



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:26 AM GMT

PDB ID : 2AES
Title : Crystal Structure of Human M340H-Beta1,4-Galactosyltransferase-I (M340H-B4Gal-T1) in Complex with GlcNAc-beta1,2-Man-alpha1,3-Man-beta-O
R
Authors : Ramasamy, V.; Ramakrishnan, B.; Boeggeman, E.; Ratner, D.M.; Seeberger, P.H.; Qasba, P.K.
Deposited on : 2005-07-23
Resolution : 2.00 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

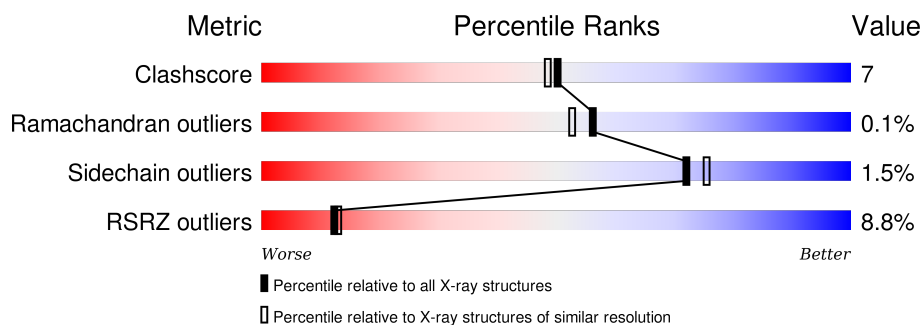
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (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. 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	287	<div> <div>10%</div> <div>79%</div> <div>15%</div> <div>5%</div> </div>
1	B	287	<div> <div>6%</div> <div>83%</div> <div>11%</div> <div>5%</div> </div>
1	C	287	<div> <div>9%</div> <div>76%</div> <div>19%</div> <div>5%</div> </div>

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	SO4	B	429	-	-	-	X
4	SO4	B	434	-	-	-	X
7	MES	C	436	-	-	-	X
8	GOL	A	417	-	-	-	X
8	GOL	A	420	-	-	-	X
8	GOL	A	421	-	-	-	X
8	GOL	B	415	-	-	-	X

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 7401 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 Beta-1,4-galactosyltransferase 1.

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 (3-MER).

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

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

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

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



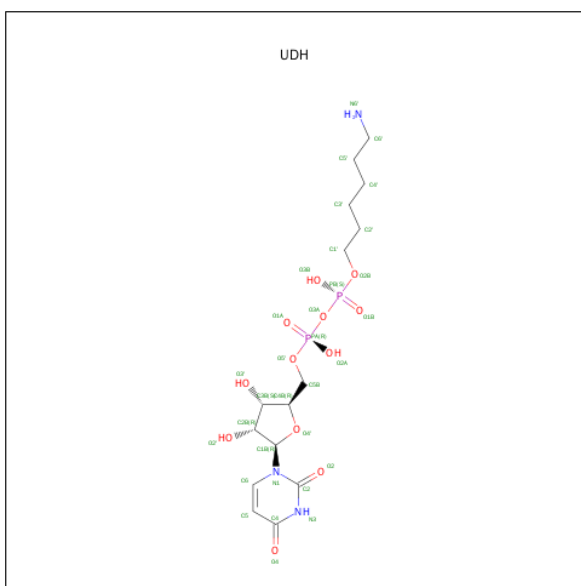
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O S 5 4 1	0	0
4	A	1	Total O S 5 4 1	0	0
4	A	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0
4	C	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0
4	C	1	Total O S 5 4 1	0	0

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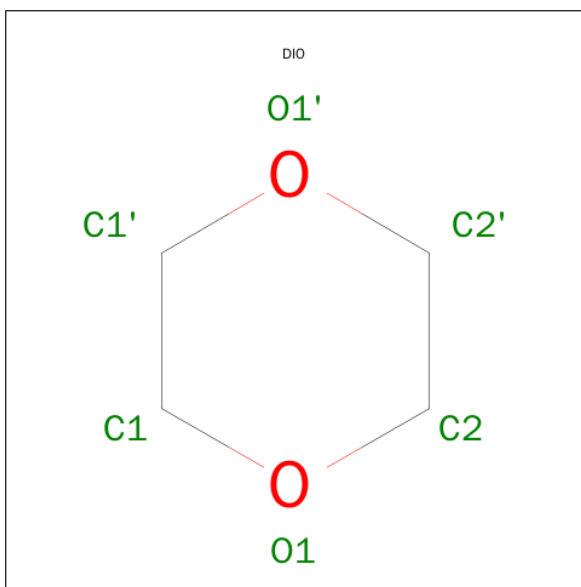
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	C	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		

