



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:35 PM UTC

PDB ID : 6AKE / pdb_00006ake
Title : Crystal structure of mouse claudin-3 in complex with C-terminal fragment of Clostridium perfringens enterotoxin
Authors : Nakamura, S.; Irie, K.; Fujiyoshi, Y.
Deposited on : 2018-08-31
Resolution : 3.60 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

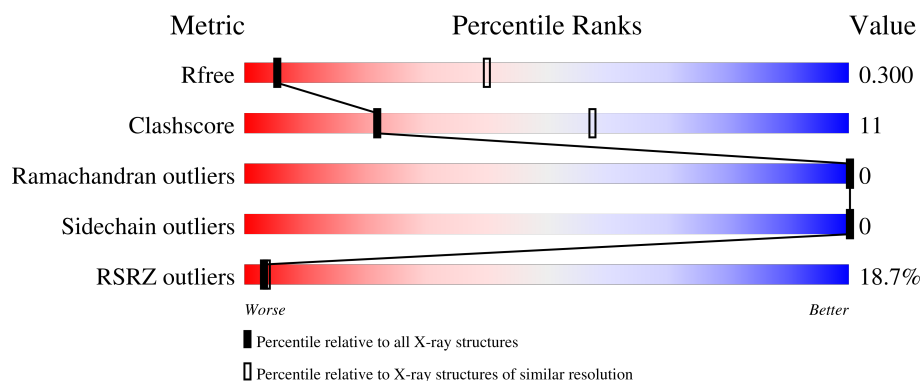
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.60 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	1747 (3.70-3.50)
Clashscore	190562	1827 (3.70-3.50)
Ramachandran outliers	187476	1773 (3.70-3.50)
Sidechain outliers	187428	1772 (3.70-3.50)
RSRZ outliers	180081	1745 (3.70-3.50)

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	190	15% 83% 11% 6%
1	C	190	30% 69% 18% 12%
2	B	119	6% 87% 12% .
2	D	119	13% 76% 19% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4306 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 Claudin-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	179	Total	C	N	O	S	0	0	0
			1278	834	206	226	12			
1	C	168	Total	C	N	O	S	0	0	0
			1211	795	193	211	12			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	GLY	-	see sequence details	UNP Q9Z0G9
A	-5	HIS	-	see sequence details	UNP Q9Z0G9
A	-4	MET	-	see sequence details	UNP Q9Z0G9
A	-3	ALA	-	see sequence details	UNP Q9Z0G9
A	-2	SER	-	see sequence details	UNP Q9Z0G9
A	-1	GLY	-	see sequence details	UNP Q9Z0G9
A	0	SER	-	see sequence details	UNP Q9Z0G9
A	103	ALA	CYS	engineered mutation	UNP Q9Z0G9
A	106	ALA	CYS	engineered mutation	UNP Q9Z0G9
A	181	ALA	CYS	engineered mutation	UNP Q9Z0G9
A	182	ALA	CYS	engineered mutation	UNP Q9Z0G9
C	-6	GLY	-	see sequence details	UNP Q9Z0G9
C	-5	HIS	-	see sequence details	UNP Q9Z0G9
C	-4	MET	-	see sequence details	UNP Q9Z0G9
C	-3	ALA	-	see sequence details	UNP Q9Z0G9
C	-2	SER	-	see sequence details	UNP Q9Z0G9
C	-1	GLY	-	see sequence details	UNP Q9Z0G9
C	0	SER	-	see sequence details	UNP Q9Z0G9
C	103	ALA	CYS	engineered mutation	UNP Q9Z0G9
C	106	ALA	CYS	engineered mutation	UNP Q9Z0G9
C	181	ALA	CYS	engineered mutation	UNP Q9Z0G9
C	182	ALA	CYS	engineered mutation	UNP Q9Z0G9

- Molecule 2 is a protein called Heat-labile enterotoxin B chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	117	Total 918	C 586	N 152	O 179	S 1	0	0	0
2	D	115	Total 899	C 575	N 146	O 177	S 1	0	0	0

