor DCC
R_{free} test set	228 reflections (4.58%)	wwPDB-VP
Wilson B-factor (Å ²)	122.5	Xtriage
Anisotropy	0.019	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 177.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.38$, $\langle L^2 \rangle = 0.21$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.83	EDS
Total number of atoms	1798	wwPDB-VP
Average B, all atoms (Å ²)	124.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.30% 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	1.11	3/914 (0.3%)	1.55	20/1233 (1.6%)
1	B	1.06	3/917 (0.3%)	1.49	20/1240 (1.6%)
All	All	1.08	6/1831 (0.3%)	1.52	40/2473 (1.6%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	188	PRO	N-CD	-17.81	1.22	1.47
1	B	82	TRP	CA-C	12.36	1.68	1.52
1	A	107	LEU	N-CA	8.51	1.57	1.46
1	B	184	ARG	CZ-NH1	-7.77	1.21	1.32
1	B	82	TRP	N-CA	6.47	1.53	1.46
1	A	107	LEU	CA-C	-6.37	1.44	1.53

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	188	PRO	N-CA-CB	-14.50	89.02	103.08
1	A	188	PRO	CA-N-CD	11.30	127.82	112.00
1	B	161	MET	N-CA-C	-9.73	99.53	111.40
1	B	163	GLN	N-CA-C	-9.71	99.71	112.34
1	B	133	LEU	N-CA-C	-9.66	94.39	109.76
1	A	214	MET	N-CA-C	9.35	121.08	111.07
1	A	187	TYR	C-N-CD	-8.96	88.28	125.00
1	B	157	CYS	N-CA-C	-8.72	101.78	111.28
1	B	165	LEU	N-CA-C	-8.65	95.16	108.31
1	B	160	GLN	N-CA-C	-8.56	102.86	113.38
1	A	88	VAL	N-CA-C	8.07	120.58	108.80
1	A	85	ASP	N-CA-C	7.80	121.28	111.69
1	A	139	SER	N-CA-C	-7.68	98.78	110.30
1	B	192	ARG	N-CA-C	-7.14	103.43	111.14
1	B	184	ARG	NE-CZ-NH2	6.80	125.32	119.20
1	A	97	LEU	N-CA-C	6.70	118.24	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	103	LEU	N-CA-C	-6.48	103.49	111.40
1	A	210	LYS	N-CA-C	-6.47	104.43	112.90
1	B	137	LEU	N-CA-C	6.35	120.31	111.74
1	B	184	ARG	NH1-CZ-NH2	-6.33	111.07	119.30
1	A	152	LEU	N-CA-C	-6.18	104.46	111.07
1	B	210	LYS	N-CA-C	-5.79	104.88	111.14
1	A	174	CYS	N-CA-C	5.61	118.23	108.76
1	B	136	ALA	N-CA-C	-5.59	102.37	110.64
1	A	159	TYR	N-CA-C	-5.57	105.20	111.28
1	B	90	VAL	CB-CA-C	-5.55	104.59	111.08
1	B	154	ASP	CA-C-N	5.54	125.20	119.05
1	B	154	ASP	C-N-CA	5.54	125.20	119.05
1	B	85	ASP	N-CA-C	5.46	122.43	110.80
1	A	100	ALA	N-CA-C	-5.44	105.78	112.90
1	A	182	LEU	CD1-CG-CD2	-5.43	98.86	110.80
1	A	106	TYR	CA-C-N	5.36	133.38	121.80
1	A	106	TYR	C-N-CA	5.36	133.38	121.80
1	A	186	ALA	CB-CA-C	-5.34	99.79	110.42
1	B	101	GLN	N-CA-C	-5.34	106.61	113.23
1	B	108	GLU	CB-CA-C	-5.31	110.47	116.63
1	A	165	LEU	N-CA-C	-5.27	106.69	113.01
1	A	187	TYR	CA-C-N	5.20	125.74	120.38
1	A	187	TYR	C-N-CA	5.20	125.74	120.38
1	B	178	LEU	N-CA-C	-5.17	101.74	109.95

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	899	0	876	214	1
1	B	899	0	881	194	4
All	All	1798	0	1757	408	5

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 116.