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



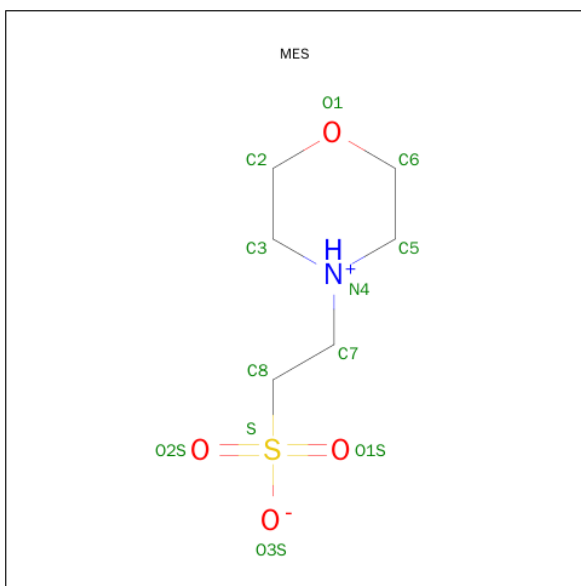
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			32	15	3	12	2		
5	B	1	Total	C	N	O	P	0	0
			32	15	3	12	2		
5	C	1	Total	C	N	O	P	0	0
			32	15	3	12	2		

- Molecule 6 is 1,4-DIETHYLENE DIOXIDE (three-letter code: DIO) (formula: $C_4H_8O_2$).



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

- Molecule 7 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	C	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- 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	A	1	Total	C	O	0	0
			6	3	3		
8	A	1	Total	C	O	0	0
			6	3	3		

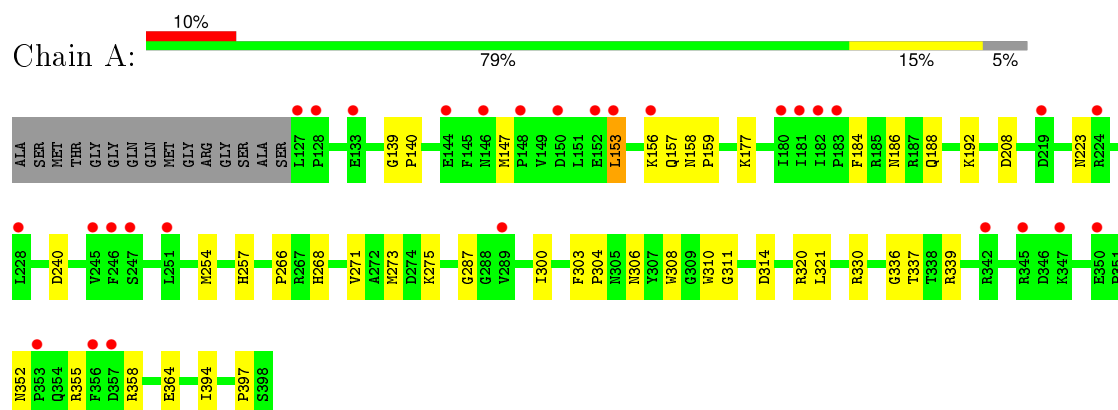
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	142	Total	O	0	0
			142	142		
9	B	157	Total	O	0	0
			157	157		
9	C	125	Total	O	0	0
			125	125		

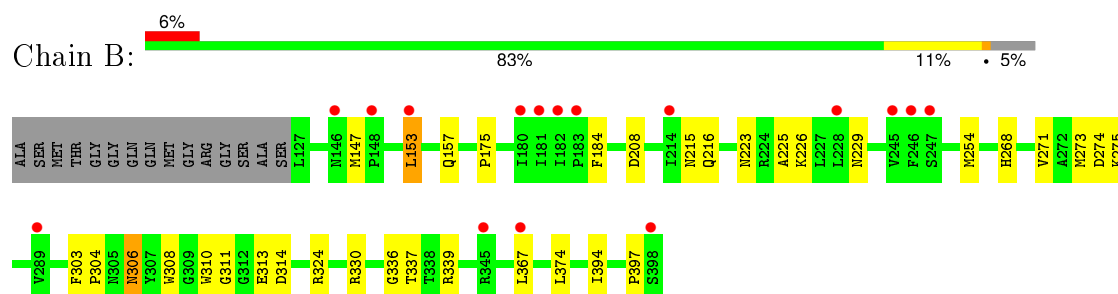
3 Residue-property plots [i](#)

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.

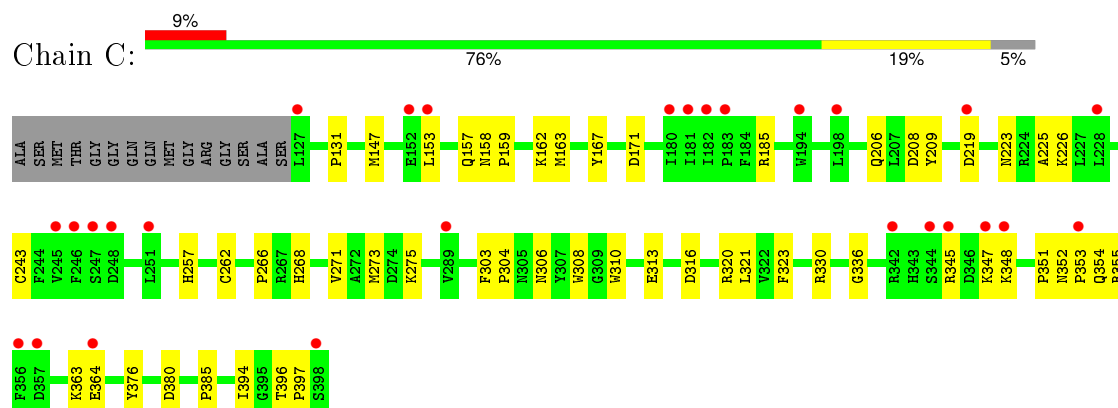
- Molecule 1: Beta-1,4-galactosyltransferase 1



