



# Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 07:55 pm BST

PDB ID : 3P2M  
Title : Crystal Structure of a Novel Esterase Rv0045c from Mycobacterium tuberculosis  
Authors : Zheng, X.D.; Guo, J.; Xu, L.; Li, H.; Zhang, D.; Zhang, K.; Sun, F.; Wen, T.; Liu, S.; Pang, H.  
Deposited on : 2010-10-03  
Resolution : 2.80 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 : 2.11  
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) : 2.11

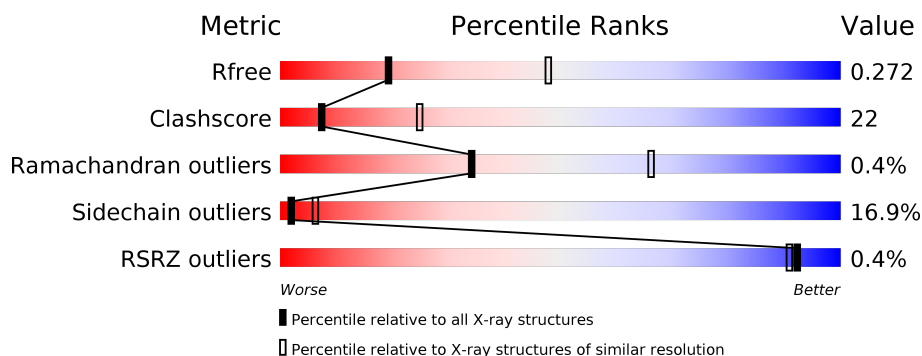
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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  $\geq 3$ , 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	330	<div> <div></div> <div>54%</div> <div>25%</div> <div>7%</div> <div>15%</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2172 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 POSSIBLE HYDROLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	281	Total	C	N	O	S	0	0	0
			2136	1340	391	402	3			

There are 33 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	EXPRESSION TAG	UNP P71702
A	2	GLY	-	EXPRESSION TAG	UNP P71702
A	3	SER	-	EXPRESSION TAG	UNP P71702
A	4	SER	-	EXPRESSION TAG	UNP P71702
A	5	HIS	-	EXPRESSION TAG	UNP P71702
A	6	HIS	-	EXPRESSION TAG	UNP P71702
A	7	HIS	-	EXPRESSION TAG	UNP P71702
A	8	HIS	-	EXPRESSION TAG	UNP P71702
A	9	HIS	-	EXPRESSION TAG	UNP P71702
A	10	HIS	-	EXPRESSION TAG	UNP P71702
A	11	SER	-	EXPRESSION TAG	UNP P71702
A	12	SER	-	EXPRESSION TAG	UNP P71702
A	13	GLY	-	EXPRESSION TAG	UNP P71702
A	14	LEU	-	EXPRESSION TAG	UNP P71702
A	15	VAL	-	EXPRESSION TAG	UNP P71702
A	16	PRO	-	EXPRESSION TAG	UNP P71702
A	17	ARG	-	EXPRESSION TAG	UNP P71702
A	18	GLY	-	EXPRESSION TAG	UNP P71702
A	19	SER	-	EXPRESSION TAG	UNP P71702
A	20	HIS	-	EXPRESSION TAG	UNP P71702
A	21	MET	-	EXPRESSION TAG	UNP P71702
A	22	ALA	-	EXPRESSION TAG	UNP P71702
A	23	SER	-	EXPRESSION TAG	UNP P71702
A	24	MET	-	EXPRESSION TAG	UNP P71702
A	25	THR	-	EXPRESSION TAG	UNP P71702
A	26	GLY	-	EXPRESSION TAG	UNP P71702
A	27	GLY	-	EXPRESSION TAG	UNP P71702

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Chain	Residue	Modelled	Actual	Comment	Reference
A	28	GLN	-	EXPRESSION TAG	UNP P71702
A	29	GLN	-	EXPRESSION TAG	UNP P71702
A	30	MET	-	EXPRESSION TAG	UNP P71702
A	31	GLY	-	EXPRESSION TAG	UNP P71702
A	32	ARG	-	EXPRESSION TAG	UNP P71702
A	33	VAL	-	EXPRESSION TAG	UNP P71702

