



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:37 AM UTC

PDB ID : 2D7I / pdb_00002d7i
Title : Crystal structure of pp-GalNAc-T10 with UDP, GalNAc and Mn²⁺
Authors : Kubota, T.; Shiba, T.; Sugioka, S.; Kato, R.; Wakatsuki, S.; Narimatsu, H.
Deposited on : 2005-11-21
Resolution : 2.50 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) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

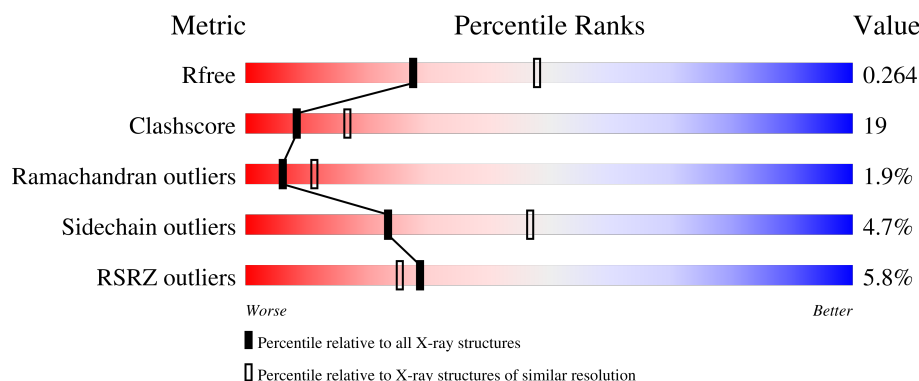
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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

Mol	Chain	Length	Quality of chain
1	A	570	5% 58% 32% 6%
2	B	2	50% 50%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NAG	A	704	X	-	-	-

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 4691 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 Polypeptide N-acetylgalactosaminyltransferase 10.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	536	4357	2762	778	789	28	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	34	HIS	-	expression tag	UNP Q86SR1
A	35	HIS	-	expression tag	UNP Q86SR1
A	36	HIS	-	expression tag	UNP Q86SR1
A	37	HIS	-	expression tag	UNP Q86SR1
A	38	HIS	-	expression tag	UNP Q86SR1
A	39	HIS	-	expression tag	UNP Q86SR1

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



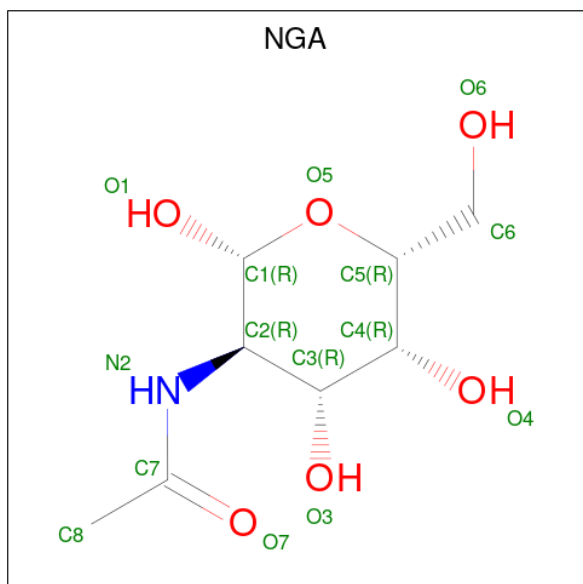
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	B	2	28	16	2	10	0	0	0

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-galactopyranose (CCD ID: NGA) (formula: $C_8H_{15}NO_6$).

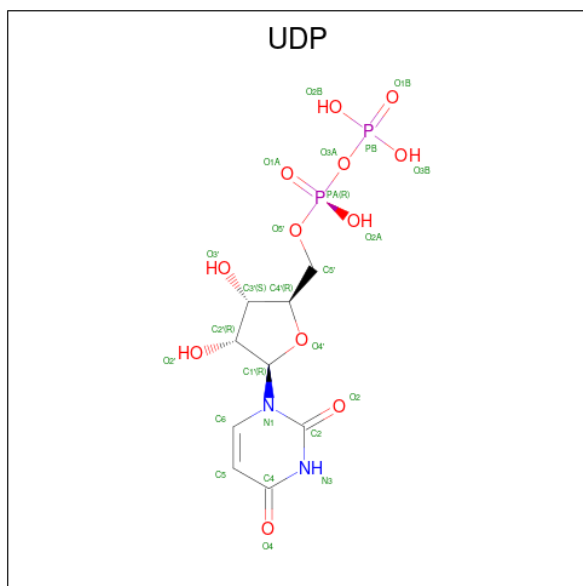


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			15	8	1	6		

- Molecule 5 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mn	0	0
			1	1		

- Molecule 6 is URIDINE-5'-DIPHOSPHATE (CCD ID: UDP) (formula: C₉H₁₄N₂O₁₂P₂).



