



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:55 PM GMT

PDB ID : 1R9G
Title : Three-dimensional Structure of YaaE from Bacillus subtilis
Authors : Bauer, J.A.; Bennett, E.M.; Begley, T.P.; Ealick, S.E.
Deposited on : 2003-10-29
Resolution : 2.50 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

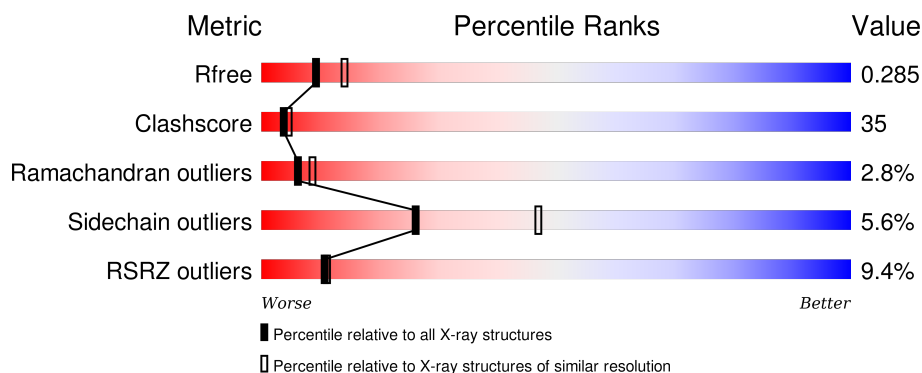
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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 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	216	
1	B	216	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3293 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 Hypothetical protein yaaE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	196	Total	C	N	O	S	Se	0	0	0
			1507	954	262	283	3	5			
1	B	196	Total	C	N	O	S	Se	0	0	0
			1507	954	262	283	3	5			

There are 50 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MSE	-	CLONING ARTIFACT	UNP P37528
A	-18	GLY	-	CLONING ARTIFACT	UNP P37528
A	-17	SER	-	CLONING ARTIFACT	UNP P37528
A	-16	SER	-	CLONING ARTIFACT	UNP P37528
A	-15	HIS	-	EXPRESSION TAG	UNP P37528
A	-14	HIS	-	EXPRESSION TAG	UNP P37528
A	-13	HIS	-	EXPRESSION TAG	UNP P37528
A	-12	HIS	-	EXPRESSION TAG	UNP P37528
A	-11	HIS	-	EXPRESSION TAG	UNP P37528
A	-10	HIS	-	EXPRESSION TAG	UNP P37528
A	-9	SER	-	CLONING ARTIFACT	UNP P37528
A	-8	SER	-	CLONING ARTIFACT	UNP P37528
A	-7	GLY	-	CLONING ARTIFACT	UNP P37528
A	-6	LEU	-	CLONING ARTIFACT	UNP P37528
A	-5	VAL	-	CLONING ARTIFACT	UNP P37528
A	-4	PRO	-	CLONING ARTIFACT	UNP P37528
A	-3	ARG	-	CLONING ARTIFACT	UNP P37528
A	-2	GLY	-	CLONING ARTIFACT	UNP P37528
A	-1	SER	-	CLONING ARTIFACT	UNP P37528
A	0	HIS	-	CLONING ARTIFACT	UNP P37528
A	1	MSE	MET	MODIFIED RESIDUE	UNP P37528
A	52	MSE	MET	MODIFIED RESIDUE	UNP P37528
A	62	MSE	MET	MODIFIED RESIDUE	UNP P37528
A	75	MSE	MET	MODIFIED RESIDUE	UNP P37528
A	186	MSE	MET	MODIFIED RESIDUE	UNP P37528

