



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 26, 2026 – 10:35 AM EDT

PDB ID : 7ZOH / pdb_00007zoh
Title : Carbohydrate binding domain CBM92-B from a multi-catalytic glucanase-chitinase from Chitinophaga pinensis DSM 2588
Authors : Mazurkewich, S.; McKee, L.S.; Lu, Z.; Branden, G.; Larsbrink, J.
Deposited on : 2022-04-25
Resolution : 1.56 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
Mogul : 2022.3.0, CSD as543be (2022)
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

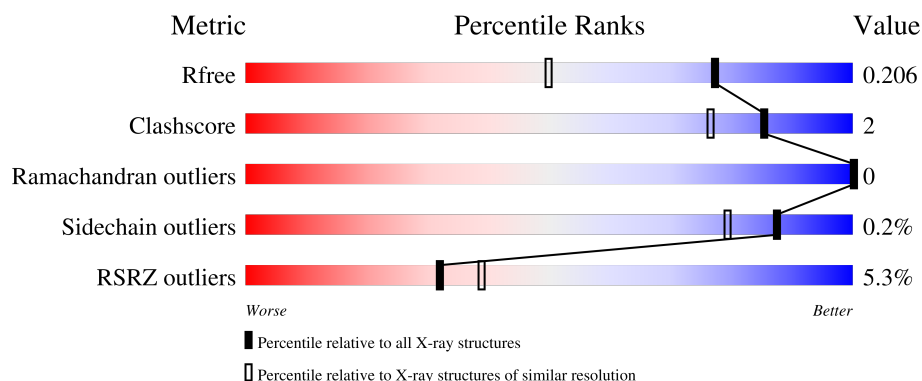
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.56 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	2145 (1.56-1.56)
Clashscore	190562	2189 (1.56-1.56)
Ramachandran outliers	187476	2153 (1.56-1.56)
Sidechain outliers	187428	2150 (1.56-1.56)
RSRZ outliers	180081	2146 (1.56-1.56)

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	147	3% 79% 6% 15%
1	B	147	5% 78% 7% 15%
1	C	147	5% 86% • 12%
1	D	147	5% 79% 9% 12%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4296 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 Glycoside hydrolase family 18.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	125	Total	C	N	O	S	0	0	0
			964	596	171	191	6			
1	B	125	Total	C	N	O	S	0	2	0
			982	606	177	193	6			
1	C	130	Total	C	N	O	S	0	2	0
			1015	627	180	202	6			
1	D	129	Total	C	N	O	S	0	3	0
			1022	629	184	203	6			

There are 92 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	874	MET	-	initiating methionine	UNP A0A979GQH9
A	875	ALA	-	expression tag	UNP A0A979GQH9
A	876	SER	-	expression tag	UNP A0A979GQH9
A	877	MET	-	expression tag	UNP A0A979GQH9
A	878	THR	-	expression tag	UNP A0A979GQH9
A	879	GLY	-	expression tag	UNP A0A979GQH9
A	880	GLY	-	expression tag	UNP A0A979GQH9
A	881	GLN	-	expression tag	UNP A0A979GQH9
A	882	GLN	-	expression tag	UNP A0A979GQH9
A	883	MET	-	expression tag	UNP A0A979GQH9
A	884	GLY	-	expression tag	UNP A0A979GQH9
A	885	ARG	-	expression tag	UNP A0A979GQH9
A	886	GLY	-	expression tag	UNP A0A979GQH9
A	887	SER	-	expression tag	UNP A0A979GQH9
A	959	SNN	ASP	conflict	UNP A0A979GQH9
A	1013	LEU	-	expression tag	UNP A0A979GQH9
A	1014	GLU	-	expression tag	UNP A0A979GQH9
A	1015	HIS	-	expression tag	UNP A0A979GQH9
A	1016	HIS	-	expression tag	UNP A0A979GQH9
A	1017	HIS	-	expression tag	UNP A0A979GQH9
A	1018	HIS	-	expression tag	UNP A0A979GQH9

