



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:03 AM UTC

PDB ID : 2VH2 / pdb_00002vh2
Title : Crystal structure of cell division protein FtsQ from *Yersinia enterocolitica*
Authors : van den Ent, F.; Vinkenvleugel, T.; Ind, A.; West, P.; Veprintsev, D.; Nanninga, N.; den Blaauwen, T.; Lowe, J.
Deposited on : 2007-11-16
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

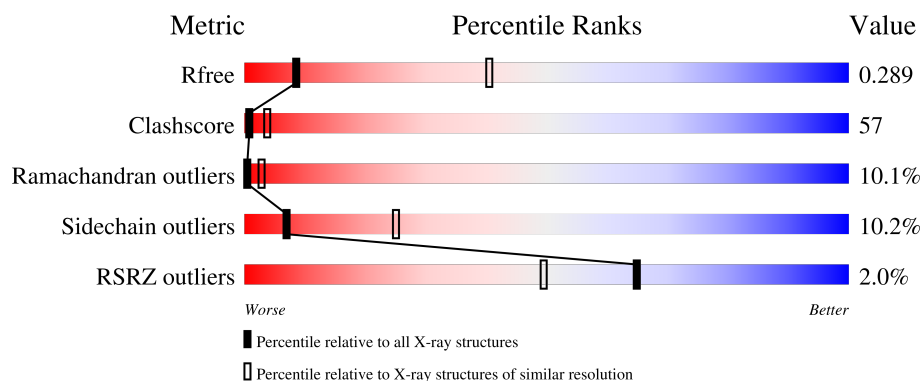
1 Overall quality at a glance

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	255	
1	B	255	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3282 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.

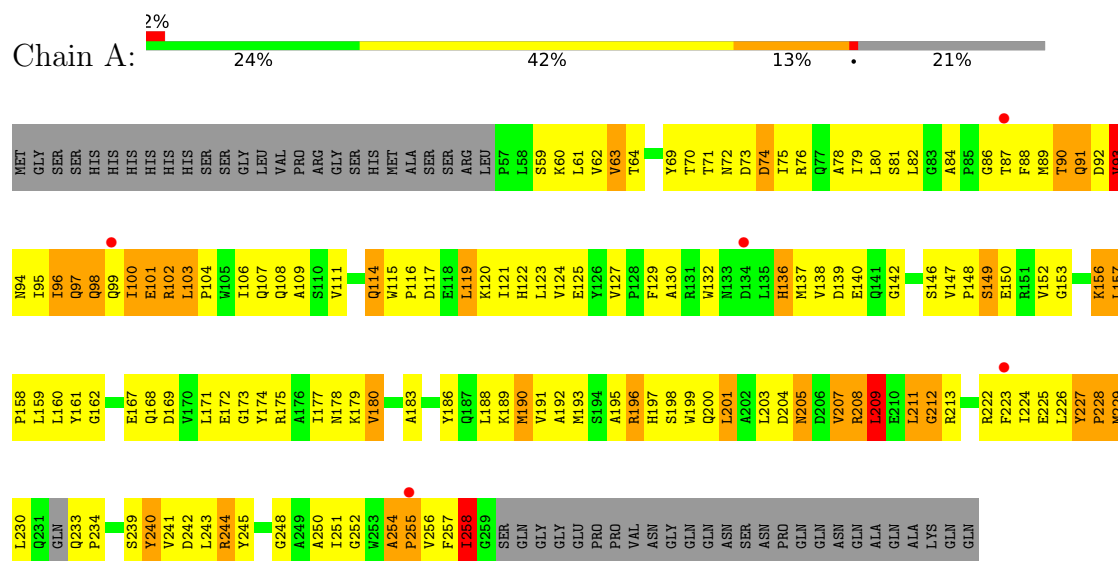
- Molecule 1 is a protein called CELL DIVISION PROTEIN FTSQ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	202	Total	C	N	O	S	0	0	1
			1639	1042	294	297	6			
1	B	202	Total	C	N	O	S	0	0	1
			1643	1045	297	295	6			

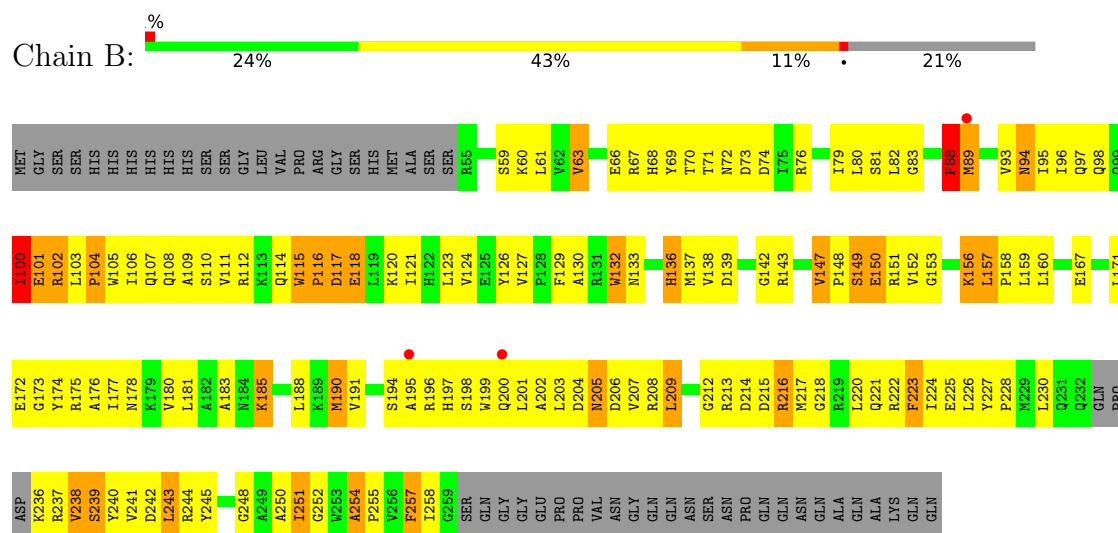
3 Residue-property plots

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: CELL DIVISION PROTEIN FTSQ