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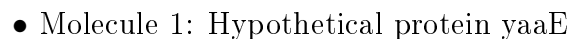
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Chain	Residue	Modelled	Actual	Comment	Reference
B	-19	MSE	-	CLONING ARTIFACT	UNP P37528
B	-18	GLY	-	CLONING ARTIFACT	UNP P37528
B	-17	SER	-	CLONING ARTIFACT	UNP P37528
B	-16	SER	-	CLONING ARTIFACT	UNP P37528
B	-15	HIS	-	EXPRESSION TAG	UNP P37528
B	-14	HIS	-	EXPRESSION TAG	UNP P37528
B	-13	HIS	-	EXPRESSION TAG	UNP P37528
B	-12	HIS	-	EXPRESSION TAG	UNP P37528
B	-11	HIS	-	EXPRESSION TAG	UNP P37528
B	-10	HIS	-	EXPRESSION TAG	UNP P37528
B	-9	SER	-	CLONING ARTIFACT	UNP P37528
B	-8	SER	-	CLONING ARTIFACT	UNP P37528
B	-7	GLY	-	CLONING ARTIFACT	UNP P37528
B	-6	LEU	-	CLONING ARTIFACT	UNP P37528
B	-5	VAL	-	CLONING ARTIFACT	UNP P37528
B	-4	PRO	-	CLONING ARTIFACT	UNP P37528
B	-3	ARG	-	CLONING ARTIFACT	UNP P37528
B	-2	GLY	-	CLONING ARTIFACT	UNP P37528
B	-1	SER	-	CLONING ARTIFACT	UNP P37528
B	0	HIS	-	CLONING ARTIFACT	UNP P37528
B	1	MSE	MET	MODIFIED RESIDUE	UNP P37528
B	52	MSE	MET	MODIFIED RESIDUE	UNP P37528
B	62	MSE	MET	MODIFIED RESIDUE	UNP P37528
B	75	MSE	MET	MODIFIED RESIDUE	UNP P37528
B	186	MSE	MET	MODIFIED RESIDUE	UNP P37528

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	148	Total O 148 148	0	0
2	B	131	Total O 131 131	0	0

- Molecule 1: Hypothetical protein yaaE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	45.50 Å 80.87 Å 115.39 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50 46.97 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.8 (50.00-2.50) 99.8 (46.97-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.95 (at 2.51 Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.240 , 0.279 0.250 , 0.285	Depositor DCC
R_{free} test set	758 reflections (5.21%)	DCC
Wilson B-factor (Å ²)	49.3	Xtriage
Anisotropy	0.257	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 50.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 28442 reflections (0.004%)	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	3293	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.47% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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

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	0.43	0/1529	0.62	0/2058
1	B	0.40	0/1529	0.63	0/2058
All	All	0.41	0/3058	0.63	0/4116

There are no bond length outliers.

There are no bond angle outliers.

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	1507	0	1512	111	0
1	B	1507	0	1512	105	0
2	A	148	0	0	30	0
2	B	131	0	0	18	0
All	All	3293	0	3024	212	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 35.

All (212) 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:B:32:ARG:HH11	1:B:32:ARG:HB2	1.03	1.07

