



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 1, 2021 – 04:08 PM EDT

PDB ID : 1XTM  
Title : Crystal structure of the double mutant Y88H-P104H of a SOD-like protein from *Bacillus subtilis*.  
Authors : Calderone, V.; Mangani, S.; Banci, L.; Benvenuti, M.; Bertini, I.; Fantoni, A.; Viezzoli, M.S.  
Deposited on : 2004-10-22  
Resolution : 1.60 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](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.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.23.2
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)	:	2.23.2

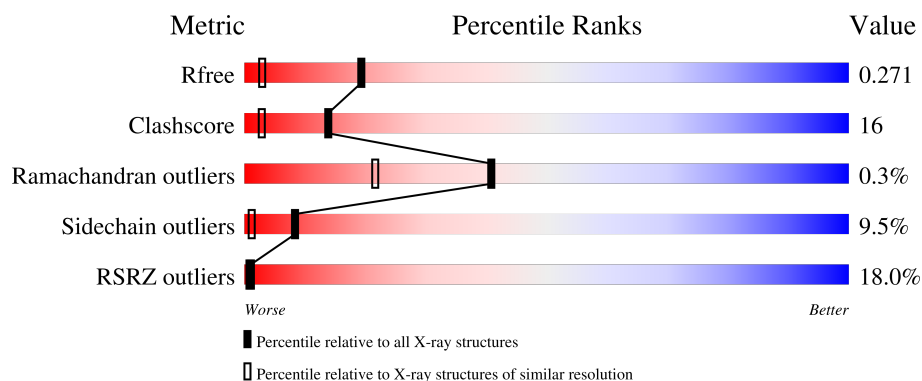
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.60 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	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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  $\geq 3$ , 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	175	<div> <div>25%</div> <div>60%</div> <div>22%</div> <div>• •</div> <div>13%</div> </div>
1	B	175	<div> <div>6%</div> <div>63%</div> <div>20%</div> <div>•</div> <div>13%</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2507 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 superoxide dismutase-like protein yojM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	153	Total	C	N	O	S	0	0	0
			1140	705	206	225	4			
1	A	152	Total	C	N	O	S	0	0	0
			1134	702	205	223	4			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	88	HIS	TYR	engineered mutation	UNP O31851
B	104	HIS	PRO	engineered mutation	UNP O31851
A	88	HIS	TYR	engineered mutation	UNP O31851
A	104	HIS	PRO	engineered mutation	UNP O31851

- Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Cu	0	0
			1	1		
2	A	1	Total	Cu	0	0
			1	1		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	4	Total	Zn	0	0
			4	4		
3	A	1	Total	Zn	0	0
			1	1		

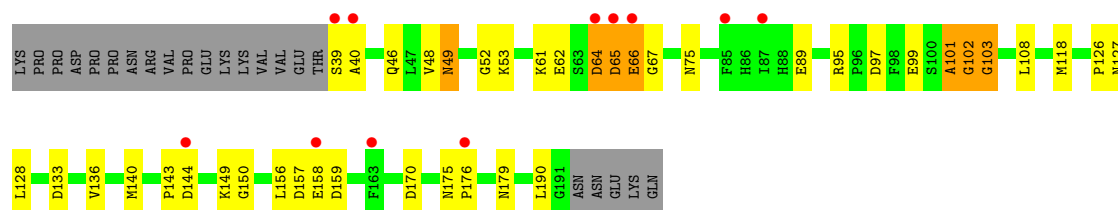
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	142	Total 142	O 142	0	0
4	A	84	Total 84	O 84	0	0

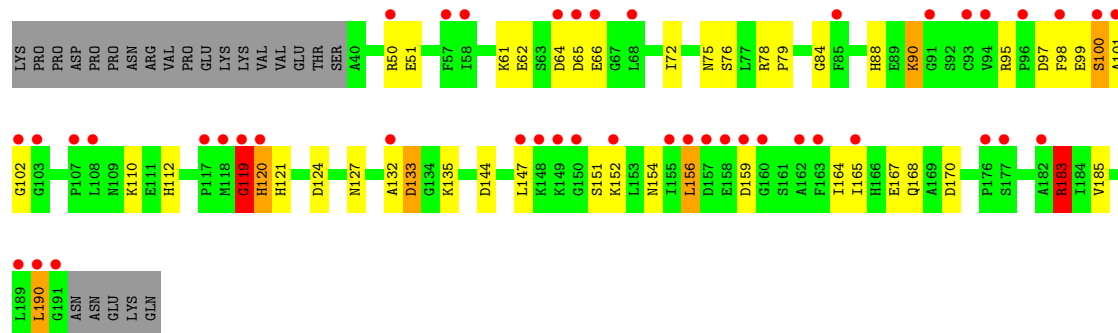
### 3 Residue-property plots [i](#)