- Molecule 1: Beta-1,4-galactosyltransferase 1



- Molecule 1: Beta-1,4-galactosyltransferase 1



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	107.16Å 192.99Å 143.43Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.00 31.54 – 2.00	Depositor EDS
% Data completeness (in resolution range)	97.5 (40.00-2.00) 97.0 (31.54-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.91 (at 2.00Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.204 , 0.232 0.199 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	29.0	Xtriage
Anisotropy	0.797	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 50.5	EDS
Estimated twinning fraction	0.000 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.017 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 97848 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7401	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 4.03% 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, 0.375 respectively for untwinned datasets, and 0.333, 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, DIO, NAG, MN, BMA, SO4, MES, 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	0.68	2/3089 (0.1%)
1	B	0.47	0/2274	0.69	2/3089 (0.1%)
1	C	0.42	0/2274	0.66	1/3089 (0.0%)
All	All	0.45	0/6822	0.68	5/9267 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	208	ASP	N-CA-C	-5.53	96.06	111.00
1	B	311	GLY	N-CA-C	5.51	126.88	113.10
1	C	208	ASP	N-CA-C	-5.27	96.77	111.00
1	A	208	ASP	N-CA-C	-5.16	97.07	111.00
1	A	311	GLY	N-CA-C	5.13	125.93	113.10

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	2212	0	2157	29	0
1	B	2212	0	2157	30	0
1	C	2212	0	2157	36	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	37	0	33	0	0
2	B	37	0	33	0	0
2	C	37	0	33	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
4	A	20	0	0	0	0
4	B	30	0	0	0	0
4	C	15	0	0	0	0
5	A	32	0	25	1	0
5	B	32	0	25	0	0
5	C	32	0	25	1	0
6	C	6	0	8	2	0
7	C	12	0	13	1	0
8	A	18	0	24	2	0
8	B	24	0	32	4	0
8	C	6	0	8	1	0
9	A	142	0	0	2	0
9	B	157	0	0	3	0
9	C	125	0	0	1	0
All	All	7401	0	6730	96	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 7.

