



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 4, 2017 – 03:46 PM EDT

PDB ID : 2B0L
Title : C-terminal DNA binding domain of transcriptional pleiotropic repressor CodY.
Authors : Levnikov, V.M.; Blagova, E.; Joseph, P.; Sonenshein, A.L.; Wilkinson, A.J.
Deposited on : unknown
Resolution : 2.90 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

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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-20030345
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-20030345

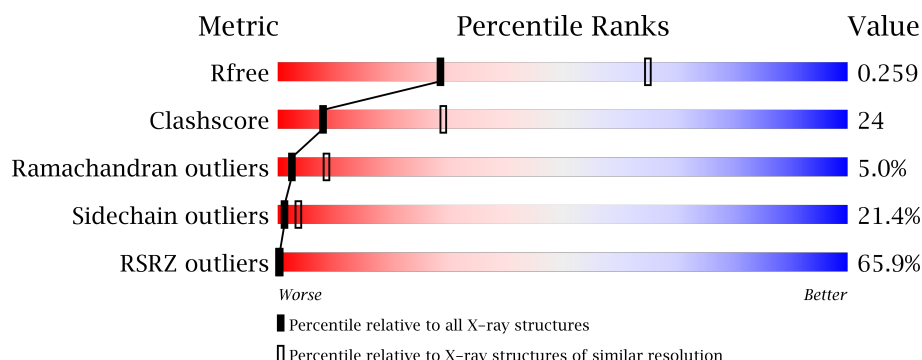
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1586 (2.90-2.90)
Clashscore	112137	1807 (2.90-2.90)
Ramachandran outliers	110173	1768 (2.90-2.90)
Sidechain outliers	110143	1770 (2.90-2.90)
RSRZ outliers	101464	1596 (2.90-2.90)

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 ≥ 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	102	<div> <div>32%</div> <div> <div>55%</div> <div>30%</div> <div>9%</div> <div>• •</div> </div> </div>
1	B	102	<div> <div>68%</div> <div> <div>50%</div> <div>33%</div> <div>10%</div> <div>7%</div> </div> </div>
1	C	102	<div> <div>85%</div> <div> <div>46%</div> <div>35%</div> <div>10%</div> <div>•</div> <div>8%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2300 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 GTP-sensing transcriptional pleiotropic repressor codY.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	98	Total	C	N	O	S	0	2	0
			783	493	141	145	4			
1	B	95	Total	C	N	O	S	4	4	0
			768	485	135	143	5			
1	C	94	Total	C	N	O	S	107	0	0
			730	461	128	138	3			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	158	GLY	-	CLONING ARTIFACT	UNP P39779
A	159	SER	-	CLONING ARTIFACT	UNP P39779
A	160	SER	-	CLONING ARTIFACT	UNP P39779
A	161	HIS	-	EXPRESSION TAG	UNP P39779
A	162	HIS	-	EXPRESSION TAG	UNP P39779
A	163	HIS	-	EXPRESSION TAG	UNP P39779
A	164	HIS	-	EXPRESSION TAG	UNP P39779
A	165	HIS	-	EXPRESSION TAG	UNP P39779
A	166	HIS	-	EXPRESSION TAG	UNP P39779
A	167	MET	-	INITIATING METHIONINE	UNP P39779
B	158	GLY	-	CLONING ARTIFACT	UNP P39779
B	159	SER	-	CLONING ARTIFACT	UNP P39779
B	160	SER	-	CLONING ARTIFACT	UNP P39779
B	161	HIS	-	EXPRESSION TAG	UNP P39779
B	162	HIS	-	EXPRESSION TAG	UNP P39779
B	163	HIS	-	EXPRESSION TAG	UNP P39779
B	164	HIS	-	EXPRESSION TAG	UNP P39779
B	165	HIS	-	EXPRESSION TAG	UNP P39779
B	166	HIS	-	EXPRESSION TAG	UNP P39779
B	167	MET	-	INITIATING METHIONINE	UNP P39779
C	158	GLY	-	CLONING ARTIFACT	UNP P39779
C	159	SER	-	CLONING ARTIFACT	UNP P39779
C	160	SER	-	CLONING ARTIFACT	UNP P39779

