



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 02:33 AM UTC

PDB ID : 4UDI / pdb_00004udi
Title : Crystal structure of b-1,4-mannopyranosyl-chitobiose phosphorylase at 1.85 Angstrom from unknown human gut bacteria (Uhgb_MP)
Authors : Ladeveze, S.; Cioci, G.; Potocki-Veronese, G.; Tranier, S.; Mourey, L.
Deposited on : 2014-12-10
Resolution : 1.80 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

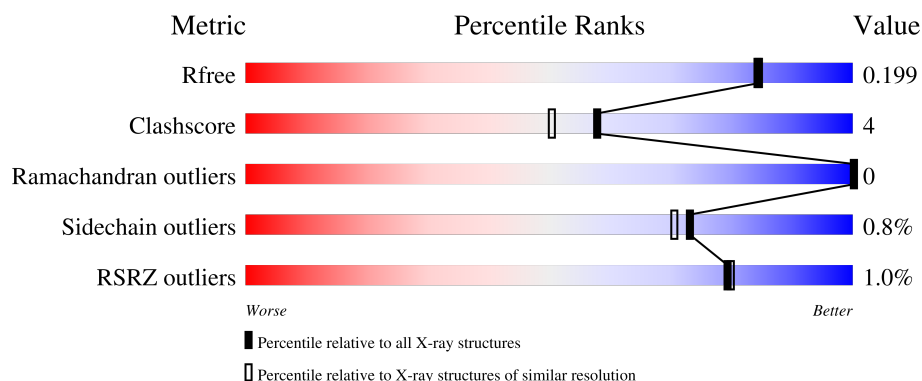
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	80% 12% 7%
1	B	347	84% 8% 7%
1	C	347	80% 12% 7%
1	D	347	85% 7% 8%
1	E	347	84% 8% 8%

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Mol	Chain	Length	Quality of chain
1	F	347	% 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
3	GOL	A	332	-	X	-	-
3	GOL	A	335	-	X	-	-
5	EDO	A	337	-	-	X	-
5	EDO	A	339	-	-	X	-
5	EDO	B	339	-	-	X	-
5	EDO	C	336	-	-	X	-
5	EDO	C	339	-	-	X	-
5	EDO	C	340	-	-	X	-

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 16842 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	321	Total	C	N	O	S	0	5	0
			2601	1683	436	465	17			
1	B	321	Total	C	N	O	S	0	2	0
			2591	1674	435	465	17			
1	C	321	Total	C	N	O	S	0	4	0
			2595	1679	434	465	17			
1	D	320	Total	C	N	O	S	0	1	0
			2558	1652	429	460	17			
1	E	320	Total	C	N	O	S	0	1	0
			2556	1653	428	459	16			
1	F	320	Total	C	N	O	S	0	2	0
			2566	1658	429	462	17			

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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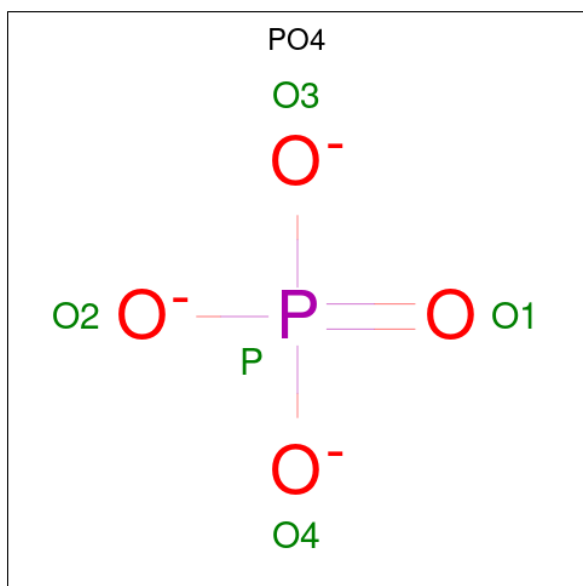
Chain	Residue	Modelled	Actual	Comment	Reference
C	0	HIS	-	expression tag	UNP D9ZDQ9
D	-19	MET	-	expression tag	UNP D9ZDQ9
D	-18	GLY	-	expression tag	UNP D9ZDQ9
D	-17	SER	-	expression tag	UNP D9ZDQ9
D	-16	SER	-	expression tag	UNP D9ZDQ9
D	-15	HIS	-	expression tag	UNP D9ZDQ9
D	-14	HIS	-	expression tag	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).



