



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:23 AM UTC

PDB ID : 2GEM / pdb\_00002gem  
Title : 2.1A crystal structure of Salmonella typhimurium YeaZ, a putative Gram-negative RPF, form-A  
Authors : Nichols, C.E.; Stammers, D.K.  
Deposited on : 2006-03-20  
Resolution : 2.10 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](mailto: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>.

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

MolProbity : 4-5-2 with Phenix2.0  
Xtrriage (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

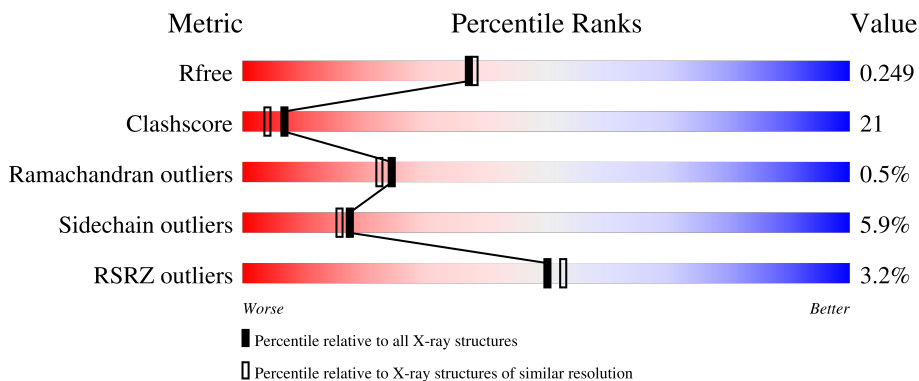
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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  $\geq 3$ , 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	231	 3% 66% 23% 5% 6%
1	B	231	 3% 63% 27% 6% 6%

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3625 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 Putative Gram negative resuscitation promoting factor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	216	1611	1014	280	308	9	0	0	0
1	B	216	1611	1014	280	308	9	0	0	0

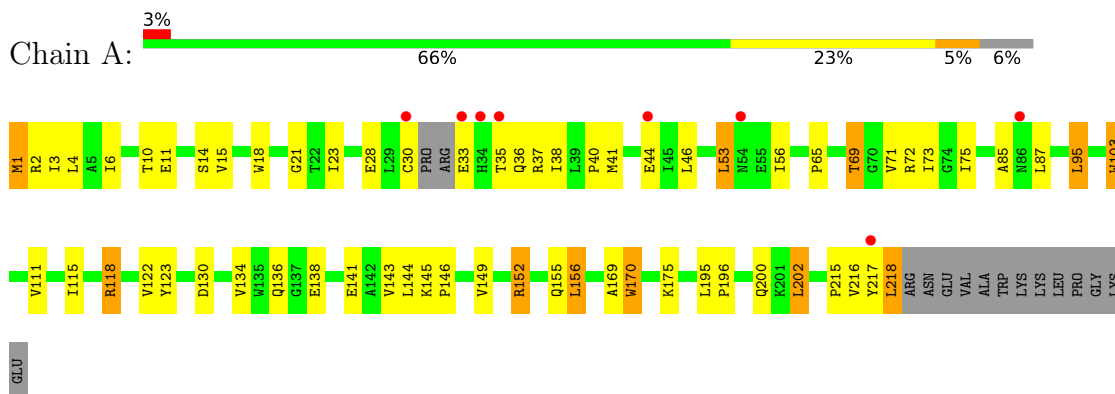
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	209	209	209	0	0
2	B	194	194	194	0	0