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Chain	Residue	Modelled	Actual	Comment	Reference
C	161	HIS	-	EXPRESSION TAG	UNP P39779
C	162	HIS	-	EXPRESSION TAG	UNP P39779
C	163	HIS	-	EXPRESSION TAG	UNP P39779
C	164	HIS	-	EXPRESSION TAG	UNP P39779
C	165	HIS	-	EXPRESSION TAG	UNP P39779
C	166	HIS	-	EXPRESSION TAG	UNP P39779
C	167	MET	-	INITIATING METHIONINE	UNP P39779

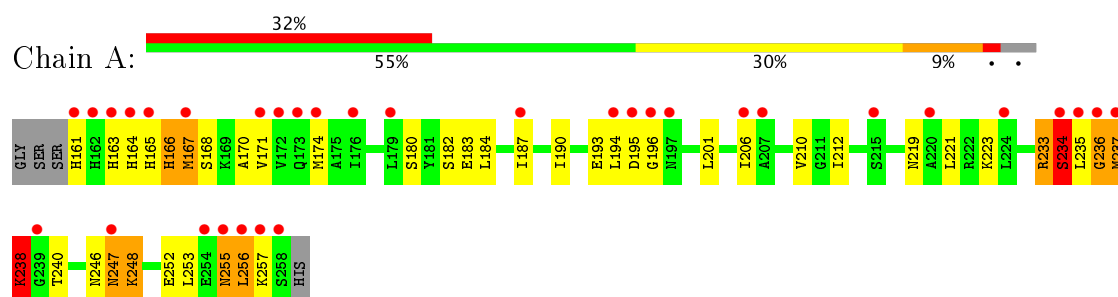
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	18	Total O 18 18	0	0
2	B	1	Total O 1 1	0	0

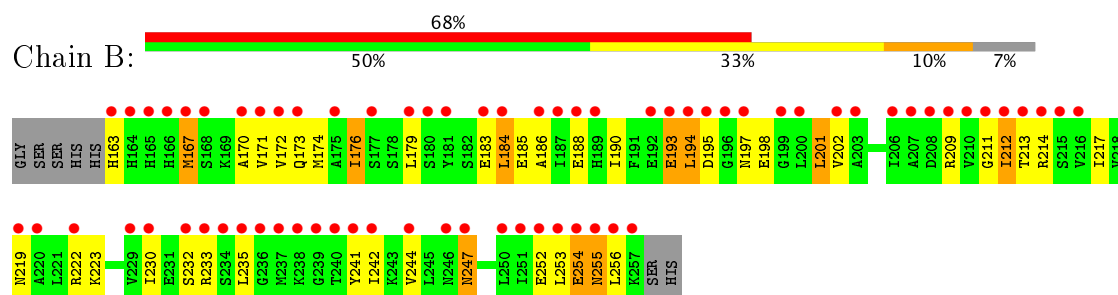
3 Residue-property plots

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.

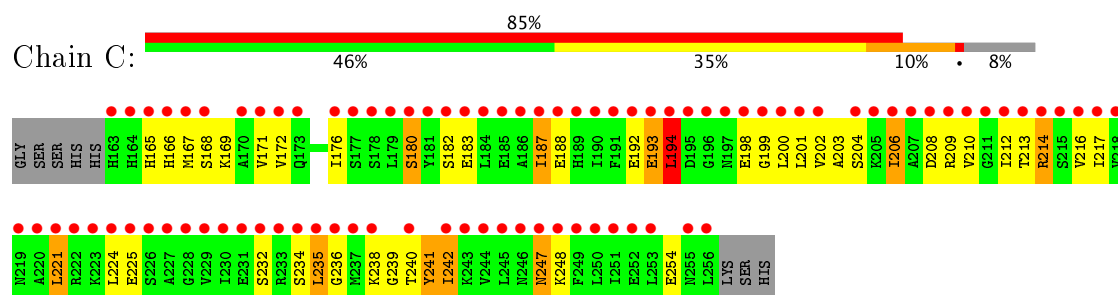
- Molecule 1: GTP-sensing transcriptional pleiotropic repressor codY



