



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:33 AM UTC

PDB ID : 6UB8 / pdb_00006ub8
Title : Crystal structure of a GH128 (subgroup VI) exo-beta-1,3-glucanase from Aureobasidium namibiae (AnGH128_VI)
Authors : Santos, C.R.; Vieira, P.S.; Domingues, M.N.; Cordeiro, R.L.; Tomazini, A.; Murakami, M.T.
Deposited on : 2019-09-11
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
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

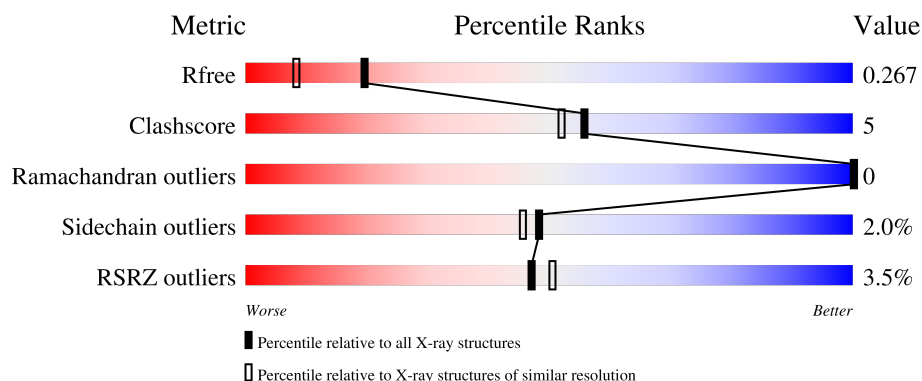
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	266	2% 79% 14% • 6%
1	B	266	% 82% 11% 6%
1	C	266	3% 80% 12% • 6%
1	D	266	3% 79% 12% • 6%
1	E	266	% 82% 10% • 7%

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Mol	Chain	Length	Quality of chain
1	F	266	
1	G	266	
1	H	266	
1	I	266	
1	J	266	

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	GOL	G	302	-	-	X	-
2	GOL	I	302	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 21733 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 Glyco_hydro_cc domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	249	Total	C	N	O	S	0	0	0
			2040	1305	354	365	16			
1	B	249	Total	C	N	O	S	0	1	0
			2047	1312	354	365	16			
1	C	249	Total	C	N	O	S	0	0	0
			2040	1305	354	365	16			
1	D	249	Total	C	N	O	S	0	0	0
			2040	1305	354	365	16			
1	E	248	Total	C	N	O	S	0	0	0
			2036	1303	353	364	16			
1	F	248	Total	C	N	O	S	0	0	0
			2036	1303	353	364	16			
1	G	249	Total	C	N	O	S	0	0	0
			2040	1305	354	365	16			
1	H	248	Total	C	N	O	S	0	0	0
			2036	1303	353	364	16			
1	I	249	Total	C	N	O	S	0	0	0
			2040	1305	354	365	16			
1	J	249	Total	C	N	O	S	0	0	0
			2040	1305	354	365	16			

There are 200 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP A0A074W9U7
A	-18	GLY	-	expression tag	UNP A0A074W9U7
A	-17	SER	-	expression tag	UNP A0A074W9U7
A	-16	SER	-	expression tag	UNP A0A074W9U7
A	-15	HIS	-	expression tag	UNP A0A074W9U7
A	-14	HIS	-	expression tag	UNP A0A074W9U7
A	-13	HIS	-	expression tag	UNP A0A074W9U7
A	-12	HIS	-	expression tag	UNP A0A074W9U7
A	-11	HIS	-	expression tag	UNP A0A074W9U7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	HIS	-	expression tag	UNP A0A074W9U7
A	-9	SER	-	expression tag	UNP A0A074W9U7
A	-8	SER	-	expression tag	UNP A0A074W9U7
A	-7	GLY	-	expression tag	UNP A0A074W9U7
A	-6	LEU	-	expression tag	UNP A0A074W9U7
A	-5	VAL	-	expression tag	UNP A0A074W9U7
A	-4	PRO	-	expression tag	UNP A0A074W9U7
A	-3	ALA	-	expression tag	UNP A0A074W9U7
A	-2	GLY	-	expression tag	UNP A0A074W9U7
A	-1	SER	-	expression tag	UNP A0A074W9U7
A	0	HIS	-	expression tag	UNP A0A074W9U7
B	-19	MET	-	initiating methionine	UNP A0A074W9U7
B	-18	GLY	-	expression tag	UNP A0A074W9U7
B	-17	SER	-	expression tag	UNP A0A074W9U7
B	-16	SER	-	expression tag	UNP A0A074W9U7
B	-15	HIS	-	expression tag	UNP A0A074W9U7
B	-14	HIS	-	expression tag	UNP A0A074W9U7
B	-13	HIS	-	expression tag	UNP A0A074W9U7
B	-12	HIS	-	expression tag	UNP A0A074W9U7
B	-11	HIS	-	expression tag	UNP A0A074W9U7
B	-10	HIS	-	expression tag	UNP A0A074W9U7
B	-9	SER	-	expression tag	UNP A0A074W9U7
B	-8	SER	-	expression tag	UNP A0A074W9U7
B	-7	GLY	-	expression tag	UNP A0A074W9U7
B	-6	LEU	-	expression tag	UNP A0A074W9U7
B	-5	VAL	-	expression tag	UNP A0A074W9U7
B	-4	PRO	-	expression tag	UNP A0A074W9U7
B	-3	ALA	-	expression tag	UNP A0A074W9U7
B	-2	GLY	-	expression tag	UNP A0A074W9U7
B	-1	SER	-	expression tag	UNP A0A074W9U7
B	0	HIS	-	expression tag	UNP A0A074W9U7
C	-19	MET	-	initiating methionine	UNP A0A074W9U7
C	-18	GLY	-	expression tag	UNP A0A074W9U7
C	-17	SER	-	expression tag	UNP A0A074W9U7
C	-16	SER	-	expression tag	UNP A0A074W9U7
C	-15	HIS	-	expression tag	UNP A0A074W9U7
C	-14	HIS	-	expression tag	UNP A0A074W9U7
C	-13	HIS	-	expression tag	UNP A0A074W9U7
C	-12	HIS	-	expression tag	UNP A0A074W9U7
C	-11	HIS	-	expression tag	UNP A0A074W9U7
C	-10	HIS	-	expression tag	UNP A0A074W9U7
C	-9	SER	-	expression tag	UNP A0A074W9U7

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-8	SER	-	expression tag	UNP A0A074W9U7
C	-7	GLY	-	expression tag	UNP A0A074W9U7
C	-6	LEU	-	expression tag	UNP A0A074W9U7
C	-5	VAL	-	expression tag	UNP A0A074W9U7
C	-4	PRO	-	expression tag	UNP A0A074W9U7
C	-3	ALA	-	expression tag	UNP A0A074W9U7
C	-2	GLY	-	expression tag	UNP A0A074W9U7
C	-1	SER	-	expression tag	UNP A0A074W9U7
C	0	HIS	-	expression tag	UNP A0A074W9U7
D	-19	MET	-	initiating methionine	UNP A0A074W9U7
D	-18	GLY	-	expression tag	UNP A0A074W9U7
D	-17	SER	-	expression tag	UNP A0A074W9U7
D	-16	SER	-	expression tag	UNP A0A074W9U7
D	-15	HIS	-	expression tag	UNP A0A074W9U7
D	-14	HIS	-	expression tag	UNP A0A074W9U7
D	-13	HIS	-	expression tag	UNP A0A074W9U7
D	-12	HIS	-	expression tag	UNP A0A074W9U7
D	-11	HIS	-	expression tag	UNP A0A074W9U7
D	-10	HIS	-	expression tag	UNP A0A074W9U7
D	-9	SER	-	expression tag	UNP A0A074W9U7
D	-8	SER	-	expression tag	UNP A0A074W9U7
D	-7	GLY	-	expression tag	UNP A0A074W9U7
D	-6	LEU	-	expression tag	UNP A0A074W9U7
D	-5	VAL	-	expression tag	UNP A0A074W9U7
D	-4	PRO	-	expression tag	UNP A0A074W9U7
D	-3	ALA	-	expression tag	UNP A0A074W9U7
D	-2	GLY	-	expression tag	UNP A0A074W9U7
D	-1	SER	-	expression tag	UNP A0A074W9U7
D	0	HIS	-	expression tag	UNP A0A074W9U7
E	-19	MET	-	initiating methionine	UNP A0A074W9U7
E	-18	GLY	-	expression tag	UNP A0A074W9U7
E	-17	SER	-	expression tag	UNP A0A074W9U7
E	-16	SER	-	expression tag	UNP A0A074W9U7
E	-15	HIS	-	expression tag	UNP A0A074W9U7
E	-14	HIS	-	expression tag	UNP A0A074W9U7
E	-13	HIS	-	expression tag	UNP A0A074W9U7
E	-12	HIS	-	expression tag	UNP A0A074W9U7
E	-11	HIS	-	expression tag	UNP A0A074W9U7
E	-10	HIS	-	expression tag	UNP A0A074W9U7
E	-9	SER	-	expression tag	UNP A0A074W9U7
E	-8	SER	-	expression tag	UNP A0A074W9U7
E	-7	GLY	-	expression tag	UNP A0A074W9U7