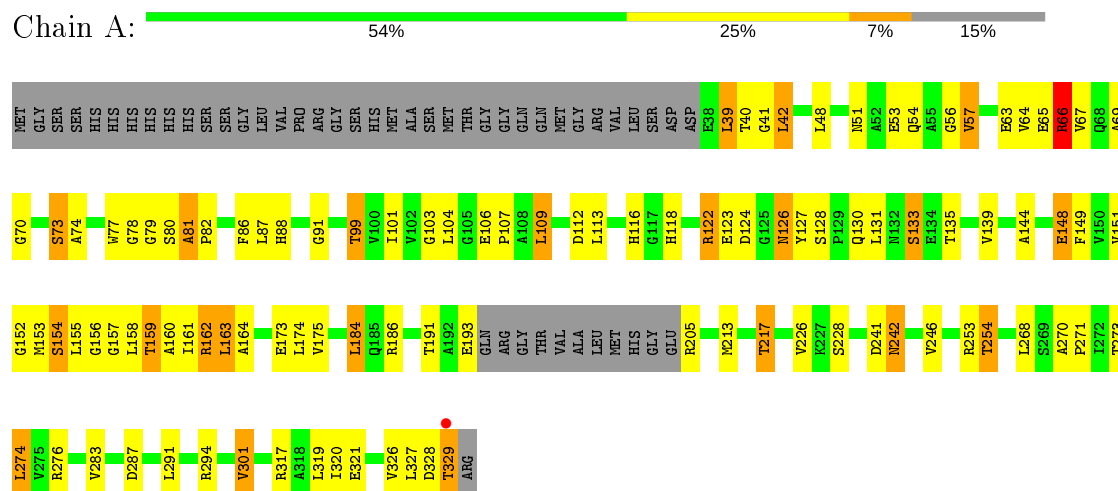
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	36	Total O 36 36	0	0

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

These plots are drawn for all protein, RNA and DNA 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: POSSIBLE HYDROLASE



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	73.47Å 73.47Å 48.06Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	38.35 – 2.80 38.35 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (38.35-2.80) 100.0 (38.35-2.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.46 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, $R_{free}$	0.217 , 0.286 0.211 , 0.272	Depositor DCC
$R_{free}$ test set	336 reflections (4.70%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	57.1	Xtriage
Anisotropy	0.261	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 36.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.028 for -h,-k,l 0.046 for h,-h-k,-l 0.029 for -k,-h,-l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	2172	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.00% 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	0.72	0/2183	0.84	2/2973 (0.1%)

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	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	41	GLY	N-CA-C	-5.53	99.28	113.10
1	A	66	ARG	NE-CZ-NH1	5.21	122.90	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	123	GLU	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	2136	0	2084	94	0
2	A	36	0	0	3	0
All	All	2172	0	2084	94	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 22.