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: Hypothetical superoxide dismutase-like protein yojM



- Molecule 1: Hypothetical superoxide dismutase-like protein yojM



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	52.46Å 104.35Å 58.76Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.60 – 1.60 25.60 – 1.60	Depositor EDS
% Data completeness (in resolution range)	100.0 (25.60-1.60) 99.5 (25.60-1.60)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.78 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.2.0003	Depositor
R, $R_{free}$	0.253 , 0.266 0.260 , 0.271	Depositor DCC
$R_{free}$ test set	6323 reflections (7.72%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	20.2	Xtriage
Anisotropy	0.495	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 51.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.44$ , $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	2507	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.23% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CU, ZN

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	1.11	2/1158 (0.2%)	1.08	10/1563 (0.6%)
1	B	0.93	2/1164 (0.2%)	1.06	7/1571 (0.4%)
All	All	1.03	4/2322 (0.2%)	1.07	17/3134 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	3
All	All	0	5

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	119	GLY	C-N	-25.11	0.76	1.34
1	A	120	HIS	C-N	19.14	1.78	1.34
1	B	143	PRO	C-N	13.94	1.66	1.34
1	B	144	ASP	C-N	7.17	1.50	1.34

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	119	GLY	O-C-N	-11.82	103.78	122.70
1	A	120	HIS	CA-C-N	-10.75	93.55	117.20
1	A	119	GLY	C-N-CA	10.40	147.71	121.70
1	A	120	HIS	CA-C-O	9.39	139.82	120.10
1	B	102	GLY	N-CA-C	-8.86	90.94	113.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	119	GLY	CA-C-N	8.80	136.56	117.20
1	A	124	ASP	CB-CG-OD2	7.40	124.96	118.30
1	B	103	GLY	N-CA-C	-7.05	95.46	113.10
1	B	101	ALA	N-CA-C	6.80	129.37	111.00
1	A	190	LEU	CA-CB-CG	6.56	130.39	115.30
1	A	65	ASP	CB-CG-OD2	6.42	124.07	118.30
1	B	157	ASP	CB-CG-OD2	6.24	123.92	118.30
1	B	158	GLU	C-N-CA	5.70	135.96	121.70
1	A	144	ASP	CB-CG-OD2	5.31	123.08	118.30
1	A	183	ARG	CB-CA-C	5.31	121.02	110.40
1	B	144	ASP	CB-CA-C	5.30	121.01	110.40
1	B	64	ASP	CB-CG-OD2	5.16	122.95	118.30

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	119	GLY	Peptide
1	A	132	ALA	Peptide
1	B	101	ALA	Peptide
1	B	102	GLY	Peptide
1	B	190	LEU	Peptide

## 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	1134	0	1074	47	0
1	B	1140	0	1081	23	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
3	B	4	0	0	0	0
4	A	84	0	0	7	0
4	B	142	0	0	3	0
All	All	2507	0	2155	70	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 16.

