



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:28 PM GMT

PDB ID : 1X8C  
Title : Crystal structure of the SeMet-derivative copper homeostasis protein (cutCm) with calcium binding from *Shigella flexneri* 2a str. 301  
Authors : Wang, D.C.; Zhu, D.Y.  
Deposited on : 2004-08-18  
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  
<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  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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) : trunk26865

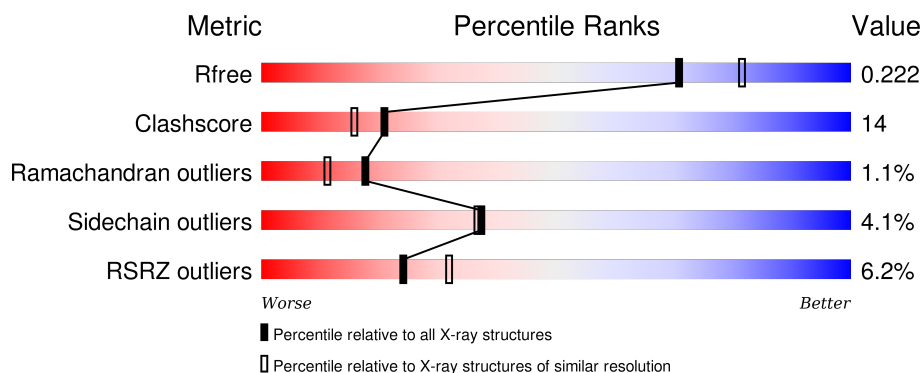
# 1 Overall quality at a glance

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))
$R_{free}$	91344	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (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 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	256	
1	B	256	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3947 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 Copper homeostasis protein cutC.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	234	Total	C	N	O	S	Se	0	0	0
			1765	1108	317	324	6	10			
1	B	238	Total	C	N	O	S	Se	0	0	0
			1799	1128	322	333	6	10			

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	MODIFIED RESIDUE	UNP P67825
A	11	MSE	MET	MODIFIED RESIDUE	UNP P67825
A	103	MSE	MET	MODIFIED RESIDUE	UNP P67825
A	106	MSE	MET	MODIFIED RESIDUE	UNP P67825
A	110	MSE	MET	MODIFIED RESIDUE	UNP P67825
A	126	MSE	MET	MODIFIED RESIDUE	UNP P67825
A	162	MSE	MET	MODIFIED RESIDUE	UNP P67825
A	174	MSE	MET	MODIFIED RESIDUE	UNP P67825
A	207	MSE	MET	MODIFIED RESIDUE	UNP P67825
A	216	MSE	MET	MODIFIED RESIDUE	UNP P67825
A	238	MSE	MET	MODIFIED RESIDUE	UNP P67825
A	249	LEU	-	EXPRESSION TAG	UNP P67825
A	250	GLU	-	EXPRESSION TAG	UNP P67825
A	251	HIS	-	EXPRESSION TAG	UNP P67825
A	252	HIS	-	EXPRESSION TAG	UNP P67825
A	253	HIS	-	EXPRESSION TAG	UNP P67825
A	254	HIS	-	EXPRESSION TAG	UNP P67825
A	255	HIS	-	EXPRESSION TAG	UNP P67825
A	256	HIS	-	EXPRESSION TAG	UNP P67825
B	1	MSE	MET	MODIFIED RESIDUE	UNP P67825
B	11	MSE	MET	MODIFIED RESIDUE	UNP P67825
B	103	MSE	MET	MODIFIED RESIDUE	UNP P67825
B	106	MSE	MET	MODIFIED RESIDUE	UNP P67825
B	110	MSE	MET	MODIFIED RESIDUE	UNP P67825
B	126	MSE	MET	MODIFIED RESIDUE	UNP P67825

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Chain	Residue	Modelled	Actual	Comment	Reference
B	162	MSE	MET	MODIFIED RESIDUE	UNP P67825
B	174	MSE	MET	MODIFIED RESIDUE	UNP P67825
B	207	MSE	MET	MODIFIED RESIDUE	UNP P67825
B	216	MSE	MET	MODIFIED RESIDUE	UNP P67825
B	238	MSE	MET	MODIFIED RESIDUE	UNP P67825
B	249	LEU	-	EXPRESSION TAG	UNP P67825
B	250	GLU	-	EXPRESSION TAG	UNP P67825
B	251	HIS	-	EXPRESSION TAG	UNP P67825
B	252	HIS	-	EXPRESSION TAG	UNP P67825
B	253	HIS	-	EXPRESSION TAG	UNP P67825
B	254	HIS	-	EXPRESSION TAG	UNP P67825
B	255	HIS	-	EXPRESSION TAG	UNP P67825
B	256	HIS	-	EXPRESSION TAG	UNP P67825

