



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:22 PM UTC

PDB ID : 1VZU / pdb_00001vzu
Title : Roles of active site tryptophans in substrate binding and catalysis by ALPHA-1,3 GALACTOSYLTRANSFERASE
Authors : Zhang, Y.; Deshpande, A.; Xie, Z.; Natesh, R.; Acharya, K.R.; Brew, K.
Deposited on : 2004-05-27
Resolution : 1.97 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

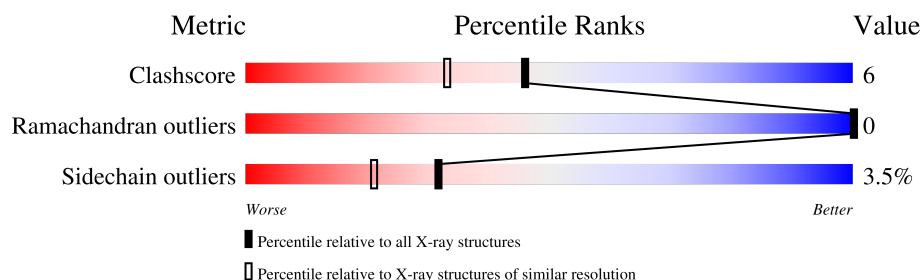
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.97 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	190562	1534 (1.98-1.98)
Ramachandran outliers	187476	1518 (1.98-1.98)
Sidechain outliers	187428	1518 (1.98-1.98)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	289	 80% 17% ..
1	B	289	 80% 18% ..
2	C	2	 50% 50%
2	D	2	 50% 50%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 5788 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 N-ACETYLLACTOSAMINIDE ALPHA-1,3-GALACTOSYL TRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	287	Total	C	N	O	S	0	4	0
			2414	1583	396	423	12			
1	B	287	Total	C	N	O	S	0	2	0
			2401	1571	395	423	12			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	314	TYR	TRP	engineered mutation	UNP P14769
B	1314	TYR	TRP	engineered mutation	UNP P14769

- Molecule 2 is an oligosaccharide called beta-D-galactopyranose-(1-4)-beta-D-glucopyranose.

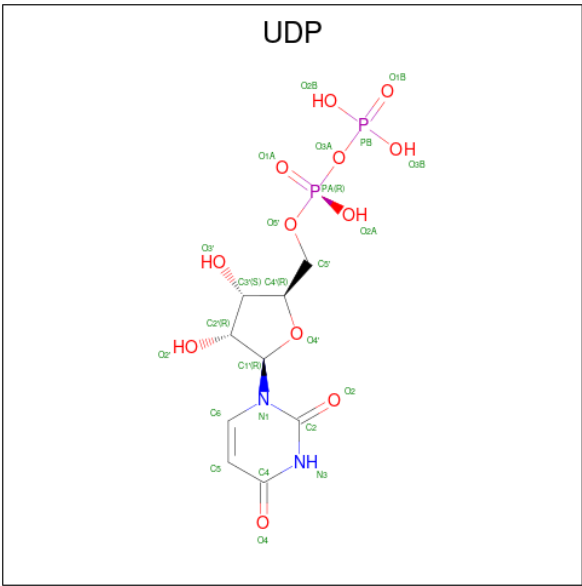


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	2	Total	C	O	0	0	0
			23	12	11			
2	D	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 3 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

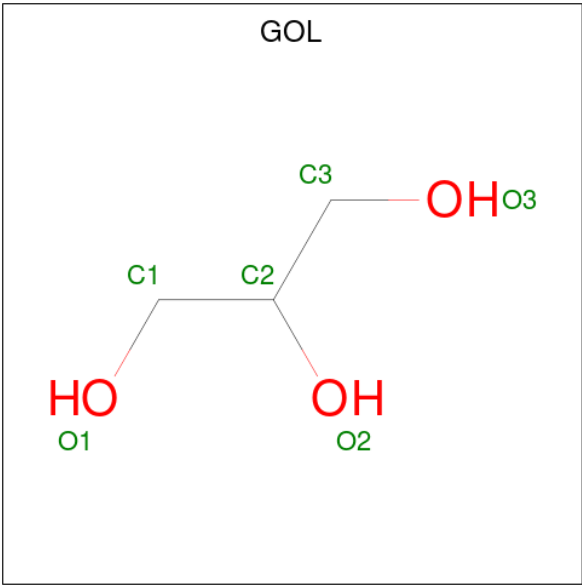
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mn	0	0
			1	1		
3	B	1	Total	Mn	0	0
			1	1		

- Molecule 4 is URIDINE-5'-DIPHOSPHATE (CCD ID: UDP) (formula: C₉H₁₄N₂O₁₂P₂).



