



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 08:29 PM UTC

PDB ID : 4UDG / pdb_00004udg
Title : Crystal structure of b-1,4-mannopyranosyl-chitobiose phosphorylase at 1.60 Angstrom in complex with N-acetylglucosamine and inorganic phosphate
Authors : Ladeveze, S.; Cioci, G.; Potocki-Veronese, G.; Tranier, S.; Mourey, L.
Deposited on : 2014-12-10
Resolution : 1.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
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

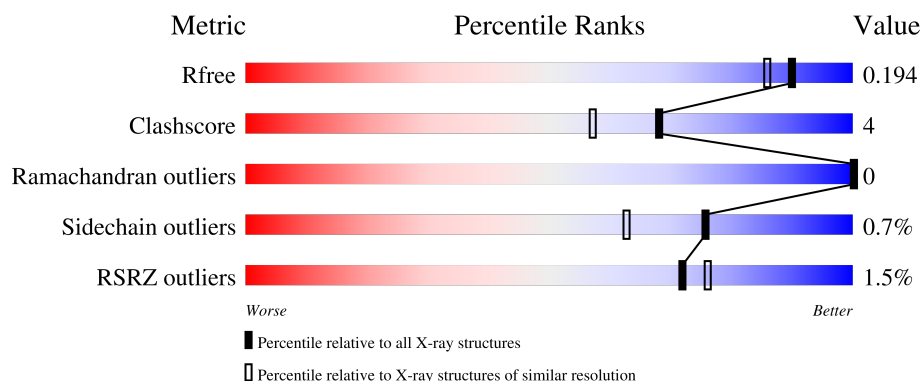
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4673 (1.60-1.60)
Clashscore	190562	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)
RSRZ outliers	180081	4672 (1.60-1.60)

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	347	
1	B	347	
1	C	347	
1	D	347	
1	E	347	

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Mol	Chain	Length	Quality of chain
1	F	347	 2% 83% 9% 8%

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
4	GOL	A	330	-	X	-	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 17360 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 UHGB_MP.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	320	Total	C	N	O	S	0	9	0
			2623	1699	437	470	17			
1	B	321	Total	C	N	O	S	0	7	0
			2633	1700	443	473	17			
1	C	320	Total	C	N	O	S	0	4	0
			2581	1671	431	462	17			
1	D	320	Total	C	N	O	S	0	9	0
			2640	1707	443	474	16			
1	E	321	Total	C	N	O	S	0	7	0
			2632	1699	442	474	17			
1	F	320	Total	C	N	O	S	0	4	0
			2600	1678	437	469	16			

There are 120 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP D9ZDQ9
A	-18	GLY	-	expression tag	UNP D9ZDQ9
A	-17	SER	-	expression tag	UNP D9ZDQ9
A	-16	SER	-	expression tag	UNP D9ZDQ9
A	-15	HIS	-	expression tag	UNP D9ZDQ9
A	-14	HIS	-	expression tag	UNP D9ZDQ9
A	-13	HIS	-	expression tag	UNP D9ZDQ9
A	-12	HIS	-	expression tag	UNP D9ZDQ9
A	-11	HIS	-	expression tag	UNP D9ZDQ9
A	-10	HIS	-	expression tag	UNP D9ZDQ9
A	-9	SER	-	expression tag	UNP D9ZDQ9
A	-8	SER	-	expression tag	UNP D9ZDQ9
A	-7	GLY	-	expression tag	UNP D9ZDQ9
A	-6	LEU	-	expression tag	UNP D9ZDQ9
A	-5	VAL	-	expression tag	UNP D9ZDQ9
A	-4	PRO	-	expression tag	UNP D9ZDQ9
A	-3	ARG	-	expression tag	UNP D9ZDQ9

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP D9ZDQ9
A	-1	SER	-	expression tag	UNP D9ZDQ9
A	0	HIS	-	expression tag	UNP D9ZDQ9
B	-19	MET	-	expression tag	UNP D9ZDQ9
B	-18	GLY	-	expression tag	UNP D9ZDQ9
B	-17	SER	-	expression tag	UNP D9ZDQ9
B	-16	SER	-	expression tag	UNP D9ZDQ9
B	-15	HIS	-	expression tag	UNP D9ZDQ9
B	-14	HIS	-	expression tag	UNP D9ZDQ9
B	-13	HIS	-	expression tag	UNP D9ZDQ9
B	-12	HIS	-	expression tag	UNP D9ZDQ9
B	-11	HIS	-	expression tag	UNP D9ZDQ9
B	-10	HIS	-	expression tag	UNP D9ZDQ9
B	-9	SER	-	expression tag	UNP D9ZDQ9
B	-8	SER	-	expression tag	UNP D9ZDQ9
B	-7	GLY	-	expression tag	UNP D9ZDQ9
B	-6	LEU	-	expression tag	UNP D9ZDQ9
B	-5	VAL	-	expression tag	UNP D9ZDQ9
B	-4	PRO	-	expression tag	UNP D9ZDQ9
B	-3	ARG	-	expression tag	UNP D9ZDQ9
B	-2	GLY	-	expression tag	UNP D9ZDQ9
B	-1	SER	-	expression tag	UNP D9ZDQ9
B	0	HIS	-	expression tag	UNP D9ZDQ9
C	-19	MET	-	expression tag	UNP D9ZDQ9
C	-18	GLY	-	expression tag	UNP D9ZDQ9
C	-17	SER	-	expression tag	UNP D9ZDQ9
C	-16	SER	-	expression tag	UNP D9ZDQ9
C	-15	HIS	-	expression tag	UNP D9ZDQ9
C	-14	HIS	-	expression tag	UNP D9ZDQ9
C	-13	HIS	-	expression tag	UNP D9ZDQ9
C	-12	HIS	-	expression tag	UNP D9ZDQ9
C	-11	HIS	-	expression tag	UNP D9ZDQ9
C	-10	HIS	-	expression tag	UNP D9ZDQ9
C	-9	SER	-	expression tag	UNP D9ZDQ9
C	-8	SER	-	expression tag	UNP D9ZDQ9
C	-7	GLY	-	expression tag	UNP D9ZDQ9
C	-6	LEU	-	expression tag	UNP D9ZDQ9
C	-5	VAL	-	expression tag	UNP D9ZDQ9
C	-4	PRO	-	expression tag	UNP D9ZDQ9
C	-3	ARG	-	expression tag	UNP D9ZDQ9
C	-2	GLY	-	expression tag	UNP D9ZDQ9
C	-1	SER	-	expression tag	UNP D9ZDQ9

