



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 16, 2024 – 04:33 AM EDT

PDB ID : 1ZBO
Title : X-Ray Crystal Structure of Protein BPP1347 from Bordetella parapertussis.
Northeast Structural Genomics Consortium Target BoR27.
Authors : Forouhar, F.; Yong, W.; Conover, K.; Acton, T.B.; Montelione, G.T.; Tong,
L.; Hunt, J.F.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2005-04-08
Resolution : 2.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

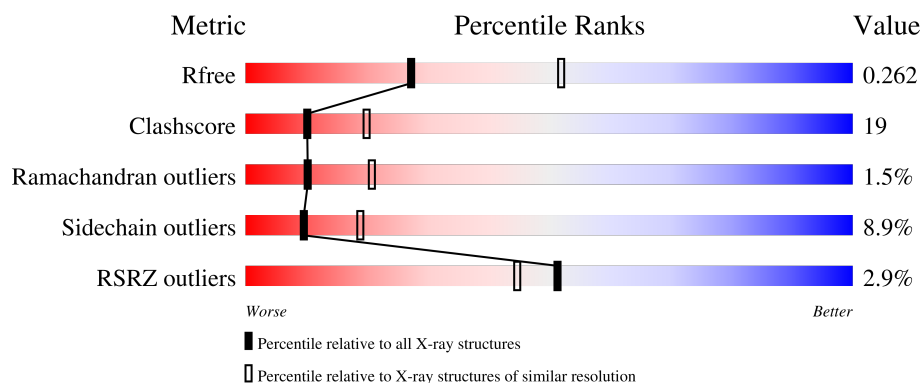
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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	210	<div> <div> <div></div> <div>57%</div> <div>33%</div> <div>6%</div> </div> </div>
1	B	210	<div> <div> <div>4%</div> <div>53%</div> <div>36%</div> <div>6%</div> <div>5%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3141 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 BPP1347.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	197	Total	C	N	O	S	Se	0	0	0
			1525	971	277	267	4	6			
1	B	200	Total	C	N	O	S	Se	0	0	0
			1547	983	282	272	4	6			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	31	MSE	MET	MODIFIED RESIDUE	UNP Q7WAM7
A	68	MSE	MET	MODIFIED RESIDUE	UNP Q7WAM7
A	78	MSE	MET	MODIFIED RESIDUE	UNP Q7WAM7
A	148	MSE	MET	MODIFIED RESIDUE	UNP Q7WAM7
A	150	MSE	MET	MODIFIED RESIDUE	UNP Q7WAM7
A	169	MSE	MET	MODIFIED RESIDUE	UNP Q7WAM7
A	203	LEU	-	CLONING ARTIFACT	UNP Q7WAM7
A	204	GLU	-	CLONING ARTIFACT	UNP Q7WAM7
A	205	HIS	-	EXPRESSION TAG	UNP Q7WAM7
A	206	HIS	-	EXPRESSION TAG	UNP Q7WAM7
A	207	HIS	-	EXPRESSION TAG	UNP Q7WAM7
A	208	HIS	-	EXPRESSION TAG	UNP Q7WAM7
A	209	HIS	-	EXPRESSION TAG	UNP Q7WAM7
A	210	HIS	-	EXPRESSION TAG	UNP Q7WAM7
B	31	MSE	MET	MODIFIED RESIDUE	UNP Q7WAM7
B	68	MSE	MET	MODIFIED RESIDUE	UNP Q7WAM7
B	78	MSE	MET	MODIFIED RESIDUE	UNP Q7WAM7
B	148	MSE	MET	MODIFIED RESIDUE	UNP Q7WAM7
B	150	MSE	MET	MODIFIED RESIDUE	UNP Q7WAM7
B	169	MSE	MET	MODIFIED RESIDUE	UNP Q7WAM7
B	203	LEU	-	CLONING ARTIFACT	UNP Q7WAM7
B	204	GLU	-	CLONING ARTIFACT	UNP Q7WAM7
B	205	HIS	-	EXPRESSION TAG	UNP Q7WAM7
B	206	HIS	-	EXPRESSION TAG	UNP Q7WAM7
B	207	HIS	-	EXPRESSION TAG	UNP Q7WAM7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	208	HIS	-	EXPRESSION TAG	UNP Q7WAM7
B	209	HIS	-	EXPRESSION TAG	UNP Q7WAM7
B	210	HIS	-	EXPRESSION TAG	UNP Q7WAM7

