



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2018 – 08:26 PM EST

PDB ID : 5X87  
Title : Crystal structure of a bacterial Bestrophin homolog from *Klebsiella pneumoniae* with a mutation L177T  
Authors : Zhang, Y.; Chen, S.; Yang, T.  
Deposited on : 2017-03-01  
Resolution : 3.14 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

<http://wwpdb.org/validation/2016/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)	:	1.9-1692
EDS	:	rb-20030736
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	rb-20030736

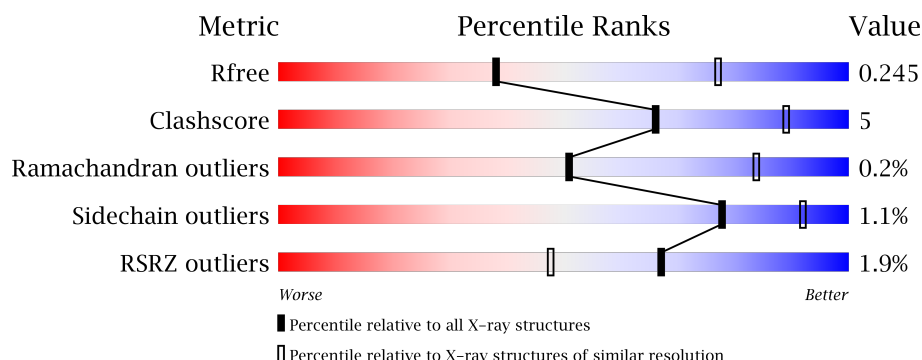
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.14 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))
$R_{free}$	100719	1234 (3.18-3.10)
Clashscore	112137	1345 (3.18-3.10)
Ramachandran outliers	110173	1301 (3.18-3.10)
Sidechain outliers	110143	1301 (3.18-3.10)
RSRZ outliers	101464	1240 (3.18-3.10)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	297	
1	B	297	
1	C	297	
1	D	297	
1	E	297	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ZN	A	304	-	-	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10678 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 Bestrophin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	268	Total	C	N	O	S	0	0	0
			2119	1371	359	380	9			
1	B	268	Total	C	N	O	S	0	0	0
			2130	1378	360	383	9			
1	C	269	Total	C	N	O	S	0	0	0
			2109	1360	362	378	9			
1	D	268	Total	C	N	O	S	0	0	0
			2113	1365	359	380	9			
1	E	270	Total	C	N	O	S	0	0	0
			2151	1394	364	384	9			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP W1ELP7
A	-1	ASN	-	expression tag	UNP W1ELP7
A	0	ALA	-	expression tag	UNP W1ELP7
A	177	THR	LEU	engineered mutation	UNP W1ELP7
B	-2	SER	-	expression tag	UNP W1ELP7
B	-1	ASN	-	expression tag	UNP W1ELP7
B	0	ALA	-	expression tag	UNP W1ELP7
B	177	THR	LEU	engineered mutation	UNP W1ELP7
C	-2	SER	-	expression tag	UNP W1ELP7
C	-1	ASN	-	expression tag	UNP W1ELP7
C	0	ALA	-	expression tag	UNP W1ELP7
C	177	THR	LEU	engineered mutation	UNP W1ELP7
D	-2	SER	-	expression tag	UNP W1ELP7
D	-1	ASN	-	expression tag	UNP W1ELP7
D	0	ALA	-	expression tag	UNP W1ELP7
D	177	THR	LEU	engineered mutation	UNP W1ELP7
E	-2	SER	-	expression tag	UNP W1ELP7
E	-1	ASN	-	expression tag	UNP W1ELP7
E	0	ALA	-	expression tag	UNP W1ELP7

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Chain	Residue	Modelled	Actual	Comment	Reference
E	177	THR	LEU	engineered mutation	UNP W1ELP7

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	3	Total Zn 3 3	0	0
2	A	4	Total Zn 4 4	0	0
2	D	3	Total Zn 3 3	0	0
2	C	3	Total Zn 3 3	0	0
2	E	2	Total Zn 2 2	0	0


- Molecule 3 is water.