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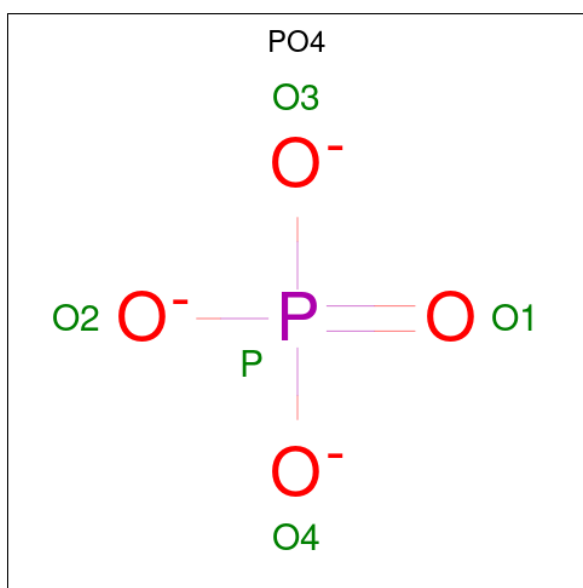
Chain	Residue	Modelled	Actual	Comment	Reference
C	0	HIS	-	expression tag	UNP D9ZDQ9
D	-19	MET	-	expression tag	UNP D9ZDQ9
D	-18	GLY	-	expression tag	UNP D9ZDQ9
D	-17	SER	-	expression tag	UNP D9ZDQ9
D	-16	SER	-	expression tag	UNP D9ZDQ9
D	-15	HIS	-	expression tag	UNP D9ZDQ9
D	-14	HIS	-	expression tag	UNP D9ZDQ9
D	-13	HIS	-	expression tag	UNP D9ZDQ9
D	-12	HIS	-	expression tag	UNP D9ZDQ9
D	-11	HIS	-	expression tag	UNP D9ZDQ9
D	-10	HIS	-	expression tag	UNP D9ZDQ9
D	-9	SER	-	expression tag	UNP D9ZDQ9
D	-8	SER	-	expression tag	UNP D9ZDQ9
D	-7	GLY	-	expression tag	UNP D9ZDQ9
D	-6	LEU	-	expression tag	UNP D9ZDQ9
D	-5	VAL	-	expression tag	UNP D9ZDQ9
D	-4	PRO	-	expression tag	UNP D9ZDQ9
D	-3	ARG	-	expression tag	UNP D9ZDQ9
D	-2	GLY	-	expression tag	UNP D9ZDQ9
D	-1	SER	-	expression tag	UNP D9ZDQ9
D	0	HIS	-	expression tag	UNP D9ZDQ9
E	-19	MET	-	expression tag	UNP D9ZDQ9
E	-18	GLY	-	expression tag	UNP D9ZDQ9
E	-17	SER	-	expression tag	UNP D9ZDQ9
E	-16	SER	-	expression tag	UNP D9ZDQ9
E	-15	HIS	-	expression tag	UNP D9ZDQ9
E	-14	HIS	-	expression tag	UNP D9ZDQ9
E	-13	HIS	-	expression tag	UNP D9ZDQ9
E	-12	HIS	-	expression tag	UNP D9ZDQ9
E	-11	HIS	-	expression tag	UNP D9ZDQ9
E	-10	HIS	-	expression tag	UNP D9ZDQ9
E	-9	SER	-	expression tag	UNP D9ZDQ9
E	-8	SER	-	expression tag	UNP D9ZDQ9
E	-7	GLY	-	expression tag	UNP D9ZDQ9
E	-6	LEU	-	expression tag	UNP D9ZDQ9
E	-5	VAL	-	expression tag	UNP D9ZDQ9
E	-4	PRO	-	expression tag	UNP D9ZDQ9
E	-3	ARG	-	expression tag	UNP D9ZDQ9
E	-2	GLY	-	expression tag	UNP D9ZDQ9
E	-1	SER	-	expression tag	UNP D9ZDQ9
E	0	HIS	-	expression tag	UNP D9ZDQ9
F	-19	MET	-	expression tag	UNP D9ZDQ9

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-18	GLY	-	expression tag	UNP D9ZDQ9
F	-17	SER	-	expression tag	UNP D9ZDQ9
F	-16	SER	-	expression tag	UNP D9ZDQ9
F	-15	HIS	-	expression tag	UNP D9ZDQ9
F	-14	HIS	-	expression tag	UNP D9ZDQ9
F	-13	HIS	-	expression tag	UNP D9ZDQ9
F	-12	HIS	-	expression tag	UNP D9ZDQ9
F	-11	HIS	-	expression tag	UNP D9ZDQ9
F	-10	HIS	-	expression tag	UNP D9ZDQ9
F	-9	SER	-	expression tag	UNP D9ZDQ9
F	-8	SER	-	expression tag	UNP D9ZDQ9
F	-7	GLY	-	expression tag	UNP D9ZDQ9
F	-6	LEU	-	expression tag	UNP D9ZDQ9
F	-5	VAL	-	expression tag	UNP D9ZDQ9
F	-4	PRO	-	expression tag	UNP D9ZDQ9
F	-3	ARG	-	expression tag	UNP D9ZDQ9
F	-2	GLY	-	expression tag	UNP D9ZDQ9
F	-1	SER	-	expression tag	UNP D9ZDQ9
F	0	HIS	-	expression tag	UNP D9ZDQ9

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



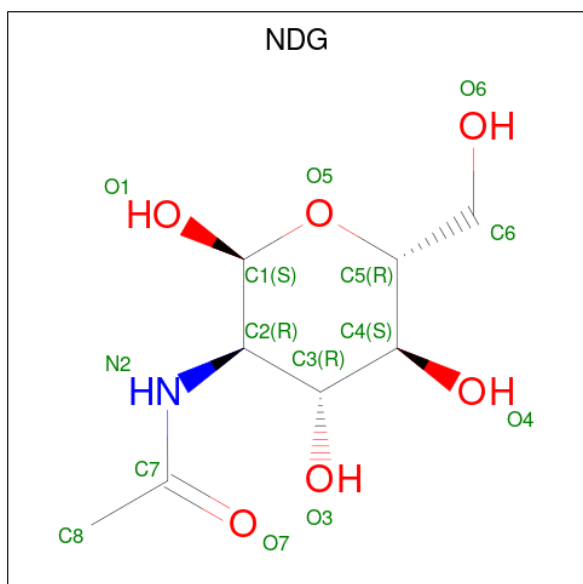
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		

