



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:16 PM GMT

PDB ID : 1O7Q  
Title : Roles of Individual Residues of Alpha-1,3 Galactosyltransferases in Substrate Binding and Catalysis  
Authors : Zhang, Y.; Swaminathan, G.J.; Deshpande, A.; Natesh, R.; Xie, Z.; Acharya, K.R.; Brew, K.  
Deposited on : 2002-11-12  
Resolution : 1.30 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](mailto: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.

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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) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

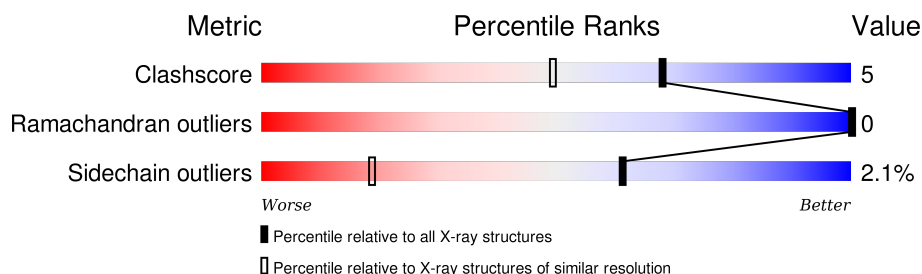
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.30 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	1031 (1.32-1.28)
Ramachandran outliers	100387	1504 (1.34-1.26)
Sidechain outliers	100360	1503 (1.34-1.26)

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  $\geq 3$ , 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	289	
1	B	289	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GOL	B	1378	-	-	X	-
5	UDP	A	1375	-	-	X	-

## 2 Entry composition [i](#)

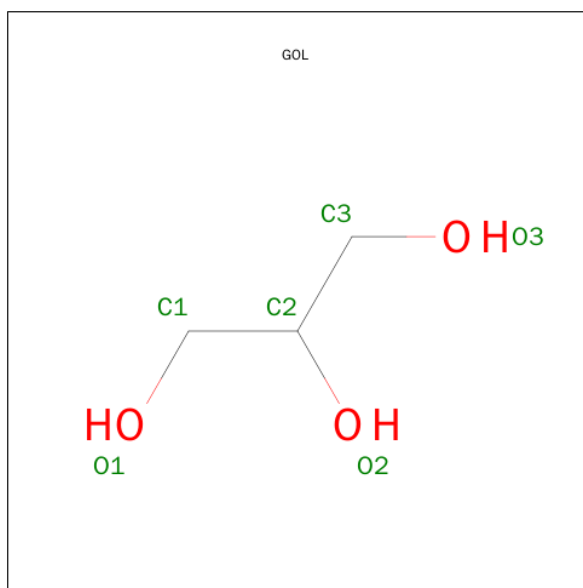
There are 6 unique types of molecules in this entry. The entry contains 10790 atoms, of which 4791 are hydrogens and 0 are deuteriums.

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 N-ACETYLLACTOSAMINIDE ALPHA-1,3-GALACTOSYL TRANSFERASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	287	Total	C	H	N	O	S	0	10	0
			4823	1592	2394	397	424	16			
1	B	287	Total	C	H	N	O	S	0	11	0
			4831	1596	2397	395	427	16			

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

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

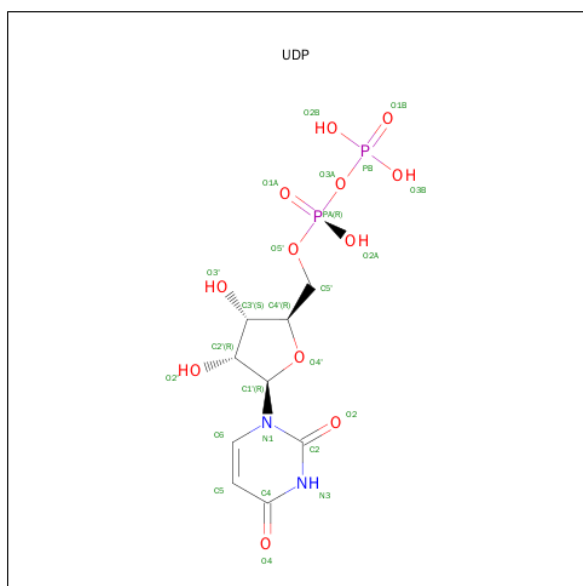
- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

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

- Molecule 4 is a polymer of unknown type called SUGAR (2-MER).

