



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:10 PM UTC

PDB ID : 3E7D / pdb_00003e7d
Title : Crystal Structure of Precorrin-8X Methyl Mutase CbiC/CobH from *Brucella melitensis*
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2008-08-18
Resolution : 1.80 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

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

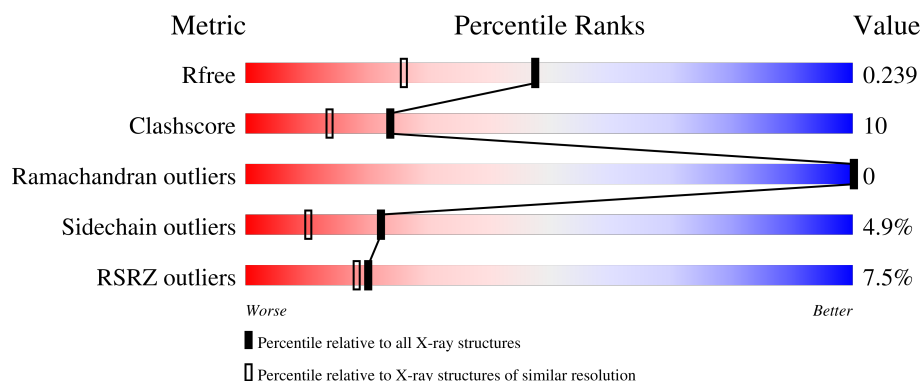
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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}	180053	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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 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	212	6% 83% 12% 5%
1	B	212	7% 81% 13% • •
1	C	212	5% 71% 22% • •
1	D	212	12% 71% 23% • •

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6557 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 CobH, precorrin-8X methylmutase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	202	Total	C	N	O	S	0	0	0
			1469	926	265	272	6			
1	B	204	Total	C	N	O	S	0	0	0
			1487	936	273	272	6			
1	C	203	Total	C	N	O	S	0	0	0
			1483	934	268	275	6			
1	D	204	Total	C	N	O	S	0	0	0
			1470	926	268	270	6			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP Q57CK9
A	-2	PRO	-	expression tag	UNP Q57CK9
A	-1	GLY	-	expression tag	UNP Q57CK9
A	0	SER	-	expression tag	UNP Q57CK9
B	-3	GLY	-	expression tag	UNP Q57CK9
B	-2	PRO	-	expression tag	UNP Q57CK9
B	-1	GLY	-	expression tag	UNP Q57CK9
B	0	SER	-	expression tag	UNP Q57CK9
C	-3	GLY	-	expression tag	UNP Q57CK9
C	-2	PRO	-	expression tag	UNP Q57CK9
C	-1	GLY	-	expression tag	UNP Q57CK9
C	0	SER	-	expression tag	UNP Q57CK9
D	-3	GLY	-	expression tag	UNP Q57CK9
D	-2	PRO	-	expression tag	UNP Q57CK9
D	-1	GLY	-	expression tag	UNP Q57CK9
D	0	SER	-	expression tag	UNP Q57CK9

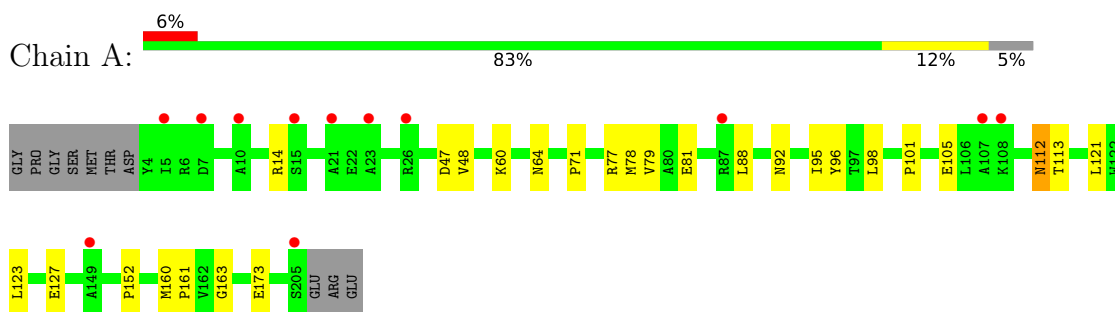
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	195	Total 195	O 195	0	0
2	B	173	Total 173	O 173	0	0
2	C	165	Total 165	O 165	0	0
2	D	115	Total 115	O 115	0	0