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

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

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

- Molecule 6 is water.

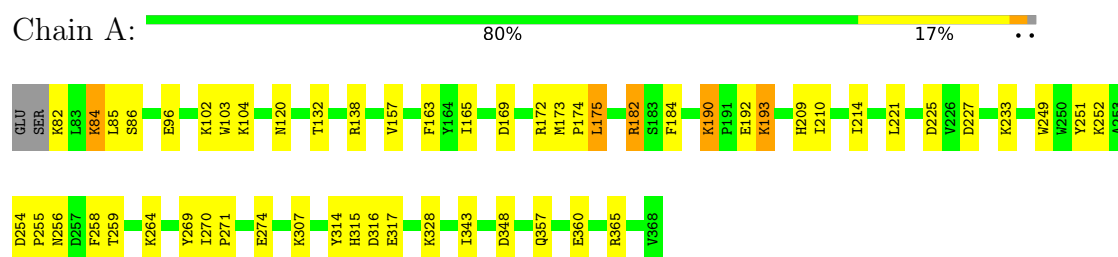
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	419	Total	O	0	0
			419	419		
6	B	444	Total	O	0	0
			444	444		

3 Residue-property plots [i](#)

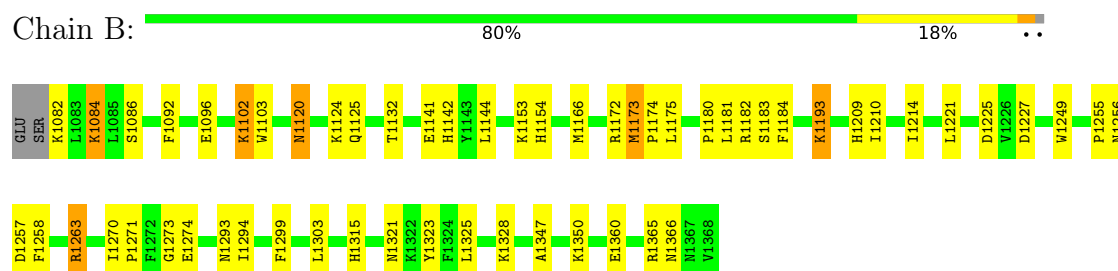
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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



- Molecule 2: beta-D-galactopyranose-(1-4)-beta-D-glucopyranose



- Molecule 2: beta-D-galactopyranose-(1-4)-beta-D-glucopyranose



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	44.86Å 93.68Å 93.71Å 90.00° 99.00° 90.00°	Depositor
Resolution (Å)	50.00 – 1.97	Depositor
% Data completeness (in resolution range)	97.5 (50.00-1.97)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.154 , 0.202	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5788	wwPDB-VP
Average B, all atoms (Å ²)	16.0	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, MN, UDP, GOL, BGC

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	0.55	0/2507	1.39	9/3395 (0.3%)
1	B	0.54	0/2484	1.42	18/3362 (0.5%)
All	All	0.54	0/4991	1.41	27/6757 (0.4%)

