



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 04:18 PM GMT

PDB ID : 2AE7
Title : Crystal Structure of Human M340H-Beta1,4-Galactosyltransferase-I(M340H-B4GAL-T1) in Complex with Pentasaccharide
Authors : Ramasamy, V.; Ramakrishnan, B.; Boeggeman, E.; Ratner, D.M.; Seeberger, P.H.; Qasba, P.K.
Deposited on : 2005-07-21
Resolution : 2.00 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

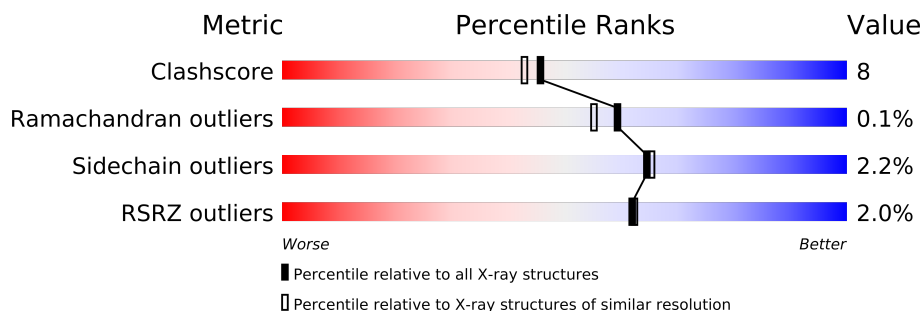
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.00 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))
Clashscore	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)
RSRZ outliers	66119	4890 (2.00-2.00)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	287	
1	B	287	
1	C	287	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
5	SO4	A	421	-	X
5	SO4	A	429	-	X
5	SO4	B	422	-	X
5	SO4	C	431	-	X
8	GOL	A	415	-	X
8	GOL	B	413	-	X
8	GOL	B	414	-	X
8	GOL	B	418	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
8	GOL	C	416	-	X

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 7413 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Beta-1,4-galactosyltransferase1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	272	Total	C	N	O	S	0	0	0
			2212	1417	385	399	11			
1	B	272	Total	C	N	O	S	0	0	0
			2212	1417	385	399	11			
1	C	272	Total	C	N	O	S	0	0	0
			2212	1417	385	399	11			

There are 51 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	112	ALA	-	SEE REMARK 999	UNP P15291
A	113	SER	-	SEE REMARK 999	UNP P15291
A	114	MET	-	SEE REMARK 999	UNP P15291
A	115	THR	-	SEE REMARK 999	UNP P15291
A	116	GLY	-	SEE REMARK 999	UNP P15291
A	117	GLY	-	SEE REMARK 999	UNP P15291
A	118	GLN	-	SEE REMARK 999	UNP P15291
A	119	GLN	-	SEE REMARK 999	UNP P15291
A	120	MET	-	SEE REMARK 999	UNP P15291
A	121	GLY	-	SEE REMARK 999	UNP P15291
A	122	ARG	-	SEE REMARK 999	UNP P15291
A	123	GLY	-	SEE REMARK 999	UNP P15291
A	124	SER	-	SEE REMARK 999	UNP P15291
A	125	ALA	-	SEE REMARK 999	UNP P15291
A	337	THR	ARG	ENGINEERED	UNP P15291
A	338	THR	CYS	ENGINEERED	UNP P15291
A	340	HIS	MET	ENGINEERED	UNP P15291
B	112	ALA	-	SEE REMARK 999	UNP P15291
B	113	SER	-	SEE REMARK 999	UNP P15291
B	114	MET	-	SEE REMARK 999	UNP P15291
B	115	THR	-	SEE REMARK 999	UNP P15291
B	116	GLY	-	SEE REMARK 999	UNP P15291
B	117	GLY	-	SEE REMARK 999	UNP P15291

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Chain	Residue	Modelled	Actual	Comment	Reference
B	118	GLN	-	SEE REMARK 999	UNP P15291
B	119	GLN	-	SEE REMARK 999	UNP P15291
B	120	MET	-	SEE REMARK 999	UNP P15291
B	121	GLY	-	SEE REMARK 999	UNP P15291
B	122	ARG	-	SEE REMARK 999	UNP P15291
B	123	GLY	-	SEE REMARK 999	UNP P15291
B	124	SER	-	SEE REMARK 999	UNP P15291
B	125	ALA	-	SEE REMARK 999	UNP P15291
B	337	THR	ARG	ENGINEERED	UNP P15291
B	338	THR	CYS	ENGINEERED	UNP P15291
B	340	HIS	MET	ENGINEERED	UNP P15291
C	112	ALA	-	SEE REMARK 999	UNP P15291
C	113	SER	-	SEE REMARK 999	UNP P15291
C	114	MET	-	SEE REMARK 999	UNP P15291
C	115	THR	-	SEE REMARK 999	UNP P15291
C	116	GLY	-	SEE REMARK 999	UNP P15291
C	117	GLY	-	SEE REMARK 999	UNP P15291
C	118	GLN	-	SEE REMARK 999	UNP P15291
C	119	GLN	-	SEE REMARK 999	UNP P15291
C	120	MET	-	SEE REMARK 999	UNP P15291
C	121	GLY	-	SEE REMARK 999	UNP P15291
C	122	ARG	-	SEE REMARK 999	UNP P15291
C	123	GLY	-	SEE REMARK 999	UNP P15291
C	124	SER	-	SEE REMARK 999	UNP P15291
C	125	ALA	-	SEE REMARK 999	UNP P15291
C	337	THR	ARG	ENGINEERED	UNP P15291
C	338	THR	CYS	ENGINEERED	UNP P15291
C	340	HIS	MET	ENGINEERED	UNP P15291