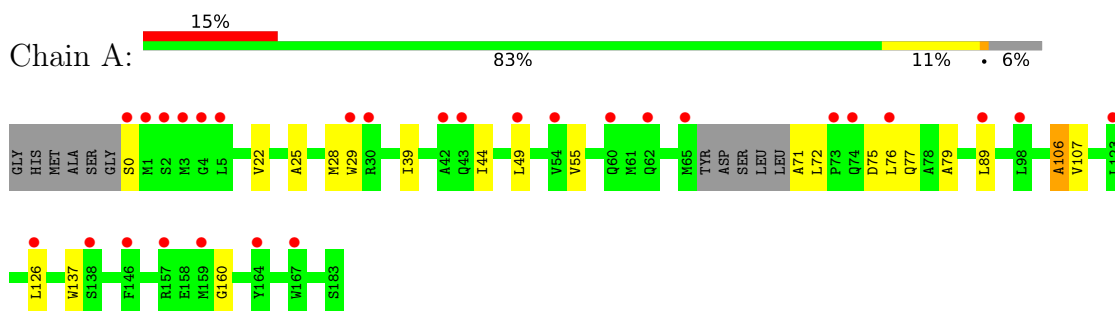
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	201	GLY	-	see sequence details	UNP P01558
B	202	SER	-	see sequence details	UNP P01558
B	313	ALA	SER	engineered mutation	UNP P01558
D	201	GLY	-	see sequence details	UNP P01558
D	202	SER	-	see sequence details	UNP P01558
D	313	ALA	SER	engineered mutation	UNP P01558

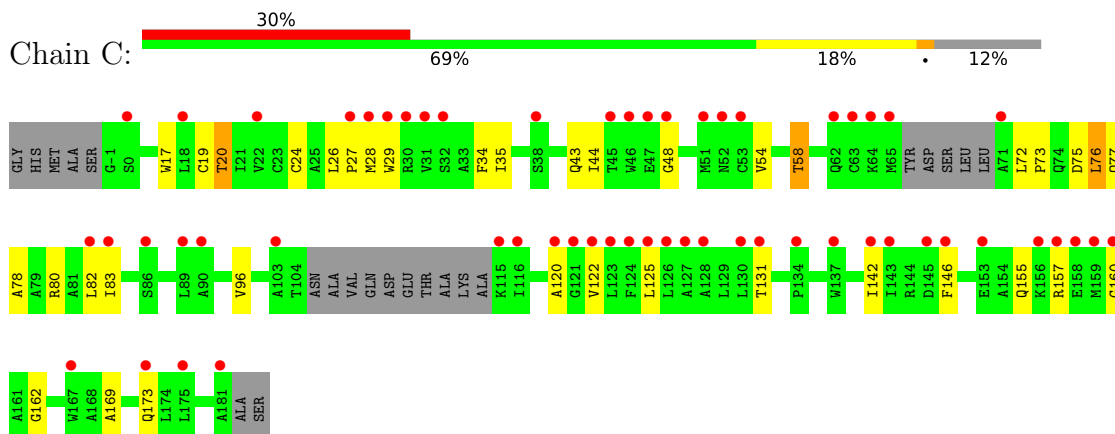
3 Residue-property plots [i](#)

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.