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

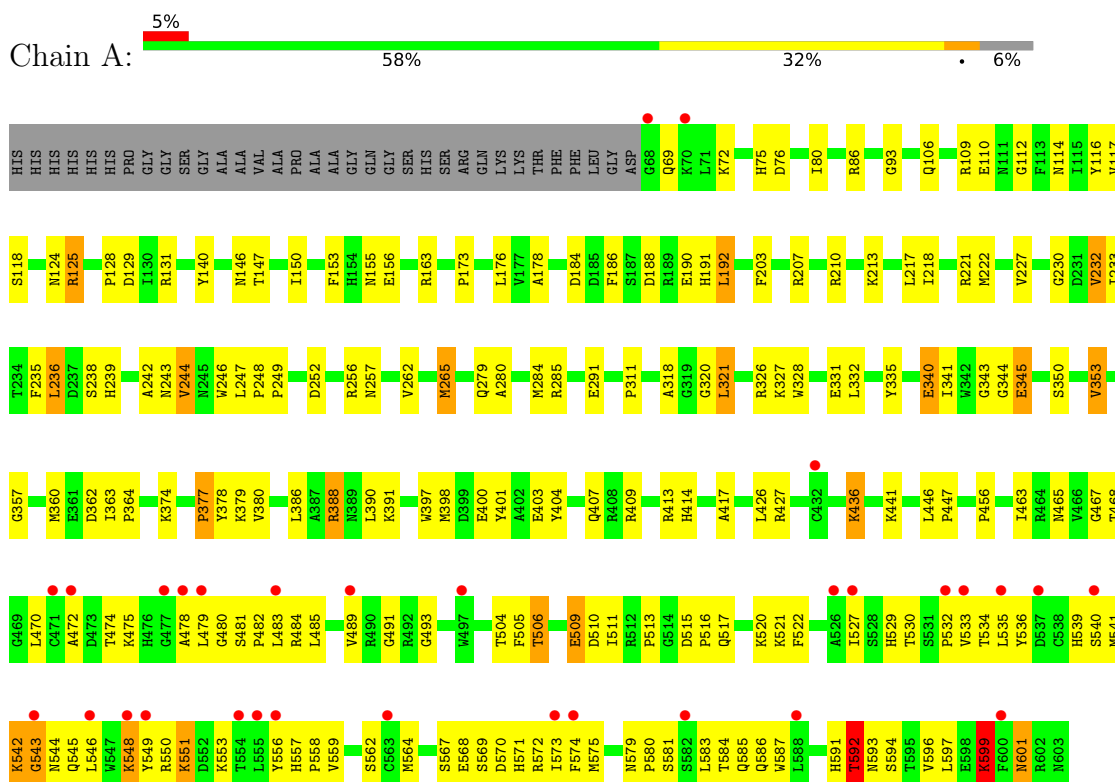
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	237	Total	O	0	0
			237	237		

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Polypeptide N-acetylgalactosaminyltransferase 10



- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	79.86Å 133.38Å 138.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.50 40.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (40.00-2.50) 99.3 (40.00-2.50)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.39 (at 2.51Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.215 , 0.267 0.220 , 0.264	Depositor DCC
R_{free} test set	2543 reflections (9.84%)	wwPDB-VP
Wilson B-factor (Å ²)	51.1	Xtriage
Anisotropy	0.557	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 45.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.014 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.029 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4691	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.89% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NGA, MN, UDP, 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	0.44	0/4481	0.92	20/6073 (0.3%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	203	PHE	CA-C-N	6.50	126.73	119.32
1	A	203	PHE	C-N-CA	6.50	126.73	119.32
1	A	409	ARG	CA-C-N	6.49	127.00	119.47
1	A	409	ARG	C-N-CA	6.49	127.00	119.47
1	A	357	GLY	N-CA-C	6.21	122.99	114.92
1	A	291	GLU	N-CA-C	-6.04	105.08	112.88
1	A	353	VAL	CB-CA-C	-5.99	104.05	112.14
1	A	601	ASN	N-CA-C	-5.92	105.65	114.64
1	A	265	MET	N-CA-C	-5.81	101.02	110.20
1	A	112	GLY	N-CA-C	-5.78	107.35	115.32
1	A	340	GLU	N-CA-C	5.47	122.44	110.80
1	A	592	THR	N-CA-C	5.43	115.40	108.24
1	A	244	VAL	N-CA-C	5.37	116.62	109.37
1	A	436	LYS	N-CA-C	-5.36	105.33	111.07
1	A	318	ALA	N-CA-C	-5.36	105.44	111.28
1	A	190	GLU	N-CA-C	5.33	117.09	111.28
1	A	398	MET	N-CA-C	5.28	119.59	113.20
1	A	344	GLY	N-CA-C	5.15	121.93	114.67
1	A	374	LYS	N-CA-C	-5.11	105.79	112.23
1	A	279	GLN	N-CA-C	-5.08	103.83	110.53

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	4357	0	4228	167	0
2	B	28	0	25	3	0
3	A	28	0	26	3	0
4	A	15	0	15	1	0
5	A	1	0	0	0	0
6	A	25	0	11	0	0
7	A	237	0	0	11	0
All	All	4691	0	4305	167	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 19.