- Molecule 2 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	2	Total	C	N	O	0	0
			26	14	1	11		
2	C	2	Total	C	N	O	0	0
			26	14	1	11		

There are 34 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	112	ALA	-	SEE REMARK 999	UNP P15291
A	113	SER	-	SEE REMARK 999	UNP P15291

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Chain	Residue	Modelled	Actual	Comment	Reference
A	114	MET	-	SEE REMARK 999	UNP P15291
A	115	THR	-	SEE REMARK 999	UNP P15291
A	116	GLY	-	SEE REMARK 999	UNP P15291
A	117	GLY	-	SEE REMARK 999	UNP P15291
A	118	GLN	-	SEE REMARK 999	UNP P15291
A	119	GLN	-	SEE REMARK 999	UNP P15291
A	120	MET	-	SEE REMARK 999	UNP P15291
A	121	GLY	-	SEE REMARK 999	UNP P15291
A	122	ARG	-	SEE REMARK 999	UNP P15291
A	123	GLY	-	SEE REMARK 999	UNP P15291
A	124	SER	-	SEE REMARK 999	UNP P15291
A	125	ALA	-	SEE REMARK 999	UNP P15291
A	337	THR	ARG	ENGINEERED	UNP P15291
A	338	THR	CYS	ENGINEERED	UNP P15291
A	340	HIS	MET	ENGINEERED	UNP P15291
C	112	ALA	-	SEE REMARK 999	UNP P15291
C	113	SER	-	SEE REMARK 999	UNP P15291
C	114	MET	-	SEE REMARK 999	UNP P15291
C	115	THR	-	SEE REMARK 999	UNP P15291
C	116	GLY	-	SEE REMARK 999	UNP P15291
C	117	GLY	-	SEE REMARK 999	UNP P15291
C	118	GLN	-	SEE REMARK 999	UNP P15291
C	119	GLN	-	SEE REMARK 999	UNP P15291
C	120	MET	-	SEE REMARK 999	UNP P15291
C	121	GLY	-	SEE REMARK 999	UNP P15291
C	122	ARG	-	SEE REMARK 999	UNP P15291
C	123	GLY	-	SEE REMARK 999	UNP P15291
C	124	SER	-	SEE REMARK 999	UNP P15291
C	125	ALA	-	SEE REMARK 999	UNP P15291
C	337	THR	ARG	ENGINEERED	UNP P15291
C	338	THR	CYS	ENGINEERED	UNP P15291
C	340	HIS	MET	ENGINEERED	UNP P15291

- Molecule 3 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	B	3	Total	C	N	O	0	0
			37	20	1	16		

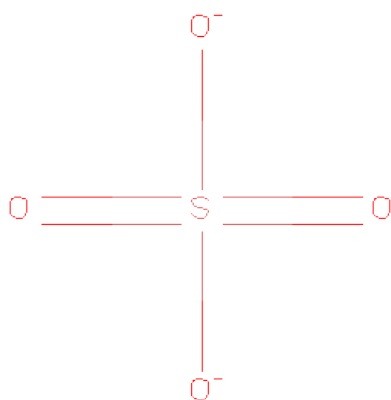
There are 17 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	112	ALA	-	SEE REMARK 999	UNP P15291
B	113	SER	-	SEE REMARK 999	UNP P15291
B	114	MET	-	SEE REMARK 999	UNP P15291
B	115	THR	-	SEE REMARK 999	UNP P15291
B	116	GLY	-	SEE REMARK 999	UNP P15291
B	117	GLY	-	SEE REMARK 999	UNP P15291
B	118	GLN	-	SEE REMARK 999	UNP P15291
B	119	GLN	-	SEE REMARK 999	UNP P15291
B	120	MET	-	SEE REMARK 999	UNP P15291
B	121	GLY	-	SEE REMARK 999	UNP P15291
B	122	ARG	-	SEE REMARK 999	UNP P15291
B	123	GLY	-	SEE REMARK 999	UNP P15291
B	124	SER	-	SEE REMARK 999	UNP P15291
B	125	ALA	-	SEE REMARK 999	UNP P15291
B	337	THR	ARG	ENGINEERED	UNP P15291
B	338	THR	CYS	ENGINEERED	UNP P15291
B	340	HIS	MET	ENGINEERED	UNP P15291

- Molecule 4 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

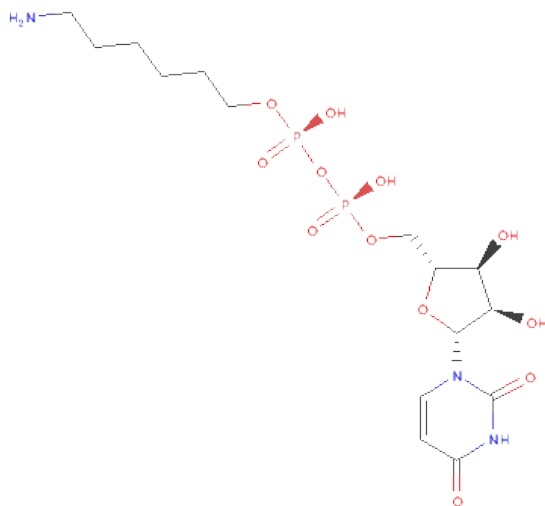
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Mn 1 1	0	0
4	A	1	Total Mn 1 1	0	0
4	C	1	Total Mn 1 1	0	0

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



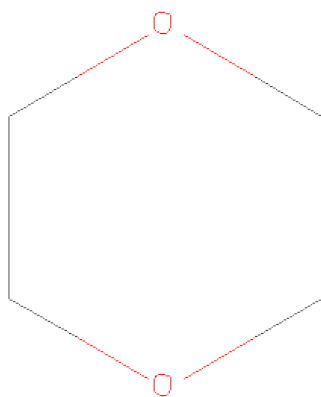
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	C	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		
5	C	1	Total	O	S	0	0
			5	4	1		
5	C	1	Total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

- Molecule 6 is 6-AMINOHEXYL-URIDINE-C1,5'-DIPHOSPHATE (three-letter code: UDH) (formula: $C_{15}H_{27}N_3O_{12}P_2$).