All (408) 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:136:ALA:C	1:A:137:LEU:HD22	1.41	1.45
1:A:133:LEU:O	1:A:135:GLN:CB	1.80	1.27
1:B:147:ILE:HB	1:B:179:LEU:CD1	1.62	1.27
1:A:210:LYS:HD2	1:A:213:VAL:CG1	1.67	1.23
1:B:142:CYS:HA	1:B:174:CYS:SG	1.81	1.20
1:A:134:CYS:CA	1:A:135:GLN:HB3	1.75	1.16
1:A:179:LEU:CD2	1:A:187:TYR:CD2	2.28	1.15
1:A:136:ALA:O	1:A:137:LEU:HD22	1.42	1.14
1:A:179:LEU:HD22	1:A:187:TYR:CD2	1.84	1.13
1:A:137:LEU:HD13	1:A:137:LEU:N	1.59	1.12
1:B:107:LEU:HD22	1:B:107:LEU:N	1.61	1.11
1:B:90:VAL:C	1:B:134:CYS:HG	1.57	1.11
1:B:147:ILE:HB	1:B:179:LEU:HD12	1.15	1.11
1:A:89:CYS:SG	1:A:143:ARG:HG3	1.89	1.10
1:B:155:PRO:HA	1:B:158:LYS:CB	1.81	1.10
1:A:97:LEU:HD23	1:A:98:VAL:N	1.67	1.09
1:A:207:ARG:HG2	1:A:211:GLU:HG3	1.34	1.09
1:B:155:PRO:CA	1:B:158:LYS:HB3	1.82	1.09
1:A:147:ILE:HG22	1:A:182:LEU:CD2	1.84	1.07
1:A:134:CYS:HA	1:A:135:GLN:CB	1.74	1.07
1:A:210:LYS:HD2	1:A:213:VAL:HG12	1.32	1.07
1:A:81:ARG:HG3	1:A:141:HIS:NE2	1.69	1.06
1:B:107:LEU:H	1:B:107:LEU:CD2	1.61	1.06
1:A:136:ALA:CB	1:A:137:LEU:HD13	1.84	1.04
1:A:136:ALA:C	1:A:137:LEU:CD2	2.30	1.04
1:B:106:TYR:CE2	1:B:206:PHE:HB3	1.93	1.03
1:A:210:LYS:HD2	1:A:213:VAL:HG11	1.42	1.02
1:B:160:GLN:HE21	1:B:160:GLN:HA	1.22	1.02
1:A:163:GLN:HE21	1:A:163:GLN:N	1.58	1.01
1:A:136:ALA:HB1	1:A:137:LEU:HD13	1.37	1.00
1:A:187:TYR:O	1:A:188:PRO:O	1.79	1.00
1:A:136:ALA:O	1:A:138:SER:OG	1.80	1.00
1:B:161:MET:O	1:B:164:ALA:N	1.93	1.00
1:A:133:LEU:O	1:A:135:GLN:HB2	1.57	1.00
1:A:187:TYR:CD2	1:A:188:PRO:HD2	1.96	1.00
1:B:160:GLN:HE21	1:B:160:GLN:CA	1.72	0.99
1:A:137:LEU:HD23	1:A:137:LEU:O	1.62	0.99
1:A:81:ARG:CD	1:A:220:LEU:HD13	1.93	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:179:LEU:CD2	1:A:187:TYR:HD2	1.76	0.98
1:B:147:ILE:HD12	1:B:179:LEU:HD11	1.44	0.98
1:A:136:ALA:CB	1:A:137:LEU:CD1	2.42	0.98
1:A:134:CYS:CA	1:A:135:GLN:CB	2.35	0.97
1:A:133:LEU:C	1:A:135:GLN:HB2	1.90	0.96
1:A:137:LEU:CD2	1:A:137:LEU:C	2.39	0.96
1:B:132:GLU:CG	1:B:135:GLN:HE21	1.78	0.96
1:A:134:CYS:HA	1:A:135:GLN:HB3	0.96	0.95
1:A:136:ALA:HB3	1:A:137:LEU:HD11	1.46	0.95
1:B:155:PRO:HA	1:B:158:LYS:HB3	0.97	0.95
1:A:179:LEU:HD21	1:A:187:TYR:CD2	1.99	0.94
1:A:210:LYS:HA	1:A:213:VAL:HB	1.49	0.94
1:A:89:CYS:SG	1:A:143:ARG:CG	2.54	0.94
1:B:106:TYR:CD2	1:B:206:PHE:HB3	2.02	0.94
1:B:147:ILE:HB	1:B:179:LEU:HD11	1.50	0.94
1:A:210:LYS:O	1:A:213:VAL:HG12	1.68	0.93
1:B:90:VAL:C	1:B:134:CYS:SG	2.50	0.93
1:A:94:GLU:HG2	1:A:95:GLU:HG3	1.47	0.93
1:B:92:HIS:CD2	1:B:97:LEU:HA	2.04	0.93
1:B:97:LEU:C	1:B:97:LEU:HD23	1.94	0.93
1:A:210:LYS:CD	1:A:213:VAL:HG11	1.99	0.92
1:A:137:LEU:HD13	1:A:137:LEU:H	1.25	0.92
1:A:82:TRP:H	1:A:82:TRP:CD1	1.76	0.92
1:B:147:ILE:CD1	1:B:179:LEU:HD11	2.00	0.92
1:A:208:GLN:O	1:A:212:ALA:N	2.03	0.91
1:B:107:LEU:HD22	1:B:107:LEU:H	0.77	0.91
1:A:210:LYS:CD	1:A:213:VAL:CG1	2.47	0.91
