



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:21 PM UTC

PDB ID : 3FHV / pdb_00003fhv
Title : Structural basis of Salmonella typhi type IVb PilS and cystic fibrosis trans-membrane conductance regulator (CFTR) interaction
Authors : Balakrishna, A.M.; Swaminathan, K.
Deposited on : 2008-12-10
Resolution : 1.90 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

<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

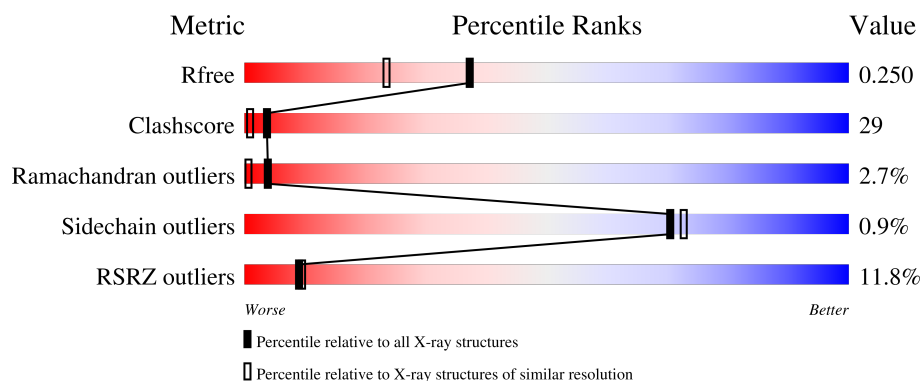
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	180053	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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 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	156	10% 72% 23% ..
1	B	156	10% 74% 21% ..
2	C	10	60% 40% 40% 20%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	DTY	C	109	-	-	X	-
2	DPR	C	111	-	-	X	-
2	DGL	C	116	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2474 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 Prepilin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	150	Total	C	N	O	S	0	0	0
			1054	639	183	225	7			
1	B	150	Total	C	N	O	S	0	0	0
			1054	639	183	225	7			

- Molecule 2 is a protein called cfr peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	10	Total	C	N	O	0	0	0
			88	51	15	22			

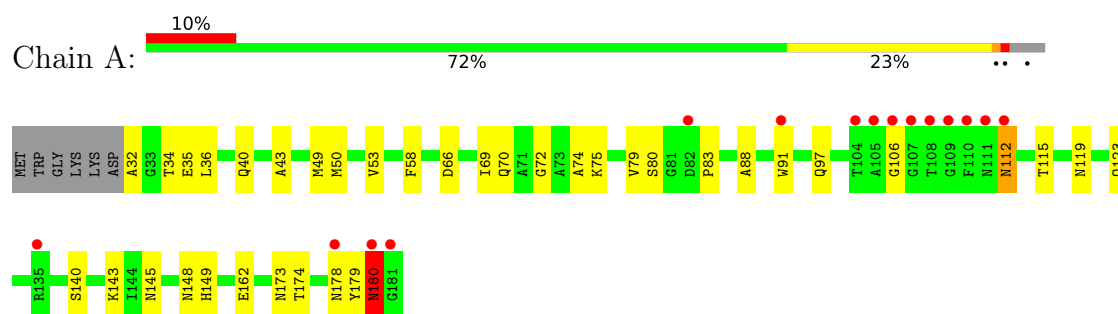
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	142	Total	O	0	0
			142	142		
3	B	130	Total	O	0	0
			130	130		
3	C	6	Total	O	0	0
			6	6		

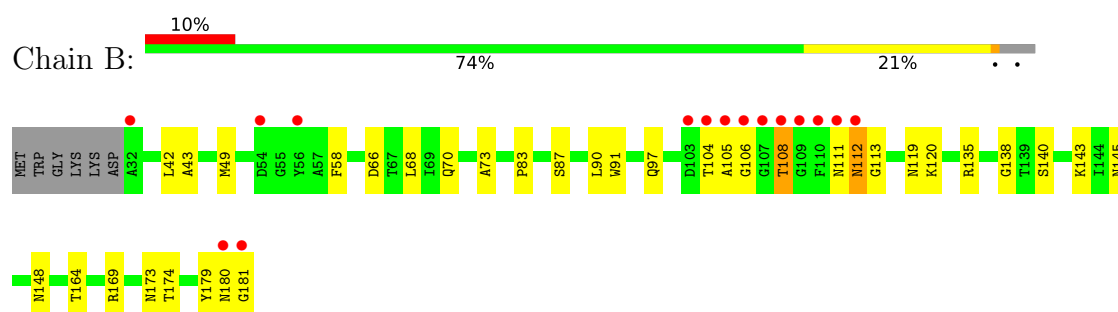
3 Residue-property plots [i](#)

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.