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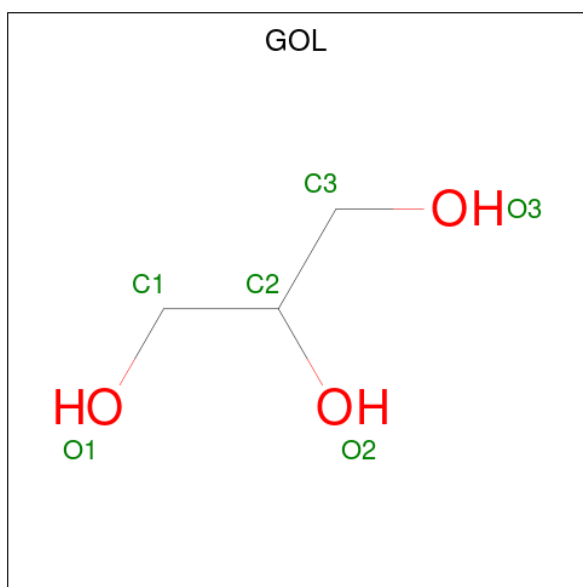
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	O	P	0	0
			5	4	1		
2	D	1	Total	O	P	0	0
			5	4	1		
2	E	1	Total	O	P	0	0
			5	4	1		
2	F	1	Total	O	P	0	0
			5	4	1		

- Molecule 3 is 2-acetamido-2-deoxy- α -D-glucopyranose (CCD ID: NDG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			15	8	1	6		
3	B	1	Total	C	N	O	0	0
			15	8	1	6		
3	C	1	Total	C	N	O	0	0
			15	8	1	6		
3	D	1	Total	C	N	O	0	0
			15	8	1	6		
3	E	1	Total	C	N	O	0	0
			15	8	1	6		
3	F	1	Total	C	N	O	0	0
			15	8	1	6		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 5 is POTASSIUM ION (CCD ID: K) (formula: K).

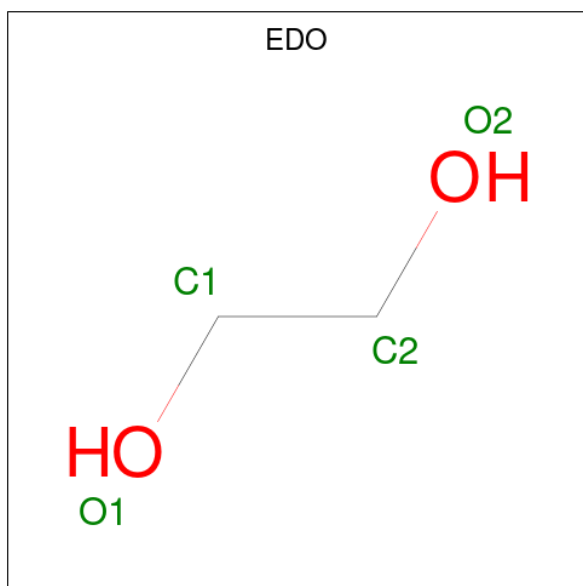
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	K	0	0
			1	1		
5	B	1	Total	K	0	0
			1	1		
5	C	1	Total	K	0	0
			1	1		
5	D	1	Total	K	0	0
			1	1		
5	E	1	Total	K	0	0
			1	1		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	F	1	Total K 1 1	0	0

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	E	1	Total C O 4 2 2	0	0
6	E	1	Total C O 4 2 2	0	0
6	F	1	Total C O 4 2 2	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	277	Total O 277 277	0	0
7	B	277	Total O 277 277	0	0
7	C	262	Total O 262 262	0	0

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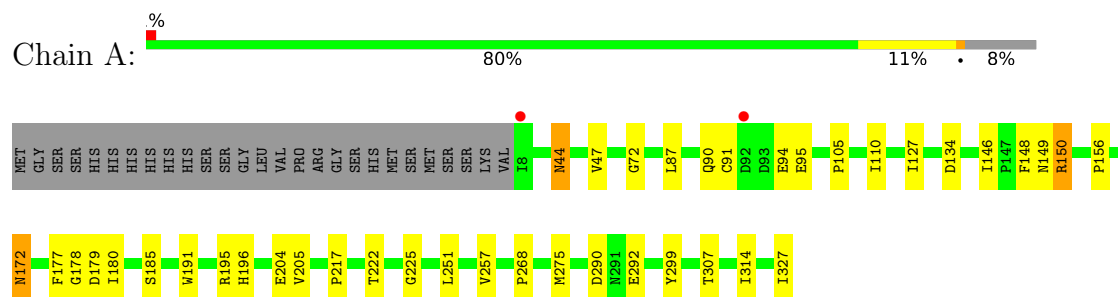
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	D	228	Total 228	O 228	0	0
7	E	239	Total 239	O 239	0	0
7	F	180	Total 180	O 180	0	0

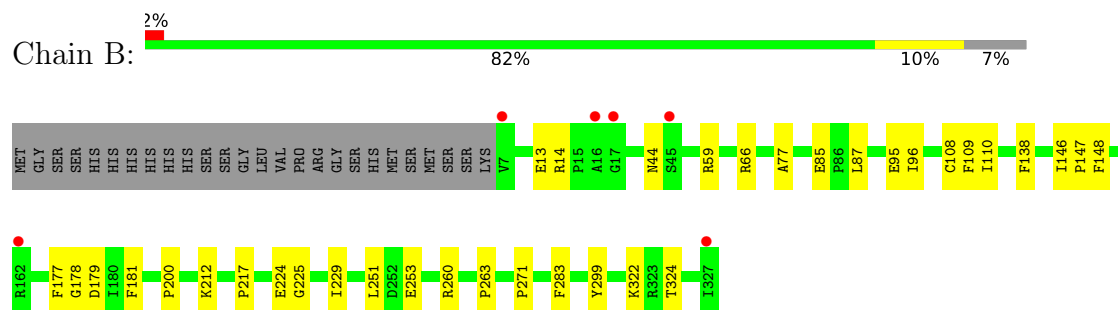
3 Residue-property plots [i](#)

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.

