



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:16 PM UTC

PDB ID : 2ZVZ / pdb_00002zvz
Title : Structure of the periplasmic domain of MotB from Salmonella (crystal form III)
Authors : Imada, K.; Kojima, S.; Namba, K.; Homma, S.
Deposited on : 2008-11-26
Resolution : 2.40 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

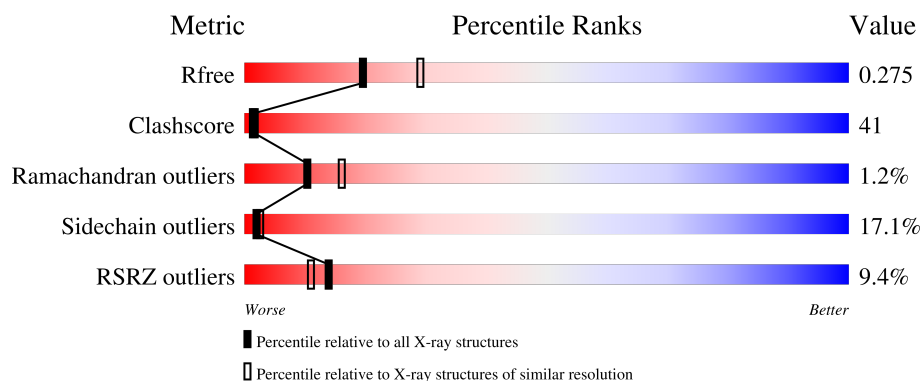
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	183	5% 38% 39% 14% 7%
1	B	183	11% 37% 37% 13% 12%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2844 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 Chemotaxis protein motB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	170	Total	C	N	O	S	0	0	0
			1339	831	255	248	5			
1	B	161	Total	C	N	O	S	0	0	0
			1269	787	243	235	4			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	277	HIS	-	expression tag	UNP P55892
A	278	HIS	-	expression tag	UNP P55892
A	279	HIS	-	expression tag	UNP P55892
A	280	HIS	-	expression tag	UNP P55892
A	281	HIS	-	expression tag	UNP P55892
B	277	HIS	-	expression tag	UNP P55892
B	278	HIS	-	expression tag	UNP P55892
B	279	HIS	-	expression tag	UNP P55892
B	280	HIS	-	expression tag	UNP P55892
B	281	HIS	-	expression tag	UNP P55892

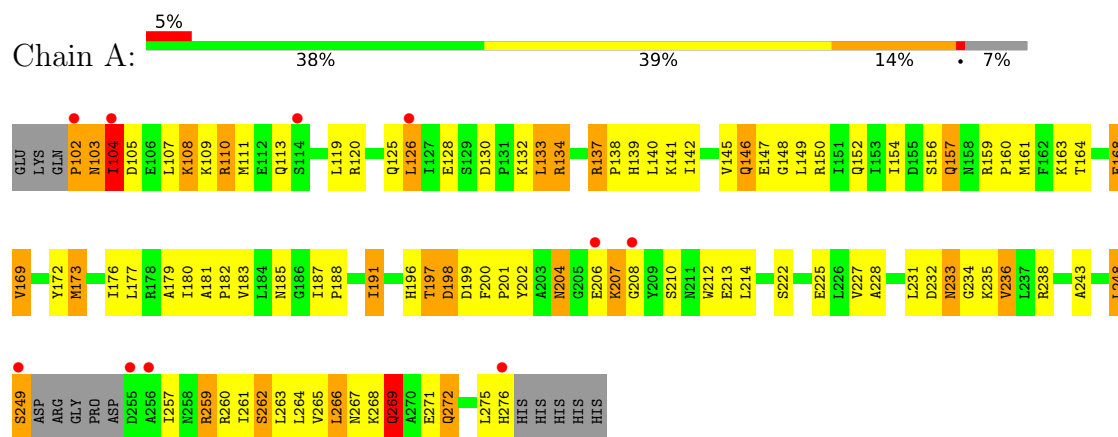
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	146	Total	O	0	0
			146	146		
2	B	90	Total	O	0	0
			90	90		

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: Chemotaxis protein motB



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	45.20Å 46.24Å 173.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.68 – 2.40 44.68 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.2 (44.68-2.40) 99.3 (44.68-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.21 (at 2.39Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.245 , 0.275 0.246 , 0.275	Depositor DCC
R_{free} test set	724 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	42.0	Xtriage
Anisotropy	0.609	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 59.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.042 for k,h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	2844	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.52% 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	0.61	0/1355	1.28	20/1822 (1.1%)
1	B	0.51	0/1282	1.17	13/1721 (0.8%)
All	All	0.56	0/2637	1.23	33/3543 (0.9%)

There are no bond length outliers.