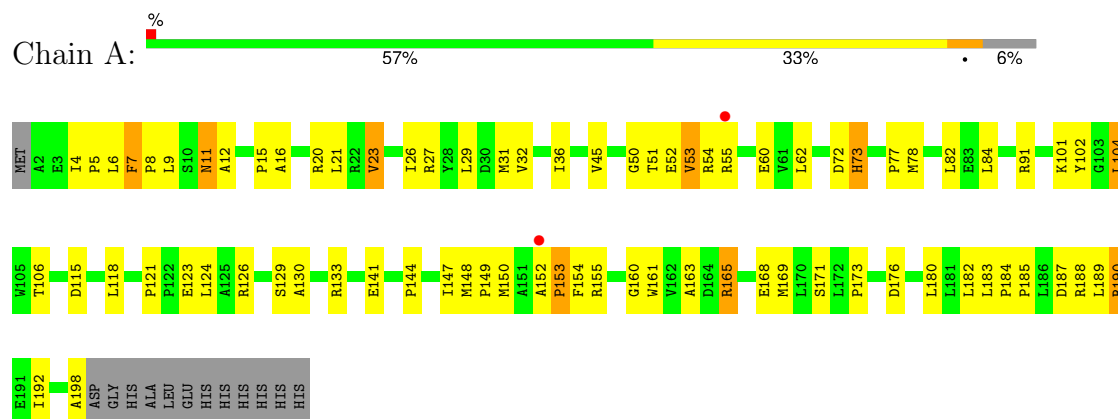
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	42	Total 42	O 42	0	0
2	B	27	Total 27	O 27	0	0

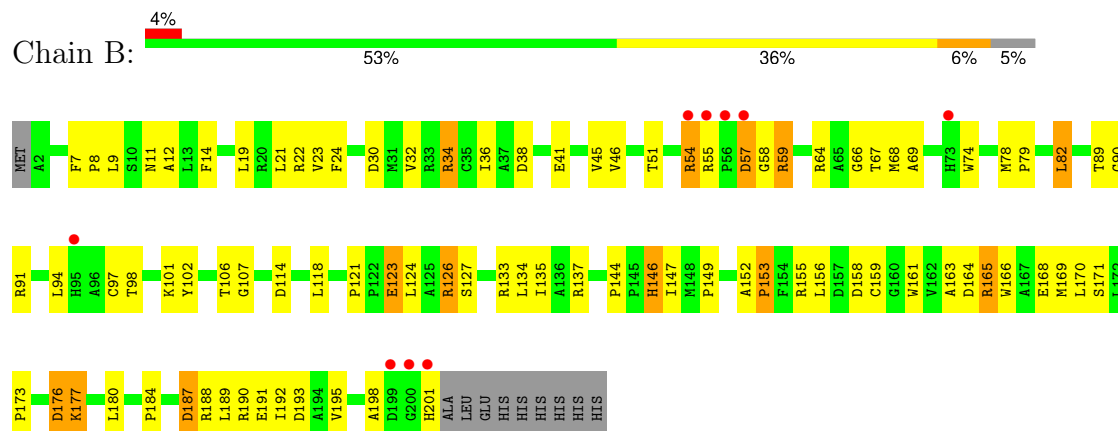
3 Residue-property plots

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: hypothetical protein BPP1347



• Molecule 1: hypothetical protein BPP1347



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	70.02Å 101.38Å 122.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.53 – 2.60 29.53 – 2.60	Depositor EDS
% Data completeness (in resolution range)	96.5 (29.53-2.60) 99.5 (29.53-2.60)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.78 (at 2.61Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.240 , 0.297 0.248 , 0.262	Depositor DCC
R_{free} test set	2425 reflections (9.42%)	wwPDB-VP
Wilson B-factor (Å ²)	77.5	Xtriage
Anisotropy	0.170	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 34.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3141	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.51% 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	0.45	0/1556	0.63	0/2110
1	B	0.37	0/1579	0.55	0/2141
All	All	0.41	0/3135	0.59	0/4251

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	1525	0	1549	57	0
1	B	1547	0	1563	70	0
2	A	42	0	0	5	0
2	B	27	0	0	2	0
All	All	3141	0	3112	117	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 19.