• Molecule 1: CELL DIVISION PROTEIN FTSQ



4 Data and refinement statistics

Property	Value	Source
Space group	P 62	Depositor
Cell constants a, b, c, α , β , γ	160.76Å 160.76Å 54.58Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	35.00 – 3.40 35.00 – 3.40	Depositor EDS
% Data completeness (in resolution range)	99.4 (35.00-3.40) 99.3 (35.00-3.40)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.52 (at 3.18Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.277 , 0.323 0.282 , 0.289	Depositor DCC
R_{free} test set	659 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	135.4	Xtriage
Anisotropy	0.420	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 98.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.054 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3282	wwPDB-VP
Average B, all atoms (Å ²)	151.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.73% 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	1/1675 (0.1%)	1.12	13/2268 (0.6%)
1	B	0.63	1/1678 (0.1%)	1.09	12/2271 (0.5%)
All	All	0.64	2/3353 (0.1%)	1.11	25/4539 (0.6%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	258	ILE	C-N	-6.05	1.24	1.33
1	A	258	ILE	C-N	-5.67	1.25	1.33

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	212	GLY	N-CA-C	9.99	123.34	111.35
1	B	115	TRP	N-CA-C	8.70	124.56	113.25
1	B	205	ASN	N-CA-C	-8.49	95.45	108.96
1	B	254	ALA	CA-C-N	8.35	130.28	119.84
1	B	254	ALA	C-N-CA	8.35	130.28	119.84
1	A	70	THR	N-CA-C	7.64	121.11	108.96
1	A	114	GLN	N-CA-C	-7.07	97.71	108.96
1	A	93	VAL	N-CA-C	-6.67	104.29	112.76
1	A	180	VAL	N-CA-C	6.17	116.35	110.42
1	A	254	ALA	N-CA-C	5.96	117.53	109.24
1	B	157	LEU	CA-C-N	5.94	127.27	119.84
1	B	157	LEU	C-N-CA	5.94	127.27	119.84
1	A	74	ASP	N-CA-C	-5.90	105.18	112.38
1	A	254	ALA	CA-C-N	5.86	127.17	119.84
1	A	254	ALA	C-N-CA	5.86	127.17	119.84
1	B	70	THR	N-CA-C	5.85	119.07	109.76
1	A	157	LEU	CA-C-N	5.82	127.12	119.84
1	A	157	LEU	C-N-CA	5.82	127.12	119.84
1	A	256	VAL	N-CA-C	5.78	116.75	110.21
1	B	223	PHE	N-CA-C	5.78	118.32	111.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	205	ASN	N-CA-C	-5.57	103.14	110.33
1	B	118	GLU	N-CA-C	5.38	118.70	110.20
1	B	133	ASN	N-CA-C	-5.29	99.52	110.80
1	B	132	TRP	N-CA-C	5.27	118.17	111.69
1	B	66	GLU	N-CA-C	5.01	117.36	109.39

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	1639	0	1638	194	0
1	B	1643	0	1650	192	0
All	All	3282	0	3288	377	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 57.

