



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 01:31 PM UTC

PDB ID : 8JR5 / pdb_00008jr5
Title : Crystal structure of Hendra Virus attachment(G) glycoprotein mutant S586N
Authors : Li, Y.H.; Huang, X.Y.; Xu, J.J.; Chen, W.
Deposited on : 2023-06-16
Resolution : 3.30 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

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

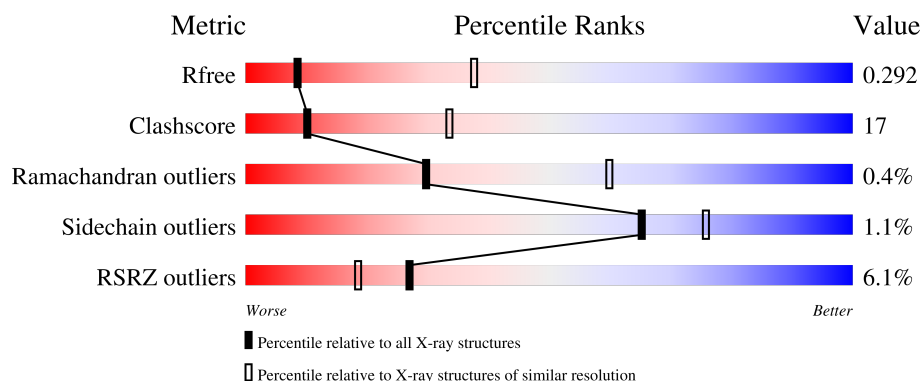
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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))
R_{free}	180053	1169 (3.32-3.28)
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)

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	416	7% 62% 33% ..
1	B	416	5% 64% 32% ..
2	C	2	100%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6451 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 Glycoprotein G.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	402	Total	C	N	O	S	0	0	0
			3182	2032	533	598	19			
1	B	403	Total	C	N	O	S	0	0	0
			3185	2033	534	599	19			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	586	ASN	SER	engineered mutation	UNP O89343
B	586	ASN	SER	engineered mutation	UNP O89343

- 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
2	C	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆) (labeled as "Ligand of Interest" by depositor).