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- Molecule 1: GTP-sensing transcriptional pleiotropic repressor codY



4 Data and refinement statistics

Property	Value	Source
Space group	P 4 2 2	Depositor
Cell constants a, b, c, α , β , γ	68.13 Å 68.13 Å 164.90 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.90 9.99 – 2.90	Depositor EDS
% Data completeness (in resolution range)	100.0 (10.00-2.90) 93.8 (9.99-2.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.64 (at 2.89 Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.202 , 0.250 0.212 , 0.259	Depositor DCC
R_{free} test set	393 reflections (4.90%)	DCC
Wilson B-factor (Å ²)	77.5	Xtriage
Anisotropy	0.152	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.18 , 142.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	2300	wwPDB-VP
Average B, all atoms (Å ²)	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.57% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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

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	0.55	0/794	0.70	0/1065
1	B	0.47	0/780	0.70	0/1045
1	C	0.24	0/738	0.43	1/991 (0.1%)
All	All	0.44	0/2312	0.63	1/3101 (0.0%)

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	3
1	B	0	1
1	C	0	2
All	All	0	6

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	208	ASP	CB-CG-OD2	5.21	122.99	118.30

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	161	HIS	Peptide
1	A	233	ARG	Peptide
1	A	234	SER	Peptide
1	B	193	GLU	Peptide
1	C	210	VAL	Peptide
1	C	236	GLY	Peptide

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	783	0	799	48	0
1	B	768	0	795	44	0
1	C	730	0	753	23	0
2	A	18	0	0	0	0
2	B	1	0	0	0	0
All	All	2300	0	2347	103	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 24.