All (377) 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:102:ARG:C	1:A:104:PRO:HA	1.80	1.07
1:A:245:TYR:HE2	1:A:250:ALA:HB2	1.19	1.04
1:B:104:PRO:HG2	1:B:106:ILE:HD13	1.38	1.03
1:B:71:THR:HB	1:B:74:ASP:OD1	1.59	1.03
1:B:177:ILE:HD11	1:B:216:ARG:NH1	1.71	1.03
1:B:177:ILE:HD11	1:B:216:ARG:HH12	1.17	1.03
1:A:61:LEU:HD11	1:A:76:ARG:HH11	1.29	0.97
1:B:198:SER:HA	1:B:212:GLY:O	1.64	0.96
1:B:88:PHE:CD2	1:B:88:PHE:O	2.18	0.96
1:B:97:GLN:O	1:B:100:ILE:HG22	1.68	0.93
1:B:157:LEU:HD22	1:B:158:PRO:HD2	1.51	0.93
1:A:201:LEU:HD21	1:A:211:LEU:HD23	1.50	0.91
1:A:196:ARG:HB2	1:B:68:HIS:NE2	1.88	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:196:ARG:HB2	1:B:68:HIS:CE1	2.08	0.87
1:B:238:VAL:HG12	1:B:239:SER:H	1.38	0.86
1:B:114:GLN:HB2	1:B:118:GLU:HG3	1.56	0.85
1:A:245:TYR:CE2	1:A:250:ALA:HB2	2.12	0.82
1:B:203:LEU:HB2	1:B:207:VAL:HG12	1.61	0.82
1:B:200:GLN:NE2	1:B:208:ARG:HD2	1.94	0.82
1:B:88:PHE:O	1:B:88:PHE:HD2	1.59	0.81
1:A:61:LEU:CD1	1:A:76:ARG:HH11	1.92	0.81
1:B:61:LEU:HD11	1:B:76:ARG:HH11	1.44	0.80
1:B:101:GLU:O	1:B:103:LEU:N	2.14	0.80
1:B:245:TYR:HE2	1:B:250:ALA:HB2	1.47	0.80
1:A:103:LEU:N	1:A:104:PRO:HA	1.98	0.79
1:A:157:LEU:HD22	1:A:158:PRO:HD2	1.64	0.79
1:B:101:GLU:HG3	1:B:102:ARG:H	1.47	0.79
1:B:71:THR:HG22	1:B:73:ASP:H	1.48	0.78
1:B:104:PRO:HB2	1:B:105:TRP:CE3	2.19	0.78
1:B:114:GLN:HB2	1:B:118:GLU:CG	2.14	0.77
1:B:101:GLU:C	1:B:103:LEU:H	1.91	0.76
1:B:79:ILE:CD1	1:B:96:ILE:HG23	2.14	0.76
1:B:227:TYR:O	1:B:230:LEU:HG	1.86	0.76
1:B:207:VAL:HG22	1:B:208:ARG:H	1.50	0.76
1:A:198:SER:HA	1:A:212:GLY:O	1.86	0.75
1:B:108:GLN:O	1:B:124:VAL:HG22	1.85	0.75
1:A:245:TYR:HD1	1:B:72:ASN:ND2	1.84	0.75
1:A:115:TRP:O	1:A:116:PRO:C	2.31	0.74
1:B:130:ALA:HB1	1:B:160:LEU:HD13	1.69	0.74
1:B:88:PHE:O	1:B:89:MET:O	2.07	0.73
1:A:101:GLU:O	1:A:102:ARG:C	2.31	0.72
1:A:245:TYR:CD1	1:B:72:ASN:ND2	2.57	0.72
1:A:79:ILE:HD11	1:A:95:ILE:HD12	1.71	0.72
1:A:104:PRO:HG2	1:A:146:SER:H	1.55	0.71
1:A:71:THR:HB	1:A:74:ASP:OD1	1.90	0.71
1:A:213:ARG:HH12	1:B:152:VAL:HG12	1.54	0.71
1:A:188:LEU:HD21	1:A:190:MET:O	1.90	0.71
1:A:64:THR:CG2	1:A:120:LYS:HE3	2.20	0.71
1:A:89:MET:HG3	1:A:90:THR:N	2.04	0.71
1:B:188:LEU:HD23	1:B:188:LEU:C	2.16	0.71
1:A:188:LEU:HD23	1:A:189:LYS:N	2.05	0.70
1:B:79:ILE:HD13	1:B:96:ILE:HG23	1.74	0.70
1:A:233:GLN:CD	1:A:234:PRO:HD2	2.16	0.69
1:A:188:LEU:HD23	1:A:188:LEU:C	2.17	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:239:SER:O	1:A:240:TYR:HB3	1.92	0.69
1:A:208:ARG:HG2	1:A:209:LEU:H	1.56	0.69
1:A:188:LEU:HD23	1:A:190:MET:N	2.07	0.69
1:B:100:ILE:HD13	1:B:100:ILE:O	1.93	0.69
1:A:186:TYR:HB3	1:A:203:LEU:HD12	1.75	0.68
1:B:97:GLN:HE21	1:B:110:SER:HA	1.58	0.66
1:A:208:ARG:CG	1:A:209:LEU:H	2.07	0.66
1:B:130:ALA:HA	1:B:157:LEU:HD11	1.77	0.66
1:A:132:TRP:CH2	1:A:167:GLU:HA	2.29	0.66
1:B:101:GLU:O	1:B:104:PRO:HD2	1.95	0.66
1:B:104:PRO:CG	1:B:106:ILE:HD13	2.20	0.66
1:B:130:ALA:O	1:B:138:VAL:HG12	1.96	0.66
1:A:103:LEU:N	1:A:104:PRO:CA	2.59	0.66
1:B:241:VAL:HG12	1:B:243:LEU:CD1	2.25	0.65
1:A:79:ILE:CD1	1:A:95:ILE:HG23	2.25	0.65
1:A:213:ARG:NH1	1:B:152:VAL:HA	2.12	0.65
1:A:222:ARG:NH1	1:A:248:GLY:C	2.54	0.65
1:A:257:PHE:HD1	1:A:258:ILE:N	1.94	0.65
1:B:114:GLN:CB	1:B:118:GLU:HG3	2.27	0.65
1:B:61:LEU:CD1	1:B:76:ARG:HH11	2.10	0.65
1:A:61:LEU:N	1:A:61:LEU:HD12	2.12	0.64
1:B:132:TRP:CH2	1:B:167:GLU:HA	2.32	0.64
1:A:148:PRO:C	1:A:150:GLU:H	2.06	0.64
1:B:101:GLU:HG3	1:B:102:ARG:N	2.13	0.64
1:B:202:ALA:O	1:B:203:LEU:HD12	1.98	0.64
1:B:177:ILE:CD1	1:B:216:ARG:HH12	2.02	0.64
1:B:181:LEU:CD1	1:B:188:LEU:HD12	2.28	0.64
1:A:159:LEU:C	1:A:160:LEU:HD12	2.23	0.63
1:B:93:VAL:HG23	1:B:94:ASN:H	1.62	0.63
1:A:61:LEU:CD1	1:A:76:ARG:NH1	2.59	0.63
1:B:180:VAL:O	1:B:183:ALA:HB3	1.99	0.63
1:B:129:PHE:C	1:B:157:LEU:HD21	2.24	0.63
1:A:201:LEU:CD2	1:A:211:LEU:HD23	2.26	0.63
1:A:257:PHE:CD1	1:A:258:ILE:N	2.67	0.63
1:B:138:VAL:HG22	1:B:139:ASP:N	2.14	0.63
1:A:79:ILE:CG2	1:A:80:LEU:N	2.61	0.62
1:A:233:GLN:OE1	1:A:234:PRO:HD2	1.97	0.62
1:B:203:LEU:HD22	1:B:207:VAL:HG11	1.81	0.62
1:A:227:TYR:C	1:A:229:MET:H	2.08	0.62
1:A:79:ILE:HG23	1:A:80:LEU:N	2.14	0.62