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1019	HIS	-	expression tag	UNP A0A979GQH9
A	1020	HIS	-	expression tag	UNP A0A979GQH9
B	874	MET	-	initiating methionine	UNP A0A979GQH9
B	875	ALA	-	expression tag	UNP A0A979GQH9
B	876	SER	-	expression tag	UNP A0A979GQH9
B	877	MET	-	expression tag	UNP A0A979GQH9
B	878	THR	-	expression tag	UNP A0A979GQH9
B	879	GLY	-	expression tag	UNP A0A979GQH9
B	880	GLY	-	expression tag	UNP A0A979GQH9
B	881	GLN	-	expression tag	UNP A0A979GQH9
B	882	GLN	-	expression tag	UNP A0A979GQH9
B	883	MET	-	expression tag	UNP A0A979GQH9
B	884	GLY	-	expression tag	UNP A0A979GQH9
B	885	ARG	-	expression tag	UNP A0A979GQH9
B	886	GLY	-	expression tag	UNP A0A979GQH9
B	887	SER	-	expression tag	UNP A0A979GQH9
B	959	SNN	ASP	conflict	UNP A0A979GQH9
B	1013	LEU	-	expression tag	UNP A0A979GQH9
B	1014	GLU	-	expression tag	UNP A0A979GQH9
B	1015	HIS	-	expression tag	UNP A0A979GQH9
B	1016	HIS	-	expression tag	UNP A0A979GQH9
B	1017	HIS	-	expression tag	UNP A0A979GQH9
B	1018	HIS	-	expression tag	UNP A0A979GQH9
B	1019	HIS	-	expression tag	UNP A0A979GQH9
B	1020	HIS	-	expression tag	UNP A0A979GQH9
C	874	MET	-	initiating methionine	UNP A0A979GQH9
C	875	ALA	-	expression tag	UNP A0A979GQH9
C	876	SER	-	expression tag	UNP A0A979GQH9
C	877	MET	-	expression tag	UNP A0A979GQH9
C	878	THR	-	expression tag	UNP A0A979GQH9
C	879	GLY	-	expression tag	UNP A0A979GQH9
C	880	GLY	-	expression tag	UNP A0A979GQH9
C	881	GLN	-	expression tag	UNP A0A979GQH9
C	882	GLN	-	expression tag	UNP A0A979GQH9
C	883	MET	-	expression tag	UNP A0A979GQH9
C	884	GLY	-	expression tag	UNP A0A979GQH9
C	885	ARG	-	expression tag	UNP A0A979GQH9
C	886	GLY	-	expression tag	UNP A0A979GQH9
C	887	SER	-	expression tag	UNP A0A979GQH9
C	959	SNN	ASP	conflict	UNP A0A979GQH9
C	1013	LEU	-	expression tag	UNP A0A979GQH9
C	1014	GLU	-	expression tag	UNP A0A979GQH9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1015	HIS	-	expression tag	UNP A0A979GQH9
C	1016	HIS	-	expression tag	UNP A0A979GQH9
C	1017	HIS	-	expression tag	UNP A0A979GQH9
C	1018	HIS	-	expression tag	UNP A0A979GQH9
C	1019	HIS	-	expression tag	UNP A0A979GQH9
C	1020	HIS	-	expression tag	UNP A0A979GQH9
D	874	MET	-	initiating methionine	UNP A0A979GQH9
D	875	ALA	-	expression tag	UNP A0A979GQH9
D	876	SER	-	expression tag	UNP A0A979GQH9
D	877	MET	-	expression tag	UNP A0A979GQH9
D	878	THR	-	expression tag	UNP A0A979GQH9
D	879	GLY	-	expression tag	UNP A0A979GQH9
D	880	GLY	-	expression tag	UNP A0A979GQH9
D	881	GLN	-	expression tag	UNP A0A979GQH9
D	882	GLN	-	expression tag	UNP A0A979GQH9
D	883	MET	-	expression tag	UNP A0A979GQH9
D	884	GLY	-	expression tag	UNP A0A979GQH9
D	885	ARG	-	expression tag	UNP A0A979GQH9
D	886	GLY	-	expression tag	UNP A0A979GQH9
D	887	SER	-	expression tag	UNP A0A979GQH9
D	959	SNN	ASP	conflict	UNP A0A979GQH9
D	1013	LEU	-	expression tag	UNP A0A979GQH9
D	1014	GLU	-	expression tag	UNP A0A979GQH9
D	1015	HIS	-	expression tag	UNP A0A979GQH9
D	1016	HIS	-	expression tag	UNP A0A979GQH9
D	1017	HIS	-	expression tag	UNP A0A979GQH9
D	1018	HIS	-	expression tag	UNP A0A979GQH9
D	1019	HIS	-	expression tag	UNP A0A979GQH9
D	1020	HIS	-	expression tag	UNP A0A979GQH9