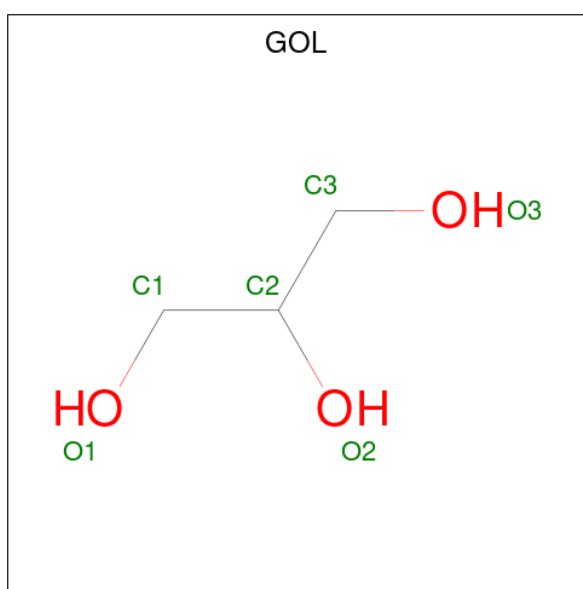
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	P	0	0
			5	4	1		
2	B	1	Total	O	P	0	0
			5	4	1		

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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 GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

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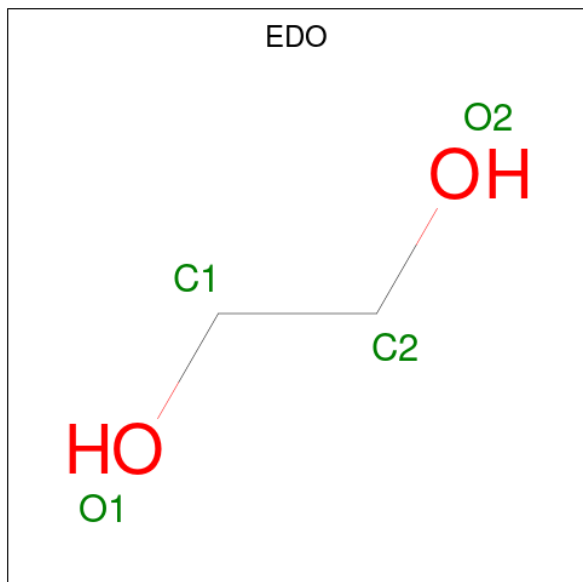
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0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

- Molecule 4 is POTASSIUM ION (CCD ID: K) (formula: K).

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

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



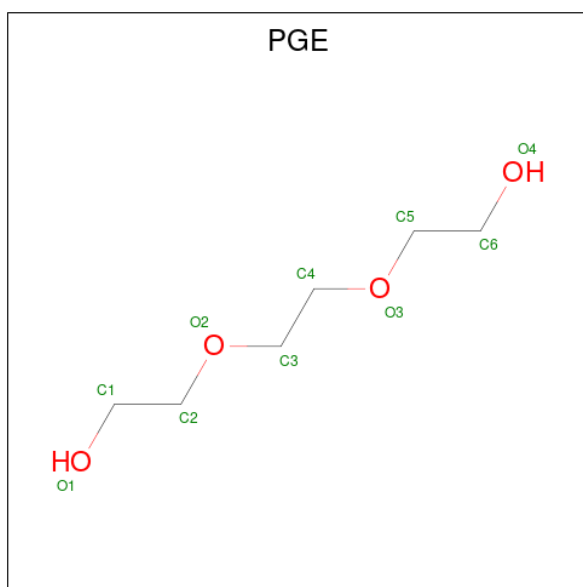
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0

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

- Molecule 6 is TRIETHYLENE GLYCOL (CCD ID: PGE) (formula: C₆H₁₄O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			10	6	4		
6	B	1	Total	C	O	0	0
			10	6	4		
6	C	1	Total	C	O	0	0
			10	6	4		