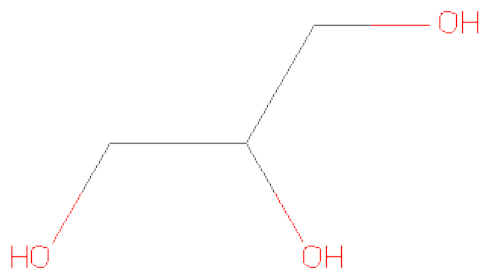
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	P	0	0
			32	15	3	12	2		
6	B	1	Total	C	N	O	P	0	0
			32	15	3	12	2		
6	C	1	Total	C	N	O	P	0	0
			32	15	3	12	2		

- Molecule 7 is 1,4-DIETHYLENE DIOXIDE (three-letter code: DIO) (formula: $C_4H_8O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	C	1	Total	C	O	0	0
			6	4	2		

- Molecule 8 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	157	Total	O	0	0
			157	157		
9	B	192	Total	O	0	0
			192	192		

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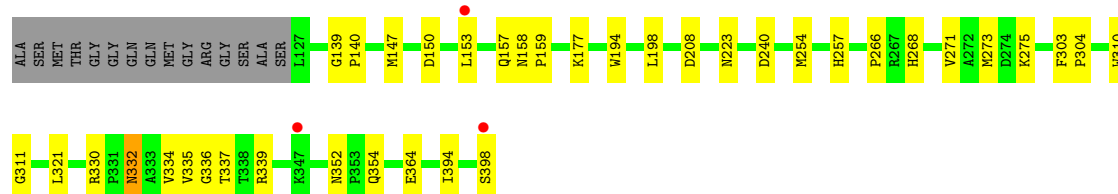
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	C	122	Total 122	O 122	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

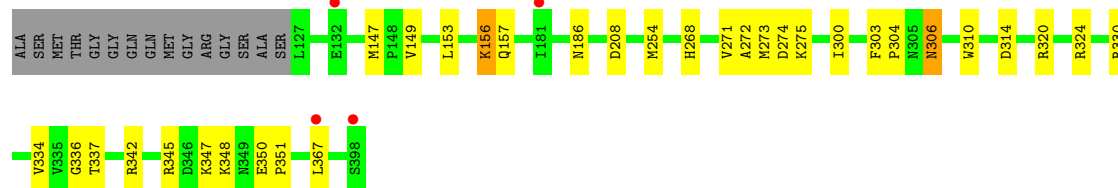
• Molecule 1: Beta-1,4-galactosyltransferase1

Chain A: 



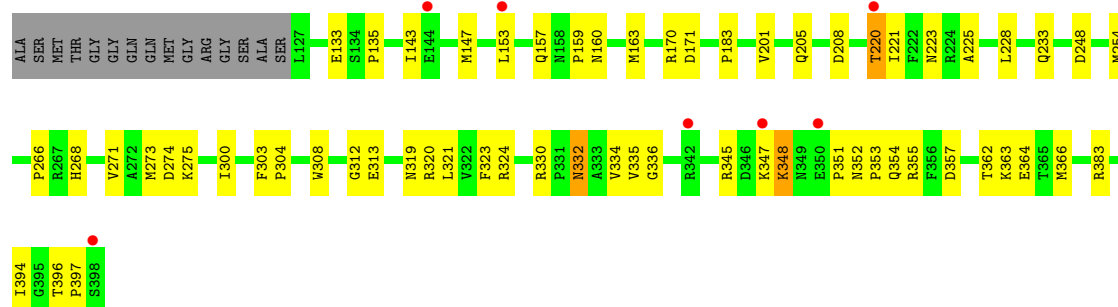
• Molecule 1: Beta-1,4-galactosyltransferase1

Chain B: 



• Molecule 1: Beta-1,4-galactosyltransferase1

Chain C: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	106.98Å 195.17Å 143.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.00 36.67 – 1.98	Depositor EDS
% Data completeness (in resolution range)	96.5 (40.00-2.00) 96.3 (36.67-1.98)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 1.98Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.219 , 0.246 0.216 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	26.3	Xtriage
Anisotropy	0.737	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 41.2	EDS
Estimated twinning fraction	0.000 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.020 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Outliers	0 of 101271 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7413	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.12% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, DIO, NAG, MN, SO4, UDH, MAN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.44	0/2274	0.71	2/3089 (0.1%)
1	B	0.46	0/2274	0.70	1/3089 (0.0%)
1	C	0.42	0/2274	0.66	1/3089 (0.0%)
All	All	0.44	0/6822	0.69	4/9267 (0.0%)

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
3	B	1	0

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	208	ASP	N-CA-C	-5.80	95.33	111.00
1	C	208	ASP	N-CA-C	-5.30	96.69	111.00
1	A	311	GLY	N-CA-C	5.11	125.88	113.10
1	A	208	ASP	N-CA-C	-5.03	97.41	111.00

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	B	409	MAN	C1

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2212	0	2157	32	0
1	B	2212	0	2157	32	0
1	C	2212	0	2157	52	0
2	A	26	0	24	0	0
2	C	26	0	24	0	0
3	B	37	0	33	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
5	A	20	0	0	0	0
5	B	30	0	0	0	0
5	C	20	0	0	0	0
6	A	32	0	25	0	0
6	B	32	0	25	4	0
6	C	32	0	25	2	0
7	C	6	0	8	2	0
8	A	6	0	8	1	0
8	B	30	0	40	6	0
8	C	6	0	8	1	0
9	A	157	0	0	0	0
9	B	192	0	0	3	0
9	C	122	0	0	1	0
All	All	7413	0	6691	113	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 8.