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1256	ASN	CA-CB-CG	-9.09	103.51	112.60
1	A	138	ARG	CD-NE-CZ	7.79	135.31	124.40
1	A	270	ILE	O-C-N	7.37	127.81	120.59
1	B	1092	PHE	CA-CB-CG	-7.15	106.65	113.80
1	A	315	HIS	CA-CB-CG	-6.85	106.95	113.80
1	A	256	ASN	CA-CB-CG	-6.81	105.79	112.60
1	B	1270	ILE	O-C-N	6.58	127.04	120.59
1	A	96	GLU	CB-CG-CD	6.55	123.74	112.60
1	B	1293	ASN	CA-CB-CG	-6.32	106.28	112.60
1	A	120	ASN	CA-CB-CG	-6.32	106.28	112.60
1	B	1183	SER	CA-C-N	-6.11	112.39	122.11
1	B	1183	SER	C-N-CA	-6.11	112.39	122.11
1	B	1257	ASP	CA-CB-CG	-6.00	106.60	112.60
1	B	1315	HIS	CA-CB-CG	-5.86	107.94	113.80
1	A	270	ILE	CA-C-O	-5.86	116.89	119.94
1	B	1209	HIS	CA-CB-CG	-5.58	108.22	113.80
1	B	1273	GLY	CA-C-N	-5.49	113.97	122.59
1	B	1273	GLY	C-N-CA	-5.49	113.97	122.59
1	B	1096	GLU	CB-CG-CD	5.32	121.64	112.60
1	A	209	HIS	CA-CB-CG	-5.18	108.62	113.80
1	B	1365	ARG	CA-C-N	-5.18	113.35	123.13
1	B	1365	ARG	C-N-CA	-5.18	113.35	123.13

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1325	LEU	CA-C-O	5.17	125.90	120.42
1	A	365	ARG	CA-C-O	-5.16	115.33	121.16
1	B	1120	ASN	CA-CB-CG	-5.12	107.48	112.60
1	B	1321	ASN	CA-CB-CG	-5.11	107.49	112.60
1	B	1263	ARG	NE-CZ-NH1	-5.02	116.48	121.50

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2414	0	2376	32	0
1	B	2401	0	2361	26	0
2	C	23	0	17	1	0
2	D	23	0	21	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	25	0	11	0	0
4	B	25	0	11	0	0
5	A	6	0	8	1	0
5	B	6	0	8	0	0
6	A	419	0	0	4	0
6	B	444	0	0	1	0
All	All	5788	0	4813	57	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 6.