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-6	LEU	-	expression tag	UNP A0A074W9U7
E	-5	VAL	-	expression tag	UNP A0A074W9U7
E	-4	PRO	-	expression tag	UNP A0A074W9U7
E	-3	ALA	-	expression tag	UNP A0A074W9U7
E	-2	GLY	-	expression tag	UNP A0A074W9U7
E	-1	SER	-	expression tag	UNP A0A074W9U7
E	0	HIS	-	expression tag	UNP A0A074W9U7
F	-19	MET	-	initiating methionine	UNP A0A074W9U7
F	-18	GLY	-	expression tag	UNP A0A074W9U7
F	-17	SER	-	expression tag	UNP A0A074W9U7
F	-16	SER	-	expression tag	UNP A0A074W9U7
F	-15	HIS	-	expression tag	UNP A0A074W9U7
F	-14	HIS	-	expression tag	UNP A0A074W9U7
F	-13	HIS	-	expression tag	UNP A0A074W9U7
F	-12	HIS	-	expression tag	UNP A0A074W9U7
F	-11	HIS	-	expression tag	UNP A0A074W9U7
F	-10	HIS	-	expression tag	UNP A0A074W9U7
F	-9	SER	-	expression tag	UNP A0A074W9U7
F	-8	SER	-	expression tag	UNP A0A074W9U7
F	-7	GLY	-	expression tag	UNP A0A074W9U7
F	-6	LEU	-	expression tag	UNP A0A074W9U7
F	-5	VAL	-	expression tag	UNP A0A074W9U7
F	-4	PRO	-	expression tag	UNP A0A074W9U7
F	-3	ALA	-	expression tag	UNP A0A074W9U7
F	-2	GLY	-	expression tag	UNP A0A074W9U7
F	-1	SER	-	expression tag	UNP A0A074W9U7
F	0	HIS	-	expression tag	UNP A0A074W9U7
G	-19	MET	-	initiating methionine	UNP A0A074W9U7
G	-18	GLY	-	expression tag	UNP A0A074W9U7
G	-17	SER	-	expression tag	UNP A0A074W9U7
G	-16	SER	-	expression tag	UNP A0A074W9U7
G	-15	HIS	-	expression tag	UNP A0A074W9U7
G	-14	HIS	-	expression tag	UNP A0A074W9U7
G	-13	HIS	-	expression tag	UNP A0A074W9U7
G	-12	HIS	-	expression tag	UNP A0A074W9U7
G	-11	HIS	-	expression tag	UNP A0A074W9U7
G	-10	HIS	-	expression tag	UNP A0A074W9U7
G	-9	SER	-	expression tag	UNP A0A074W9U7
G	-8	SER	-	expression tag	UNP A0A074W9U7
G	-7	GLY	-	expression tag	UNP A0A074W9U7
G	-6	LEU	-	expression tag	UNP A0A074W9U7
G	-5	VAL	-	expression tag	UNP A0A074W9U7

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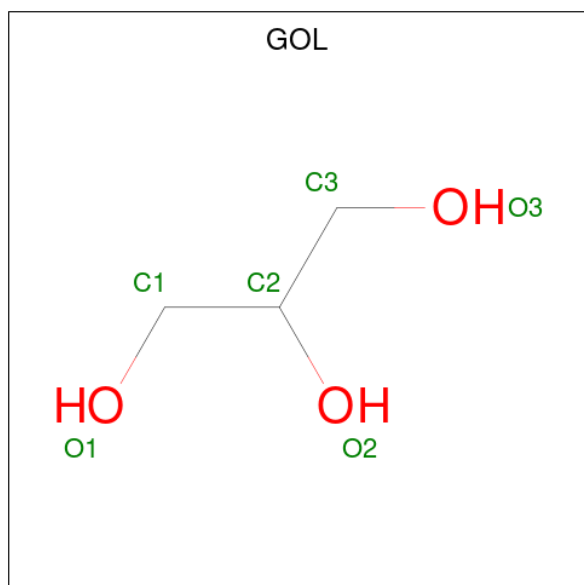
Chain	Residue	Modelled	Actual	Comment	Reference
G	-4	PRO	-	expression tag	UNP A0A074W9U7
G	-3	ALA	-	expression tag	UNP A0A074W9U7
G	-2	GLY	-	expression tag	UNP A0A074W9U7
G	-1	SER	-	expression tag	UNP A0A074W9U7
G	0	HIS	-	expression tag	UNP A0A074W9U7
H	-19	MET	-	initiating methionine	UNP A0A074W9U7
H	-18	GLY	-	expression tag	UNP A0A074W9U7
H	-17	SER	-	expression tag	UNP A0A074W9U7
H	-16	SER	-	expression tag	UNP A0A074W9U7
H	-15	HIS	-	expression tag	UNP A0A074W9U7
H	-14	HIS	-	expression tag	UNP A0A074W9U7
H	-13	HIS	-	expression tag	UNP A0A074W9U7
H	-12	HIS	-	expression tag	UNP A0A074W9U7
H	-11	HIS	-	expression tag	UNP A0A074W9U7
H	-10	HIS	-	expression tag	UNP A0A074W9U7
H	-9	SER	-	expression tag	UNP A0A074W9U7
H	-8	SER	-	expression tag	UNP A0A074W9U7
H	-7	GLY	-	expression tag	UNP A0A074W9U7
H	-6	LEU	-	expression tag	UNP A0A074W9U7
H	-5	VAL	-	expression tag	UNP A0A074W9U7
H	-4	PRO	-	expression tag	UNP A0A074W9U7
H	-3	ALA	-	expression tag	UNP A0A074W9U7
H	-2	GLY	-	expression tag	UNP A0A074W9U7
H	-1	SER	-	expression tag	UNP A0A074W9U7
H	0	HIS	-	expression tag	UNP A0A074W9U7
I	-19	MET	-	initiating methionine	UNP A0A074W9U7
I	-18	GLY	-	expression tag	UNP A0A074W9U7
I	-17	SER	-	expression tag	UNP A0A074W9U7
I	-16	SER	-	expression tag	UNP A0A074W9U7
I	-15	HIS	-	expression tag	UNP A0A074W9U7
I	-14	HIS	-	expression tag	UNP A0A074W9U7
I	-13	HIS	-	expression tag	UNP A0A074W9U7
I	-12	HIS	-	expression tag	UNP A0A074W9U7
I	-11	HIS	-	expression tag	UNP A0A074W9U7
I	-10	HIS	-	expression tag	UNP A0A074W9U7
I	-9	SER	-	expression tag	UNP A0A074W9U7
I	-8	SER	-	expression tag	UNP A0A074W9U7
I	-7	GLY	-	expression tag	UNP A0A074W9U7
I	-6	LEU	-	expression tag	UNP A0A074W9U7
I	-5	VAL	-	expression tag	UNP A0A074W9U7
I	-4	PRO	-	expression tag	UNP A0A074W9U7
I	-3	ALA	-	expression tag	UNP A0A074W9U7

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Chain	Residue	Modelled	Actual	Comment	Reference
I	-2	GLY	-	expression tag	UNP A0A074W9U7
I	-1	SER	-	expression tag	UNP A0A074W9U7
I	0	HIS	-	expression tag	UNP A0A074W9U7
J	-19	MET	-	initiating methionine	UNP A0A074W9U7
J	-18	GLY	-	expression tag	UNP A0A074W9U7
J	-17	SER	-	expression tag	UNP A0A074W9U7
J	-16	SER	-	expression tag	UNP A0A074W9U7
J	-15	HIS	-	expression tag	UNP A0A074W9U7
J	-14	HIS	-	expression tag	UNP A0A074W9U7
J	-13	HIS	-	expression tag	UNP A0A074W9U7
J	-12	HIS	-	expression tag	UNP A0A074W9U7
J	-11	HIS	-	expression tag	UNP A0A074W9U7
J	-10	HIS	-	expression tag	UNP A0A074W9U7
J	-9	SER	-	expression tag	UNP A0A074W9U7
J	-8	SER	-	expression tag	UNP A0A074W9U7
J	-7	GLY	-	expression tag	UNP A0A074W9U7
J	-6	LEU	-	expression tag	UNP A0A074W9U7
J	-5	VAL	-	expression tag	UNP A0A074W9U7
J	-4	PRO	-	expression tag	UNP A0A074W9U7
J	-3	ALA	-	expression tag	UNP A0A074W9U7
J	-2	GLY	-	expression tag	UNP A0A074W9U7
J	-1	SER	-	expression tag	UNP A0A074W9U7
J	0	HIS	-	expression tag	UNP A0A074W9U7

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	A	1	Total C O 6 3 3	0	0
2	B	1	Total C O 6 3 3	0	0
2	B	1	Total C O 6 3 3	0	0
2	B	1	Total C O 6 3 3	0	0
2	C	1	Total C O 6 3 3	0	0
2	C	1	Total C O 6 3 3	0	0
2	D	1	Total C O 6 3 3	0	0
2	D	1	Total C O 6 3 3	0	0
2	D	1	Total C O 6 3 3	0	0
2	E	1	Total C O 6 3 3	0	0
2	F	1	Total C O 6 3 3	0	0
2	F	1	Total C O 6 3 3	0	0
2	F	1	Total C O 6 3 3	0	0
2	F	1	Total C O 6 3 3	0	0
2	F	1	Total C O 6 3 3	0	0
2	G	1	Total C O 6 3 3	0	0
2	G	1	Total C O 6 3 3	0	0
2	G	1	Total C O 6 3 3	0	0
2	G	1	Total C O 6 3 3	0	0
2	G	1	Total C O 6 3 3	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	G	1	Total C O 6 3 3	0	0
2	H	1	Total C O 6 3 3	0	0
2	H	1	Total C O 6 3 3	0	0
2	I	1	Total C O 6 3 3	0	0
2	I	1	Total C O 6 3 3	0	0
2	I	1	Total C O 6 3 3	0	0
2	I	1	Total C O 6 3 3	0	0
2	J	1	Total C O 6 3 3	0	0
2	J	1	Total C O 6 3 3	0	0
2	J	1	Total C O 6 3 3	0	0
2	J	1	Total C O 6 3 3	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	108	Total O 108 108	0	0
3	B	141	Total O 141 141	0	0
3	C	97	Total O 97 97	0	0
3	D	93	Total O 93 93	0	0
3	E	104	Total O 104 104	0	0
3	F	132	Total O 132 132	0	0
3	G	132	Total O 132 132	0	0
3	H	83	Total O 83 83	0	0

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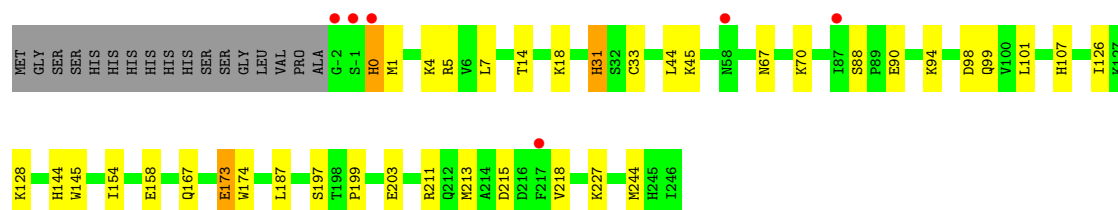
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	I	113	Total 113	O 113	0	0
3	J	137	Total 137	O 137	0	0

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.