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	248	LEU	N-CA-C	12.07	128.99	108.23
1	A	108	LYS	N-CA-C	-9.74	100.81	112.89
1	A	257	ILE	N-CA-C	-8.67	100.26	111.09
1	B	133	LEU	N-CA-C	-7.66	102.57	112.23
1	A	173	MET	N-CA-C	-7.12	103.45	111.14
1	B	191	ILE	N-CA-C	6.92	119.50	108.85
1	A	204	ASN	N-CA-C	-6.66	104.41	112.54
1	B	105	ASP	N-CA-C	-6.55	102.78	111.24
1	A	111	MET	N-CA-C	-6.48	104.30	111.82
1	A	248	LEU	CA-CB-CG	-6.31	94.21	116.30
1	A	128	GLU	N-CA-C	6.12	117.95	111.28
1	B	243	ALA	CB-CA-C	-6.06	109.60	116.63
1	B	266	LEU	N-CA-C	6.04	119.36	109.76
1	B	169	VAL	N-CA-C	6.03	117.82	109.80
1	A	154	ILE	N-CA-C	5.92	117.88	108.89
1	A	104	ILE	N-CA-C	-5.86	107.57	113.20
1	A	266	LEU	N-CA-C	5.84	119.04	109.76
1	A	130	ASP	CA-C-N	-5.81	113.86	120.89
1	A	130	ASP	C-N-CA	-5.81	113.86	120.89
1	A	156	SER	N-CA-C	-5.70	100.84	109.85
1	B	225	GLU	N-CA-C	-5.64	104.75	111.69
1	A	110	ARG	N-CA-C	-5.61	107.08	114.31
1	B	156	SER	N-CA-C	-5.60	101.62	109.69
1	A	197	THR	N-CA-C	5.50	118.36	109.40
1	B	268	LYS	N-CA-C	5.46	117.99	111.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	267	ASN	N-CA-C	-5.39	103.57	110.53
1	B	187	ILE	CA-C-N	5.30	125.62	119.47
1	B	187	ILE	C-N-CA	5.30	125.62	119.47
1	A	191	ILE	N-CA-C	5.23	117.39	108.86
1	B	173	MET	N-CA-C	-5.18	105.63	111.28
1	A	183	VAL	N-CA-C	-5.16	104.52	111.44
1	A	185	ASN	N-CA-C	-5.14	106.96	113.18
1	A	126	LEU	N-CA-C	-5.09	105.62	111.07

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1339	0	1380	89	0
1	B	1269	0	1307	130	0
2	A	146	0	0	30	0
2	B	90	0	0	8	0
All	All	2844	0	2687	217	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 41.