1:A:137:LEU:HD22	1:A:137:LEU:N	1.75	0.89
1:A:81:ARG:HG3	1:A:141:HIS:HE2	1.36	0.89
1:A:133:LEU:O	1:A:135:GLN:CG	2.19	0.89
1:B:178:LEU:C	1:B:179:LEU:HD13	1.97	0.89
1:A:81:ARG:HD2	1:A:220:LEU:HD13	1.51	0.89
1:A:136:ALA:HB3	1:A:137:LEU:CD1	1.99	0.89
1:B:104:VAL:O	1:B:107:LEU:HD23	1.73	0.88
1:A:147:ILE:HG22	1:A:182:LEU:HD23	1.57	0.86
1:A:133:LEU:C	1:A:135:GLN:CB	2.46	0.86
1:A:207:ARG:O	1:A:211:GLU:N	2.08	0.86
1:B:201:GLY:HA3	1:B:208:GLN:NE2	1.90	0.85
1:A:179:LEU:HD21	1:A:187:TYR:HD2	1.35	0.85
1:B:160:GLN:HA	1:B:160:GLN:NE2	1.91	0.84
1:A:92:HIS:HD2	1:A:93:SER:O	1.59	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:TRP:CD1	1:A:82:TRP:N	2.41	0.83
1:A:208:GLN:HA	1:A:211:GLU:HB2	1.61	0.83
1:A:137:LEU:N	1:A:137:LEU:CD1	2.30	0.82
1:A:187:TYR:HE2	1:A:191:LEU:HD12	1.42	0.82
1:A:94:GLU:C	1:A:95:GLU:HG3	2.04	0.82
1:A:187:TYR:OH	1:A:191:LEU:HB3	1.79	0.82
1:B:106:TYR:HD2	1:B:206:PHE:HD2	1.24	0.82
1:A:81:ARG:HD3	1:A:220:LEU:HD13	1.62	0.82
1:A:137:LEU:HD22	1:A:137:LEU:C	1.98	0.81
1:B:161:MET:O	1:B:164:ALA:CB	2.28	0.81
1:A:191:LEU:N	1:A:191:LEU:HD23	1.95	0.81
1:B:147:ILE:CB	1:B:179:LEU:CD1	2.56	0.81
1:A:90:VAL:HA	1:A:144:VAL:HB	1.62	0.80
1:B:142:CYS:CA	1:B:174:CYS:SG	2.68	0.80
1:A:163:GLN:HA	1:A:166:THR:OG1	1.82	0.80
1:B:103:LEU:O	1:B:107:LEU:HD21	1.81	0.80
1:B:106:TYR:HD2	1:B:206:PHE:CD2	2.00	0.80
1:A:137:LEU:CD2	1:A:137:LEU:O	2.30	0.80
1:A:92:HIS:CD2	1:A:93:SER:O	2.35	0.79
1:A:82:TRP:H	1:A:82:TRP:HD1	1.29	0.79
1:A:156:TRP:CE2	1:A:160:GLN:NE2	2.51	0.78
1:B:163:GLN:C	1:B:163:GLN:CD	2.51	0.78
1:A:136:ALA:O	1:A:137:LEU:CD2	2.30	0.78
1:B:149:PRO:HA	1:B:182:LEU:HD11	1.65	0.78
1:A:133:LEU:O	1:A:135:GLN:HB3	1.67	0.78
1:A:84:LYS:HB2	1:A:141:HIS:CE1	2.20	0.77
1:A:162:LEU:O	1:A:166:THR:OG1	2.02	0.77
1:B:86:TYR:CD1	1:B:132:GLU:HB2	2.19	0.77
1:B:155:PRO:O	1:B:158:LYS:O	2.03	0.77
1:A:156:TRP:CZ2	1:A:160:GLN:NE2	2.53	0.76
1:A:136:ALA:O	1:A:138:SER:N	2.18	0.76
1:A:209:VAL:O	1:A:213:VAL:HB	1.86	0.75
1:A:147:ILE:HG22	1:A:182:LEU:HD21	1.69	0.75
1:B:148:THR:H	1:B:151:PHE:HB3	1.50	0.75
1:A:187:TYR:CG	1:A:188:PRO:HD2	2.22	0.75
1:B:165:LEU:N	1:B:165:LEU:HD23	2.01	0.75
1:B:198:ASP:OD1	1:B:200:ARG:NE	2.20	0.75
1:B:179:LEU:HD13	1:B:179:LEU:N	2.02	0.75
1:A:174:CYS:HB2	1:A:216:TYR:CZ	2.22	0.74
1:B:103:LEU:O	1:B:107:LEU:CD2	2.35	0.74
1:A:163:GLN:NE2	1:A:163:GLN:CA	2.49	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:207:ARG:O	1:A:211:GLU:HB2	1.87	0.74
1:B:106:TYR:HD1	1:B:106:TYR:O	1.69	0.74
1:A:89:CYS:SG	1:A:89:CYS:O	2.46	0.73
1:A:136:ALA:O	1:A:137:LEU:C	2.31	0.73
1:A:191:LEU:C	1:A:193:PHE:N	2.45	0.73
1:A:163:GLN:HA	1:A:163:GLN:NE2	2.03	0.72
1:B:159:TYR:HD1	1:B:162:LEU:HD22	1.54	0.72
1:B:148:THR:HB	1:B:149:PRO:HD2	1.72	0.72
1:A:163:GLN:HE21	1:A:163:GLN:CA	2.02	0.72
1:B:94:GLU:HG2	1:B:95:GLU:H	1.55	0.71
1:B:90:VAL:HG22	1:B:103:LEU:HD23	1.72	0.71