All (167) 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:75:HIS:HE1	1:A:131:ARG:HH12	1.09	0.92
1:A:155:ASN:HD21	1:A:188:ASP:H	1.15	0.90
1:A:262:VAL:HB	1:A:363:ILE:HD11	1.58	0.84
1:A:567:SER:HB3	1:A:571:HIS:HB2	1.59	0.83
1:A:481:SER:HB3	1:A:484:ARG:HH22	1.45	0.81
1:A:75:HIS:CE1	1:A:131:ARG:HH12	1.98	0.78
1:A:218:ILE:HG13	1:A:345:GLU:HG2	1.65	0.77
1:A:542:LYS:HB3	1:A:546:LEU:HD22	1.66	0.76
1:A:262:VAL:HB	1:A:363:ILE:CD1	2.17	0.75
1:A:218:ILE:HD12	1:A:218:ILE:H	1.51	0.74
1:A:583:LEU:H	1:A:583:LEU:HD23	1.54	0.73
1:A:485:LEU:HD13	1:A:573:ILE:HD11	1.72	0.71
1:A:265:MET:HE1	1:A:362:ASP:HB3	1.71	0.71
1:A:509:GLU:HG2	7:A:908:HOH:O	1.90	0.70
1:A:388:ARG:HB2	7:A:895:HOH:O	1.92	0.68
1:A:474:THR:HG23	1:A:483:LEU:HD12	1.75	0.68
1:A:544:ASN:HA	1:A:559:VAL:HG11	1.76	0.68
1:A:593:ASN:OD1	1:A:596:VAL:HG23	1.94	0.68
1:A:489:VAL:HG13	1:A:493:GLY:HA3	1.76	0.68
1:A:114:ASN:CG	1:A:117:VAL:HG23	2.20	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:551:LYS:HZ3	1:A:551:LYS:HA	1.59	0.67
1:A:218:ILE:HG13	1:A:345:GLU:CG	2.25	0.66
1:A:482:PRO:HD2	1:A:572:ARG:HH21	1.62	0.64
1:A:548:LYS:HB3	1:A:556:TYR:HB3	1.80	0.63
1:A:114:ASN:C	1:A:114:ASN:HD22	2.06	0.63
1:A:363:ILE:HD12	1:A:363:ILE:N	2.13	0.63
1:A:75:HIS:HE1	1:A:131:ARG:NH1	1.92	0.63
1:A:467:GLY:HA2	1:A:583:LEU:HD12	1.80	0.63
1:A:86:ARG:HH21	2:B:1:NAG:C8	2.12	0.63
1:A:533:VAL:HG22	1:A:575:MET:CE	2.30	0.62
1:A:217:LEU:HG	1:A:218:ILE:HD12	1.82	0.62
1:A:539:HIS:O	1:A:539:HIS:ND1	2.33	0.62
1:A:465:ASN:HB2	1:A:587:TRP:CZ3	2.35	0.62
1:A:221:ARG:HD2	1:A:235:PHE:CD2	2.35	0.61
1:A:207:ARG:HH12	1:A:227:VAL:CG1	2.13	0.61
1:A:599:LYS:HZ2	1:A:599:LYS:HB2	1.64	0.61
1:A:548:LYS:HA	7:A:920:HOH:O	2.00	0.61
1:A:529:HIS:HB2	1:A:562:SER:HB3	1.81	0.61
1:A:328:TRP:CH2	1:A:332:LEU:HD21	2.36	0.60
1:A:335:TYR:CE2	1:A:345:GLU:HB2	2.36	0.60
1:A:475:LYS:H	1:A:484:ARG:NH1	1.99	0.60
1:A:86:ARG:HH21	2:B:1:NAG:H82	1.66	0.60
1:A:539:HIS:O	1:A:541:MET:N	2.31	0.59
1:A:551:LYS:HA	1:A:551:LYS:NZ	2.18	0.59
1:A:114:ASN:HD21	1:A:116:TYR:HB2	1.67	0.59
1:A:218:ILE:HD12	1:A:218:ILE:N	2.17	0.58
1:A:388:ARG:HD2	1:A:417:ALA:HA	1.84	0.58
1:A:548:LYS:HZ2	1:A:601:ASN:HB2	1.69	0.58
1:A:446:LEU:N	1:A:447:PRO:HD2	2.20	0.57
1:A:155:ASN:HD21	1:A:188:ASP:N	1.95	0.56
1:A:109:ARG:HB3	1:A:109:ARG:NH1	2.21	0.56
1:A:530:THR:HA	1:A:575:MET:O	2.06	0.56
1:A:553:LYS:O	1:A:586:GLN:HA	2.06	0.56
1:A:76:ASP:O	1:A:80:ILE:HD13	2.07	0.55
1:A:207:ARG:HH12	1:A:227:VAL:HG11	1.70	0.55
1:A:441:LYS:HE2	7:A:798:HOH:O	2.06	0.55
1:A:548:LYS:HE2	1:A:549:TYR:N	2.22	0.55
1:A:564:MET:HE3	1:A:573:ILE:HG21	1.90	0.54
1:A:265:MET:HE3	1:A:364:PRO:N	2.23	0.54
1:A:482:PRO:CD	1:A:572:ARG:HH21	2.21	0.54
1:A:401:TYR:OH	1:A:436:LYS:HD2	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:327:LYS:O	1:A:331:GLU:HG3	2.09	0.53
1:A:482:PRO:HB2	1:A:572:ARG:HH21	1.74	0.53
1:A:109:ARG:HB3	1:A:109:ARG:HH11	1.73	0.53
1:A:541:MET:C	1:A:543:GLY:H	2.17	0.52
1:A:218:ILE:H	1:A:218:ILE:CD1	2.22	0.52
1:A:568:GLU:C	1:A:570:ASP:H	2.16	0.52
1:A:550:ARG:HB2	7:A:866:HOH:O	2.07	0.52
1:A:332:LEU:O	1:A:426:LEU:HD21	2.09	0.52
1:A:591:HIS:O	1:A:592:THR:HG23	2.09	0.52
1:A:156:GLU:HG2	1:A:238:SER:OG	2.10	0.51
1:A:86:ARG:NH2	2:B:1:NAG:H82	2.26	0.51
1:A:533:VAL:HG22	1:A:575:MET:HE3	1.93	0.51
1:A:265:MET:CE	1:A:362:ASP:HB3	2.41	0.50
1:A:489:VAL:HG13	1:A:493:GLY:CA	2.41	0.50
1:A:446:LEU:C	1:A:446:LEU:HD23	2.37	0.50
1:A:539:HIS:H	1:A:545:GLN:HE22	1.60	0.50
1:A:599:LYS:HZ1	1:A:599:LYS:HA	1.76	0.50
1:A:583:LEU:HD23	1:A:583:LEU:N	2.26	0.49
1:A:178:ALA:HA	3:A:704:NAG:H83	1.93	0.49
1:A:118:SER:OG	1:A:163:ARG:NH1	2.46	0.49
1:A:280:ALA:O	1:A:284:MET:HE2	2.13	0.49
1:A:504:THR:HG21	1:A:516:PRO:HA	1.95	0.48
1:A:557:HIS:CD2	1:A:559:VAL:HG12	2.48	0.48