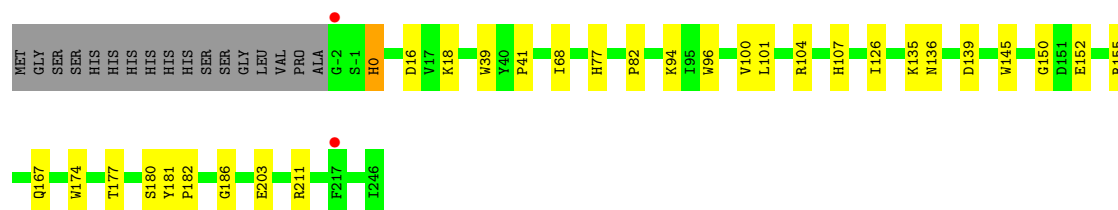
- Molecule 1: Glyco_hydro_cc domain-containing protein

Chain A: 



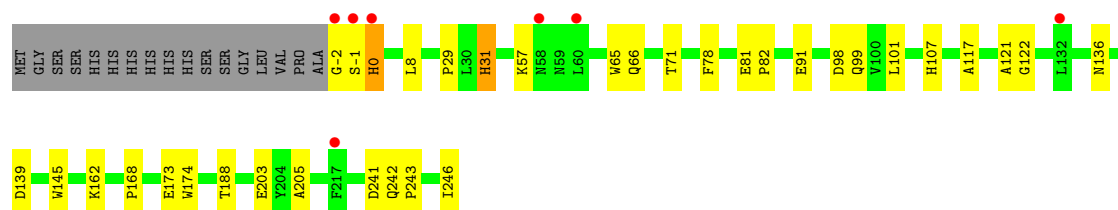
- Molecule 1: Glyco_hydro_cc domain-containing protein

Chain B: 



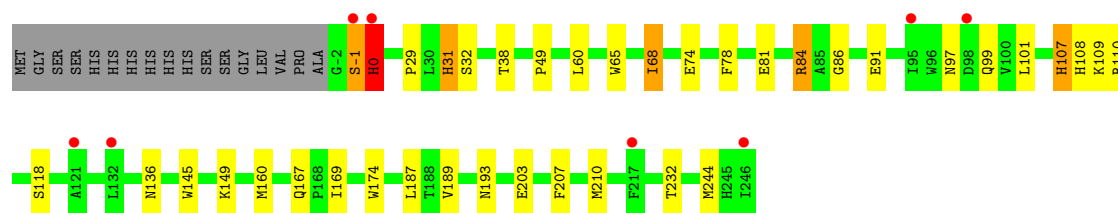
- Molecule 1: Glyco_hydro_cc domain-containing protein

Chain C: 

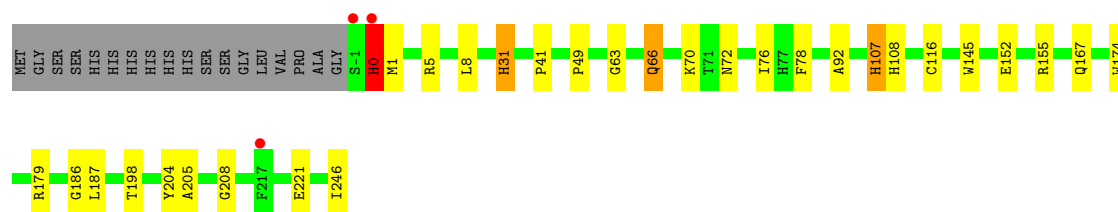
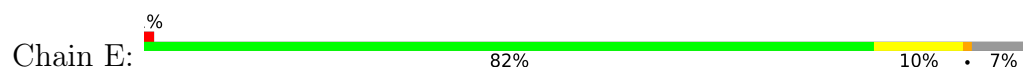


- Molecule 1: Glyco_hydro_cc domain-containing protein

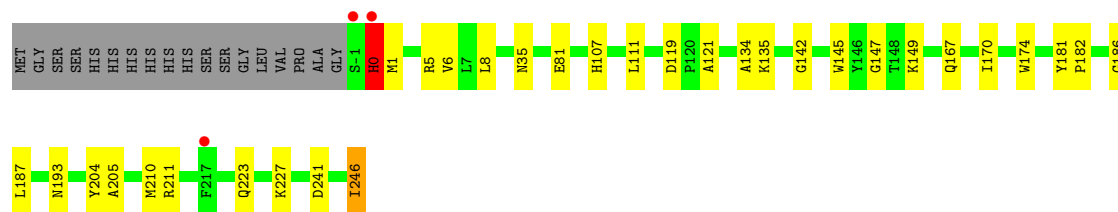
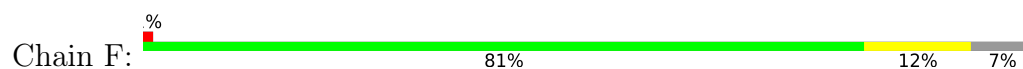
Chain D: 



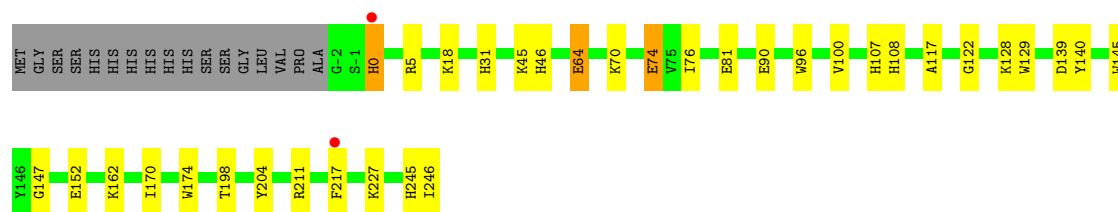
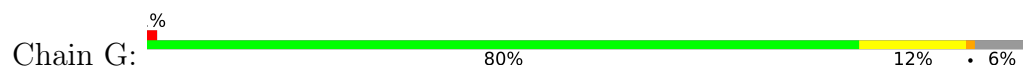
- Molecule 1: Glyco_hydro_cc domain-containing protein



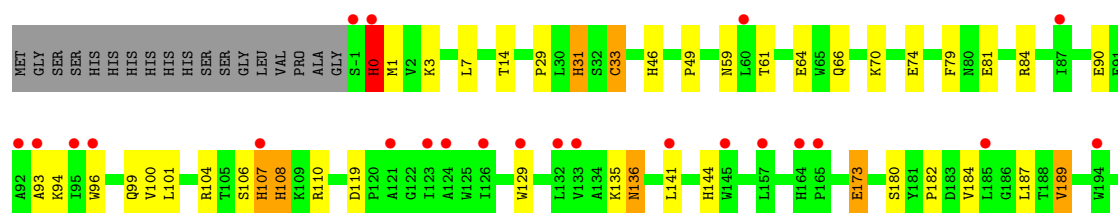
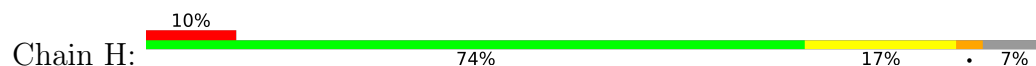
- Molecule 1: Glyco_hydro_cc domain-containing protein



- Molecule 1: Glyco_hydro_cc domain-containing protein

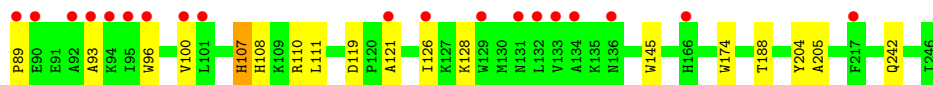
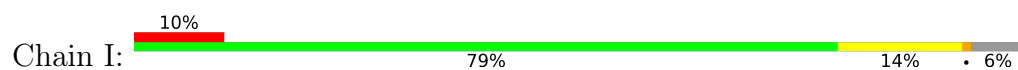


- Molecule 1: Glyco_hydro_cc domain-containing protein

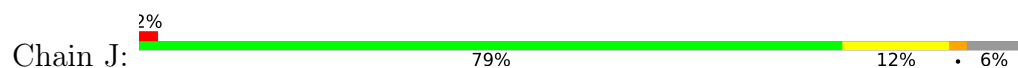




- Molecule 1: Glyco_hydro_cc domain-containing protein



- Molecule 1: Glyco_hydro_cc domain-containing protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	117.99Å 151.86Å 166.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.21 – 1.90 48.21 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.3 (48.21-1.90) 99.3 (48.21-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.08 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.8.0253	Depositor
R, R_{free}	0.222 , 0.264 0.228 , 0.267	Depositor DCC
R_{free} test set	11663 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	28.6	Xtriage
Anisotropy	0.240	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 25.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	21733	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.48% 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:
GOL

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	1.24	6/2108 (0.3%)	1.46	7/2864 (0.2%)
1	B	1.27	7/2120 (0.3%)	1.45	9/2880 (0.3%)
1	C	1.21	3/2108 (0.1%)	1.47	14/2864 (0.5%)
1	D	1.22	8/2108 (0.4%)	1.46	7/2864 (0.2%)
1	E	1.24	6/2104 (0.3%)	1.48	13/2859 (0.5%)
1	F	1.26	5/2104 (0.2%)	1.47	10/2859 (0.3%)
1	G	1.28	7/2108 (0.3%)	1.54	13/2864 (0.5%)
1	H	1.27	6/2104 (0.3%)	1.57	19/2859 (0.7%)
1	I	1.27	6/2108 (0.3%)	1.46	7/2864 (0.2%)
1	J	1.27	12/2108 (0.6%)	1.49	12/2864 (0.4%)
All	All	1.25	66/21080 (0.3%)	1.49	111/28641 (0.4%)