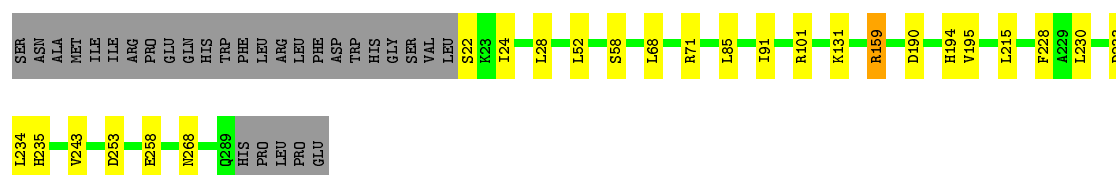
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	6	Total O 6 6	0	0
3	B	5	Total O 5 5	0	0
3	C	10	Total O 10 10	0	0
3	D	10	Total O 10 10	0	0
3	E	10	Total O 10 10	0	0

### 3 Residue-property plots


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.

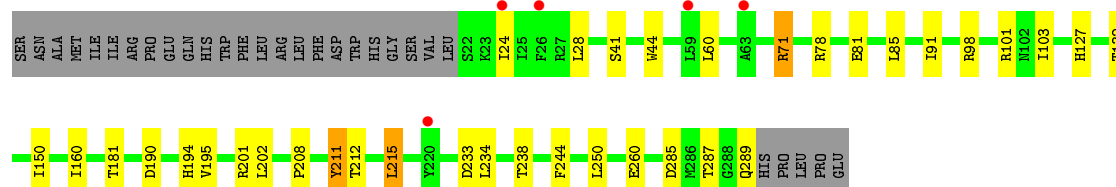
#### • Molecule 1: Bestrophin

Chain A: 



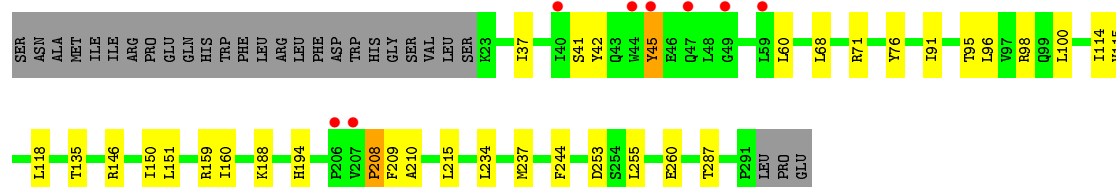
#### • Molecule 1: Bestrophin

Chain B: 




#### • Molecule 1: Bestrophin

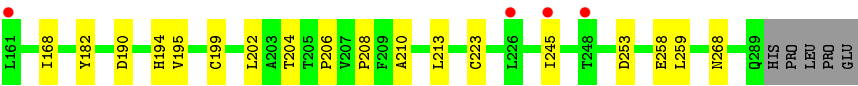
Chain C: 



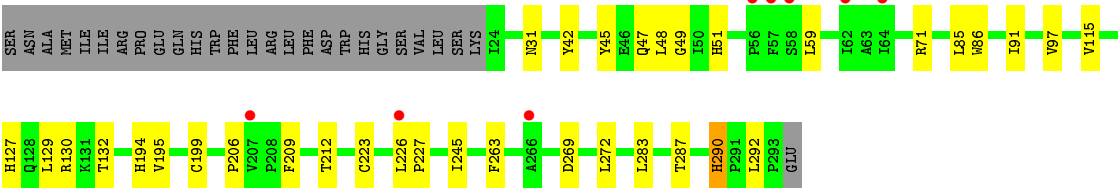
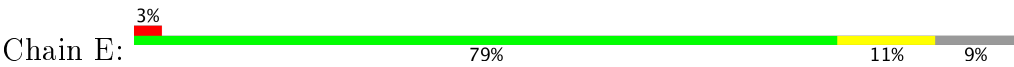
#### • Molecule 1: Bestrophin

Chain D: 





● Molecule 1: Bestrophin



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	114.48Å 160.57Å 162.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	65.73 – 3.14 114.27 – 3.14	Depositor EDS
% Data completeness (in resolution range)	99.9 (65.73-3.14) 89.6 (114.27-3.14)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.51 (at 3.13Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, $R_{free}$	0.199 , 0.237 0.201 , 0.245	Depositor DCC
$R_{free}$ test set	1843 reflections (3.75%)	DCC
Wilson B-factor (Å <sup>2</sup> )	73.9	Xtriage
Anisotropy	0.023	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.27 , 52.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.025 for -h,l,k	Xtriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	10678	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	92.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.29% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.23	0/2165	0.37	0/2949
1	B	0.26	0/2176	0.42	0/2963
1	C	0.23	0/2154	0.37	0/2936
1	D	0.23	0/2158	0.37	0/2940
1	E	0.28	0/2200	0.43	0/2998
All	All	0.25	0/10853	0.39	0/14786

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 [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	2119	0	2147	22	1
1	B	2130	0	2166	25	0
1	C	2109	0	2130	24	0
1	D	2113	0	2140	24	0
1	E	2151	0	2191	26	1
2	A	4	0	0	0	0
2	B	3	0	0	0	0
2	C	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	3	0	0	0	0
2	E	2	0	0	0	0
3	A	6	0	0	0	0
3	B	5	0	0	0	0
3	C	10	0	0	0	0
3	D	10	0	0	1	0
3	E	10	0	0	0	0
All	All	10678	0	10774	98	1

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 5.