1:B:97:LEU:C	1:B:97:LEU:CD2	2.64	0.71
1:B:164:ALA:O	1:B:165:LEU:HG	1.91	0.71
1:B:147:ILE:CB	1:B:179:LEU:HD11	2.21	0.70
1:B:97:LEU:HD21	1:B:101:GLN:OE1	1.92	0.70
1:A:88:VAL:HG23	1:A:131:SER:HA	1.73	0.70
1:A:136:ALA:HB1	1:A:137:LEU:CD1	2.11	0.70
1:B:97:LEU:HD23	1:B:97:LEU:O	1.92	0.70
1:B:149:PRO:HA	1:B:182:LEU:CD1	2.22	0.69
1:A:155:PRO:O	1:A:158:LYS:HB3	1.91	0.69
1:B:106:TYR:HE2	1:B:206:PHE:HB3	1.54	0.69
1:A:82:TRP:N	1:A:82:TRP:HD1	1.84	0.69
1:A:107:LEU:HB2	1:A:131:SER:OG	1.93	0.69
1:A:210:LYS:HD2	1:A:210:LYS:O	1.92	0.69
1:B:147:ILE:CB	1:B:179:LEU:HD12	2.09	0.69
1:A:190:GLU:N	1:A:190:GLU:OE1	2.26	0.68
1:A:92:HIS:CD2	1:A:92:HIS:C	2.71	0.68
1:B:84:LYS:HB2	1:B:141:HIS:CE1	2.29	0.68
1:A:190:GLU:CD	1:A:190:GLU:H	2.02	0.68
1:B:148:THR:H	1:B:151:PHE:CB	2.06	0.67
1:A:147:ILE:HB	1:A:179:LEU:HG	1.75	0.67
1:B:201:GLY:HA3	1:B:208:GLN:HE21	1.58	0.67
1:B:164:ALA:CB	1:B:165:LEU:HD23	2.25	0.67
1:A:81:ARG:HD3	1:A:220:LEU:CB	2.25	0.66
1:A:89:CYS:SG	1:A:143:ARG:HG2	2.35	0.66
1:A:163:GLN:CA	1:A:166:THR:OG1	2.43	0.66
1:A:188:PRO:HB2	1:A:190:GLU:OE1	1.96	0.66
1:B:103:LEU:HA	1:B:206:PHE:CE2	2.29	0.66
1:A:210:LYS:CD	1:A:213:VAL:HG12	2.16	0.66
1:A:208:GLN:CA	1:A:211:GLU:HB2	2.27	0.66
1:A:207:ARG:O	1:A:211:GLU:CB	2.44	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:LYS:HA	1:A:213:VAL:CB	2.23	0.65
1:B:132:GLU:HG3	1:B:135:GLN:HE21	1.60	0.65
1:B:178:LEU:O	1:B:179:LEU:HD13	1.95	0.65
1:A:210:LYS:CA	1:A:213:VAL:HB	2.25	0.65
1:B:148:THR:HA	1:B:180:SER:O	1.96	0.65
1:A:136:ALA:C	1:A:137:LEU:HD13	2.22	0.65
1:B:162:LEU:C	1:B:164:ALA:N	2.46	0.65
1:B:92:HIS:NE2	1:B:97:LEU:HA	2.11	0.65
1:B:94:GLU:HG2	1:B:95:GLU:N	2.11	0.65
1:B:104:VAL:C	1:B:107:LEU:HD23	2.21	0.65
1:B:160:GLN:HE21	1:B:160:GLN:C	2.04	0.65
1:A:94:GLU:HG2	1:A:95:GLU:CG	2.22	0.65
1:B:182:LEU:HD22	1:B:186:ALA:HB3	1.78	0.65
1:B:132:GLU:CG	1:B:135:GLN:NE2	2.58	0.64
1:A:191:LEU:O	1:A:192:ARG:C	2.41	0.64
1:A:81:ARG:NH1	1:A:217:LEU:HD22	2.12	0.64
1:B:149:PRO:HD3	1:B:181:GLY:C	2.23	0.64
1:A:182:LEU:HD12	1:A:182:LEU:O	1.98	0.64
1:A:198:ASP:OD1	1:A:200:ARG:HB2	1.96	0.63
1:A:135:GLN:OE1	1:A:136:ALA:N	2.30	0.62
1:A:187:TYR:O	1:A:188:PRO:C	2.41	0.62
1:B:132:GLU:HG2	1:B:135:GLN:HE21	1.63	0.62
1:A:189:PRO:O	1:A:192:ARG:N	2.32	0.62
1:A:81:ARG:HD3	1:A:220:LEU:CD1	2.28	0.62
1:A:191:LEU:O	1:A:193:PHE:N	2.33	0.62
1:A:133:LEU:O	1:A:135:GLN:HG2	1.97	0.62
1:A:187:TYR:CE2	1:A:191:LEU:HD12	2.30	0.62
1:A:81:ARG:HH12	1:A:217:LEU:HD22	1.64	0.62
1:B:149:PRO:HB3	1:B:182:LEU:HG	1.82	0.62
1:A:153:GLN:O	1:A:155:PRO:HD3	2.00	0.62
1:B:106:TYR:CD2	1:B:206:PHE:HD2	2.12	0.62
1:B:163:GLN:OE1	1:B:164:ALA:HA	2.00	0.62
1:B:92:HIS:CD2	1:B:97:LEU:CA	2.83	0.61
1:A:190:GLU:C	1:A:191:LEU:HD23	2.26	0.61
1:A:89:CYS:O	1:A:144:VAL:N	2.30	0.60
1:A:162:LEU:C	1:A:166:THR:OG1	2.44	0.60
1:A:179:LEU:CD2	1:A:187:TYR:CE2	2.84	0.60