All (217) 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:182:PRO:HD2	2:A:319:HOH:O	1.48	1.08
1:A:201:PRO:HD2	1:A:204:ASN:HD21	1.03	1.07
1:B:122:ASP:O	1:B:125:GLN:HG2	1.56	1.04
1:B:125:GLN:HG3	1:B:126:LEU:H	1.17	1.02
1:A:201:PRO:HD2	1:A:204:ASN:ND2	1.78	0.98
1:B:275:LEU:HB2	2:B:72:HOH:O	1.63	0.97
1:B:125:GLN:HG3	1:B:126:LEU:N	1.73	0.97
1:A:197:THR:HG21	1:A:214:LEU:HD23	1.46	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:204:ASN:HB2	2:A:318:HOH:O	1.69	0.93
1:A:199:ASP:OD1	1:A:259:ARG:HD3	1.70	0.90
1:A:200:PHE:HB3	1:A:204:ASN:HD22	1.40	0.87
1:A:272:GLN:HA	1:A:272:GLN:HE21	1.36	0.87
1:A:168:GLU:HA	1:A:168:GLU:OE1	1.73	0.86
1:A:248:LEU:O	1:A:249:SER:HB2	1.80	0.82
1:B:127:ILE:HG21	1:B:137:ARG:HG3	1.61	0.81
1:A:207:LYS:HG3	1:A:208:GLY:N	1.95	0.81
1:A:119:LEU:HD22	1:A:187:ILE:HD11	1.64	0.79
1:A:268:LYS:O	1:A:272:GLN:HG2	1.82	0.79
1:B:198:ASP:HB2	1:B:259:ARG:HG3	1.65	0.78
1:A:110:ARG:HG3	2:A:340:HOH:O	1.84	0.78
1:A:200:PHE:HB3	1:A:204:ASN:ND2	1.99	0.77
1:A:181:ALA:N	2:A:319:HOH:O	2.18	0.77
1:B:108:LYS:HA	1:B:111:MET:HE2	1.66	0.76
1:B:112:GLU:OE2	1:B:267:ASN:HB2	1.86	0.75
1:B:105:ASP:O	1:B:109:LYS:HG3	1.89	0.73
1:A:169:VAL:HG23	1:A:173:MET:HB3	1.69	0.73
1:B:233:ASN:ND2	1:B:233:ASN:N	2.36	0.73
1:B:272:GLN:NE2	1:B:277:HIS:HB2	2.05	0.71
1:A:227:VAL:HA	2:A:347:HOH:O	1.91	0.69
1:A:157:GLN:HG3	2:A:308:HOH:O	1.92	0.69
1:B:197:THR:OG1	1:B:211:ASN:HB3	1.92	0.69
1:B:272:GLN:HE22	1:B:277:HIS:CG	2.12	0.68
1:B:271:GLU:O	1:B:275:LEU:HG	1.92	0.68
1:B:161:MET:HB3	1:B:173:MET:CG	2.23	0.68
1:B:108:LYS:HG2	1:B:111:MET:CE	2.25	0.67
1:B:127:ILE:CG2	1:B:137:ARG:HG3	2.25	0.67
1:B:272:GLN:NE2	1:B:272:GLN:HA	2.09	0.66
1:B:104:ILE:CA	1:B:107:LEU:HB3	2.25	0.66
1:B:272:GLN:HE21	1:B:277:HIS:HB2	1.60	0.66
1:A:262:SER:HB3	2:A:339:HOH:O	1.96	0.66
1:A:201:PRO:CD	1:A:204:ASN:HD21	1.94	0.65
1:A:179:ALA:C	2:A:319:HOH:O	2.38	0.65
1:B:139:HIS:HE1	1:B:156:SER:HB3	1.60	0.65
1:A:210:SER:C	2:A:323:HOH:O	2.40	0.65
1:B:104:ILE:HA	1:B:107:LEU:HB3	1.79	0.65
1:B:257:ILE:HD13	1:B:257:ILE:H	1.62	0.65
1:B:269:GLN:NE2	2:B:294:HOH:O	2.30	0.65
1:A:275:LEU:O	1:A:276:HIS:C	2.39	0.65
1:B:140:LEU:O	1:B:141:LYS:HG2	1.96	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:136:LEU:C	1:B:138:PRO:HD2	2.23	0.64
1:A:119:LEU:HD22	1:A:187:ILE:CD1	2.28	0.64
1:B:197:THR:OG1	1:B:243:ALA:HA	1.99	0.63
1:B:223:ARG:HG3	1:B:236:VAL:HG21	1.81	0.63