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
1	D	0	2
1	F	0	1
1	J	0	1
All	All	0	4

All (66) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	0	HIS	CE1-NE2	-10.74	1.21	1.32
1	A	0	HIS	CE1-NE2	-8.41	1.24	1.32
1	I	-2	GLY	N-CA	7.93	1.58	1.45
1	H	107	HIS	CD2-NE2	7.64	1.46	1.37
1	J	110	ARG	C-O	7.33	1.32	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	I	0	HIS	CE1-NE2	7.29	1.39	1.32
1	C	0	HIS	CA-CB	7.08	1.64	1.53
1	E	92	ALA	C-O	7.00	1.32	1.24
1	E	208	GLY	C-O	6.97	1.33	1.23
1	F	111	LEU	C-O	6.92	1.32	1.24
1	G	0	HIS	CD2-NE2	6.84	1.45	1.37
1	G	0	HIS	CA-CB	6.81	1.63	1.53
1	H	33	CYS	C-O	6.72	1.32	1.23
1	B	41	PRO	C-O	-6.67	1.18	1.24
1	C	0	HIS	CE1-NE2	-6.66	1.25	1.32
1	D	107	HIS	CE1-NE2	6.47	1.39	1.32
1	E	0	HIS	CA-CB	6.39	1.62	1.53
1	B	94	LYS	C-O	-6.37	1.16	1.24
1	D	81	GLU	N-CA	6.32	1.50	1.46
1	I	121	ALA	C-O	6.30	1.31	1.24
1	J	70	LYS	C-O	-6.23	1.15	1.23
1	G	81	GLU	N-CA	6.17	1.50	1.46
1	H	107	HIS	ND1-CE1	6.17	1.38	1.32
1	G	0	HIS	ND1-CE1	5.94	1.38	1.32
1	D	232	THR	C-O	5.93	1.31	1.23
1	I	75	VAL	C-O	-5.86	1.18	1.24
1	J	245	HIS	CE1-NE2	5.82	1.38	1.32
1	I	107	HIS	CE1-NE2	5.81	1.38	1.32
1	G	46	HIS	CE1-NE2	5.79	1.38	1.32
1	D	32	SER	C-O	5.77	1.30	1.23
1	F	210	MET	C-O	5.76	1.30	1.23
1	B	68	ILE	C-O	5.74	1.30	1.24
1	A	44	LEU	C-O	5.73	1.31	1.24
1	B	77	HIS	CE1-NE2	5.72	1.38	1.32
1	H	81	GLU	N-CA	5.70	1.50	1.46
1	F	6	VAL	C-O	5.67	1.31	1.24
1	D	78	PHE	C-O	5.65	1.29	1.23
1	A	215	ASP	C-O	5.64	1.29	1.24
1	F	223	GLN	C-O	5.63	1.30	1.23
1	A	199	PRO	C-O	-5.63	1.17	1.24
1	I	111	LEU	C-O	5.44	1.29	1.23
1	J	79	PHE	C-O	5.42	1.30	1.23
1	J	220	PRO	C-O	-5.42	1.17	1.24
1	J	124	ALA	C-O	5.39	1.30	1.24
1	G	129	TRP	C-O	5.38	1.30	1.24
1	J	232	THR	C-O	5.33	1.31	1.23
1	E	107	HIS	CE1-NE2	5.27	1.37	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	J	5	ARG	C-O	5.27	1.30	1.24
1	J	12	THR	C-O	5.26	1.30	1.24
1	H	0	HIS	ND1-CE1	5.25	1.37	1.32
1	E	116	CYS	C-O	5.22	1.30	1.23
1	A	5	ARG	C-O	5.16	1.30	1.24
1	J	174	TRP	C-O	5.15	1.30	1.23
1	A	0	HIS	C-O	-5.14	1.17	1.23
1	F	81	GLU	N-CA	5.14	1.50	1.46
1	H	208	GLY	C-O	5.08	1.30	1.23
1	B	150	GLY	C-O	5.07	1.29	1.23
1	D	29	PRO	C-O	-5.07	1.18	1.23
1	J	50	PHE	C-O	5.06	1.30	1.24
1	B	203	GLU	C-O	5.06	1.29	1.23
1	C	0	HIS	ND1-CE1	5.05	1.37	1.32
1	D	169	ILE	C-O	5.04	1.29	1.24
1	E	76	ILE	C-O	5.04	1.29	1.24
1	B	186	GLY	C-O	5.03	1.29	1.23
1	J	201	VAL	N-CA	5.03	1.52	1.46
1	D	193	ASN	C-O	5.01	1.30	1.24

All (111) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	0	HIS	CA-CB-CG	19.24	133.04	113.80
1	E	0	HIS	CA-CB-CG	14.14	127.94	113.80
1	C	0	HIS	CA-CB-CG	13.95	127.75	113.80
1	J	0	HIS	CA-CB-CG	12.46	126.26	113.80
1	F	0	HIS	CA-CB-CG	12.05	125.85	113.80
1	D	0	HIS	CA-CB-CG	11.32	125.12	113.80
1	A	0	HIS	CA-CB-CG	11.28	125.08	113.80
1	H	107	HIS	CB-CA-C	10.51	127.64	110.09
1	H	0	HIS	CA-CB-CG	9.34	123.14	113.80
1	H	107	HIS	CE1-NE2-CD2	-9.04	99.96	109.00
1	E	198	THR	CB-CA-C	8.40	118.90	110.33
1	J	81	GLU	CB-CA-C	8.40	117.16	111.20
1	F	167	GLN	CB-CA-C	-8.08	96.86	109.55
1	G	0	HIS	ND1-CG-CD2	-7.96	98.14	106.10
1	F	121	ALA	CA-C-N	7.41	128.02	119.94
1	F	121	ALA	C-N-CA	7.41	128.02	119.94
1	C	0	HIS	ND1-CG-CD2	-7.21	98.89	106.10
1	H	107	HIS	CG-CD2-NE2	7.17	114.37	107.20
1	E	41	PRO	CB-CA-C	7.03	117.78	111.87

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	0	HIS	CA-CB-CG	6.93	120.73	113.80
1	I	18	LYS	CB-CA-C	-6.72	99.26	110.68
1	C	98	ASP	CB-CA-C	-6.69	99.55	110.79
1	B	18	LYS	CB-CA-C	-6.69	100.11	110.81
1	H	198	THR	CB-CA-C	6.53	116.64	110.17
1	E	0	HIS	CB-CA-C	6.51	120.46	109.84
1	B	104	ARG	NE-CZ-NH2	6.40	124.96	119.20
1	H	173	GLU	CB-CA-C	-6.39	101.39	110.34
1	D	0	HIS	CB-CA-C	6.39	120.08	109.53
1	C	71	THR	CA-CB-OG1	-6.38	100.02	109.60
1	A	18	LYS	CA-C-N	6.26	129.51	120.38
1	A	18	LYS	C-N-CA	6.26	129.51	120.38
1	I	188	THR	N-CA-C	-6.19	104.44	111.07
1	H	14	THR	N-CA-C	-6.17	104.64	111.36
1	D	108	HIS	CA-CB-CG	-6.10	107.70	113.80
1	B	177	THR	CA-CB-OG1	-6.07	100.49	109.60
1	F	119	ASP	CB-CA-C	6.07	118.26	109.09
1	A	14	THR	N-CA-C	-6.06	104.68	111.28
1	H	107	HIS	CA-CB-CG	6.05	119.85	113.80
1	J	211	ARG	CB-CG-CD	-6.01	97.47	111.30
1	F	147	GLY	CA-C-O	-5.99	115.06	121.35
1	I	93	ALA	CA-C-N	5.98	128.22	120.44
1	I	93	ALA	C-N-CA	5.98	128.22	120.44
1	E	66	GLN	N-CA-C	-5.94	104.80	111.28
1	C	0	HIS	CB-CA-C	5.94	119.99	109.72
1	G	108	HIS	CA-CB-CG	-5.93	107.87	113.80
1	B	16	ASP	CA-CB-CG	5.87	118.47	112.60
1	H	189	VAL	CA-C-O	-5.84	115.20	121.27
1	J	51	ARG	CB-CA-C	-5.77	104.46	110.17
1	F	134	ALA	CA-C-N	5.76	128.00	120.28
1	F	134	ALA	C-N-CA	5.76	128.00	120.28
1	E	167	GLN	CB-CA-C	-5.70	99.51	108.91
1	J	119	ASP	CA-CB-CG	5.70	118.30	112.60
1	F	193	ASN	CA-CB-CG	-5.69	106.91	112.60
1	E	78	PHE	CA-C-N	5.68	128.78	120.71
1	E	78	PHE	C-N-CA	5.68	128.78	120.71
1	G	0	HIS	CB-CA-C	5.65	119.50	109.72
1	C	-1	SER	CA-C-N	5.64	130.78	123.00
1	C	-1	SER	C-N-CA	5.64	130.78	123.00
1	D	109	LYS	CB-CA-C	5.64	118.95	109.48
1	A	244	MET	CB-CA-C	5.62	119.13	109.51
1	C	188	THR	N-CA-C	-5.62	105.07	111.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	184	VAL	N-CA-C	-5.61	105.25	110.53
1	J	64	GLU	N-CA-C	-5.60	105.07	111.07
1	E	186	GLY	CA-C-O	-5.56	115.33	121.05
1	C	0	HIS	CG-CD2-NE2	5.53	112.73	107.20
1	I	188	THR	CB-CA-C	5.53	119.56	110.88
1	B	82	PRO	CB-CA-C	-5.52	104.01	111.85
1	C	121	ALA	CA-C-N	5.51	126.05	119.94
1	C	121	ALA	C-N-CA	5.51	126.05	119.94
1	G	0	HIS	CG-CD2-NE2	5.50	112.70	107.20
1	H	141	LEU	CA-C-N	5.49	127.23	120.92
1	H	141	LEU	C-N-CA	5.49	127.23	120.92
1	H	93	ALA	CA-C-N	5.46	127.53	120.44
1	H	93	ALA	C-N-CA	5.46	127.53	120.44
1	E	108	HIS	CA-CB-CG	-5.44	108.36	113.80
1	E	0	HIS	ND1-CG-CD2	-5.44	100.66	106.10
1	J	167	GLN	O-C-N	-5.42	118.06	121.85
1	A	173	GLU	CB-CA-C	-5.38	100.81	109.89
1	G	198	THR	CB-CA-C	5.37	115.49	110.17
1	G	217	PHE	CB-CA-C	5.37	120.52	110.70
1	I	62	GLY	CA-C-N	5.36	125.93	119.98
1	I	62	GLY	C-N-CA	5.36	125.93	119.98
1	C	29	PRO	CB-CA-C	5.34	117.78	110.52
1	G	139	ASP	CA-CB-CG	-5.34	107.26	112.60
1	G	0	HIS	CE1-NE2-CD2	-5.32	103.69	109.00
1	A	4	LYS	CB-CA-C	5.30	121.92	110.41
1	D	74	GLU	CB-CG-CD	5.29	121.59	112.60
1	J	73	GLU	CA-C-O	-5.29	115.57	121.81
1	H	79	PHE	CA-CB-CG	5.28	119.08	113.80
1	J	108	HIS	CA-CB-CG	-5.28	108.52	113.80
1	B	104	ARG	NE-CZ-NH1	-5.28	116.22	121.50
1	E	66	GLN	CB-CA-C	5.23	119.47	110.79
1	G	64	GLU	N-CA-C	-5.22	105.59	111.28
1	E	1	MET	N-CA-CB	5.22	117.56	109.48
1	H	108	HIS	CA-CB-CG	-5.19	108.61	113.80
1	G	0	HIS	CB-CG-CD2	5.17	137.93	131.20
1	H	129	TRP	CA-C-O	-5.15	115.41	120.82
1	F	241	ASP	CA-CB-CG	5.12	117.72	112.60
1	B	39	TRP	CA-C-N	5.09	130.62	120.94
1	B	39	TRP	C-N-CA	5.09	130.62	120.94
1	H	107	HIS	N-CA-C	-5.09	106.76	113.17
1	J	138	PRO	CA-C-O	-5.09	115.57	122.08
1	H	107	HIS	CB-CG-CD2	5.07	137.80	131.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	110	ARG	CB-CA-C	5.06	118.77	109.46
1	C	162	LYS	CA-C-N	5.05	127.05	120.28
1	C	162	LYS	C-N-CA	5.05	127.05	120.28
1	G	162	LYS	CA-C-N	5.04	126.99	120.44
1	G	162	LYS	C-N-CA	5.04	126.99	120.44
1	D	160	MET	CA-C-O	-5.03	115.09	120.42
1	J	94	LYS	CA-C-N	5.03	127.35	120.46
1	J	94	LYS	C-N-CA	5.03	127.35	120.46