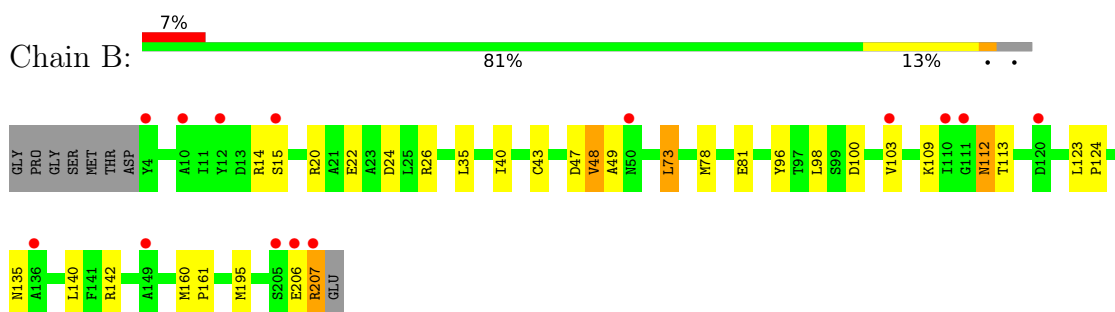
3 Residue-property plots [i](#)

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.

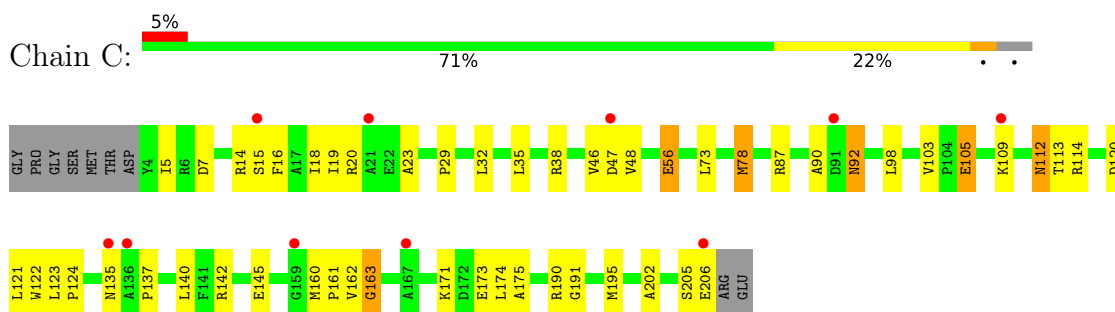
- Molecule 1: CobH, precorrin-8X methylmutase