1:B:196:ARG:O	1:B:198:SER:N	2.32	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:ASN:HB3	1:A:207:VAL:HG12	1.83	0.61
1:A:62:VAL:O	1:A:120:LYS:HD2	2.01	0.61
1:B:226:LEU:N	1:B:226:LEU:HD12	2.15	0.61
1:B:237:ARG:HG2	1:B:237:ARG:HH21	1.65	0.61
1:B:241:VAL:HG12	1:B:243:LEU:HD11	1.82	0.61
1:B:79:ILE:HG23	1:B:80:LEU:N	2.16	0.61
1:B:127:VAL:O	1:B:139:ASP:HB3	2.01	0.61
1:B:148:PRO:C	1:B:150:GLU:H	2.09	0.61
1:A:188:LEU:HD23	1:A:190:MET:H	1.65	0.60
1:B:242:ASP:OD1	1:B:244:ARG:HD2	2.00	0.60
1:A:188:LEU:HD11	1:A:191:VAL:CG2	2.31	0.60
1:B:101:GLU:HG3	1:B:102:ARG:HG3	1.83	0.60
1:A:64:THR:HG21	1:A:120:LYS:HE3	1.84	0.59
1:A:188:LEU:CD2	1:A:190:MET:N	2.65	0.59
1:A:203:LEU:HB2	1:A:207:VAL:HG13	1.84	0.59
1:B:82:LEU:N	1:B:82:LEU:HD22	2.17	0.59
1:A:130:ALA:HB1	1:A:160:LEU:HD13	1.85	0.59
1:A:130:ALA:CA	1:A:157:LEU:HD11	2.33	0.59
1:B:130:ALA:CA	1:B:157:LEU:HD11	2.33	0.59
1:B:200:GLN:HE21	1:B:208:ARG:HD2	1.68	0.59
1:A:173:GLY:C	1:A:177:ILE:HD12	2.28	0.58
1:A:100:ILE:HG13	1:A:109:ALA:H	1.69	0.58
1:B:93:VAL:HG23	1:B:94:ASN:N	2.19	0.58
1:A:64:THR:HG23	1:A:120:LYS:HE3	1.86	0.58
1:A:227:TYR:O	1:A:229:MET:N	2.37	0.58
1:A:102:ARG:HB3	1:A:104:PRO:HA	1.86	0.57
1:B:138:VAL:HG23	1:B:143:ARG:O	2.04	0.57
1:B:190:MET:SD	1:B:191:VAL:N	2.77	0.57
1:B:152:VAL:HG23	1:B:153:GLY:H	1.68	0.57
1:A:245:TYR:HE2	1:A:250:ALA:CB	2.07	0.57
1:A:245:TYR:HD1	1:B:72:ASN:HD21	1.53	0.57
1:B:94:ASN:O	1:B:98:GLN:HB2	2.04	0.57
1:A:152:VAL:HG23	1:A:153:GLY:H	1.69	0.57
1:B:205:ASN:O	1:B:207:VAL:N	2.38	0.57
1:A:203:LEU:N	1:A:203:LEU:HD22	2.20	0.57
1:B:223:PHE:CE1	1:B:227:TYR:HB2	2.40	0.57
1:A:98:GLN:O	1:A:101:GLU:HB3	2.05	0.56
1:A:71:THR:HG22	1:A:73:ASP:H	1.70	0.56
1:A:86:GLY:O	1:A:88:PHE:N	2.39	0.56
1:B:103:LEU:HB2	1:B:104:PRO:CD	2.34	0.56
1:A:96:ILE:O	1:A:98:GLN:N	2.39	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:205:ASN:O	1:B:205:ASN:OD1	2.23	0.56
1:A:156:LYS:HE2	1:A:156:LYS:HA	1.85	0.56
1:A:127:VAL:O	1:A:139:ASP:HB3	2.06	0.56
1:A:92:ASP:C	1:A:94:ASN:H	2.14	0.56
1:A:201:LEU:HD23	1:A:201:LEU:N	2.21	0.56
1:A:223:PHE:HE1	1:A:241:VAL:HG21	1.70	0.56
1:B:138:VAL:HG22	1:B:139:ASP:H	1.71	0.56
1:B:173:GLY:O	1:B:174:TYR:C	2.48	0.56
1:B:200:GLN:HE21	1:B:208:ARG:CG	2.19	0.56
1:A:102:ARG:H	1:A:102:ARG:HD3	1.71	0.55
1:A:92:ASP:C	1:A:94:ASN:N	2.60	0.55
1:A:188:LEU:HD21	1:A:190:MET:C	2.30	0.55
1:B:123:LEU:HD12	1:B:123:LEU:O	2.07	0.55
1:B:130:ALA:HA	1:B:157:LEU:CD1	2.36	0.55
1:A:180:VAL:O	1:A:183:ALA:HB3	2.07	0.55
1:A:226:LEU:HD12	1:A:226:LEU:H	1.71	0.55
1:B:106:ILE:N	1:B:106:ILE:HD12	2.21	0.55
1:B:199:TRP:NE1	1:B:216:ARG:HD3	2.21	0.55
1:B:215:ASP:OD2	1:B:218:GLY:HA3	2.05	0.55
1:B:97:GLN:NE2	1:B:110:SER:HA	2.21	0.55
1:A:95:ILE:O	1:A:96:ILE:O	2.25	0.55
1:A:130:ALA:HA	1:A:157:LEU:CD1	2.36	0.55
1:B:157:LEU:HD22	1:B:158:PRO:CD	2.33	0.55
1:A:100:ILE:HD11	1:A:108:GLN:HA	1.89	0.55
1:A:213:ARG:NH1	1:B:152:VAL:HG12	2.22	0.55
1:A:130:ALA:HA	1:A:157:LEU:HD11	1.89	0.54
1:A:63:VAL:HA	1:A:121:ILE:O	2.08	0.54
1:B:101:GLU:C	1:B:103:LEU:N	2.62	0.54
1:B:217:MET:O	1:B:220:LEU:HB3	2.07	0.54
1:B:79:ILE:CG2	1:B:80:LEU:N	2.70	0.54
1:A:100:ILE:O	1:A:102:ARG:N	2.40	0.54
1:A:82:LEU:N	1:A:82:LEU:HD22	2.23	0.54
1:A:98:GLN:O	1:A:101:GLU:N	2.38	0.54
1:B:61:LEU:HD12	1:B:61:LEU:N	2.22	0.54
1:A:98:GLN:O	1:A:101:GLU:CB	2.55	0.54
1:A:111:VAL:HG13	1:A:121:ILE:HG13	1.90	0.54
1:A:84:ALA:HB1	1:A:89:MET:SD	2.48	0.54
1:A:124:VAL:O	1:A:124:VAL:HG23	2.07	0.53
1:A:188:LEU:HD11	1:A:191:VAL:HG22	1.90	0.53
1:B:174:TYR:C	1:B:174:TYR:CD2	2.86	0.53
1:B:71:THR:HG22	1:B:73:ASP:N	2.22	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:ILE:HG12	1:A:99:GLN:OE1	2.08	0.53
1:B:188:LEU:HD23	1:B:190:MET:N	2.23	0.53
1:A:76:ARG:O	1:A:79:ILE:HG22	2.08	0.53
1:B:199:TRP:CD1	1:B:216:ARG:HD3	2.43	0.53
1:B:185:LYS:O	1:B:185:LYS:HG2	2.09	0.53
1:B:139:ASP:O	1:B:142:GLY:N	2.41	0.53
1:B:61:LEU:CD1	1:B:76:ARG:NH1	2.71	0.52
1:B:112:ARG:HG2	1:B:120:LYS:HB3	1.91	0.52
1:A:208:ARG:HG2	1:A:209:LEU:N	2.25	0.52
1:B:156:LYS:HE2	1:B:156:LYS:HA	1.91	0.52
1:A:208:ARG:CG	1:A:209:LEU:N	2.72	0.52
1:B:238:VAL:HG12	1:B:239:SER:N	2.15	0.52