All (57) 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:254:ASP:HB3	6:A:2266:HOH:O	1.86	0.74
1:A:165:ILE:HD13	1:A:173:MET:HE1	1.72	0.70
1:A:317:GLU:OE1	2:C:2:GAL:H62	1.93	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:271:PRO:HG2	1:A:274:GLU:OE1	1.95	0.67
1:B:1166:MET:N	1:B:1166:MET:HE2	2.15	0.62
1:A:190:LYS:HD3	1:A:192:GLU:OE1	2.02	0.60
1:A:255:PRO:HA	1:A:258:PHE:CD1	2.36	0.59
1:A:84:LYS:HE2	1:A:85:LEU:N	2.19	0.57
1:B:1255:PRO:HA	1:B:1258:PHE:CD1	2.40	0.55
1:B:1173:MET:HE1	1:B:1175:LEU:HD23	1.88	0.55
1:B:1263:ARG:HD2	6:B:4279:HOH:O	2.07	0.55
1:A:193:LYS:HD2	6:A:2191:HOH:O	2.09	0.53
1:A:132:THR:HG23	1:A:221:LEU:HD11	1.91	0.53
1:A:193:LYS:HE3	6:A:2191:HOH:O	2.08	0.53
1:B:1102:LYS:O	1:B:1102:LYS:HG2	2.09	0.52
1:A:210:ILE:HA	1:A:214:ILE:HB	1.93	0.51
1:B:1132:THR:HG23	1:B:1221:LEU:HD11	1.93	0.51
1:A:255:PRO:HA	1:A:258:PHE:CE1	2.47	0.50
1:A:252:LYS:HE2	1:A:348:ASP:OD2	2.13	0.49
1:B:1141:GLU:OE2	1:B:1172:ARG:HD2	2.12	0.49
1:B:1271:PRO:HG2	1:B:1274:GLU:CD	2.38	0.49
1:B:1084[A]:LYS:HE2	1:B:1086:SER:OG	2.13	0.49
1:B:1255:PRO:HA	1:B:1258:PHE:CE1	2.49	0.48
1:A:264:LYS:HA	1:A:269:TYR:CG	2.48	0.48
1:A:175:LEU:HD23	1:A:175:LEU:HA	1.59	0.47
1:A:163:PHE:HB2	1:A:184[B]:PHE:HB3	1.95	0.47
1:A:251:TYR:CE1	1:A:252:LYS:HG3	2.50	0.47
1:A:103:TRP:CE2	1:A:328:LYS:HB3	2.50	0.47
1:A:84:LYS:HE2	1:A:85:LEU:H	1.77	0.47
1:B:1210:ILE:HA	1:B:1214:ILE:HB	1.97	0.46
1:A:360:GLU:OE1	5:A:1372:GOL:H2	2.16	0.46
1:B:1193:LYS:O	1:B:1193:LYS:HE3	2.15	0.46
1:A:271:PRO:HG2	1:A:274:GLU:CD	2.40	0.46
1:B:1120:ASN:O	1:B:1124:LYS:HG3	2.17	0.45
1:A:157:VAL:HG21	1:A:233:LYS:HE3	1.98	0.45
1:B:1175:LEU:HD23	1:B:1175:LEU:HA	1.79	0.44
1:A:103:TRP:O	1:A:104:LYS:HB2	2.17	0.44
1:B:1294:ILE:HG12	1:B:1323:TYR:CE1	2.53	0.44
1:A:307:LYS:HE2	6:A:2322:HOH:O	2.17	0.43
1:B:1180:PRO:O	1:B:1181:LEU:HB2	2.17	0.43
1:B:1103:TRP:CE2	1:B:1328:LYS:HB3	2.54	0.43
1:B:1347:ALA:O	1:B:1350:LYS:HD3	2.19	0.43
1:B:1142:HIS:HE1	1:B:1360:GLU:OE2	2.02	0.42
1:A:249[B]:TRP:HD1	1:A:343:ILE:HD11	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:259:THR:HB	1:A:314:TYR:CE1	2.54	0.42
1:B:1153:LYS:HE2	1:B:1154:HIS:CE1	2.55	0.42
1:A:84:LYS:NZ	1:A:86:SER:H	2.18	0.41
1:B:1225:ASP:HB3	1:B:1227:ASP:OD1	2.20	0.41
1:A:169:ASP:OD2	1:A:172:ARG:HD3	2.19	0.41
1:A:175:LEU:CD1	1:B:1173:MET:HE2	2.51	0.41
1:A:173:MET:HE3	1:A:174:PRO:O	2.21	0.41
1:B:1173:MET:HE1	1:B:1175:LEU:CD2	2.51	0.41
1:B:1144:LEU:HD23	1:B:1174:PRO:HD2	2.02	0.40
1:A:225:ASP:HB3	1:A:227:ASP:OD1	2.22	0.40
1:B:1249:TRP:CG	2:D:1:BGC:H3	2.56	0.40
1:B:1299:PHE:CE2	1:B:1303:LEU:HD11	2.56	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	289/289 (100%)	283 (98%)	6 (2%)	0	100	100
1	B	287/289 (99%)	283 (99%)	4 (1%)	0	100	100
All	All	576/578 (100%)	566 (98%)	10 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

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

The Analysed column shows the number of residues for which the sidechain conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	262/260 (101%)	253 (97%)	9 (3%)	32	22
1	B	260/260 (100%)	249 (96%)	11 (4%)	26	15
All	All	522/520 (100%)	502 (96%)	20 (4%)	32	19

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	82	LYS
1	A	84	LYS
1	A	102	LYS
1	A	175	LEU
1	A	182	ARG
1	A	190	LYS
1	A	193	LYS
1	A	316	ASP
1	A	357	GLN
1	B	1082	LYS
1	B	1084[A]	LYS
1	B	1084[B]	LYS
1	B	1102	LYS
1	B	1125	GLN
1	B	1173	MET
1	B	1182	ARG
1	B	1184[A]	PHE
1	B	1184[B]	PHE
1	B	1193	LYS
1	B	1366	ASN

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

Mol	Chain	Res	Type
1	A	296	GLN
1	A	313	GLN
1	B	1142	HIS
1	B	1231	GLN
1	B	1296	GLN
1	B	1313	GLN
1	B	1342	HIS

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 ⓘ

4 monosaccharides 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	BGC	C	1	2	12,12,12	0.93	0	17,17,17	0.77	0
2	GAL	C	2	2	11,11,12	1.08	1 (9%)	15,15,17	1.75	2 (13%)
2	BGC	D	1	2	12,12,12	0.95	0	17,17,17	1.09	1 (5%)
2	GAL	D	2	2	11,11,12	1.11	1 (9%)	15,15,17	1.03	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	BGC	C	1	2	-	1/2/22/22	0/1/1/1
2	GAL	C	2	2	-	0/2/19/22	0/1/1/1
2	BGC	D	1	2	-	0/2/22/22	0/1/1/1
2	GAL	D	2	2	-	0/2/19/22	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	2	GAL	C4-C5	2.64	1.58	1.53
2	D	2	GAL	C4-C5	2.38	1.58	1.53

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	2	GAL	C3-C4-C5	-4.32	102.39	110.23
2	C	2	GAL	C1-O5-C5	3.27	116.57	112.19
2	D	1	BGC	C4-C3-C2	-2.09	107.15	110.83

There are no chirality outliers.