- Molecule 1: CobH, precorrin-8X methylmutase

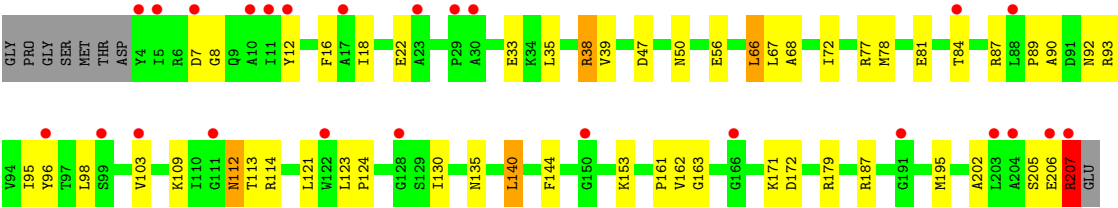


- Molecule 1: CobH, precorrin-8X methylmutase



- Molecule 1: CobH, precorrin-8X methylmutase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	57.72Å 68.93Å 103.20Å 90.00° 95.88° 90.00°	Depositor
Resolution (Å)	50.00 – 1.80 50.00 – 1.80	Depositor EDS
% Data completeness (in resolution range)	91.2 (50.00-1.80) 91.2 (50.00-1.80)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 1.81Å)	Xtriage
Refinement program	REFMAC refmac_5.4.0067	Depositor
R, R_{free}	0.232 , 0.260 0.238 , 0.239	Depositor DCC
R_{free} test set	3392 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	24.3	Xtriage
Anisotropy	0.105	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 48.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6557	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.69% 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	0.76	2/1492 (0.1%)	1.04	2/2028 (0.1%)
1	B	0.79	1/1510 (0.1%)	0.98	0/2051
1	C	0.98	5/1507 (0.3%)	1.07	2/2048 (0.1%)
1	D	0.83	0/1494	1.10	4/2033 (0.2%)
All	All	0.85	8/6003 (0.1%)	1.05	8/8160 (0.1%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	171	LYS	C-O	-7.35	1.15	1.24
1	A	77	ARG	C-O	-6.35	1.16	1.24
1	C	7	ASP	C-O	-5.82	1.17	1.24
1	C	175	ALA	C-O	-5.64	1.17	1.24
1	A	79	VAL	C-O	-5.51	1.18	1.24
1	B	142	ARG	C-O	-5.44	1.17	1.24
1	C	174	LEU	C-O	-5.23	1.18	1.24
1	C	5	ILE	CA-CB	-5.12	1.48	1.54

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	163	GLY	N-CA-C	6.58	120.32	110.97
1	D	68	ALA	N-CA-C	-5.69	106.19	113.02
1	C	163	GLY	N-CA-C	5.66	119.01	110.97
1	D	205	SER	N-CA-C	5.50	115.64	107.88
1	D	130	ILE	CB-CA-C	-5.40	103.58	110.98
1	C	173	GLU	CA-CB-CG	-5.26	103.58	114.10
1	D	207	ARG	NE-CZ-NH2	5.17	123.85	119.20
1	A	79	VAL	N-CA-C	-5.09	105.74	110.53

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	1469	0	1496	18	0
1	B	1487	0	1518	26	0
1	C	1483	0	1507	39	0
1	D	1470	0	1479	41	0
2	A	195	0	0	1	0
2	B	173	0	0	1	0
2	C	165	0	0	1	0
2	D	115	0	0	1	0
All	All	6557	0	6000	120	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 10.