All (94) 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:54:GLN:NE2	1:A:99:THR:HG21	1.39	1.35
1:A:40:THR:O	1:A:40:THR:HG22	1.53	1.06
1:A:54:GLN:HE22	1:A:99:THR:HG21	1.31	0.92
1:A:81:ALA:CB	1:A:107:PRO:HB2	2.01	0.90
1:A:164:ALA:CB	1:A:268:LEU:HD21	2.03	0.88
1:A:81:ALA:HB1	1:A:107:PRO:HB2	1.55	0.88
1:A:54:GLN:HE21	1:A:99:THR:HG21	1.12	0.88
1:A:81:ALA:HB1	1:A:82:PRO:HA	1.58	0.86
1:A:164:ALA:HB1	1:A:268:LEU:HD21	1.59	0.84
1:A:126:ASN:CG	1:A:131:LEU:HD12	1.98	0.83
1:A:126:ASN:C	1:A:126:ASN:HD22	1.84	0.81
1:A:40:THR:HG23	1:A:42:LEU:HB2	1.63	0.80
1:A:54:GLN:HE21	1:A:99:THR:CG2	1.93	0.76
1:A:66:ARG:NH2	1:A:118:HIS:HD2	1.85	0.74
1:A:51:ASN:OD1	1:A:99:THR:HG22	1.88	0.73
1:A:40:THR:O	1:A:40:THR:CG2	2.30	0.71
1:A:88:HIS:CE1	1:A:113:LEU:H	2.08	0.71
1:A:79:GLY:HA2	1:A:80:SER:C	2.12	0.70
1:A:66:ARG:NH2	1:A:118:HIS:CD2	2.60	0.69
1:A:155:LEU:O	1:A:159:THR:CG2	2.43	0.67
1:A:66:ARG:HD2	1:A:73:SER:OG	1.95	0.67
1:A:174:LEU:HD23	1:A:175:VAL:N	2.10	0.65
1:A:148:GLU:HB3	1:A:149:PHE:CD1	2.32	0.65
1:A:81:ALA:HB1	1:A:82:PRO:CA	2.26	0.65
1:A:173:GLU:HG3	1:A:174:LEU:N	2.13	0.64
1:A:54:GLN:NE2	1:A:99:THR:CG2	2.35	0.64
1:A:66:ARG:HH11	1:A:66:ARG:HG2	1.62	0.64
1:A:81:ALA:HB2	1:A:107:PRO:HB2	1.79	0.64
1:A:91:GLY:H	1:A:116:HIS:HE1	1.48	0.62
1:A:155:LEU:O	1:A:159:THR:HG23	2.00	0.61
1:A:126:ASN:ND2	1:A:128:SER:H	2.00	0.59
1:A:88:HIS:HE1	1:A:113:LEU:H	1.49	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:GLU:HB3	1:A:78:GLY:HA2	1.85	0.59
1:A:164:ALA:HB1	1:A:268:LEU:CD2	2.32	0.58
1:A:186:ARG:HH22	1:A:287:ASP:CG	2.06	0.58
1:A:158:LEU:HD12	1:A:184:LEU:HD12	1.86	0.58
1:A:213:MET:O	1:A:217:THR:HG23	2.04	0.57
1:A:66:ARG:HH11	1:A:66:ARG:CG	2.17	0.57
1:A:321:GLU:HG2	2:A:347:HOH:O	2.05	0.56
1:A:164:ALA:HB3	1:A:268:LEU:HD21	1.85	0.56
1:A:126:ASN:HA	1:A:254:THR:HG22	1.87	0.55
1:A:122:ARG:HG3	1:A:127:TYR:CE2	2.42	0.54
1:A:39:LEU:HD23	1:A:39:LEU:N	2.23	0.54
1:A:69:ALA:O	1:A:70:GLY:C	2.45	0.52
1:A:82:PRO:HG2	1:A:144:ALA:CB	2.40	0.52
1:A:122:ARG:NH2	1:A:124:ASP:OD2	2.42	0.51
1:A:133:SER:CB	1:A:163:LEU:HD22	2.40	0.51
1:A:82:PRO:HG2	1:A:144:ALA:HB2	1.92	0.51
1:A:57:VAL:HG22	2:A:355:HOH:O	2.10	0.51
1:A:130:GLN:HA	1:A:162:ARG:HG3	1.93	0.51
1:A:66:ARG:HH22	1:A:118:HIS:HD2	1.57	0.51
1:A:173:GLU:HG3	1:A:174:LEU:H	1.74	0.51
1:A:155:LEU:O	1:A:159:THR:HG22	2.10	0.51
1:A:133:SER:HB3	1:A:163:LEU:HD22	1.93	0.50
1:A:276:ARG:NH1	1:A:283:VAL:O	2.44	0.50
1:A:126:ASN:ND2	1:A:126:ASN:C	2.58	0.50
1:A:133:SER:HA	1:A:163:LEU:HD22	1.94	0.49
1:A:88:HIS:HE1	1:A:112:ASP:HA	1.76	0.49
1:A:154:SER:HB3	1:A:155:LEU:H	1.37	0.49
1:A:65:GLU:HG3	1:A:67:VAL:HG23	1.95	0.48
1:A:56:GLY:CA	1:A:317:ARG:HH21	2.26	0.48
1:A:39:LEU:N	1:A:39:LEU:CD2	2.76	0.48
1:A:69:ALA:HB3	1:A:139:VAL:HG13	1.95	0.48
1:A:77:TRP:CD2	1:A:101:ILE:HG21	2.49	0.47
1:A:48:LEU:HA	1:A:48:LEU:HD23	1.48	0.47
1:A:148:GLU:HB3	1:A:149:PHE:CE1	2.49	0.47
1:A:103:GLY:HA3	1:A:320:ILE:HD13	1.97	0.46
1:A:191:THR:C	1:A:193:GLU:H	2.19	0.46
1:A:74:ALA:HB1	1:A:109:LEU:HD22	1.98	0.46
1:A:326:VAL:C	1:A:328:ASP:H	2.20	0.45
1:A:126:ASN:OD1	1:A:131:LEU:HD12	2.16	0.45
1:A:186:ARG:NH2	1:A:287:ASP:OD1	2.49	0.45
1:A:152:GLY:HA3	1:A:157:GLY:HA2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:ALA:CB	1:A:82:PRO:CA	2.94	0.44
1:A:66:ARG:CZ	1:A:118:HIS:CD2	3.00	0.44
1:A:242:ASN:N	1:A:242:ASN:OD1	2.51	0.44
1:A:149:PHE:CD2	1:A:173:GLU:HB3	2.53	0.44
1:A:113:LEU:O	1:A:116:HIS:HB2	2.16	0.43
1:A:131:LEU:HD23	1:A:131:LEU:HA	1.78	0.43
1:A:79:GLY:CA	1:A:80:SER:C	2.85	0.43
1:A:270:ALA:HA	1:A:271:PRO:HD3	1.92	0.43
1:A:79:GLY:HA3	1:A:81:ALA:CB	2.49	0.43
1:A:327:LEU:C	1:A:329:THR:H	2.22	0.42
1:A:63:GLU:CB	1:A:78:GLY:HA2	2.48	0.42
1:A:133:SER:HA	1:A:163:LEU:CD2	2.50	0.42
1:A:205:ARG:HB3	2:A:349:HOH:O	2.19	0.41
1:A:158:LEU:O	1:A:161:ILE:HB	2.20	0.41
1:A:174:LEU:HD23	1:A:175:VAL:H	1.84	0.41
1:A:274:LEU:HD13	1:A:301:VAL:CG1	2.51	0.41
1:A:66:ARG:NH1	1:A:66:ARG:HG2	2.33	0.41
1:A:156:GLY:HA2	1:A:159:THR:HG23	2.02	0.41
1:A:82:PRO:HD2	1:A:144:ALA:HA	2.01	0.41
1:A:87:LEU:HD12	1:A:160:ALA:HB2	2.03	0.41
1:A:86:PHE:CD1	1:A:151:VAL:HB	2.56	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	277/330 (84%)	262 (95%)	14 (5%)	1 (0%)	34 66