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:32:ARG:NH1	1:B:32:ARG:HB2	1.74	1.03
1:B:165:LEU:HD22	1:B:186:MSE:HE1	1.36	1.02
1:A:52:MSE:HE2	1:A:84:ILE:HG13	1.46	0.97
1:B:75:MSE:HE3	1:B:164:PHE:HE1	1.30	0.95
1:B:32:ARG:H	1:B:35:GLN:HE21	1.14	0.94
1:A:145:ASN:HD22	1:A:145:ASN:H	1.10	0.93
1:B:174:THR:HB	2:B:257:HOH:O	1.68	0.92
1:A:102:VAL:HB	2:A:235:HOH:O	1.72	0.90
1:B:31:LYS:H	1:B:35:GLN:NE2	1.71	0.89
1:A:130:THR:HG22	2:A:278:HOH:O	1.77	0.84
1:A:146:VAL:HG11	2:A:235:HOH:O	1.80	0.80
1:B:145:ASN:H	1:B:145:ASN:HD22	1.29	0.80
1:A:17:ILE:HD11	1:A:29:VAL:HG22	1.64	0.80
1:B:183:PHE:HA	1:B:186:MSE:HE3	1.63	0.78
1:B:75:MSE:HE3	1:B:164:PHE:CE1	2.16	0.78
1:A:160:LYS:HB3	2:A:309:HOH:O	1.83	0.77
1:B:69:ALA:HB2	1:B:75:MSE:HE1	1.64	0.77
1:B:181:GLN:HG2	2:B:245:HOH:O	1.85	0.76
1:B:9:LEU:HD21	1:B:31:LYS:HE3	1.66	0.76
1:A:69:ALA:CB	1:A:75:MSE:HE1	2.16	0.75
1:B:32:ARG:CB	1:B:32:ARG:HH11	1.92	0.73
1:A:61:PHE:HA	2:A:302:HOH:O	1.89	0.72
1:B:69:ALA:CB	1:B:75:MSE:HE1	2.21	0.69
1:A:4:ILE:HB	1:A:20:ILE:HD13	1.73	0.69
1:B:69:ALA:HA	1:B:75:MSE:HE2	1.74	0.69
1:B:28:LEU:HD23	1:B:29:VAL:N	2.08	0.69
1:B:165:LEU:HD22	1:B:186:MSE:CE	2.19	0.69
1:A:2:LEU:H	1:A:2:LEU:HD22	1.57	0.69
1:A:69:ALA:HB2	1:A:75:MSE:HE1	1.75	0.68
1:A:149:LEU:HD12	2:A:276:HOH:O	1.93	0.67
1:B:2:LEU:HG	2:B:269:HOH:O	1.94	0.67
1:A:145:ASN:H	1:A:145:ASN:ND2	1.90	0.67
1:A:102:VAL:CG2	1:A:139:ILE:HG23	2.24	0.67
1:A:132:VAL:HG23	2:A:278:HOH:O	1.94	0.67
1:B:145:ASN:N	1:B:145:ASN:HD22	1.93	0.66
1:B:102:VAL:CG2	1:B:139:ILE:HG23	2.26	0.66
1:B:32:ARG:N	1:B:35:GLN:HE21	1.89	0.65
1:B:2:LEU:HD23	1:B:3:THR:N	2.12	0.65
1:A:34:GLU:CD	1:A:34:GLU:H	2.00	0.64
1:A:47:GLY:N	1:A:80:ALA:HB3	2.13	0.64
1:A:10:GLN:HB2	1:A:55:LEU:HD12	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:52:MSE:O	1:A:56:ILE:HG23	1.97	0.64
1:A:63:GLU:HB2	1:A:64:PRO:HD3	1.80	0.63
1:A:62:MSE:HG3	2:A:287:HOH:O	1.99	0.62
1:B:89:ILE:C	1:B:91:GLY:H	2.03	0.62
1:A:145:ASN:N	1:A:145:ASN:HD22	1.83	0.62
1:B:89:ILE:O	1:B:91:GLY:N	2.33	0.62
1:A:16:HIS:HB2	2:A:333:HOH:O	1.98	0.62
1:A:2:LEU:HD21	1:A:188:GLU:HG2	1.82	0.62
1:B:102:VAL:HG21	1:B:139:ILE:HG23	1.81	0.62
1:A:6:VAL:HG11	2:A:333:HOH:O	1.99	0.61
1:B:1:MSE:HE3	1:B:2:LEU:H	1.65	0.61
1:A:46:GLY:CA	1:A:80:ALA:HB3	2.31	0.61