All (120) 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:D:206:GLU:O	1:D:207:ARG:HB2	1.41	1.15
1:A:112:ASN:HD22	1:A:113:THR:H	1.11	0.96
1:C:87:ARG:NH1	1:C:205:SER:HB3	1.82	0.94
1:C:112:ASN:HD22	1:C:113:THR:H	1.21	0.89
1:B:24:ASP:OD1	1:B:26:ARG:HD3	1.76	0.85
1:B:20:ARG:HH12	1:B:207:ARG:NH2	1.76	0.83
1:C:121:LEU:HD12	1:C:121:LEU:N	1.95	0.81
1:A:112:ASN:ND2	1:A:113:THR:H	1.82	0.77
1:D:153:LYS:HE3	1:D:179:ARG:O	1.88	0.73
1:A:112:ASN:HD22	1:A:113:THR:N	1.86	0.73
1:C:112:ASN:ND2	1:C:113:THR:H	1.87	0.71
1:B:206:GLU:O	1:B:207:ARG:CB	2.39	0.71
1:D:12:TYR:CZ	1:D:16:PHE:HE1	2.09	0.71
1:C:47:ASP:OD1	1:C:47:ASP:C	2.35	0.70
1:D:39:VAL:HG13	1:D:195:MET:HG2	1.75	0.68
1:C:112:ASN:HD22	1:C:113:THR:N	1.89	0.68
1:D:93:ARG:HG2	1:D:95:ILE:HD11	1.76	0.66
1:A:78:MET:HE1	1:B:81:GLU:HB3	1.76	0.66
1:C:56:GLU:OE2	2:C:390:HOH:O	2.13	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:12:TYR:CZ	1:D:16:PHE:CE1	2.84	0.66
1:C:105:GLU:OE2	1:C:109:LYS:HE3	1.95	0.65
1:D:206:GLU:O	1:D:207:ARG:CB	2.29	0.65
1:D:161:PRO:HD2	1:D:171:LYS:HD3	1.79	0.64
1:C:121:LEU:N	1:C:121:LEU:CD1	2.61	0.63
1:C:92:ASN:H	1:C:92:ASN:HD22	1.47	0.62
1:A:78:MET:HA	1:A:78:MET:HE2	1.79	0.62
1:D:18:ILE:O	1:D:22:GLU:HG3	1.99	0.62
1:D:12:TYR:CE1	1:D:16:PHE:HE1	2.18	0.62
1:B:112:ASN:HD22	1:B:113:THR:H	1.48	0.61
1:D:66:LEU:HD21	1:D:72:ILE:HD11	1.83	0.61
1:D:12:TYR:CE1	1:D:16:PHE:CE1	2.90	0.60
1:B:40:ILE:HG12	1:B:48:VAL:HG12	1.84	0.59
1:C:16:PHE:O	1:C:20:ARG:HG3	2.03	0.58
1:A:101:PRO:HG2	2:A:303:HOH:O	2.02	0.57
1:B:206:GLU:O	1:B:207:ARG:HB2	2.04	0.57
1:B:112:ASN:ND2	1:B:113:THR:H	2.03	0.57
1:C:47:ASP:OD1	1:C:48:VAL:N	2.37	0.57
1:D:89:PRO:HD2	1:D:92:ASN:ND2	2.18	0.57
1:D:123:LEU:HB2	1:D:124:PRO:HD3	1.87	0.57
1:C:92:ASN:H	1:C:92:ASN:ND2	2.03	0.56
1:A:88:LEU:HB3	1:A:92:ASN:HD22	1.72	0.55
1:C:78:MET:HE1	1:D:78:MET:HE2	1.88	0.55
1:D:103:VAL:HG12	1:D:114:ARG:HG2	1.87	0.55
1:D:93:ARG:CG	1:D:95:ILE:HD11	2.37	0.54
1:D:112:ASN:HD22	1:D:113:THR:H	1.55	0.53
1:C:103:VAL:HG12	1:C:114:ARG:HG2	1.90	0.53
1:D:121:LEU:N	1:D:121:LEU:HD12	2.22	0.53
1:B:207:ARG:HG3	2:B:329:HOH:O	2.07	0.53
1:D:50:ASN:OD1	2:D:276:HOH:O	2.19	0.53
1:D:140:LEU:HD22	1:D:144:PHE:CZ	2.45	0.52
1:C:78:MET:HE2	1:C:78:MET:HA	1.90	0.52
1:D:12:TYR:CE2	1:D:16:PHE:HE1	2.27	0.52
1:B:35:LEU:C	1:B:35:LEU:HD23	2.35	0.52
1:C:47:ASP:OD1	1:C:190:ARG:NH1	2.41	0.51
1:C:105:GLU:HA	1:C:105:GLU:OE1	2.10	0.51
1:A:173:GLU:HA	1:A:173:GLU:OE1	2.11	0.50
1:B:20:ARG:NH1	1:B:207:ARG:NH2	2.54	0.50
1:D:35:LEU:HA	1:D:202:ALA:CB	2.41	0.50
1:A:96:TYR:CZ	1:A:98:LEU:HB3	2.47	0.49
1:C:191:GLY:O	1:C:195:MET:HE3	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:89:PRO:HD2	1:D:92:ASN:HD21	1.76	0.49
1:D:96:TYR:CZ	1:D:98:LEU:HB3	2.48	0.49
1:D:90:ALA:C	1:D:92:ASN:N	2.70	0.49
1:D:38:ARG:HH22	1:D:207:ARG:HD2	1.77	0.48
1:B:112:ASN:HD22	1:B:113:THR:N	2.11	0.47