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	81	ALA

### 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	219/258 (85%)	182 (83%)	37 (17%)	<b>2</b> <b>6</b>

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	LEU
1	A	42	LEU
1	A	53	GLU
1	A	57	VAL
1	A	64	VAL
1	A	66	ARG
1	A	73	SER
1	A	99	THR
1	A	104	LEU
1	A	106	GLU
1	A	109	LEU
1	A	122	ARG
1	A	126	ASN
1	A	133	SER
1	A	135	THR
1	A	148	GLU
1	A	153	MET
1	A	154	SER
1	A	159	THR
1	A	162	ARG
1	A	163	LEU
1	A	184	LEU
1	A	217	THR
1	A	226	VAL
1	A	228	SER
1	A	241	ASP

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Mol	Chain	Res	Type
1	A	242	ASN
1	A	246	VAL
1	A	253	ARG
1	A	254	THR
1	A	273	THR
1	A	274	LEU
1	A	291	LEU
1	A	294	ARG
1	A	301	VAL
1	A	319	LEU
1	A	329	THR

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	54	GLN
1	A	88	HIS
1	A	116	HIS
1	A	118	HIS
1	A	126	ASN
1	A	223	HIS
1	A	235	HIS

### 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 carbohydrates in this entry.

### 5.6 Ligand geometry ⓘ

There are no ligands in this entry.

## 5.7 Other polymers

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	281/330 (85%)	-0.23	1 (0%) 92 91	32, 52, 71, 82	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	329	THR	2.3

### 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.