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:A:174[B]:MET:HE1	1:B:174[B]:MET:CA	1.62	1.28
1:A:174[B]:MET:CE	1:B:174[B]:MET:HA	1.70	1.19
1:A:174[B]:MET:CE	1:B:174[B]:MET:CA	2.28	1.07
1:A:174[B]:MET:HE1	1:B:174[B]:MET:CB	1.86	1.04
1:A:170:ALA:O	1:A:174[A]:MET:HG3	1.59	1.01
1:C:206:ILE:HD13	1:C:206:ILE:N	1.74	1.00
1:A:174[B]:MET:CE	1:B:174[B]:MET:CB	2.44	0.96
1:A:190:ILE:HG23	1:A:201:LEU:HD13	1.47	0.94
1:A:256:LEU:C	1:A:256:LEU:HD23	1.85	0.94
1:A:174[B]:MET:HE1	1:B:174[B]:MET:HB3	1.50	0.94
1:A:174[B]:MET:CE	1:B:174[B]:MET:HB3	2.02	0.89
1:A:237:MET:CG	1:A:238:LYS:H	1.88	0.85
1:A:174[B]:MET:HE2	1:B:174[B]:MET:HA	1.61	0.83
1:A:247:ASN:HD22	1:A:247:ASN:H	1.22	0.83
1:A:235:LEU:O	1:A:237:MET:HG2	1.80	0.81
1:A:237:MET:HG3	1:A:238:LYS:H	1.46	0.79
1:A:247:ASN:HD22	1:A:247:ASN:N	1.81	0.79
1:A:237:MET:HG3	1:A:238:LYS:HG2	1.68	0.75
1:B:172:VAL:O	1:B:176:ILE:HG22	1.85	0.75
1:C:247:ASN:H	1:C:247:ASN:HD22	1.34	0.74
1:B:170:ALA:O	1:B:174[B]:MET:HG3	1.92	0.69
1:B:186:ALA:O	1:B:190:ILE:HD12	1.92	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:188:GLU:O	1:C:192:GLU:HG2	1.94	0.67
1:C:206:ILE:HD13	1:C:206:ILE:H	1.57	0.66
1:B:190:ILE:HD11	1:B:217:ILE:HG23	1.79	0.64
1:B:214:ARG:HG2	1:B:214:ARG:O	1.97	0.63
1:C:167:MET:O	1:C:171:VAL:HG23	1.98	0.63
1:A:237:MET:HG3	1:A:238:LYS:CG	2.28	0.63
1:C:187:ILE:HG13	1:C:224:LEU:HD11	1.79	0.63
1:A:164[B]:HIS:CE1	1:A:167:MET:HB2	2.35	0.61
1:A:247:ASN:H	1:A:247:ASN:ND2	1.96	0.61
1:C:200:LEU:HA	1:C:241:TYR:HA	1.82	0.60
1:A:174[B]:MET:HE2	1:B:174[B]:MET:CB	2.31	0.60
1:A:235:LEU:O	1:A:236:GLY:C	2.41	0.59
1:C:172:VAL:O	1:C:176:ILE:HG13	2.03	0.59
1:A:164[B]:HIS:CG	1:A:164[B]:HIS:O	2.57	0.58
1:B:185:GLU:O	1:B:188:GLU:HB3	2.04	0.57
1:B:232:SER:HB3	1:B:242:ILE:HD13	1.85	0.57
1:A:190:ILE:HG23	1:A:201:LEU:CD1	2.30	0.57
1:A:237:MET:CG	1:A:238:LYS:N	2.62	0.57
1:C:206:ILE:N	1:C:206:ILE:CD1	2.48	0.57
1:C:247:ASN:N	1:C:247:ASN:HD22	2.02	0.57
1:B:230:ILE:HG22	1:B:244:VAL:HA	1.86	0.56
1:A:256:LEU:C	1:A:256:LEU:CD2	2.60	0.56
1:B:253:LEU:O	1:B:256:LEU:HG	2.05	0.56
1:B:253:LEU:HD12	1:B:256:LEU:CD1	2.37	0.55
1:B:201:LEU:HD12	1:B:202:VAL:H	1.72	0.55
1:C:193:GLU:O	1:C:194:LEU:C	2.46	0.54
1:A:165:HIS:O	1:A:166:HIS:C	2.46	0.54
1:A:256:LEU:HD23	1:A:256:LEU:O	2.07	0.54
1:A:180:SER:N	1:A:183:GLU:OE2	2.31	0.54
1:C:209:ARG:HG2	1:C:209:ARG:O	2.07	0.54
1:A:165:HIS:O	1:A:167:MET:N	2.42	0.53
1:A:165:HIS:O	1:A:168:SER:N	2.41	0.53
1:B:183:GLU:OE1	1:B:223:LYS:NZ	2.39	0.53
1:B:253:LEU:O	1:B:255:ASN:N	2.42	0.52
1:B:172:VAL:HG11	1:B:252:GLU:HB3	1.91	0.52
1:A:246:ASN:OD1	1:A:246:ASN:C	2.49	0.51
1:B:170:ALA:O	1:B:174[A]:MET:HG3	2.12	0.50
1:A:165:HIS:HB3	1:A:248:LYS:HD3	1.95	0.49
1:B:167[A]:MET:O	1:B:170:ALA:HB3	2.11	0.49
1:A:252:GLU:O	1:A:255:ASN:HB3	2.13	0.48
1:A:174[B]:MET:SD	1:B:174[B]:MET:HB3	2.54	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:198:GLU:HA	1:B:242:ILE:O	2.13	0.48
1:C:206:ILE:CD1	1:C:206:ILE:H	2.21	0.47
1:B:171:VAL:HA	1:B:174[A]:MET:SD	2.54	0.47
1:B:194:LEU:HG	1:B:195:ASP:H	1.79	0.47
1:C:232:SER:HA	1:C:241:TYR:O	2.15	0.47
1:C:168:SER:O	1:C:172:VAL:HG23	2.15	0.47
1:B:179:LEU:HD23	1:B:223:LYS:HD3	1.98	0.46
1:B:184:LEU:HD23	1:B:184:LEU:HA	1.82	0.46
1:B:253:LEU:HD12	1:B:256:LEU:HD11	1.98	0.46
1:B:194:LEU:HG	1:B:195:ASP:N	2.31	0.46
1:A:256:LEU:HD23	1:A:257:LYS:N	2.31	0.46
1:C:203:ALA:HB3	1:C:214:ARG:NE	2.31	0.45
1:B:214:ARG:CG	1:B:214:ARG:O	2.65	0.45
1:A:187:ILE:HD13	1:A:190:ILE:HD12	1.98	0.45
1:A:194:LEU:HD23	1:A:196:GLY:O	2.17	0.45
1:A:174[B]:MET:HE2	1:B:174[B]:MET:CG	2.47	0.44
1:B:253:LEU:HD12	1:B:256:LEU:HD12	1.99	0.44
1:A:174[B]:MET:HE2	1:B:174[B]:MET:CA	2.27	0.44
1:C:241:TYR:C	1:C:241:TYR:CD2	2.91	0.44
1:B:184:LEU:HD22	1:B:256:LEU:HD13	2.00	0.44
1:A:206:ILE:O	1:A:210:VAL:HG23	2.18	0.43
1:B:184:LEU:CD2	1:B:256:LEU:HD13	2.49	0.43
1:C:234:SER:O	1:C:235:LEU:C	2.57	0.42
1:C:199:GLY:O	1:C:242:ILE:N	2.51	0.42
1:C:234:SER:O	1:C:235:LEU:O	2.38	0.42
1:A:183:GLU:OE2	1:A:223:LYS:NZ	2.53	0.42
1:B:253:LEU:O	1:B:254:GLU:C	2.58	0.42
1:A:246:ASN:OD1	1:A:248:LYS:N	2.53	0.42
1:B:233:ARG:C	1:B:235:LEU:H	2.23	0.41
1:B:211:GLY:C	1:B:212:ILE:HG13	2.34	0.41
1:A:171:VAL:HA	1:A:174[A]:MET:SD	2.60	0.41
1:C:202:VAL:HG22	1:C:239:GLY:CA	2.51	0.41
1:A:233:ARG:O	1:A:234:SER:C	2.59	0.41
1:C:180:SER:H	1:C:183:GLU:HB2	1.85	0.41
1:B:232:SER:HA	1:B:241:TYR:O	2.20	0.40
1:A:247:ASN:N	1:A:247:ASN:ND2	2.53	0.40
1:B:176:ILE:HD13	1:B:176:ILE:C	2.42	0.40
1:C:221:LEU:O	1:C:225:GLU:HB2	2.21	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	98/102 (96%)	79 (81%)	13 (13%)	6 (6%)	2	5
1	B	97/102 (95%)	81 (84%)	12 (12%)	4 (4%)	3	13
1	C	92/102 (90%)	73 (79%)	15 (16%)	4 (4%)	3	12
All	All	287/306 (94%)	233 (81%)	40 (14%)	14 (5%)	2	9