All (113) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:337:THR:HG21	1:C:135:PRO:HB2	1.57	0.86
1:B:324:ARG:HH12	1:B:367:LEU:HD21	1.43	0.84
1:C:332:ASN:ND2	1:C:335:VAL:H	1.82	0.78
1:B:324:ARG:NH1	1:B:367:LEU:HD21	1.98	0.78
1:A:332:ASN:HD22	1:A:334:VAL:H	1.31	0.77

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:320:ARG:O	1:C:324:ARG:HG2	1.86	0.75
1:A:150:ASP:OD1	1:A:153:LEU:HD13	1.89	0.73
1:A:254:MET:HE3	1:A:337:THR:HG22	1.71	0.73
1:C:300:ILE:HG22	1:C:324:ARG:HG3	1.72	0.72
1:C:153:LEU:O	1:C:157:GLN:HG3	1.90	0.71
1:C:332:ASN:HD21	1:C:335:VAL:H	1.38	0.71
1:A:273:MET:CE	1:A:275:LYS:HE2	2.21	0.70
1:A:254:MET:HE3	1:A:337:THR:CG2	2.23	0.69
1:B:149:VAL:HA	8:B:418:GOL:H32	1.74	0.69
1:B:186:ASN:ND2	8:B:414:GOL:H32	2.10	0.67
1:B:273:MET:CE	1:B:275:LYS:HE2	2.24	0.67
1:C:300:ILE:HG22	1:C:324:ARG:CG	2.25	0.66
1:A:332:ASN:ND2	1:A:335:VAL:H	1.93	0.66
1:C:308:TRP:HE1	7:C:433:DIO:H1'1	1.64	0.63
1:B:268:HIS:HB3	1:B:330:ARG:HG2	1.79	0.63
1:B:342:ARG:HE	1:C:170:ARG:NH1	1.96	0.63
1:A:332:ASN:HD21	1:A:335:VAL:H	1.45	0.62
1:A:268:HIS:HB3	1:A:330:ARG:HG2	1.82	0.61
1:B:351:PRO:HG3	6:B:401:UDH:H6'1	1.83	0.61
1:A:273:MET:HE1	1:A:275:LYS:HE2	1.82	0.61
1:C:266:PRO:HG2	1:C:321:LEU:HD22	1.82	0.60
1:C:345:ARG:HG2	6:C:403:UDH:H4'2	1.83	0.60
1:A:157:GLN:C	1:A:159:PRO:HD3	2.22	0.59
1:A:139:GLY:HA2	1:A:257:HIS:HD2	1.68	0.58
1:B:306:ASN:HD22	1:B:306:ASN:H	1.50	0.58
1:A:332:ASN:HD22	1:A:334:VAL:N	2.01	0.58
1:C:273:MET:CE	1:C:275:LYS:HE2	2.34	0.58
1:C:268:HIS:HB3	1:C:330:ARG:HG2	1.86	0.57
1:A:254:MET:CE	1:A:337:THR:HG22	2.34	0.57
8:B:412:GOL:H31	9:B:445:HOH:O	2.06	0.56
1:C:323:PHE:CE1	1:C:363:LYS:HB2	2.41	0.55
1:B:254:MET:HB3	9:B:568:HOH:O	2.07	0.55
1:B:156:LYS:N	1:B:156:LYS:HE2	2.21	0.55
1:B:345:ARG:HE	6:B:401:UDH:H5B	1.73	0.54
1:C:274:ASP:OD1	1:C:275:LYS:N	2.41	0.54
1:C:160:ASN:OD1	1:C:170:ARG:NH2	2.41	0.53
1:C:157:GLN:C	1:C:159:PRO:HD3	2.28	0.53
1:B:310:TRP:CD1	8:B:412:GOL:H12	2.44	0.53
1:C:271:VAL:CG1	1:C:330:ARG:HD3	2.40	0.52
1:A:398:SER:OXT	1:B:348:LYS:HG3	2.08	0.52
1:B:271:VAL:HG22	1:B:336:GLY:HA3	1.90	0.52
1:A:332:ASN:ND2	1:A:334:VAL:H	2.03	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:306:ASN:ND2	1:B:306:ASN:H	2.08	0.51
1:A:352:ASN:OD1	1:A:354:GLN:HG2	2.11	0.51
1:A:254:MET:CE	1:A:339:ARG:HG2	2.42	0.50
1:C:332:ASN:HD22	1:C:334:VAL:N	2.10	0.50
1:B:306:ASN:HD22	1:B:306:ASN:N	2.08	0.50
1:C:352:ASN:OD1	1:C:354:GLN:HG2	2.12	0.49
1:C:271:VAL:HG12	1:C:330:ARG:HD3	1.95	0.48
1:C:273:MET:HE1	1:C:275:LYS:HE2	1.95	0.48
1:B:273:MET:HE1	1:B:275:LYS:HE2	1.94	0.48
1:A:139:GLY:HA2	1:A:257:HIS:CD2	2.47	0.48
1:C:271:VAL:HG22	1:C:336:GLY:HA3	1.94	0.48
1:C:324:ARG:HD3	1:C:366:MET:CE	2.44	0.48