All (96) 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:268:HIS:HD2	1:A:330:ARG:HE	1.22	0.83
1:A:273:MET:HE1	1:A:275:LYS:HE2	1.63	0.81
1:A:273:MET:CE	1:A:275:LYS:HE2	2.12	0.79
1:B:324:ARG:HH11	1:B:367:LEU:HD21	1.50	0.76
1:C:306:ASN:HB3	6:C:435:DIO:H11	1.67	0.76
1:B:273:MET:CE	1:B:275:LYS:HE2	2.16	0.76
1:B:153:LEU:O	1:B:157:GLN:HG3	1.85	0.75
1:B:324:ARG:NH1	1:B:367:LEU:HD21	2.07	0.69
1:C:308:TRP:HE1	6:C:435:DIO:H12	1.57	0.69
1:A:268:HIS:CD2	1:A:330:ARG:HE	2.10	0.68
1:A:254:MET:SD	1:A:337:THR:HG22	2.35	0.67
1:A:153:LEU:O	1:A:157:GLN:HG3	1.97	0.65
1:C:268:HIS:HB3	1:C:330:ARG:HG2	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:266:PRO:HG2	1:C:321:LEU:HD22	1.81	0.61
1:B:268:HIS:HB3	1:B:330:ARG:HG2	1.82	0.61
8:B:414:GOL:H31	9:B:508:HOH:O	2.01	0.59
1:B:303:PHE:HB3	1:B:304:PRO:HD2	1.86	0.58
1:C:153:LEU:O	1:C:157:GLN:HG3	2.03	0.57
1:B:273:MET:HE2	1:B:275:LYS:HE2	1.84	0.57
1:B:306:ASN:HD22	1:B:306:ASN:H	1.50	0.57
1:A:303:PHE:HB3	1:A:304:PRO:HD2	1.88	0.56
1:B:324:ARG:HH11	1:B:367:LEU:CD2	2.19	0.55
1:C:266:PRO:CG	1:C:321:LEU:HD22	2.36	0.55
1:B:271:VAL:HG22	1:B:336:GLY:HA3	1.88	0.55
1:C:323:PHE:CE1	1:C:363:LYS:HB2	2.43	0.53
1:A:268:HIS:HB3	1:A:330:ARG:HG2	1.90	0.53
1:C:352:ASN:OD1	1:C:354:GLN:HG2	2.08	0.53
1:A:157:GLN:C	1:A:159:PRO:HD3	2.29	0.53
1:B:306:ASN:HB2	8:B:419:GOL:H2	1.91	0.53
1:B:273:MET:HE1	1:B:275:LYS:HE2	1.90	0.52
1:C:157:GLN:C	1:C:159:PRO:HD3	2.30	0.52
1:A:306:ASN:ND2	1:A:358:ARG:HE	2.07	0.52
1:A:268:HIS:HE1	1:A:287:GLY:O	1.94	0.51
1:C:273:MET:HE2	1:C:275:LYS:HG2	1.93	0.50
1:C:396:THR:HB	1:C:397:PRO:HD2	1.93	0.50
1:B:306:ASN:ND2	1:B:306:ASN:H	2.09	0.50
1:C:223:ASN:HB3	1:C:394:ILE:CD1	2.41	0.50
1:B:306:ASN:HD22	1:B:306:ASN:N	2.09	0.50
1:C:376:TYR:O	7:C:436:MES:H71	2.12	0.50
1:C:162:LYS:HE2	9:C:589:HOH:O	2.11	0.50
1:C:303:PHE:HB3	1:C:304:PRO:HD2	1.94	0.49
1:C:347:LYS:O	1:C:348:LYS:HB2	2.12	0.49
1:B:254:MET:SD	1:B:337:THR:HG22	2.53	0.48
1:B:216:GLN:OE1	1:B:394:ILE:HD11	2.14	0.47
1:B:254:MET:SD	1:B:339:ARG:HG3	2.54	0.47
1:A:352:ASN:O	1:A:355:ARG:HG2	2.14	0.47
1:C:345:ARG:HG2	5:C:403:UDH:H4'2	1.95	0.47
1:A:158:ASN:N	1:A:159:PRO:HD3	2.29	0.47
1:A:266:PRO:HG2	1:A:321:LEU:HD22	1.96	0.47
1:C:163:MET:CE	1:C:385:PRO:HG3	2.45	0.47
1:C:158:ASN:N	1:C:159:PRO:HD3	2.30	0.46
1:A:177:LYS:NZ	1:A:240:ASP:OD2	2.47	0.46
1:B:175:PRO:HD2	9:B:680:HOH:O	2.15	0.45
1:B:184:PHE:O	1:B:215:ASN:HA	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:308:TRP:O	1:C:352:ASN:HB2	2.16	0.45
1:C:352:ASN:O	1:C:355:ARG:HG2	2.15	0.45
1:C:243:CYS:SG	1:C:262:CYS:N	2.90	0.45
1:A:188:GLN:O	1:A:192:LYS:HG3	2.17	0.45
1:C:219:ASP:O	1:C:348:LYS:HE2	2.17	0.45
1:A:266:PRO:CG	1:A:321:LEU:HD22	2.47	0.44
1:A:257:HIS:HE1	9:A:580:HOH:O	2.00	0.44
8:A:417:GOL:H31	9:A:472:HOH:O	2.17	0.44
1:B:306:ASN:CB	8:B:419:GOL:H2	2.48	0.44
1:B:271:VAL:CG2	1:B:336:GLY:HA3	2.47	0.44
1:A:254:MET:SD	1:A:337:THR:CG2	3.06	0.43
1:C:226:LYS:HD3	1:C:394:ILE:HB	2.00	0.43
1:C:351:PRO:O	1:C:353:PRO:HD3	2.18	0.43
1:A:223:ASN:HB3	1:A:394:ILE:CD1	2.48	0.43
1:C:310:TRP:CE2	8:C:418:GOL:H12	2.53	0.43
1:C:223:ASN:HB3	1:C:394:ILE:HD12	2.00	0.43
1:C:304:PRO:HB3	1:C:320:ARG:HH21	1.84	0.43
1:A:300:ILE:HB	1:A:320:ARG:HB3	1.99	0.43
1:A:184:PHE:HA	5:A:399:UDH:O2	2.18	0.43
1:C:163:MET:HE1	1:C:385:PRO:HG3	2.00	0.43
1:C:225:ALA:HB2	1:C:313:GLU:HG3	2.00	0.43
1:A:308:TRP:CG	1:A:397:PRO:HG3	2.55	0.42
1:C:271:VAL:HG22	1:C:336:GLY:HA3	2.00	0.42
1:A:310:TRP:CE2	8:A:417:GOL:H12	2.54	0.42
1:C:131:PRO:HD2	1:C:206:GLN:HE22	1.84	0.42
1:B:229:ASN:HB3	1:B:374:LEU:HD22	2.02	0.42
1:A:139:GLY:HA3	1:A:140:PRO:HD3	1.91	0.41
1:B:223:ASN:HB3	1:B:394:ILE:HD12	2.02	0.41
1:B:274:ASP:O	1:C:171:ASP:HB3	2.20	0.41
1:B:367:LEU:HB2	9:B:630:HOH:O	2.20	0.41
1:B:226:LYS:HD3	1:B:394:ILE:HB	2.02	0.41
1:A:254:MET:SD	1:A:339:ARG:HG3	2.61	0.41
1:B:314:ASP:OD1	1:B:314:ASP:N	2.53	0.41
1:B:310:TRP:NE1	8:B:414:GOL:H12	2.36	0.41
1:A:271:VAL:HG22	1:A:336:GLY:HA3	2.03	0.40
1:A:156:LYS:O	1:A:159:PRO:HG3	2.21	0.40
1:B:225:ALA:HB2	1:B:313:GLU:HG3	2.03	0.40
1:A:314:ASP:OD1	1:A:314:ASP:N	2.52	0.40
1:C:273:MET:HE2	1:C:275:LYS:CG	2.52	0.40
1:C:167:TYR:HB3	1:C:209:TYR:CE2	2.57	0.40
1:C:316:ASP:O	1:C:320:ARG:HG3	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:308:TRP:CD1	1:B:397:PRO:HB3	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	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%)	262 (97%)	7 (3%)	1 (0%)	39	33
All	All	810/861 (94%)	789 (97%)	20 (2%)	1 (0%)	56	53

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	185	ARG

5.3.2 Protein sidechains [i](#)

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%)	239 (98%)	4 (2%)	70	73
1	B	243/252 (96%)	240 (99%)	3 (1%)	78	81