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	0	HIS	Peptide
1	D	86	GLY	Peptide
1	F	0	HIS	Peptide
1	J	0	HIS	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	2040	0	1966	23	0
1	B	2047	0	1971	15	0
1	C	2040	0	1966	22	0
1	D	2040	0	1966	23	0
1	E	2036	0	1963	18	0
1	F	2036	0	1963	19	0
1	G	2040	0	1966	18	0
1	H	2036	0	1963	40	0
1	I	2040	0	1966	32	0
1	J	2040	0	1966	22	0
2	A	18	0	24	0	0
2	B	18	0	24	0	0
2	C	12	0	16	3	0
2	D	18	0	24	1	0
2	E	6	0	8	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	24	0	32	0	0
2	G	42	0	56	7	0
2	H	12	0	16	2	0
2	I	24	0	32	4	0
2	J	24	0	32	2	0
3	A	108	0	0	4	0
3	B	141	0	0	4	0
3	C	97	0	0	0	0
3	D	93	0	0	1	0
3	E	104	0	0	0	0
3	F	132	0	0	2	0
3	G	132	0	0	1	0
3	H	83	0	0	2	0
3	I	113	0	0	1	0
3	J	137	0	0	3	0
All	All	21733	0	19920	192	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 5.

All (192) 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:H:107:HIS:CE1	1:I:0:HIS:NE2	2.09	1.21
1:A:107:HIS:HE2	1:G:0:HIS:CE1	1.59	1.20
1:C:0:HIS:CE1	1:D:107:HIS:HE2	1.58	1.19
1:C:107:HIS:HE2	1:D:0:HIS:CE1	1.63	1.15
1:E:0:HIS:CE1	1:J:107:HIS:NE2	2.18	1.12
1:A:107:HIS:NE2	1:G:0:HIS:CE1	2.15	1.12
1:B:107:HIS:NE2	1:F:0:HIS:CE1	2.20	1.10
3:A:484:HOH:O	1:G:74:GLU:HG3	1.56	1.06
1:H:135:LYS:C	1:H:136:ASN:HD22	1.66	1.02
1:H:0:HIS:CE1	1:I:107:HIS:NE2	2.32	0.98
1:C:107:HIS:NE2	1:D:0:HIS:CE1	2.32	0.97
1:H:107:HIS:CE1	1:I:0:HIS:CE1	2.52	0.97
1:C:0:HIS:CE1	1:D:107:HIS:NE2	2.33	0.97
1:E:107:HIS:NE2	1:J:0:HIS:CE1	2.38	0.92
1:H:107:HIS:NE2	1:I:0:HIS:NE2	2.17	0.91
1:J:141:LEU:HD23	1:J:169:ILE:CD1	2.08	0.83
1:A:0:HIS:CE1	1:G:107:HIS:NE2	2.48	0.82
2:G:301:GOL:O1	2:G:303:GOL:H11	1.78	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:141:LEU:HD23	1:J:169:ILE:HD11	1.61	0.82
1:D:38:THR:HG22	1:D:68:ILE:HD11	1.62	0.80
1:H:107:HIS:HE1	1:I:0:HIS:NE2	1.76	0.78
2:G:302:GOL:H32	1:I:1:MET:HE2	1.64	0.78
1:B:0:HIS:CE1	1:F:107:HIS:NE2	2.53	0.76
1:C:66:GLN:NE2	1:D:-1:SER:OG	2.18	0.76
1:H:107:HIS:CE1	1:I:0:HIS:HE2	1.88	0.72
1:C:57:LYS:NZ	1:C:91:GLU:OE2	2.22	0.71
1:H:136:ASN:HD22	1:H:136:ASN:N	1.91	0.69
1:C:0:HIS:NE2	1:D:107:HIS:NE2	2.38	0.69
1:A:211:ARG:HD3	3:A:454:HOH:O	1.93	0.68
1:B:107:HIS:CE1	1:F:0:HIS:CE1	2.83	0.67
1:A:0:HIS:CE1	1:G:107:HIS:CE1	2.83	0.67
1:G:45:LYS:HE3	3:G:414:HOH:O	1.95	0.66
1:A:0:HIS:HE1	1:G:107:HIS:NE2	1.89	0.66
1:H:59:ASN:HA	1:H:64:GLU:HG2	1.78	0.65
1:B:211:ARG:HD3	3:B:403:HOH:O	1.95	0.65
1:F:187:LEU:C	1:F:187:LEU:HD23	2.23	0.63
1:A:145:TRP:O	1:A:174:TRP:HA	2.00	0.62
1:A:107:HIS:CE1	1:G:0:HIS:CE1	2.87	0.62
1:D:84:ARG:HG3	1:D:84:ARG:HH11	1.64	0.61
1:H:0:HIS:CE1	1:I:107:HIS:CE1	2.88	0.61
1:G:64:GLU:OE2	2:G:302:GOL:H2	2.01	0.61
1:H:107:HIS:NE2	1:I:0:HIS:CE1	2.66	0.60
1:C:241:ASP:OD2	2:C:302:GOL:O3	2.20	0.60
1:H:0:HIS:HE1	1:I:107:HIS:NE2	1.93	0.60
1:G:145:TRP:CH2	1:G:147:GLY:HA3	2.36	0.60
1:B:0:HIS:CE1	1:F:107:HIS:CE1	2.90	0.59
1:H:135:LYS:C	1:H:136:ASN:ND2	2.51	0.58
1:E:0:HIS:CE1	1:J:107:HIS:CE1	2.92	0.58
1:I:37:ASN:OD1	2:I:302:GOL:H31	2.05	0.57
1:A:67:ASN:HA	1:A:70:LYS:HE3	1.87	0.56
1:H:224:LEU:HB3	1:H:235:MET:HE2	1.88	0.55
1:H:59:ASN:HA	1:H:64:GLU:CG	2.36	0.54
1:E:31:HIS:O	1:E:49:PRO:HD2	2.07	0.54
1:E:63:GLY:HA3	1:F:1:MET:CE	2.38	0.54
2:D:402:GOL:H2	1:H:1:MET:HE2	1.89	0.54
1:F:149:LYS:HE3	3:F:498:HOH:O	2.08	0.53
1:J:155:ARG:NH1	3:J:401:HOH:O	2.24	0.53
1:I:96:TRP:HA	1:I:100:VAL:HB	1.88	0.53
1:E:107:HIS:NE2	1:J:0:HIS:ND1	2.57	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:301:GOL:C1	2:G:303:GOL:H11	2.38	0.53
1:D:38:THR:CG2	1:D:68:ILE:HD11	2.35	0.52
1:A:94:LYS:HE2	1:A:98:ASP:OD2	2.08	0.52
1:C:101:LEU:HD11	1:C:136:ASN:HB3	1.91	0.52
1:I:84:ARG:NH1	1:I:119:ASP:OD2	2.41	0.52
1:H:84:ARG:NH2	1:H:119:ASP:OD2	2.42	0.52
1:F:186:GLY:HA2	1:F:246:ILE:HD11	1.91	0.52
1:A:187:LEU:HD23	1:A:187:LEU:C	2.34	0.52
1:B:101:LEU:HD11	1:B:136:ASN:HB3	1.92	0.52
1:H:187:LEU:C	1:H:187:LEU:HD23	2.35	0.52
1:C:101:LEU:HD11	1:C:136:ASN:CB	2.40	0.51
1:J:141:LEU:HB3	1:J:169:ILE:HD13	1.91	0.51
1:C:107:HIS:NE2	1:D:0:HIS:NE2	2.50	0.51
1:H:144:HIS:CD2	1:H:173:GLU:HB2	2.46	0.51
1:H:180:SER:OG	1:H:182:PRO:HD2	2.12	0.50
1:H:31:HIS:O	1:H:49:PRO:HD2	2.12	0.50
1:H:101:LEU:HD11	1:H:136:ASN:CB	2.42	0.50
1:E:145:TRP:O	1:E:174:TRP:HA	2.12	0.50
1:G:90:GLU:OE2	1:G:128:LYS:HE3	2.11	0.49
1:G:117:ALA:O	1:G:122:GLY:HA3	2.11	0.49
1:I:145:TRP:O	1:I:174:TRP:HA	2.12	0.49
1:G:145:TRP:O	1:G:174:TRP:HA	2.13	0.49
1:A:144:HIS:HA	1:A:173:GLU:O	2.13	0.49
1:J:141:LEU:HD23	1:J:169:ILE:HD12	1.95	0.49
1:B:139:ASP:O	1:B:167:GLN:HB2	2.13	0.49
1:F:145:TRP:O	1:F:174:TRP:HA	2.13	0.48
1:A:1:MET:HE2	2:H:501:GOL:H11	1.94	0.48
2:G:302:GOL:H32	1:I:1:MET:CE	2.39	0.48
1:J:145:TRP:O	1:J:174:TRP:HA	2.13	0.48
1:D:84:ARG:HG3	1:D:84:ARG:NH1	2.28	0.48
1:C:173:GLU:OE1	2:C:301:GOL:H31	2.14	0.48
1:B:155:ARG:NH2	3:B:406:HOH:O	2.41	0.48
1:H:106:SER:OG	1:H:107:HIS:HD2	1.96	0.48
1:E:0:HIS:NE2	1:J:107:HIS:NE2	2.56	0.48
1:J:241:ASP:OD2	2:J:301:GOL:O1	2.28	0.48
1:A:90:GLU:OE2	1:A:128:LYS:HE2	2.14	0.47
1:I:68:ILE:O	1:I:71:THR:HG22	2.14	0.47
1:J:211:ARG:HD3	3:J:416:HOH:O	2.13	0.47
1:D:60:LEU:O	1:D:99:GLN:NE2	2.48	0.47