1:A:456:PRO:HG3	3:A:701:NAG:H3	1.94	0.48
1:A:69:GLN:H	1:A:69:GLN:CD	2.22	0.48
1:A:252:ASP:O	1:A:256:ARG:HG2	2.14	0.48
1:A:465:ASN:HD21	1:A:584:THR:HA	1.79	0.48
1:A:532:PRO:HA	1:A:574:PHE:HA	1.96	0.48
1:A:594:SER:HA	1:A:597:LEU:HB2	1.96	0.48
1:A:579:ASN:C	1:A:581:SER:H	2.21	0.47
1:A:69:GLN:HB3	7:A:746:HOH:O	2.14	0.47
1:A:482:PRO:HB2	1:A:572:ARG:NH2	2.30	0.46
1:A:350:SER:OG	1:A:360:MET:HE1	2.16	0.46
1:A:478:ALA:HB1	1:A:522:PHE:CZ	2.51	0.46
1:A:548:LYS:NZ	1:A:601:ASN:HB2	2.31	0.46
1:A:146:ASN:OD1	1:A:230:GLY:HA2	2.16	0.46
1:A:153:PHE:CE1	1:A:192:LEU:HG	2.51	0.46
1:A:505:PHE:CD2	1:A:511:ILE:HG12	2.50	0.46
1:A:184:ASP:HB3	1:A:210:ARG:HG2	1.98	0.46
1:A:534:THR:HG23	7:A:903:HOH:O	2.15	0.46
1:A:380:VAL:O	1:A:380:VAL:HG23	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:184:ASP:OD1	1:A:210:ARG:NH1	2.46	0.45
1:A:515:ASP:C	1:A:517:GLN:H	2.24	0.45
1:A:109:ARG:NH1	7:A:727:HOH:O	2.48	0.45
1:A:557:HIS:NE2	1:A:559:VAL:HG12	2.31	0.45
1:A:544:ASN:ND2	7:A:907:HOH:O	2.48	0.45
1:A:401:TYR:HA	1:A:404:TYR:CD1	2.52	0.45
1:A:140:TYR:CZ	1:A:249:PRO:HB3	2.52	0.45
1:A:506:THR:OG1	1:A:510:ASP:OD2	2.29	0.45
1:A:114:ASN:ND2	1:A:117:VAL:HG23	2.32	0.44
1:A:341:ILE:HD12	1:A:341:ILE:N	2.31	0.44
1:A:343:GLY:HA3	1:A:386:LEU:CD1	2.48	0.44
1:A:247:LEU:HB3	1:A:248:PRO:HD3	1.99	0.44
1:A:114:ASN:ND2	1:A:117:VAL:H	2.16	0.44
1:A:556:TYR:CE2	1:A:558:PRO:HB3	2.53	0.44
1:A:403:GLU:O	1:A:407:GLN:HG3	2.18	0.43
1:A:513:PRO:HG2	1:A:520:LYS:HB2	2.00	0.43
1:A:147:THR:HB	1:A:232:VAL:HG22	2.00	0.43
1:A:599:LYS:CB	1:A:599:LYS:NZ	2.81	0.43
1:A:482:PRO:HD2	1:A:572:ARG:NH2	2.29	0.43
1:A:534:THR:HG22	1:A:535:LEU:N	2.33	0.43
1:A:377:PRO:HG2	1:A:378:TYR:H	1.82	0.43
1:A:489:VAL:HG12	1:A:491:GLY:H	1.83	0.43
1:A:186:PHE:CD1	1:A:213:LYS:HA	2.54	0.43
1:A:129:ASP:OD2	1:A:129:ASP:C	2.61	0.43
1:A:320:GLY:C	1:A:321:LEU:HD22	2.44	0.43
1:A:262:VAL:CB	1:A:363:ILE:HD11	2.40	0.42
1:A:233:ILE:HD11	1:A:326:ARG:HG3	2.02	0.42
1:A:479:LEU:HD23	1:A:480:GLY:N	2.34	0.42
1:A:548:LYS:HD2	1:A:601:ASN:HA	2.01	0.42
1:A:599:LYS:HB2	1:A:599:LYS:NZ	2.31	0.42
1:A:69:GLN:HG2	1:A:69:GLN:O	2.19	0.42
1:A:191:HIS:ND1	1:A:192:LEU:HD13	2.34	0.42
1:A:463:ILE:HB	1:A:472:ALA:HB3	2.01	0.42
1:A:534:THR:CG2	1:A:535:LEU:N	2.82	0.42
1:A:114:ASN:C	1:A:114:ASN:ND2	2.77	0.42
1:A:456:PRO:HG3	3:A:701:NAG:H83	2.01	0.42
1:A:150:ILE:HD12	1:A:150:ILE:N	2.34	0.42
1:A:153:PHE:HE1	1:A:192:LEU:HG	1.85	0.42
1:A:520:LYS:C	1:A:522:PHE:H	2.27	0.42
1:A:489:VAL:HG13	1:A:493:GLY:N	2.34	0.42
1:A:548:LYS:NZ	1:A:601:ASN:CB	2.83	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:515:ASP:C	1:A:517:GLN:N	2.77	0.42
1:A:265:MET:HE3	1:A:363:ILE:C	2.45	0.41
1:A:236:LEU:HD12	1:A:321:LEU:HD13	2.01	0.41
1:A:397:TRP:HA	1:A:427:ARG:NH1	2.35	0.41
1:A:124:ASN:O	1:A:125:ARG:C	2.62	0.41
1:A:482:PRO:CB	1:A:572:ARG:HH21	2.32	0.41
1:A:242:ALA:HB1	1:A:246:TRP:CD1	2.56	0.41
1:A:243:ASN:OD1	1:A:244:VAL:HG12	2.20	0.41
1:A:173:PRO:HB2	1:A:176:LEU:HD12	2.03	0.41
1:A:72:LYS:NZ	7:A:923:HOH:O	2.53	0.41
1:A:285:ARG:HD2	7:A:781:HOH:O	2.20	0.41
1:A:362:ASP:O	1:A:364:PRO:HD3	2.20	0.41
1:A:106:GLN:NE2	1:A:109:ARG:HG3	2.36	0.41
1:A:93:GLY:HA3	1:A:128:PRO:HD3	2.03	0.41
1:A:481:SER:HB3	1:A:484:ARG:NH2	2.26	0.41
1:A:534:THR:HG22	1:A:536:TYR:N	2.36	0.41
1:A:207:ARG:NH1	1:A:227:VAL:HG11	2.35	0.40
1:A:222:MET:HE2	1:A:335:TYR:O	2.21	0.40
1:A:569:SER:OG	1:A:571:HIS:HD2	2.05	0.40
1:A:413:ARG:HH11	1:A:413:ARG:HG2	1.86	0.40
1:A:468:THR:C	1:A:470:LEU:H	2.29	0.40
1:A:345:GLU:OE2	4:A:702:NGA:H62	2.22	0.40
1:A:529:HIS:ND1	1:A:530:THR:HG23	2.35	0.40
1:A:579:ASN:HB3	1:A:585:GLN:OE1	2.22	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	534/570 (94%)	478 (90%)	46 (9%)	10 (2%)	6 11