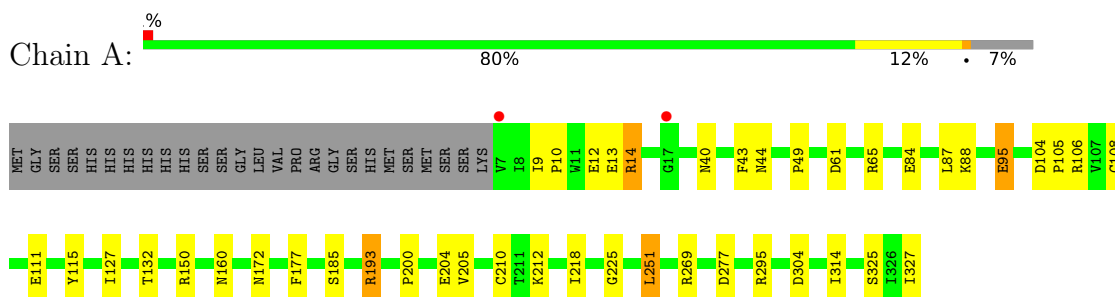
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	205	Total	O	0	0
			205	205		
7	B	213	Total	O	0	0
			213	213		
7	C	229	Total	O	0	0
			229	229		
7	D	144	Total	O	0	0
			144	144		
7	E	143	Total	O	0	0
			143	143		
7	F	129	Total	O	0	0
			129	129		

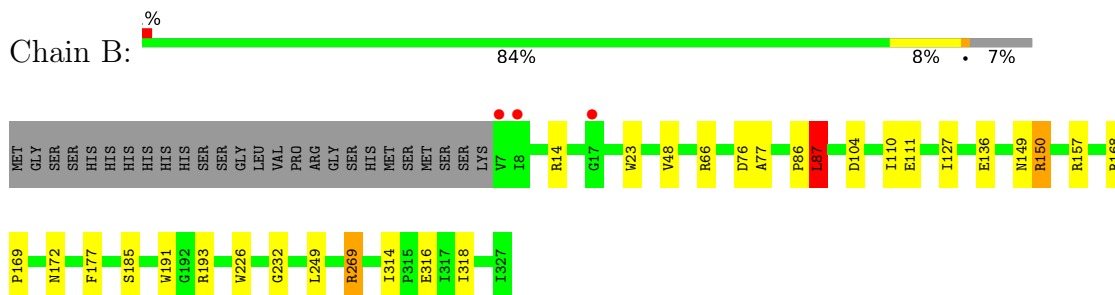
3 Residue-property plots [i](#)

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.