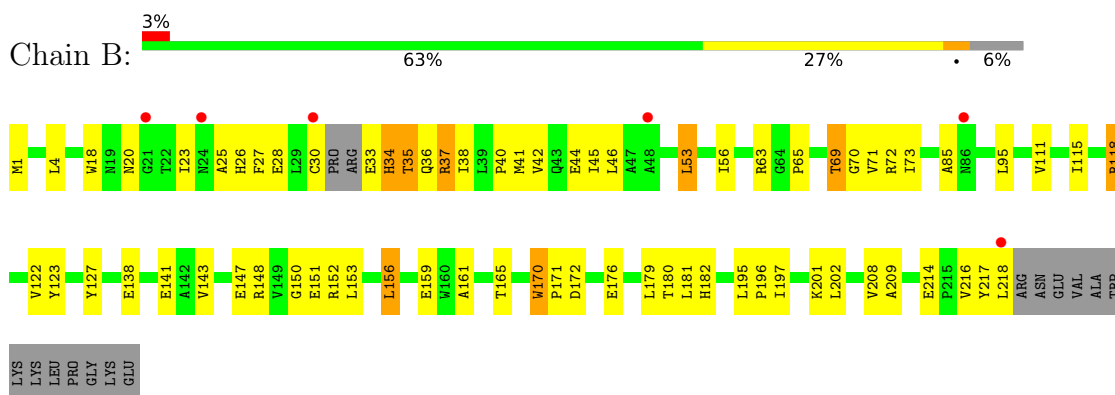
### 3 Residue-property plots [i](#)

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: Putative Gram negative resuscitation promoting factor



- Molecule 1: Putative Gram negative resuscitation promoting factor



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	40.70Å 40.70Å 464.83Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.11 – 2.10 29.11 – 2.10	Depositor EDS
% Data completeness (in resolution range)	84.1 (29.11-2.10) 84.2 (29.11-2.10)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.42 (at 2.10Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.220 , 0.248 0.220 , 0.249	Depositor DCC
$R_{free}$ test set	2323 reflections (9.86%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	37.8	Xtrriage
Anisotropy	0.237	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 71.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.42$ , $\langle L^2 \rangle = 0.25$	Xtrriage
Estimated twinning fraction	0.097 for -h,-k,l	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3625	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	42.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.05% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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 [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	0.41	0/1641	0.87	4/2232 (0.2%)
1	B	0.40	0/1641	0.92	5/2232 (0.2%)
All	All	0.40	0/3282	0.90	9/4464 (0.2%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	170	TRP	CA-C-N	6.26	127.66	119.84
1	B	170	TRP	C-N-CA	6.26	127.66	119.84
1	B	214	GLU	N-CA-C	6.04	113.57	108.13
1	A	170	TRP	CA-C-N	5.92	127.24	119.84
1	A	170	TRP	C-N-CA	5.92	127.24	119.84
1	A	111	VAL	N-CA-C	5.86	117.19	108.46
1	B	111	VAL	N-CA-C	5.67	116.91	108.46
1	B	25	ALA	N-CA-C	5.21	117.73	109.50
1	A	14	SER	N-CA-C	5.06	116.88	109.24

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	1611	0	1592	68	0
1	B	1611	0	1592	68	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	209	0	0	3	0
2	B	194	0	0	8	0
All	All	3625	0	3184	136	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 21.