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	243/252 (96%)	239 (98%)	4 (2%)	70	73
All	All	729/756 (96%)	718 (98%)	11 (2%)	72	75

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	147	MET
1	A	153	LEU
1	A	186	ASN
1	A	364	GLU
1	B	147	MET
1	B	153	LEU
1	B	306	ASN
1	C	147	MET
1	C	257	HIS
1	C	364	GLU
1	C	380	ASP

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	186	ASN
1	A	257	HIS
1	A	268	HIS
1	B	186	ASN
1	B	306	ASN
1	C	206	GLN
1	C	257	HIS
1	C	377	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 ⓘ

9 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	14,14,15	1.60	3 (21%)	15,19,21	1.18	1 (6%)
2	MAN	A	406	2	11,11,12	3.17	4 (36%)	14,15,17	1.56	2 (14%)
2	BMA	A	407	2	12,12,12	2.13	3 (25%)	17,17,17	0.85	0
2	NAG	B	408	2	14,14,15	1.72	5 (35%)	15,19,21	1.11	1 (6%)
2	MAN	B	409	2	11,11,12	3.27	4 (36%)	14,15,17	1.44	2 (14%)
2	BMA	B	410	2	12,12,12	1.98	3 (25%)	17,17,17	0.75	0
2	NAG	C	411	2	14,14,15	1.81	5 (35%)	15,19,21	1.16	1 (6%)
2	MAN	C	412	2	11,11,12	3.14	5 (45%)	14,15,17	1.65	3 (21%)
2	BMA	C	413	2	12,12,12	2.06	4 (33%)	17,17,17	0.72	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	-	0/2/19/22	0/1/1/1
2	BMA	A	407	2	-	0/2/22/22	0/1/1/1
2	NAG	B	408	2	-	0/6/23/26	0/1/1/1
2	MAN	B	409	2	-	0/2/19/22	0/1/1/1
2	BMA	B	410	2	-	0/2/22/22	0/1/1/1
2	NAG	C	411	2	-	0/6/23/26	0/1/1/1
2	MAN	C	412	2	-	0/2/19/22	0/1/1/1
2	BMA	C	413	2	-	0/2/22/22	0/1/1/1

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	413	BMA	C3-C2	2.05	1.57	1.52
2	C	412	MAN	C1-C2	2.11	1.57	1.52
2	B	408	NAG	O5-C1	2.20	1.47	1.43
2	B	408	NAG	C4-C3	2.20	1.58	1.52
2	B	410	BMA	C4-C3	2.33	1.58	1.52
2	B	409	MAN	C4-C5	2.37	1.58	1.53
2	A	405	NAG	C2-N2	2.42	1.50	1.46
2	B	408	NAG	C1-C2	2.46	1.55	1.52
2	C	411	NAG	C4-C3	2.47	1.58	1.52
2	C	411	NAG	C2-N2	2.49	1.50	1.46
2	A	406	MAN	C4-C5	2.53	1.58	1.53
2	B	408	NAG	C2-N2	2.57	1.50	1.46
2	C	413	BMA	C4-C3	2.58	1.59	1.52
2	A	405	NAG	C4-C3	2.61	1.59	1.52
2	C	412	MAN	C4-C5	2.62	1.58	1.53
2	C	411	NAG	O5-C1	2.65	1.48	1.43
2	C	411	NAG	C1-C2	2.70	1.56	1.52
2	A	405	NAG	O5-C5	2.73	1.49	1.43
2	A	407	BMA	C4-C3	2.81	1.59	1.52
2	C	411	NAG	O5-C5	3.37	1.50	1.43
2	B	408	NAG	O5-C5	3.39	1.50	1.43
2	B	410	BMA	O5-C1	3.47	1.49	1.43
2	C	413	BMA	C4-C5	3.73	1.61	1.53
2	A	407	BMA	O5-C1	3.74	1.50	1.43
2	C	413	BMA	O5-C1	3.75	1.50	1.43
2	B	410	BMA	C4-C5	3.90	1.61	1.53
2	A	407	BMA	C4-C5	4.14	1.61	1.53
2	A	406	MAN	O5-C5	4.43	1.53	1.43
2	C	412	MAN	O5-C5	4.47	1.53	1.43
2	B	409	MAN	O5-C5	4.47	1.53	1.43
2	C	412	MAN	C2-C3	5.86	1.60	1.52
2	A	406	MAN	C2-C3	5.98	1.60	1.52
2	C	412	MAN	O5-C1	6.24	1.54	1.43
2	B	409	MAN	O5-C1	6.27	1.54	1.43
2	A	406	MAN	O5-C1	6.29	1.54	1.43
2	B	409	MAN	C2-C3	6.57	1.61	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	405	NAG	C4-C3-C2	-3.42	105.92	111.23
2	B	408	NAG	C4-C3-C2	-3.22	106.23	111.23
2	C	411	NAG	C4-C3-C2	-3.10	106.42	111.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	406	MAN	O2-C2-C3	-2.38	105.33	110.12
2	C	412	MAN	O2-C2-C3	-2.34	105.40	110.12
2	B	409	MAN	O2-C2-C3	-2.15	105.80	110.12
2	C	412	MAN	C1-C2-C3	2.23	112.18	109.54
2	B	409	MAN	C1-O5-C5	3.88	117.18	112.25
2	A	406	MAN	C1-O5-C5	4.06	117.40	112.25
2	C	412	MAN	C1-O5-C5	4.27	117.67	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.6 Ligand geometry [i](#)