1:A:145:LEU:HB2	1:A:177:PRO:HA	1.84	0.60
1:A:187:TYR:C	1:A:188:PRO:O	2.39	0.60
1:A:104:VAL:O	1:A:108:GLU:HG2	2.01	0.60
1:B:188:PRO:HG2	1:B:191:LEU:HD12	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:88:VAL:HG12	1:B:89:CYS:N	2.17	0.60
1:B:147:ILE:HG12	1:B:151:PHE:CE1	2.37	0.59
1:A:133:LEU:C	1:A:135:GLN:HB3	2.22	0.59
1:B:155:PRO:C	1:B:158:LYS:HB3	2.27	0.59
1:A:191:LEU:N	1:A:191:LEU:CD2	2.65	0.59
1:B:106:TYR:C	1:B:106:TYR:CD1	2.80	0.59
1:B:160:GLN:O	1:B:163:GLN:HG3	2.02	0.59
1:B:89:CYS:C	1:B:134:CYS:SG	2.86	0.59
1:A:81:ARG:HD3	1:A:220:LEU:HB3	1.84	0.59
1:B:160:GLN:HE22	1:B:163:GLN:HG2	1.66	0.59
1:A:97:LEU:HD23	1:A:98:VAL:H	1.65	0.59
1:A:162:LEU:C	1:A:166:THR:HG1	2.11	0.59
1:B:155:PRO:HA	1:B:158:LYS:CG	2.31	0.59
1:A:189:PRO:C	1:A:191:LEU:N	2.56	0.59
1:A:132:GLU:OE1	1:A:138:SER:HB3	2.01	0.59
1:B:135:GLN:O	1:B:136:ALA:C	2.45	0.58
1:A:157:CYS:HA	1:A:160:GLN:HG3	1.84	0.58
1:B:104:VAL:O	1:B:107:LEU:CD2	2.51	0.58
1:B:198:ASP:CG	1:B:200:ARG:HE	2.12	0.58
1:B:162:LEU:C	1:B:164:ALA:H	2.10	0.58
1:A:210:LYS:HD3	1:A:213:VAL:HG11	1.83	0.58
1:B:201:GLY:O	1:B:202:PRO:C	2.47	0.58
1:A:187:TYR:OH	1:A:191:LEU:CB	2.50	0.58
1:A:97:LEU:HD23	1:A:98:VAL:CA	2.34	0.57
1:A:202:PRO:O	1:A:203:ASP:HB2	2.04	0.57
1:B:161:MET:O	1:B:164:ALA:CA	2.51	0.57
1:A:182:LEU:HD12	1:A:182:LEU:C	2.29	0.57
1:A:97:LEU:CD2	1:A:98:VAL:N	2.57	0.57
1:A:94:GLU:CG	1:A:95:GLU:HG3	2.28	0.56
1:A:157:CYS:O	1:A:160:GLN:HB2	2.05	0.56
1:B:106:TYR:HD1	1:B:106:TYR:C	2.11	0.56
1:A:163:GLN:N	1:A:163:GLN:NE2	2.41	0.56
1:A:107:LEU:HD23	1:A:210:LYS:HB2	1.87	0.56
1:B:178:LEU:O	1:B:179:LEU:CD1	2.53	0.56
1:A:150:GLY:O	1:A:154:ASP:N	2.37	0.55
1:B:160:GLN:CA	1:B:160:GLN:NE2	2.48	0.55
1:A:157:CYS:HA	1:A:160:GLN:CG	2.36	0.55
1:B:182:LEU:HD22	1:B:186:ALA:CB	2.37	0.55
1:A:155:PRO:HA	1:A:158:LYS:CB	2.37	0.55
1:B:159:TYR:HA	1:B:162:LEU:HB2	1.88	0.55
1:B:164:ALA:C	1:B:165:LEU:CG	2.80	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:164:ALA:C	1:B:165:LEU:HG	2.31	0.55
1:B:137:LEU:HD13	1:B:137:LEU:O	2.07	0.54
1:B:182:LEU:HB3	1:B:187:TYR:HE1	1.73	0.54
1:B:158:LYS:O	1:B:159:TYR:HB3	2.06	0.54
1:A:155:PRO:HA	1:A:158:LYS:HB3	1.88	0.54
1:B:179:LEU:HB3	1:B:187:TYR:CZ	2.43	0.54
1:B:185:ALA:O	1:B:186:ALA:HB3	2.07	0.54
1:B:81:ARG:O	1:B:82:TRP:HD1	1.91	0.54
1:B:132:GLU:CD	1:B:135:GLN:HE21	2.15	0.54
1:B:201:GLY:CA	1:B:208:GLN:NE2	2.69	0.54
1:A:81:ARG:CD	1:A:220:LEU:CD1	2.78	0.54
1:A:134:CYS:N	1:A:135:GLN:CB	2.71	0.54
1:B:90:VAL:CG2	1:B:103:LEU:HD23	2.38	0.53
1:B:154:ASP:C	1:B:154:ASP:OD1	2.50	0.53
1:B:165:LEU:O	1:B:166:THR:C	2.51	0.53
1:B:183:SER:O	1:B:184:ARG:C	2.51	0.53
1:B:163:GLN:O	1:B:166:THR:OG1	2.27	0.53
1:A:97:LEU:HD23	1:A:97:LEU:C	2.31	0.53
1:B:87:ASP:O	1:B:142:CYS:HB2	2.09	0.53
1:B:92:HIS:HD2	1:B:96:ASP:C	2.16	0.53
1:A:210:LYS:O	1:A:210:LYS:CD	2.57	0.53
1:A:179:LEU:HD21	1:A:187:TYR:CE2	2.41	0.52