All (1) torsion outliers are listed below:

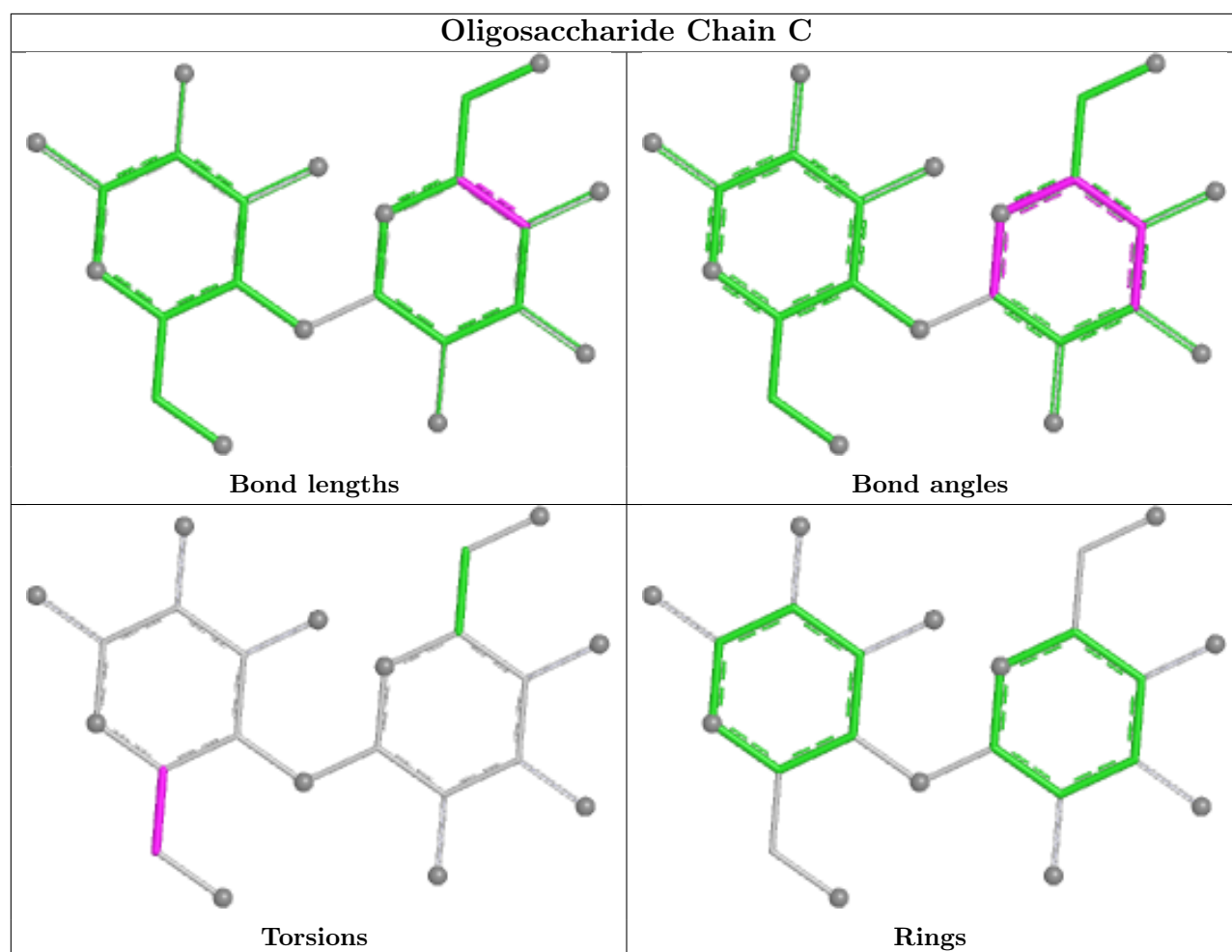
Mol	Chain	Res	Type	Atoms
2	C	1	BGC	O5-C5-C6-O6