1:H:3:LYS:HG2	1:H:29:PRO:HA	1.95	0.47
1:C:145:TRP:O	1:C:174:TRP:HA	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:5:ARG:HG2	1:E:204:TYR:CE1	2.49	0.47
1:F:142:GLY:HA2	1:F:170:ILE:O	2.14	0.47
1:A:227:LYS:HE2	3:A:490:HOH:O	2.14	0.47
1:B:107:HIS:NE2	1:F:0:HIS:NE2	2.60	0.47
1:A:154:ILE:O	1:A:158:GLU:HG3	2.15	0.47
1:H:101:LEU:HD11	1:H:136:ASN:HB3	1.96	0.47
1:D:101:LEU:HD11	1:D:136:ASN:CB	2.45	0.46
1:F:35:ASN:C	1:F:35:ASN:OD1	2.57	0.46
1:G:245:HIS:HE1	2:G:301:GOL:O1	1.99	0.46
1:E:107:HIS:CE1	1:J:0:HIS:ND1	2.83	0.46
1:E:179:ARG:O	1:E:221:GLU:HG3	2.16	0.46
1:H:59:ASN:HD22	1:H:64:GLU:CG	2.29	0.46
1:H:215:ASP:HB2	3:H:642:HOH:O	2.16	0.46
1:E:63:GLY:HA3	1:F:1:MET:HE2	1.98	0.46
1:H:189:VAL:HG22	1:H:244:MET:HB3	1.98	0.46
1:D:65:TRP:CZ2	1:D:99:GLN:HG2	2.52	0.45
1:F:211:ARG:HD2	1:F:227:LYS:O	2.16	0.45
1:H:96:TRP:HA	1:H:100:VAL:HB	1.99	0.45
1:H:74:GLU:HB3	1:I:74:GLU:HG2	1.98	0.45
1:J:187:LEU:C	1:J:187:LEU:HD23	2.41	0.45
1:G:211:ARG:HD2	1:G:227:LYS:O	2.15	0.45
1:D:207:PHE:CZ	1:D:210:MET:HE1	2.51	0.45
1:H:211:ARG:HD2	1:H:227:LYS:O	2.17	0.45
1:A:101:LEU:HD23	1:A:101:LEU:HA	1.77	0.45
1:C:8:LEU:HG	1:C:205:ALA:HB1	1.98	0.44
2:G:302:GOL:C3	1:I:1:MET:HE2	2.43	0.44
1:H:66:GLN:O	1:H:70:LYS:HG3	2.17	0.44
1:B:101:LEU:HD11	1:B:136:ASN:CB	2.47	0.44
1:C:81:GLU:N	1:C:82:PRO:CD	2.80	0.44
1:E:63:GLY:HA3	1:F:1:MET:HE3	1.99	0.44
1:I:6:VAL:HB	1:I:205:ALA:HB2	1.99	0.44
1:D:145:TRP:O	1:D:174:TRP:HA	2.18	0.44
1:H:110:ARG:NH2	1:I:108:HIS:O	2.48	0.44
1:A:45:LYS:CE	3:A:413:HOH:O	2.65	0.44
1:I:82:PRO:HA	1:I:85:ALA:HB3	2.00	0.44
2:C:302:GOL:O2	2:I:302:GOL:H12	2.18	0.43
1:H:209:CYS:SG	1:H:235:MET:CE	3.06	0.43
1:C:117:ALA:O	1:C:122:GLY:HA3	2.18	0.43
1:J:179:ARG:O	1:J:221:GLU:HG3	2.18	0.43
1:E:8:LEU:HG	1:E:205:ALA:HB1	2.00	0.43
1:E:72:ASN:OD1	1:J:31:HIS:HB2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:149:LYS:CE	3:F:498:HOH:O	2.63	0.43
1:B:181:TYR:HB3	1:B:182:PRO:HD3	2.01	0.43
1:E:152:GLU:OE1	1:E:155:ARG:NH2	2.44	0.43
1:I:37:ASN:OD1	2:I:302:GOL:C3	2.65	0.43
3:B:500:HOH:O	1:E:66:GLN:HG2	2.19	0.43
1:D:187:LEU:C	1:D:187:LEU:HD23	2.43	0.43
1:C:31:HIS:NE2	1:C:203:GLU:OE1	2.50	0.43
1:D:167:GLN:NE2	3:D:510:HOH:O	2.49	0.43
1:I:37:ASN:HD21	2:I:302:GOL:H11	1.83	0.43
1:C:139:ASP:O	1:C:168:PRO:HD2	2.19	0.43
1:E:187:LEU:C	1:E:187:LEU:HD23	2.43	0.43
1:F:8:LEU:HG	1:F:205:ALA:HB1	2.00	0.43
1:I:126:ILE:HD12	1:I:126:ILE:HA	1.82	0.43
1:D:31:HIS:O	1:D:49:PRO:HD2	2.19	0.42
1:C:107:HIS:CE1	1:D:0:HIS:CE1	3.05	0.42
1:J:79:PHE:O	1:J:114:PRO:HA	2.19	0.42
1:H:107:HIS:HE1	1:I:0:HIS:CD2	2.34	0.42
1:H:110:ARG:HG3	1:I:110:ARG:CZ	2.49	0.42
1:I:66:GLN:HB3	1:I:70:LYS:HE3	2.01	0.42
1:I:86:GLY:HA2	3:I:491:HOH:O	2.17	0.42
1:I:89:PRO:HG2	1:I:128:LYS:HD3	2.00	0.42
1:A:31:HIS:NE2	1:A:203:GLU:OE1	2.49	0.42
1:G:5:ARG:HG2	1:G:204:TYR:CE1	2.54	0.42
1:G:140:TYR:HB3	1:G:170:ILE:HG13	2.02	0.42
1:H:106:SER:OG	1:H:107:HIS:CD2	2.71	0.42
1:I:5:ARG:HG2	1:I:204:TYR:CE1	2.53	0.42
1:J:18:LYS:HG3	3:J:502:HOH:O	2.20	0.42
1:B:135:LYS:HB2	3:B:485:HOH:O	2.19	0.42
1:I:29:PRO:HG3	1:I:242:GLN:HB2	2.00	0.42
1:H:104:ARG:O	1:H:108:HIS:HA	2.20	0.42
1:J:78:PHE:HD2	1:J:114:PRO:HD3	1.85	0.42
1:B:145:TRP:O	1:B:174:TRP:HA	2.20	0.41
1:F:181:TYR:N	1:F:182:PRO:CD	2.84	0.41
1:C:-2:GLY:HA2	1:I:62:GLY:HA2	2.02	0.41
1:D:31:HIS:NE2	1:D:203:GLU:OE1	2.50	0.41
1:D:189:VAL:HG22	1:D:244:MET:HB3	2.02	0.41
1:C:242:GLN:HA	1:C:243:PRO:C	2.44	0.41
1:G:96:TRP:HA	1:G:100:VAL:HB	2.02	0.41
1:B:96:TRP:HA	1:B:100:VAL:HB	2.03	0.41
1:B:126:ILE:HD12	1:B:126:ILE:HA	1.88	0.41
1:A:1:MET:HE2	2:H:501:GOL:C1	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:ILE:HD12	1:A:126:ILE:HA	1.87	0.41
1:J:10:ASP:OD2	1:J:12:THR:HG23	2.21	0.41
1:A:7:LEU:O	1:A:33:CYS:HA	2.21	0.40
1:J:237:LYS:HG2	2:J:301:GOL:O1	2.21	0.40
1:C:65:TRP:CZ2	1:C:99:GLN:HG2	2.56	0.40
1:F:5:ARG:HG2	1:F:204:TYR:CE1	2.56	0.40
1:D:101:LEU:HD11	1:D:136:ASN:CG	2.47	0.40
1:H:7:LEU:O	1:H:33:CYS:HA	2.21	0.40
1:A:213:MET:HE3	1:A:218:VAL:HG23	2.02	0.40
1:H:46:HIS:HA	3:H:604:HOH:O	2.21	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	247/266 (93%)	235 (95%)	12 (5%)	0	100	100
1	B	248/266 (93%)	241 (97%)	7 (3%)	0	100	100
1	C	247/266 (93%)	238 (96%)	9 (4%)	0	100	100
1	D	247/266 (93%)	236 (96%)	11 (4%)	0	100	100
1	E	246/266 (92%)	233 (95%)	13 (5%)	0	100	100
1	F	246/266 (92%)	236 (96%)	10 (4%)	0	100	100
1	G	247/266 (93%)	237 (96%)	10 (4%)	0	100	100
1	H	246/266 (92%)	238 (97%)	8 (3%)	0	100	100
1	I	247/266 (93%)	237 (96%)	10 (4%)	0	100	100
1	J	247/266 (93%)	237 (96%)	10 (4%)	0	100	100
All	All	2468/2660 (93%)	2368 (96%)	100 (4%)	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	221/235 (94%)	216 (98%)	5 (2%)	44	40
1	B	222/235 (94%)	220 (99%)	2 (1%)	70	73
1	C	221/235 (94%)	218 (99%)	3 (1%)	59	59
1	D	221/235 (94%)	213 (96%)	8 (4%)	31	23
1	E	221/235 (94%)	217 (98%)	4 (2%)	51	50
1	F	221/235 (94%)	218 (99%)	3 (1%)	59	59
1	G	221/235 (94%)	214 (97%)	7 (3%)	34	27
1	H	221/235 (94%)	214 (97%)	7 (3%)	34	27
1	I	221/235 (94%)	219 (99%)	2 (1%)	70	73
1	J	221/235 (94%)	218 (99%)	3 (1%)	59	59
All	All	2211/2350 (94%)	2167 (98%)	44 (2%)	48	46