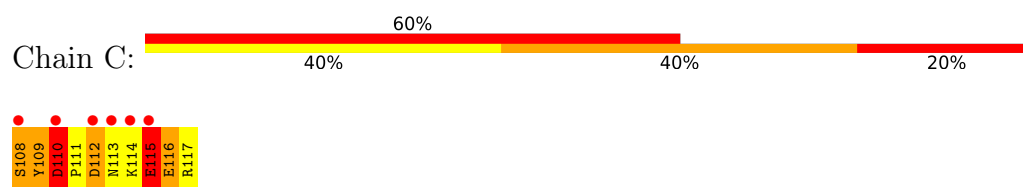
• Molecule 1: Prepilin



• Molecule 1: Prepilin



• Molecule 2: cftr peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	77.80Å 114.37Å 31.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 1.90 8.00 – 1.90	Depositor EDS
% Data completeness (in resolution range)	88.6 (8.00-1.90) 96.9 (8.00-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.88 (at 1.70Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.207 , 0.248 0.217 , 0.250	Depositor DCC
R_{free} test set	2252 reflections (7.25%)	wwPDB-VP
Wilson B-factor (Å ²)	12.8	Xtriage
Anisotropy	0.220	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.52 , 67.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	2474	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.59% 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

Bond lengths and bond angles in the following residue types are not validated in this section: DPR, DTY, DAR, DGL

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	0.41	0/1068	0.84	2/1449 (0.1%)
1	B	0.42	0/1068	0.86	1/1449 (0.1%)
2	C	0.09	0/45	0.35	0/55
All	All	0.41	0/2181	0.84	3/2953 (0.1%)

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
2	C	0	5

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	140	SER	N-CA-C	-6.08	105.18	112.59
1	A	119	ASN	N-CA-C	5.72	119.29	110.42
1	B	140	SER	N-CA-C	-5.65	104.80	112.26

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	C	108	SER	Peptide
2	C	109	DTY	Peptide
2	C	110	ASP	Peptide
2	C	115	GLU	Peptide

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Mol	Chain	Res	Type	Group
2	C	116	DGL	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	1054	0	1009	52	0
1	B	1054	0	1009	54	6
2	C	88	0	70	67	6
3	A	142	0	0	1	0
3	B	130	0	0	2	0
3	C	6	0	0	1	0
All	All	2474	0	2088	124	6

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 29.