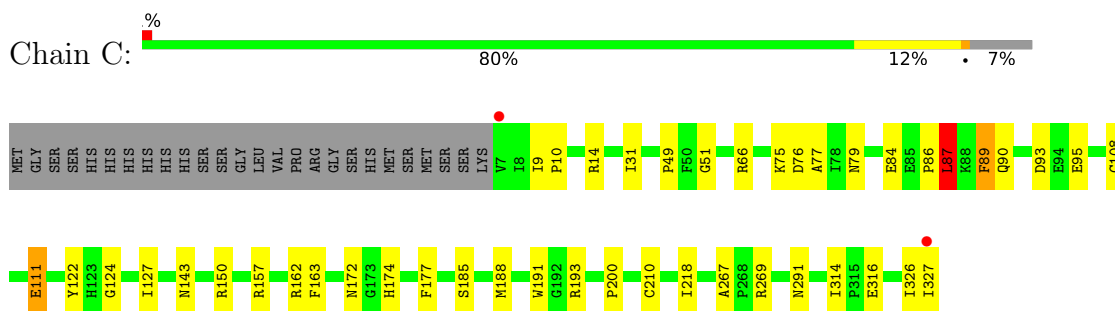
• Molecule 1: UHGB_MP



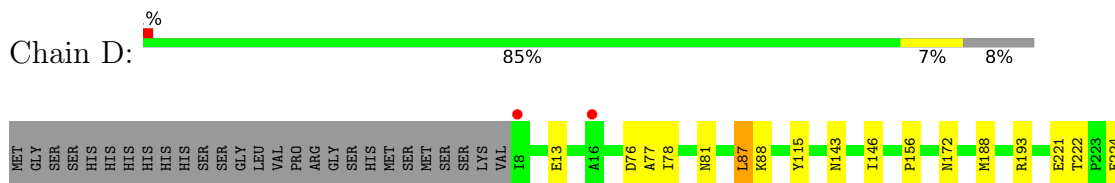
• Molecule 1: UHGB_MP

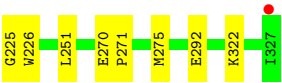


• Molecule 1: UHGB_MP

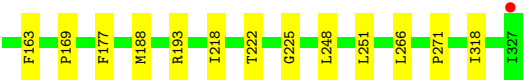
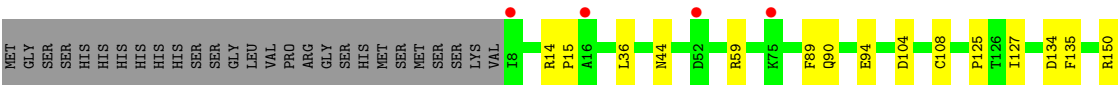
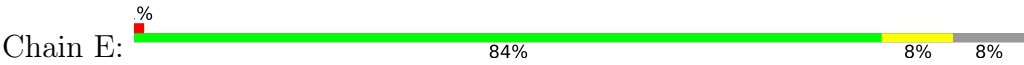


• Molecule 1: UHGB_MP

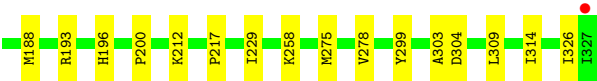
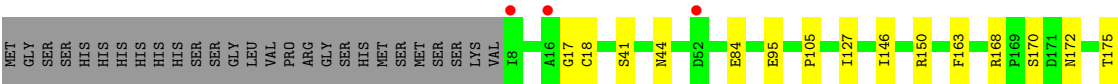
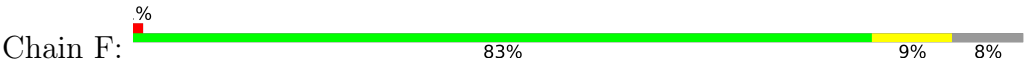




● Molecule 1: UHGB_MP



● 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	84.12Å 141.21Å 176.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	110.20 – 1.80 110.20 – 1.80	Depositor EDS
% Data completeness (in resolution range)	98.0 (110.20-1.80) 98.1 (110.20-1.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 1.81Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.155 , 0.190 0.170 , 0.199	Depositor DCC
R_{free} test set	9557 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	22.9	Xtriage
Anisotropy	0.779	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 39.0	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	0.97	EDS
Total number of atoms	16842	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.04% 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: GOL, K, PGE, PO4, EDO

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	1.39	7/2696 (0.3%)	1.23	11/3676 (0.3%)
1	B	1.46	12/2683 (0.4%)	1.15	3/3657 (0.1%)
1	C	1.45	15/2693 (0.6%)	1.18	8/3672 (0.2%)
1	D	1.10	5/2647 (0.2%)	1.02	4/3610 (0.1%)
1	E	1.10	3/2648 (0.1%)	1.13	8/3613 (0.2%)
1	F	1.21	4/2658 (0.2%)	1.09	9/3625 (0.2%)
All	All	1.30	46/16025 (0.3%)	1.14	43/21853 (0.2%)

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	B	0	1
1	C	0	1
All	All	0	2

The worst 5 of 46 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	326	ILE	C-N	-27.40	0.94	1.33
1	B	87	LEU	C-N	-15.78	1.10	1.33
1	C	87	LEU	C-N	-15.20	1.11	1.33
1	C	326	ILE	C-N	-13.37	1.14	1.33
1	C	90	GLN	C-N	-12.04	1.18	1.33

The worst 5 of 43 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	90	GLN	O-C-N	-22.56	96.99	123.16
1	C	90	GLN	O-C-N	-19.77	100.68	123.22
1	A	13	GLU	O-C-N	17.31	142.53	122.79
1	F	17	GLY	CA-C-N	13.43	143.07	121.86
1	F	17	GLY	C-N-CA	13.43	143.07	121.86

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	87	LEU	Mainchain
1	C	87	LEU	Mainchain