All (98) 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:235:HIS:ND1	1:E:51:HIS:HD2	1.03	1.49
1:A:235:HIS:ND1	1:E:51:HIS:CD2	1.90	1.37
1:A:235:HIS:CE1	1:E:51:HIS:CD2	2.36	1.12
1:A:235:HIS:CE1	1:E:51:HIS:HD2	1.72	1.06
1:B:127:HIS:HD1	1:B:132:THR:HG1	1.13	0.92
1:B:71:ARG:NH1	1:B:260:GLU:OE2	2.12	0.81
1:A:258:GLU:OE1	1:A:268:ASN:ND2	2.21	0.72
1:A:235:HIS:CG	1:E:51:HIS:CD2	2.77	0.71
1:E:47:GLN:HG2	1:E:47:GLN:O	1.90	0.71
1:D:78:ARG:NH2	1:D:202:LEU:O	2.26	0.68
1:B:41:SER:HA	1:B:44:TRP:HD1	1.59	0.66
1:A:230:LEU:HB3	1:A:234:LEU:HD12	1.78	0.65
1:B:71:ARG:HH22	1:B:212:THR:HG22	1.62	0.65
1:D:258:GLU:OE1	1:D:268:ASN:ND2	2.30	0.64
1:E:48:LEU:N	1:E:48:LEU:HD23	2.13	0.63
1:A:68:LEU:HD23	1:A:215:LEU:HD13	1.81	0.61
1:A:91:ILE:HD13	1:E:194:HIS:HA	1.82	0.61
1:C:91:ILE:HD13	1:D:194:HIS:HA	1.82	0.61
1:C:194:HIS:HA	1:E:91:ILE:HD13	1.82	0.61
1:B:103:ILE:HD13	1:B:181:THR:HB	1.83	0.59
1:B:24:ILE:HG22	1:B:28:LEU:HG	1.84	0.58
1:B:78:ARG:NH2	1:B:202:LEU:O	2.36	0.58
1:B:190:ASP:OD1	1:D:98:ARG:NH2	2.38	0.57
1:C:208:PRO:O	1:C:210:ALA:N	2.38	0.57
1:B:194:HIS:HA	1:D:91:ILE:HD13	1.85	0.56
1:A:194:HIS:HA	1:B:91:ILE:HD13	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:71:ARG:NH2	1:D:253:ASP:OD1	2.32	0.55
1:B:287:THR:OG1	1:B:287:THR:O	2.22	0.54
1:C:98:ARG:NH2	1:D:190:ASP:OD1	2.41	0.54
1:E:31:ASN:ND2	1:E:223:CYS:O	2.33	0.54
1:C:115:VAL:HG21	1:C:287:THR:HG21	1.91	0.52
1:C:68:LEU:HD23	1:C:215:LEU:HD13	1.91	0.52
1:B:211:TYR:O	1:B:215:LEU:HB2	2.10	0.52
1:B:24:ILE:HD12	1:B:250:LEU:HB3	1.90	0.52
1:C:42:TYR:HA	1:C:45:TYR:CD1	2.45	0.52
1:C:71:ARG:NH1	1:C:260:GLU:OE2	2.42	0.52
1:D:31:ASN:ND2	1:D:223:CYS:O	2.43	0.51
1:B:208:PRO:HG2	1:D:259:LEU:HD21	1.92	0.51
1:E:42:TYR:HA	1:E:45:TYR:CD1	2.46	0.50
1:A:28:LEU:HD22	1:A:243:VAL:HG13	1.94	0.50
1:C:71:ARG:NH2	1:C:253:ASP:OD1	2.43	0.50
1:B:60:LEU:HD13	1:D:245:ILE:HG12	1.94	0.49
1:A:22:SER:HB2	1:A:24:ILE:HG22	1.94	0.49
1:D:146:ARG:NH1	1:D:149:GLU:OE1	2.42	0.49
1:C:96:LEU:HD23	1:C:118:LEU:HD11	1.94	0.48
1:D:108:HIS:ND1	3:D:402:HOH:O	2.35	0.48
1:C:150:ILE:HA	1:C:159:ARG:HG2	1.95	0.48
1:D:139:ARG:HG3	1:D:147:VAL:HG11	1.96	0.48
1:A:159:ARG:NH2	1:B:289:GLN:O	2.46	0.48
1:A:52:LEU:HD12	1:A:228:PHE:HB3	1.96	0.48