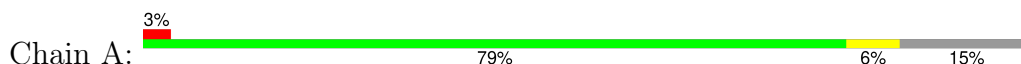
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	86	Total O 86 86	0	0
2	B	82	Total O 82 82	0	0
2	C	81	Total O 81 81	0	0
2	D	64	Total O 64 64	0	0

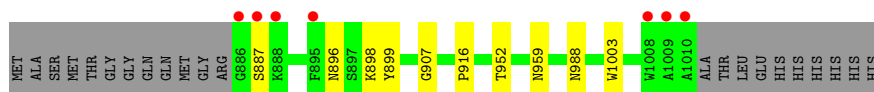
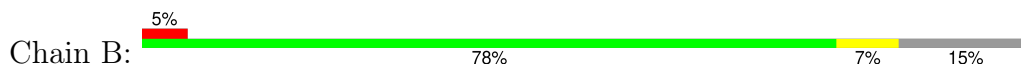
3 Residue-property plots [i](#)

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.

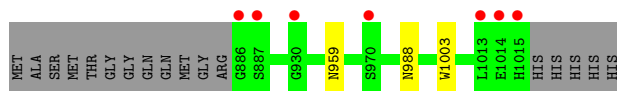
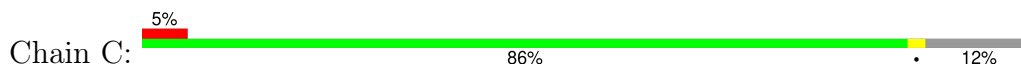
- Molecule 1: Glycoside hydrolase family 18



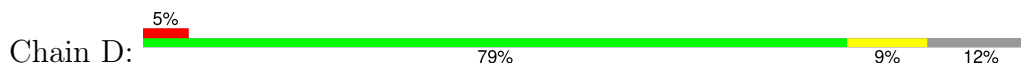
- Molecule 1: Glycoside hydrolase family 18



- Molecule 1: Glycoside hydrolase family 18



- Molecule 1: Glycoside hydrolase family 18



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	34.04Å 59.06Å 68.03Å 90.98° 92.55° 105.64°	Depositor
Resolution (Å)	33.97 – 1.56 33.97 – 1.56	Depositor EDS
% Data completeness (in resolution range)	95.3 (33.97-1.56) 95.5 (33.97-1.56)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.31 (at 1.56Å)	Xtriage
Refinement program	PHENIX 1.15.2_3472	Depositor
R, R_{free}	0.175 , 0.206 0.175 , 0.206	Depositor DCC
R_{free} test set	4977 reflections (6.83%)	wwPDB-VP
Wilson B-factor (Å ²)	24.4	Xtriage
Anisotropy	0.187	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 35.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.034 for -h,h+k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4296	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.08% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SNN

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	0.27	0/978	0.52	0/1327
1	B	0.33	0/996	0.58	0/1348
1	C	0.27	0/1030	0.51	0/1397
1	D	0.25	0/1037	0.49	0/1406
All	All	0.28	0/4041	0.52	0/5478

There are no bond length outliers.