1:B:306:ASN:HB2	8:B:417:GOL:H2	1.94	0.48
1:C:273:MET:HE2	1:C:275:LYS:HE2	1.95	0.48
1:C:347:LYS:O	1:C:348:LYS:HB2	2.13	0.48
1:A:303:PHE:HB3	1:A:304:PRO:HD2	1.96	0.48
1:B:345:ARG:NE	6:B:401:UDH:H5B	2.29	0.47
1:C:271:VAL:CG2	1:C:336:GLY:HA3	2.44	0.47
1:B:274:ASP:O	1:C:171:ASP:HB3	2.14	0.47
1:C:332:ASN:HD22	1:C:334:VAL:H	1.63	0.47
1:C:304:PRO:HB3	1:C:320:ARG:HH21	1.79	0.47
1:B:310:TRP:NE1	8:B:412:GOL:H12	2.30	0.47
1:A:139:GLY:HA3	1:A:140:PRO:HD3	1.81	0.46
1:A:271:VAL:HG22	1:A:336:GLY:HA3	1.97	0.46
1:C:300:ILE:HG22	1:C:324:ARG:HG2	1.98	0.46
1:C:248:ASP:HB3	6:C:403:UDH:O3'	2.15	0.46
1:A:398:SER:HB2	1:B:347:LYS:HB2	1.98	0.46
1:C:396:THR:HB	1:C:397:PRO:HD2	1.98	0.45
1:A:254:MET:HE1	1:A:339:ARG:HG2	1.98	0.45
1:C:163:MET:CE	1:C:383:ARG:O	2.65	0.45
1:A:153:LEU:O	1:A:157:GLN:HG3	2.16	0.45
1:A:266:PRO:HG2	1:A:321:LEU:HD22	1.99	0.45
1:A:194:TRP:CZ2	1:A:198:LEU:HG	2.52	0.44
8:C:416:GOL:H31	9:C:829:HOH:O	2.17	0.44
1:C:266:PRO:CG	1:C:321:LEU:HD22	2.45	0.44
1:C:352:ASN:O	1:C:355:ARG:HG2	2.17	0.44
1:C:351:PRO:O	1:C:353:PRO:HD3	2.16	0.44
1:C:303:PHE:HB3	1:C:304:PRO:HD2	2.00	0.44
1:A:254:MET:SD	1:A:337:THR:HG22	2.57	0.44
1:A:273:MET:HE2	1:A:275:LYS:HE2	1.98	0.43
1:B:334:VAL:O	1:B:337:THR:HG23	2.18	0.43
1:C:223:ASN:HB3	1:C:394:ILE:CD1	2.48	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:220:THR:HG23	1:C:221:ILE:N	2.33	0.43
1:C:183:PRO:HD3	1:C:228:LEU:HD21	2.01	0.43
1:C:324:ARG:HD3	1:C:366:MET:HE2	2.00	0.43
1:B:303:PHE:HB3	1:B:304:PRO:HD2	2.00	0.43
1:B:272:ALA:HB3	1:B:337:THR:HA	2.01	0.42
1:C:271:VAL:HG11	1:C:330:ARG:HB3	2.01	0.42
1:B:314:ASP:OD1	1:B:314:ASP:N	2.52	0.42
1:C:319:ASN:HB2	1:C:362:THR:HG21	2.00	0.42
1:C:273:MET:HE2	1:C:275:LYS:HG2	2.02	0.42
1:A:310:TRP:CE2	8:A:415:GOL:H12	2.55	0.42
1:C:163:MET:HE3	1:C:383:ARG:O	2.19	0.42
1:B:351:PRO:CG	6:B:401:UDH:H6'1	2.49	0.41
1:C:308:TRP:NE1	7:C:433:DIO:H1'1	2.33	0.41
1:A:158:ASN:N	1:A:159:PRO:HD3	2.36	0.41
1:C:143:ILE:HG13	1:C:254:MET:HG2	2.03	0.41
1:C:201:VAL:O	1:C:205:GLN:HG3	2.21	0.41
1:C:345:ARG:HG3	1:C:345:ARG:HH11	1.86	0.41
1:A:223:ASN:HB3	1:A:394:ILE:CD1	2.51	0.41
1:B:306:ASN:ND2	1:B:306:ASN:N	2.68	0.40
1:C:220:THR:CG2	1:C:221:ILE:N	2.84	0.40
1:B:300:ILE:HB	1:B:320:ARG:HB3	2.03	0.40
1:C:225:ALA:HB2	1:C:313:GLU:HG3	2.03	0.40
1:B:157:GLN:NE2	9:B:492:HOH:O	2.53	0.40
1:A:177:LYS:NZ	1:A:240:ASP:OD2	2.55	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	270/287 (94%)	263 (97%)	7 (3%)	0	100	100
1	B	270/287 (94%)	264 (98%)	6 (2%)	0	100	100
1	C	270/287 (94%)	261 (97%)	8 (3%)	1 (0%)	43	36

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	810/861 (94%)	788 (97%)	21 (3%)	1 (0%)	59	55