All (136) 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:118:ARG:HH11	1:A:118:ARG:HB3	1.27	0.96
1:B:118:ARG:HH11	1:B:118:ARG:HB3	1.39	0.86
1:B:28:GLU:OE1	1:B:37:ARG:HD3	1.76	0.85
1:A:28:GLU:OE1	1:A:37:ARG:HD3	1.85	0.77
1:B:35:THR:HA	1:B:38:ILE:HD13	1.67	0.77
1:A:35:THR:HA	1:A:38:ILE:HD13	1.68	0.75
1:A:33:GLU:HG2	1:A:36:GLN:HE21	1.53	0.73
1:B:115:ILE:HB	1:B:123:TYR:HB2	1.69	0.73
1:A:134:VAL:HG13	1:A:200:GLN:HE22	1.55	0.71
1:A:33:GLU:HG2	1:A:36:GLN:NE2	2.06	0.70
1:A:152:ARG:HH11	1:A:156:LEU:HD11	1.58	0.68
1:A:152:ARG:HH11	1:A:156:LEU:CD1	2.08	0.67
1:A:195:LEU:HB2	1:A:196:PRO:HD3	1.79	0.65
1:A:75:ILE:HD12	1:A:215:PRO:HG3	1.80	0.64
1:B:179:LEU:HB2	1:B:181:LEU:HD21	1.79	0.64
1:B:35:THR:CA	1:B:38:ILE:HD13	2.28	0.64
1:A:46:LEU:HD11	1:A:53:LEU:HD23	1.80	0.63
1:A:53:LEU:HD12	1:A:85:ALA:CB	2.29	0.63
1:A:33:GLU:HG2	1:A:36:GLN:HG3	1.81	0.63
1:A:40:PRO:O	1:A:44:GLU:HG3	1.99	0.62
1:A:1:MET:HE1	1:A:3:ILE:HD11	1.81	0.62
1:B:37:ARG:O	1:B:40:PRO:HG2	2.00	0.61
1:A:1:MET:CE	1:A:202:LEU:HD12	2.30	0.61
1:B:197:ILE:O	1:B:201:LYS:HG3	2.01	0.61
1:B:150:GLY:HA2	1:B:153:LEU:HD12	1.84	0.60
1:A:138:GLU:O	1:A:141:GLU:HG3	2.02	0.60
1:A:33:GLU:OE2	1:A:35:THR:HG22	2.02	0.60
1:B:4:LEU:HB2	1:B:56:ILE:HD13	1.84	0.60
1:B:118:ARG:HB3	1:B:118:ARG:NH1	2.11	0.60
1:A:38:ILE:HD12	1:A:38:ILE:N	2.16	0.60
1:B:18:TRP:CE3	1:B:23:ILE:HD11	2.38	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:38:ILE:HD12	1:B:38:ILE:H	1.67	0.59
1:B:35:THR:HG23	1:B:36:GLN:H	1.68	0.58
1:B:18:TRP:HB2	1:B:195:LEU:HD13	1.85	0.58
1:A:37:ARG:O	1:A:41:MET:HG3	2.04	0.58
1:A:118:ARG:HB3	1:A:118:ARG:NH1	2.08	0.58
1:A:1:MET:HE2	1:A:202:LEU:HD12	1.87	0.57
1:B:217:TYR:HB3	2:B:278:HOH:O	2.04	0.57
1:A:4:LEU:HB2	1:A:56:ILE:HD13	1.86	0.57
1:B:65:PRO:HD2	1:B:95:LEU:CD2	2.34	0.57
1:B:152:ARG:NH1	1:B:156:LEU:HD13	2.19	0.56
1:B:143:VAL:O	1:B:143:VAL:HG13	2.05	0.56
1:B:138:GLU:O	1:B:141:GLU:HG3	2.06	0.56
1:A:71:VAL:HG13	1:A:72:ARG:N	2.20	0.56
1:B:38:ILE:HD12	1:B:38:ILE:N	2.20	0.56
1:A:218:LEU:HD12	1:A:218:LEU:N	2.21	0.55
1:A:115:ILE:HB	1:A:123:TYR:HB2	1.88	0.55
1:A:118:ARG:HH11	1:A:118:ARG:CB	2.11	0.55
1:A:152:ARG:NH1	1:A:156:LEU:CD1	2.70	0.54
1:A:11:GLU:CD	1:A:11:GLU:H	2.15	0.54
1:A:33:GLU:CG	1:A:36:GLN:HE21	2.18	0.54
1:B:63:ARG:HH22	1:B:138:GLU:CD	2.15	0.54
1:A:122:VAL:HG11	1:A:149:VAL:HG21	1.89	0.54
1:A:6:ILE:HG12	1:A:15:VAL:HG22	1.89	0.54
1:B:63:ARG:NH1	2:B:332:HOH:O	2.40	0.54
1:A:21:GLY:O	1:A:23:ILE:HG13	2.08	0.53
1:A:33:GLU:HG2	1:A:36:GLN:CG	2.38	0.53
1:B:216:VAL:HG22	1:B:218:LEU:HD12	1.89	0.52
1:A:37:ARG:O	1:A:40:PRO:HG2	2.09	0.52
1:B:23:ILE:HD12	1:B:195:LEU:HD13	1.92	0.52
1:B:65:PRO:HD2	1:B:95:LEU:HD22	1.92	0.52
1:B:195:LEU:HB2	1:B:196:PRO:HD3	1.90	0.52
1:B:69:THR:O	1:B:73:ILE:HG12	2.09	0.52
1:B:35:THR:HG23	1:B:36:GLN:N	2.25	0.51
1:A:38:ILE:HD12	1:A:38:ILE:H	1.75	0.51
1:B:53:LEU:HD12	1:B:85:ALA:CB	2.41	0.51
1:B:218:LEU:HD12	1:B:218:LEU:N	2.25	0.51
1:A:152:ARG:HA	1:A:155:GLN:OE1	2.10	0.51
1:A:169:ALA:HB2	2:A:312:HOH:O	2.11	0.51