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	2601	0	2465	34	0
1	B	2591	0	2460	24	0
1	C	2595	0	2462	28	0
1	D	2558	0	2400	12	0
1	E	2556	0	2403	13	0
1	F	2566	0	2409	15	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	24	0	32	5	0
3	B	36	0	48	6	0
3	C	36	0	48	6	0
3	D	24	0	32	2	0
3	E	12	0	16	2	0
3	F	30	0	40	3	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
5	A	28	0	41	12	0
5	B	8	0	12	6	0
5	C	24	0	36	20	0
5	D	4	0	6	0	0
5	E	12	0	18	1	0
5	F	8	0	12	2	0
6	A	10	0	14	3	0
6	B	10	0	14	0	0
6	C	10	0	14	1	0
7	A	205	0	0	8	1
7	B	213	0	0	1	0
7	C	229	0	0	6	1
7	D	144	0	0	1	0
7	E	143	0	0	1	0
7	F	129	0	0	1	0
All	All	16842	0	14982	138	1

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.

The worst 5 of 138 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:277:ASP:OD1	5:C:336:EDO:H21	1.70	0.90
1:C:269:ARG:HD3	3:C:335:GOL:H12	1.54	0.87
1:A:14[A]:ARG:HH11	1:A:14[A]:ARG:HG3	1.40	0.87
1:D:172:ASN:HB2	3:D:332:GOL:O1	1.74	0.87
1:C:163:PHE:HZ	5:C:340:EDO:H11	1.45	0.82

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:505:HOH:O	7:C:390:HOH:O[1_455]	1.79	0.41

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	324/347 (93%)	305 (94%)	19 (6%)	0	100	100
1	B	321/347 (92%)	298 (93%)	23 (7%)	0	100	100
1	C	323/347 (93%)	302 (94%)	21 (6%)	0	100	100
1	D	319/347 (92%)	302 (95%)	17 (5%)	0	100	100
1	E	319/347 (92%)	302 (95%)	17 (5%)	0	100	100
1	F	320/347 (92%)	304 (95%)	16 (5%)	0	100	100
All	All	1926/2082 (92%)	1813 (94%)	113 (6%)	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	276/298 (93%)	273 (99%)	3 (1%)	65	60
1	B	277/298 (93%)	275 (99%)	2 (1%)	76	73
1	C	277/298 (93%)	275 (99%)	2 (1%)	76	73
1	D	270/298 (91%)	268 (99%)	2 (1%)	76	73
1	E	269/298 (90%)	266 (99%)	3 (1%)	65	60
1	F	271/298 (91%)	269 (99%)	2 (1%)	76	73
All	All	1640/1788 (92%)	1626 (99%)	14 (1%)	73	67

5 of 14 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	13	GLU
1	D	87	LEU
1	F	168	ARG
1	E	266	LEU
1	F	95	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	160	ASN
1	D	44	ASN
1	D	139	HIS
1	D	79	ASN
1	B	280	ASN

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](#)

Of 63 ligands modelled in this entry, 6 are monoatomic - leaving 57 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
5	EDO	F	337	-	3,3,3	0.14	0	2,2,2	0.53	0
6	PGE	C	341	-	9,9,9	1.00	0	8,8,8	1.57	2 (25%)
5	EDO	B	346	-	3,3,3	0.65	0	2,2,2	0.24	0
5	EDO	A	337	-	3,3,3	0.99	0	2,2,2	0.48	0
5	EDO	A	569	-	3,3,3	0.65	0	2,2,2	0.23	0
5	EDO	C	340	-	3,3,3	0.53	0	2,2,2	0.59	0
5	EDO	C	347	-	3,3,3	0.63	0	2,2,2	0.78	0
5	EDO	E	573	-	3,3,3	0.41	0	2,2,2	0.94	0
3	GOL	F	329	-	5,5,5	0.29	0	5,5,5	0.65	0
3	GOL	C	335	-	5,5,5	1.03	0	5,5,5	0.87	0
3	GOL	A	329	-	5,5,5	0.41	0	5,5,5	0.49	0
3	GOL	F	334	-	5,5,5	0.39	0	5,5,5	0.91	0
2	PO4	A	328	-	4,4,4	1.36	0	6,6,6	1.17	1 (16%)
3	GOL	B	334	-	5,5,5	0.33	0	5,5,5	1.90	2 (40%)
5	EDO	D	339	-	3,3,3	0.26	0	2,2,2	0.55	0
6	PGE	B	341	-	9,9,9	0.73	0	8,8,8	0.68	0
5	EDO	A	348	-	3,3,3	0.64	0	2,2,2	0.32	0
2	PO4	B	328	-	4,4,4	0.96	0	6,6,6	0.90	0
3	GOL	E	329	-	5,5,5	0.26	0	5,5,5	0.41	0
2	PO4	E	328	-	4,4,4	1.62	1 (25%)	6,6,6	1.46	2 (33%)
3	GOL	B	332	-	5,5,5	0.63	0	5,5,5	1.43	1 (20%)
3	GOL	F	332	-	5,5,5	0.39	0	5,5,5	0.80	0
5	EDO	C	343	-	3,3,3	0.30	0	2,2,2	1.21	0
3	GOL	F	330	-	5,5,5	0.38	0	5,5,5	0.30	0
5	EDO	E	570	-	3,3,3	0.49	0	2,2,2	0.63	0
3	GOL	B	331	-	5,5,5	0.44	0	5,5,5	1.54	2 (40