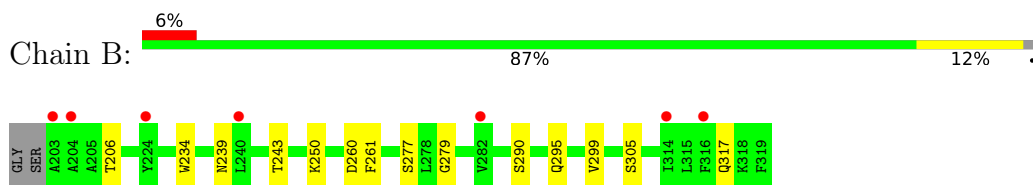
- Molecule 1: Claudin-3



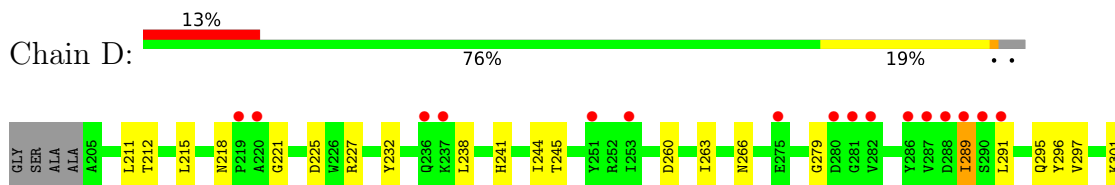
- Molecule 1: Claudin-3



- Molecule 2: Heat-labile enterotoxin B chain



- Molecule 2: Heat-labile enterotoxin B chain





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	92.60Å 66.63Å 111.54Å 90.00° 97.55° 90.00°	Depositor
Resolution (Å)	45.90 – 3.60 45.90 – 3.60	Depositor EDS
% Data completeness (in resolution range)	94.0 (45.90-3.60) 94.0 (45.90-3.60)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.95 (at 3.57Å)	Xtriage
Refinement program	REFMAC 5, PHENIX 1.13_2998	Depositor
R, R_{free}	0.264 , 0.283 (Not available) , 0.300	Depositor DCC
R_{free} test set	756 reflections (4.74%)	wwPDB-VP
Wilson B-factor (Å ²)	147.3	Xtriage
Anisotropy	0.068	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 69.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.84	EDS
Total number of atoms	4306	wwPDB-VP
Average B, all atoms (Å ²)	142.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.35% 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	0.80	0/1299	1.10	4/1778 (0.2%)
1	C	0.79	0/1231	1.07	4/1682 (0.2%)
2	B	0.75	0/939	0.89	0/1281
2	D	0.78	0/920	0.99	1/1257 (0.1%)
All	All	0.78	0/4389	1.03	9/5998 (0.2%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	106	ALA	N-CA-C	8.26	120.20	111.03
1	A	107	VAL	N-CA-C	-7.51	93.71	109.34
1	A	29	TRP	N-CA-C	7.46	119.31	111.03
1	C	58	THR	N-CA-C	5.67	119.06	111.24
2	D	289	ILE	N-CA-C	5.64	116.69	109.30
1	A	77	GLN	N-CA-C	5.49	117.35	111.36
1	C	54	VAL	N-CA-C	5.48	114.69	106.42
1	C	20	THR	N-CA-CB	5.08	117.69	110.16
1	C	76	LEU	N-CA-C	5.01	119.39	113.12

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	1278	0	1316	31	0
1	C	1211	0	1259	37	0
2	B	918	0	874	9	0
2	D	899	0	844	24	0
All	All	4306	0	4293	92	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 11.

