



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 02:41 PM UTC

PDB ID : 8C7T / pdb_00008c7t
Title : Unliganded transcriptional pleiotropic repressor CodY from *Enterococcus faecalis*
Authors : Hainzl, T.; Sauer-Eriksson, A.E.
Deposited on : 2023-01-17
Resolution : 2.21 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

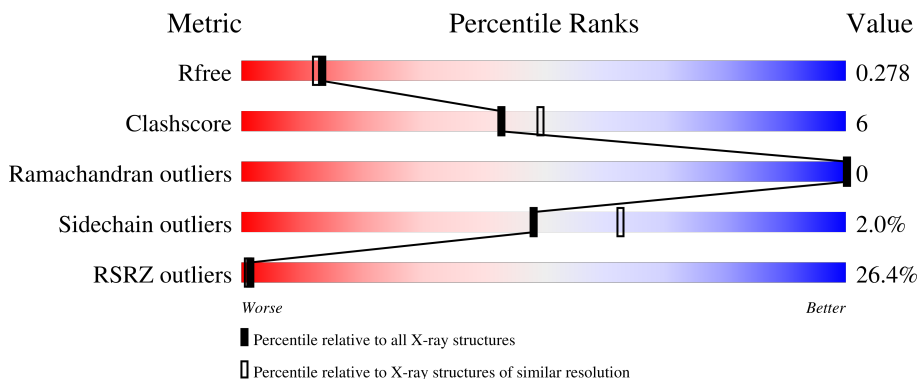
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.21 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}	180053	7682 (2.24-2.20)
Clashscore	190562	8402 (2.24-2.20)
Ramachandran outliers	187476	8303 (2.24-2.20)
Sidechain outliers	187428	8304 (2.24-2.20)
RSRZ outliers	180081	7683 (2.24-2.20)

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 ≥ 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	262	35% 84% 15%
1	B	262	26% 81% 18%
1	C	262	10% 55% 13% 32%
1	D	262	26% 86% 14%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7659 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	261	Total	C	N	O	S	0	0	0
			2031	1290	336	397	8			
1	B	262	Total	C	N	O	S	0	1	0
			2048	1301	339	400	8			
1	C	178	Total	C	N	O	S	0	0	0
			1398	893	228	270	7			
1	D	262	Total	C	N	O	S	0	0	0
			2041	1296	338	399	8			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP A0A1B4XP18
A	0	ALA	-	expression tag	UNP A0A1B4XP18
B	-1	GLY	-	expression tag	UNP A0A1B4XP18
B	0	ALA	-	expression tag	UNP A0A1B4XP18
C	-1	GLY	-	expression tag	UNP A0A1B4XP18
C	0	ALA	-	expression tag	UNP A0A1B4XP18
D	-1	GLY	-	expression tag	UNP A0A1B4XP18
D	0	ALA	-	expression tag	UNP A0A1B4XP18

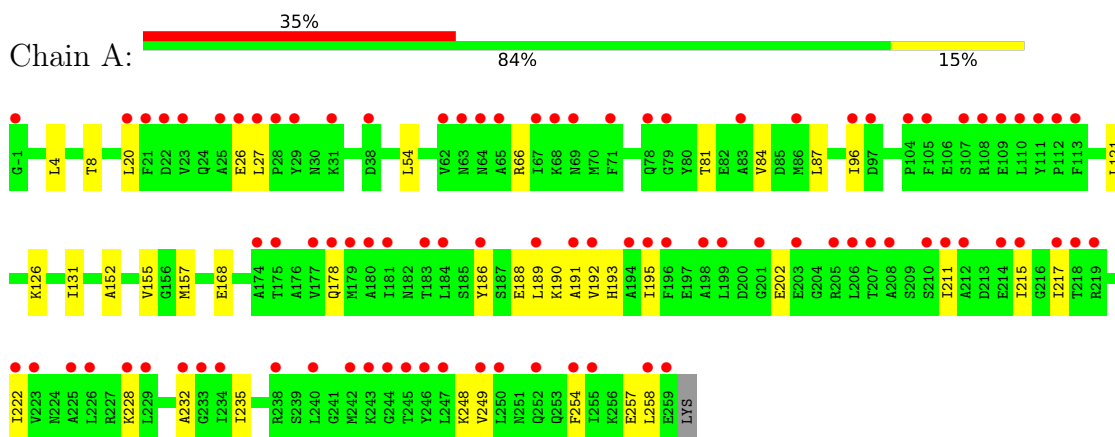
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	12	Total	O	0	0
			12	12		
2	B	53	Total	O	0	0
			53	53		
2	C	21	Total	O	0	0
			21	21		
2	D	55	Total	O	0	0
			55	55		