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	166	HIS
1	A	237	MET
1	B	254	GLU
1	A	234	SER
1	A	236	GLY
1	A	255	ASN
1	B	194	LEU
1	B	247	ASN
1	C	194	LEU
1	C	235	LEU
1	A	238	LYS
1	B	197	ASN
1	C	193	GLU
1	C	180	SER

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	87/88 (99%)	71 (82%)	16 (18%)	2	6
1	B	86/88 (98%)	71 (83%)	15 (17%)	2	7
1	C	81/88 (92%)	58 (72%)	23 (28%)	0	1
All	All	254/264 (96%)	200 (79%)	54 (21%)	1	3

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	163	HIS
1	A	167	MET
1	A	182	SER
1	A	184	LEU
1	A	193	GLU
1	A	195	ASP
1	A	212	ILE
1	A	219	ASN
1	A	221	LEU
1	A	234	SER
1	A	238	LYS
1	A	240	THR
1	A	247	ASN
1	A	248	LYS
1	A	253	LEU
1	A	256	LEU
1	B	163	HIS
1	B	167[A]	MET
1	B	167[B]	MET
1	B	173	GLN
1	B	176	ILE
1	B	184	LEU
1	B	193	GLU
1	B	201	LEU
1	B	209	ARG
1	B	212	ILE
1	B	213	THR
1	B	219	ASN
1	B	222	ARG
1	B	247	ASN
1	B	255	ASN
1	C	165	HIS
1	C	166	HIS
1	C	169	LYS