1:B:245:TYR:HB2	1:B:248:GLY:O	2.10	0.52
1:A:95:ILE:O	1:A:96:ILE:C	2.52	0.52
1:B:79:ILE:HD13	1:B:96:ILE:CG2	2.39	0.52
1:B:225:GLU:O	1:B:228:PRO:HD2	2.10	0.52
1:A:173:GLY:O	1:A:177:ILE:HD12	2.10	0.51
1:A:174:TYR:C	1:A:174:TYR:CD2	2.88	0.51
1:A:175:ARG:HG3	1:A:175:ARG:HH21	1.76	0.51
1:A:174:TYR:CE2	1:A:178:ASN:HB2	2.45	0.51
1:B:200:GLN:HE22	1:B:208:ARG:HD2	1.70	0.51
1:A:174:TYR:HE2	1:A:178:ASN:CG	2.19	0.51
1:B:71:THR:O	1:B:74:ASP:N	2.43	0.51
1:B:188:LEU:HD21	1:B:190:MET:O	2.11	0.51
1:A:188:LEU:CD2	1:A:190:MET:H	2.23	0.51
1:B:60:LYS:C	1:B:61:LEU:HD12	2.36	0.51
1:B:111:VAL:HG13	1:B:121:ILE:CD1	2.41	0.51
1:B:242:ASP:C	1:B:243:LEU:HD12	2.36	0.51
1:A:96:ILE:O	1:A:97:GLN:C	2.53	0.50
1:A:160:LEU:HD21	1:A:174:TYR:HD1	1.75	0.50
1:B:127:VAL:O	1:B:139:ASP:CB	2.58	0.50
1:A:100:ILE:CG1	1:A:108:GLN:HA	2.41	0.50
1:A:102:ARG:HB3	1:A:104:PRO:CA	2.41	0.50
1:A:119:LEU:O	1:A:121:ILE:HD12	2.11	0.50
1:B:160:LEU:HD11	1:B:174:TYR:CD1	2.47	0.50
1:B:209:LEU:N	1:B:209:LEU:CD1	2.74	0.50
1:A:101:GLU:O	1:A:102:ARG:O	2.29	0.50
1:A:60:LYS:C	1:A:61:LEU:HD12	2.37	0.50
1:A:79:ILE:CG2	1:A:80:LEU:H	2.24	0.49
1:B:114:GLN:HG3	1:B:118:GLU:HG3	1.94	0.49
1:B:220:LEU:O	1:B:221:GLN:C	2.55	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:188:LEU:C	1:A:188:LEU:CD2	2.85	0.49
1:A:227:TYR:CE1	1:A:230:LEU:HD11	2.47	0.49
1:B:101:GLU:CD	1:B:108:GLN:HA	2.38	0.49
1:A:81:SER:HB2	1:A:98:GLN:NE2	2.28	0.49
1:A:94:ASN:O	1:A:97:GLN:HB3	2.13	0.49
1:B:200:GLN:HE21	1:B:208:ARG:CD	2.26	0.49
1:A:203:LEU:N	1:A:203:LEU:CD2	2.76	0.49
1:B:76:ARG:O	1:B:79:ILE:HG22	2.13	0.48
1:A:148:PRO:C	1:A:150:GLU:N	2.69	0.48
1:B:207:VAL:HG22	1:B:208:ARG:N	2.22	0.48
1:A:79:ILE:HD13	1:A:95:ILE:HG23	1.96	0.48
1:B:243:LEU:HD12	1:B:243:LEU:N	2.28	0.48
1:A:174:TYR:CD2	1:A:178:ASN:HB2	2.49	0.48
1:A:127:VAL:O	1:A:127:VAL:HG13	2.12	0.48
1:A:243:LEU:C	1:A:245:TYR:H	2.21	0.48
1:B:196:ARG:HB3	1:B:213:ARG:NH1	2.29	0.48
1:A:173:GLY:O	1:A:174:TYR:C	2.57	0.48
1:A:257:PHE:CD1	1:A:257:PHE:C	2.92	0.48
1:B:115:TRP:O	1:B:116:PRO:C	2.54	0.48
1:B:221:GLN:O	1:B:224:ILE:CG1	2.62	0.48
1:B:241:VAL:HG12	1:B:243:LEU:HD12	1.95	0.48
1:B:237:ARG:HG2	1:B:237:ARG:NH2	2.28	0.47
1:B:104:PRO:HB2	1:B:105:TRP:CZ3	2.49	0.47
1:B:114:GLN:HB2	1:B:118:GLU:HG2	1.94	0.47
1:B:171:LEU:O	1:B:172:GLU:C	2.57	0.47
1:B:251:ILE:HG22	1:B:252:GLY:N	2.29	0.47
1:A:79:ILE:HD11	1:A:95:ILE:CD1	2.43	0.47
1:A:227:TYR:C	1:A:229:MET:N	2.69	0.47
1:B:174:TYR:CE2	1:B:178:ASN:HB2	2.50	0.47
1:A:72:ASN:O	1:A:75:ILE:HB	2.15	0.47
1:A:127:VAL:O	1:A:139:ASP:CB	2.62	0.47
1:A:75:ILE:O	1:A:78:ALA:N	2.33	0.47
1:B:101:GLU:CG	1:B:102:ARG:H	2.12	0.47
1:B:159:LEU:C	1:B:160:LEU:HD12	2.39	0.47
1:B:254:ALA:HA	1:B:255:PRO:HD3	1.58	0.47
1:A:200:GLN:C	1:A:201:LEU:HD23	2.40	0.47
1:B:103:LEU:HB2	1:B:104:PRO:HD3	1.96	0.47
1:A:92:ASP:HB2	1:A:111:VAL:HB	1.97	0.46
1:A:148:PRO:O	1:A:150:GLU:N	2.42	0.46
1:B:245:TYR:HE2	1:B:250:ALA:CB	2.22	0.46
1:B:257:PHE:H	1:B:257:PHE:HD2	1.62	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:ASN:C	1:A:207:VAL:H	2.24	0.46
1:B:241:VAL:O	1:B:243:LEU:CD1	2.63	0.46
1:A:160:LEU:HA	1:A:191:VAL:O	2.14	0.46
1:A:222:ARG:O	1:A:223:PHE:C	2.57	0.46
1:A:114:GLN:C	1:A:116:PRO:HD2	2.40	0.46
1:B:177:ILE:O	1:B:178:ASN:C	2.56	0.46
1:B:181:LEU:HD12	1:B:188:LEU:HD12	1.98	0.46
1:B:188:LEU:C	1:B:188:LEU:CD2	2.87	0.46
1:B:227:TYR:HB3	1:B:228:PRO:HD3	1.97	0.46
1:A:75:ILE:O	1:A:76:ARG:C	2.59	0.46
1:B:101:GLU:CD	1:B:109:ALA:H	2.23	0.46
1:B:132:TRP:O	1:B:160:LEU:O	2.34	0.46
1:A:160:LEU:HD11	1:A:174:TYR:CD1	2.51	0.46
1:A:226:LEU:HD12	1:A:226:LEU:N	2.30	0.46
1:A:254:ALA:HA	1:A:255:PRO:HD3	1.66	0.46
1:B:226:LEU:N	1:B:226:LEU:CD1	2.78	0.46
1:A:91:GLN:HE21	1:A:91:GLN:HB2	1.56	0.45
1:B:152:VAL:HG23	1:B:153:GLY:N	2.30	0.45
1:A:199:TRP:H	1:A:199:TRP:CD1	2.32	0.45
1:B:240:TYR:CD1	1:B:240:TYR:C	2.95	0.45
1:A:201:LEU:HD21	1:A:211:LEU:CD2	2.34	0.45
1:B:174:TYR:CD2	1:B:178:ASN:HB2	2.52	0.45
1:A:121:ILE:HD12	1:A:121:ILE:N	2.32	0.45
1:A:178:ASN:O	1:A:179:LYS:C	2.60	0.45
1:A:129:PHE:C	1:A:157:LEU:HD21	2.42	0.45
1:A:73:ASP:C	1:A:75:ILE:N	2.70	0.44