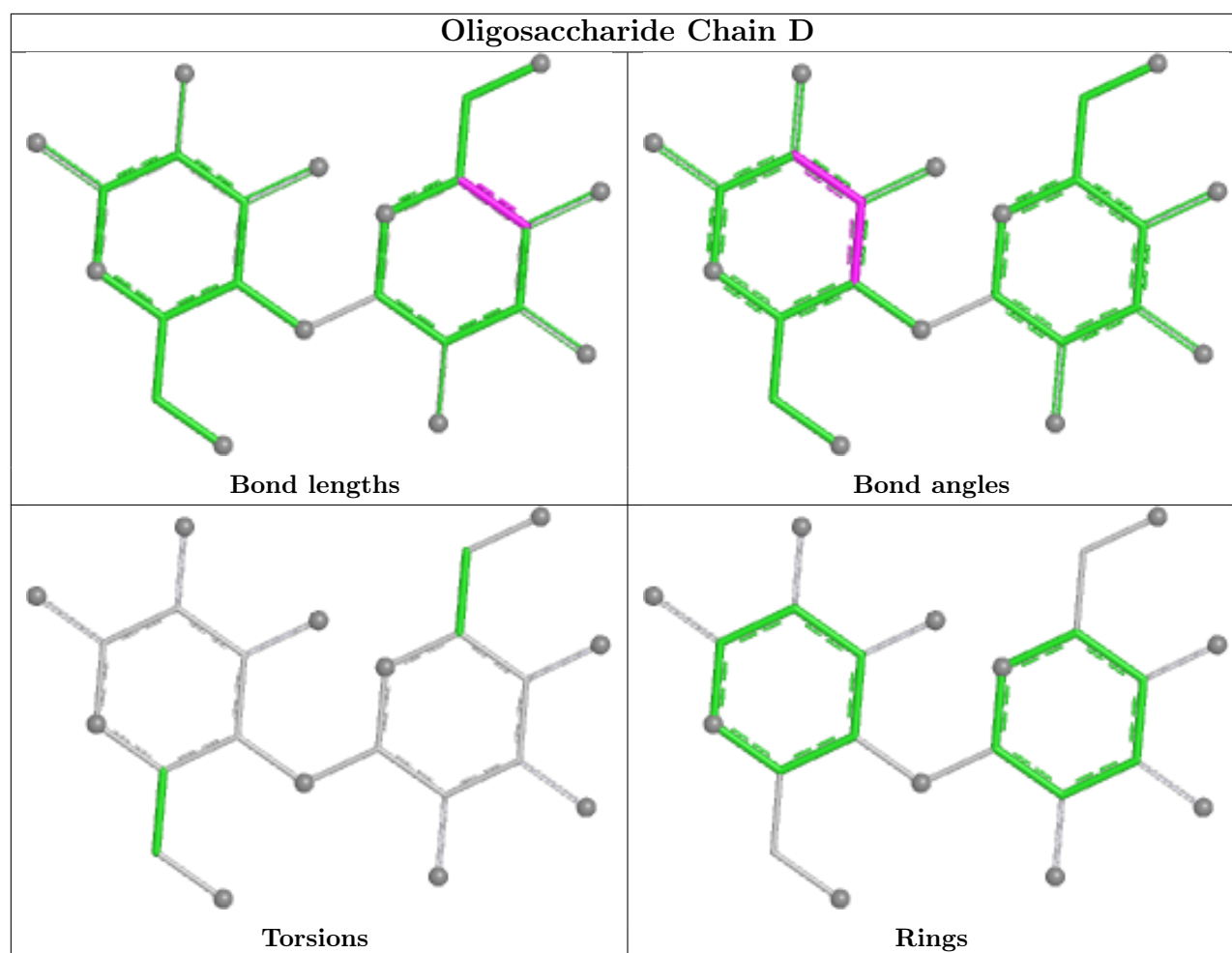
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1	BGC	1	0
2	C	2	GAL	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.