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

- Molecule 5 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: C<sub>9</sub>H<sub>14</sub>N<sub>2</sub>O<sub>12</sub>P<sub>2</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
5	A	1	Total	C	N	O	P	0	0
			25	9	2	12	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
5	B	1	Total	C	N	O	P	0	0
			25	9	2	12	2		

- Molecule 6 is water.

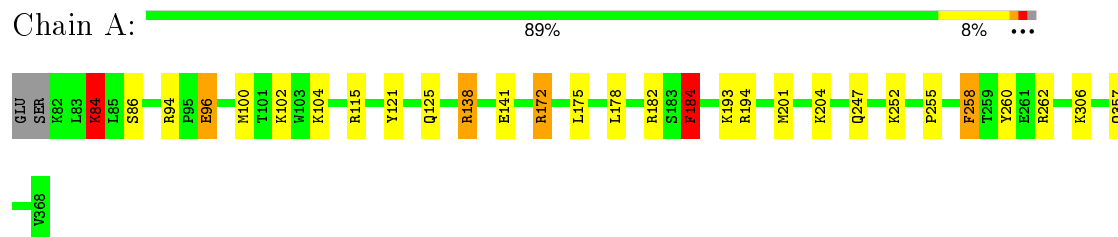
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	468	Total	O	0	0
			468	468		
6	B	484	Total	O	0	0
			484	484		

### 3 Residue-property plots [i](#)

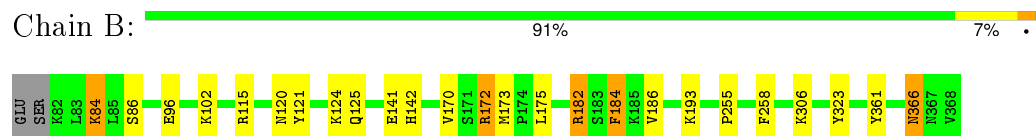
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

- Molecule 1: N-ACETYLLACTOSAMINIDE ALPHA-1,3-GALACTOSYLTRANSFERASE



- Molecule 1: N-ACETYLLACTOSAMINIDE ALPHA-1,3-GALACTOSYLTRANSFERASE



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	45.33 Å 94.68 Å 95.01 Å 90.00° 99.03° 90.00°	Depositor
Resolution (Å)	40.00 – 1.30	Depositor
% Data completeness (in resolution range)	94.9 (40.00-1.30)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	SHELXL-97	Depositor
R, $R_{free}$	0.116 , 0.154	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	10790	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	14.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GAL, GOL, UDP, MN, NAG

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	0.64	2/2544 (0.1%)	1.10	17/3444 (0.5%)
1	B	0.59	0/2554	1.09	9/3457 (0.3%)
All	All	0.62	2/5098 (0.0%)	1.10	26/6901 (0.4%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	96	GLU	CD-OE1	-9.08	1.15	1.25
1	A	96	GLU	CD-OE2	-6.41	1.18	1.25

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	182	ARG	NE-CZ-NH2	9.31	124.95	120.30
1	B	184[A]	PHE	CB-CG-CD1	-9.20	114.36	120.80
1	B	184[B]	PHE	CB-CG-CD1	-9.20	114.36	120.80
1	A	96	GLU	OE1-CD-OE2	-8.86	112.67	123.30
1	B	323	TYR	CB-CG-CD1	8.73	126.24	121.00
1	A	260	TYR	CB-CG-CD2	6.93	125.16	121.00
1	A	184[A]	PHE	CB-CG-CD1	-6.89	115.98	120.80
1	A	184[B]	PHE	CB-CG-CD1	-6.89	115.98	120.80
1	A	138[A]	ARG	CD-NE-CZ	6.57	132.80	123.60
1	A	138[B]	ARG	CD-NE-CZ	6.57	132.80	123.60
1	A	94	ARG	NE-CZ-NH2	-6.54	117.03	120.30
1	A	96	GLU	CG-CD-OE2	6.44	131.19	118.30
1	A	115	ARG	NE-CZ-NH1	6.34	123.47	120.30
1	A	141	GLU	OE1-CD-OE2	-6.28	115.76	123.30
1	A	84	LYS	CD-CE-NZ	6.22	126.00	111.70
1	B	96	GLU	CB-CG-CD	6.05	130.55	114.20
1	B	115	ARG	NE-CZ-NH1	5.82	123.21	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	194	ARG	CD-NE-CZ	5.54	131.35	123.60
1	A	258	PHE	CB-CG-CD2	5.49	124.64	120.80
1	B	361	TYR	CB-CG-CD2	5.39	124.23	121.00
1	A	138[A]	ARG	NE-CZ-NH2	5.18	122.89	120.30
1	A	138[B]	ARG	NE-CZ-NH2	5.18	122.89	120.30
1	A	172	ARG	CA-CB-CG	-5.15	102.08	113.40
1	B	366	ASN	CA-CB-CG	-5.12	102.13	113.40
1	B	172	ARG	NE-CZ-NH1	5.12	122.86	120.30
1	A	172	ARG	CD-NE-CZ	5.05	130.67	123.60