All (70) 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:120:HIS:C	1:A:121:HIS:N	1.78	1.37
1:A:119:GLY:O	1:A:120:HIS:N	1.57	1.26
1:A:119:GLY:C	1:A:120:HIS:CA	2.14	1.15
1:A:119:GLY:CA	1:A:120:HIS:N	2.15	1.09
1:A:183:ARG:HG3	4:A:540:HOH:O	1.56	1.04
1:A:101:ALA:HB1	1:A:102:GLY:HA2	1.46	0.97
1:A:99:GLU:O	1:A:101:ALA:HA	1.65	0.96
1:A:120:HIS:CA	1:A:121:HIS:N	2.41	0.83
1:A:95:ARG:HG3	4:A:544:HOH:O	1.80	0.81
1:A:119:GLY:C	1:A:120:HIS:N	0.76	0.81
1:A:97:ASP:OD2	1:A:99:GLU:HG3	1.88	0.74
1:A:127:ASN:HD21	1:A:170:ASP:CB	2.00	0.73
1:A:154:ASN:ND2	1:A:156:LEU:O	2.22	0.72
1:A:51:GLU:HB3	4:A:577:HOH:O	1.93	0.69
1:A:127:ASN:HD21	1:A:170:ASP:HB3	1.59	0.66
1:A:102:GLY:O	1:A:183:ARG:NH1	2.28	0.66
1:A:133:ASP:HB3	1:A:135:LYS:H	1.61	0.66
1:B:118:MET:HA	1:B:118:MET:CE	2.26	0.65
1:A:167:GLU:HG3	1:A:168:GLN:HG3	1.81	0.63
1:B:66:GLU:HG2	1:B:67:GLY:N	2.14	0.62
1:A:154:ASN:CG	1:A:156:LEU:O	2.39	0.61
1:B:127:ASN:HD21	1:B:170:ASP:HB3	1.65	0.61
1:A:97:ASP:OD2	1:A:99:GLU:CG	2.48	0.60
1:B:49:ASN:C	1:B:49:ASN:HD22	2.06	0.59
1:A:120:HIS:HA	1:A:121:HIS:N	2.15	0.59
1:B:62:GLU:HG2	4:B:525:HOH:O	2.04	0.58
1:B:118:MET:HA	1:B:118:MET:HE2	1.84	0.57
1:B:65:ASP:OD2	1:B:66:GLU:HB2	2.04	0.57
1:A:101:ALA:HB1	1:A:102:GLY:CA	2.29	0.57
1:A:164:ILE:HG21	1:A:183:ARG:HB3	1.87	0.56
1:A:88:HIS:CG	1:A:164:ILE:HD13	2.41	0.56
1:B:39:SER:O	1:B:40:ALA:HB3	2.08	0.54
1:A:90:LYS:HD3	1:A:159:ASP:OD2	2.07	0.54
1:A:151:SER:HB3	1:A:154:ASN:HB2	1.89	0.53
1:B:66:GLU:HG2	1:B:67:GLY:H	1.72	0.53
1:A:127:ASN:HD21	1:A:170:ASP:H	1.55	0.53
1:A:110:LYS:NZ	4:A:582:HOH:O	2.41	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:GLY:HA2	1:A:127:ASN:HD22	1.76	0.51
1:A:51:GLU:CB	4:A:577:HOH:O	2.55	0.51
1:A:127:ASN:ND2	1:A:170:ASP:H	2.08	0.50
1:A:98:PHE:C	1:A:100:SER:H	2.14	0.50
1:A:62:GLU:CG	1:A:147:LEU:HD12	2.41	0.49
1:A:112:HIS:HB2	1:A:121:HIS:CE1	2.47	0.49
1:A:95:ARG:N	4:A:544:HOH:O	2.46	0.49
1:A:99:GLU:C	1:A:101:ALA:HA	2.32	0.49
1:A:50:ARG:NH1	4:A:533:HOH:O	2.22	0.48
1:A:101:ALA:CB	1:A:102:GLY:HA2	2.30	0.46
1:B:150:GLY:HA2	4:B:638:HOH:O	2.15	0.46
1:B:126:PRO:HD2	1:B:140:MET:SD	2.56	0.46
1:A:90:LYS:HB2	1:A:101:ALA:HB3	1.97	0.46
1:B:89:GLU:HG2	1:B:103:GLY:H	1.80	0.46
1:B:97:ASP:OD1	1:B:99:GLU:OE1	2.33	0.46
1:B:48:VAL:HG22	4:B:543:HOH:O	2.17	0.44
1:A:88:HIS:CD2	1:A:164:ILE:HD13	2.53	0.44
1:B:64:ASP:HA	1:B:65:ASP:OD1	2.18	0.43
1:A:62:GLU:HG3	1:A:147:LEU:HD12	2.01	0.43
1:B:66:GLU:CG	1:B:67:GLY:N	2.78	0.43
1:B:89:GLU:CD	1:B:103:GLY:H	2.21	0.43
1:A:165:ILE:HB	1:A:185:VAL:HB	2.01	0.43
1:B:127:ASN:HD21	1:B:170:ASP:CB	2.30	0.42
1:B:49:ASN:ND2	1:B:53:LYS:H	2.17	0.42
1:B:39:SER:O	1:B:40:ALA:CB	2.67	0.42
1:A:78:ARG:HA	1:A:79:PRO:HD3	1.87	0.42
1:A:72:ILE:HD11	1:A:165:ILE:HD11	2.03	0.41
1:A:98:PHE:C	1:A:100:SER:N	2.73	0.41
1:B:128:LEU:HD22	1:B:136:VAL:HG21	2.03	0.41
1:B:52:GLY:HA2	1:B:95:ARG:HD3	2.03	0.41
1:A:127:ASN:ND2	1:A:170:ASP:HB3	2.31	0.41
1:B:175:ASN:HA	1:B:176:PRO:HA	1.79	0.40
1:A:127:ASN:HD21	1:A:170:ASP:HB2	1.81	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	150/175 (86%)	136 (91%)	13 (9%)	1 (1%)	22	7
1	B	151/175 (86%)	141 (93%)	10 (7%)	0	100	100
All	All	301/350 (86%)	277 (92%)	23 (8%)	1 (0%)	41	21