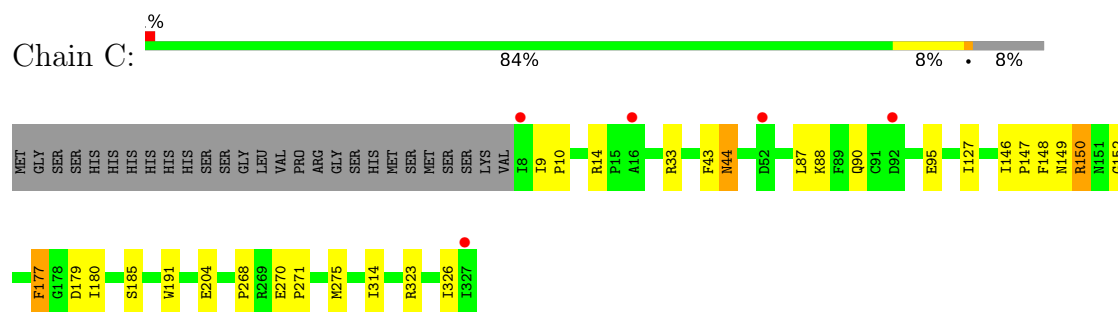
• Molecule 1: UHGB_MP



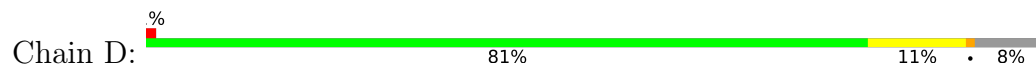
• Molecule 1: UHGB_MP

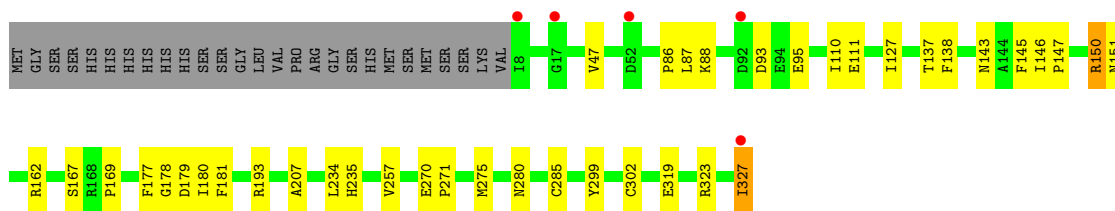


• Molecule 1: UHGB_MP

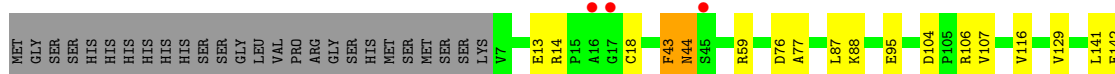
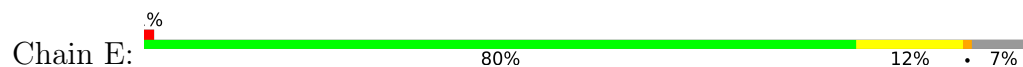


• Molecule 1: UHGB_MP

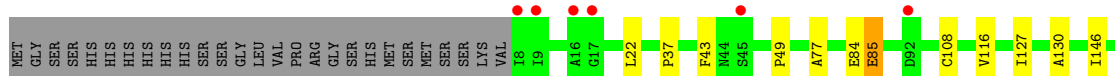
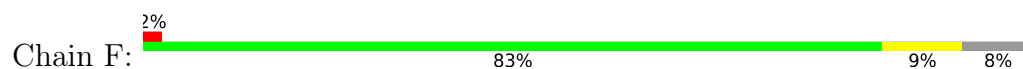




• Molecule 1: UHGB_MP



• Molecule 1: UHGB_MP



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.82Å 140.91Å 168.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	108.12 – 1.60 108.12 – 1.60	Depositor EDS
% Data completeness (in resolution range)	99.8 (108.12-1.60) 99.8 (108.12-1.60)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.154 , 0.183 0.168 , 0.194	Depositor DCC
R_{free} test set	13154 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	19.4	Xtriage
Anisotropy	0.217	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 34.2	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	0.97	EDS
Total number of atoms	17360	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.19% 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, K, PO4, EDO, NDG

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	1.36	19/2721 (0.7%)	1.05	5/3711 (0.1%)
1	B	1.29	15/2722 (0.6%)	1.05	2/3711 (0.1%)
1	C	1.35	15/2676 (0.6%)	1.07	2/3649 (0.1%)
1	D	1.23	11/2729 (0.4%)	1.01	3/3720 (0.1%)
1	E	1.26	14/2721 (0.5%)	1.06	2/3709 (0.1%)
1	F	1.19	4/2692 (0.1%)	1.04	0/3669
All	All	1.28	78/16261 (0.5%)	1.05	14/22169 (0.1%)