1:D:121:LEU:N	1:D:121:LEU:CD1	2.77	0.47
1:B:206:GLU:O	1:B:207:ARG:HB3	2.15	0.46
1:A:105:GLU:OE1	1:A:105:GLU:HA	2.16	0.46
1:C:160:MET:N	1:C:161:PRO:CD	2.79	0.46
1:C:15:SER:O	1:C:19:ILE:HG13	2.17	0.45
1:C:23:ALA:HA	1:C:46:VAL:HG13	1.98	0.45
1:D:96:TYR:CE1	1:D:98:LEU:HB3	2.51	0.45
1:C:98:LEU:HD11	1:C:114:ARG:HB3	1.98	0.45
1:A:47:ASP:C	1:A:47:ASP:OD1	2.59	0.45
1:C:14:ARG:NH1	1:C:18:ILE:HD11	2.31	0.45
1:B:96:TYR:CZ	1:B:98:LEU:HB3	2.51	0.45
1:C:87:ARG:HH12	1:C:205:SER:HB3	1.77	0.45
1:D:47:ASP:OD1	1:D:47:ASP:C	2.59	0.45
1:A:81:GLU:OE1	1:B:81:GLU:OE2	2.35	0.44
1:C:23:ALA:HA	1:C:46:VAL:CG1	2.46	0.44
1:B:73:LEU:N	1:B:73:LEU:CD1	2.81	0.44
1:B:78:MET:HA	1:B:78:MET:HE2	1.99	0.44
1:D:38:ARG:NH2	1:D:207:ARG:HD2	2.33	0.44
1:D:112:ASN:ND2	1:D:113:THR:H	2.14	0.44
1:A:92:ASN:ND2	1:A:92:ASN:H	2.16	0.44
1:A:160:MET:N	1:A:161:PRO:CD	2.80	0.44
1:B:43:CYS:SG	1:B:195:MET:HE2	2.58	0.44
1:C:73:LEU:HD23	1:C:122:TRP:HB3	1.98	0.44
1:A:71:PRO:HB2	1:A:95:ILE:HD12	1.99	0.43
1:C:142:ARG:O	1:C:145:GLU:HB2	2.18	0.43
1:A:123:LEU:HD23	1:A:123:LEU:HA	1.78	0.43
1:B:43:CYS:CB	1:B:195:MET:HE2	2.48	0.43
1:D:77:ARG:O	1:D:81:GLU:HG3	2.19	0.43
1:C:38:ARG:HD2	1:C:202:ALA:HB2	2.01	0.43
1:D:67:LEU:HD23	1:D:67:LEU:HA	1.93	0.43
1:D:172:ASP:CG	1:D:187:ARG:HH21	2.27	0.43
1:B:47:ASP:OD1	1:B:47:ASP:C	2.62	0.42
1:C:47:ASP:CG	1:C:190:ARG:HH12	2.24	0.42
1:D:162:VAL:HG22	1:D:163:GLY:N	2.34	0.42
1:B:22:GLU:OE2	1:D:22:GLU:OE2	2.37	0.42
1:B:160:MET:N	1:B:161:PRO:CD	2.83	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:90:ALA:H	1:C:92:ASN:ND2	2.18	0.42
1:C:123:LEU:HD23	1:C:123:LEU:HA	1.79	0.42
1:D:7:ASP:O	1:D:8:GLY:C	2.62	0.42
1:A:60:LYS:NZ	1:A:64:ASN:HD21	2.18	0.41
1:C:206:GLU:O	1:C:206:GLU:HG3	2.19	0.41
1:C:120:ASP:C	1:C:121:LEU:HD12	2.44	0.41
1:C:135:ASN:O	1:C:137:PRO:HD3	2.21	0.41
1:B:100:ASP:HB3	1:B:103:VAL:HG23	2.03	0.41
1:C:123:LEU:HB2	1:C:124:PRO:HD3	2.02	0.41
1:A:127:GLU:HB2	1:A:152:PRO:HB2	2.02	0.41
1:C:29:PRO:HG2	1:C:32:LEU:HB2	2.02	0.41
1:D:93:ARG:CG	1:D:95:ILE:CD1	2.99	0.41
1:D:123:LEU:HD23	1:D:123:LEU:HA	1.93	0.41
1:C:35:LEU:C	1:C:35:LEU:HD23	2.45	0.41
1:D:84:THR:O	1:D:87:ARG:HB2	2.20	0.40
1:B:123:LEU:N	1:B:124:PRO:CD	2.84	0.40
1:C:162:VAL:HG22	1:C:163:GLY:N	2.36	0.40
1:B:40:ILE:HD11	1:B:49:ALA:HB2	2.02	0.40
1:B:20:ARG:HH12	1:B:207:ARG:HH21	1.63	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	200/212 (94%)	195 (98%)	5 (2%)	0	100	100
1	B	202/212 (95%)	198 (98%)	4 (2%)	0	100	100
1	C	201/212 (95%)	195 (97%)	6 (3%)	0	100	100
1	D	202/212 (95%)	193 (96%)	9 (4%)	0	100	100
All	All	805/848 (95%)	781 (97%)	24 (3%)	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	144/155 (93%)	140 (97%)	4 (3%)	38	26
1	B	145/155 (94%)	136 (94%)	9 (6%)	16	6
1	C	146/155 (94%)	140 (96%)	6 (4%)	27	15
1	D	141/155 (91%)	132 (94%)	9 (6%)	16	6
All	All	576/620 (93%)	548 (95%)	28 (5%)	22	10