There are no bond angle outliers.

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	964	0	878	6	0
1	B	982	0	898	5	0
1	C	1015	0	924	1	0
1	D	1022	0	928	8	0
2	A	86	0	0	1	0
2	B	82	0	0	0	0
2	C	81	0	0	0	0
2	D	64	0	0	1	0
All	All	4296	0	3628	17	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 2.

All (17) 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:901[B]:ASN:OD1	1:D:903:ARG:NH1	2.34	0.61
1:D:904:ASN:OD1	1:D:938:ASN:HB3	2.01	0.60
1:D:997:ARG:HG2	2:D:1136:HOH:O	2.10	0.52
1:A:952:THR:HG21	1:B:907:GLY:HA2	1.94	0.50
1:A:988:ASN:HB2	1:A:1003:TRP:CE2	2.47	0.50
1:D:887:SER:HB3	1:D:1010:ALA:HB1	1.94	0.49
1:B:988:ASN:HB2	1:B:1003:TRP:CE2	2.49	0.47
1:D:901[B]:ASN:ND2	1:D:912:ASP:OD1	2.40	0.47
1:B:899:TYR:CE1	1:B:916:PRO:HD3	2.51	0.46
1:A:946:ASN:HB2	1:A:961:TRP:CE2	2.52	0.45
1:C:988:ASN:HB2	1:C:1003:TRP:CE2	2.53	0.44
1:A:907:GLY:HA2	1:B:952:THR:HG21	2.00	0.43
1:A:997:ARG:HG2	2:A:1128:HOH:O	2.18	0.43
1:B:896:ASN:ND2	1:B:898[B]:LYS:HD2	2.35	0.42
1:D:988:ASN:HB2	1:D:1003:TRP:CE2	2.55	0.41
1:D:971:ASP:OD1	1:D:1005[A]:ARG:HD3	2.21	0.40
1:A:957:ALA:HB2	1:D:887:SER:CB	2.51	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	121/147 (82%)	115 (95%)	6 (5%)	0	100	100
1	B	123/147 (84%)	118 (96%)	5 (4%)	0	100	100
1	C	128/147 (87%)	122 (95%)	6 (5%)	0	100	100
1	D	128/147 (87%)	120 (94%)	8 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	500/588 (85%)	475 (95%)	25 (5%)	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	99/116 (85%)	99 (100%)	0	100	100
1	B	101/116 (87%)	100 (99%)	1 (1%)	68	47
1	C	105/116 (90%)	105 (100%)	0	100	100
1	D	106/116 (91%)	106 (100%)	0	100	100
All	All	411/464 (89%)	410 (100%)	1 (0%)	87	79

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	887	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	906	GLN
1	A	929	ASN
1	A	991	GLN
1	A	1007	ASN
1	B	954	ASN
1	B	991	GLN
1	C	917	GLN
1	C	991	GLN
1	D	1007	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.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
1	SNN	C	959	1	7,8,8	1.38	1 (14%)	10,11,11	3.25	5 (50%)
1	SNN	A	959	1	7,8,8	1.87	2 (28%)	10,11,11	2.33	5 (50%)
1	SNN	D	959	1	7,8,8	1.47	2 (28%)	10,11,11	2.81	5 (50%)
1	SNN	B	959	1	7,8,8	1.77	2 (28%)	10,11,11	2.72	6 (60%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	SNN	C	959	1	-	-	0/1/1/1
1	SNN	A	959	1	-	-	0/1/1/1
1	SNN	D	959	1	-	-	0/1/1/1
1	SNN	B	959	1	-	-	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	959	SNN	C-N1	-4.23	1.32	1.37
1	B	959	SNN	C-N1	-3.71	1.32	1.37
1	D	959	SNN	C-N1	-2.90	1.33	1.37
1	C	959	SNN	C-N1	-2.80	1.34	1.37
1	B	959	SNN	C5-N1	-2.55	1.33	1.37
1	A	959	SNN	C5-N1	-2.39	1.33	1.37
1	D	959	SNN	C5-N1	-2.25	1.33	1.37