All (92) 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:71:ALA:CB	1:A:76:LEU:HD23	1.15	1.63
1:A:71:ALA:CB	1:A:76:LEU:CD2	1.77	1.52
1:A:71:ALA:HB1	1:A:76:LEU:CD2	1.44	1.33
2:D:212:THR:OG1	2:D:245:THR:HG23	1.32	1.28
1:A:71:ALA:HB3	1:A:76:LEU:CD2	1.54	1.19
2:D:212:THR:OG1	2:D:245:THR:CG2	1.92	1.16
1:A:71:ALA:CB	1:A:76:LEU:HD21	1.75	1.13
2:D:212:THR:HG1	2:D:245:THR:HG23	0.82	0.98
1:A:71:ALA:HB3	1:A:76:LEU:HD23	1.02	0.96
2:D:212:THR:OG1	2:D:245:THR:CB	2.15	0.94
1:A:71:ALA:HB1	1:A:76:LEU:HD23	0.98	0.86
1:A:71:ALA:HB3	1:A:76:LEU:CG	2.09	0.81
1:A:106:ALA:HA	1:C:35:ILE:CG2	2.11	0.81
1:C:78:ALA:O	1:C:82:LEU:CB	2.31	0.78
1:C:72:LEU:HD23	1:C:75:ASP:CG	2.10	0.77
2:D:260:ASP:OD1	2:D:279:GLY:N	2.20	0.74
1:C:78:ALA:O	1:C:82:LEU:HB3	1.87	0.73
1:C:76:LEU:CD1	1:C:80:ARG:HG3	2.17	0.73
1:C:76:LEU:HD11	1:C:80:ARG:HG3	1.71	0.72
2:D:212:THR:OG1	2:D:245:THR:OG1	2.07	0.72
1:A:106:ALA:HA	1:C:35:ILE:HG22	1.71	0.72
1:A:28:MET:O	1:A:160:GLY:HA3	1.89	0.71
1:A:71:ALA:HB2	1:A:76:LEU:HD21	1.74	0.69
2:B:206:THR:HG22	2:B:239:ASN:HB2	1.74	0.67
1:C:43:GLN:HE22	2:D:218:ASN:HD22	1.41	0.67
1:C:73:PRO:HA	1:C:77:GLN:HG2	1.77	0.66
1:C:72:LEU:HB3	1:C:75:ASP:HB2	1.77	0.65
1:C:155:GLN:HE21	2:D:225:ASP:HA	1.60	0.65
1:A:49:LEU:HD22	1:A:79:ALA:HB3	1.79	0.63
2:D:212:THR:CB	2:D:245:THR:HG1	2.10	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:ALA:HA	1:C:35:ILE:HG21	1.82	0.62
1:C:76:LEU:HD12	1:C:76:LEU:O	1.99	0.61
1:A:49:LEU:HD22	1:A:79:ALA:CB	2.31	0.59
1:C:78:ALA:O	1:C:82:LEU:HB2	2.00	0.59
1:A:72:LEU:H	1:A:76:LEU:HD23	1.67	0.58
1:C:35:ILE:HD11	1:C:44:ILE:HD11	1.87	0.57
1:A:39:ILE:HD13	2:B:317:GLN:HE22	1.70	0.57
1:C:35:ILE:HD11	1:C:44:ILE:CD1	2.34	0.57
1:A:106:ALA:CA	1:C:35:ILE:HG21	2.35	0.56
1:C:77:GLN:HA	1:C:77:GLN:OE1	2.07	0.55
1:C:24:CYS:HB2	1:C:83:ILE:HG21	1.90	0.53
1:A:71:ALA:HB3	1:A:76:LEU:HG	1.90	0.53
2:D:212:THR:HG1	2:D:245:THR:CG2	1.73	0.51
1:A:49:LEU:CD2	1:A:79:ALA:CB	2.88	0.51
1:C:76:LEU:HD12	1:C:80:ARG:HG3	1.93	0.50
1:A:75:ASP:CG	1:A:137:TRP:CZ3	2.90	0.49
1:C:27:PRO:HA	1:C:48:GLY:HA3	1.95	0.49
2:D:225:ASP:OD1	2:D:227:ARG:NH1	2.46	0.49
1:C:122:VAL:HA	1:C:125:LEU:HB2	1.95	0.49
2:D:291:LEU:HD22	2:D:296:TYR:CE2	2.48	0.49
2:B:234:TRP:CG	2:B:305:SER:HG	2.30	0.48
1:A:49:LEU:CD2	1:A:79:ALA:HB1	2.44	0.48
2:D:211:LEU:HD11	2:D:316:PHE:CG	2.48	0.48
1:C:34:PHE:C	1:C:35:ILE:HG13	2.38	0.48
1:C:72:LEU:HD12	1:C:73:PRO:HD2	1.95	0.48
2:D:212:THR:OG1	2:D:245:THR:N	2.46	0.48
1:C:142:ILE:HG23	1:C:157:ARG:O	2.14	0.48
2:D:263:ILE:HD13	2:D:289:ILE:HD12	1.95	0.47
1:A:22:VAL:O	1:A:25:ALA:N	2.47	0.47