Of 29 ligands modelled in this entry, 3 are monoatomic - leaving 26 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	UDH	A	399	3	25,33,33	1.72	3 (12%)	33,47,47	2.72	8 (24%)
8	GOL	A	417	-	5,5,5	0.43	0	5,5,5	0.72	0
8	GOL	A	420	-	5,5,5	0.21	0	5,5,5	0.62	0
8	GOL	A	421	-	5,5,5	0.37	0	5,5,5	0.60	0
4	SO4	A	422	-	4,4,4	3.24	2 (50%)	6,6,6	0.98	0
4	SO4	A	423	-	4,4,4	3.19	2 (50%)	6,6,6	1.01	0
4	SO4	A	424	-	4,4,4	3.27	2 (50%)	6,6,6	0.97	0
4	SO4	A	433	-	4,4,4	3.27	2 (50%)	6,6,6	0.97	0
5	UDH	B	401	3	25,33,33	1.70	4 (16%)	33,47,47	2.65	9 (27%)
8	GOL	B	414	-	5,5,5	0.38	0	5,5,5	0.70	0
8	GOL	B	415	-	5,5,5	0.30	0	5,5,5	0.62	0
8	GOL	B	416	-	5,5,5	0.39	0	5,5,5	0.49	0
8	GOL	B	419	-	5,5,5	0.35	0	5,5,5	0.62	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SO4	B	425	-	4,4,4	3.25	2 (50%)	6,6,6	1.00	0
4	SO4	B	426	-	4,4,4	3.23	2 (50%)	6,6,6	0.95	0
4	SO4	B	428	-	4,4,4	3.17	2 (50%)	6,6,6	1.02	0
4	SO4	B	429	-	4,4,4	3.19	2 (50%)	6,6,6	0.97	0
4	SO4	B	432	-	4,4,4	3.28	2 (50%)	6,6,6	0.96	0
4	SO4	B	434	-	4,4,4	3.32	2 (50%)	6,6,6	0.97	0
5	UDH	C	403	3	25,33,33	1.69	4 (16%)	33,47,47	2.68	7 (21%)
8	GOL	C	418	-	5,5,5	0.26	0	5,5,5	0.64	0
4	SO4	C	427	-	4,4,4	3.18	2 (50%)	6,6,6	0.89	0
4	SO4	C	430	-	4,4,4	3.24	2 (50%)	6,6,6	0.93	0
4	SO4	C	431	-	4,4,4	3.32	2 (50%)	6,6,6	0.93	0
6	DIO	C	435	-	6,6,6	0.88	0	6,6,6	0.69	0
7	MES	C	436	-	11,12,12	1.47	2 (18%)	14,16,16	0.81	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	UDH	A	399	3	-	0/21/41/41	0/2/2/2
8	GOL	A	417	-	-	0/4/4/4	0/0/0/0
8	GOL	A	420	-	-	0/4/4/4	0/0/0/0
8	GOL	A	421	-	-	0/4/4/4	0/0/0/0
4	SO4	A	422	-	-	0/0/0/0	0/0/0/0
4	SO4	A	423	-	-	0/0/0/0	0/0/0/0
4	SO4	A	424	-	-	0/0/0/0	0/0/0/0
4	SO4	A	433	-	-	0/0/0/0	0/0/0/0
5	UDH	B	401	3	-	0/21/41/41	0/2/2/2
8	GOL	B	414	-	-	0/4/4/4	0/0/0/0
8	GOL	B	415	-	-	0/4/4/4	0/0/0/0
8	GOL	B	416	-	-	0/4/4/4	0/0/0/0
8	GOL	B	419	-	-	0/4/4/4	0/0/0/0
4	SO4	B	425	-	-	0/0/0/0	0/0/0/0
4	SO4	B	426	-	-	0/0/0/0	0/0/0/0
4	SO4	B	428	-	-	0/0/0/0	0/0/0/0
4	SO4	B	429	-	-	0/0/0/0	0/0/0/0
4	SO4	B	432	-	-	0/0/0/0	0/0/0/0
4	SO4	B	434	-	-	0/0/0/0	0/0/0/0
5	UDH	C	403	3	-	0/21/41/41	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	GOL	C	418	-	-	0/4/4/4	0/0/0/0
4	SO4	C	427	-	-	0/0/0/0	0/0/0/0
4	SO4	C	430	-	-	0/0/0/0	0/0/0/0
4	SO4	C	431	-	-	0/0/0/0	0/0/0/0
6	DIO	C	435	-	-	0/0/6/6	0/1/1/1
7	MES	C	436	-	-	0/6/14/14	0/1/1/1