All (78) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	95	GLU	C-N	-16.02	1.14	1.33
1	C	326	ILE	C-N	-15.70	1.11	1.33
1	E	326	ILE	C-N	9.83	1.47	1.33
1	D	150	ARG	C-O	-9.37	1.15	1.24
1	A	222	THR	N-CA	7.86	1.51	1.46
1	E	179	ASP	C-O	-7.26	1.15	1.23
1	A	314	ILE	CA-CB	-7.17	1.50	1.54
1	A	179	ASP	C-O	-6.95	1.15	1.23
1	B	283	PHE	C-O	-6.88	1.18	1.24
1	E	225	GLY	C-O	6.84	1.29	1.23
1	A	91	CYS	C-O	-6.82	1.16	1.23
1	B	179	ASP	C-O	-6.80	1.15	1.23
1	C	150	ARG	C-O	-6.78	1.17	1.24
1	A	156	PRO	CA-C	6.64	1.58	1.52
1	B	108	CYS	C-O	-6.60	1.15	1.23
1	B	66	ARG	C-O	-6.48	1.16	1.23
1	A	150	ARG	C-O	-6.42	1.17	1.24
1	A	110	ILE	C-O	-6.34	1.17	1.24
1	B	178	GLY	C-O	-6.33	1.15	1.23
1	A	90	GLN	C-O	-6.31	1.16	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	179	ASP	C-O	-6.23	1.16	1.23
1	C	90	GLN	C-O	-6.18	1.16	1.24
1	C	268	PRO	C-O	-6.13	1.16	1.23
1	E	270	GLU	C-O	-6.13	1.15	1.23
1	A	307	THR	C-O	-6.08	1.16	1.24
1	D	179	ASP	C-O	-5.99	1.16	1.23
1	D	178	GLY	C-O	-5.92	1.16	1.23
1	A	178	GLY	C-O	-5.91	1.16	1.23
1	C	314	ILE	CA-CB	-5.89	1.51	1.54
1	C	180	ILE	C-O	-5.87	1.17	1.24
1	B	59	ARG	C-O	-5.87	1.17	1.24
1	B	271	PRO	C-O	-5.84	1.16	1.24
1	A	72	GLY	N-CA	5.82	1.51	1.45
1	E	177	PHE	C-O	-5.82	1.17	1.23
1	A	195	ARG	C-O	-5.81	1.16	1.23
1	D	271	PRO	N-CD	5.74	1.55	1.47
1	E	309	LEU	C-O	-5.72	1.17	1.23
1	B	44	ASN	N-CA	5.70	1.52	1.46
1	C	44	ASN	N-CA	5.66	1.52	1.46
1	F	49	PRO	C-O	-5.65	1.17	1.23
1	D	86	PRO	C-O	5.62	1.30	1.23
1	C	14	ARG	C-N	-5.61	1.26	1.33
1	E	271	PRO	N-CD	5.55	1.55	1.47
1	B	13	GLU	C-N	-5.52	1.24	1.33
1	E	150	ARG	C-O	-5.49	1.18	1.24
1	C	271	PRO	N-CD	5.48	1.55	1.47
1	F	179	ASP	C-O	-5.45	1.17	1.23
1	C	270	GLU	C-O	-5.45	1.16	1.23
1	B	271	PRO	N-CD	5.45	1.55	1.47
1	F	302	CYS	N-CA	5.44	1.52	1.46
1	D	302	CYS	C-O	5.43	1.30	1.24
1	E	271	PRO	C-O	-5.38	1.17	1.24
1	D	143	ASN	C-O	-5.34	1.17	1.23
1	B	110	ILE	C-O	-5.33	1.18	1.24
1	E	205	VAL	C-O	-5.26	1.18	1.23
1	C	177	PHE	C-O	-5.25	1.17	1.23
1	B	260	ARG	C-O	-5.23	1.17	1.23
1	B	181	PHE	C-O	-5.20	1.17	1.23
1	D	180	ILE	C-O	-5.18	1.17	1.24
1	F	178	GLY	C-O	-5.18	1.17	1.23
1	D	271	PRO	C-O	-5.18	1.17	1.24
1	A	134	ASP	C-O	-5.17	1.17	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	33	ARG	C-O	-5.17	1.17	1.24
1	C	323	ARG	C-O	-5.17	1.17	1.24
1	E	180	ILE	C-O	-5.14	1.18	1.24
1	C	149	ASN	C-O	-5.14	1.17	1.23
1	E	270	GLU	C-N	5.13	1.40	1.33
1	A	149	ASN	C-O	-5.13	1.17	1.23
1	D	151	ASN	C-O	-5.13	1.17	1.23
1	A	196	HIS	N-CA	-5.11	1.39	1.46
1	A	44	ASN	C-O	5.10	1.30	1.24
1	E	44	ASN	N-CA	5.10	1.52	1.46
1	B	96	ILE	C-O	-5.07	1.18	1.24
1	E	104	ASP	CA-CB	5.06	1.61	1.53
1	A	180	ILE	C-O	-5.04	1.18	1.24
1	A	172	ASN	C-O	-5.04	1.17	1.24
1	D	167	SER	C-N	-5.02	1.29	1.33
1	B	109	PHE	C-O	-5.01	1.17	1.24

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	43	PHE	CA-C-N	-7.69	111.40	122.94
1	E	43	PHE	C-N-CA	-7.69	111.40	122.94
1	A	217	PRO	CB-CA-C	-6.74	102.42	111.12
1	A	268	PRO	O-C-N	6.21	129.75	122.98
1	A	95	GLU	O-C-N	-6.10	114.30	122.23
1	D	280	ASN	N-CA-C	5.46	119.16	111.52
1	A	94	GLU	CA-C-N	-5.35	111.42	120.58
1	A	94	GLU	C-N-CA	-5.35	111.42	120.58
1	B	253	GLU	CA-C-N	-5.33	114.53	120.45
1	B	253	GLU	C-N-CA	-5.33	114.53	120.45
1	C	43	PHE	CA-C-N	-5.11	115.28	122.94
1	C	43	PHE	C-N-CA	-5.11	115.28	122.94
1	D	270	GLU	CA-C-N	-5.09	113.68	119.28
1	D	270	GLU	C-N-CA	-5.09	113.68	119.28

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2623	0	2495	21	0
1	B	2633	0	2498	14	0
1	C	2581	0	2443	16	0
1	D	2640	0	2506	31	0
1	E	2632	0	2493	25	0
1	F	2600	0	2459	19	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
2	E	5	0	0	0	0
2	F	5	0	0	0	0
3	A	15	0	12	0	0
3	B	15	0	12	0	0
3	C	15	0	12	0	0
3	D	15	0	12	0	0
3	E	15	0	12	0	0
3	F	15	0	12	0	0
4	A	6	0	8	0	0
4	B	12	0	16	0	0
4	C	6	0	8	0	0
4	D	6	0	8	0	0
4	E	6	0	8	0	0
4	F	6	0	8	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
5	E	1	0	0	0	0
5	F	1	0	0	0	0
6	A	8	0	12	0	0
6	E	8	0	12	0	0
6	F	4	0	6	0	0
7	A	277	0	0	2	0
7	B	277	0	0	1	0
7	C	262	0	0	2	0
7	D	228	0	0	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	E	239	0	0	2	0
7	F	180	0	0	0	0
All	All	17360	0	15052	111	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 4.