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	C	332	-	5,5,5	0.54	0	5,5,5	0.83	0
3	GOL	D	332	-	5,5,5	0.67	0	5,5,5	2.02	2 (40%)
5	EDO	F	570	-	3,3,3	0.39	0	2,2,2	0.82	0
5	EDO	A	339	-	3,3,3	0.22	0	2,2,2	0.43	0
3	GOL	B	335	-	5,5,5	0.67	0	5,5,5	1.19	0
5	EDO	A	345	-	3,3,3	0.50	0	2,2,2	1.37	0
3	GOL	B	329	-	5,5,5	0.30	0	5,5,5	0.63	0
3	GOL	F	335	-	5,5,5	0.61	0	5,5,5	0.50	0
5	EDO	B	339	-	3,3,3	0.60	0	2,2,2	0.65	0
3	GOL	A	335	-	5,5,5	1.06	0	5,5,5	1.67	2 (40%)
3	GOL	D	330	-	5,5,5	0.31	0	5,5,5	0.95	0
5	EDO	C	336	-	3,3,3	0.39	0	2,2,2	1.14	0
3	GOL	C	330	-	5,5,5	0.48	0	5,5,5	0.73	0
5	EDO	C	339	-	3,3,3	0.14	0	2,2,2	0.91	0
3	GOL	B	330	-	5,5,5	0.89	0	5,5,5	0.56	0
3	GOL	D	331	-	5,5,5	0.46	0	5,5,5	1.16	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
5	EDO	F	337	-	-	0/1/1/1	-
6	PGE	C	341	-	-	2/7/7/7	-
5	EDO	B	346	-	-	0/1/1/1	-
5	EDO	A	337	-	-	0/1/1/1	-
5	EDO	A	569	-	-	1/1/1/1	-
5	EDO	C	340	-	-	0/1/1/1	-
5	EDO	C	347	-	-	1/1/1/1	-
5	EDO	E	573	-	-	1/1/1/1	-
3	GOL	F	329	-	-	0/4/4/4	-
3	GOL	C	335	-	-	4/4/4/4	-
3	GOL	A	329	-	-	0/4/4/4	-
3							