1:A:217:TYR:C	1:A:218:LEU:HD12	2.36	0.51
1:A:69:THR:O	1:A:73:ILE:HG12	2.10	0.51
1:B:23:ILE:HD12	1:B:195:LEU:CD1	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2:ARG:HA	1:A:18:TRP:O	2.12	0.50
1:A:35:THR:CA	1:A:38:ILE:HD13	2.40	0.50
1:B:34:HIS:O	1:B:36:GLN:N	2.45	0.50
1:B:71:VAL:HG13	1:B:72:ARG:N	2.27	0.49
1:B:181:LEU:HD22	1:B:181:LEU:N	2.26	0.49
1:A:65:PRO:HD2	1:A:95:LEU:HD22	1.94	0.49
1:A:72:ARG:HH11	1:A:72:ARG:HG3	1.78	0.49
1:B:38:ILE:H	1:B:38:ILE:CD1	2.25	0.49
1:B:41:MET:O	1:B:45:ILE:HD13	2.13	0.49
1:B:179:LEU:HB2	1:B:181:LEU:CD2	2.41	0.49
1:B:26:HIS:HB2	2:B:404:HOH:O	2.12	0.48
1:A:143:VAL:HG13	1:A:143:VAL:O	2.12	0.48
1:A:134:VAL:HG12	1:A:136:GLN:NE2	2.29	0.48
1:B:216:VAL:HG22	1:B:218:LEU:CD1	2.44	0.48
1:A:152:ARG:NH1	1:A:156:LEU:HD11	2.26	0.47
1:B:70:GLY:HA3	2:B:419:HOH:O	2.15	0.47
1:B:40:PRO:O	1:B:44:GLU:HG3	2.14	0.47
1:A:53:LEU:HD12	1:A:85:ALA:HB2	1.97	0.47
1:A:1:MET:HE2	1:A:202:LEU:CD1	2.44	0.47
1:A:130:ASP:OD1	1:A:130:ASP:C	2.58	0.47
1:B:37:ARG:HA	1:B:37:ARG:HE	1.79	0.46
1:B:27:PHE:CD1	1:B:27:PHE:C	2.94	0.46
1:B:42:VAL:O	1:B:46:LEU:HG	2.15	0.46
1:A:10:THR:HG22	2:A:292:HOH:O	2.17	0.45
1:A:38:ILE:H	1:A:38:ILE:CD1	2.29	0.45
1:A:38:ILE:N	1:A:38:ILE:CD1	2.80	0.44
1:A:134:VAL:HG13	1:A:200:GLN:NE2	2.29	0.44
1:B:127:TYR:OH	1:B:138:GLU:HA	2.17	0.44
1:B:143:VAL:HG12	2:B:314:HOH:O	2.17	0.44
1:A:71:VAL:CG1	1:A:72:ARG:N	2.80	0.44
1:A:134:VAL:HG12	1:A:136:GLN:HE21	1.81	0.44
1:B:218:LEU:N	1:B:218:LEU:CD1	2.81	0.44
1:B:53:LEU:HD22	1:B:53:LEU:HA	1.87	0.44
1:A:216:VAL:HG22	1:A:218:LEU:CD1	2.47	0.43
1:A:53:LEU:HD13	1:A:87:LEU:HD11	2.01	0.43
1:B:18:TRP:HA	1:B:23:ILE:HD13	2.00	0.43
1:B:118:ARG:HH11	1:B:118:ARG:CB	2.21	0.43
1:B:147:GLU:HG2	2:B:279:HOH:O	2.17	0.43
1:A:1:MET:HE1	1:A:3:ILE:CD1	2.47	0.43
1:A:103:TRP:CD1	1:A:103:TRP:C	2.96	0.43
1:B:33:GLU:CD	1:B:36:GLN:NE2	2.76	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:46:LEU:HD11	1:B:53:LEU:HD23	2.01	0.42
1:B:33:GLU:CD	1:B:36:GLN:HE21	2.28	0.42
1:A:2:ARG:NH1	2:A:435:HOH:O	2.52	0.42
1:A:122:VAL:HG12	1:A:144:LEU:O	2.18	0.42
1:B:208:VAL:CG1	1:B:209:ALA:N	2.82	0.42
1:A:145:LYS:O	1:A:146:PRO:C	2.62	0.42
1:B:1:MET:N	1:B:20:ASN:OD1	2.52	0.42
1:B:44:GLU:HB3	2:B:241:HOH:O	2.20	0.42
1:B:161:ALA:HA	1:B:182:HIS:O	2.20	0.41
1:B:165:THR:HB	2:B:307:HOH:O	2.20	0.41
1:B:172:ASP:O	1:B:176:GLU:HG2	2.21	0.41
1:A:1:MET:HE1	1:A:202:LEU:HD12	2.01	0.41
1:B:152:ARG:NH1	1:B:156:LEU:CD1	2.83	0.41
1:B:34:HIS:O	1:B:35:THR:C	2.64	0.41
1:B:159:GLU:HB3	1:B:180:THR:HB	2.02	0.41
1:B:208:VAL:HG12	1:B:209:ALA:N	2.35	0.41
1:A:3:ILE:HB	1:A:18:TRP:HB3	2.03	0.41
1:B:148:ARG:O	1:B:151:GLU:HB3	2.21	0.41
1:B:170:TRP:HA	1:B:171:PRO:HD2	1.91	0.41
1:A:72:ARG:HG3	1:A:72:ARG:NH1	2.35	0.40
1:A:146:PRO:HG2	1:A:170:TRP:CZ2	2.56	0.40
1:A:175:LYS:HB2	1:A:175:LYS:HE2	1.89	0.40
1:B:38:ILE:N	1:B:38:ILE:CD1	2.85	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	212/231 (92%)	203 (96%)	9 (4%)	0	100 100
1	B	212/231 (92%)	203 (96%)	7 (3%)	2 (1%)	14 10