All (39) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	430	SO4	O3-S	-4.66	1.30	1.47
4	B	434	SO4	O3-S	-4.58	1.30	1.47
4	B	432	SO4	O3-S	-4.58	1.30	1.47
4	A	433	SO4	O3-S	-4.56	1.31	1.47
4	C	431	SO4	O3-S	-4.52	1.31	1.47
4	A	422	SO4	O3-S	-4.46	1.31	1.47
4	B	425	SO4	O3-S	-4.45	1.31	1.47
4	A	424	SO4	O3-S	-4.45	1.31	1.47
4	B	428	SO4	O3-S	-4.43	1.31	1.47
4	A	423	SO4	O3-S	-4.40	1.31	1.47
4	B	429	SO4	O3-S	-4.37	1.31	1.47
4	C	427	SO4	O3-S	-4.35	1.31	1.47
4	B	426	SO4	O3-S	-3.87	1.33	1.47
5	C	403	UDH	C6-C5	-2.13	1.33	1.38
5	B	401	UDH	PB-O2B	-2.05	1.49	1.59
5	C	403	UDH	PB-O1B	2.21	1.59	1.51
5	A	399	UDH	PB-O1B	2.27	1.59	1.51
7	C	436	MES	C5-N4	2.60	1.54	1.46
5	B	401	UDH	PB-O1B	2.63	1.60	1.51
7	C	436	MES	C3-N4	2.74	1.54	1.46
4	C	430	SO4	O1-S	4.40	1.62	1.47
4	B	428	SO4	O1-S	4.45	1.62	1.47
5	C	403	UDH	C4-N3	4.46	1.41	1.33
4	C	427	SO4	O1-S	4.53	1.62	1.47
4	A	423	SO4	O1-S	4.54	1.62	1.47
4	B	429	SO4	O1-S	4.56	1.62	1.47
4	B	432	SO4	O1-S	4.58	1.62	1.47
4	A	422	SO4	O1-S	4.59	1.62	1.47
4	A	433	SO4	O1-S	4.60	1.62	1.47
5	A	399	UDH	C4-N3	4.62	1.41	1.33
5	B	401	UDH	C4-N3	4.64	1.41	1.33
4	B	425	SO4	O1-S	4.65	1.63	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	401	UDH	C6-N1	4.65	1.42	1.35
4	A	424	SO4	O1-S	4.69	1.63	1.47
4	B	434	SO4	O1-S	4.74	1.63	1.47
4	C	431	SO4	O1-S	4.77	1.63	1.47
5	C	403	UDH	C6-N1	5.03	1.42	1.35
4	B	426	SO4	O1-S	5.16	1.64	1.47
5	A	399	UDH	C6-N1	5.29	1.43	1.35

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	401	UDH	O4'-C4B-C3B	-3.40	98.29	105.15
5	A	399	UDH	O4'-C4B-C3B	-3.38	98.33	105.15
5	B	401	UDH	C5-C4-N3	-3.21	114.89	123.12
5	A	399	UDH	C5-C4-N3	-3.19	114.94	123.12
5	C	403	UDH	C5-C4-N3	-3.14	115.08	123.12
5	C	403	UDH	O4'-C4B-C3B	-3.05	98.99	105.15
5	C	403	UDH	PB-O3A-PA	-2.86	124.70	132.73
5	B	401	UDH	PB-O3A-PA	-2.72	125.10	132.73
5	A	399	UDH	C3'-C2'-C1'	-2.31	103.16	113.47
5	B	401	UDH	C3'-C2'-C1'	-2.24	103.46	113.47
5	A	399	UDH	C4'-C5'-C6'	-2.23	103.50	114.01
5	C	403	UDH	C5'-C4'-C3'	-2.19	103.22	114.53
5	C	403	UDH	C4'-C5'-C6'	-2.18	103.72	114.01
5	B	401	UDH	C5'-C4'-C3'	-2.18	103.28	114.53
5	A	399	UDH	C5'-C4'-C3'	-2.15	103.41	114.53
5	A	399	UDH	C4'-C3'-C2'	-2.07	103.86	114.53
5	B	401	UDH	C4'-C5'-C6'	-2.06	104.29	114.01
5	B	401	UDH	C4'-C3'-C2'	-2.02	104.09	114.53
5	B	401	UDH	O3B-PB-O1B	2.18	124.33	112.53
5	A	399	UDH	O3B-PB-O1B	2.46	125.85	112.53
5	C	403	UDH	O3B-PB-O1B	2.70	127.15	112.53
5	B	401	UDH	C4-N3-C2	12.45	126.47	114.14
5	C	403	UDH	C4-N3-C2	12.59	126.61	114.14
5	A	399	UDH	C4-N3-C2	12.99	127.01	114.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

8 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	399	UDH	1	0
8	A	417	GOL	2	0
8	B	414	GOL	2	0
8	B	419	GOL	2	0
5	C	403	UDH	1	0
8	C	418	GOL	1	0
6	C	435	DIO	2	0
7	C	436	MES	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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%)	0.31	29 (10%) 8 8	23, 32, 49, 60	0
1	B	272/287 (94%)	0.12	16 (5%) 26 27	20, 29, 43, 53	0
1	C	272/287 (94%)	0.37	27 (9%) 9 10	24, 36, 56, 64	0
All	All	816/861 (94%)	0.27	72 (8%) 12 13	20, 32, 51, 64	0