All (117) 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:68:MSE:HB3	1:B:89:THR:HG23	1.45	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:VAL:HG13	1:A:54:ARG:H	1.29	0.93
1:A:11:ASN:HD22	1:A:12:ALA:N	1.72	0.86
1:A:26:ILE:HA	1:A:29:LEU:HD23	1.58	0.86
1:B:68:MSE:HB3	1:B:89:THR:CG2	2.07	0.85
1:A:11:ASN:HD22	1:A:12:ALA:H	1.27	0.83
1:A:163:ALA:HA	1:A:192:ILE:HD13	1.62	0.80
1:A:53:VAL:HG13	1:A:54:ARG:N	2.01	0.75
1:B:173:PRO:HG2	1:B:176:ASP:HB2	1.69	0.74
1:A:5:PRO:HB2	1:A:31:MSE:HE2	1.69	0.73
1:B:184:PRO:HG2	1:B:187:ASP:HB2	1.72	0.72
1:A:72:ASP:O	1:A:73:HIS:HB3	1.90	0.72
1:A:183:LEU:HD12	1:A:184:PRO:HD2	1.72	0.71
1:A:23:VAL:HG13	1:A:82:LEU:HB3	1.74	0.69
1:B:11:ASN:HB2	2:B:221:HOH:O	1.93	0.69
1:A:115:ASP:HB2	2:A:241:HOH:O	1.93	0.68
1:B:121:PRO:HB2	1:B:123:GLU:HG3	1.77	0.67
1:A:163:ALA:HA	1:A:192:ILE:CD1	2.25	0.67
1:B:124:LEU:HD13	1:B:189:LEU:HD22	1.77	0.66
1:B:45:VAL:HG12	1:B:66:GLY:HA3	1.77	0.66
1:B:180:LEU:O	1:B:188:ARG:HD3	1.96	0.66
1:B:163:ALA:HB1	1:B:188:ARG:HG2	1.78	0.65
1:B:68:MSE:O	1:B:89:THR:HG22	1.98	0.64
1:A:133:ARG:NH2	1:B:198:ALA:HA	2.13	0.64
1:B:159:CYS:HA	1:B:189:LEU:HD12	1.81	0.63
1:B:41:GLU:HB2	1:B:69:ALA:O	1.99	0.62
1:A:45:VAL:HG11	1:A:62:LEU:HD13	1.80	0.62
1:A:51:THR:HG21	1:A:55:ARG:HG2	1.82	0.61
1:B:12:ALA:HB2	1:B:46:VAL:HB	1.83	0.61
1:A:198:ALA:C	1:B:133:ARG:HE	2.04	0.60
1:B:94:LEU:HD11	1:B:97:CYS:HB3	1.83	0.60
1:B:23:VAL:HG11	1:B:32:VAL:HG21	1.83	0.60
1:B:97:CYS:HB2	1:B:106:THR:O	2.02	0.59
1:B:9:LEU:HD11	1:B:21:LEU:HD11	1.85	0.59
1:A:144:PRO:HD2	1:A:147:ILE:HD12	1.86	0.58
1:B:189:LEU:O	1:B:189:LEU:HD23	2.04	0.58
1:A:169:MSE:HE2	1:B:171:SER:OG	2.04	0.58
1:A:11:ASN:ND2	1:A:12:ALA:N	2.49	0.57
1:A:124:LEU:HD11	1:A:190:ARG:HG2	1.86	0.57
1:B:78:MSE:HB2	1:B:79:PRO:HD2	1.87	0.57
1:B:32:VAL:O	1:B:36:ILE:HG13	2.05	0.56
1:B:14:PHE:HZ	1:B:177:LYS:HG3	1.71	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:94:LEU:HD11	1:B:107:GLY:HA3	1.87	0.56
1:A:152:ALA:HB3	1:A:153:PRO:HD3	1.87	0.56
1:B:176:ASP:O	1:B:180:LEU:HG	2.06	0.55
1:A:198:ALA:O	1:B:133:ARG:NE	2.41	0.54
1:B:19:LEU:HG	1:B:21:LEU:CD2	2.38	0.54
1:B:163:ALA:CB	1:B:188:ARG:HG2	2.38	0.53
1:B:19:LEU:HG	1:B:21:LEU:HD22	1.89	0.53
1:A:8:PRO:O	1:A:9:LEU:HD23	2.07	0.53
1:B:163:ALA:HA	1:B:192:ILE:HD13	1.91	0.52
1:B:191:GLU:O	1:B:195:VAL:HG23	2.09	0.52
1:B:59:ARG:HE	1:B:59:ARG:H	1.58	0.52
1:A:149:PRO:HG3	1:B:171:SER:HB3	1.92	0.52
1:A:6:LEU:O	1:A:31:MSE:HE1	2.10	0.51
1:A:77:PRO:HD2	2:A:231:HOH:O	2.11	0.51
1:B:152:ALA:HB3	1:B:153:PRO:HD3	1.92	0.51
1:B:67:THR:HA	1:B:91:ARG:HA	1.92	0.51
1:B:30:ASP:HB3	1:B:34:ARG:NH1	2.26	0.50