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	540	SER
1	A	599	LYS
1	A	340	GLU
1	A	527	ILE
1	A	542	LYS
1	A	377	PRO
1	A	580	PRO
1	A	506	THR
1	A	521	LYS
1	A	543	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	471/494 (95%)	449 (95%)	22 (5%)	23	47

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	110	GLU
1	A	125	ARG
1	A	192	LEU
1	A	232	VAL
1	A	236	LEU
1	A	239	HIS
1	A	257	ASN
1	A	311	PRO
1	A	321	LEU
1	A	345	GLU
1	A	353	VAL
1	A	379	LYS
1	A	388	ARG
1	A	390	LEU
1	A	391	LYS

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Mol	Chain	Res	Type
1	A	400	GLU
1	A	414	HIS
1	A	509	GLU
1	A	548	LYS
1	A	551	LYS
1	A	592	THR
1	A	599	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	75	HIS
1	A	85	GLN
1	A	106	GLN
1	A	114	ASN
1	A	134	ASN
1	A	155	ASN
1	A	170	ASN
1	A	245	ASN
1	A	271	HIS
1	A	304	GLN
1	A	407	GLN
1	A	423	GLN
1	A	465	ASN
1	A	476	HIS
1	A	545	GLN
1	A	571	HIS
1	A	576	ASN
1	A	579	ASN

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

2 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	NAG	B	1	1,2	14,14,15	0.65	0	17,19,21	1.11	2 (11%)
2	NAG	B	2	2	14,14,15	0.70	0	17,19,21	0.79	1 (5%)

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	B	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	B	2	2	-	4/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	NAG	C2-N2-C7	-2.73	119.24	122.90
2	B	2	NAG	C2-N2-C7	-2.21	119.94	122.90
2	B	1	NAG	C3-C4-C5	-2.17	106.29	110.23

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	2	NAG	O7-C7-N2-C2
2	B	2	NAG	C8-C7-N2-C2
2	B	2	NAG	O5-C5-C6-O6
2	B	1	NAG	C8-C7-N2-C2
2	B	1	NAG	O7-C7-N2-C2