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Mol	Chain	Res	Type
1	C	182	SER
1	C	187	ILE
1	C	194	LEU
1	C	198	GLU
1	C	201	LEU
1	C	204	SER
1	C	206	ILE
1	C	212	ILE
1	C	213	THR
1	C	214	ARG
1	C	216	VAL
1	C	217	ILE
1	C	221	LEU
1	C	238	LYS
1	C	240	THR
1	C	241	TYR
1	C	242	ILE
1	C	247	ASN
1	C	248	LYS
1	C	254	GLU

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

Mol	Chain	Res	Type
1	A	247	ASN
1	B	247	ASN
1	C	247	ASN

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 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 ⓘ

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, 95th 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 ²)	Q<0.9
1	A	98/102 (96%)	1.98	33 (33%) 0 0	10, 32, 79, 93	0
1	B	95/102 (93%)	3.25	69 (72%) 0 0	41, 75, 111, 116	1 (1%)
1	C	94/102 (92%)	5.50	87 (92%) 0 0	103, 127, 152, 154	37 (39%)
All	All	287/306 (93%)	3.56	189 (65%) 0 0	10, 79, 148, 154	38 (13%)

All (189) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	243	LYS	23.3
1	C	246	ASN	15.7
1	C	251	ILE	15.2
1	C	211	GLY	11.2
1	B	206	ILE	10.5
1	C	163	HIS	10.2
1	C	250	LEU	10.1
1	C	256	LEU	10.0
1	A	258	SER	9.8
1	C	237	MET	9.8
1	C	187	ILE	9.4
1	B	234	SER	9.4
1	C	164	HIS	9.3
1	C	216	VAL	8.8
1	C	232	SER	8.8
1	B	235	LEU	8.6
1	B	255	ASN	8.4
1	B	207	ALA	8.3
1	C	210	VAL	8.3
1	B	241	TYR	8.2
1	C	197	ASN	8.2
1	C	247	ASN	8.1
1	C	242	ILE	7.9

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Mol	Chain	Res	Type	RSRZ
1	C	185	GLU	7.7
1	C	255	ASN	7.7
1	C	215	SER	7.6
1	C	180	SER	7.5
1	C	193	GLU	7.5
1	C	182	SER	7.4
1	C	202	VAL	7.4
1	C	235	LEU	7.2
1	A	161	HIS	7.2
1	C	178	SER	6.9
1	C	244	VAL	6.9
1	B	250	LEU	6.7
1	C	168	SER	6.7
1	C	252	GLU	6.6
1	C	234	SER	6.3
1	A	164[A]	HIS	6.2
1	B	236	GLY	6.2
1	A	194	LEU	6.1
1	B	256	LEU	6.1
1	C	225	GLU	6.1
1	C	177	SER	6.0
1	C	188	GLU	5.9
1	A	196	GLY	5.9
1	A	162	HIS	5.9
1	B	163	HIS	5.8
1	B	165	HIS	5.8
1	C	181	TYR	5.8
1	C	200	LEU	5.7
1	C	194	LEU	5.7
1	C	230	ILE	5.7
1	C	184	LEU	5.7
1	C	186	ALA	5.5
1	B	194	LEU	5.5
1	A	234	SER	5.5
1	B	164	HIS	5.4
1	B	210	VAL	5.3
1	C	248	LYS	5.3
1	A	165	HIS	5.3
1	A	207	ALA	5.2
1	C	183	GLU	5.2
1	C	209	ARG	5.1
1	C	219	ASN	5.1