1:A:40:ASP:OD1	1:A:191:LYS:HE3	2.00	0.60
1:A:85:LEU:HD22	1:A:100:LEU:HB2	1.84	0.60
1:B:75:MSE:CE	1:B:164:PHE:HE1	2.11	0.60
1:A:195:LEU:O	1:A:195:LEU:HD12	2.02	0.59
1:B:89:ILE:O	1:B:89:ILE:HG13	2.02	0.59
1:B:50:THR:HB	2:B:231:HOH:O	2.02	0.59
1:B:127:GLU:HG3	2:B:288:HOH:O	2.02	0.59
1:A:85:LEU:O	1:A:85:LEU:HD23	2.03	0.59
1:A:111:ARG:NH1	1:A:135:ARG:HD3	2.17	0.58
1:A:75:MSE:HE3	1:A:164:PHE:CE1	2.39	0.58
1:B:106:ARG:NE	2:B:233:HOH:O	2.37	0.58
1:B:131:GLY:HA2	2:B:209:HOH:O	2.03	0.58
1:A:69:ALA:HA	1:A:75:MSE:CE	2.34	0.58
1:B:165:LEU:CD2	1:B:186:MSE:HE1	2.23	0.57
1:B:44:LEU:N	1:B:44:LEU:HD12	2.19	0.57
1:B:17:ILE:HD11	1:B:29:VAL:HG22	1.87	0.57
1:A:47:GLY:H	1:A:80:ALA:HB3	1.70	0.57
1:B:13:VAL:HG12	2:B:225:HOH:O	2.05	0.57
1:A:17:ILE:HG13	2:A:333:HOH:O	2.05	0.56
1:B:106:ARG:HG3	1:B:106:ARG:HH11	1.70	0.56
1:B:94:ASN:HB3	1:B:95:PRO:CD	2.35	0.56
1:A:75:MSE:HE3	1:A:164:PHE:HE1	1.70	0.56
1:B:11:GLY:O	1:B:12:ALA:HB3	2.06	0.56
1:B:69:ALA:HA	1:B:75:MSE:CE	2.35	0.56
1:A:141:GLU:HG3	2:A:239:HOH:O	2.07	0.55
1:B:51:THR:O	1:B:55:LEU:HD23	2.06	0.55
1:B:128:PRO:HD2	2:B:288:HOH:O	2.06	0.55
1:B:192:GLN:N	1:B:192:GLN:OE1	2.32	0.55
1:A:171:PRO:O	1:A:174:THR:HG22	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1:MSE:HE3	1:B:1:MSE:N	2.21	0.55
1:A:174:THR:HG23	1:A:176:ASP:H	1.72	0.55
1:B:1:MSE:HG3	2:B:252:HOH:O	2.06	0.55
1:A:56:ILE:HD11	1:A:62:MSE:CE	2.37	0.55
1:B:124:GLY:C	1:B:125:LEU:HD12	2.28	0.55
1:B:31:LYS:H	1:B:35:GLN:HE22	1.54	0.54
1:A:32:ARG:N	1:A:35:GLN:HE21	2.06	0.54
1:A:181:GLN:HG2	2:A:253:HOH:O	2.06	0.54
1:B:63:GLU:HB2	1:B:64:PRO:HD3	1.89	0.54
1:A:6:VAL:HG21	2:A:333:HOH:O	2.08	0.54
1:A:90:ALA:HA	1:B:112:GLN:CD	2.28	0.54
1:B:86:ALA:HB1	1:B:102:VAL:O	2.08	0.54
1:A:55:LEU:HA	2:A:342:HOH:O	2.07	0.54
1:B:8:GLY:HA3	2:B:262:HOH:O	2.07	0.54
1:A:23:CYS:HB3	1:A:181:GLN:HE22	1.74	0.54
1:B:1:MSE:HE3	1:B:1:MSE:H3	1.74	0.53
1:B:143:GLY:O	1:B:146:VAL:HG13	2.08	0.53
1:A:64:PRO:HB2	2:A:302:HOH:O	2.08	0.53
1:B:73:LYS:HE3	2:B:293:HOH:O	2.09	0.53
1:A:49:SER:HB3	1:A:106:ARG:HE	1.73	0.53
1:B:99:LEU:HB3	1:B:164:PHE:CZ	2.44	0.52
1:B:172:GLU:N	1:B:172:GLU:OE1	2.39	0.52
1:A:57:ASP:HB2	2:A:313:HOH:O	2.09	0.52
1:A:17:ILE:HD11	1:A:29:VAL:CG2	2.35	0.52
1:B:106:ARG:NH1	1:B:106:ARG:HG3	2.24	0.52
1:B:149:LEU:HD11	1:B:160:LYS:HB2	1.91	0.52