1:A:146:GLN:HB2	2:A:331:HOH:O	1.98	0.62
1:B:143:ASP:OD2	1:B:143:ASP:N	2.31	0.62
1:B:161:MET:HB3	1:B:173:MET:HG2	1.81	0.62
1:A:259:ARG:HG2	2:A:35:HOH:O	1.98	0.62
1:A:126:LEU:HD21	1:A:179:ALA:O	1.99	0.62
1:B:104:ILE:N	1:B:107:LEU:HB2	2.15	0.62
1:B:161:MET:SD	1:B:173:MET:HG2	2.40	0.62
1:A:197:THR:CG2	1:A:214:LEU:HD23	2.27	0.62
1:B:232:ASP:CG	1:B:233:ASN:N	2.57	0.61
1:A:204:ASN:HB3	1:A:207:LYS:HE2	1.82	0.61
1:B:123:LEU:HD13	1:B:142:ILE:HD13	1.81	0.61
1:B:233:ASN:ND2	1:B:233:ASN:H	1.97	0.61
1:B:233:ASN:N	1:B:233:ASN:HD22	1.97	0.61
1:A:157:GLN:HB2	2:A:290:HOH:O	1.99	0.61
1:B:108:LYS:HG2	1:B:111:MET:HE2	1.81	0.61
2:A:328:HOH:O	1:B:224:ARG:HG2	2.00	0.61
1:A:201:PRO:CD	1:A:204:ASN:ND2	2.57	0.61
1:A:212:TRP:N	2:A:323:HOH:O	2.34	0.61
1:B:137:ARG:N	1:B:138:PRO:HD2	2.16	0.61
1:B:275:LEU:O	1:B:276:HIS:HB2	1.99	0.60
1:B:174:ARG:HG3	1:B:225:GLU:OE1	2.00	0.60
1:B:198:ASP:HB2	1:B:259:ARG:CG	2.31	0.60
1:A:248:LEU:O	1:A:249:SER:CB	2.50	0.60
1:A:231:LEU:O	2:A:347:HOH:O	2.15	0.60
1:B:126:LEU:HA	1:B:129:SER:OG	2.03	0.59
1:B:106:GLU:HA	1:B:109:LYS:HB2	1.83	0.59
1:A:105:ASP:HA	2:A:343:HOH:O	2.04	0.58
1:A:272:GLN:HE21	1:A:272:GLN:CA	2.11	0.58
1:B:130:ASP:HB3	1:B:133:LEU:HG	1.86	0.58
1:B:232:ASP:CG	1:B:233:ASN:H	2.12	0.58
1:A:147:GLU:O	1:A:266:LEU:HA	2.04	0.57
1:A:200:PHE:HB2	2:A:318:HOH:O	2.04	0.57
1:B:231:LEU:HD21	1:B:235:LYS:HB3	1.87	0.57
1:B:108:LYS:HA	1:B:111:MET:CE	2.35	0.57
1:A:168:GLU:OE1	1:A:168:GLU:CA	2.50	0.56
1:B:237:LEU:O	1:B:237:LEU:HD23	2.06	0.56
1:A:238:ARG:HA	1:B:212:TRP:HZ3	1.72	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:228:ALA:HB3	2:A:336:HOH:O	2.06	0.55
1:B:123:LEU:HD12	1:B:149:LEU:HD21	1.89	0.55
1:A:108:LYS:HB2	2:A:343:HOH:O	2.07	0.54
1:B:272:GLN:NE2	1:B:277:HIS:CG	2.74	0.54
1:B:188:PRO:HD2	2:B:98:HOH:O	2.06	0.54
1:B:223:ARG:NH2	1:B:224:ARG:HG3	2.23	0.54
1:A:147:GLU:HB2	1:A:264:LEU:HD21	1.89	0.54
1:A:269:GLN:NE2	2:A:333:HOH:O	2.40	0.54
1:B:119:LEU:HG	1:B:149:LEU:HD22	1.90	0.53
1:B:139:HIS:HE1	1:B:156:SER:CB	2.20	0.53
1:B:139:HIS:CE1	1:B:156:SER:HB3	2.42	0.53
1:B:161:MET:HB3	1:B:173:MET:HG3	1.90	0.53
1:B:255:ASP:C	1:B:257:ILE:HD13	2.34	0.53
1:B:140:LEU:C	1:B:141:LYS:HG2	2.33	0.53
1:A:272:GLN:HA	1:A:272:GLN:NE2	2.14	0.53
1:B:198:ASP:O	1:B:199:ASP:C	2.51	0.53
1:B:119:LEU:HG	1:B:149:LEU:CD2	2.38	0.53
1:A:102:PRO:O	1:A:103:ASN:HB2	2.08	0.53
1:B:145:VAL:HG22	1:B:148:GLY:O	2.09	0.53
1:B:162:PHE:HE2	1:B:261:ILE:HD11	1.74	0.52