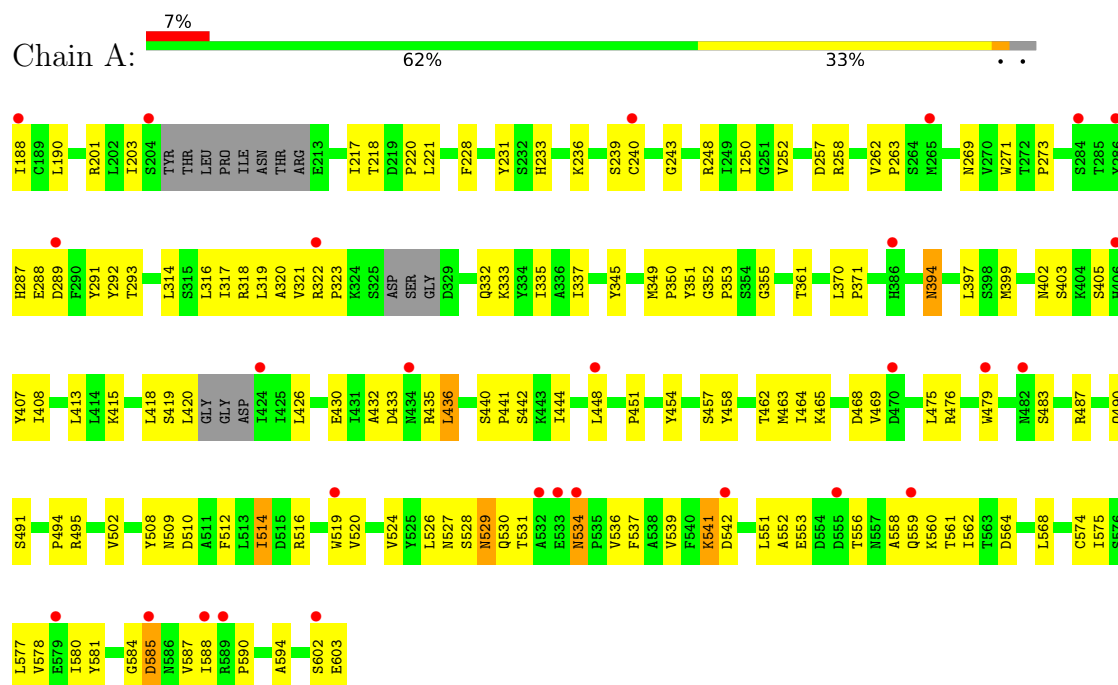


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		

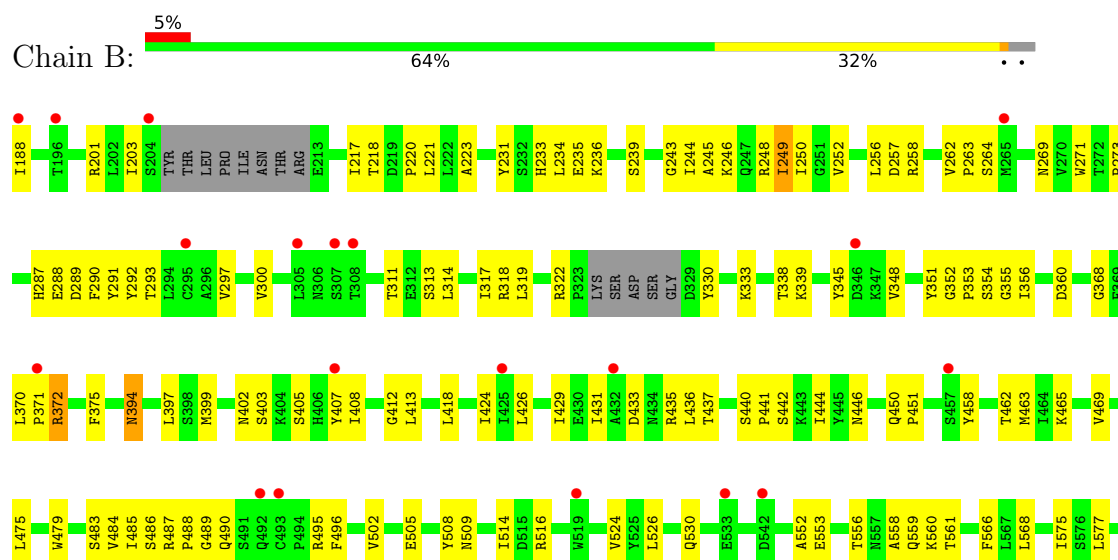
3 Residue-property plots [i](#)

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 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: Glycoprotein G



• Molecule 1: Glycoprotein G





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

Chain C:  100%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.93Å 63.66Å 121.62Å 90.00° 97.69° 90.00°	Depositor
Resolution (Å)	41.59 – 3.30 41.59 – 3.30	Depositor EDS
% Data completeness (in resolution range)	98.5 (41.59-3.30) 98.5 (41.59-3.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.08 (at 3.32Å)	Xtriage
Refinement program	PHENIX (1.18.2_3874: ???)	Depositor
R, R_{free}	0.258 , 0.290 0.258 , 0.292	Depositor DCC
R_{free} test set	968 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	87.0	Xtriage
Anisotropy	0.600	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 44.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	6451	wwPDB-VP
Average B, all atoms (Å ²)	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.62% 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: 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	0/3257	1.06	7/4432 (0.2%)
1	B	0.63	0/3261	1.00	4/4438 (0.1%)
All	All	0.64	0/6518	1.03	11/8870 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	529	ASN	CB-CA-C	-7.37	98.25	110.19
1	A	529	ASN	CA-CB-CG	6.88	119.47	112.60
1	A	559	GLN	N-CA-C	6.76	118.65	111.28
1	B	394	ASN	CB-CG-ND2	6.19	125.68	116.40
1	A	514	ILE	N-CA-C	6.18	116.34	110.53
1	A	394	ASN	CB-CG-ND2	6.18	125.66	116.40
1	B	559	GLN	N-CA-C	5.61	117.39	111.28
1	A	534	ASN	N-CA-C	5.29	121.51	109.81
1	A	394	ASN	CB-CG-OD1	-5.08	110.63	120.80
1	B	311	THR	CA-C-N	-5.02	115.74	122.42
1	B	311	THR	C-N-CA	-5.02	115.74	122.42

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	372	ARG	Sidechain

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	3182	0	3120	110	0
1	B	3185	0	3124	112	0
2	C	28	0	25	0	0
3	A	28	0	26	0	0
3	B	28	0	26	1	0
All	All	6451	0	6321	219	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 17.