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	398	SER	4.7
1	A	228	LEU	4.0
1	C	182	ILE	3.8
1	A	347	LYS	3.8
1	C	347	LYS	3.7
1	B	398	SER	3.7
1	A	181	ILE	3.6
1	C	181	ILE	3.6
1	C	153	LEU	3.6
1	A	153	LEU	3.4
1	A	180	ILE	3.4
1	C	180	ILE	3.2
1	A	356	PHE	3.2
1	C	228	LEU	3.2
1	A	148	PRO	3.2
1	A	289	VAL	3.2
1	B	182	ILE	3.1
1	A	245	VAL	3.0
1	A	353	PRO	3.0
1	A	152	GLU	3.0
1	C	245	VAL	2.9
1	A	251	LEU	2.8
1	B	183	PRO	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	219	ASP	2.7
1	B	289	VAL	2.7
1	C	183	PRO	2.7
1	B	153	LEU	2.7
1	B	180	ILE	2.6
1	B	245	VAL	2.6
1	C	246	PHE	2.6
1	A	156	LYS	2.6
1	A	350	GLU	2.6
1	A	146	ASN	2.6
1	B	247	SER	2.6
1	A	246	PHE	2.5
1	C	289	VAL	2.5
1	B	228	LEU	2.5
1	A	182	ILE	2.5
1	A	357	ASP	2.5
1	B	181	ILE	2.4
1	C	127	LEU	2.4
1	C	353	PRO	2.4
1	A	219	ASP	2.4
1	C	247	SER	2.4
1	C	356	PHE	2.3
1	B	146	ASN	2.3
1	B	246	PHE	2.3
1	C	344	SER	2.3
1	C	345	ARG	2.3
1	A	342	ARG	2.3
1	A	247	SER	2.2
1	C	198	LEU	2.2
1	B	345	ARG	2.2
1	C	194	TRP	2.2
1	B	367	LEU	2.2
1	C	348	LYS	2.2
1	A	128	PRO	2.2
1	A	345	ARG	2.2
1	A	224	ARG	2.2
1	C	357	ASP	2.2
1	C	152	GLU	2.2
1	A	133	GLU	2.1
1	B	148	PRO	2.1
1	B	214	ILE	2.1
1	C	364	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	127	LEU	2.1
1	C	342	ARG	2.1
1	A	150	ASP	2.1
1	A	144	GLU	2.0
1	C	251	LEU	2.0
1	A	183	PRO	2.0
1	C	248	ASP	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](#)

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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	NAG	C	411	14/15	0.84	0.19	1.40	56,60,61,61	0
2	NAG	A	405	14/15	0.90	0.17	1.13	41,46,48,49	0
2	NAG	B	408	14/15	0.92	0.14	0.77	34,41,43,44	0
2	MAN	C	412	11/12	0.90	0.31	-	64,68,69,69	0
2	BMA	C	413	12/12	0.70	0.42	-	68,69,70,70	11
2	MAN	A	406	11/12	0.87	0.30	-	53,60,61,62	0
2	BMA	B	410	12/12	0.82	0.29	-	52,55,56,57	11
2	MAN	B	409	11/12	0.89	0.28	-	49,54,57,59	0
2	BMA	A	407	12/12	0.78	0.47	-	64,67,70,71	0

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. 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	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
7	MES	C	436	12/12	0.66	0.32	10.80	67,69,78,79	0
8	GOL	B	415	6/6	0.83	0.24	7.14	45,50,53,54	0
4	SO4	B	434	5/5	0.84	0.30	5.70	96,96,97,97	0
8	GOL	A	421	6/6	0.70	0.28	4.05	58,61,61,62	0
8	GOL	A	420	6/6	0.89	0.32	3.39	53,59,60,60	0
4	SO4	B	429	5/5	0.91	0.26	2.34	63,66,67,67	0
8	GOL	A	417	6/6	0.89	0.28	2.04	39,46,47,52	0
8	GOL	B	414	6/6	0.90	0.19	0.98	33,40,43,45	0
8	GOL	B	419	6/6	0.71	0.23	0.96	55,58,58,61	0
8	GOL	C	418	6/6	0.92	0.21	0.92	47,50,51,53	0
6	DIO	C	435	6/6	0.74	0.24	0.60	76,76,77,77	0
4	SO4	A	423	5/5	0.94	0.15	0.46	60,63,63,63	0
4	SO4	C	430	5/5	0.93	0.15	0.00	84,84,85,85	0
4	SO4	C	427	5/5	0.95	0.13	-0.01	49,53,54,55	0
4	SO4	B	428	5/5	0.91	0.16	-0.05	82,83,83,84	0
5	UDH	A	399	32/32	0.94	0.16	-0.21	29,34,55,57	0
5	UDH	B	401	32/32	0.96	0.14	-0.37	20,26,55,57	0
5	UDH	C	403	32/32	0.93	0.16	-0.48	31,37,56,58	0
8	GOL	B	416	6/6	0.96	0.11	-0.48	24,26,29,30	0
4	SO4	A	422	5/5	0.91	0.14	-0.56	90,91,91,91	0
4	SO4	A	424	5/5	0.85	0.30	-	88,88,89,89	0
4	SO4	B	426	5/5	0.85	0.41	-	83,83,84,86	0
3	MN	B	402	1/1	1.00	0.08	-	26,26,26,26	0
3	MN	A	400	1/1	1.00	0.08	-	33,33,33,33	0
4	SO4	C	431	5/5	0.79	0.29	-	95,96,96,96	0
4	SO4	B	432	5/5	0.96	0.24	-	88,88,88,89	0
3	MN	C	404	1/1	0.98	0.07	-	35,35,35,35	0
4	SO4	B	425	5/5	0.89	0.23	-	78,78,78,78	0
4	SO4	A	433	5/5	0.87	0.31	-	98,99,99,99	0

6.5 Other polymers

There are no such residues in this entry.