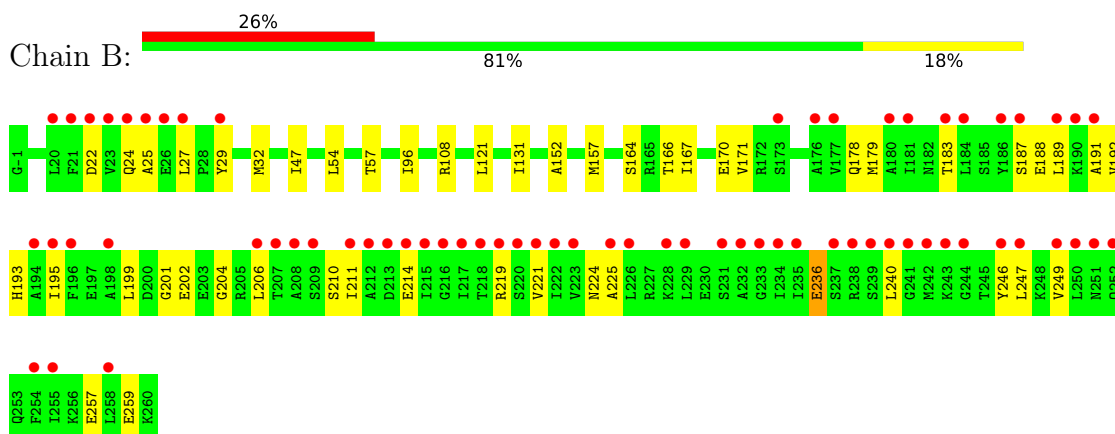
3 Residue-property plots [i](#)

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.

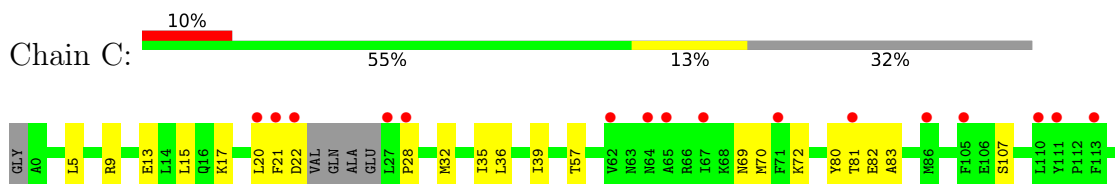
- Molecule 1: GTP-sensing transcriptional pleiotropic repressor CodY