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Mol	Chain	Res	Type	RSRZ
1	C	207	ALA	5.1
1	C	165	HIS	5.0
1	B	195	ASP	5.0
1	B	247	ASN	4.9
1	C	208	ASP	4.9
1	C	236	GLY	4.8
1	C	220	ALA	4.8
1	B	240	THR	4.8
1	A	256	LEU	4.8
1	C	221	LEU	4.7
1	A	163	HIS	4.7
1	B	253	LEU	4.6
1	C	206	ILE	4.6
1	C	240	THR	4.6
1	C	171	VAL	4.6
1	C	204	SER	4.6
1	C	176	ILE	4.6
1	C	212	ILE	4.6
1	B	167[A]	MET	4.5
1	C	218	VAL	4.5
1	C	214	ARG	4.5
1	A	237	MET	4.5
1	B	239	GLY	4.5
1	C	217	ILE	4.5
1	B	257	LYS	4.5
1	C	198	GLU	4.4
1	C	249	PHE	4.4
1	A	235	LEU	4.3
1	C	222	ARG	4.3
1	C	229	VAL	4.2
1	C	201	LEU	4.2
1	C	227	ALA	4.1
1	B	232	SER	4.1
1	B	187	ILE	4.1
1	B	172	VAL	4.0
1	B	216	VAL	4.0
1	B	215	SER	3.9
1	B	220	ALA	3.9
1	B	180	SER	3.8
1	B	209	ARG	3.7
1	B	212	ILE	3.7
1	C	196	GLY	3.6

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Mol	Chain	Res	Type	RSRZ
1	C	189	HIS	3.6
1	C	245	LEU	3.6
1	A	174[A]	MET	3.6
1	C	223	LYS	3.5
1	B	238	LYS	3.5
1	B	233	ARG	3.5
1	C	170	ALA	3.5
1	C	228	GLY	3.4
1	B	254	GLU	3.4
1	C	195	ASP	3.4
1	A	236	GLY	3.4
1	B	193	GLU	3.4
1	C	226	SER	3.3
1	B	208	ASP	3.3
1	A	257	LYS	3.3
1	B	251	ILE	3.3
1	C	224	LEU	3.3
1	B	202	VAL	3.2
1	B	230	ILE	3.1
1	B	173	GLN	3.1
1	A	172	VAL	3.0
1	A	247	ASN	3.0
1	C	238	LYS	3.0
1	B	222	ARG	3.0
1	C	231	GLU	3.0
1	A	167	MET	3.0
1	C	253	LEU	3.0
1	B	183	GLU	3.0
1	B	181	TYR	2.9
1	B	184	LEU	2.9
1	B	179	LEU	2.9
1	B	188	GLU	2.8
1	C	190	ILE	2.8
1	B	197	ASN	2.8
1	B	229	VAL	2.7
1	B	211	GLY	2.7
1	B	192	GLU	2.7
1	C	205	LYS	2.7
1	C	179	LEU	2.7
1	C	166	HIS	2.6
1	C	199	GLY	2.6
1	B	242	ILE	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	254	GLU	2.6
1	C	167	MET	2.6
1	A	215	SER	2.6
1	C	213	THR	2.5
1	B	189	HIS	2.5
1	B	244	VAL	2.5
1	C	172	VAL	2.5
1	A	239	GLY	2.5
1	A	224	LEU	2.5
1	B	186	ALA	2.5
1	B	196	GLY	2.5
1	B	199	GLY	2.5
1	B	175	ALA	2.4
1	A	187	ILE	2.4
1	B	171	VAL	2.4
1	A	255	ASN	2.4
1	C	191	PHE	2.4
1	C	173	GLN	2.4
1	B	166[A]	HIS	2.3
1	C	192	GLU	2.3
1	B	246	ASN	2.3
1	B	252	GLU	2.3
1	A	220	ALA	2.3
1	B	214	ARG	2.3
1	B	170	ALA	2.3
1	B	203	ALA	2.3
1	B	213	THR	2.2
1	A	179	LEU	2.2
1	B	237	MET	2.2
1	B	200	LEU	2.2
1	A	206	ILE	2.2
1	B	168[A]	SER	2.2
1	A	173	GLN	2.2
1	B	219	ASN	2.1
1	C	233	ARG	2.1
1	B	177	SER	2.0
1	A	171	VAL	2.0
1	A	197	ASN	2.0
1	A	176	ILE	2.0
1	A	195	ASP	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.