1:C:58:THR:HG22	1:C:58:THR:O	2.14	0.47
1:A:106:ALA:CA	1:C:35:ILE:CG2	2.87	0.46
1:A:72:LEU:N	1:A:76:LEU:HD23	2.31	0.46
2:B:250:LYS:HA	2:B:290:SER:HA	1.98	0.45
2:D:212:THR:HG23	2:D:244:ILE:HG23	1.98	0.45
1:C:19:CYS:HB3	1:C:169:ALA:HB2	1.99	0.45
2:D:241:HIS:CD2	2:D:297:VAL:HG22	2.52	0.44
1:A:75:ASP:OD1	1:A:137:TRP:CH2	2.70	0.44
1:A:75:ASP:OD2	1:A:137:TRP:CZ3	2.71	0.44
2:D:266:ASN:HB3	2:D:295:GLN:HB3	2.01	0.43
1:A:44:ILE:HB	1:A:55:VAL:CG1	2.48	0.43
1:C:20:THR:HB	1:C:131:THR:HG21	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:291:LEU:HD22	2:D:296:TYR:CD2	2.54	0.43
2:D:232:TYR:CE2	2:D:238:LEU:HD13	2.54	0.43
2:B:239:ASN:OD1	2:B:299:VAL:HG22	2.20	0.42
2:B:260:ASP:OD1	2:B:279:GLY:N	2.52	0.42
2:D:301:LYS:HA	2:D:312:TYR:OH	2.19	0.42
1:C:96:VAL:HG12	1:C:120:ALA:HB2	2.01	0.42
2:D:212:THR:CG2	2:D:244:ILE:HG23	2.50	0.42
2:D:295:GLN:C	2:D:296:TYR:CD1	2.98	0.42
2:B:243:THR:HG23	2:B:295:GLN:HE22	1.85	0.42
2:D:215:LEU:O	2:D:221:GLY:HA2	2.20	0.42
1:A:89:LEU:HD21	1:A:126:LEU:HB2	2.01	0.41
1:C:20:THR:HG23	1:C:169:ALA:HB1	2.03	0.41
1:C:29:TRP:O	1:C:160:GLY:N	2.53	0.41
1:C:76:LEU:HD11	1:C:80:ARG:CG	2.47	0.41
2:B:206:THR:HG22	2:B:239:ASN:HD22	1.85	0.41
1:A:0:SER:HB3	1:C:146:PHE:CE1	2.56	0.41
2:B:261:PHE:CE1	2:B:277:SER:HB3	2.55	0.41
1:A:72:LEU:HB3	1:A:76:LEU:HB3	2.02	0.41
1:C:26:LEU:HD22	1:C:28:MET:HE2	2.04	0.40
1:C:17:TRP:HA	1:C:173:GLN:NE2	2.36	0.40
1:C:28:MET:O	1:C:162:GLY:N	2.55	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	175/190 (92%)	169 (97%)	6 (3%)	0	100	100
1	C	162/190 (85%)	161 (99%)	1 (1%)	0	100	100
2	B	115/119 (97%)	115 (100%)	0	0	100	100
2	D	113/119 (95%)	111 (98%)	2 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	565/618 (91%)	556 (98%)	9 (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	127/144 (88%)	127 (100%)	0	100	100
1	C	122/144 (85%)	122 (100%)	0	100	100
2	B	98/102 (96%)	98 (100%)	0	100	100
2	D	96/102 (94%)	96 (100%)	0	100	100
All	All	443/492 (90%)	443 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	52	ASN
1	A	102	GLN
1	A	105	ASN
2	B	218	ASN
2	B	239	ASN
2	B	295	GLN
1	C	74	GLN
1	C	173	GLN
2	D	218	ASN
2	D	241	HIS
2	D	276	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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	179/190 (94%)	0.82	28 (15%) 5 5	84, 135, 202, 228	0
1	C	168/190 (88%)	1.59	57 (33%) 1 1	88, 134, 200, 264	0
2	B	117/119 (98%)	0.30	7 (5%) 27 17	114, 144, 174, 204	0
2	D	115/119 (96%)	0.59	16 (13%) 6 6	122, 147, 177, 212	0
All	All	579/618 (93%)	0.89	108 (18%) 3 4	84, 141, 196, 264	0