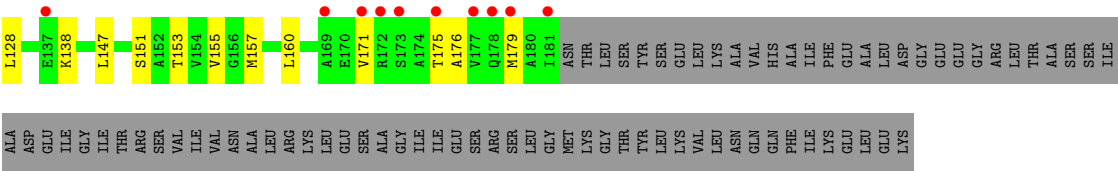


- Molecule 1: GTP-sensing transcriptional pleiotropic repressor CodY

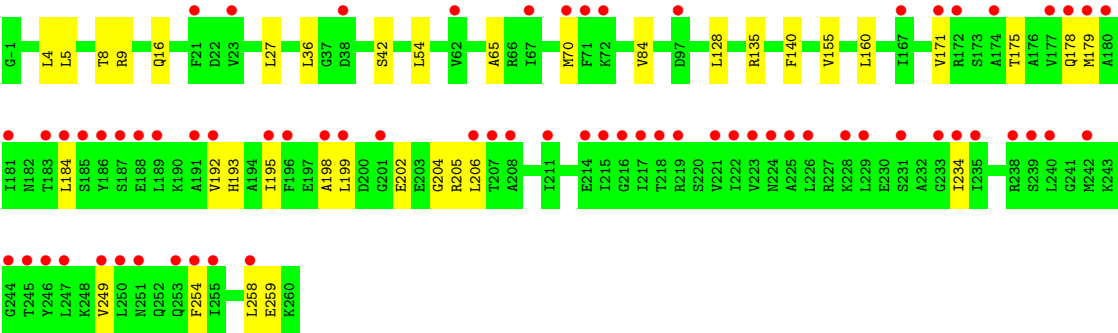
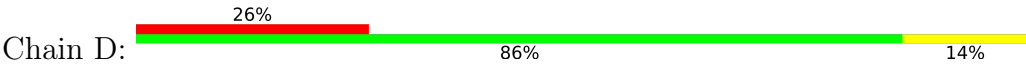


- Molecule 1: GTP-sensing transcriptional pleiotropic repressor CodY





● Molecule 1: GTP-sensing transcriptional pleiotropic repressor CodY



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	32.99Å 171.24Å 215.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	35.87 – 2.21 35.87 – 2.21	Depositor EDS
% Data completeness (in resolution range)	99.8 (35.87-2.21) 99.8 (35.87-2.21)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.49 (at 2.22Å)	Xtriage
Refinement program	PHENIX 1.17.1 _3660	Depositor
R, R_{free}	0.236 , 0.273 0.245 , 0.278	Depositor DCC
R_{free} test set	3191 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	41.9	Xtriage
Anisotropy	0.371	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 39.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7659	wwPDB-VP
Average B, all atoms (Å ²)	77.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.91% 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 [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	0.50	0/2057	0.68	2/2777 (0.1%)
1	B	0.60	1/2074 (0.0%)	0.77	0/2798
1	C	0.51	0/1418	0.64	0/1916
1	D	0.57	0/2067	0.69	0/2788
All	All	0.55	1/7616 (0.0%)	0.70	2/10279 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	47	ILE	CG1-CD1	-5.04	1.32	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	ARG	N-CA-C	-5.46	99.50	108.41
1	A	186	TYR	N-CA-CB	5.27	117.65	110.01

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	2031	0	2076	22	0
1	B	2048	0	2097	33	0
1	C	1398	0	1424	26	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2041	0	2089	25	0
2	A	12	0	0	0	0
2	B	53	0	0	0	0
2	C	21	0	0	0	0
2	D	55	0	0	0	0
All	All	7659	0	7686	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 6.