1:A:101:LYS:HE2	1:A:102:TYR:CE2	2.46	0.50
1:A:27:ARG:O	1:A:104:LEU:HD11	2.12	0.50
1:A:7:PHE:CE2	1:A:21:LEU:HD21	2.47	0.49
1:A:165:ARG:O	1:A:168:GLU:HG2	2.13	0.49
1:A:50:GLY:HA3	1:A:60:GLU:HG3	1.94	0.49
1:B:23:VAL:HG22	1:B:82:LEU:HB2	1.94	0.49
1:B:166:TRP:O	1:B:170:LEU:HG	2.12	0.49
1:A:155:ARG:HB2	1:A:161:TRP:CG	2.48	0.48
1:B:146:HIS:CD2	1:B:147:ILE:HG13	2.48	0.48
1:B:134:LEU:HD13	1:B:169:MSE:HE1	1.95	0.48
1:A:4:ILE:HB	1:A:5:PRO:HD2	1.95	0.48
1:B:54:ARG:CB	1:B:57:ASP:HB2	2.44	0.47
1:B:144:PRO:HB2	1:B:146:HIS:CD2	2.49	0.47
1:A:184:PRO:HG2	1:A:187:ASP:OD2	2.14	0.47
1:B:155:ARG:HB2	1:B:161:TRP:CG	2.49	0.47
1:B:126:ARG:HG2	1:B:127:SER:H	1.78	0.47
1:B:36:ILE:HG23	1:B:74:TRP:CZ3	2.49	0.47
1:B:177:LYS:HE3	2:B:222:HOH:O	2.14	0.47
1:A:72:ASP:O	1:A:73:HIS:CB	2.63	0.46
1:A:180:LEU:O	1:A:188:ARG:HD3	2.16	0.46
1:A:52:GLU:HB3	1:B:22:ARG:HH12	1.81	0.46
1:A:130:ALA:HA	1:A:133:ARG:NH2	2.30	0.46
1:B:54:ARG:HB2	1:B:57:ASP:HB2	1.98	0.45
1:A:171:SER:HB3	1:B:149:PRO:HG3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:24:PHE:CD1	1:B:24:PHE:N	2.84	0.45
1:B:101:LYS:HG2	1:B:102:TYR:CD1	2.52	0.45
1:B:135:ILE:HD12	1:B:156:LEU:HD21	1.99	0.45
1:B:11:ASN:OD1	1:B:12:ALA:N	2.49	0.45
1:A:91:ARG:HD3	2:A:211:HOH:O	2.17	0.45
1:B:97:CYS:HB3	1:B:107:GLY:HA3	1.99	0.45
1:B:118:LEU:O	1:B:158:ASP:HA	2.16	0.45
1:B:190:ARG:O	1:B:193:ASP:HB3	2.17	0.45
1:A:118:LEU:HD23	2:A:246:HOH:O	2.16	0.44
1:A:173:PRO:HG2	1:A:176:ASP:HB2	2.00	0.44
1:A:148:MSE:HE1	2:A:244:HOH:O	2.18	0.43
1:A:118:LEU:HD21	1:A:185:PRO:HB2	2.00	0.43
1:B:164:ASP:O	1:B:168:GLU:HG3	2.17	0.43
1:B:34:ARG:O	1:B:38:ASP:HB2	2.19	0.43
1:A:53:VAL:CG1	1:A:54:ARG:N	2.73	0.42
1:A:160:GLY:HA2	1:A:188:ARG:NH2	2.34	0.42
1:B:98:THR:N	1:B:106:THR:O	2.48	0.42
1:B:192:ILE:HG13	1:B:193:ASP:N	2.33	0.42
1:A:129:SER:HB2	1:B:201:HIS:O	2.20	0.42
1:B:89:THR:OG1	1:B:90:GLY:N	2.53	0.42
1:A:15:PRO:O	1:A:16:ALA:HB3	2.20	0.42
1:A:20:ARG:NH1	1:A:148:MSE:O	2.53	0.41
1:B:68:MSE:HB2	1:B:90:GLY:O	2.20	0.41
1:A:154:PHE:HA	1:A:161:TRP:CZ2	2.55	0.41
1:A:77:PRO:HG2	1:A:78:MSE:H	1.85	0.41
1:A:32:VAL:O	1:A:36:ILE:HG13	2.20	0.41
1:A:188:ARG:HG3	1:A:188:ARG:HH11	1.86	0.41
1:B:8:PRO:HA	1:B:45:VAL:O	2.21	0.41
1:A:176:ASP:HB3	1:B:137:ARG:HH11	1.86	0.41
1:B:165:ARG:O	1:B:168:GLU:HB2	2.20	0.41
1:A:198:ALA:O	1:B:133:ARG:HG2	2.22	0.40
1:B:68:MSE:C	1:B:89:THR:HG22	2.41	0.40
1:A:106:THR:O	1:A:106:THR:HG23	2.20	0.40
1:A:121:PRO:HB2	1:A:123:GLU:OE1	2.21	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	195/210 (93%)	180 (92%)	12 (6%)	3 (2%)	10	21
1	B	198/210 (94%)	174 (88%)	21 (11%)	3 (2%)	10	21
All	All	393/420 (94%)	354 (90%)	33 (8%)	6 (2%)	10	21