1:A:92:ASP:HA	1:A:95:ILE:HG12	1.99	0.44
1:A:103:LEU:HD12	1:A:106:ILE:HD13	1.99	0.44
1:A:139:ASP:O	1:A:140:GLU:C	2.59	0.44
1:B:63:VAL:HG13	1:B:121:ILE:HG22	1.98	0.44
1:A:227:TYR:H	1:A:228:PRO:CD	2.29	0.44
1:B:71:THR:O	1:B:72:ASN:C	2.59	0.44
1:B:194:SER:O	1:B:195:ALA:C	2.59	0.44
1:A:152:VAL:HG23	1:A:153:GLY:N	2.33	0.44
1:B:148:PRO:C	1:B:150:GLU:N	2.74	0.44
1:B:129:PHE:O	1:B:157:LEU:HD21	2.17	0.44
1:B:130:ALA:C	1:B:157:LEU:HD11	2.42	0.44
1:B:209:LEU:HD21	1:B:223:PHE:CE2	2.53	0.44
1:A:81:SER:C	1:A:82:LEU:HD22	2.42	0.44
1:A:114:GLN:HB3	1:A:116:PRO:HD2	1.99	0.44
1:A:223:PHE:O	1:A:224:ILE:C	2.61	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:ASP:HA	1:A:95:ILE:CG1	2.48	0.43
1:A:102:ARG:CA	1:A:104:PRO:HA	2.45	0.43
1:B:157:LEU:HD13	1:B:157:LEU:C	2.43	0.43
1:B:175:ARG:HH21	1:B:175:ARG:HG3	1.83	0.43
1:A:171:LEU:O	1:A:172:GLU:C	2.60	0.43
1:B:81:SER:C	1:B:82:LEU:HD22	2.42	0.43
1:A:242:ASP:C	1:A:243:LEU:HD12	2.42	0.43
1:B:59:SER:HB3	1:B:117:ASP:O	2.18	0.43
1:A:139:ASP:O	1:A:142:GLY:N	2.50	0.43
1:B:221:GLN:O	1:B:224:ILE:HG12	2.19	0.43
1:B:226:LEU:CD1	1:B:226:LEU:H	2.32	0.43
1:A:82:LEU:HD21	1:A:98:GLN:NE2	2.33	0.43
1:A:162:GLY:HA2	1:A:193:MET:O	2.18	0.43
1:B:121:ILE:HG22	1:B:121:ILE:O	2.18	0.43
1:A:223:PHE:CE1	1:A:241:VAL:HG21	2.52	0.43
1:B:114:GLN:CG	1:B:118:GLU:HG3	2.49	0.43
1:B:174:TYR:HD2	1:B:174:TYR:O	2.01	0.43
1:B:188:LEU:HD21	1:B:191:VAL:HG23	2.01	0.43
1:A:91:GLN:CG	1:A:93:VAL:HG23	2.49	0.42
1:B:71:THR:O	1:B:74:ASP:HB2	2.18	0.42
1:B:88:PHE:CD2	1:B:88:PHE:C	2.88	0.42
1:B:106:ILE:N	1:B:106:ILE:CD1	2.82	0.42
1:A:196:ARG:NE	1:B:68:HIS:CE1	2.87	0.42
1:A:201:LEU:HD23	1:A:209:LEU:O	2.19	0.42
1:B:111:VAL:HG13	1:B:121:ILE:HD11	2.00	0.42
1:A:132:TRP:HB3	1:A:136:HIS:HB3	2.01	0.42
1:A:156:LYS:HE2	1:A:156:LYS:CA	2.50	0.42
1:A:207:VAL:O	1:A:207:VAL:HG22	2.17	0.42
1:B:132:TRP:HB3	1:B:136:HIS:HB3	2.00	0.42
1:B:243:LEU:C	1:B:245:TYR:H	2.26	0.42
1:A:73:ASP:C	1:A:75:ILE:H	2.27	0.42
1:B:103:LEU:CB	1:B:104:PRO:CD	2.98	0.42
1:B:105:TRP:O	1:B:126:TYR:HD1	2.02	0.42
1:B:245:TYR:CE2	1:B:250:ALA:HB2	2.39	0.42
1:A:203:LEU:O	1:A:204:ASP:C	2.62	0.42
1:A:95:ILE:HG12	1:A:95:ILE:H	1.66	0.42
1:B:199:TRP:CE2	1:B:216:ARG:HD3	2.55	0.42
1:A:71:THR:O	1:A:72:ASN:C	2.63	0.41
1:A:97:GLN:CA	1:A:100:ILE:HG22	2.50	0.41
1:A:102:ARG:CB	1:A:104:PRO:HA	2.50	0.41
1:A:161:TYR:HB2	1:A:192:ALA:CB	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:93:VAL:O	1:B:95:ILE:N	2.53	0.41
1:B:238:VAL:HG13	1:B:251:ILE:HG23	2.03	0.41
1:A:129:PHE:HB3	1:A:139:ASP:HA	2.02	0.41
1:A:199:TRP:CD1	1:A:199:TRP:N	2.89	0.41
1:A:59:SER:HB3	1:A:88:PHE:CE1	2.56	0.41
1:A:130:ALA:C	1:A:157:LEU:HD11	2.44	0.41
1:A:239:SER:O	1:A:240:TYR:CB	2.65	0.41
1:B:203:LEU:HD22	1:B:207:VAL:CG1	2.50	0.41
1:A:177:ILE:O	1:A:178:ASN:C	2.64	0.41
1:A:195:ALA:C	1:A:197:HIS:H	2.28	0.41
1:B:138:VAL:CG2	1:B:139:ASP:N	2.83	0.41
1:B:220:LEU:O	1:B:222:ARG:N	2.54	0.41
1:A:69:TYR:HD2	1:A:125:GLU:OE2	2.04	0.41
1:A:91:GLN:O	1:A:94:ASN:HB3	2.21	0.41
1:A:138:VAL:HG22	1:A:139:ASP:N	2.35	0.41
1:B:173:GLY:O	1:B:176:ALA:N	2.54	0.41
1:B:216:ARG:O	1:B:217:MET:C	2.63	0.41
1:A:111:VAL:CG1	1:A:119:LEU:HD22	2.51	0.41
1:B:207:VAL:HG13	1:B:208:ARG:N	2.34	0.41
1:A:64:THR:OG1	1:A:122:HIS:CD2	2.74	0.40
1:A:168:GLN:O	1:A:169:ASP:C	2.64	0.40
1:A:240:TYR:CZ	1:A:252:GLY:HA3	2.56	0.40
1:B:226:LEU:HD12	1:B:226:LEU:H	1.83	0.40
1:A:79:ILE:HG22	1:A:80:LEU:H	1.87	0.40
1:B:111:VAL:HG13	1:B:121:ILE:HD12	2.03	0.40
1:B:240:TYR:CE2	1:B:252:GLY:HA3	2.57	0.40
1:A:242:ASP:OD1	1:A:244:ARG:HD2	2.22	0.40
1:B:188:LEU:HD21	1:B:190:MET:C	2.46	0.40
1:A:79:ILE:C	1:A:81:SER:N	2.79	0.40
1:A:132:TRP:O	1:A:160:LEU:O	2.39	0.40
1:B:147:VAL:HA	1:B:148:PRO:HD3	1.96	0.40
1:B:196:ARG:HD3	1:B:213:ARG:HH12	1.87	0.40
1:A:81:SER:HB2	1:A:98:GLN:HE21	1.85	0.40
1:B:69:TYR:HB2	1:B:105:TRP:NE1	2.37	0.40
1:B:150:GLU:HB3	1:B:151:ARG:H	1.62	0.40
1:B:160:LEU:HD12	1:B:160:LEU:N	2.37	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	198/255 (78%)	129 (65%)	50 (25%)	19 (10%)	0	3
1	B	198/255 (78%)	134 (68%)	43 (22%)	21 (11%)	0	2
All	All	396/510 (78%)	263 (66%)	93 (24%)	40 (10%)	0	3