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	ARG
1	A	48	VAL
1	A	112	ASN
1	A	121	LEU
1	B	14	ARG
1	B	15	SER
1	B	48	VAL
1	B	73	LEU
1	B	109	LYS
1	B	112	ASN
1	B	135	ASN
1	B	140	LEU
1	B	207	ARG
1	C	56	GLU
1	C	78	MET
1	C	92	ASN
1	C	105	GLU
1	C	112	ASN
1	C	140	LEU
1	D	33	GLU
1	D	38	ARG
1	D	56	GLU

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Mol	Chain	Res	Type
1	D	66	LEU
1	D	109	LYS
1	D	112	ASN
1	D	135	ASN
1	D	140	LEU
1	D	207	ARG

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

Mol	Chain	Res	Type
1	A	64	ASN
1	A	92	ASN
1	A	112	ASN
1	B	92	ASN
1	B	112	ASN
1	B	135	ASN
1	C	27	HIS
1	C	41	HIS
1	C	92	ASN
1	C	112	ASN
1	D	41	HIS
1	D	50	ASN
1	D	92	ASN
1	D	112	ASN
1	D	135	ASN

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	202/212 (95%)	0.79	12 (5%) 28 27	17, 24, 34, 39	0
1	B	204/212 (96%)	0.79	14 (6%) 23 21	16, 23, 35, 42	0
1	C	203/212 (95%)	0.83	10 (4%) 35 33	18, 24, 35, 39	0
1	D	204/212 (96%)	1.13	25 (12%) 8 6	18, 25, 36, 49	0
All	All	813/848 (95%)	0.89	61 (7%) 20 18	16, 24, 35, 49	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	4	TYR	4.2
1	A	108	LYS	4.1
1	D	204	ALA	4.0
1	D	207	ARG	3.6
1	D	12	TYR	3.6
1	A	15	SER	3.5
1	B	207	ARG	3.4
1	D	203	LEU	3.3
1	A	10	ALA	3.3
1	A	107	ALA	3.2
1	D	10	ALA	3.1
1	B	12	TYR	3.1
1	C	91	ASP	3.0
1	D	166	GLY	3.0
1	B	110	ILE	3.0
1	D	5	ILE	3.0
1	A	26	ARG	2.9
1	A	23	ALA	2.8
1	B	10	ALA	2.8
1	A	205	SER	2.7
1	A	87	ARG	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	136	ALA	2.7
1	D	29	PRO	2.7
1	B	4	TYR	2.7
1	D	128	GLY	2.6
1	D	96	TYR	2.5
1	A	7	ASP	2.5
1	D	7	ASP	2.5
1	C	47	ASP	2.5
1	D	17	ALA	2.5
1	D	23	ALA	2.5
1	D	88	LEU	2.4
1	C	167	ALA	2.4
1	D	191	GLY	2.4
1	C	109	LYS	2.4
1	D	84	THR	2.4
1	D	111	GLY	2.4
1	C	135	ASN	2.4
1	B	206	GLU	2.3
1	C	15	SER	2.3
1	C	159	GLY	2.3
1	B	50	ASN	2.3
1	A	21	ALA	2.2
1	B	205	SER	2.2
1	A	5	ILE	2.2
1	D	11	ILE	2.2
1	C	21	ALA	2.2
1	B	103	VAL	2.2
1	B	149	ALA	2.2
1	B	111	GLY	2.2
1	D	99	SER	2.1
1	A	149	ALA	2.1
1	D	122	TRP	2.1
1	D	103	VAL	2.1
1	D	30	ALA	2.1
1	D	206	GLU	2.1
1	D	150	GLY	2.1
1	C	136	ALA	2.1
1	B	120	ASP	2.1
1	B	15	SER	2.0
1	C	206	GLU	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.