1:E:115:VAL:HG21	1:E:287:THR:HG21	1.97	0.47
1:B:150:ILE:HD13	1:B:160:ILE:HG12	1.96	0.47
1:C:146:ARG:HH11	1:C:146:ARG:HA	1.80	0.47
1:D:150:ILE:HD13	1:D:160:ILE:HG12	1.97	0.47
1:E:130:ARG:NH1	1:E:269:ASP:OD2	2.47	0.47
1:E:42:TYR:HA	1:E:45:TYR:CE1	2.50	0.46
1:E:97:VAL:HG21	1:E:283:LEU:HD22	1.97	0.46
1:A:85:LEU:HD22	1:A:195:VAL:HA	1.97	0.46
1:C:188:LYS:HD3	1:C:188:LYS:HA	1.78	0.46
1:A:190:ASP:OD1	1:B:98:ARG:NH2	2.49	0.46
1:B:208:PRO:O	1:B:211:TYR:HB2	2.15	0.46
1:C:100:LEU:HD11	1:C:114:ILE:HD12	1.97	0.46
1:E:85:LEU:HD22	1:E:195:VAL:HA	1.98	0.46
1:C:208:PRO:C	1:C:210:ALA:H	2.20	0.45
1:E:129:LEU:HD21	1:E:199:CYS:HB3	1.98	0.45
1:D:204:THR:O	1:D:206:PRO:HD3	2.16	0.45
1:E:71:ARG:NH1	1:E:212:THR:OG1	2.39	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:71:ARG:NH2	1:A:253:ASP:OD1	2.42	0.45
1:D:129:LEU:HD21	1:D:199:CYS:HB3	1.98	0.45
1:B:85:LEU:HD22	1:B:195:VAL:HA	1.98	0.45
1:E:127:HIS:ND1	1:E:132:THR:OG1	2.44	0.45
1:B:127:HIS:ND1	1:B:132:THR:OG1	2.24	0.44
1:A:58:SER:HA	1:E:59:LEU:HD13	1.99	0.44
1:C:135:THR:HG23	1:C:151:LEU:HD11	2.00	0.44
1:C:60:LEU:HD13	1:E:245:ILE:HG12	2.00	0.43
1:E:86:TRP:CE2	1:E:272:LEU:HD11	2.54	0.43
1:C:150:ILE:HD13	1:C:160:ILE:HG12	2.00	0.43
1:A:52:LEU:HA	1:B:234:LEU:HD22	2.00	0.43
1:D:127:HIS:ND1	1:D:132:THR:OG1	2.43	0.43
1:D:168:ILE:HG22	1:D:182:TYR:HD1	1.84	0.43
1:D:210:ALA:HA	1:D:213:LEU:HB3	2.00	0.43
1:C:255:LEU:HD13	1:D:210:ALA:HB3	2.01	0.43
1:B:234:LEU:O	1:B:238:THR:OG1	2.37	0.43
1:B:285:ASP:HB3	1:B:289:GLN:O	2.19	0.43
1:C:95:THR:OG1	1:C:188:LYS:HD2	2.20	0.42
1:D:89:VAL:HG23	1:D:195:VAL:HG11	2.01	0.42
1:A:235:HIS:HB2	1:E:49:GLY:O	2.20	0.42
1:C:76:TYR:HE1	1:D:208:PRO:HD2	1.85	0.42
1:C:37:ILE:O	1:C:41:SER:OG	2.24	0.42
1:B:81:GLU:OE2	1:B:201:ARG:NH2	2.49	0.41
1:E:129:LEU:HB3	1:E:263:PHE:HB3	2.02	0.41
1:C:234:LEU:HB3	1:C:237:MET:HB2	2.02	0.41
1:A:101:ARG:HB3	1:A:101:ARG:HE	1.60	0.41
1:A:234:LEU:O	1:A:235:HIS:HB2	2.21	0.41
1:D:168:ILE:HG22	1:D:182:TYR:CD1	2.56	0.41
1:E:226:LEU:N	1:E:227:PRO:HD2	2.36	0.41
1:E:292:LEU:HD23	1:E:292:LEU:HA	1.92	0.41
1:C:234:LEU:HA	1:D:51:HIS:HB2	2.03	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:LYS:CB	1:E:290:HIS:O[4_545]	2.17	0.03