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	133	ASP

### 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	120/143 (84%)	109 (91%)	11 (9%)	9	1
1	B	121/143 (85%)	109 (90%)	12 (10%)	8	1
All	All	241/286 (84%)	218 (90%)	23 (10%)	8	1

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	46	GLN
1	B	49	ASN
1	B	61	LYS
1	B	65	ASP
1	B	66	GLU

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Mol	Chain	Res	Type
1	B	75	ASN
1	B	108	LEU
1	B	133	ASP
1	B	149	LYS
1	B	156	LEU
1	B	159	ASP
1	B	179	ASN
1	A	61	LYS
1	A	64	ASP
1	A	66	GLU
1	A	75	ASN
1	A	76	SER
1	A	90	LYS
1	A	100	SER
1	A	152	LYS
1	A	156	LEU
1	A	183	ARG
1	A	190	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	49	ASN
1	B	75	ASN
1	B	127	ASN
1	A	75	ASN
1	A	127	ASN
1	A	179	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 7 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	2
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	120:HIS	C	121:HIS	N	1.78
1	B	143:PRO	C	144:ASP	N	1.66
1	A	119:GLY	C	120:HIS	N	0.76

## 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, 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	152/175 (86%)	1.59	44 (28%) 0 0	17, 32, 48, 51	0
1	B	153/175 (87%)	0.55	11 (7%) 15 14	11, 17, 30, 43	0
All	All	305/350 (87%)	1.07	55 (18%) 1 1	11, 25, 44, 51	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	65	ASP	8.0
1	A	108	LEU	6.8
1	B	64	ASP	6.6
1	A	64	ASP	6.4
1	A	118	MET	5.9
1	B	65	ASP	5.4
1	B	39	SER	5.3
1	A	101	ALA	4.7
1	A	190	LEU	4.3
1	A	150	GLY	4.2
1	A	66	GLU	4.1
1	A	102	GLY	4.0
1	A	156	LEU	4.0
1	A	158	GLU	4.0
1	A	191	GLY	3.9
1	A	162	ALA	3.8
1	A	149	LYS	3.8
1	A	182	ALA	3.7
1	A	132	ALA	3.6
1	A	103	GLY	3.6
1	A	120	HIS	3.5
1	A	100	SER	3.3
1	A	117	PRO	3.2
1	A	50	ARG	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	160	GLY	3.0
1	A	163	PHE	2.9
1	A	119	GLY	2.9
1	A	93	CYS	2.8
1	A	91	GLY	2.8
1	A	157	ASP	2.8
1	B	158	GLU	2.8
1	A	152	LYS	2.7
1	B	176	PRO	2.7
1	B	87	ILE	2.6
1	B	66	GLU	2.5
1	A	189	LEU	2.5
1	A	147	LEU	2.5
1	A	176	PRO	2.5
1	A	159	ASP	2.4
1	A	148	LYS	2.4
1	A	165	ILE	2.4
1	B	85	PHE	2.3
1	A	58	ILE	2.2
1	B	144	ASP	2.2
1	A	57	PHE	2.2
1	B	40	ALA	2.2
1	A	98	PHE	2.1
1	A	155	ILE	2.1
1	A	107	PRO	2.1
1	A	94	VAL	2.1
1	B	163	PHE	2.0
1	A	68	LEU	2.0
1	A	177	SER	2.0
1	A	85	PHE	2.0
1	A	96	PRO	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 monosaccharides in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	ZN	B	505	1/1	0.95	0.15	26,26,26,26	1
2	CU	A	503	1/1	0.96	0.04	41,41,41,41	0
3	ZN	A	504	1/1	0.96	0.05	33,33,33,33	1
3	ZN	B	506	1/1	0.99	0.09	21,21,21,21	1
2	CU	B	500	1/1	0.99	0.04	23,23,23,23	0
3	ZN	B	502	1/1	1.00	0.04	21,21,21,21	0
3	ZN	B	501	1/1	1.00	0.05	20,20,20,20	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.