All (111) 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:E:107:VAL:HG22	1:E:116[A]:VAL:HG12	1.45	0.94
1:B:87:LEU:HD11	1:B:138:PHE:CD1	2.12	0.85
1:D:285:CYS:HB3	7:D:587:HOH:O	1.78	0.83
1:E:129[A]:VAL:HG13	1:E:141:LEU:HB2	1.62	0.82
1:C:44:ASN:OD1	7:C:419:HOH:O	1.99	0.79
1:E:44:ASN:OD1	7:E:419:HOH:O	2.04	0.75
1:D:234[B]:LEU:HD12	1:D:235:HIS:N	2.03	0.73
1:A:146[A]:ILE:HD11	1:E:181:PHE:CB	2.19	0.72
1:D:257[A]:VAL:HG12	7:D:536:HOH:O	1.90	0.70
1:D:137[A]:THR:HG23	7:D:598:HOH:O	1.94	0.68
1:D:234[B]:LEU:CD1	7:D:507:HOH:O	2.40	0.67
1:A:146[A]:ILE:HG22	7:A:625:HOH:O	1.95	0.64
1:F:225:GLY:HA2	1:F:251:LEU:HG	1.80	0.64
1:A:47[B]:VAL:HG21	1:A:299:TYR:CE1	2.35	0.61
1:E:257:VAL:HG12	1:E:327:ILE:HD11	1.83	0.61
1:F:22:LEU:HD22	1:F:265:LEU:HB3	1.83	0.60
1:D:47[A]:VAL:HG11	1:D:299:TYR:CE1	2.36	0.60
1:D:275:MET:CE	1:F:275:MET:HE3	2.30	0.60
1:D:234[B]:LEU:HD12	7:D:507:HOH:O	2.01	0.59
1:A:275:MET:HE2	1:C:275:MET:HE2	1.84	0.58
1:D:137[B]:THR:HG23	7:D:446:HOH:O	2.03	0.58
1:F:187:ASP:O	1:F:188:MET:HB2	2.04	0.57
1:A:275:MET:HE1	1:C:275:MET:CE	2.34	0.57
1:A:146[A]:ILE:HD11	1:E:181:PHE:HB2	1.86	0.57
1:F:248:LEU:HD11	1:F:318:ILE:HG23	1.87	0.56
1:D:275:MET:HE1	1:F:275:MET:HE3	1.87	0.55
1:C:127[B]:ILE:HD13	1:C:152:GLY:HA3	1.89	0.55
1:A:275:MET:CE	1:C:275:MET:CE	2.85	0.54
1:D:234[A]:LEU:HD22	7:D:507:HOH:O	2.07	0.54
1:E:275:MET:HB2	1:F:275:MET:HE2	1.90	0.54
1:F:257:VAL:HG12	1:F:327:ILE:HD11	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:127[A]:ILE:HD11	1:F:150:ARG:HA	1.89	0.53
1:E:129[A]:VAL:HG12	1:E:142:GLU:O	2.09	0.52
1:C:204:GLU:OE2	7:C:614:HOH:O	2.19	0.52
1:D:162:ARG:NH2	1:D:193[A]:ARG:NH1	2.58	0.52
1:F:127[A]:ILE:CD1	1:F:150:ARG:HA	2.40	0.51
1:B:146[A]:ILE:CG2	1:D:169:PRO:HG3	2.41	0.51
1:D:127[A]:ILE:CD1	1:D:150:ARG:HA	2.40	0.51
1:D:110:ILE:HG22	1:D:111:GLU:HG3	1.93	0.51
1:E:43:PHE:CZ	1:E:59:ARG:CZ	2.94	0.51
1:B:224:GLU:OE1	1:B:322:LYS:NZ	2.39	0.50
1:C:146:ILE:CG2	1:C:147:PRO:HD2	2.42	0.50
1:B:146[B]:ILE:HG22	7:B:595:HOH:O	2.11	0.50
1:E:220[A]:VAL:CG2	1:E:227:LEU:HB3	2.42	0.50
1:A:257:VAL:HG12	1:A:327:ILE:HD11	1.93	0.50
1:D:323:ARG:NE	7:D:669:HOH:O	2.37	0.50
1:B:225:GLY:HA2	1:B:251:LEU:HG	1.95	0.49
1:A:44:ASN:ND2	7:A:419:HOH:O	2.14	0.49
1:A:290:ASP:OD1	1:A:292:GLU:HG2	2.13	0.49
1:C:146:ILE:HG22	1:C:147:PRO:HD2	1.94	0.49
1:E:106:ARG:NH1	1:E:285[B]:CYS:SG	2.86	0.49
1:F:108:CYS:HB3	1:F:218:ILE:HG13	1.94	0.48
1:E:87:LEU:HD23	1:E:88:LYS:N	2.29	0.47
1:A:225:GLY:HA2	1:A:251[A]:LEU:HG	1.98	0.46
1:D:93:ASP:OD1	1:D:95:GLU:HB3	2.15	0.46
1:D:137[B]:THR:HG23	1:D:137[B]:THR:O	2.15	0.46
1:D:127[A]:ILE:HD11	1:D:150:ARG:HA	1.98	0.46
1:E:234:LEU:C	1:E:234:LEU:HD23	2.40	0.46
1:D:87[B]:LEU:HD11	1:D:138:PHE:CD1	2.50	0.46
1:C:87:LEU:HD23	1:C:87:LEU:C	2.41	0.45
1:C:127[A]:ILE:HD11	1:C:150:ARG:HA	1.98	0.45
1:E:87:LEU:HD23	1:E:87:LEU:C	2.42	0.45
1:F:85[A]:GLU:H	1:F:85[A]:GLU:CD	2.23	0.45
1:D:127[B]:ILE:HD12	1:D:145:PHE:CD2	2.52	0.45
1:A:275:MET:CE	1:C:275:MET:HE2	2.47	0.45
1:C:148:PHE:HE2	1:F:177:PHE:CE1	2.35	0.44
1:A:146[A]:ILE:HD11	1:E:181:PHE:HB3	1.97	0.44
1:E:116[A]:VAL:HG23	1:E:116[A]:VAL:O	2.17	0.44
1:A:148:PHE:CE1	1:A:172:ASN:HB2	2.52	0.43
1:E:187:ASP:O	1:E:188:MET:HB2	2.17	0.43
1:A:204:GLU:HG2	1:A:205:VAL:HG23	2.00	0.43
1:D:162:ARG:NH2	1:D:193[A]:ARG:HH12	2.15	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:177:PHE:C	1:B:177:PHE:CD1	2.97	0.43
1:B:217:PRO:HD2	1:B:229[B]:ILE:HB	2.01	0.43
1:A:185:SER:HB2	1:A:191:TRP:CD2	2.54	0.43
1:D:87[A]:LEU:HD23	1:D:88:LYS:N	2.34	0.43
1:A:87:LEU:C	1:A:87:LEU:HD23	2.44	0.42
1:A:127[A]:ILE:CD1	1:A:150:ARG:HA	2.49	0.42
1:D:207:ALA:HB1	1:D:234[A]:LEU:HD23	2.00	0.42
1:F:222:THR:HG21	1:F:318:ILE:HD11	2.00	0.42
1:B:146[A]:ILE:HG22	1:B:147:PRO:O	2.19	0.42
1:E:221:GLU:HB2	1:E:226:TRP:CZ3	2.55	0.42
1:F:146:ILE:HD13	1:F:146:ILE:HG21	1.85	0.42
1:B:263:PRO:HD3	1:B:324:THR:HB	2.02	0.42
1:B:200:PRO:HG3	1:B:212:LYS:HA	2.00	0.42
1:A:146[A]:ILE:HD13	1:A:146[A]:ILE:HG21	1.75	0.42
1:C:127[A]:ILE:CD1	1:C:150:ARG:HA	2.50	0.42
1:F:77:ALA:HB1	1:F:299:TYR:OH	2.20	0.41
1:B:148:PHE:HE2	1:D:177:PHE:CE1	2.39	0.41
1:C:9:ILE:HB	1:C:10:PRO:HD2	2.02	0.41
1:E:14:ARG:HG2	1:E:18:CYS:HB3	2.01	0.41
1:B:146[A]:ILE:HG21	1:D:169:PRO:HG3	2.02	0.41
1:E:43:PHE:HZ	1:E:59:ARG:CZ	2.34	0.41
1:F:37:PRO:HD2	1:F:84:GLU:HG2	2.02	0.41
1:F:116:VAL:CG1	1:F:130:ALA:HB3	2.50	0.41
1:E:216:GLY:HA3	1:E:217:PRO:HD3	1.95	0.41
1:E:220[A]:VAL:HG23	1:E:227:LEU:HB3	2.03	0.41
1:C:185:SER:HB2	1:C:191:TRP:CD2	2.56	0.41
1:D:146:ILE:CG2	1:D:147:PRO:HD2	2.51	0.41
1:D:257[B]:VAL:HG12	1:D:327:ILE:HD13	2.03	0.41
1:C:177:PHE:C	1:C:177:PHE:CD1	2.99	0.41
1:D:87[A]:LEU:HD23	1:D:87[A]:LEU:C	2.46	0.41
1:E:76:ASP:O	1:E:77:ALA:HB3	2.20	0.41
1:F:43:PHE:CD2	1:F:304:ASP:HA	2.56	0.41
1:B:77:ALA:HB1	1:B:299:TYR:OH	2.20	0.40
1:E:77:ALA:HB1	1:E:299:TYR:OH	2.21	0.40
1:A:44:ASN:HD22	1:A:105:PRO:HD2	1.87	0.40
1:B:146[B]:ILE:HD11	1:D:181:PHE:HB2	2.03	0.40
1:A:177:PHE:CD1	1:A:177:PHE:C	2.99	0.40
1:E:235:HIS:HE1	7:E:607:HOH:O	2.05	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	327/347 (94%)	311 (95%)	16 (5%)	0	100	100
1	B	326/347 (94%)	308 (94%)	18 (6%)	0	100	100
1	C	322/347 (93%)	304 (94%)	18 (6%)	0	100	100
1	D	327/347 (94%)	313 (96%)	14 (4%)	0	100	100
1	E	326/347 (94%)	312 (96%)	14 (4%)	0	100	100
1	F	322/347 (93%)	305 (95%)	17 (5%)	0	100	100
All	All	1950/2082 (94%)	1853 (95%)	97 (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	281/298 (94%)	281 (100%)	0	100	100
1	B	282/298 (95%)	279 (99%)	3 (1%)	65	46
1	C	274/298 (92%)	272 (99%)	2 (1%)	76	63
1	D	282/298 (95%)	280 (99%)	2 (1%)	76	63
1	E	282/298 (95%)	279 (99%)	3 (1%)	65	46
1	F	277/298 (93%)	275 (99%)	2 (1%)	76	63
All	All	1678/1788 (94%)	1666 (99%)	12 (1%)	76	63