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:B:202:GLU:O	1:B:249:VAL:HG23	1.81	0.81
1:C:157:MET:CE	1:D:160:LEU:HD12	2.13	0.78
1:B:188:GLU:HB3	1:B:225:ALA:HB2	1.68	0.75
1:B:22:ASP:CG	1:B:25:ALA:HB2	2.14	0.73
1:B:187:SER:HB3	1:B:221:VAL:HG11	1.76	0.68
1:D:205:ARG:O	1:D:206:LEU:HD12	1.94	0.66
1:C:39:ILE:HG22	1:C:147:LEU:HD21	1.77	0.65
1:D:198:ALA:HB3	1:D:206:LEU:HD21	1.81	0.63
1:B:29:TYR:CD1	1:B:32:MET:HE2	2.33	0.63
1:B:187:SER:HB3	1:B:221:VAL:CG1	2.29	0.63
1:A:254:PHE:HE1	1:A:258:LEU:HD22	1.64	0.63
1:B:188:GLU:OE1	1:B:224:ASN:ND2	2.32	0.62
1:B:27:LEU:HB3	1:B:54:LEU:O	2.01	0.60
1:C:157:MET:HE2	1:D:160:LEU:HD12	1.81	0.60
1:C:176:ALA:HB2	1:D:179:MET:HE1	1.83	0.60
1:C:157:MET:HE3	1:D:160:LEU:HD12	1.83	0.58
1:A:190:LYS:HE3	1:A:215:ILE:HD11	1.85	0.57
1:D:4:LEU:O	1:D:8:THR:HG23	2.05	0.57
1:A:131:ILE:HD11	1:A:155:VAL:HG21	1.86	0.56
1:A:178:GLN:NE2	1:A:257:GLU:OE1	2.38	0.56
1:D:175:THR:HA	1:D:178:GLN:OE1	2.05	0.56
1:D:135:ARG:HG3	1:D:140:PHE:CZ	2.41	0.56
1:C:36:LEU:HD11	1:C:155:VAL:HG21	1.88	0.55
1:D:254:PHE:HE1	1:D:258:LEU:HD22	1.71	0.55
1:A:121:LEU:HD23	1:A:152:ALA:HB1	1.89	0.54
1:B:199:LEU:HD11	1:B:204:GLY:CA	2.38	0.54
1:B:188:GLU:O	1:B:192:VAL:HG23	2.07	0.54
1:B:188:GLU:O	1:B:191:ALA:HB3	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:LEU:HB3	1:A:54:LEU:O	2.09	0.53
1:A:87:LEU:HD12	1:A:87:LEU:N	2.24	0.52
1:C:175:THR:O	1:C:179:MET:HG2	2.09	0.52
1:B:240:LEU:HD11	1:B:246:TYR:HB2	1.91	0.51
1:B:24:GLN:HG3	1:B:24:GLN:O	2.11	0.50
1:D:199:LEU:HD11	1:D:204:GLY:N	2.27	0.50
1:B:166:THR:HG22	1:B:170:GLU:HG3	1.94	0.50
1:A:4:LEU:O	1:A:8:THR:HG23	2.13	0.49
1:C:107:SER:O	1:C:107:SER:OG	2.26	0.49
1:B:164:SER:O	1:B:167:ILE:HG22	2.13	0.48
1:B:201:GLY:O	1:B:202:GLU:HB2	2.12	0.48
1:C:175:THR:OG1	1:D:175:THR:HG21	2.13	0.48
1:C:5:LEU:O	1:C:9:ARG:HG2	2.15	0.47
1:A:168:GLU:HG3	1:B:171:VAL:HG11	1.97	0.47
1:C:21:PHE:O	1:C:22:ASP:HB3	2.15	0.47
1:A:81:THR:O	1:A:84:VAL:HG22	2.15	0.46
1:C:36:LEU:HD22	1:C:151:SER:HB3	1.97	0.46
1:B:192:VAL:HA	1:B:195:ILE:HD12	1.97	0.46
1:B:57:THR:O	1:B:57:THR:HG23	2.15	0.46
1:B:211:ILE:HD12	1:B:214:GLU:OE1	2.16	0.46
1:B:121:LEU:HD21	1:B:131:ILE:HG13	1.97	0.46
1:B:178:GLN:NE2	1:B:257:GLU:OE1	2.49	0.45
1:A:192:VAL:HA	1:A:195:ILE:HD12	1.98	0.45
1:C:20:LEU:O	1:C:28:PRO:HB3	2.17	0.45
1:B:96:ILE:HG12	1:B:108:ARG:HG2	1.98	0.45
1:D:184:LEU:HD11	1:D:192:VAL:HG21	1.98	0.45
1:A:191:ALA:HB1	1:A:222:ILE:HD13	1.99	0.45
1:B:199:LEU:HD11	1:B:204:GLY:C	2.42	0.45
1:A:191:ALA:O	1:A:195:ILE:HG13	2.18	0.44
1:B:199:LEU:HD11	1:B:204:GLY:N	2.32	0.44
1:C:157:MET:HE1	1:D:128:LEU:HD11	1.99	0.44
1:A:235:ILE:HA	1:A:248:LYS:O	2.17	0.44
1:D:36:LEU:HD11	1:D:155:VAL:CG2	2.48	0.44
1:C:157:MET:HE1	1:D:128:LEU:CD1	2.47	0.43
1:C:57:THR:O	1:C:57:THR:HG23	2.17	0.43
1:C:171:VAL:HB	1:D:171:VAL:HG21	2.00	0.43
1:C:153:THR:HG22	1:D:16:GLN:NE2	2.34	0.43
1:A:188:GLU:OE1	1:A:228:LYS:NZ	2.46	0.43
1:B:166:THR:HG22	1:B:170:GLU:CD	2.43	0.43
1:C:15:LEU:HD21	1:C:35:ILE:HD13	2.00	0.43
1:C:69:ASN:HA	1:C:72:LYS:HE3	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:192:VAL:CG2	1:B:225:ALA:HB1	2.49	0.42
1:B:179:MET:O	1:B:183:THR:HG23	2.19	0.42
1:C:32:MET:HB3	1:C:32:MET:HE2	1.74	0.42
1:D:192:VAL:HA	1:D:195:ILE:HD12	2.01	0.42
1:D:234:ILE:HG22	1:D:234:ILE:O	2.19	0.42
1:D:202:GLU:O	1:D:249:VAL:HG23	2.19	0.42
1:B:189:LEU:C	1:B:189:LEU:HD23	2.45	0.42
1:C:157:MET:HE2	1:C:157:MET:HB3	1.92	0.42
1:A:121:LEU:CD2	1:A:152:ALA:HB1	2.49	0.42
1:A:189:LEU:C	1:A:189:LEU:HD23	2.45	0.42
1:B:236:GLU:O	1:B:247:LEU:HA	2.19	0.42
1:D:65:ALA:HB1	1:D:70:MET:HE2	2.01	0.42
1:C:81:THR:OG1	1:C:82:GLU:OE1	2.37	0.41
1:A:202:GLU:O	1:A:249:VAL:HG23	2.21	0.41
1:A:211:ILE:O	1:A:215:ILE:HG22	2.21	0.41
1:A:157:MET:HG3	1:B:157:MET:HG3	2.03	0.41
1:A:232:ALA:HB1	1:B:179:MET:HE2	2.03	0.41
1:B:121:LEU:HD23	1:B:152:ALA:HB1	2.01	0.41
1:D:42:SER:OG	1:D:135:ARG:HD3	2.21	0.41
1:A:26:GLU:O	1:A:27:LEU:HD23	2.21	0.41
1:C:13:GLU:OE1	1:C:17:LYS:HE2	2.21	0.41
1:C:80:TYR:O	1:C:83:ALA:HB3	2.22	0.40
1:C:128:LEU:HD21	1:C:160:LEU:HD13	2.03	0.40
1:D:5:LEU:O	1:D:9:ARG:HG2	2.21	0.40
1:D:27:LEU:HB3	1:D:54:LEU:O	2.22	0.40