1:A:187:TYR:CZ	1:A:191:LEU:HB2	2.44	0.52
1:B:188:PRO:HG2	1:B:191:LEU:CD1	2.39	0.52
1:A:162:LEU:O	1:A:166:THR:N	2.43	0.52
1:B:101:GLN:HG3	1:B:133:LEU:HD11	1.90	0.52
1:B:101:GLN:HG2	1:B:133:LEU:HD13	1.91	0.52
1:A:92:HIS:CD2	1:A:92:HIS:O	2.63	0.52
1:A:156:TRP:O	1:A:159:TYR:N	2.43	0.52
1:A:179:LEU:HD23	1:A:182:LEU:HD21	1.91	0.52
1:A:189:PRO:HD2	1:A:190:GLU:CD	2.35	0.52
1:B:164:ALA:C	1:B:165:LEU:HD23	2.35	0.52
1:B:161:MET:O	1:B:164:ALA:HB2	2.09	0.52
1:A:156:TRP:O	1:A:157:CYS:C	2.51	0.51
1:B:97:LEU:CD2	1:B:101:GLN:OE1	2.56	0.51
1:B:94:GLU:CG	1:B:95:GLU:N	2.73	0.51
1:B:147:ILE:CG1	1:B:179:LEU:HD11	2.40	0.51
1:B:159:TYR:CD1	1:B:162:LEU:HD22	2.40	0.51
1:B:160:GLN:NE2	1:B:160:GLN:O	2.44	0.51
1:B:104:VAL:C	1:B:107:LEU:CD2	2.84	0.51
1:B:205:GLY:HA2	1:B:208:GLN:NE2	2.26	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:160:GLN:HE22	1:B:163:GLN:CG	2.24	0.51
1:A:183:SER:OG	1:A:185:ALA:HB3	2.10	0.51
1:A:189:PRO:O	1:A:191:LEU:N	2.44	0.51
1:A:81:ARG:HD3	1:A:220:LEU:CG	2.41	0.50
1:A:88:VAL:HG21	1:A:107:LEU:HD13	1.94	0.50
1:A:189:PRO:O	1:A:190:GLU:C	2.55	0.50
1:B:185:ALA:O	1:B:186:ALA:CB	2.58	0.50
1:A:81:ARG:CG	1:A:141:HIS:HE2	2.17	0.50
1:B:132:GLU:HG2	1:B:135:GLN:NE2	2.23	0.50
1:B:164:ALA:HB3	1:B:165:LEU:HD23	1.93	0.50
1:A:191:LEU:C	1:A:193:PHE:H	2.17	0.50
1:B:107:LEU:HA	1:B:210:LYS:HE3	1.92	0.50
1:A:189:PRO:HD2	1:A:190:GLU:OE1	2.13	0.49
1:B:159:TYR:CD1	1:B:162:LEU:CD2	2.95	0.49
1:B:149:PRO:HD3	1:B:181:GLY:O	2.13	0.49
1:A:107:LEU:CD2	1:A:210:LYS:HB2	2.43	0.49
1:A:147:ILE:CG2	1:A:182:LEU:CD2	2.73	0.49
1:B:92:HIS:HD2	1:B:97:LEU:HA	1.71	0.49
1:B:147:ILE:HA	1:B:151:PHE:CG	2.48	0.49
1:A:188:PRO:HG2	1:A:191:LEU:HG	1.94	0.49
1:A:138:SER:O	1:A:139:SER:C	2.55	0.49
1:A:187:TYR:CZ	1:A:191:LEU:CB	2.95	0.49
1:A:207:ARG:C	1:A:211:GLU:HB2	2.38	0.49
1:B:161:MET:O	1:B:164:ALA:HB3	2.09	0.49
1:B:142:CYS:CB	1:B:174:CYS:SG	3.01	0.49
1:A:183:SER:C	1:A:185:ALA:H	2.21	0.48
1:A:174:CYS:HB2	1:A:216:TYR:CE1	2.47	0.48
1:B:213:VAL:O	1:B:216:TYR:N	2.44	0.48
1:A:162:LEU:H	1:A:162:LEU:HD12	1.78	0.48
1:B:106:TYR:HD2	1:B:206:PHE:HB3	1.71	0.48
1:A:94:GLU:O	1:A:95:GLU:HG3	2.14	0.47
1:A:210:LYS:HD2	1:A:210:LYS:C	2.37	0.47
1:B:132:GLU:OE2	1:B:135:GLN:HG3	2.14	0.47
1:B:163:GLN:OE1	1:B:164:ALA:CA	2.63	0.47
1:B:164:ALA:HB3	1:B:165:LEU:CD2	2.43	0.47
1:B:163:GLN:CD	1:B:163:GLN:O	2.56	0.47
1:A:134:CYS:N	1:A:135:GLN:HB3	2.27	0.47
1:A:189:PRO:C	1:A:191:LEU:H	2.21	0.47
1:B:161:MET:HE1	1:B:188:PRO:HG3	1.95	0.47
1:B:81:ARG:O	1:B:81:ARG:HG2	2.15	0.47
1:B:104:VAL:HA	1:B:107:LEU:HD23	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:183:SER:C	1:B:185:ALA:N	2.68	0.47
1:A:132:GLU:HG2	1:A:133:LEU:N	2.29	0.47
1:B:89:CYS:HG	1:B:134:CYS:CB	2.27	0.47
1:B:159:TYR:CE1	1:B:162:LEU:HD23	2.50	0.47
1:B:101:GLN:CG	1:B:133:LEU:CD1	2.93	0.46
1:A:214:MET:O	1:A:215:ARG:C	2.58	0.46
