



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:11 AM GMT

PDB ID : 2ZVZ
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

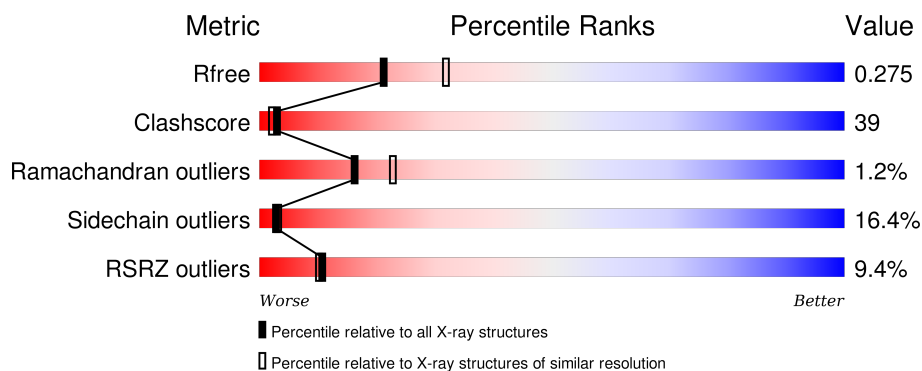
1 Overall quality at a glance

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}	91344	2919 (2.40-2.40)
Clashscore	102246	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	
1	B	183	

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

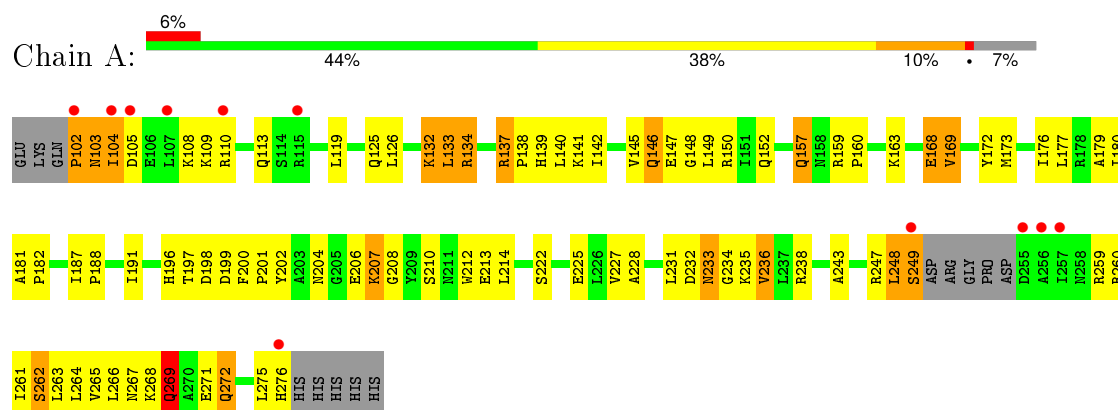
- Molecule 2 is water.

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

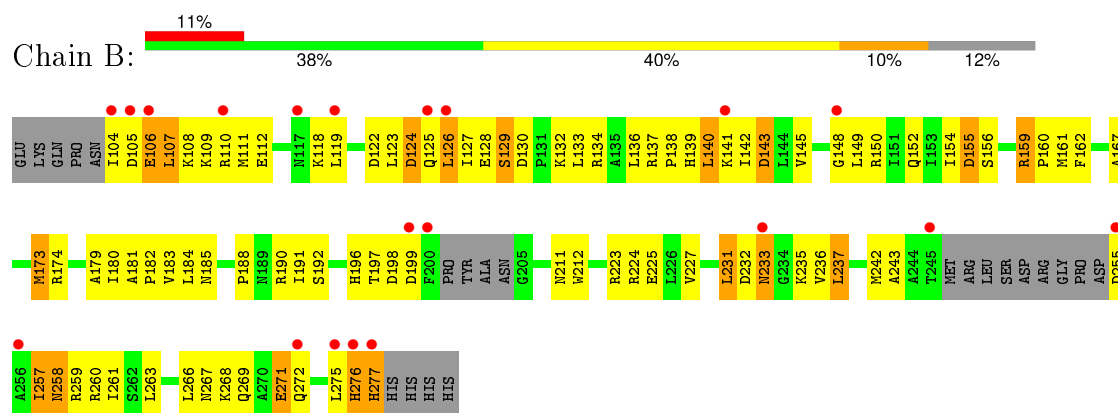
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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



• 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	45.20 Å 46.24 Å 173.70 Å 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 2.39 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.245 , 0.275 0.246 , 0.275	Depositor DCC
R_{free} test set	723 reflections (4.87%)	DCC
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 , 63.5	EDS
Estimated twinning fraction	0.042 for k,h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 14836 reflections	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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.47	0/1355	0.84	2/1822 (0.1%)
1	B	0.39	0/1282	0.77	0/1721
All	All	0.43	0/2637	0.81	2/3543 (0.1%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	248	LEU	CA-CB-CG	-9.17	94.21	115.30
1	A	248	LEU	N-CA-C	6.66	128.99	111.00

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	86	0
1	B	1269	0	1307	122	0
2	A	146	0	0	30	0
2	B	90	0	0	8	0
All	All	2844	0	2687	206	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 39.