5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	GOL	A	1372	-	5,5,5	1.13	0	5,5,5	0.87	0
5	GOL	B	2372	-	5,5,5	1.18	1 (20%)	5,5,5	0.93	0
4	UDP	A	1371	3	25,26,26	1.21	2 (8%)	38,40,40	1.35	4 (10%)
4	UDP	B	2371	3	25,26,26	1.27	3 (12%)	38,40,40	1.36	5 (13%)

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	GOL	A	1372	-	-	2/4/4/4	-
5	GOL	B	2372	-	-	0/4/4/4	-
4	UDP	A	1371	3	-	1/16/32/32	0/2/2/2
4	UDP	B	2371	3	-	2/16/32/32	0/2/2/2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1371	UDP	C4-N3	3.11	1.43	1.38
4	B	2371	UDP	C4-N3	2.92	1.43	1.38
4	A	1371	UDP	PB-O1B	2.77	1.59	1.50
4	B	2371	UDP	PB-O1B	2.47	1.58	1.50
4	B	2371	UDP	PA-O3A	2.41	1.62	1.59
5	B	2372	GOL	C3-C2	-2.10	1.43	1.51

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1371	UDP	O4-C4-N3	-4.92	112.14	119.27
4	B	2371	UDP	O2-C2-N1	-3.38	118.39	122.80
4	B	2371	UDP	C4-N3-C2	-3.24	122.59	126.61
4	A	1371	UDP	C4-N3-C2	-3.10	122.76	126.61
4	B	2371	UDP	O4-C4-N3	-3.04	114.87	119.27
4	A	1371	UDP	O3B-PB-O2B	2.95	118.86	107.80
4	B	2371	UDP	O3B-PB-O2B	2.71	117.97	107.80
4	B	2371	UDP	O3'-C3'-C2'	-2.32	104.37	111.82
4	A	1371	UDP	O4'-C4'-C3'	-2.08	101.02	105.15

There are no chirality outliers.

All (5) torsion outliers are listed below:

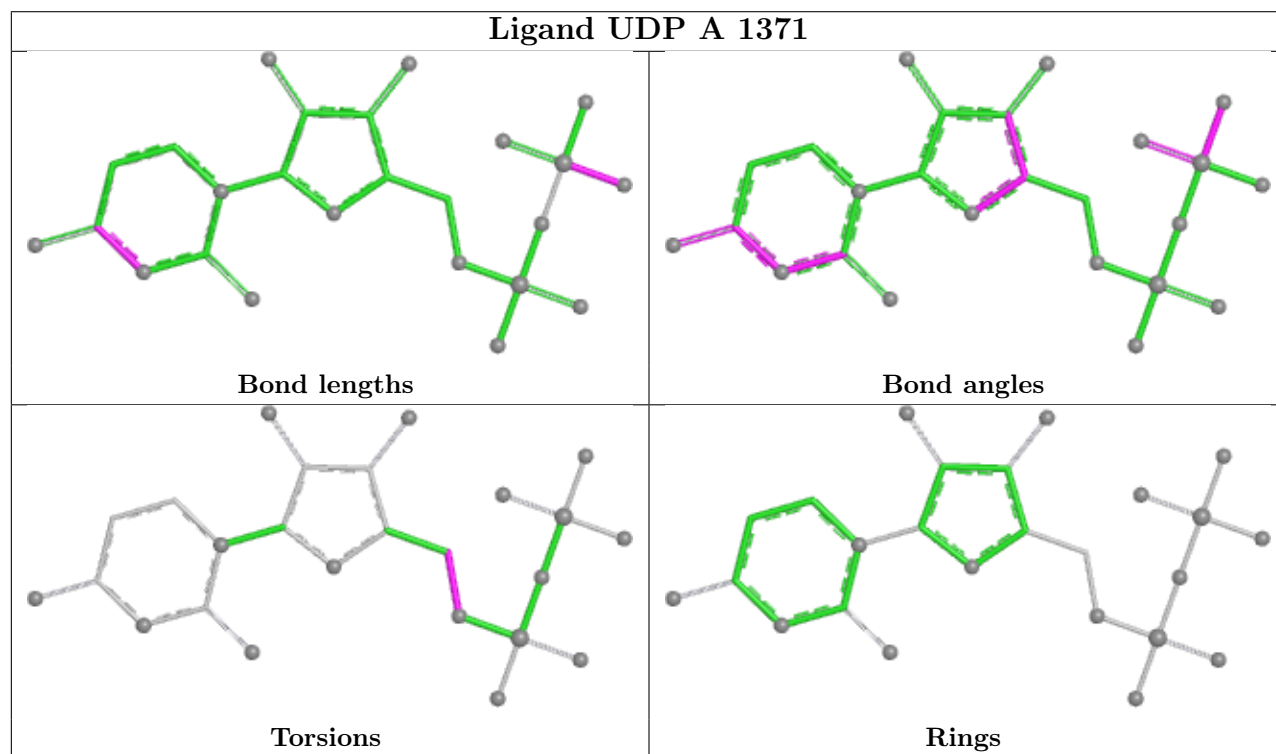
Mol	Chain	Res	Type	Atoms
5	A	1372	GOL	C1-C2-C3-O3
5	A	1372	GOL	O2-C2-C3-O3
4	A	1371	UDP	C4'-C5'-O5'-PA
4	B	2371	UDP	C4'-C5'-O5'-PA
4	B	2371	UDP	C3'-C4'-C5'-O5'