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	F	332	-	-	4/4/4/4	-
5	EDO	C	343	-	-	0/1/1/1	-
3	GOL	F	330	-	-	2/4/4/4	-
5	EDO	E	570	-	-	0/1/1/1	-
3	GOL	B	331	-	-	0/4/4/4	-
3	GOL	A	332	-	-	4/4/4/4	-
3	GOL	A	330	-	-	2/4/4/4	-
5	EDO	A	597	-	-	1/1/1/1	-
5	EDO	C	344	-	-	1/1/1/1	-
3	GOL	E	332	-	-	4/4/4/4	-
3	GOL	C	329	-	-	0/4/4/4	-
3	GOL	D	329	-	-	0/4/4/4	-
3	GOL	C	334	-	-	4/4/4/4	-
5	EDO	A	346	-	-	1/1/1/1	-
5	EDO	E	569	-	-	1/1/1/1	-
6	PGE	A	341	-	-	2/7/7/7	-
3	GOL	C	331	-	-	2/4/4/4	-
3	GOL	C	332	-	-	4/4/4/4	-
3	GOL	D	332	-	-	3/4/4/4	-
5	EDO	F	570	-	-	1/1/1/1	-
5	EDO	A	339	-	-	0/1/1/1	-
3	GOL	B	335	-	-	4/4/4/4	-
5	EDO	A	345	-	-	0/1/1/1	-
3	GOL	B	329	-	-	2/4/4/4	-
3	GOL	F	335	-	-	0/4/4/4	-
5	EDO	B	339	-	-	0/1/1/1	-
3	GOL	A	335	-	-	4/4/4/4	-
3	GOL	D	330	-	-	0/4/4/4	-
5	EDO	C	336	-	-	0/1/1/1	-
3	GOL	C	330	-	-	0/4/4/4	-
5	EDO	C	339	-	-	0/1/1/1	-
3	GOL	B	330	-	-	2/4/4/4	-
3	GOL	D	331	-	-	2/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	328	PO4	P-O1	2.97	1.57	1.50

The worst 5 of 18 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	334	GOL	O3-C3-C2	-3.65	93.93	110.38
3	A	332	GOL	C3-C2-C1	-3.29	99.74	111.80
3	D	332	GOL	O1-C1-C2	-3.26	95.70	110.38
3	E	332	GOL	O2-C2-C3	3.17	122.32	109.18
6	C	341	PGE	O3-C5-C6	3.16	124.02	110.11

There are no chirality outliers.

5 of 70 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	330	GOL	C1-C2-C3-O3
3	A	332	GOL	O1-C1-C2-C3
3	A	335	GOL	O1-C1-C2-C3
3	A	335	GOL	C1-C2-C3-O3
3	B	334	GOL	O1-C1-C2-C3

There are no ring outliers.

28 monomers are involved in 68 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	F	337	EDO	2	0
6	C	341	PGE	1	0
5	B	346	EDO	1	0
5	A	337	EDO	5	0
5	C	340	EDO	4	0
5	C	347	EDO	3	0
5	E	573	EDO	1	0
3	C	335	GOL	2	0
3	B	334	GOL	1	0
3	B	332	GOL	3	0
3	F	332	GOL	3	0
5	C	343	EDO	3	0
3	A	332	GOL	2	0
3	A	330	GOL	1	0
5	A	597	EDO	1	0
3	E	332	GOL	2	0
5	A	346	EDO	2	0
6	A	341	PGE	3	0
3	C	331	GOL	1	0
3	C	332	GOL	2	0
3	D	332	GOL	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	339	EDO	4	0
3	B	335	GOL	2	0
5	B	339	EDO	5	0
3	A	335	GOL	2	0
5	C	336	EDO	6	0
3	C	330	GOL	1	0
5	C	339	EDO	4	0

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	C	4
1	B	1
1	F	1

The worst 5 of 6 chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	C	89:PHE	C	90:GLN	N	1.20
1	C	90:GLN	C	91:CYS	N	1.17
1	C	326:ILE	C	327:ILE	N	1.14
1	C	87:LEU	C	88:LYS	N	1.11
1	B	87:LEU	C	88:LYS	N	1.10

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	321/347 (92%)	-0.47	2 (0%) 85 86	11, 26, 43, 66	5 (1%)
1	B	321/347 (92%)	-0.50	3 (0%) 81 81	15, 25, 39, 57	2 (0%)
1	C	321/347 (92%)	-0.54	2 (0%) 85 86	13, 25, 42, 63	4 (1%)
1	D	320/347 (92%)	-0.13	3 (0%) 81 81	19, 39, 59, 82	1 (0%)
1	E	320/347 (92%)	-0.05	5 (1%) 70 71	17, 40, 63, 81	1 (0%)
1	F	320/347 (92%)	-0.10	4 (1%) 75 75	18, 40, 60, 82	2 (0%)
All	All	1923/2082 (92%)	-0.30	19 (0%) 79 80	11, 32, 58, 82	15 (0%)

The worst 5 of 19 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	8	ILE	5.1
1	F	8	ILE	3.9
1	C	7	VAL	3.8
1	A	7	VAL	3.7
1	F	327	ILE	3.4