All (40) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	87	THR
1	A	96	ILE
1	A	97	GLN
1	A	117	ASP
1	A	227	TYR
1	B	88	PHE
1	B	89	MET
1	B	100	ILE
1	B	102	ARG
1	B	104	PRO
1	B	197	HIS
1	B	206	ASP
1	A	102	ARG
1	A	208	ARG
1	B	67	ARG
1	B	101	GLU
1	B	216	ARG
1	A	98	GLN
1	A	101	GLU
1	A	149	SER
1	A	209	LEU
1	A	225	GLU
1	A	228	PRO
1	A	240	TYR
1	A	244	ARG

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Mol	Chain	Res	Type
1	B	94	ASN
1	B	117	ASP
1	B	185	LYS
1	B	149	SER
1	B	150	GLU
1	A	103	LEU
1	B	214	ASP
1	A	147	VAL
1	A	196	ARG
1	B	83	GLY
1	B	239	SER
1	A	255	PRO
1	B	116	PRO
1	B	147	VAL
1	B	238	VAL

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	176/220 (80%)	156 (89%)	20 (11%)	5	21
1	B	176/220 (80%)	160 (91%)	16 (9%)	9	30
All	All	352/440 (80%)	316 (90%)	36 (10%)	7	25

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	VAL
1	A	90	THR
1	A	91	GLN
1	A	93	VAL
1	A	100	ILE
1	A	107	GLN
1	A	119	LEU
1	A	123	LEU