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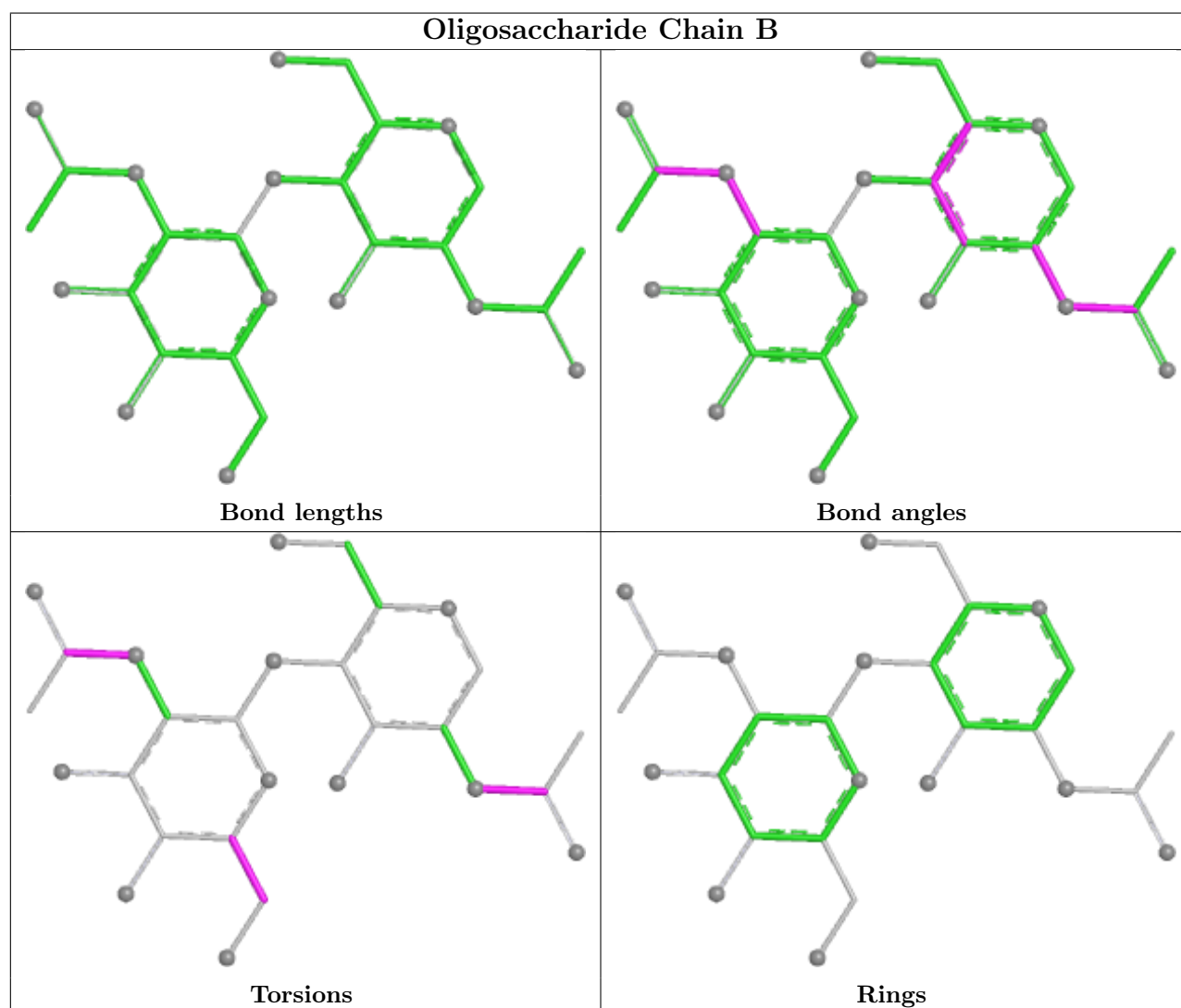
Mol	Chain	Res	Type	Atoms
2	B	2	NAG	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1	NAG	3	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 5 ligands modelled in this entry, 1 is 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	$\# Z > 2$
6	UDP	A	705	5	25,26,26	0.97	1 (4%)	38,40,40	1.48	5 (13%)
3	NAG	A	701	1	14,14,15	0.70	0	17,19,21	0.67	0
3	NAG	A	704	1	14,14,15	0.73	0	17,19,21	0.95	2 (11%)
4	NGA	A	702	-	15,15,15	0.61	0	21,21,21	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	UDP	A	705	5	-	1/16/32/32	0/2/2/2
3	NAG	A	701	1	-	5/6/23/26	0/1/1/1
3	NAG	A	704	1	1/1/5/7	2/6/23/26	0/1/1/1
4	NGA	A	702	-	-	1/6/26/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	705	UDP	C6-C5	2.28	1.40	1.35

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	705	UDP	C4-N3-C2	-4.90	120.53	126.61
6	A	705	UDP	C5-C4-N3	3.96	120.34	114.80
6	A	705	UDP	N3-C2-N1	3.72	119.73	114.89
3	A	704	NAG	C2-N2-C7	-2.41	119.67	122.90
6	A	705	UDP	O4-C4-C5	-2.37	121.08	125.16
6	A	705	UDP	O2B-PB-O3A	2.11	111.72	104.64
3	A	704	NAG	O5-C1-C2	-2.11	108.03	111.29

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	704	NAG	C1

All (9) torsion outliers are listed below:

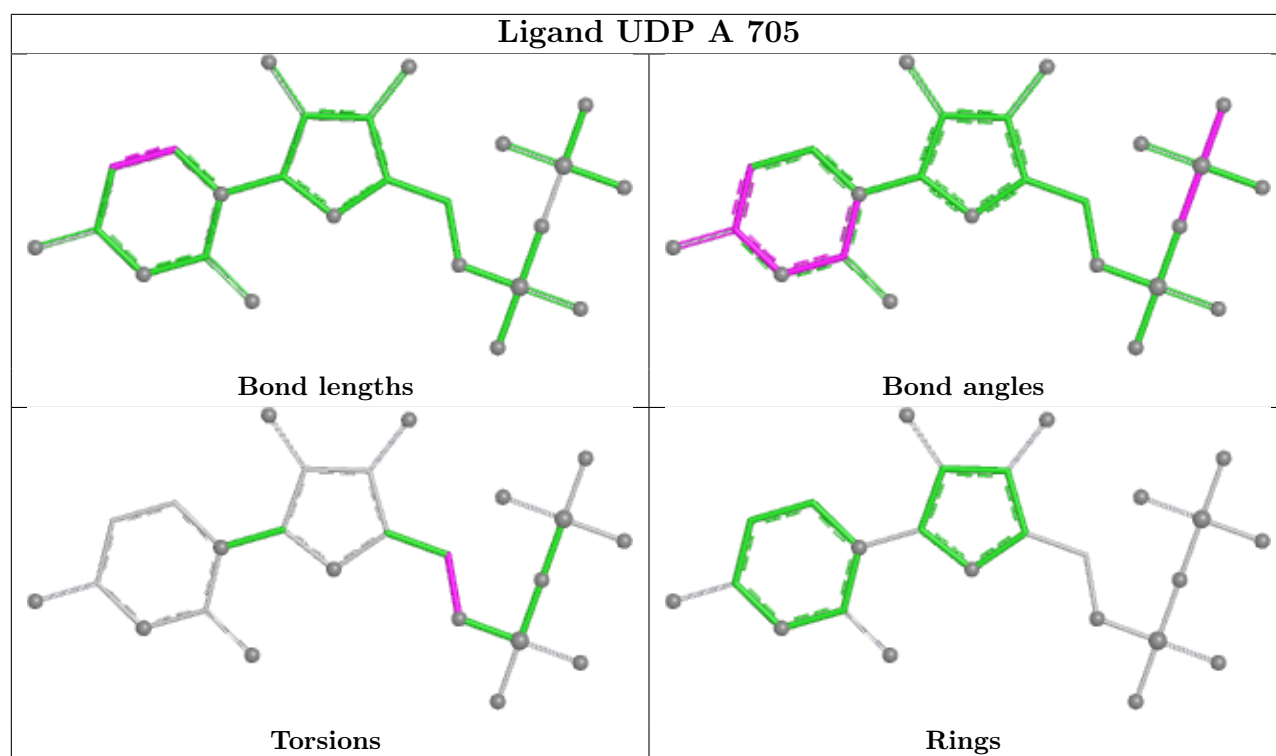
Mol	Chain	Res	Type	Atoms
3	A	701	NAG	C8-C7-N2-C2
3	A	701	NAG	O7-C7-N2-C2
3	A	704	NAG	C8-C7-N2-C2
3	A	704	NAG	O7-C7-N2-C2
3	A	701	NAG	O5-C5-C6-O6
3	A	701	NAG	C4-C5-C6-O6
4	A	702	NGA	O5-C5-C6-O6
3	A	701	NAG	C3-C2-N2-C7
6	A	705	UDP	C4'-C5'-O5'-PA

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	701	NAG	2	0
3	A	704	NAG	1	0
4	A	702	NGA	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 ⓘ

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	536/570 (94%)	0.27	31 (5%) 29 25	29, 53, 109, 126	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	477	GLY	4.4
1	A	532	PRO	4.1
1	A	556	TYR	3.9
1	A	573	ILE	3.5
1	A	554	THR	3.3
1	A	533	VAL	3.1
1	A	527	ILE	3.0
1	A	555	LEU	2.9
1	A	574	PHE	2.9
1	A	497	TRP	2.8
1	A	479	LEU	2.6
1	A	543	GLY	2.6
1	A	489	VAL	2.6
1	A	472	ALA	2.6
1	A	535	LEU	2.6
1	A	600	PHE	2.6
1	A	432	CYS	2.6
1	A	471	CYS	2.5
1	A	563	CYS	2.5
1	A	548	LYS	2.5
1	A	582	SER	2.4
1	A	549	TYR	2.3
1	A	68	GLY	2.3
1	A	540	SER	2.3
1	A	70	LYS	2.2
1	A	546	LEU	2.2
1	A	537	ASP	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	588	LEU	2.2
1	A	478	ALA	2.1
1	A	483	LEU	2.1
1	A	526	ALA	2.1