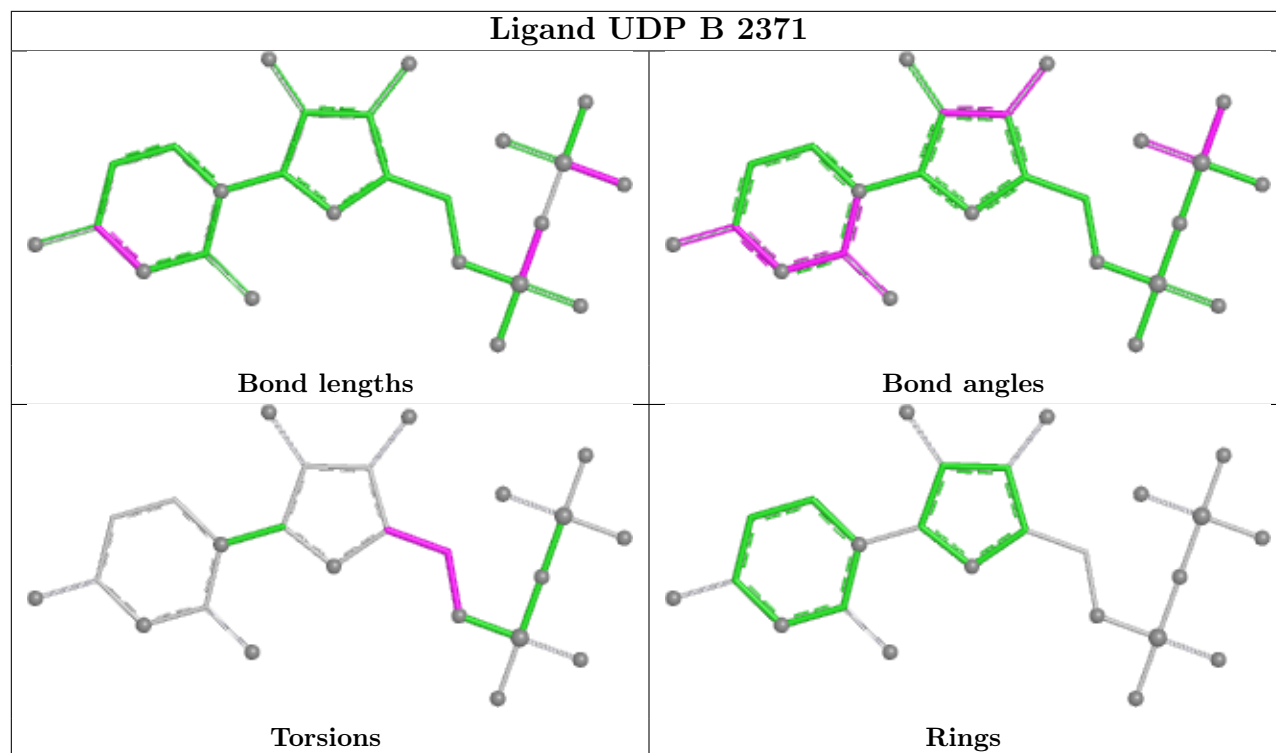
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1372	GOL	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.