1:A:140:LEU:HD21	1:A:176:ILE:HG21	1.90	0.52
1:B:272:GLN:NE2	1:B:277:HIS:CB	2.72	0.52
1:A:191:ILE:HD13	1:A:263:LEU:HD21	1.92	0.52
1:B:154:ILE:HG22	1:B:155:ASP:N	2.25	0.51
1:B:161:MET:CB	1:B:173:MET:HG2	2.41	0.51
1:A:137:ARG:HD2	2:A:329:HOH:O	2.11	0.51
1:A:148:GLY:HA3	1:A:265:VAL:O	2.11	0.51
1:A:233:ASN:OD1	1:A:233:ASN:N	2.31	0.51
1:B:162:PHE:CE2	1:B:261:ILE:HD11	2.46	0.51
1:A:200:PHE:CB	2:A:318:HOH:O	2.60	0.50
1:B:142:ILE:HG22	1:B:149:LEU:HD11	1.94	0.50
1:B:160:PRO:HB3	2:B:323:HOH:O	2.11	0.50
1:B:136:LEU:O	1:B:137:ARG:C	2.55	0.50
1:A:213:GLU:OE1	2:A:323:HOH:O	2.19	0.50
1:A:232:ASP:O	1:A:235:LYS:HG2	2.12	0.49
1:A:103:ASN:OD1	1:A:104:ILE:N	2.44	0.49
1:B:191:ILE:CG2	1:B:263:LEU:HD23	2.42	0.49
1:B:191:ILE:HD13	1:B:263:LEU:HD21	1.94	0.49
1:A:137:ARG:HG2	1:A:138:PRO:HD3	1.94	0.49
1:B:127:ILE:HD13	1:B:140:LEU:HD13	1.93	0.49
1:A:236:VAL:HG22	2:A:71:HOH:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:276:HIS:C	1:A:276:HIS:ND1	2.71	0.49
1:B:191:ILE:HG22	1:B:192:SER:N	2.28	0.49
1:B:136:LEU:HD21	1:B:159:ARG:HH11	1.77	0.49
1:A:202:TYR:HA	2:A:352:HOH:O	2.12	0.48
1:B:137:ARG:N	1:B:138:PRO:CD	2.75	0.48
1:B:255:ASP:C	1:B:257:ILE:H	2.21	0.48
1:A:134:ARG:HG3	1:A:137:ARG:NH2	2.28	0.48
1:B:260:ARG:HH11	1:B:260:ARG:HG2	1.77	0.48
1:A:152:GLN:HG2	1:A:262:SER:OG	2.14	0.48
1:B:191:ILE:HG21	1:B:263:LEU:CD2	2.44	0.48
1:A:248:LEU:HD23	1:A:248:LEU:HA	1.35	0.48
1:B:150:ARG:NH1	1:B:152:GLN:OE1	2.47	0.48
1:A:260:ARG:C	1:A:261:ILE:HD12	2.38	0.48
1:B:196:HIS:CE1	1:B:242:MET:HG3	2.49	0.48
1:A:132:LYS:C	1:A:133:LEU:HD23	2.39	0.47
1:A:225:GLU:HG3	2:A:336:HOH:O	2.14	0.47
1:B:184:LEU:C	1:B:235:LYS:HZ1	2.17	0.47
1:B:255:ASP:CA	1:B:257:ILE:HD13	2.45	0.47
1:B:275:LEU:O	1:B:276:HIS:CB	2.63	0.47
1:B:190:ARG:HG2	1:B:235:LYS:O	2.15	0.47
1:A:139:HIS:ND1	1:A:161:MET:HE3	2.30	0.47
1:A:187:ILE:HB	1:A:188:PRO:HD2	1.97	0.46
1:B:197:THR:OG1	1:B:243:ALA:CA	2.63	0.46
1:B:142:ILE:C	1:B:143:ASP:OD2	2.58	0.46
1:A:212:TRP:CE2	1:A:243:ALA:HB3	2.51	0.46
1:A:213:GLU:HB3	2:A:334:HOH:O	2.15	0.46
1:B:122:ASP:O	1:B:125:GLN:CG	2.47	0.45
1:A:137:ARG:C	1:A:139:HIS:H	2.23	0.45
1:A:234:GLY:HA2	2:A:320:HOH:O	2.16	0.45
1:A:177:LEU:HD13	1:A:222:SER:HA	1.99	0.45
1:B:104:ILE:HG23	1:B:105:ASP:N	2.31	0.45
1:B:268:LYS:O	1:B:272:GLN:HG2	2.16	0.45
1:A:180:ILE:N	2:A:319:HOH:O	2.49	0.45
1:B:181:ALA:HB3	1:B:182:PRO:HD3	1.99	0.45
1:B:104:ILE:CA	1:B:107:LEU:CB	2.92	0.45