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	424/462 (92%)	406 (96%)	16 (4%)	2 (0%)	24 22

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	35	THR
1	B	34	HIS

### 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	160/173 (92%)	149 (93%)	11 (7%)	14 12
1	B	160/173 (92%)	152 (95%)	8 (5%)	22 22
All	All	320/346 (92%)	301 (94%)	19 (6%)	18 16

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	30	CYS
1	A	53	LEU
1	A	69	THR
1	A	95	LEU
1	A	103	TRP
1	A	118	ARG
1	A	152	ARG
1	A	156	LEU
1	A	202	LEU
1	A	218	LEU
1	B	30	CYS
1	B	37	ARG
1	B	53	LEU
1	B	69	THR
1	B	118	ARG

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Mol	Chain	Res	Type
1	B	122	VAL
1	B	156	LEU
1	B	202	LEU

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	24	ASN
1	A	36	GLN
1	A	132	GLN
1	A	136	GLN
1	A	182	HIS
1	A	200	GLN
1	B	36	GLN
1	B	132	GLN
1	B	155	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

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	216/231 (93%)	0.18	8 (3%) 45 47	22, 36, 66, 98	0
1	B	216/231 (93%)	0.23	6 (2%) 55 57	19, 38, 66, 109	0
All	All	432/462 (93%)	0.21	14 (3%) 50 53	19, 38, 66, 109	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	35	THR	4.0
1	B	30	CYS	3.9
1	A	30	CYS	3.6
1	A	86	ASN	2.7
1	B	48	ALA	2.6
1	A	34	HIS	2.6
1	B	24	ASN	2.5
1	A	217	TYR	2.5
1	A	33	GLU	2.5
1	B	218	LEU	2.5
1	B	21	GLY	2.3
1	B	86	ASN	2.2
1	A	44	GLU	2.1
1	A	54	ASN	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

There are no ligands in this entry.

## 6.5 Other polymers

There are no such residues in this entry.