All (206) 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
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: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:HE2	1.76	0.67
1:B:127:ILE:CG2	1:B:137:ARG:HG3	2.25	0.67
1:B:108:LYS:HG2	1:B:111:MET:CE	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:139:HIS:HE1	1:B:156:SER:HB3	1.60	0.65
1:B:104:ILE:HA	1:B:107:LEU:HB3	1.79	0.65
1:B:257:ILE:H	1:B:257:ILE:HD13	1.62	0.65
1:B:269:GLN:NE2	2:B:294:HOH:O	2.30	0.65
1:B:140:LEU:O	1:B:141:LYS:HG2	1.96	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: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:179:ALA:C	2:A:319:HOH:O	2.38	0.61
1:A:157:GLN:HB2	2:A:290:HOH:O	1.99	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:B:137:ARG:N	1:B:138:PRO:HD2	2.16	0.61
1:A:212:TRP:N	2:A:323:HOH:O	2.34	0.61
1:B:275:LEU:O	1:B:276:HIS:HB2	1.99	0.60
1:B:198:ASP:HB2	1:B:259:ARG:CG	2.31	0.60
1:B:174:ARG:HG3	1:B:225:GLU:OE1	2.00	0.60
1:A:248:LEU:O	1:A:249:SER:CB	2.50	0.60
1:A:210:SER:C	2:A:323:HOH:O	2.40	0.60
1:A:275:LEU:O	1:A:276:HIS:C	2.39	0.60
1:A:231:LEU:O	2:A:347:HOH:O	2.15	0.60
1:B:136:LEU:C	1:B:138:PRO:HD2	2.23	0.59
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:CA	1:A:272:GLN:HE21	2.11	0.58
1:B:130:ASP:HB3	1:B:133:LEU:HG	1.86	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:232:ASP:CG	1:B:233:ASN:N	2.57	0.57
1:A:168:GLU:CA	1:A:168:GLU:OE1	2.50	0.56
1:B:108:LYS:HA	1:B:111:MET:HE2	1.86	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
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:223:ARG:NH2	1:B:224:ARG:HG3	2.23	0.54
1:B:188:PRO:HD2	2:B:98:HOH:O	2.06	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:CE1	1:B:156:SER:HB3	2.42	0.53
1:B:139:HIS:HE1	1:B:156:SER:CB	2.20	0.53
1:B:161:MET:HB3	1:B:173:MET:HG3	1.90	0.53
1:B:232:ASP:CG	1:B:233:ASN:H	2.12	0.53
1:A:272:GLN:HA	1:A:272:GLN:NE2	2.14	0.53
1:B:119:LEU:HG	1:B:149:LEU:CD2	2.38	0.53
1:B:145:VAL:HG22	1:B:148:GLY:O	2.09	0.53
1:A:102:PRO:O	1:A:103:ASN:HB2	2.08	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:B:185:ASN:HA	1:B:235:LYS:HZ3	1.75	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:B:162:PHE:CE2	1:B:261:ILE:HD11	2.46	0.51
1:A:148:GLY:HA3	1:A:265:VAL:O	2.11	0.51
1:A:233:ASN:N	1:A:233:ASN:OD1	2.31	0.51
1:A:137:ARG:HD2	2:A:329:HOH:O	2.11	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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:191:ILE:HD13	1:B:263:LEU:HD21	1.94	0.49
1:B:191:ILE:CG2	1:B:263:LEU:HD23	2.42	0.49
1:B:127:ILE:HD13	1:B:140:LEU:HD13	1.93	0.49
1:A:137:ARG:HG2	1:A:138:PRO:HD3	1.94	0.49
1:A:236:VAL:HG22	2:A:71:HOH:O	2.12	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:HD13	2.34	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:B:140:LEU:C	1:B:141:LYS:HG2	2.33	0.48
1:B:198:ASP:O	1:B:199:ASP:C	2.51	0.48
1:A:248:LEU:HA	1:A:248:LEU:HD23	1.35	0.48
1:B:150:ARG:NH1	1:B:152:GLN:OE1	2.47	0.48
1:B:196:HIS:CE1	1:B:242:MET:HG3	2.49	0.48
1:A:225:GLU:HG3	2:A:336:HOH:O	2.14	0.47
1:B:275:LEU:O	1:B:276:HIS:CB	2.63	0.47
1:B:255:ASP:CA	1:B:257:ILE:HD13	2.45	0.47
1:B:190:ARG:HG2	1:B:235:LYS:O	2.15	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: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:B:136:LEU:O	1:B:137:ARG:C	2.55	0.45
1:A:234:GLY:HA2	2:A:320:HOH:O	2.16	0.45
1:B:268:LYS:O	1:B:272:GLN:HG2	2.16	0.45
1:B:104:ILE:HG23	1:B:105:ASP:N	2.31	0.45
1:A:177:LEU:HD13	1:A:222:SER:HA	1.99	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:104:ILE:N	1:B:107:LEU:CB	2.80	0.44
1:B:180:ILE:O	1:B:184:LEU:HG	2.16	0.44
1:B:179:ALA:O	1:B:182:PRO:HD2	2.17	0.44
1:B:156:SER:OG	1:B:159:ARG:HB2	2.18	0.44
1:B:104:ILE:HA	1:B:107:LEU:CB	2.46	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:A:276:HIS:ND1	1:A:276:HIS:C	2.71	0.44
1:B:192:SER:HB2	1:B:266:LEU:HD11	2.00	0.44
1:B:184:LEU:O	1:B:235:LYS:NZ	2.33	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:NH1	1:B:260:ARG:HG2	2.33	0.44
1:B:255:ASP:C	1:B:257:ILE:H	2.21	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:B:197:THR:HG1	1:B:211:ASN:HB3	1.83	0.43
1:A:260:ARG:C	1:A:261:ILE:HD12	2.38	0.43
1:B:125:GLN:HG3	1:B:126:LEU:HD23	2.00	0.43
1:B:191:ILE:HG23	1:B:263:LEU:HD23	2.00	0.43
1:A:177:LEU:CD1	1:A:222:SER:HA	2.48	0.43
1:A:132:LYS:C	1:A:133:LEU:HD23	2.39	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:B:277:HIS:HB2	2:B:72:HOH:O	2.19	0.42
1:B:124:ASP:N	1:B:124:ASP:OD2	2.51	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:B:128:GLU:HB2	1:B:137:ARG:HH22	1.83	0.41
1:A:172:TYR:O	1:A:176:ILE:HG13	2.20	0.41
1:B:132:LYS:HB2	2:B:83:HOH:O	2.19	0.41
1:B:272:GLN:HE22	1:B:277:HIS:CD2	2.37	0.41
1:B:232:ASP:HB3	1:B:235:LYS:HB2	2.01	0.41
1:B:154:ILE:HG22	1:B:155:ASP:O	2.21	0.41
1:B:123:LEU:O	1:B:126:LEU:HG	2.20	0.41
1:B:142:ILE:C	1:B:143:ASP:OD2	2.58	0.41
1:B:125:GLN:CG	1:B:126:LEU:HD23	2.50	0.41
1:A:196:HIS:O	1:A:259:ARG:HA	2.20	0.41
1:B:128:GLU:O	1:B:134:ARG:NH2	2.43	0.41
1:A:137:ARG:C	1:A:139:HIS:H	2.23	0.41
1:A:134:ARG:HG2	1:A:134:ARG:O	2.19	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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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%)	11	13
1	B	155/183 (85%)	142 (92%)	12 (8%)	1 (1%)	30	43
All	All	321/366 (88%)	295 (92%)	22 (7%)	4 (1%)	16	23