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

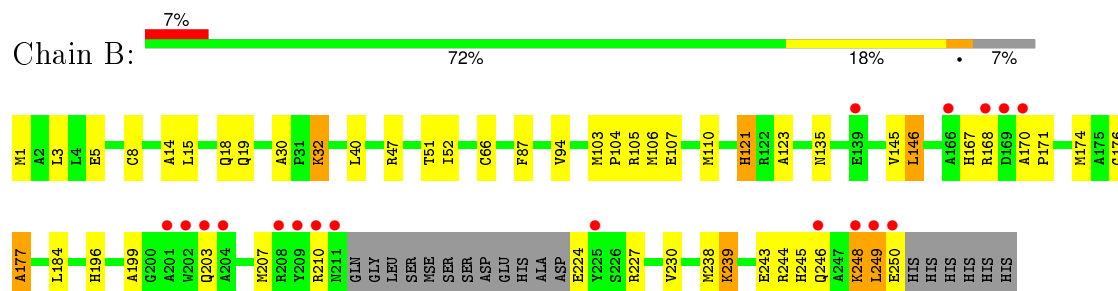
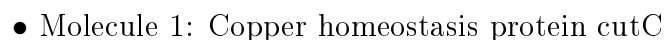
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Ca 1 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	191	Total O 191 191	0	0
3	B	191	Total O 191 191	0	0



- Molecule 1: Copper homeostasis protein cutC



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	74.58 Å 97.58 Å 131.85 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	59.26 – 2.10 59.25 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.4 (59.26-2.10) 99.6 (59.25-2.10)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	8.73 (at 2.10 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.196 , 0.222 0.193 , 0.222	Depositor DCC
$R_{free}$ test set	1377 reflections (4.86%)	DCC
Wilson B-factor (Å <sup>2</sup> )	14.9	Xtriage
Anisotropy	0.489	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 53.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 28343 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3947	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.80% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: CA

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	0.30	0/1786	0.60	1/2400 (0.0%)
1	B	0.30	0/1820	0.59	0/2446
All	All	0.30	0/3606	0.59	1/4846 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	170	ALA	C-N-CD	5.06	139.03	128.40

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	1765	0	1781	59	0
1	B	1799	0	1810	47	0
2	B	1	0	0	0	0
3	A	191	0	0	3	0
3	B	191	0	0	1	1
All	All	3947	0	3591	103	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 14.