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
5	EDO	A	345	4/4	0.76	0.17	56,58,59,64	0
5	EDO	C	343	4/4	0.76	0.16	45,48,53,54	0
6	PGE	A	341	10/10	0.77	0.19	44,58,69,73	0
5	EDO	E	573	4/4	0.81	0.15	54,59,61,65	0
6	PGE	B	341	10/10	0.81	0.17	43,62,67,69	0
5	EDO	A	348	4/4	0.82	0.12	53,57,58,62	0
6	PGE	C	341	10/10	0.82	0.15	36,58,62,62	0
5	EDO	A	597	4/4	0.83	0.13	56,57,59,67	0
5	EDO	E	570	4/4	0.83	0.15	57,58,60,63	0
3	GOL	B	331	6/6	0.84	0.16	54,55,58,64	0
3	GOL	E	332	6/6	0.85	0.15	47,49,51,57	0
3	GOL	C	331	6/6	0.85	0.16	53,56,59,61	0
3	GOL	F	335	6/6	0.86	0.14	58,66,67,71	0
3	GOL	C	335	6/6	0.86	0.13	36,46,52,53	0
5	EDO	F	570	4/4	0.86	0.13	52,61,64,69	0
3	GOL	D	331	6/6	0.86	0.13	59,60,61,62	0
3	GOL	A	332	6/6	0.86	0.14	46,52,56,58	0
3	GOL	F	332	6/6	0.86	0.13	50,52,54,57	0
3	GOL	D	332	6/6	0.88	0.13	40,48,55,60	0
3	GOL	A	330	6/6	0.88	0.12	43,53,55,57	0
5	EDO	C	344	4/4	0.88	0.13	52,55,61,61	0
3	GOL	C	332	6/6	0.89	0.12	36,46,58,62	0
3	GOL	F	334	6/6	0.90	0.10	56,58,62,62	0
5	EDO	E	569	4/4	0.90	0.12	55,60,61,65	0
3	GOL	B	334	6/6	0.90	0.10	43,48,51,52	0
3	GOL	B	332	6/6	0.90	0.12	40,49,55,58	0
3	GOL	B	330	6/6	0.91	0.11	39,48,53,58	0
5	EDO	A	569	4/4	0.			

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GOL	A	335	6/6	0.93	0.11	33,44,47,53	0
5	EDO	F	337	4/4	0.93	0.25	40,43,43,46	0
5	EDO	B	346	4/4	0.93	0.26	37,37,42,43	0
3	GOL	E	329	6/6	0.93	0.09	45,46,47,50	0
3	GOL	D	330	6/6	0.93	0.09	55,58,59,60	0
5	EDO	D	339	4/4	0.93	0.09	51,55,55,56	0
4	K	D	333	1/1	0.94	0.17	60,60,60,60	0
3	GOL	C	330	6/6	0.94	0.08	41,44,46,49	0
5	EDO	C	347	4/4	0.94	0.12	37,48,52,52	0
5	EDO	B	339	4/4	0.94	0.20	33,34,35,35	0
3	GOL	A	329	6/6	0.94	0.08	29,33,34,36	0
5	EDO	C	340	4/4	0.94	0.19	35,43,44,44	0
5	EDO	A	339	4/4	0.95	0.24	32,33,33,33	0
4	K	A	333	1/1	0.95	0.25	50,50,50,50	0
3	GOL	C	329	6/6	0.95	0.08	27,31,31,35	0
4	K	C	333	1/1	0.96	0.20	50,50,50,50	0
5	EDO	A	337	4/4	0.97	0.18	26,27,27,30	0
5	EDO	C	336	4/4	0.97	0.14	28,28,30,30	0
5	EDO	C	339	4/4	0.97	0.21	29,32,34,35	0
4	K	B	333	1/1	0.97	0.18	48,48,48,48	0
2	PO4	D	328	5/5	0.99	0.07	30,30,32,35	0
2	PO4	E	328	5/5	0.99	0.06	32,34,34,36	0
2	PO4	F	328	5/5	0.99	0.07	29,29,34,35	0
2	PO4	A	328	5/5	0.99	0.06	23,24,25,26	0
2	PO4	B	328	5/5	0.99	0.05	21,21,24,24	0
2	PO4	C	328	5/5	1.00	0.03	23,23,24,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.