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	14	ARG
1	B	85	GLU
1	B	95	GLU
1	C	88	LYS
1	C	95	GLU
1	D	319	GLU
1	D	327	ILE
1	E	13	GLU
1	E	95	GLU
1	E	309	LEU
1	F	85[A]	GLU
1	F	85[B]	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	79	ASN
1	A	280	ASN
1	B	172	ASN
1	C	44	ASN
1	C	172	ASN
1	E	44	ASN
1	E	140	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 30 ligands modelled in this entry, 6 are monoatomic - leaving 24 for Mogul analysis.

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$
4	GOL	D	330	-	5,5,5	0.45	0	5,5,5	0.56	0
3	NDG	A	329	-	15,15,15	0.86	0	21,21,21	1.25	2 (9%)
4	GOL	C	330	-	5,5,5	0.87	0	5,5,5	0.95	0
4	GOL	B	330	-	5,5,5	1.16	0	5,5,5	1.41	0
4	GOL	E	330	-	5,5,5	1.33	0	5,5,5	1.62	1 (20%)
2	PO4	B	328	-	4,4,4	1.17	0	6,6,6	0.95	0
4	GOL	A	330	-	5,5,5	1.59	2 (40%)	5,5,5	1.65	2 (40%)
2	PO4	D	328	-	4,4,4	0.48	0	6,6,6	0.39	0
2	PO4	F	328	-	4,4,4	0.72	0	6,6,6	0.79	0
4	GOL	F	330	-	5,5,5	1.01	0	5,5,5	0.85	0
2	PO4	E	328	-	4,4,4	0.44	0	6,6,6	0.88	0
6	EDO	E	569	-	3,3,3	0.32	0	2,2,2	0.72	0
2	PO4	A	328	-	4,4,4	1.20	0	6,6,6	0.58	0
4	GOL	B	331	-	5,5,5	0.39	0	5,5,5	0.59	0
3	NDG	D	329	-	15,15,15	0.66	0	21,21,21	1.33	2 (9%)
3	NDG	B	329	-	15,15,15	0.66	0	21,21,21	1.01	1 (4%)
6	EDO	F	569	-	3,3,3	0.28	0	2,2,2	0.46	0
6	EDO	E	570	-	3,3,3	0.82	0	2,2,2	0.54	0
2	PO4	C	328	-	4,4,4	1.25	0	6,6,6	0.40	0
6	EDO	A	597	-	3,3,3	0.49	0	2,2,2	0.18	0
6	EDO	A	569	-	3,3,3	0.56	0	2,2,2	0.07	0
3	NDG	E	329	-	15,15,15	0.64	0	21,21,21	1.25	3 (14%)
3	NDG	C	329	-	15,15,15	0.62	0	21,21,21	1.10	1 (4%)
3	NDG	F	329	-	15,15,15	0.73	0	21,21,21	1.06	1 (4%)

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
6	EDO	E	569	-	-	1/1/1/1	-
4	GOL	B	331	-	-	2/4/4/4	-
3	NDG	A	329	-	-	0/6/26/26	0/1/1/1
4	GOL	D	330	-	-	4/4/4/4	-
6	EDO	A	597	-	-	1/1/1/1	-
6	EDO	A	569	-	-	0/1/1/1	-
4	GOL	C	330	-	-	2/4/4/4	-
4	GOL	B	330	-	-	2/4/4/4	-
3	NDG	E	329	-	-	0/6/26/26	0/1/1/1
4	GOL	E	330	-	-	4/4/4/4	-
3	NDG	C	329	-	-	0/6/26/26	0/1/1/1
3	NDG	D	329	-	-	0/6/26/26	0/1/1/1
4	GOL	A	330	-	-	4/4/4/4	-
3	NDG	F	329	-	-	0/6/26/26	0/1/1/1
3	NDG	B	329	-	-	0/6/26/26	0/1/1/1
6	EDO	F	569	-	-	1/1/1/1	-
6	EDO	E	570	-	-	1/1/1/1	-
4	GOL	F	330	-	-	3/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	330	GOL	O1-C1	2.29	1.52	1.42
4	A	330	GOL	O2-C2	2.15	1.49	1.43