1:A:49:SER:H	1:A:106:ARG:NH2	2.07	0.52
1:B:91:GLY:O	1:B:92:SER:HB3	2.10	0.52
1:B:89:ILE:C	1:B:91:GLY:N	2.63	0.52
1:A:53:ARG:NE	1:A:96:HIS:ND1	2.57	0.51
1:A:177:HIS:HE1	2:A:308:HOH:O	1.93	0.51
1:A:145:ASN:N	1:A:145:ASN:ND2	2.56	0.51
1:A:63:GLU:HG2	2:A:303:HOH:O	2.11	0.51
1:B:48:GLU:OE2	1:B:52:MSE:HB2	2.11	0.51
1:B:28:LEU:HD23	1:B:28:LEU:C	2.30	0.50
1:B:158:ALA:HA	1:B:166:GLY:O	2.11	0.50
1:A:2:LEU:N	1:A:2:LEU:HD22	2.26	0.50
1:A:31:LYS:H	1:A:35:GLN:NE2	2.10	0.50
1:A:23:CYS:HB3	1:A:181:GLN:NE2	2.25	0.50
1:A:123:LYS:HE3	2:A:285:HOH:O	2.11	0.50
1:A:90:ALA:N	1:B:112:GLN:HE22	2.09	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:101:ASN:HD22	1:A:145:ASN:HD21	1.59	0.50
1:B:160:LYS:HE3	2:B:241:HOH:O	2.12	0.50
1:B:50:THR:HG22	1:B:50:THR:O	2.11	0.49
1:A:89:ILE:HG22	1:A:92:SER:HB3	1.95	0.49
1:A:160:LYS:HD3	2:A:309:HOH:O	2.12	0.49
1:A:49:SER:HB3	1:A:106:ARG:NE	2.27	0.49
1:A:32:ARG:H	1:A:35:GLN:HE21	1.60	0.48
1:A:155:ARG:NH1	1:A:155:ARG:HG2	2.27	0.48
1:A:106:ARG:HH11	1:A:106:ARG:HG2	1.78	0.48
1:A:69:ALA:HA	1:A:75:MSE:HE2	1.94	0.48
1:A:53:ARG:NH2	1:A:94:ASN:ND2	2.62	0.48
1:B:29:VAL:HG23	2:B:283:HOH:O	2.13	0.48
1:B:101:ASN:HD22	1:B:145:ASN:HD21	1.62	0.47
1:B:2:LEU:HD23	1:B:2:LEU:C	2.35	0.47
1:B:102:VAL:HG21	1:B:139:ILE:CG2	2.44	0.47
1:A:135:ARG:NH1	2:A:230:HOH:O	2.47	0.47
1:B:52:MSE:HE3	1:B:84:ILE:HG13	1.96	0.47
1:B:31:LYS:N	1:B:35:GLN:NE2	2.53	0.47
1:A:69:ALA:CA	1:A:75:MSE:HE1	2.44	0.47
1:B:51:THR:O	1:B:54:ARG:HB2	2.14	0.47
1:A:32:ARG:H	1:A:35:GLN:NE2	2.12	0.47
1:B:52:MSE:CE	1:B:84:ILE:HG13	2.45	0.47
1:B:73:LYS:HG2	1:B:190:TYR:CE2	2.50	0.47
1:B:91:GLY:O	1:B:92:SER:CB	2.64	0.46
1:A:75:MSE:CE	1:A:164:PHE:HE1	2.28	0.46
1:A:52:MSE:CE	1:A:84:ILE:HG13	2.33	0.46
1:A:53:ARG:HH22	1:A:94:ASN:HD21	1.64	0.46
1:B:46:GLY:HA3	1:B:80:ALA:HB3	1.97	0.46
1:A:44:LEU:HD22	1:A:44:LEU:N	2.30	0.46
1:A:155:ARG:HG2	1:A:155:ARG:HH11	1.80	0.46
1:A:165:LEU:HB2	1:A:186:MSE:HE1	1.96	0.46
1:A:69:ALA:HA	1:A:75:MSE:HE1	1.97	0.46
1:A:111:ARG:HA	1:A:114:ASP:OD2	2.15	0.46
1:B:122:ILE:HG12	1:B:150:SER:HB2	1.99	0.45
1:B:13:VAL:HG13	1:B:14:ARG:N	2.30	0.45
1:A:43:ILE:O	1:A:45:PRO:HD3	2.16	0.45
1:A:196:VAL:HG12	1:A:196:VAL:OXT	2.16	0.45
1:A:33:PRO:HB3	2:A:302:HOH:O	2.17	0.45
1:A:111:ARG:HH11	1:A:135:ARG:HD3	1.81	0.45