All (124) 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:120:LYS:NZ	2:C:115:GLU:HG3	1.21	1.44
1:B:120:LYS:NZ	2:C:115:GLU:CG	1.98	1.24
1:A:75:LYS:CD	2:C:112:ASP:OD2	1.87	1.21
1:A:75:LYS:HD3	2:C:112:ASP:OD2	1.05	1.20
1:A:75:LYS:HD3	2:C:112:ASP:CG	1.76	1.09
2:C:109:DTY:CG	2:C:110:ASP:H	1.69	1.04
1:B:120:LYS:HZ3	2:C:115:GLU:CG	1.60	1.02
1:B:120:LYS:HZ1	2:C:115:GLU:CG	1.77	0.94
2:C:109:DTY:CD2	2:C:110:ASP:H	1.81	0.93
1:B:164:THR:H	1:B:173:ASN:HD21	1.03	0.93
1:B:97:GLN:HB3	2:C:116:DGL:O	1.73	0.88
1:A:75:LYS:HA	2:C:108:SER:OG	1.73	0.87
1:B:120:LYS:HZ1	2:C:115:GLU:HG2	1.41	0.85
1:B:120:LYS:HD2	2:C:113:ASN:HD22	1.39	0.85
1:B:120:LYS:CD	2:C:113:ASN:HD22	1.92	0.83
1:B:120:LYS:CE	2:C:115:GLU:HG3	2.06	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:120:LYS:HD2	2:C:113:ASN:ND2	1.93	0.82
1:B:164:THR:N	1:B:173:ASN:HD21	1.78	0.82
1:B:143:LYS:HE2	1:B:148:ASN:ND2	1.96	0.81
1:A:69:ILE:O	2:C:108:SER:N	2.14	0.80
1:A:72:GLY:HA2	2:C:108:SER:HA	1.62	0.79
1:A:115:THR:HG22	1:A:178:ASN:ND2	1.99	0.78
1:B:145:ASN:HD21	1:B:173:ASN:HD22	1.32	0.77
1:A:75:LYS:CA	2:C:108:SER:OG	2.31	0.76
1:A:50:MET:HE1	1:A:58:PHE:HE2	1.50	0.76
1:A:75:LYS:HG2	1:B:91:TRP:CZ2	2.23	0.74
1:B:119:ASN:ND2	2:C:117:DAR:CG	2.52	0.73
2:C:111:DPR:O	2:C:112:ASP:CB	2.36	0.73
1:A:75:LYS:CG	2:C:112:ASP:OD2	2.37	0.72
2:C:109:DTY:CG	2:C:110:ASP:N	2.35	0.71
2:C:109:DTY:CD2	2:C:110:ASP:N	2.51	0.71
1:B:119:ASN:OD1	1:B:174:THR:HG23	1.90	0.71
1:A:106:GLY:H	1:A:112:ASN:ND2	1.88	0.70
1:B:164:THR:H	1:B:173:ASN:ND2	1.85	0.70
1:B:112:ASN:CG	1:B:113:GLY:H	2.00	0.69
1:B:119:ASN:HD21	2:C:117:DAR:HG2	1.58	0.68
1:B:120:LYS:CG	2:C:113:ASN:HD22	2.07	0.67
1:B:66:ASP:HB2	1:B:87:SER:HB3	1.76	0.67
1:B:120:LYS:CE	2:C:115:GLU:CG	2.70	0.67
1:B:66:ASP:O	1:B:70:GLN:HG2	1.94	0.66
1:A:70:GLN:O	2:C:110:ASP:O	2.14	0.65
1:B:119:ASN:ND2	2:C:117:DAR:HG3	2.11	0.65
2:C:114:LYS:O	2:C:115:GLU:HG2	1.97	0.64
1:A:50:MET:HE2	1:A:50:MET:HA	1.77	0.64
1:B:68:LEU:HG	1:B:73:ALA:HB3	1.79	0.63
1:A:106:GLY:H	1:A:112:ASN:HD22	1.46	0.63
1:A:66:ASP:O	1:A:70:GLN:HG2	1.99	0.63
1:A:149:HIS:HE1	1:A:162:GLU:OE1	1.82	0.62
1:A:72:GLY:HA2	2:C:108:SER:CA	2.30	0.61
1:A:50:MET:O	1:A:53:VAL:HG22	2.01	0.61
1:A:75:LYS:HB2	2:C:112:ASP:OD2	2.01	0.61
1:B:119:ASN:ND2	2:C:117:DAR:HG2	2.14	0.60
1:B:169:ARG:O	1:B:169:ARG:HG2	2.03	0.59
1:A:69:ILE:HG22	2:C:109:DTY:HA	1.85	0.59
1:B:120:LYS:HZ3	2:C:115:GLU:HG3	0.66	0.59
1:B:97:GLN:CB	2:C:116:DGL:O	2.48	0.58
1:A:143:LYS:HE2	1:A:148:ASN:ND2	2.18	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:91:TRP:CZ3	1:A:97:GLN:HB2	2.41	0.56
1:A:32:ALA:HA	1:A:35:GLU:OE1	2.04	0.56
1:A:49:MET:HE3	1:A:50:MET:HE3	1.85	0.56