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

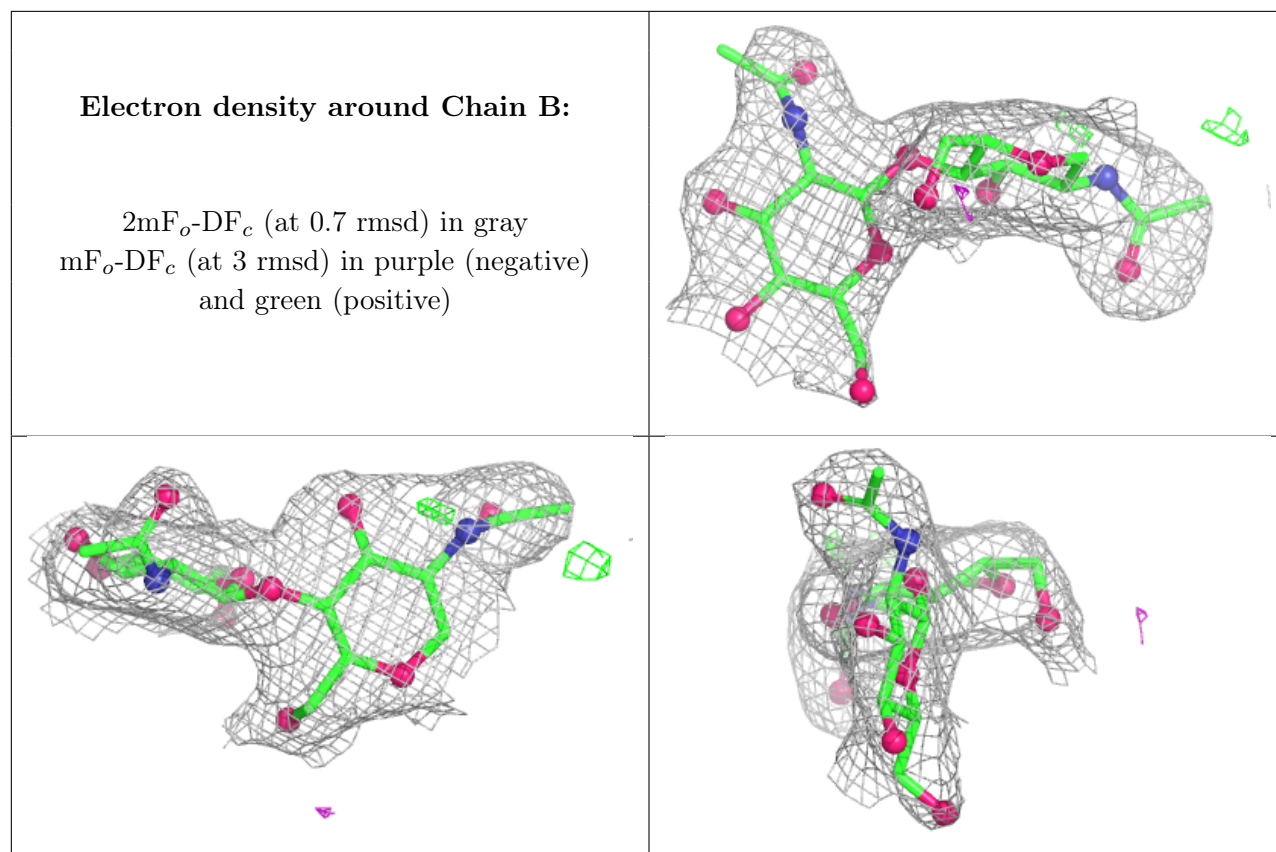
There are no non-standard protein/DNA/RNA residues in this entry.

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. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	B	2	14/15	0.85	0.11	71,74,75,75	0
2	NAG	B	1	14/15	0.91	0.10	52,57,61,66	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



6.4 Ligands [i](#)

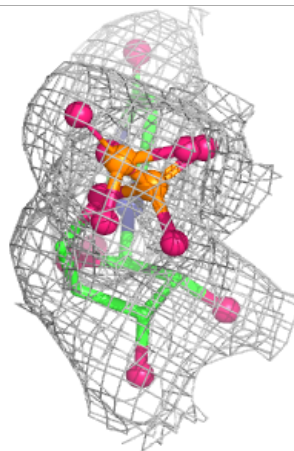
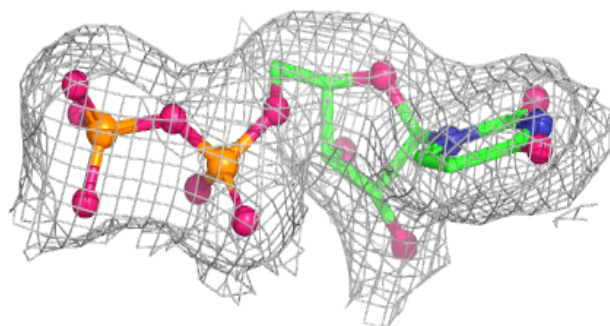
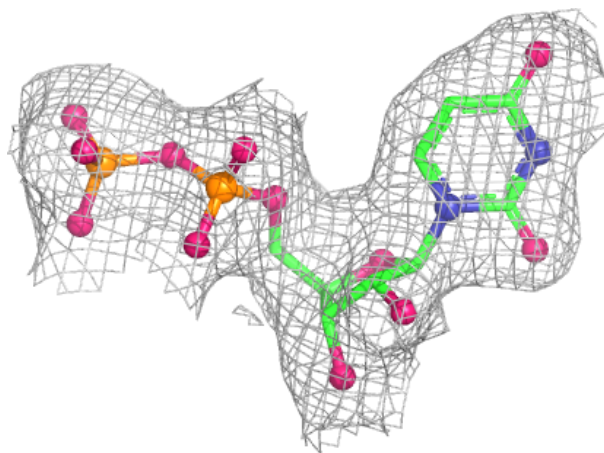
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	A	704	14/15	0.61	0.15	82,84,87,88	0
3	NAG	A	701	14/15	0.78	0.12	92,95,97,97	0
4	NGA	A	702	15/15	0.92	0.12	53,57,59,59	0
6	UDP	A	705	25/25	0.98	0.05	33,36,39,39	0
5	MN	A	706	1/1	1.00	0.02	30,30,30,30	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around UDP A 705:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.