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	312	GLY

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	243/252 (96%)	240 (99%)	3 (1%)	82	84
1	B	243/252 (96%)	238 (98%)	5 (2%)	66	67
1	C	243/252 (96%)	235 (97%)	8 (3%)	50	46
All	All	729/756 (96%)	713 (98%)	16 (2%)	64	65

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	147	MET
1	A	332	ASN
1	A	364	GLU
1	B	147	MET
1	B	153	LEU
1	B	156	LYS
1	B	306	ASN
1	B	350	GLU
1	C	133	GLU
1	C	147	MET
1	C	220	THR
1	C	233	GLN
1	C	332	ASN
1	C	348	LYS
1	C	357	ASP
1	C	364	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	157	GLN
1	A	186	ASN
1	A	295	GLN
1	A	332	ASN
1	B	186	ASN
1	B	306	ASN
1	C	295	GLN
1	C	332	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

7 carbohydrates 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	NAG	A	405	2	12,14,15	1.27	2 (16%)	15,19,21	0.96	0
2	MAN	A	406	2	12,12,12	1.64	3 (25%)	17,17,17	0.93	0
3	NAG	B	407	3	12,14,15	1.54	2 (16%)	15,19,21	1.16	2 (13%)
3	MAN	B	408	3	10,11,12	2.02	5 (50%)	11,15,17	1.09	1 (9%)
3	MAN	B	409	3	12,12,12	2.59	5 (41%)	17,17,17	0.74	0
2	NAG	C	410	2	12,14,15	1.29	1 (8%)	15,19,21	0.96	0
2	MAN	C	411	2	12,12,12	1.61	2 (16%)	17,17,17	0.80	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	NAG	A	405	2	-	0/6/23/26	0/1/1/1
2	MAN	A	406	2	-	0/2/22/22	0/1/1/1
3	NAG	B	407	3	-	0/6/23/26	0/1/1/1
3	MAN	B	408	3	-	0/2/19/22	0/1/1/1
3	MAN	B	409	3	1/1/5/5	0/2/22/22	0/1/1/1
2	NAG	C	410	2	-	0/6/23/26	0/1/1/1
2	MAN	C	411	2	-	0/2/22/22	0/1/1/1

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	409	MAN	C4-C5	4.74	1.63	1.53
3	B	409	MAN	C1-C2	4.57	1.61	1.52
3	B	409	MAN	O5-C1	3.59	1.50	1.43
3	B	408	MAN	C3-C2	3.37	1.60	1.52
2	A	405	NAG	C2-N2	3.26	1.50	1.46
2	C	410	NAG	C2-N2	3.18	1.50	1.46
3	B	407	NAG	C4-C5	3.09	1.59	1.53
2	A	406	MAN	C1-C2	3.08	1.58	1.52
2	C	411	MAN	O5-C1	3.06	1.49	1.43
3	B	408	MAN	O2-C2	2.97	1.50	1.43
3	B	409	MAN	C3-C2	2.96	1.60	1.52
3	B	407	NAG	C2-N2	2.67	1.49	1.46
3	B	409	MAN	O5-C5	2.61	1.50	1.44
3	B	408	MAN	C4-C5	2.56	1.58	1.53
3	B	408	MAN	O5-C5	2.51	1.50	1.45
3	B	408	MAN	C4-C3	2.46	1.58	1.52
2	A	406	MAN	C4-C5	2.44	1.58	1.53
2	A	406	MAN	O5-C1	2.36	1.47	1.43
2	C	411	MAN	C1-C2	2.23	1.57	1.52
2	A	405	NAG	C8-C7	2.10	1.55	1.50

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	407	NAG	O5-C5-C4	-2.49	107.49	110.65
3	B	407	NAG	O5-C5-C6	2.32	109.42	106.98
3	B	408	MAN	C3-C4-C5	-2.26	106.17	110.20

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	B	409	MAN	C1

There are no torsion outliers.

There are no ring outliers.