There are no symmetry-related clashes.

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	259/262 (99%)	255 (98%)	4 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	261/262 (100%)	256 (98%)	5 (2%)	0	100	100
1	C	174/262 (66%)	171 (98%)	3 (2%)	0	100	100
1	D	260/262 (99%)	257 (99%)	3 (1%)	0	100	100
All	All	954/1048 (91%)	939 (98%)	15 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	221/222 (100%)	216 (98%)	5 (2%)	44	58
1	B	223/222 (100%)	217 (97%)	6 (3%)	39	52
1	C	153/222 (69%)	151 (99%)	2 (1%)	61	75
1	D	222/222 (100%)	219 (99%)	3 (1%)	59	73
All	All	819/888 (92%)	803 (98%)	16 (2%)	48	63

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	LEU
1	A	96	ILE
1	A	126	LYS
1	A	193	HIS
1	A	217	ILE
1	B	193	HIS
1	B	206	LEU
1	B	210	SER
1	B	219	ARG
1	B	236	GLU
1	B	259	GLU
1	C	70	MET
1	C	138	LYS

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Mol	Chain	Res	Type
1	D	84	VAL
1	D	193	HIS
1	D	259	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	163	GLN
1	C	16	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 oligosaccharides 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	261/262 (99%)	1.48	91 (34%)	1 0	32, 76, 131, 156	0
1	B	262/262 (100%)	1.10	68 (25%)	1 1	21, 55, 171, 190	1 (0%)
1	C	178/262 (67%)	0.87	26 (14%)	6 4	35, 56, 135, 158	0
1	D	262/262 (100%)	1.08	69 (26%)	1 1	29, 57, 145, 158	0
All	All	963/1048 (91%)	1.15	254 (26%)	1 1	21, 59, 152, 190	1 (0%)