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	HIS
1	A	88	SER
1	A	99	GLN
1	A	167	GLN
1	A	197	SER
1	B	152	GLU
1	B	180	SER
1	C	31	HIS
1	C	78	PHE
1	C	246	ILE
1	D	-1	SER
1	D	31	HIS
1	D	68	ILE
1	D	84	ARG

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Mol	Chain	Res	Type
1	D	91	GLU
1	D	97	ASN
1	D	118	SER
1	D	149	LYS
1	E	0	HIS
1	E	31	HIS
1	E	70	LYS
1	E	246	ILE
1	F	0	HIS
1	F	135	LYS
1	F	246	ILE
1	G	18	LYS
1	G	31	HIS
1	G	70	LYS
1	G	74	GLU
1	G	76	ILE
1	G	152	GLU
1	G	246	ILE
1	H	0	HIS
1	H	31	HIS
1	H	61	THR
1	H	90	GLU
1	H	94	LYS
1	H	99	GLN
1	H	136	ASN
1	I	31	HIS
1	I	57	LYS
1	J	0	HIS
1	J	31	HIS
1	J	120	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	0	HIS
1	A	46	HIS
1	A	166	HIS
1	B	0	HIS
1	B	46	HIS
1	B	66	GLN
1	B	212	GLN
1	C	37	ASN

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Mol	Chain	Res	Type
1	C	55	HIS
1	C	66	GLN
1	C	97	ASN
1	C	167	GLN
1	D	0	HIS
1	D	46	HIS
1	D	58	ASN
1	E	0	HIS
1	E	46	HIS
1	F	0	HIS
1	F	46	HIS
1	F	166	HIS
1	G	0	HIS
1	G	37	ASN
1	G	223	GLN
1	H	0	HIS
1	H	107	HIS
1	H	136	ASN
1	I	46	HIS
1	J	0	HIS
1	J	66	GLN
1	J	161	HIS

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](#)

33 ligands 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	$\# Z > 2$
2	GOL	E	301	-	5,5,5	0.15	0	5,5,5	0.56	0
2	GOL	G	304	-	5,5,5	0.18	0	5,5,5	0.28	0
2	GOL	D	403	-	5,5,5	0.11	0	5,5,5	0.32	0
2	GOL	C	302	-	5,5,5	0.45	0	5,5,5	0.70	0
2	GOL	J	303	-	5,5,5	0.08	0	5,5,5	0.51	0
2	GOL	G	303	-	5,5,5	0.20	0	5,5,5	0.61	0
2	GOL	C	301	-	5,5,5	0.19	0	5,5,5	0.39	0
2	GOL	I	304	-	5,5,5	0.39	0	5,5,5	0.68	0
2	GOL	F	302	-	5,5,5	0.15	0	5,5,5	0.35	0
2	GOL	A	301	-	5,5,5	0.12	0	5,5,5	0.39	0
2	GOL	G	301	-	5,5,5	0.36	0	5,5,5	0.46	0
2	GOL	D	402	-	5,5,5	0.24	0	5,5,5	0.31	0
2	GOL	G	302	-	5,5,5	0.38	0	5,5,5	1.07	0
2	GOL	J	304	-	5,5,5	0.27	0	5,5,5	0.51	0
2	GOL	A	302	-	5,5,5	0.31	0	5,5,5	0.32	0
2	GOL	B	301	-	5,5,5	0.17	0	5,5,5	0.51	0
2	GOL	B	302	-	5,5,5	0.24	0	5,5,5	0.54	0
2	GOL	F	304	-	5,5,5	0.19	0	5,5,5	0.37	0
2	GOL	J	301	-	5,5,5	0.11	0	5,5,5	0.48	0
2	GOL	D	401	-	5,5,5	0.35	0	5,5,5	1.24	1 (20%)
2	GOL	G	307	-	5,5,5	0.20	0	5,5,5	0.52	0
2	GOL	F	301	-	5,5,5	0.24	0	5,5,5	0.62	0
2	GOL	A	303	-	5,5,5	0.14	0	5,5,5	0.25	0
2	GOL	G	305	-	5,5,5	0.13	0	5,5,5	0.43	0
2	GOL	F	303	-	5,5,5	0.27	0	5,5,5	0.44	0
2	GOL	J	302	-	5,5,5	0.13	0	5,5,5	0.31	0
2	GOL	B	303	-	5,5,5	0.16	0	5,5,5	0.42	0
2	GOL	I	301	-	5,5,5	0.51	0	5,5,5	0.99	0
2	GOL	H	501	-	5,5,5	0.26	0	5,5,5	0.42	0
2	GOL	I	302	-	5,5,5	0.29	0	5,5,5	0.64	0
2	GOL	H	502	-	5,5,5	0.22	0	5,5,5	0.39	0
2	GOL	G	306	-	5,5,5	0.14	0	5,5,5	0.38	0
2	GOL	I	303	-	5,5,5	0.28	0	5,5,5	0.85	0

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
2	GOL	E	301	-	-	3/4/4/4	-
2	GOL	G	304	-	-	2/4/4/4	-
2	GOL	D	403	-	-	4/4/4/4	-
2	GOL	C	302	-	-	0/4/4/4	-
2	GOL	J	303	-	-	2/4/4/4	-
2	GOL	G	303	-	-	3/4/4/4	-
2	GOL	C	301	-	-	4/4/4/4	-
2	GOL	I	304	-	-	0/4/4/4	-
2	GOL	F	302	-	-	2/4/4/4	-
2	GOL	A	301	-	-	0/4/4/4	-
2	GOL	G	301	-	-	1/4/4/4	-
2	GOL	D	402	-	-	0/4/4/4	-
2	GOL	G	302	-	-	4/4/4/4	-
2	GOL	J	304	-	-	4/4/4/4	-
2	GOL	A	302	-	-	0/4/4/4	-
2	GOL	B	301	-	-	0/4/4/4	-
2	GOL	B	302	-	-	2/4/4/4	-
2	GOL	F	304	-	-	0/4/4/4	-
2	GOL	J	301	-	-	4/4/4/4	-
2	GOL	D	401	-	-	4/4/4/4	-
2	GOL	G	307	-	-	2/4/4/4	-
2	GOL	F	301	-	-	0/4/4/4	-
2	GOL	A	303	-	-	2/4/4/4	-
2	GOL	G	305	-	-	0/4/4/4	-
2	GOL	F	303	-	-	0/4/4/4	-
2	GOL	J	302	-	-	2/4/4/4	-
2	GOL	B	303	-	-	2/4/4/4	-
2	GOL	I	301	-	-	2/4/4/4	-
2	GOL	H	501	-	-	0/4/4/4	-
2	GOL	I	302	-	-	4/4/4/4	-
2	GOL	H	502	-	-	2/4/4/4	-
2	GOL	G	306	-	-	2/4/4/4	-
2	GOL	I	303	-	-	2/4/4/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	D	401	GOL	O2-C2-C1	2.30	118.69	109.18

There are no chirality outliers.