## 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	266/297 (90%)	264 (99%)	2 (1%)	0	100	100
1	B	266/297 (90%)	258 (97%)	8 (3%)	0	100	100
1	C	267/297 (90%)	259 (97%)	6 (2%)	2 (1%)	25	64
1	D	266/297 (90%)	263 (99%)	3 (1%)	0	100	100
1	E	268/297 (90%)	262 (98%)	5 (2%)	1 (0%)	38	75
All	All	1333/1485 (90%)	1306 (98%)	24 (2%)	3 (0%)	51	84

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	209	PHE
1	C	208	PRO
1	E	206	PRO

### 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	229/260 (88%)	227 (99%)	2 (1%)	82	93
1	B	232/260 (89%)	226 (97%)	6 (3%)	51	81
1	C	226/260 (87%)	224 (99%)	2 (1%)	82	93
1	D	228/260 (88%)	227 (100%)	1 (0%)	93	97
1	E	235/260 (90%)	233 (99%)	2 (1%)	82	93

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1150/1300 (88%)	1137 (99%)	13 (1%)	78 92

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	159	ARG
1	A	233	ASP
1	B	71	ARG
1	B	101	ARG
1	B	211	TYR
1	B	215	LEU
1	B	233	ASP
1	B	244	PHE
1	C	45	TYR
1	C	244	PHE
1	D	126	LYS
1	E	209	PHE
1	E	290	HIS

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

Mol	Chain	Res	Type
1	B	102	ASN
1	D	268	ASN
1	E	51	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

Of 15 ligands modelled in this entry, 15 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

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 ⓘ

### 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	268/297 (90%)	-0.05	0 100 100	47, 82, 147, 188	0
1	B	268/297 (90%)	0.09	5 (1%) 67 48	50, 90, 140, 157	0
1	C	269/297 (90%)	0.01	8 (2%) 51 28	50, 83, 158, 189	0
1	D	268/297 (90%)	0.04	5 (1%) 67 48	55, 94, 141, 160	0
1	E	270/297 (90%)	0.11	8 (2%) 51 28	47, 82, 140, 186	0
All	All	1343/1485 (90%)	0.04	26 (1%) 67 48	47, 85, 145, 189	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	44	TRP	5.6
1	B	220	TYR	4.0
1	C	40	ILE	3.3
1	E	56	PRO	3.0
1	E	57	PHE	2.9
1	E	58	SER	2.9
1	B	63	ALA	2.9
1	E	207	VAL	2.7
1	B	24	ILE	2.6
1	C	47	GLN	2.6
1	B	59	LEU	2.5
1	E	226	LEU	2.5
1	C	59	LEU	2.5
1	C	206	PRO	2.4
1	C	207	VAL	2.4
1	E	266	ALA	2.4
1	B	26	PHE	2.4
1	D	58	SER	2.3
1	E	62	ILE	2.3
1	E	64	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	226	LEU	2.2
1	D	248	THR	2.2
1	C	49	GLY	2.1
1	D	161	LEU	2.1
1	D	245	ILE	2.1
1	C	45	TYR	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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	ZN	A	304	1/1	0.99	0.26	9.07	91,91,91,91	0
2	ZN	C	302	1/1	0.97	0.20	0.10	79,79,79,79	0
2	ZN	A	301	1/1	0.99	0.20	-0.92	74,74,74,74	0
2	ZN	C	301	1/1	0.98	0.28	-	61,61,61,61	0
2	ZN	D	303	1/1	0.91	0.21	-	156,156,156,156	0
2	ZN	B	301	1/1	0.96	0.21	-	137,137,137,137	0
2	ZN	D	302	1/1	0.65	0.18	-	113,113,113,113	0
2	ZN	E	301	1/1	0.97	0.25	-	63,63,63,63	0
2	ZN	E	302	1/1	0.85	0.22	-	74,74,74,74	0
2	ZN	B	302	1/1	0.99	0.19	-	84,84,84,84	0
2	ZN	D	301	1/1	0.97	0.25	-	87,87,87,87	0
2	ZN	A	302	1/1	0.98	0.19	-	77,77,77,77	0
2	ZN	A	303	1/1	0.98	0.27	-	83,83,83,83	0
2	ZN	C	303	1/1	0.99	0.23	-	71,71,71,71	0
2	ZN	B	303	1/1	0.94	0.27	-	74,74,74,74	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.