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	53	VAL
1	A	73	HIS
1	B	57	ASP
1	A	153	PRO
1	B	153	PRO
1	B	58	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	157/162 (97%)	145 (92%)	12 (8%)	13	26
1	B	159/162 (98%)	143 (90%)	16 (10%)	7	14
All	All	316/324 (98%)	288 (91%)	28 (9%)	9	19

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	PHE
1	A	11	ASN
1	A	23	VAL
1	A	84	LEU
1	A	104	LEU
1	A	126	ARG
1	A	141	GLU
1	A	150	MSE
1	A	165	ARG
1	A	182	LEU
1	A	189	LEU
1	A	190	ARG
1	B	7	PHE
1	B	34	ARG
1	B	51	THR
1	B	54	ARG
1	B	55	ARG
1	B	59	ARG
1	B	64	ARG
1	B	82	LEU
1	B	114	ASP
1	B	123	GLU
1	B	126	ARG
1	B	146	HIS
1	B	165	ARG
1	B	176	ASP
1	B	177	LYS
1	B	187	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	108	GLN
1	B	146	HIS

5.3.3 RNA ⓘ

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 monosaccharides 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	191/210 (90%)	-0.27	2 (1%) 82 80	29, 46, 77, 90	0
1	B	194/210 (92%)	0.14	9 (4%) 32 26	37, 64, 85, 104	0
All	All	385/420 (91%)	-0.06	11 (2%) 51 45	29, 54, 84, 104	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	200	GLY	5.3
1	B	54	ARG	4.4
1	B	201	HIS	4.3
1	B	55	ARG	3.8
1	A	152	ALA	3.0
1	B	199	ASP	2.8
1	B	56	PRO	2.6
1	B	95	HIS	2.5
1	B	57	ASP	2.4
1	B	73	HIS	2.1
1	A	55	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 monosaccharides 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.