1:A:3:THR:HG22	1:A:39:VAL:HG12	1.98	0.45
1:A:174:THR:HG23	1:A:176:ASP:N	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:46:GLY:HA3	1:A:80:ALA:HB3	1.98	0.45
1:B:179:VAL:O	1:B:182:LEU:HB2	2.17	0.45
1:B:145:ASN:H	1:B:145:ASN:ND2	2.05	0.44
1:A:177:HIS:CE1	2:A:308:HOH:O	2.70	0.44
1:A:92:SER:O	1:A:93:ASP:O	2.34	0.44
1:B:62:MSE:CE	1:B:97:LEU:HD23	2.47	0.44
1:B:94:ASN:HB2	2:B:258:HOH:O	2.17	0.44
1:A:15:GLU:HG3	2:A:226:HOH:O	2.16	0.44
1:B:16:HIS:O	1:B:20:ILE:HG13	2.18	0.44
1:A:76:PHE:CZ	1:A:78:THR:HG22	2.53	0.44
1:B:66:ARG:NE	2:B:250:HOH:O	2.50	0.44
1:B:20:ILE:HD11	1:B:43:ILE:HD12	2.00	0.44
1:A:149:LEU:HD21	2:A:309:HOH:O	2.17	0.43
1:B:13:VAL:HG22	1:B:17:ILE:CD1	2.48	0.43
1:B:155:ARG:NH2	2:B:246:HOH:O	2.51	0.43
1:A:53:ARG:O	1:A:56:ILE:HG12	2.18	0.43
1:A:89:ILE:HD12	1:A:90:ALA:H	1.83	0.43
1:A:90:ALA:N	1:B:112:GLN:NE2	2.67	0.43
1:A:130:THR:HA	2:A:286:HOH:O	2.18	0.43
1:A:6:VAL:HG13	1:A:45:PRO:HG3	2.00	0.43
1:A:89:ILE:C	1:B:112:GLN:HE22	2.21	0.43
1:A:47:GLY:H	1:A:80:ALA:CB	2.32	0.42
1:B:94:ASN:HB3	1:B:95:PRO:HD2	2.00	0.42
1:A:10:GLN:HB2	1:A:55:LEU:CD1	2.47	0.42
1:B:169:PHE:O	1:B:171:PRO:HD3	2.19	0.42
1:B:1:MSE:SE	1:B:2:LEU:N	3.02	0.42
1:B:196:VAL:HG12	1:B:196:VAL:O	2.20	0.42
1:A:9:LEU:HD12	1:A:56:ILE:HG22	2.02	0.42
1:A:130:THR:HG23	2:A:286:HOH:O	2.19	0.42
1:B:32:ARG:HH12	1:B:35:GLN:CG	2.33	0.41
1:A:7:LEU:HD23	1:A:52:MSE:HE3	2.02	0.41
1:B:114:ASP:C	1:B:114:ASP:OD1	2.59	0.41
1:A:147:GLU:CD	2:A:309:HOH:O	2.59	0.41
1:B:1:MSE:CE	1:B:1:MSE:H3	2.34	0.41
1:B:1:MSE:CE	1:B:2:LEU:H	2.32	0.41
1:B:183:PHE:CA	1:B:186:MSE:HE3	2.40	0.41
1:A:89:ILE:O	1:A:90:ALA:C	2.58	0.41
1:B:142:ALA:HB1	1:B:146:VAL:CG2	2.50	0.41
1:A:53:ARG:HG2	1:A:96:HIS:HE1	1.86	0.41
1:A:10:GLN:HB3	1:A:10:GLN:HE21	1.60	0.41
1:B:11:GLY:O	1:B:12:ALA:CB	2.69	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:SER:N	1:A:106:ARG:NH2	2.69	0.41
1:A:138:HIS:HE1	2:A:261:HOH:O	2.04	0.40
1:B:187:VAL:O	1:B:190:TYR:HB3	2.21	0.40
1:B:31:LYS:HD2	2:B:253:HOH:O	2.21	0.40
1:B:160:LYS:HE2	1:B:186:MSE:SE	2.71	0.40
1:A:2:LEU:H	1:A:2:LEU:CD2	2.30	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	194/216 (90%)	169 (87%)	22 (11%)	3 (2%)	13	22
1	B	194/216 (90%)	169 (87%)	17 (9%)	8 (4%)	3	4
All	All	388/432 (90%)	338 (87%)	39 (10%)	11 (3%)	6	9