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	2429	2394	2389	18	0
1	B	2434	2397	2392	15	0
2	A	12	0	16	1	0
2	B	18	0	19	5	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	26	0	24	0	0
4	B	26	0	24	0	0
5	A	50	0	22	10	0
5	B	50	0	22	4	0
6	A	468	0	0	8	0
6	B	484	0	0	7	0
All	All	5999	4791	4908	47	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (47) 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)
5:A:1375:UDP:H6	5:A:1375:UDP:C5'	1.84	0.91
5:A:1375:UDP:O5'	5:A:1375:UDP:C6	2.34	0.80
5:A:1375:UDP:C5'	5:A:1375:UDP:C6	2.68	0.77
2:B:1378:GOL:H11	6:B:2009:HOH:O	1.85	0.76
5:A:1375:UDP:C6	5:A:1375:UDP:H3'	2.23	0.74
1:B:86:SER:HB3	2:B:1378:GOL:H12	1.69	0.73
5:A:1375:UDP:H6	5:A:1375:UDP:H5'2	1.55	0.70
1:A:100[A]:MET:HG3	6:A:2045:HOH:O	1.91	0.70
1:A:102:LYS:HE3	6:A:2333:HOH:O	1.92	0.70
5:A:1375:UDP:C6	5:A:1375:UDP:C3'	2.76	0.69
1:B:141:GLU:OE2	1:B:172:ARG:HD2	1.94	0.68
1:A:84:LYS:HD3	1:A:86[A]:SER:OG	1.94	0.68
5:B:1377:UDP:H3'	5:B:1377:UDP:H6	1.61	0.67
1:A:175:LEU:HB2	6:A:2170:HOH:O	1.95	0.65
2:B:1378:GOL:H31	6:B:2009:HOH:O	1.98	0.63
1:A:100[B]:MET:SD	1:A:104:LYS:HA	2.38	0.62
1:B:173[B]:MET:SD	1:B:186:VAL:HG21	2.43	0.58
5:A:1375:UDP:C6	5:A:1375:UDP:C4'	2.89	0.56
1:A:178:LEU:HD11	1:A:184[A]:PHE:CD1	2.41	0.56
1:B:306:LYS:HB3	5:B:1377:UDP:O3'	2.07	0.54
1:B:86:SER:HB3	2:B:1378:GOL:C1	2.37	0.54
5:B:1377:UDP:H3'	5:B:1377:UDP:C6	2.43	0.53
1:A:138[B]:ARG:HD3	6:A:2165:HOH:O	2.08	0.53
1:B:84:LYS:HG3	6:B:2002:HOH:O	2.10	0.52
1:A:193:LYS:HD2	1:A:193:LYS:N	2.25	0.52
5:A:1375:UDP:H6	5:A:1375:UDP:O5'	1.81	0.51
1:A:306:LYS:O	5:A:1375:UDP:O2'	2.29	0.51
1:B:120:ASN:O	1:B:124:LYS:HG3	2.11	0.50
1:A:255:PRO:HA	1:A:258:PHE:CD1	2.48	0.48
1:B:84:LYS:HD3	1:B:86:SER:OG	2.14	0.47
2:B:1379:GOL:H12	6:B:2482:HOH:O	2.14	0.47
1:A:172:ARG:NH2	6:A:2165:HOH:O	2.48	0.46
1:A:201[A]:MET:HE2	6:A:2222:HOH:O	2.15	0.46
1:B:142:HIS:HD2	6:B:2100:HOH:O	1.97	0.46
1:B:121:TYR:O	1:B:125:GLN:HG3	2.16	0.46
1:A:262:ARG:O	2:A:1370:GOL:H11	2.15	0.46
1:A:121:TYR:O	1:A:125:GLN:HG3	2.16	0.46
1:B:193:LYS:NZ	6:B:2221:HOH:O	2.49	0.45
5:A:1375:UDP:O5'	5:A:1375:UDP:C5	2.70	0.44
5:B:1377:UDP:C3'	5:B:1377:UDP:C6	3.01	0.44
1:A:138[A]:ARG:HG2	6:A:2443:HOH:O	2.17	0.43
1:A:204:LYS:HE2	6:A:2241:HOH:O	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:LEU:HD21	1:B:175:LEU:HD21	1.99	0.43
1:B:255:PRO:HA	1:B:258:PHE:CD1	2.53	0.43
1:B:120:ASN:ND2	6:B:2059:HOH:O	2.51	0.42
1:B:170:VAL:HG12	1:B:170:VAL:O	2.20	0.41
1:A:96:GLU:H	1:A:96:GLU:CD	2.25	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	295/289 (102%)	289 (98%)	6 (2%)	0	100	100
1	B	296/289 (102%)	290 (98%)	6 (2%)	0	100	100
All	All	591/578 (102%)	579 (98%)	12 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	267/260 (103%)	260 (97%)	7 (3%)	54	12
1	B	268/260 (103%)	262 (98%)	6 (2%)	60	17
All	All	535/520 (103%)	522 (98%)	13 (2%)	61	14