1:A:210:LYS:HA	1:A:213:VAL:CG1	2.45	0.46
1:A:97:LEU:O	1:A:98:VAL:C	2.58	0.46
1:B:91:CYS:N	1:B:134:CYS:SG	2.89	0.46
1:B:104:VAL:CA	1:B:107:LEU:HD23	2.45	0.46
1:A:93:SER:OG	1:A:94:GLU:N	2.47	0.46
1:A:104:VAL:HG12	1:A:108:GLU:OE2	2.15	0.46
1:B:84:LYS:HB2	1:B:141:HIS:ND1	2.30	0.46
1:A:107:LEU:CB	1:A:131:SER:OG	2.62	0.45
1:B:137:LEU:HA	1:B:137:LEU:HD22	1.40	0.45
1:B:88:VAL:CG1	1:B:89:CYS:N	2.80	0.45
1:A:183:SER:OG	1:A:185:ALA:CB	2.65	0.45
1:B:90:VAL:O	1:B:134:CYS:SG	2.46	0.45
1:B:201:GLY:CA	1:B:208:GLN:HE22	2.28	0.45
1:B:135:GLN:H	1:B:135:GLN:HG2	1.58	0.45
1:B:162:LEU:O	1:B:163:GLN:C	2.59	0.45
1:B:163:GLN:OE1	1:B:164:ALA:N	2.50	0.45
1:A:86:TYR:O	1:A:140:SER:HA	2.16	0.45
1:A:156:TRP:C	1:A:158:LYS:N	2.71	0.45
1:B:135:GLN:HB2	1:B:138:SER:HB3	1.99	0.45
1:A:157:CYS:O	1:A:160:GLN:CB	2.65	0.44
1:B:163:GLN:C	1:B:163:GLN:OE1	2.59	0.44
1:A:162:LEU:HD12	1:A:162:LEU:N	2.32	0.44
1:A:187:TYR:CD2	1:A:188:PRO:CD	2.86	0.44
1:A:107:LEU:HB3	1:A:129:ILE:HG21	1.98	0.44
1:A:142:CYS:SG	1:A:174:CYS:HB3	2.57	0.44
1:B:107:LEU:HB2	1:B:131:SER:OG	2.17	0.44
1:B:147:ILE:HD12	1:B:179:LEU:CD1	2.32	0.44
1:A:132:GLU:CD	1:A:138:SER:HB3	2.43	0.44
1:A:216:TYR:O	1:A:219:THR:HB	2.18	0.44
1:B:164:ALA:O	1:B:165:LEU:C	2.59	0.44
1:B:183:SER:O	1:B:185:ALA:O	2.35	0.44
1:B:101:GLN:HG3	1:B:133:LEU:CD1	2.49	0.43
1:B:147:ILE:O	1:B:179:LEU:HA	2.17	0.43
1:A:147:ILE:CG2	1:A:182:LEU:HD21	2.43	0.43
1:A:148:THR:HB	1:A:149:PRO:CD	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:90:VAL:N	1:B:134:CYS:SG	2.92	0.43
1:A:142:CYS:SG	1:A:174:CYS:CB	3.06	0.43
1:B:103:LEU:C	1:B:107:LEU:HD21	2.41	0.43
1:B:106:TYR:HE2	1:B:206:PHE:CB	2.28	0.43
1:B:154:ASP:OD1	1:B:155:PRO:HD2	2.19	0.43
1:A:86:TYR:CE1	1:A:132:GLU:HB2	2.53	0.43
1:A:94:GLU:HG2	1:A:95:GLU:N	2.32	0.43
1:A:134:CYS:CA	1:A:135:GLN:O	2.66	0.43
1:A:156:TRP:NE1	1:A:160:GLN:HE21	2.15	0.43
1:A:179:LEU:HD23	1:A:182:LEU:CD2	2.49	0.43
1:B:92:HIS:CD2	1:B:97:LEU:N	2.87	0.43
1:A:134:CYS:HA	1:A:135:GLN:O	2.19	0.43
1:A:81:ARG:HG3	1:A:141:HIS:CD2	2.47	0.42
1:B:104:VAL:HA	1:B:107:LEU:CD2	2.49	0.42
1:B:159:TYR:O	1:B:159:TYR:CG	2.72	0.42
1:B:132:GLU:HG2	1:B:133:LEU:O	2.18	0.42
1:B:176:ILE:HD11	1:B:216:TYR:CE2	2.55	0.42
1:B:81:ARG:C	1:B:82:TRP:CD1	2.98	0.42
1:B:92:HIS:NE2	1:B:97:LEU:CA	2.82	0.42
1:B:164:ALA:HB1	1:B:165:LEU:HD23	1.99	0.42
1:B:101:GLN:HG2	1:B:133:LEU:CD1	2.49	0.42
1:A:199:GLY:HA2	1:A:204:GLY:O	2.20	0.42
1:B:132:GLU:HG3	1:B:135:GLN:NE2	2.32	0.42
1:B:158:LYS:O	1:B:159:TYR:CB	2.68	0.42
1:B:159:TYR:HD1	1:B:162:LEU:CD2	2.25	0.42
1:B:89:CYS:SG	1:B:134:CYS:HB3	2.60	0.41
1:B:81:ARG:O	1:B:82:TRP:CD1	2.72	0.41
1:B:92:HIS:HD2	1:B:97:LEU:N	2.18	0.41
1:B:104:VAL:CA	1:B:107:LEU:CD2	2.99	0.41
1:B:86:TYR:CE1	1:B:132:GLU:HB2	2.54	0.41
1:B:142:CYS:SG	1:B:213:VAL:HG22	2.60	0.41
1:A:136:ALA:C	1:A:137:LEU:CD1	2.88	0.41
1:A:107:LEU:O	1:A:129:ILE:HB	2.21	0.41