5.6 Ligand geometry ⓘ

Of 28 ligands modelled in this entry, 3 are monoatomic - leaving 25 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	$\# Z > 2$
6	UDH	A	399	4	33,33,33	1.60	6 (18%)	43,47,47	1.42	6 (13%)
8	GOL	A	415	-	5,5,5	0.60	0	5,5,5	0.97	0
5	SO4	A	419	-	4,4,4	3.38	2 (50%)	6,6,6	0.93	0
5	SO4	A	420	-	4,4,4	3.33	2 (50%)	6,6,6	0.96	0
5	SO4	A	421	-	4,4,4	3.37	2 (50%)	6,6,6	0.93	0
5	SO4	A	429	-	4,4,4	3.44	2 (50%)	6,6,6	0.92	0
6	UDH	B	401	4	33,33,33	1.61	5 (15%)	43,47,47	1.38	4 (9%)
8	GOL	B	412	-	5,5,5	0.51	0	5,5,5	1.03	0
8	GOL	B	413	-	5,5,5	0.42	0	5,5,5	0.90	0
8	GOL	B	414	-	5,5,5	0.50	0	5,5,5	0.76	0
8	GOL	B	417	-	5,5,5	0.36	0	5,5,5	0.83	0
8	GOL	B	418	-	5,5,5	0.34	0	5,5,5	0.77	0
5	SO4	B	422	-	4,4,4	3.35	2 (50%)	6,6,6	0.98	0
5	SO4	B	424	-	4,4,4	3.34	2 (50%)	6,6,6	0.96	0
5	SO4	B	425	-	4,4,4	3.34	2 (50%)	6,6,6	0.93	0
5	SO4	B	428	-	4,4,4	3.47	2 (50%)	6,6,6	0.91	0
5	SO4	B	430	-	4,4,4	3.41	2 (50%)	6,6,6	0.89	0
5	SO4	B	432	-	4,4,4	3.35	2 (50%)	6,6,6	0.96	0
6	UDH	C	403	4	33,33,33	1.50	7 (21%)	43,47,47	1.37	6 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	GOL	C	416	-	5,5,5	0.44	0	5,5,5	0.84	0
5	SO4	C	423	-	4,4,4	3.30	2 (50%)	6,6,6	0.91	0
5	SO4	C	426	-	4,4,4	3.51	2 (50%)	6,6,6	0.90	0
5	SO4	C	427	-	4,4,4	3.47	2 (50%)	6,6,6	0.87	0
5	SO4	C	431	-	4,4,4	3.46	2 (50%)	6,6,6	0.92	0
7	DIO	C	433	-	6,6,6	0.89	0	6,6,6	0.65	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
6	UDH	A	399	4	-	0/23/41/41	0/2/2/2
8	GOL	A	415	-	-	0/4/4/4	0/0/0/0
5	SO4	A	419	-	-	0/0/0/0	0/0/0/0
5	SO4	A	420	-	-	0/0/0/0	0/0/0/0
5	SO4	A	421	-	-	0/0/0/0	0/0/0/0
5	SO4	A	429	-	-	0/0/0/0	0/0/0/0
6	UDH	B	401	4	-	0/23/41/41	0/2/2/2
8	GOL	B	412	-	-	0/4/4/4	0/0/0/0
8	GOL	B	413	-	-	0/4/4/4	0/0/0/0
8	GOL	B	414	-	-	0/4/4/4	0/0/0/0
8	GOL	B	417	-	-	0/4/4/4	0/0/0/0
8	GOL	B	418	-	-	0/4/4/4	0/0/0/0
5	SO4	B	422	-	-	0/0/0/0	0/0/0/0
5	SO4	B	424	-	-	0/0/0/0	0/0/0/0
5	SO4	B	425	-	-	0/0/0/0	0/0/0/0
5	SO4	B	428	-	-	0/0/0/0	0/0/0/0
5	SO4	B	430	-	-	0/0/0/0	0/0/0/0
5	SO4	B	432	-	-	0/0/0/0	0/0/0/0
6	UDH	C	403	4	-	0/23/41/41	0/2/2/2
8	GOL	C	416	-	-	0/4/4/4	0/0/0/0
5	SO4	C	423	-	-	0/0/0/0	0/0/0/0
5	SO4	C	426	-	-	0/0/0/0	0/0/0/0
5	SO4	C	427	-	-	0/0/0/0	0/0/0/0
5	SO4	C	431	-	-	0/0/0/0	0/0/0/0
7	DIO	C	433	-	-	0/0/6/6	0/1/1/1

All (46) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	427	SO4	O1-S	5.15	1.63	1.47
5	C	431	SO4	O1-S	5.10	1.63	1.47
5	A	429	SO4	O1-S	5.01	1.63	1.47
5	C	426	SO4	O3-S	-4.99	1.30	1.47
5	B	428	SO4	O3-S	-4.85	1.30	1.47
5	B	422	SO4	O1-S	4.84	1.62	1.47
5	C	426	SO4	O1-S	4.83	1.62	1.47
5	B	430	SO4	O1-S	4.80	1.62	1.47
5	A	421	SO4	O1-S	4.79	1.62	1.47
5	B	428	SO4	O1-S	4.76	1.62	1.47
5	B	430	SO4	O3-S	-4.76	1.31	1.47
5	A	419	SO4	O1-S	4.75	1.62	1.47
5	B	432	SO4	O1-S	4.74	1.62	1.47
5	A	420	SO4	O3-S	-4.72	1.31	1.47
5	A	419	SO4	O3-S	-4.71	1.31	1.47
5	C	423	SO4	O3-S	-4.70	1.31	1.47
5	B	424	SO4	O3-S	-4.69	1.31	1.47
5	B	425	SO4	O1-S	4.69	1.62	1.47
5	B	425	SO4	O3-S	-4.68	1.31	1.47
5	A	429	SO4	O3-S	-4.67	1.31	1.47
5	B	432	SO4	O3-S	-4.67	1.31	1.47
5	A	421	SO4	O3-S	-4.66	1.31	1.47
5	B	424	SO4	O1-S	4.64	1.62	1.47
5	C	431	SO4	O3-S	-4.63	1.31	1.47
5	A	420	SO4	O1-S	4.60	1.62	1.47
5	C	427	SO4	O3-S	-4.59	1.31	1.47
5	B	422	SO4	O3-S	-4.57	1.31	1.47
6	A	399	UDH	C6-N1	4.54	1.43	1.35
5	C	423	SO4	O1-S	4.53	1.61	1.47
6	C	403	UDH	C6-N1	4.29	1.42	1.35
6	B	401	UDH	C6-N1	4.21	1.42	1.35
6	B	401	UDH	C2-N1	4.04	1.42	1.38
6	B	401	UDH	PA-O3A	3.59	1.66	1.59
6	A	399	UDH	PA-O3A	3.30	1.65	1.59
6	A	399	UDH	C2-N1	2.84	1.41	1.38
6	C	403	UDH	C2-N1	2.80	1.41	1.38
6	C	403	UDH	C4-N3	2.79	1.41	1.37
6	C	403	UDH	PA-O3A	2.68	1.64	1.59
6	A	399	UDH	C2B-C1B	-2.63	1.49	1.53
6	B	401	UDH	PB-O1B	2.44	1.60	1.51
6	A	399	UDH	C4-N3	2.31	1.41	1.37
6	A	399	UDH	PB-O1B	2.22	1.59	1.51
6	C	403	UDH	PB-O2B	-2.16	1.49	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	C	403	UDH	PB-O1B	2.16	1.59	1.51
6	B	401	UDH	C4-N3	2.15	1.40	1.37
6	C	403	UDH	C6-C5	-2.02	1.32	1.36