All (108) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	47	GLU	13.0
1	C	31	VAL	10.4
2	D	282	VAL	9.5
1	C	32	SER	8.2
1	C	157	ARG	7.3
1	C	146	PHE	6.9
1	C	30	ARG	6.8
1	C	159	MET	6.1
2	D	288	ASP	6.0
2	D	289	ILE	6.0
1	C	158	GLU	6.0
1	A	2	SER	5.8
2	D	290	SER	5.5
1	C	63	CYS	5.3
1	A	4	GLY	5.2
1	C	45	THR	5.0
1	C	153	GLU	5.0
2	B	316	PHE	4.9
1	A	74	GLN	4.9
1	A	164	TYR	4.8
1	A	0	SER	4.8

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Mol	Chain	Res	Type	RSRZ
1	C	46	TRP	4.7
1	A	43	GLN	4.6
1	A	54	VAL	4.5
1	C	18	LEU	4.5
1	C	116	ILE	4.5
1	A	30	ARG	4.5
1	A	60	GLN	4.4
1	C	131	THR	4.4
1	C	53	CYS	4.1
2	B	203	ALA	4.1
1	C	115	LYS	4.0
1	A	3	MET	3.9
1	C	160	GLY	3.9
1	C	64	LYS	3.9
1	C	103	ALA	3.7
1	C	156	LYS	3.6
2	D	287	VAL	3.6
2	D	251	TYR	3.5
1	C	65	MET	3.4
1	A	62	GLN	3.4
1	C	83	ILE	3.3
2	B	282	VAL	3.3
2	D	280	ASP	3.3
1	A	42	ALA	3.2
1	C	51	MET	3.2
1	C	86	SER	3.2
1	A	5	LEU	3.1
1	C	143	ILE	3.0
1	C	123	LEU	3.0
1	C	124	PHE	3.0
1	C	127	ALA	3.0
1	C	52	ASN	3.0
1	C	167	TRP	3.0
1	C	71	ALA	3.0
1	A	1	MET	3.0
1	C	125	LEU	2.9
2	D	219	PRO	2.9
2	B	240	LEU	2.9
2	D	286	TYR	2.9
1	A	123	LEU	2.8
1	C	28	MET	2.8
1	C	142	ILE	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	167	TRP	2.8
1	C	134	PRO	2.8
1	C	62	GLN	2.8
2	D	220	ALA	2.7
1	A	76	LEU	2.7
1	C	122	VAL	2.6
2	B	204	ALA	2.6
1	C	48	GLY	2.6
1	C	173	GLN	2.6
1	A	138	SER	2.5
1	C	22	VAL	2.5
1	C	29	TRP	2.5
1	C	38	SER	2.5
2	D	291	LEU	2.5
1	A	73	PRO	2.5
1	C	145	ASP	2.5
1	A	157	ARG	2.5
1	C	175	LEU	2.5
2	D	236	GLN	2.4
1	C	128	ALA	2.4
1	C	126	LEU	2.4
2	D	237	LYS	2.4
1	A	126	LEU	2.3
1	A	146	PHE	2.3
2	B	224	TYR	2.3
1	A	29	TRP	2.3
1	C	137	TRP	2.3
1	A	49	LEU	2.3
2	B	314	ILE	2.3
1	A	65	MET	2.3
1	C	27	PRO	2.2
1	A	98	LEU	2.2
2	D	253	ILE	2.2
1	C	181	ALA	2.2
1	C	121	GLY	2.2
1	C	82	LEU	2.2
1	C	89	LEU	2.2
1	C	0	SER	2.1
2	D	275	GLU	2.1
1	A	159	MET	2.1
1	C	120	ALA	2.1
1	C	90	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
2	D	281	GLY	2.1
1	A	89	LEU	2.1
1	C	130	LEU	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.