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	84	LYS
1	A	182	ARG
1	A	184[A]	PHE
1	A	184[B]	PHE
1	A	247	GLN
1	A	252	LYS
1	A	357	GLN
1	B	84	LYS
1	B	102	LYS
1	B	182	ARG
1	B	184[A]	PHE
1	B	184[B]	PHE
1	B	366	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	296	GLN
1	A	313	GLN
1	A	357	GLN
1	B	231	GLN
1	B	296	GLN
1	B	313	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

4 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
4	GAL	A	1372	4	11,11,12	0.85	0	14,15,17	0.58	0
4	NAG	A	1373	4	15,15,15	1.05	1 (6%)	17,21,21	0.77	0
4	GAL	B	1373	4	11,11,12	0.84	0	14,15,17	0.77	0
4	NAG	B	1374	4	15,15,15	1.47	3 (20%)	17,21,21	1.06	2 (11%)

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
4	GAL	A	1372	4	-	0/2/19/22	0/1/1/1
4	NAG	A	1373	4	-	0/6/26/26	0/1/1/1
4	GAL	B	1373	4	-	0/2/19/22	0/1/1/1
4	NAG	B	1374	4	-	0/6/26/26	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1374	NAG	O5-C1	2.18	1.47	1.43
4	A	1373	NAG	O5-C1	2.25	1.47	1.43
4	B	1374	NAG	C2-N2	2.73	1.50	1.45
4	B	1374	NAG	C1-C2	3.24	1.56	1.53

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1374	NAG	C4-C3-C2	-2.20	107.38	110.43
4	B	1374	NAG	O5-C5-C4	-2.00	105.92	109.68

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

Of 11 ligands modelled in this entry, 2 are monoatomic - leaving 9 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
2	GOL	A	1369	-	5,5,5	0.43	0	5,5,5	0.79	0
2	GOL	A	1370	-	5,5,5	1.22	0	5,5,5	1.82	2 (40%)
5	UDP	A	1374	3	18,26,26	1.74	4 (22%)	26,40,40	2.74	4 (15%)
5	UDP	A	1375	-	18,26,26	2.75	6 (33%)	26,40,40	2.91	9 (34%)
2	GOL	B	1372	-	5,5,5	0.33	0	5,5,5	0.87	0
5	UDP	B	1376	3	18,26,26	1.97	4 (22%)	26,40,40	2.69	4 (15%)
5	UDP	B	1377	-	18,26,26	3.12	6 (33%)	26,40,40	2.80	9 (34%)
2	GOL	B	1378	-	5,5,5	2.24	3 (60%)	5,5,5	3.45	2 (40%)
2	GOL	B	1379	-	5,5,5	2.05	2 (40%)	5,5,5	3.52	2 (40%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	A	1369	-	-	0/4/4/4	0/0/0/0
2	GOL	A	1370	-	-	0/4/4/4	0/0/0/0
5	UDP	A	1374	3	-	0/12/32/32	0/2/2/2
5	UDP	A	1375	-	-	0/12/32/32	0/2/2/2
2	GOL	B	1372	-	-	0/4/4/4	0/0/0/0
5	UDP	B	1376	3	-	0/12/32/32	0/2/2/2
5	UDP	B	1377	-	-	0/12/32/32	0/2/2/2
2	GOL	B	1378	-	-	0/4/4/4	0/0/0/0
2	GOL	B	1379	-	-	0/4/4/4	0/0/0/0