All (254) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	254	PHE	5.4
1	B	221	VAL	5.3
1	B	234	ILE	5.1
1	D	208	ALA	5.1
1	D	186	TYR	5.1
1	B	226	LEU	5.0
1	B	222	ILE	4.9
1	B	223	VAL	4.7
1	A	20	LEU	4.6
1	A	186	TYR	4.3
1	B	27	LEU	4.3
1	D	180	ALA	4.3
1	D	245	THR	4.3
1	B	187	SER	4.2
1	A	62	VAL	4.2
1	B	240	LEU	4.1
1	A	21	PHE	4.1
1	A	67	ILE	4.1
1	A	110	LEU	4.1
1	D	189	LEU	4.1
1	D	226	LEU	4.1

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Mol	Chain	Res	Type	RSRZ
1	D	171	VAL	4.0
1	B	208	ALA	4.0
1	B	195	ILE	4.0
1	A	181	ILE	3.9
1	B	21	PHE	3.9
1	B	23	VAL	3.9
1	D	223	VAL	3.9
1	A	208	ALA	3.9
1	A	232	ALA	3.9
1	B	238	ARG	3.9
1	A	211	ILE	3.8
1	A	217	ILE	3.8
1	A	218	THR	3.8
1	B	225	ALA	3.8
1	A	108	ARG	3.8
1	A	233	GLY	3.8
1	B	250	LEU	3.8
1	B	235	ILE	3.8
1	C	137	GLU	3.8
1	A	65	ALA	3.7
1	B	191	ALA	3.7
1	D	258	LEU	3.7
1	A	26	GLU	3.7
1	D	229	LEU	3.6
1	A	180	ALA	3.6
1	D	188	GLU	3.6
1	B	233	GLY	3.6
1	D	242	MET	3.6
1	B	218	THR	3.6
1	B	217	ILE	3.6
1	B	184	LEU	3.6
1	A	23	VAL	3.6
1	A	206	LEU	3.5
1	B	212	ALA	3.5
1	B	241	GLY	3.5
1	D	254	PHE	3.5
1	A	63	ASN	3.5
1	B	186	TYR	3.5
1	B	26	GLU	3.5
1	B	229	LEU	3.5
1	D	250	LEU	3.4
1	A	113	PHE	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	245	THR	3.4
1	B	219	ARG	3.4
1	B	194	ALA	3.4
1	A	178	GLN	3.4
1	A	184	LEU	3.4
1	B	189	LEU	3.4
1	D	181	ILE	3.4
1	D	247	LEU	3.4
1	A	258	LEU	3.3
1	A	189	LEU	3.3
1	D	183	THR	3.3
1	D	70	MET	3.3
1	A	27	LEU	3.3
1	D	240	LEU	3.2
1	B	196	PHE	3.2
1	D	184	LEU	3.2
1	A	234	ILE	3.2
1	B	25	ALA	3.2
1	B	249	VAL	3.2
1	B	215	ILE	3.2
1	D	238	ARG	3.1
1	B	180	ALA	3.1
1	A	112	PRO	3.1
1	B	206	LEU	3.1
1	D	206	LEU	3.1
1	D	218	THR	3.1
1	C	177	VAL	3.1
1	D	217	ILE	3.1
1	D	255	ILE	3.1
1	C	105	PHE	3.1
1	B	242	MET	3.1
1	D	222	ILE	3.1
1	A	183	THR	3.1
1	D	221	VAL	3.1
1	A	259	GLU	3.0
1	A	105	PHE	3.0
1	C	113	PHE	3.0
1	D	233	GLY	3.0
1	D	251	ASN	3.0
1	A	68	LYS	3.0
1	D	211	ILE	3.0
1	D	215	ILE	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	-1	GLY	3.0
1	B	232	ALA	3.0
1	D	224	ASN	3.0
1	A	71	PHE	3.0
1	C	27	LEU	3.0
1	C	71	PHE	2.9
1	A	247	LEU	2.9
1	D	192	VAL	2.9
1	A	78	GLN	2.9
1	B	20	LEU	2.9
1	A	177	VAL	2.9
1	D	219	ARG	2.9
1	C	67	ILE	2.9
1	D	235	ILE	2.9
1	A	210	SER	2.9
1	D	216	GLY	2.9
1	A	219	ARG	2.9
1	B	247	LEU	2.9
1	D	195	ILE	2.9
1	D	97	ASP	2.9
1	D	246	TYR	2.8
1	A	196	PHE	2.8