All (103) 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:18:GLN:HE22	1:B:51:THR:H	1.07	0.99
1:B:203:GLN:NE2	1:B:227:ARG:HH11	1.63	0.97
1:A:18:GLN:HE22	1:A:51:THR:H	1.18	0.89
1:A:168:ARG:NE	1:A:170:ALA:HB2	1.87	0.89
1:A:135:ASN:ND2	1:A:168:ARG:HH21	1.73	0.85
1:A:1:MSE:H3	1:A:1:MSE:HE3	1.43	0.83
1:A:179:VAL:HG11	1:A:195:VAL:CG2	2.10	0.82
1:B:168:ARG:HD3	1:B:170:ALA:O	1.82	0.80
1:A:199:ALA:HB2	1:A:238:MSE:HE1	1.66	0.78
1:B:207:MSE:HE3	1:B:210:ARG:HG2	1.67	0.77
1:A:122:ARG:HD2	1:A:125:ASP:OD2	1.86	0.75
1:A:170:ALA:HB1	1:A:171:PRO:CD	2.19	0.72
1:B:18:GLN:NE2	1:B:51:THR:H	1.86	0.71
1:A:168:ARG:HE	1:A:170:ALA:HB2	1.54	0.69
1:B:203:GLN:HE22	1:B:227:ARG:HH11	1.39	0.69
1:A:179:VAL:HG11	1:A:195:VAL:HG21	1.73	0.69
1:A:170:ALA:HB1	1:A:171:PRO:HD2	1.74	0.68
1:A:168:ARG:CD	1:A:170:ALA:HB2	2.26	0.66
1:A:170:ALA:CB	1:A:171:PRO:HD2	2.26	0.65
1:A:131:LEU:HD22	1:A:167:HIS:CD2	2.31	0.65
1:A:205:SER:HB3	1:A:225:TYR:HB3	1.79	0.64
1:A:138:ALA:HB2	1:A:170:ALA:HB1	1.79	0.64
1:B:146:LEU:HG	1:B:174:MSE:HE3	1.79	0.64
1:A:125:ASP:O	1:A:126:MSE:HE2	1.99	0.63
1:B:203:GLN:HE21	1:B:227:ARG:HH11	1.45	0.63
1:B:203:GLN:NE2	1:B:227:ARG:NH1	2.43	0.62
1:B:18:GLN:HE22	1:B:51:THR:N	1.89	0.62
1:A:170:ALA:CB	1:A:171:PRO:CD	2.78	0.61
1:B:248:LYS:HE3	1:B:248:LYS:N	2.17	0.59
1:A:18:GLN:HE21	1:A:52:ILE:H	1.51	0.59
1:B:203:GLN:HE22	1:B:227:ARG:NH1	2.01	0.58
1:A:169:ASP:O	1:A:170:ALA:O	2.20	0.58
1:B:135:ASN:OD1	1:B:167:HIS:HE1	1.86	0.57
1:A:176:GLY:O	1:A:177:ALA:HB3	2.06	0.56
1:B:203:GLN:HE21	1:B:227:ARG:HD3	1.70	0.55
1:A:168:ARG:HG3	1:A:170:ALA:H	1.71	0.54
1:A:1:MSE:N	1:A:1:MSE:HE3	2.18	0.54
1:B:32:LYS:HZ2	1:B:32:LYS:HB3	1.73	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:5:GLU:OE1	1:A:196:HIS:HD2	1.90	0.53
1:B:5:GLU:OE1	1:B:196:HIS:HD2	1.90	0.53
1:B:203:GLN:HG3	1:B:227:ARG:HB2	1.89	0.53
1:A:135:ASN:HD21	1:A:167:HIS:HE1	1.56	0.53
1:B:248:LYS:C	1:B:250:GLU:H	2.11	0.53
1:A:15:LEU:O	1:A:19:GLN:HG3	2.08	0.53
1:B:15:LEU:O	1:B:19:GLN:HG3	2.09	0.52
1:A:36:LEU:HD12	1:B:30:ALA:HB2	1.91	0.52
1:B:18:GLN:HG3	1:B:52:ILE:HG23	1.91	0.52
1:A:229:ILE:C	1:A:229:ILE:HD12	2.31	0.52
1:A:135:ASN:HD22	1:A:168:ARG:HH21	1.53	0.52
1:B:176:GLY:O	1:B:177:ALA:HB3	2.10	0.52
1:A:135:ASN:HD21	1:A:167:HIS:CE1	2.28	0.51
1:A:207:MSE:SE	1:B:66:CYS:SG	3.19	0.51
1:B:18:GLN:HE21	1:B:52:ILE:H	1.58	0.51
1:A:18:GLN:NE2	1:A:51:THR:H	1.98	0.50
1:B:32:LYS:NZ	1:B:32:LYS:HB3	2.25	0.50
1:A:24:ARG:HD2	1:A:55:HIS:CE1	2.47	0.50
1:A:135:ASN:ND2	1:A:168:ARG:NH2	2.52	0.50
1:A:159:SER:HA	1:A:162:MSE:HE3	1.93	0.50
1:A:40:LEU:HD11	1:A:44:LYS:HE3	1.93	0.49
1:B:106:MSE:O	1:B:110:MSE:HG2	2.13	0.49
1:B:1:MSE:CA	1:B:246:GLN:HG3	2.42	0.49
1:A:44:LYS:O	1:A:48:GLN:HG2	2.13	0.48
1:A:18:GLN:NE2	1:A:52:ILE:H	2.10	0.48
1:A:121:HIS:HD2	1:A:123:ALA:H	1.61	0.48
1:A:176:GLY:O	1:A:177:ALA:CB	2.62	0.48
1:A:40:LEU:HG	1:B:40:LEU:HD22	1.96	0.48
1:A:138:ALA:HB2	1:A:170:ALA:CB	2.44	0.48
1:B:121:HIS:HB2	3:B:538:HOH:O	2.14	0.47
1:B:176:GLY:O	1:B:177:ALA:CB	2.61	0.47
1:A:210:ARG:HG2	1:A:225:TYR:CE2	2.49	0.47
1:A:80:ARG:HG3	3:A:390:HOH:O	2.14	0.47
1:A:179:VAL:HG11	1:A:195:VAL:HG22	1.94	0.47
1:A:121:HIS:HE1	3:A:265:HOH:O	1.98	0.47
1:A:1:MSE:H2	1:A:246:GLN:HG2	1.80	0.46
1:A:40:LEU:CD1	1:A:44:LYS:HE3	2.46	0.46
1:B:94:VAL:HG11	1:B:105:ARG:HD3	1.97	0.46
1:A:125:ASP:HA	1:A:150:GLN:HG2	1.99	0.45
1:B:244:ARG:O	1:B:248:LYS:HD2	2.16	0.45
1:A:158:LEU:HG	1:A:162:MSE:HE2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:47:ARG:HG2	1:B:87:PHE:CZ	2.52	0.45
1:A:239:LYS:HE3	3:A:331:HOH:O	2.17	0.45
1:B:239:LYS:O	1:B:243:GLU:HG3	2.18	0.44
1:B:199:ALA:HB2	1:B:238:MSE:HE1	1.99	0.44
1:B:245:HIS:O	1:B:249:LEU:HD13	2.17	0.44
1:A:47:ARG:HG2	1:A:87:PHE:CZ	2.52	0.44
1:A:203:GLN:O	1:A:226:SER:HA	2.18	0.44
1:A:1:MSE:N	1:A:246:GLN:HG2	2.32	0.44
1:B:103:MSE:O	1:B:107:GLU:HG3	2.18	0.43
1:A:58:ILE:HD11	1:A:82:VAL:HG21	1.99	0.43
1:B:1:MSE:HA	1:B:246:GLN:HG3	1.99	0.43
1:B:145:VAL:HG23	1:B:171:PRO:HB2	2.01	0.43
1:B:18:GLN:NE2	1:B:52:ILE:H	2.17	0.43
1:B:203:GLN:CG	1:B:227:ARG:HB2	2.48	0.43
1:B:103:MSE:HB2	1:B:104:PRO:HD3	2.00	0.42
1:B:248:LYS:HE3	1:B:248:LYS:CA	2.50	0.42
1:A:160:LYS:O	1:A:164:LEU:HG	2.20	0.42
1:A:9:TYR:CE2	1:A:31:PRO:HA	2.55	0.42
1:B:8:CYS:SG	1:B:14:ALA:HA	2.60	0.41
1:B:203:GLN:O	1:B:203:GLN:HG3	2.20	0.41
1:B:121:HIS:CE1	1:B:123:ALA:HB2	2.56	0.41
1:B:230:VAL:HG13	1:B:230:VAL:O	2.20	0.41
1:A:167:HIS:ND1	1:A:168:ARG:HG2	2.35	0.41
1:A:247:ALA:O	1:A:248:LYS:C	2.60	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)
3:B:601:HOH:O	3:B:601:HOH:O[4_566]	1.79	0.41