All (59) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	302	GOL	C1-C2-C3-O3
2	B	302	GOL	O2-C2-C3-O3
2	B	303	GOL	C1-C2-C3-O3
2	D	401	GOL	O1-C1-C2-C3
2	F	302	GOL	O1-C1-C2-C3
2	G	302	GOL	C1-C2-C3-O3
2	G	303	GOL	O1-C1-C2-C3
2	G	304	GOL	O1-C1-C2-C3
2	G	307	GOL	O1-C1-C2-C3
2	H	502	GOL	C1-C2-C3-O3
2	H	502	GOL	O2-C2-C3-O3
2	I	302	GOL	O1-C1-C2-C3
2	I	302	GOL	C1-C2-C3-O3
2	J	301	GOL	C1-C2-C3-O3
2	J	302	GOL	C1-C2-C3-O3
2	J	302	GOL	O2-C2-C3-O3
2	J	303	GOL	O1-C1-C2-C3
2	J	304	GOL	O1-C1-C2-C3
2	A	303	GOL	O2-C2-C3-O3
2	G	304	GOL	O1-C1-C2-O2
2	J	301	GOL	O2-C2-C3-O3
2	J	304	GOL	O1-C1-C2-O2
2	A	303	GOL	C1-C2-C3-O3
2	C	301	GOL	O1-C1-C2-C3
2	C	301	GOL	C1-C2-C3-O3
2	D	401	GOL	C1-C2-C3-O3
2	D	403	GOL	O1-C1-C2-C3
2	D	403	GOL	C1-C2-C3-O3
2	E	301	GOL	O1-C1-C2-C3
2	G	302	GOL	O1-C1-C2-C3
2	G	306	GOL	O1-C1-C2-C3
2	J	301	GOL	O1-C1-C2-C3
2	J	304	GOL	C1-C2-C3-O3
2	B	303	GOL	O2-C2-C3-O3
2	C	301	GOL	O1-C1-C2-O2
2	D	401	GOL	O2-C2-C3-O3
2	E	301	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
2	F	302	GOL	O1-C1-C2-O2
2	I	302	GOL	O1-C1-C2-O2
2	I	302	GOL	O2-C2-C3-O3
2	J	303	GOL	O1-C1-C2-O2
2	J	304	GOL	O2-C2-C3-O3
2	D	401	GOL	O1-C1-C2-O2
2	G	302	GOL	O2-C2-C3-O3
2	G	303	GOL	O1-C1-C2-O2
2	G	307	GOL	O1-C1-C2-O2
2	C	301	GOL	O2-C2-C3-O3
2	J	301	GOL	O1-C1-C2-O2
2	E	301	GOL	O2-C2-C3-O3
2	I	303	GOL	O2-C2-C3-O3
2	G	303	GOL	C1-C2-C3-O3
2	D	403	GOL	O2-C2-C3-O3
2	I	301	GOL	O1-C1-C2-C3
2	G	301	GOL	O1-C1-C2-C3
2	I	303	GOL	C1-C2-C3-O3
2	D	403	GOL	O1-C1-C2-O2
2	G	306	GOL	O1-C1-C2-O2
2	G	302	GOL	O1-C1-C2-O2
2	I	301	GOL	O1-C1-C2-O2

There are no ring outliers.

9 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	302	GOL	2	0
2	G	303	GOL	2	0
2	C	301	GOL	1	0
2	G	301	GOL	3	0
2	D	402	GOL	1	0
2	G	302	GOL	4	0
2	J	301	GOL	2	0
2	H	501	GOL	2	0
2	I	302	GOL	4	0

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	249/266 (93%)	0.33	6 (2%) 59 63	26, 37, 54, 67	0
1	B	249/266 (93%)	0.03	2 (0%) 82 85	20, 30, 42, 58	1 (0%)
1	C	249/266 (93%)	0.40	7 (2%) 55 59	27, 37, 58, 67	0
1	D	249/266 (93%)	0.46	8 (3%) 50 54	28, 38, 58, 71	0
1	E	248/266 (93%)	0.25	3 (1%) 76 79	28, 35, 47, 66	0
1	F	248/266 (93%)	0.18	3 (1%) 76 79	25, 33, 44, 59	0
1	G	249/266 (93%)	0.00	2 (0%) 82 85	23, 31, 45, 51	0
1	H	248/266 (93%)	0.86	26 (10%) 11 12	29, 41, 57, 64	0
1	I	249/266 (93%)	0.52	26 (10%) 11 12	24, 35, 57, 73	0
1	J	249/266 (93%)	0.09	4 (1%) 70 74	24, 31, 47, 70	0
All	All	2487/2660 (93%)	0.31	87 (3%) 47 50	20, 35, 55, 73	1 (0%)

All (87) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	93	ALA	4.6
1	C	0	HIS	4.4
1	C	217	PHE	4.2
1	H	0	HIS	4.0
1	E	0	HIS	4.0
1	D	0	HIS	3.8
1	F	-1	SER	3.7
1	F	0	HIS	3.6
1	J	0	HIS	3.6
1	H	92	ALA	3.5
1	I	60	LEU	3.5
1	G	0	HIS	3.4
1	I	217	PHE	3.4

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Mol	Chain	Res	Type	RSRZ
1	E	-1	SER	3.4
1	J	-2	GLY	3.4
1	E	217	PHE	3.3
1	I	89	PRO	3.2
1	J	-1	SER	3.2
1	D	217	PHE	3.2
1	I	100	VAL	3.1
1	C	-2	GLY	3.1
1	H	194	TRP	3.1
1	A	-2	GLY	3.0
1	H	107	HIS	3.0
1	I	166	HIS	3.0
1	C	-1	SER	2.9
1	H	-1	SER	2.9
1	H	132	LEU	2.9
1	H	217	PHE	2.9
1	D	132	LEU	2.9
1	I	94	LYS	2.9
1	C	132	LEU	2.8
1	D	121	ALA	2.8
1	H	157	LEU	2.8
1	I	-2	GLY	2.7
1	A	0	HIS	2.7
1	H	93	ALA	2.7
1	I	133	VAL	2.7
1	I	96	TRP	2.6
1	I	132	LEU	2.6
1	H	133	VAL	2.6
1	F	217	PHE	2.5
1	I	87	ILE	2.5
1	I	95	ILE	2.5
1	J	217	PHE	2.5
1	H	87	ILE	2.5
1	A	217	PHE	2.4
1	H	95	ILE	2.4
1	I	57	LYS	2.4
1	I	121	ALA	2.4
1	I	79	PHE	2.4
1	I	126	ILE	2.4
1	C	58	ASN	2.4
1	I	92	ALA	2.4
1	I	134	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	H	129	TRP	2.4
1	G	217	PHE	2.3
1	D	95	ILE	2.3
1	H	60	LEU	2.3
1	H	145	TRP	2.3
1	I	131	ASN	2.3
1	B	-2	GLY	2.3
1	H	124	ALA	2.2
1	H	200	TRP	2.2
1	I	90	GLU	2.2
1	H	165	PRO	2.2
1	H	121	ALA	2.2
1	H	185	LEU	2.2
1	A	58	ASN	2.1
1	D	246	ILE	2.1
1	H	123	ILE	2.1
1	H	126	ILE	2.1
1	I	129	TRP	2.1
1	H	141	LEU	2.1
1	I	136	ASN	2.1
1	C	60	LEU	2.1
1	H	199	PRO	2.1
1	I	65	TRP	2.1
1	I	101	LEU	2.1
1	D	-1	SER	2.0
1	H	164	HIS	2.0
1	H	96	TRP	2.0
1	D	98	ASP	2.0
1	A	-1	SER	2.0
1	A	87	ILE	2.0
1	B	217[A]	PHE	2.0
1	I	78	PHE	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

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	GOL	I	304	6/6	0.76	0.22	46,48,49,49	0
2	GOL	D	403	6/6	0.78	0.15	60,62,66,66	0
2	GOL	A	303	6/6	0.78	0.17	60,66,69,69	0
2	GOL	J	303	6/6	0.79	0.18	54,54,55,56	0
2	GOL	F	304	6/6	0.82	0.20	61,62,62,63	0
2	GOL	G	304	6/6	0.83	0.15	54,56,58,60	0
2	GOL	I	302	6/6	0.85	0.14	52,54,55,55	0
2	GOL	I	303	6/6	0.85	0.12	42,45,45,48	0
2	GOL	G	305	6/6	0.85	0.16	55,56,56,56	0
2	GOL	I	301	6/6	0.85	0.13	37,42,42,44	0
2	GOL	G	303	6/6	0.86	0.14	43,44,45,50	0
2	GOL	C	301	6/6	0.86	0.13	56,57,58,62	0
2	GOL	B	302	6/6	0.87	0.12	39,44,45,45	0
2	GOL	G	301	6/6	0.88	0.11	30,32,34,35	0
2	GOL	H	502	6/6	0.89	0.10	39,41,43,48	0
2	GOL	E	301	6/6	0.90	0.11	40,41,42,43	0
2	GOL	G	306	6/6	0.90	0.11	40,47,49,51	0
2	GOL	J	302	6/6	0.90	0.12	50,52,55,59	0
2	GOL	H	501	6/6	0.90	0.12	46,48,52,54	0
2	GOL	G	307	6/6	0.91	0.13	50,54,55,56	0
2	GOL	B	303	6/6	0.92	0.10	45,46,48,48	0
2	GOL	G	302	6/6	0.92	0.09	30,31,31,31	0
2	GOL	C	302	6/6	0.92	0.09	40,40,41,41	0
2	GOL	F	303	6/6	0.92	0.10	31,34,36,37	0
2	GOL	D	402	6/6	0.92	0.10	43,46,48,49	0
2	GOL	J	304	6/6	0.92	0.10	41,46,48,52	0
2	GOL	J	301	6/6	0.93	0.09	38,40,43,50	0
2	GOL	D	401	6/6	0.93	0.09	32,35,39,42	0
2	GOL	A	301	6/6	0.94	0.08	31,32,32,32	0
2	GOL	F	301	6/6	0.94	0.09	41,45,47,49	0
2	GOL	A	302	6/6	0.94	0.07	30,33,34,36	0
2	GOL	B	301	6/6	0.95	0.07	31,33,34,34	0
2	GOL	F	302	6/6	0.96	0.08	41,47,47,48	0

6.5 Other polymers [i](#)

There are no such residues in this entry.