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	401	UDH	C2-N1-C1B	-4.30	115.51	118.21
6	A	399	UDH	O4'-C4B-C3B	-3.38	98.31	105.17
6	B	401	UDH	O4'-C4B-C3B	-3.21	98.67	105.17
6	C	403	UDH	O4'-C4B-C3B	-2.92	99.26	105.17
6	A	399	UDH	C2-N1-C1B	-2.87	116.41	118.21
6	C	403	UDH	O3B-PB-O1B	2.62	126.86	112.21
6	A	399	UDH	O3B-PB-O1B	2.54	126.38	112.21
6	A	399	UDH	C2B-C1B-N1	2.48	119.64	113.26
6	C	403	UDH	PB-O3A-PA	-2.45	124.50	131.68
6	C	403	UDH	C2B-C1B-N1	2.36	119.34	113.26
6	B	401	UDH	PB-O3A-PA	-2.35	124.79	131.68
6	B	401	UDH	O3B-PB-O1B	2.34	125.28	112.21
6	C	403	UDH	C5-C6-N1	2.15	123.65	121.21
6	A	399	UDH	C5'-C4'-C3'	-2.13	103.07	114.61
6	A	399	UDH	C3'-C2'-C1'	-2.11	103.34	113.58
6	C	403	UDH	C5'-C4'-C3'	-2.09	103.31	114.61

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

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	272/287 (94%)	-0.09	3 (1%) 77 78	20, 28, 45, 54	0
1	B	272/287 (94%)	-0.08	4 (1%) 70 70	17, 25, 38, 51	0
1	C	272/287 (94%)	0.13	7 (2%) 53 53	22, 35, 56, 65	0
All	All	816/861 (94%)	-0.02	14 (1%) 62 67	17, 29, 49, 65	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	398	SER	4.5
1	C	347	LYS	3.4
1	B	398	SER	3.0
1	A	398	SER	2.7
1	A	347	LYS	2.6
1	B	367	LEU	2.5
1	C	220	THR	2.3
1	A	153	LEU	2.3
1	B	132	GLU	2.2
1	B	181	ILE	2.1
1	C	144	GLU	2.0
1	C	153	LEU	2.0
1	C	342	ARG	2.0
1	C	350	GLU	2.0

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

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

6.3 Carbohydrates ⓘ

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	MAN	B	409	12/12	0.47	30.28	49,57,57,58	12
3	MAN	B	408	11/12	0.20	8.47	34,46,50,55	0
2	MAN	C	411	12/12	0.23	1.80	54,61,64,66	0
2	MAN	A	406	12/12	0.19	1.70	39,51,54,56	0
2	NAG	A	405	14/15	0.10	-0.16	26,30,32,33	0
2	NAG	C	410	14/15	0.11	-0.25	45,48,49,51	0
3	NAG	B	407	14/15	0.09	-0.87	24,26,27,29	0

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
8	GOL	B	413	6/6	0.34	14.44	41,44,48,51	0
5	SO4	B	422	5/5	0.18	9.90	73,73,74,74	0
5	SO4	A	421	5/5	0.18	9.05	84,85,85,85	0
8	GOL	B	414	6/6	0.20	4.51	31,35,37,37	0
8	GOL	B	418	6/6	0.27	3.33	62,65,65,65	0
8	GOL	A	415	6/6	0.16	2.30	35,40,41,45	0
5	SO4	A	429	5/5	0.27	2.26	72,72,73,74	0
8	GOL	C	416	6/6	0.17	2.12	45,47,48,50	0
5	SO4	C	431	5/5	0.46	2.03	88,88,89,90	0
5	SO4	B	430	5/5	0.25	1.83	81,81,81,82	0
5	SO4	B	432	5/5	0.35	1.79	89,89,89,90	0
8	GOL	B	417	6/6	0.22	1.14	47,49,50,53	0
5	SO4	B	425	5/5	0.20	0.87	65,67,67,68	0
8	GOL	B	412	6/6	0.16	0.73	32,38,39,40	0
5	SO4	B	428	5/5	0.16	0.59	81,81,81,82	0
5	SO4	A	420	5/5	0.18	0.55	71,73,73,74	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
7	DIO	C	433	6/6	0.27	0.48	70,70,70,71	0
5	SO4	C	423	5/5	0.14	0.47	48,50,51,52	0
5	SO4	C	427	5/5	0.19	0.05	73,73,75,75	0
5	SO4	B	424	5/5	0.15	-0.07	80,80,80,81	0
6	UDH	B	401	32/32	0.15	-0.20	18,22,51,55	0
6	UDH	A	399	32/32	0.12	-0.29	26,30,53,56	0
6	UDH	C	403	32/32	0.14	-0.38	30,38,56,60	0
4	MN	A	400	1/1	0.10	-0.72	31,31,31,31	0
5	SO4	A	419	5/5	0.12	-0.82	79,79,80,80	0
5	SO4	C	426	5/5	0.13	-0.96	88,88,88,88	0
4	MN	B	402	1/1	0.10	-1.00	25,25,25,25	0
4	MN	C	404	1/1	0.07	-4.25	38,38,38,38	0

6.5 Other polymers ⓘ

There are no such residues in this entry.