## 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	230/256 (90%)	224 (97%)	3 (1%)	3 (1%)	15	9
1	B	234/256 (91%)	227 (97%)	5 (2%)	2 (1%)	21	15
All	All	464/512 (91%)	451 (97%)	8 (2%)	5 (1%)	17	11

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	170	ALA
1	A	177	ALA
1	B	177	ALA
1	A	122	ARG
1	B	249	LEU

### 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	183/192 (95%)	176 (96%)	7 (4%)	40	40
1	B	187/192 (97%)	179 (96%)	8 (4%)	35	34
All	All	370/384 (96%)	355 (96%)	15 (4%)	37	36

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MSE
1	A	27	LEU
1	A	40	LEU
1	A	116	LEU
1	A	155	LEU
1	A	184	LEU
1	A	246	GLN
1	B	3	LEU
1	B	32	LYS

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Mol	Chain	Res	Type
1	B	121	HIS
1	B	146	LEU
1	B	184	LEU
1	B	224	GLU
1	B	239	LYS
1	B	248	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	48	GLN
1	A	121	HIS
1	A	135	ASN
1	A	136	ASN
1	A	150	GLN
1	A	156	GLN
1	A	185	HIS
1	A	196	HIS
1	B	18	GLN
1	B	48	GLN
1	B	55	HIS
1	B	167	HIS
1	B	196	HIS
1	B	203	GLN

### 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 1 ligands modelled in this entry, 1 is 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	224/256 (87%)	0.16	10 (4%) 37 46	7, 14, 34, 60	0
1	B	228/256 (89%)	0.26	18 (7%) 15 21	6, 13, 43, 66	0
All	All	452/512 (88%)	0.21	28 (6%) 24 32	6, 13, 36, 66	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	225	TYR	9.5
1	A	248	LYS	6.7
1	B	210	ARG	5.8
1	A	247	ALA	5.3
1	B	250	GLU	5.3
1	B	169	ASP	4.9
1	B	248	LYS	4.4
1	A	169	ASP	3.7
1	B	225	TYR	3.6
1	B	209	TYR	3.4
1	B	204	ALA	3.4
1	B	246	GLN	3.1
1	A	246	GLN	3.0
1	B	211	ASN	3.0
1	B	202	TRP	3.0
1	B	249	LEU	2.8
1	A	170	ALA	2.6
1	B	203	GLN	2.5
1	A	226	SER	2.4
1	A	210	ARG	2.4
1	B	170	ALA	2.3
1	A	168	ARG	2.2
1	A	167	HIS	2.1
1	B	139	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	166	ALA	2.1
1	B	208	ARG	2.1
1	B	201	ALA	2.1
1	B	168	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 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( $\text{\AA}^2$ )	Q<0.9
2	CA	B	501	1/1	0.99	0.07	-	26,26,26,26	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.