1:B:124:ASP:O	1:B:125:GLN:C	2.59	0.45
1:B:104:ILE:N	1:B:107:LEU:CB	2.80	0.44
1:B:179:ALA:O	1:B:182:PRO:HD2	2.17	0.44
1:B:180:ILE:O	1:B:184:LEU:HG	2.16	0.44
1:B:104:ILE:HA	1:B:107:LEU:CB	2.46	0.44
1:B:156:SER:OG	1:B:159:ARG:HB2	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:LYS:HG3	1:A:109:LYS:O	2.17	0.44
1:B:125:GLN:O	1:B:129:SER:OG	2.28	0.44
1:B:192:SER:HB2	1:B:266:LEU:HD11	2.00	0.44
1:B:255:ASP:C	1:B:257:ILE:N	2.74	0.44
1:B:136:LEU:C	1:B:138:PRO:CD	2.88	0.44
1:B:191:ILE:HG21	1:B:263:LEU:HD21	2.00	0.44
1:A:181:ALA:HB3	1:A:182:PRO:CD	2.48	0.44
1:B:237:LEU:O	1:B:237:LEU:CD2	2.65	0.44
1:B:260:ARG:HG2	1:B:260:ARG:NH1	2.33	0.44
1:A:206:GLU:HA	1:B:227:VAL:HG11	1.98	0.44
1:B:255:ASP:O	1:B:258:ASN:ND2	2.50	0.43
1:A:198:ASP:OD1	1:A:198:ASP:C	2.61	0.43
1:B:125:GLN:HG3	1:B:126:LEU:HD23	2.00	0.43
1:A:177:LEU:CD1	1:A:222:SER:HA	2.48	0.43
1:B:191:ILE:HG23	1:B:263:LEU:HD23	2.00	0.43
1:B:276:HIS:O	1:B:277:HIS:C	2.62	0.43
1:A:146:GLN:O	1:A:267:ASN:ND2	2.47	0.43
1:B:167:ALA:HB3	2:B:10:HOH:O	2.19	0.43
1:A:198:ASP:O	1:A:198:ASP:CG	2.61	0.43
1:B:130:ASP:C	1:B:130:ASP:OD2	2.62	0.42
1:B:277:HIS:HB2	2:B:72:HOH:O	2.19	0.42
1:B:124:ASP:OD2	1:B:124:ASP:N	2.51	0.42
1:B:197:THR:OG1	1:B:243:ALA:C	2.63	0.42
1:A:159:ARG:HA	1:A:160:PRO:HD2	1.97	0.42
1:A:137:ARG:NH1	2:A:59:HOH:O	2.52	0.42
1:A:172:TYR:O	1:A:176:ILE:HG13	2.20	0.41
1:B:128:GLU:HB2	1:B:137:ARG:HH22	1.83	0.41
1:B:183:VAL:O	1:B:184:LEU:C	2.63	0.41
1:B:132:LYS:HB2	2:B:83:HOH:O	2.19	0.41
1:B:154:ILE:HG22	1:B:155:ASP:O	2.21	0.41
1:B:232:ASP:HB3	1:B:235:LYS:HB2	2.01	0.41
1:B:272:GLN:HE22	1:B:277:HIS:CD2	2.37	0.41
1:B:107:LEU:C	1:B:107:LEU:HD13	2.45	0.41
1:B:123:LEU:O	1:B:126:LEU:HG	2.20	0.41
1:A:196:HIS:O	1:A:259:ARG:HA	2.20	0.41
1:B:125:GLN:CG	1:B:126:LEU:HD23	2.50	0.41
1:B:128:GLU:O	1:B:134:ARG:NH2	2.43	0.41
1:A:134:ARG:O	1:A:134:ARG:HG2	2.19	0.41
1:B:104:ILE:C	1:B:107:LEU:HB3	2.46	0.41
1:A:147:GLU:CB	1:A:264:LEU:HD21	2.51	0.41
1:A:149:LEU:HB3	1:A:265:VAL:HB	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:185:ASN:HA	1:B:235:LYS:HZ3	1.86	0.40
1:B:258:ASN:HB2	2:B:310:HOH:O	2.21	0.40
1:A:146:GLN:O	1:A:146:GLN:HG2	2.22	0.40
1:B:191:ILE:CG2	1:B:192:SER:N	2.83	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	166/183 (91%)	153 (92%)	10 (6%)	3 (2%)	6	9
1	B	155/183 (85%)	142 (92%)	12 (8%)	1 (1%)	21	32
All	All	321/366 (88%)	295 (92%)	22 (7%)	4 (1%)	10	16