1	B	255	ILE	2.8
1	B	239	SER	2.8
1	D	244	GLY	2.8
1	A	250	LEU	2.8
1	A	25	ALA	2.8
1	A	174	ALA	2.8
1	A	242	MET	2.8
1	A	96	ILE	2.8
1	A	226	LEU	2.8
1	C	111	TYR	2.8
1	B	243	LYS	2.8
1	A	215	ILE	2.7
1	C	181	ILE	2.7
1	D	187	SER	2.7
1	D	72	LYS	2.7
1	B	198	ALA	2.7
1	C	65	ALA	2.7
1	A	97	ASP	2.7
1	A	222	ILE	2.7
1	B	213	ASP	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	28	PRO	2.7
1	A	199	LEU	2.7
1	D	225	ALA	2.7
1	B	209	SER	2.7
1	D	179	MET	2.7
1	C	178	GLN	2.7
1	B	173	SER	2.7
1	C	169	ALA	2.6
1	A	31	LYS	2.6
1	B	183	THR	2.6
1	A	192	VAL	2.6
1	C	171	VAL	2.6
1	B	220	SER	2.6
1	B	190	LYS	2.6
1	C	21	PHE	2.6
1	D	207	THR	2.6
1	B	237	SER	2.6
1	D	234	ILE	2.6
1	B	29	TYR	2.6
1	D	174	ALA	2.5
1	A	254	PHE	2.5
1	D	196	PHE	2.5
1	A	249	VAL	2.5
1	D	62	VAL	2.5
1	A	109	GLU	2.5
1	B	207	THR	2.5
1	B	246	TYR	2.5
1	C	20	LEU	2.5
1	C	62	VAL	2.5
1	A	243	LYS	2.5
1	A	104	PRO	2.5
1	D	67	ILE	2.4
1	A	64	ASN	2.4
1	A	69	ASN	2.4
1	D	231	SER	2.4
1	A	86	MET	2.4
1	B	258	LEU	2.4
1	A	246	TYR	2.4
1	B	252	GLN	2.4
1	A	191	ALA	2.4
1	B	211	ILE	2.4
1	D	23	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	179	MET	2.4
1	A	111	TYR	2.4
1	A	175	THR	2.4
1	A	223	VAL	2.4
1	A	255	ILE	2.4
1	D	172	ARG	2.4
1	D	253	GLN	2.4
1	A	38	ASP	2.3
1	C	175	THR	2.3
1	B	181	ILE	2.3
1	D	249	VAL	2.3
1	B	22	ASP	2.3
1	A	107	SER	2.3
1	D	178	GLN	2.3
1	D	199	LEU	2.3
1	A	79	GLY	2.3
1	D	71	PHE	2.3
1	A	22	ASP	2.3
1	A	198	ALA	2.3
1	D	198	ALA	2.3
1	A	214	GLU	2.3
1	D	167	ILE	2.3
1	A	229	LEU	2.2
1	D	177	VAL	2.3
1	B	251	ASN	2.2
1	C	22	ASP	2.2
1	D	38	ASP	2.2
1	A	201	GLY	2.2
1	A	238	ARG	2.2
1	C	172	ARG	2.2
1	A	212	ALA	2.2
1	A	207	THR	2.2
1	B	24	GLN	2.2
1	A	29	TYR	2.2
1	A	83	ALA	2.2
1	C	81	THR	2.2
1	A	195	ILE	2.2
1	A	240	LEU	2.2
1	C	110	LEU	2.2
1	B	244	GLY	2.2
1	C	173	SER	2.2
1	D	239	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	205	ARG	2.1
1	D	201	GLY	2.1
1	B	231	SER	2.1
1	D	191	ALA	2.1
1	C	86	MET	2.1
1	C	179	MET	2.1
1	A	228	LYS	2.1
1	B	214	GLU	2.1
1	C	28	PRO	2.1
1	B	216	GLY	2.1
1	A	225	ALA	2.1
1	D	228	LYS	2.1
1	C	64	ASN	2.1
1	B	177	VAL	2.1
1	B	228	LYS	2.0
1	B	176	ALA	2.0
1	A	252	GLN	2.0
1	A	244	GLY	2.0
1	D	21	PHE	2.0
1	A	203	GLU	2.0
1	D	185	SER	2.0
1	D	214	GLU	2.0
1	A	194	ALA	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 oligosaccharides 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.