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	959	SNN	O-C-CA	-5.99	121.88	126.19
1	B	959	SNN	CA-C-N1	5.45	111.58	107.30
1	D	959	SNN	O-C-CA	-4.88	122.68	126.19
1	D	959	SNN	CA-C-N1	4.87	111.12	107.30
1	C	959	SNN	CA-C-N1	4.76	111.04	107.30
1	A	959	SNN	CA-C-N1	4.37	110.73	107.30
1	B	959	SNN	O-C-CA	-3.98	123.33	126.19
1	C	959	SNN	C5-N1-C	-3.82	110.12	113.96
1	C	959	SNN	O5-C5-C4	-3.81	121.62	126.48
1	A	959	SNN	O-C-CA	-3.74	123.50	126.19
1	D	959	SNN	C5-N1-C	-3.15	110.80	113.96
1	B	959	SNN	C5-N1-C	-3.04	110.91	113.96
1	D	959	SNN	O5-C5-C4	-3.03	122.62	126.48
1	D	959	SNN	C4-C5-N1	2.69	111.21	108.02
1	A	959	SNN	C4-C5-N1	2.60	111.11	108.02
1	B	959	SNN	O5-C5-C4	-2.58	123.19	126.48
1	B	959	SNN	C4-C5-N1	2.54	111.04	108.02
1	A	959	SNN	O5-C5-C4	-2.50	123.29	126.48
1	C	959	SNN	C4-C5-N1	2.42	110.89	108.02
1	A	959	SNN	C5-N1-C	-2.13	111.83	113.96
1	B	959	SNN	C4-CA-C	-2.09	101.95	104.30

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 ⓘ

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	124/147 (84%)	0.28	5 (4%)	42	50	18, 26, 39, 55	0
1	B	124/147 (84%)	0.32	7 (5%)	30	36	13, 26, 42, 60	2 (1%)
1	C	129/147 (87%)	0.33	7 (5%)	31	38	10, 27, 44, 56	2 (1%)
1	D	128/147 (87%)	0.63	8 (6%)	26	31	14, 31, 50, 58	3 (2%)
All	All	505/588 (85%)	0.39	27 (5%)	32	39	10, 27, 46, 60	7 (1%)

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	904	ASN	7.2
1	B	1010	ALA	5.0
1	B	886	GLY	4.3
1	C	886	GLY	4.2
1	C	1015	HIS	3.7
1	A	1011	ALA	3.2
1	C	887	SER	3.2
1	D	1013	LEU	3.1
1	D	887	SER	3.1
1	A	928	GLY	3.1
1	C	970	SER	2.9
1	B	1009	ALA	2.9
1	A	1010	ALA	2.8
1	A	930	GLY	2.7
1	D	970	SER	2.7
1	C	1014	GLU	2.6
1	D	971	ASP	2.6
1	D	895	PHE	2.5
1	C	1013	LEU	2.5
1	C	930	GLY	2.4
1	B	887	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	929	ASN	2.4
1	D	1015	HIS	2.2
1	B	888	LYS	2.2
1	B	1008	TRP	2.1
1	D	972	GLY	2.1
1	B	895	PHE	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
1	SNN	C	959	8/8	0.92	0.09	26,30,33,33	0
1	SNN	D	959	8/8	0.92	0.08	31,37,38,38	0
1	SNN	B	959	8/8	0.94	0.08	22,28,33,35	0
1	SNN	A	959	8/8	0.96	0.07	25,29,34,37	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.