1:A:75:LYS:CB	2:C:112:ASP:OD2	2.53	0.56
1:A:49:MET:CE	1:A:50:MET:HE3	2.36	0.56
1:A:80:SER:OG	1:A:91:TRP:NE1	2.36	0.55
1:A:115:THR:CG2	1:A:178:ASN:ND2	2.67	0.55
1:B:143:LYS:HE2	1:B:148:ASN:CG	2.31	0.55
2:C:111:DPR:O	2:C:112:ASP:HB2	2.07	0.55
2:C:111:DPR:O	2:C:112:ASP:HB3	2.05	0.54
2:C:109:DTY:O	2:C:110:ASP:C	2.51	0.54
1:A:74:ALA:C	2:C:108:SER:HB3	2.34	0.53
1:B:119:ASN:ND2	3:B:224:HOH:O	2.41	0.53
2:C:115:GLU:H	2:C:116:DGL:HG3	1.74	0.53
1:B:120:LYS:CE	2:C:115:GLU:HB2	2.40	0.52
1:A:145:ASN:HD21	1:A:173:ASN:HA	1.75	0.52
2:C:114:LYS:O	2:C:115:GLU:CG	2.57	0.52
1:B:111:ASN:HB3	1:B:181:GLY:HA3	1.92	0.52
1:A:69:ILE:CG2	2:C:109:DTY:HA	2.40	0.52
1:A:43:ALA:CB	1:A:179:TYR:CZ	2.93	0.51
1:A:79:VAL:HG11	2:C:109:DTY:HB3	1.94	0.50
1:A:43:ALA:HB1	1:A:179:TYR:CZ	2.46	0.50
1:A:179:TYR:O	1:A:180:ASN:O	2.31	0.49
1:A:75:LYS:N	2:C:108:SER:CB	2.76	0.49
1:A:69:ILE:HG23	2:C:108:SER:N	2.28	0.48
2:C:115:GLU:N	2:C:116:DGL:HG3	2.28	0.48
1:B:120:LYS:CE	2:C:115:GLU:CB	2.92	0.48
1:A:145:ASN:ND2	1:A:174:THR:H	2.12	0.48
2:C:109:DTY:O	2:C:111:DPR:O	2.32	0.48
2:C:114:LYS:C	2:C:115:GLU:HG2	2.38	0.48
1:B:42:LEU:HD11	1:B:90:LEU:HD21	1.96	0.47
2:C:108:SER:HA	2:C:109:DTY:O	2.13	0.47
1:B:104:THR:HG23	1:B:105:ALA:N	2.30	0.47
1:B:112:ASN:ND2	1:B:113:GLY:H	2.11	0.47
1:B:119:ASN:HD22	2:C:117:DAR:HG3	1.77	0.47
1:B:91:TRP:HZ3	2:C:116:DGL:OE1	1.98	0.47
1:A:75:LYS:CD	2:C:112:ASP:CG	2.68	0.47
1:A:149:HIS:CE1	1:A:162:GLU:OE1	2.65	0.47
2:C:110:ASP:OD1	3:C:231:HOH:O	2.21	0.46
1:B:135:ARG:NH2	3:B:293:HOH:O	2.49	0.46
1:B:120:LYS:HE2	2:C:115:GLU:HB2	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:MET:HE3	1:A:50:MET:CE	2.45	0.46
1:A:50:MET:HE1	1:A:58:PHE:CE2	2.40	0.45
1:A:83:PRO:HB3	1:A:88:ALA:HB2	1.98	0.45
1:B:145:ASN:HD21	1:B:173:ASN:ND2	2.09	0.45
2:C:113:ASN:N	2:C:116:DGL:OE2	2.49	0.45
1:A:106:GLY:N	1:A:112:ASN:ND2	2.61	0.45
1:B:120:LYS:HE2	2:C:115:GLU:CB	2.47	0.45
1:A:72:GLY:C	2:C:108:SER:HB2	2.42	0.44
2:C:114:LYS:O	2:C:115:GLU:OE2	2.35	0.44
1:A:91:TRP:CE3	1:A:97:GLN:HA	2.53	0.44
2:C:109:DTY:O	2:C:110:ASP:O	2.35	0.44
1:A:34:THR:HG21	1:B:83:PRO:HA	2.00	0.43
1:B:104:THR:OG1	1:B:105:ALA:N	2.51	0.43
1:B:106:GLY:C	1:B:108:THR:H	2.27	0.43
1:A:180:ASN:HD22	1:A:180:ASN:HA	1.55	0.43
1:B:97:GLN:OE1	2:C:116:DGL:C	2.67	0.42
1:B:70:GLN:NE2	1:B:70:GLN:HA	2.34	0.42
1:A:36:LEU:HG	1:A:40:GLN:NE2	2.34	0.42
2:C:108:SER:CA	2:C:109:DTY:O	2.67	0.42
1:A:32:ALA:N	3:A:262:HOH:O	2.51	0.42
1:B:112:ASN:CG	1:B:113:GLY:N	2.68	0.42
1:B:49:MET:HE3	1:B:58:PHE:CE2	2.56	0.41
1:B:119:ASN:OD1	1:B:174:THR:CG2	2.63	0.41
1:B:120:LYS:HE3	2:C:113:ASN:CB	2.51	0.41
1:B:120:LYS:HE3	2:C:113:ASN:HB2	2.02	0.41
1:A:106:GLY:N	1:A:112:ASN:HD22	2.14	0.40