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1379	GOL	C3-C2	-3.24	1.39	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1378	GOL	C3-C2	-3.16	1.40	1.52
2	B	1378	GOL	O2-C2	-2.85	1.35	1.43
2	B	1379	GOL	O3-C3	-2.74	1.30	1.42
2	B	1378	GOL	O3-C3	-2.62	1.31	1.42
5	A	1374	UDP	O4'-C1'	2.01	1.43	1.41
5	A	1375	UDP	C5'-C4'	2.06	1.58	1.51
5	B	1377	UDP	C5'-C4'	2.37	1.59	1.51
5	B	1376	UDP	C6-N1	2.63	1.39	1.35
5	B	1377	UDP	O4'-C1'	2.84	1.44	1.41
5	A	1375	UDP	O4'-C1'	2.84	1.44	1.41
5	B	1376	UDP	O4'-C1'	3.28	1.45	1.41
5	A	1374	UDP	C6-C5	3.39	1.45	1.38
5	B	1377	UDP	PB-O1B	3.62	1.63	1.51
5	A	1375	UDP	PB-O1B	3.63	1.63	1.51
5	A	1374	UDP	C6-N1	3.68	1.41	1.35
5	B	1376	UDP	C6-C5	4.02	1.46	1.38
5	A	1374	UDP	C4-N3	4.21	1.40	1.33
5	B	1377	UDP	C4-N3	4.57	1.41	1.33
5	A	1375	UDP	C6-N1	4.97	1.42	1.35
5	B	1376	UDP	C4-N3	5.04	1.42	1.33
5	A	1375	UDP	C4-N3	5.35	1.43	1.33
5	A	1375	UDP	C6-C5	6.68	1.52	1.38
5	B	1377	UDP	C6-N1	6.80	1.45	1.35
5	B	1377	UDP	C6-C5	8.28	1.56	1.38

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1375	UDP	C6-C5-C4	-3.35	111.01	117.28
5	A	1374	UDP	C5-C4-N3	-3.02	115.38	123.12
5	B	1376	UDP	C5-C4-N3	-2.98	115.47	123.12
5	A	1374	UDP	PA-O3A-PB	-2.88	123.01	132.67
5	A	1375	UDP	C5-C4-N3	-2.62	116.39	123.12
5	B	1377	UDP	C5-C4-N3	-2.57	116.52	123.12
5	B	1377	UDP	C6-N1-C2	-2.51	117.22	121.28
5	B	1377	UDP	O4'-C4'-C3'	-2.44	100.23	105.15
5	B	1376	UDP	PA-O3A-PB	-2.33	124.86	132.67
5	B	1377	UDP	C6-C5-C4	-2.29	113.00	117.28
5	A	1375	UDP	O3B-PB-O1B	-2.20	103.49	110.58
5	B	1377	UDP	C2'-C3'-C4'	-2.05	98.41	102.61
5	B	1376	UDP	O3B-PB-O2B	2.03	115.11	107.38
5	A	1375	UDP	O3B-PB-O3A	2.05	114.40	105.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1375	UDP	O2'-C2'-C3'	2.06	118.54	111.83
5	A	1375	UDP	PA-O3A-PB	2.07	139.62	132.67
2	A	1370	GOL	O2-C2-C1	2.21	118.79	108.65
5	A	1374	UDP	O3B-PB-O2B	2.32	116.19	107.38
5	A	1375	UDP	O3B-PB-O2B	2.50	116.92	107.38
5	A	1375	UDP	O2A-PA-O3A	2.61	116.95	105.09
5	B	1377	UDP	O2A-PA-O3A	2.70	117.32	105.09
5	B	1377	UDP	O3B-PB-O2B	2.79	117.99	107.38
2	A	1370	GOL	O2-C2-C3	2.81	121.55	108.65
5	B	1377	UDP	O4'-C1'-N1	2.89	114.17	108.08
2	B	1379	GOL	O2-C2-C3	3.89	126.49	108.65
2	B	1378	GOL	O2-C2-C3	4.33	128.50	108.65
2	B	1378	GOL	O3-C3-C2	6.01	139.34	110.18
2	B	1379	GOL	O3-C3-C2	6.55	141.96	110.18
5	B	1377	UDP	C4-N3-C2	11.69	125.72	114.14
5	A	1375	UDP	C4-N3-C2	12.61	126.62	114.14
5	B	1376	UDP	C4-N3-C2	12.74	126.76	114.14
5	A	1374	UDP	C4-N3-C2	12.82	126.84	114.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1370	GOL	1	0
5	A	1375	UDP	10	0
5	B	1377	UDP	4	0
2	B	1378	GOL	4	0
2	B	1379	GOL	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 [i](#)

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

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.