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	103	ASN
1	A	269	GLN
1	B	183	VAL
1	A	145	VAL

5.3.2 Protein sidechains [i](#)

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	144/156 (92%)	116 (81%)	28 (19%)	1	2
1	B	136/156 (87%)	116 (85%)	20 (15%)	3	4
All	All	280/312 (90%)	232 (83%)	48 (17%)	2	2

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	102	PRO
1	A	104	ILE
1	A	107	LEU
1	A	113	GLN
1	A	120	ARG
1	A	125	GLN
1	A	133	LEU
1	A	134	ARG
1	A	137	ARG
1	A	141	LYS
1	A	142	ILE
1	A	146	GLN
1	A	150	ARG
1	A	157	GLN
1	A	163	LYS
1	A	164	THR
1	A	168	GLU
1	A	169	VAL
1	A	198	ASP
1	A	207	LYS
1	A	233	ASN
1	A	236	VAL
1	A	249	SER
1	A	259	ARG
1	A	262	SER
1	A	269	GLN
1	A	271	GLU
1	A	272	GLN
1	B	106	GLU
1	B	107	LEU
1	B	110	ARG
1	B	118	LYS
1	B	126	LEU
1	B	129	SER
1	B	133	LEU

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Mol	Chain	Res	Type
1	B	140	LEU
1	B	142	ILE
1	B	143	ASP
1	B	159	ARG
1	B	173	MET
1	B	190	ARG
1	B	233	ASN
1	B	237	LEU
1	B	257	ILE
1	B	258	ASN
1	B	268	LYS
1	B	271	GLU
1	B	276	HIS

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

Mol	Chain	Res	Type
1	A	113	GLN
1	A	146	GLN
1	A	152	GLN
1	A	258	ASN
1	A	272	GLN
1	B	139	HIS
1	B	233	ASN
1	B	258	ASN
1	B	277	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

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	170/183 (92%)	0.50	10 (5%) 28 24	23, 42, 78, 99	0
1	B	161/183 (87%)	0.92	21 (13%) 7 5	25, 53, 81, 96	0
All	All	331/366 (90%)	0.70	31 (9%) 14 11	23, 49, 81, 99	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	276	HIS	4.7
1	B	233	ASN	3.9
1	B	256	ALA	3.3
1	B	140	LEU	3.3
1	A	255	ASP	3.2
1	B	175	ASP	3.2
1	B	200	PHE	3.1
1	B	104	ILE	3.0
1	B	117	ASN	3.0
1	A	276	HIS	2.9
1	B	244	ALA	2.9
1	A	206	GLU	2.7
1	B	277	HIS	2.6
1	B	142	ILE	2.6
1	A	104	ILE	2.6
1	A	102	PRO	2.5
1	B	199	ASP	2.5
1	A	256	ALA	2.5
1	B	105	ASP	2.4
1	A	126	LEU	2.4
1	B	275	LEU	2.4
1	B	255	ASP	2.3
1	B	126	LEU	2.2
1	A	208	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	249	SER	2.2
1	B	141	LYS	2.2
1	B	197	THR	2.2
1	A	114	SER	2.2
1	B	143	ASP	2.2
1	B	274	ILE	2.1
1	B	198	ASP	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.