All (6) 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)
1:B:138:GLY:CA	2:C:115:GLU:OE2[1_556]	1.61	0.59
1:B:138:GLY:C	2:C:115:GLU:OE2[1_556]	2.02	0.18
1:B:179:TYR:CZ	2:C:114:LYS:NZ[1_556]	2.02	0.18
1:B:179:TYR:CE1	2:C:114:LYS:NZ[1_556]	2.03	0.17
1:B:43:ALA:CB	2:C:114:LYS:NZ[1_556]	2.12	0.08
1:B:179:TYR:CE1	2:C:114:LYS:CE[1_556]	2.12	0.08

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	148/156 (95%)	141 (95%)	5 (3%)	2 (1%)	9	3
1	B	148/156 (95%)	143 (97%)	2 (1%)	3 (2%)	6	1
2	C	5/10 (50%)	1 (20%)	1 (20%)	3 (60%)	0	0
All	All	301/322 (94%)	285 (95%)	8 (3%)	8 (3%)	4	1

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	180	ASN
1	B	112	ASN
2	C	110	ASP
2	C	112	ASP
1	B	180	ASN
2	C	115	GLU
1	A	112	ASN
1	B	108	THR

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	111/116 (96%)	109 (98%)	2 (2%)	51	50
1	B	111/116 (96%)	111 (100%)	0	100	100
2	C	6/6 (100%)	6 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	228/238 (96%)	226 (99%)	2 (1%)	70	73

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	123	GLN
1	A	180	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	112	ASN
1	A	123	GLN
1	A	145	ASN
1	A	149	HIS
1	A	178	ASN
1	A	180	ASN
1	B	38	ASN
1	B	70	GLN
1	B	111	ASN
1	B	112	ASN
1	B	119	ASN
1	B	173	ASN
1	B	180	ASN
2	C	113	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 non-standard protein/DNA/RNA residues are modelled in this entry.

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.

5.5 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	150/156 (96%)	0.29	15 (10%)	12 13	6, 12, 56, 90	0
1	B	150/156 (96%)	0.19	15 (10%)	12 13	7, 13, 54, 91	0
2	C	6/10 (60%)	5.31	6 (100%)	0 0	45, 56, 62, 63	0
All	All	306/322 (95%)	0.34	36 (11%)	9 9	6, 12, 58, 91	0

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	104	THR	7.9
1	A	107	GLY	7.2
1	A	105	ALA	6.8
1	A	111	ASN	6.5
2	C	112	ASP	6.4
1	A	180	ASN	6.3
1	B	108	THR	5.9
2	C	108	SER	5.9
2	C	114	LYS	5.8
1	A	181	GLY	5.8
1	A	112	ASN	5.4
1	B	109	GLY	5.0
1	A	106	GLY	4.9
2	C	113	ASN	4.9
1	B	111	ASN	4.9
1	B	110	PHE	4.6
1	A	109	GLY	4.6
2	C	110	ASP	4.6
1	A	108	THR	4.6
1	B	107	GLY	4.6
1	B	112	ASN	4.5
1	A	91	TRP	4.5
2	C	115	GLU	4.2

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Mol	Chain	Res	Type	RSRZ
1	B	180	ASN	4.2
1	A	110	PHE	4.0
1	B	104	THR	3.9
1	B	106	GLY	3.9
1	B	105	ALA	3.9
1	A	82	ASP	3.3
1	B	181	GLY	3.3
1	B	103	ASP	3.1
1	A	178	ASN	2.7
1	B	32	ALA	2.6
1	B	56	TYR	2.3
1	B	54	ASP	2.2
1	A	135	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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. The B-factors column lists the minimum, median, 95th 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	B-factors(Å ²)	Q<0.9
2	DGL	C	116	9/10	0.26	0.36	54,57,59,60	0
2	DPR	C	111	7/8	0.39	0.28	49,51,52,52	0
2	DAR	C	117	12/12	0.42	0.23	59,62,63,63	0
2	DTY	C	109	12/13	0.56	0.25	56,60,61,63	0

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.