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	90	ALA
1	A	93	ASP
1	B	90	ALA
1	B	92	SER
1	B	47	GLY
1	B	54	ARG
1	A	95	PRO
1	B	12	ALA
1	B	49	SER
1	B	89	ILE
1	B	91	GLY

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	160/171 (94%)	151 (94%)	9 (6%)	26	47
1	B	160/171 (94%)	151 (94%)	9 (6%)	26	47
All	All	320/342 (94%)	302 (94%)	18 (6%)	26	47

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	28	LEU
1	A	34	GLU
1	A	58	THR
1	A	61	PHE
1	A	145	ASN
1	A	163	GLN
1	A	192	GLN
1	A	195	LEU
1	B	1	MSE
1	B	3	THR
1	B	15	GLU
1	B	61	PHE
1	B	96	HIS
1	B	119	ASP
1	B	121	THR
1	B	145	ASN
1	B	155	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	A	35	GLN
1	A	37	ASN
1	A	94	ASN

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Mol	Chain	Res	Type
1	A	138	HIS
1	A	145	ASN
1	A	181	GLN
1	A	192	GLN
1	B	35	GLN
1	B	94	ASN
1	B	112	GLN
1	B	145	ASN

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	191/216 (88%)	0.56	15 (7%) 15 17	29, 50, 87, 115	0
1	B	191/216 (88%)	0.90	21 (10%) 7 7	28, 53, 103, 123	0
All	All	382/432 (88%)	0.73	36 (9%) 11 11	28, 51, 96, 123	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	195	LEU	20.1
1	B	196	VAL	18.4
1	B	194	ALA	14.1
1	B	60	GLN	6.3
1	A	46	GLY	4.3
1	A	90	ALA	4.1
1	B	193	LYS	4.0
1	A	106	ARG	3.8
1	B	10	GLN	3.5
1	A	91	GLY	3.4
1	A	196	VAL	3.0
1	A	93	ASP	2.9
1	B	2	LEU	2.8
1	A	92	SER	2.7
1	B	188	GLU	2.6
1	B	9	LEU	2.5
1	B	12	ALA	2.5
1	B	55	LEU	2.5
1	B	94	ASN	2.4
1	A	94	ASN	2.4
1	A	192	GLN	2.3
1	B	53	ARG	2.3
1	B	50	THR	2.3
1	A	126	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	58	THR	2.2
1	A	59	TYR	2.2
1	B	26	ALA	2.2
1	B	191	LYS	2.2
1	A	195	LEU	2.2
1	B	89	ILE	2.1
1	B	15	GLU	2.1
1	A	14	ARG	2.1
1	A	95	PRO	2.1
1	B	4	ILE	2.1
1	B	46	GLY	2.0
1	B	93	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 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.