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	329	NDG	C1-C2-N2	3.39	114.66	110.73
3	C	329	NDG	O5-C1-C2	2.85	112.38	109.52
4	A	330	GOL	O2-C2-C1	2.79	120.72	109.18
4	E	330	GOL	O2-C2-C1	2.73	120.50	109.18
3	D	329	NDG	C1-C2-N2	2.55	113.68	110.73
3	F	329	NDG	C6-C5-C4	-2.53	106.82	113.02
3	D	329	NDG	O5-C1-C2	2.51	112.04	109.52
3	E	329	NDG	C1-C2-C3	-2.26	107.46	110.54
3	A	329	NDG	O6-C6-C5	2.21	118.85	111.33
3	A	329	NDG	C1-C2-N2	2.19	113.27	110.73
3	E	329	NDG	O7-C7-N2	2.10	125.69	121.98
4	A	330	GOL	O1-C1-C2	2.03	119.51	110.38
3	B	329	NDG	O7-C7-N2	2.01	125.53	121.98

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	330	GOL	O1-C1-C2-C3
4	B	331	GOL	O1-C1-C2-O2
4	B	331	GOL	O1-C1-C2-C3
4	D	330	GOL	C1-C2-C3-O3
4	E	330	GOL	O1-C1-C2-C3
4	F	330	GOL	O1-C1-C2-C3
4	A	330	GOL	O1-C1-C2-O2
4	B	330	GOL	C1-C2-C3-O3
4	C	330	GOL	O1-C1-C2-C3
4	D	330	GOL	O1-C1-C2-C3
4	B	330	GOL	O2-C2-C3-O3
4	C	330	GOL	O1-C1-C2-O2
4	D	330	GOL	O1-C1-C2-O2
4	D	330	GOL	O2-C2-C3-O3
4	E	330	GOL	O1-C1-C2-O2
4	F	330	GOL	O1-C1-C2-O2
6	A	597	EDO	O1-C1-C2-O2
6	E	569	EDO	O1-C1-C2-O2
4	E	330	GOL	O2-C2-C3-O3
4	F	330	GOL	O2-C2-C3-O3
4	E	330	GOL	C1-C2-C3-O3
4	A	330	GOL	O2-C2-C3-O3
6	E	570	EDO	O1-C1-C2-O2
6	F	569	EDO	O1-C1-C2-O2
4	A	330	GOL	C1-C2-C3-O3

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1
1	C	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	95:GLU	C	96:ILE	N	1.14
1	C	326:ILE	C	327:ILE	N	1.11

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	320/347 (92%)	-0.31	2 (0%) 85 88	7, 17, 31, 47	9 (2%)
1	B	321/347 (92%)	-0.37	6 (1%) 66 70	7, 18, 34, 48	7 (2%)
1	C	320/347 (92%)	-0.37	5 (1%) 70 74	7, 18, 33, 51	4 (1%)
1	D	320/347 (92%)	-0.17	5 (1%) 70 74	8, 22, 38, 53	9 (2%)
1	E	321/347 (92%)	-0.15	4 (1%) 76 80	9, 21, 36, 52	7 (2%)
1	F	320/347 (92%)	-0.00	6 (1%) 66 70	10, 24, 41, 57	4 (1%)
All	All	1922/2082 (92%)	-0.23	28 (1%) 72 76	7, 20, 36, 57	40 (2%)

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	8	ILE	4.6
1	C	8	ILE	4.4
1	F	8	ILE	4.3
1	D	52	ASP	3.6
1	E	17	GLY	3.3
1	B	17	GLY	3.3
1	B	16	ALA	3.1
1	C	327	ILE	3.1
1	E	327	ILE	3.1
1	C	92	ASP	2.9
1	C	52	ASP	2.8
1	A	92	ASP	2.8
1	D	327	ILE	2.6
1	F	92	ASP	2.5
1	E	45	SER	2.5
1	B	7	VAL	2.5
1	F	16	ALA	2.5
1	B	327	ILE	2.5
1	F	9	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	16	ALA	2.3
1	D	92	ASP	2.2
1	B	45	SER	2.2
1	F	45	SER	2.2
1	B	162	ARG	2.1
1	F	17	GLY	2.1
1	E	16	ALA	2.1
1	D	17	GLY	2.1
1	D	8	ILE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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
4	GOL	D	330	6/6	0.66	0.21	42,47,50,55	0
4	GOL	B	330	6/6	0.74	0.21	38,51,53,54	0
4	GOL	F	330	6/6	0.76	0.17	41,46,47,48	0
6	EDO	E	570	4/4	0.79	0.18	41,45,46,46	0
4	GOL	E	330	6/6	0.81	0.17	35,43,46,46	0
4	GOL	A	330	6/6	0.83	0.15	28,37,40,42	0
4	GOL	B	331	6/6	0.83	0.17	46,47,51,53	0
3	NDG	B	329	15/15	0.84	0.13	29,32,33,35	0
6	EDO	A	569	4/4	0.84	0.13	49,52,56,56	0
4	GOL	C	330	6/6	0.84	0.14	32,40,43,53	0
6	EDO	E	569	4/4	0.85	0.14	45,46,47,48	0
3	NDG	A	329	15/15	0.85	0.12	21,28,32,33	0
3	NDG	C	329	15/15	0.86	0.11	26,27,29,30	0
3	NDG	F	329	15/15	0.86	0.11	28,32,33,36	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	K	C	333	1/1	0.87	0.36	60,60,60,60	0
3	NDG	D	329	15/15	0.87	0.12	24,29,33,37	0
6	EDO	F	569	4/4	0.87	0.15	50,52,54,54	0
3	NDG	E	329	15/15	0.89	0.10	24,26,28,29	0
5	K	F	333	1/1	0.89	0.33	62,62,62,62	0
5	K	B	333	1/1	0.90	0.32	52,52,52,52	0
6	EDO	A	597	4/4	0.91	0.13	42,43,51,52	0
5	K	A	333	1/1	0.94	0.33	53,53,53,53	0
5	K	D	333	1/1	0.95	0.32	58,58,58,58	0
5	K	E	333	1/1	0.96	0.31	51,51,51,51	0
2	PO4	A	328	5/5	0.99	0.03	13,14,16,17	0
2	PO4	B	328	5/5	0.99	0.03	14,15,17,19	0
2	PO4	C	328	5/5	0.99	0.03	14,15,17,18	0
2	PO4	D	328	5/5	0.99	0.03	17,17,19,21	0
2	PO4	E	328	5/5	0.99	0.03	15,17,19,20	0
2	PO4	F	328	5/5	0.99	0.03	18,18,21,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.