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Mol	Chain	Res	Type
1	A	136	HIS
1	A	137	MET
1	A	149	SER
1	A	156	LYS
1	A	190	MET
1	A	201	LEU
1	A	207	VAL
1	A	209	LEU
1	A	211	LEU
1	A	229	MET
1	A	251	ILE
1	A	258	ILE
1	B	63	VAL
1	B	88	PHE
1	B	100	ILE
1	B	107	GLN
1	B	136	HIS
1	B	137	MET
1	B	149	SER
1	B	156	LYS
1	B	190	MET
1	B	201	LEU
1	B	204	ASP
1	B	209	LEU
1	B	236	LYS
1	B	243	LEU
1	B	251	ILE
1	B	257	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	68	HIS
1	A	98	GLN
1	A	99	GLN
1	A	114	GLN
1	A	122	HIS
1	A	141	GLN
1	A	187	GLN
1	A	197	HIS
1	A	221	GLN
1	B	72	ASN

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Mol	Chain	Res	Type
1	B	97	GLN
1	B	114	GLN
1	B	200	GLN
1	B	221	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	202/255 (79%)	0.12	5 (2%) 58 43	91, 149, 192, 199	0
1	B	202/255 (79%)	0.10	3 (1%) 72 57	98, 154, 192, 199	0
All	All	404/510 (79%)	0.11	8 (1%) 65 49	91, 152, 192, 199	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	87	THR	2.8
1	B	200	GLN	2.5
1	A	134	ASP	2.3
1	B	89	MET	2.3
1	A	99	GLN	2.3
1	A	223	PHE	2.1
1	B	195	ALA	2.1
1	A	255	PRO	2.1

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.