1:A:197:VAL:HG11	1:A:209:VAL:HG13	2.03	0.41
1:B:178:LEU:HD12	1:B:197:VAL:O	2.20	0.41
1:A:207:ARG:O	1:A:211:GLU:CA	2.68	0.41
1:B:103:LEU:O	1:B:107:LEU:HD22	2.18	0.41
1:B:178:LEU:HD11	1:B:199:GLY:HA2	2.02	0.41
1:B:179:LEU:HD12	1:B:179:LEU:HA	1.43	0.41
1:A:160:GLN:H	1:A:160:GLN:HG2	1.68	0.41
1:B:107:LEU:O	1:B:108:GLU:C	2.63	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:155:PRO:HA	1:B:158:LYS:HG2	2.02	0.40

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:141:HIS:NE2	1:B:184:ARG:NH1[3_544]	1.70	0.50
1:B:83:SER:N	1:B:184:ARG:NH2[3_544]	1.90	0.30
1:B:82:TRP:CA	1:B:184:ARG:NE[3_544]	1.99	0.21
1:A:82:TRP:CB	1:A:184:ARG:O[3_554]	2.03	0.17
1:B:82:TRP:N	1:B:184:ARG:NH2[3_544]	2.13	0.07

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	106/154 (69%)	86 (81%)	14 (13%)	6 (6%)	1	13
1	B	106/154 (69%)	91 (86%)	13 (12%)	2 (2%)	6	33
All	All	212/308 (69%)	177 (84%)	27 (13%)	8 (4%)	2	21

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	137	LEU
1	A	188	PRO
1	A	186	ALA
1	B	85	ASP
1	A	192	ARG
1	A	135	GLN
1	B	202	PRO
1	A	189	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	99/128 (77%)	78 (79%)	21 (21%)	1	7
1	B	100/128 (78%)	77 (77%)	23 (23%)	1	6
All	All	199/256 (78%)	155 (78%)	44 (22%)	1	6

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	80	SER
1	A	81	ARG
1	A	82	TRP
1	A	88	VAL
1	A	89	CYS
1	A	95	GLU
1	A	97	LEU
1	A	135	GLN
1	A	137	LEU
1	A	138	SER
1	A	139	SER
1	A	140	SER
1	A	142	CYS
1	A	163	GLN
1	A	166	THR
1	A	183	SER
1	A	184	ARG
1	A	191	LEU
1	A	211	GLU
1	A	213	VAL
1	A	214	MET
1	B	85	ASP
1	B	89	CYS
1	B	93	SER
1	B	94	GLU
1	B	97	LEU
1	B	107	LEU

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Mol	Chain	Res	Type
1	B	133	LEU
1	B	134	CYS
1	B	135	GLN
1	B	137	LEU
1	B	154	ASP
1	B	155	PRO
1	B	158	LYS
1	B	160	GLN
1	B	162	LEU
1	B	163	GLN
1	B	165	LEU
1	B	174	CYS
1	B	179	LEU
1	B	180	SER
1	B	209	VAL
1	B	215	ARG
1	B	219	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	92	HIS
1	A	163	GLN
1	B	135	GLN
1	B	160	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	116/154 (75%)	0.75	14 (12%) 8 7	29, 115, 174, 199	0
1	B	114/154 (74%)	0.85	15 (13%) 7 6	32, 134, 190, 199	0
All	All	230/308 (74%)	0.80	29 (12%) 8 7	29, 123, 178, 199	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	187	TYR	4.5
1	B	214	MET	4.3
1	B	208	GLN	3.3
1	B	109	GLY	3.1
1	A	128	ALA	3.0
1	A	203	ASP	2.9
1	B	180	SER	2.8
1	B	176	ILE	2.7
1	A	82	TRP	2.6
1	A	178	LEU	2.5
1	A	193	PHE	2.4
1	A	214	MET	2.4
1	A	210	LYS	2.3
1	B	217	LEU	2.3
1	A	155	PRO	2.3
1	B	203	ASP	2.3
1	A	110	SER	2.2
1	A	133	LEU	2.2
1	B	143	ARG	2.2
1	B	216	TYR	2.1
1	B	85	ASP	2.1
1	A	217	LEU	2.1
1	B	151	PHE	2.1
1	B	166	THR	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	206	PHE	2.1
1	B	98	VAL	2.1
1	B	106	TYR	2.0
1	B	200	ARG	2.0
1	A	183	SER	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.