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%)	118 (82%)	26 (18%)	2	2

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	136/156 (87%)	116 (85%)	20 (15%)	4	4
All	All	280/312 (90%)	234 (84%)	46 (16%)	3	3

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	102	PRO
1	A	104	ILE
1	A	113	GLN
1	A	125	GLN
1	A	132	LYS
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	168	GLU
1	A	169	VAL
1	A	198	ASP
1	A	207	LYS
1	A	233	ASN
1	A	236	VAL
1	A	247	ARG
1	A	249	SER
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	124	ASP
1	B	126	LEU
1	B	129	SER
1	B	140	LEU
1	B	143	ASP
1	B	155	ASP

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Mol	Chain	Res	Type
1	B	159	ARG
1	B	173	MET
1	B	231	LEU
1	B	233	ASN
1	B	237	LEU
1	B	257	ILE
1	B	258	ASN
1	B	271	GLU
1	B	276	HIS
1	B	277	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	113	GLN
1	A	146	GLN
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 carbohydrates 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	170/183 (92%)	0.36	11 (6%) 22 22	23, 42, 78, 99	0
1	B	161/183 (87%)	0.70	20 (12%) 5 5	25, 53, 81, 96	0
All	All	331/366 (90%)	0.53	31 (9%) 11 10	23, 49, 81, 99	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	200	PHE	8.5
1	B	105	ASP	5.6
1	A	104	ILE	5.6
1	B	277	HIS	5.4
1	A	102	PRO	5.1
1	A	276	HIS	5.0
1	B	110	ARG	4.7
1	B	276	HIS	4.5
1	B	233	ASN	4.3
1	A	255	ASP	3.8
1	B	126	LEU	3.6
1	B	256	ALA	3.3
1	A	110	ARG	3.2
1	B	199	ASP	3.2
1	B	119	LEU	3.0
1	B	148	GLY	2.8
1	A	105	ASP	2.7
1	A	257	ILE	2.7
1	A	249	SER	2.7
1	B	106	GLU	2.6
1	B	245	THR	2.6
1	B	125	GLN	2.6
1	B	275	LEU	2.5
1	B	117	ASN	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	141	LYS	2.3
1	B	255	ASP	2.2
1	A	256	ALA	2.2
1	B	272	GLN	2.2
1	A	107	LEU	2.2
1	B	104	ILE	2.1
1	A	115	ARG	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 carbohydrates 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.