All (219) 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:B:360:ASP:HB3	1:B:418:LEU:HB2	1.41	1.01
1:A:552:ALA:HB1	1:A:556:THR:HG21	1.54	0.89
1:B:394:ASN:HD22	1:B:502:VAL:HG23	1.36	0.88
1:B:234:LEU:HD11	1:B:245:ALA:HB1	1.55	0.87
1:A:258:ARG:HH12	1:B:322:ARG:HD2	1.39	0.87
1:A:394:ASN:HD22	1:A:502:VAL:HG23	1.42	0.84
1:B:188:ILE:HD12	1:B:514:ILE:CG2	2.07	0.83
1:B:451:PRO:HB2	1:B:469:VAL:CG2	2.13	0.78
1:B:372:ARG:NH1	1:B:405:SER:OG	2.15	0.78
1:A:228:PHE:CD2	1:A:321:VAL:HB	2.20	0.77
1:A:451:PRO:HB2	1:A:469:VAL:CG2	2.17	0.75
1:B:488:PRO:HG3	1:B:526:LEU:HB2	1.72	0.71
1:B:486:SER:HB3	1:B:495:ARG:HG3	1.72	0.71
1:A:240:CYS:SG	1:A:588:ILE:HG13	2.29	0.70
1:A:413:LEU:HG	1:A:475:LEU:HD22	1.74	0.70
1:B:524:VAL:HG12	1:B:561:THR:HG21	1.74	0.69
1:A:454:TYR:HE1	1:A:464:ILE:HD11	1.57	0.69
1:A:435:ARG:HB2	1:A:479:TRP:CD1	2.29	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:188:ILE:HD12	1:B:514:ILE:HG21	1.73	0.68
1:B:490:GLN:HB2	1:B:505:GLU:OE1	1.94	0.67
1:B:451:PRO:HB2	1:B:469:VAL:HG22	1.76	0.67
1:A:526:LEU:HD21	1:A:558:ALA:O	1.94	0.67
1:B:217:ILE:HG23	1:B:233:HIS:CD2	2.29	0.66
1:B:188:ILE:HD12	1:B:514:ILE:HG22	1.76	0.66
1:A:432:ALA:HB2	1:A:476:ARG:HG2	1.77	0.66
1:A:257:ASP:HB3	1:A:568:LEU:HD21	1.77	0.66
1:A:190:LEU:HD22	1:A:520:VAL:HG21	1.78	0.65
1:A:318:ARG:HD3	1:A:333:LYS:HB3	1.77	0.65
1:B:394:ASN:ND2	1:B:502:VAL:HG23	2.10	0.65
1:B:441:PRO:HD2	1:B:458:TYR:CE2	2.32	0.65
1:B:489:GLY:O	1:B:530:GLN:HA	1.95	0.65
1:A:577:LEU:HD11	1:A:590:PRO:HB3	1.79	0.65
1:A:581:TYR:CZ	1:A:588:ILE:HG12	2.32	0.65
1:B:188:ILE:CD1	1:B:514:ILE:CG2	2.75	0.65
1:A:250:ILE:HG21	1:A:293:THR:HG21	1.77	0.64
1:B:188:ILE:CD1	1:B:514:ILE:HG22	2.27	0.64
1:B:394:ASN:HD22	1:B:502:VAL:CG2	2.10	0.64
1:B:435:ARG:HB2	1:B:479:TRP:CD1	2.32	0.64
1:B:318:ARG:HG2	1:B:333:LYS:HB3	1.78	0.63
1:B:250:ILE:HG21	1:B:293:THR:HG21	1.80	0.63
1:B:552:ALA:HB1	1:B:556:THR:HG21	1.81	0.63
1:A:258:ARG:HH12	1:B:322:ARG:CD	2.10	0.62
1:B:435:ARG:HD3	1:B:479:TRP:CB	2.31	0.61
1:B:220:PRO:HB2	1:B:577:LEU:HD21	1.83	0.61
1:A:441:PRO:HD2	1:A:458:TYR:CE2	2.36	0.60
1:B:577:LEU:HD11	1:B:590:PRO:HB3	1.83	0.60
1:B:413:LEU:HG	1:B:475:LEU:HD22	1.82	0.60
1:B:462:THR:HG22	1:B:487:ARG:CZ	2.31	0.60
1:A:217:ILE:HG23	1:A:233:HIS:CD2	2.37	0.60
1:B:553:GLU:H	1:B:556:THR:HG21	1.66	0.60
1:A:524:VAL:HG12	1:A:561:THR:HG21	1.83	0.59
1:A:352:GLY:HA3	1:A:442:SER:O	2.02	0.59
1:A:397:LEU:HD22	1:A:403:SER:HB2	1.86	0.58
1:A:451:PRO:HB2	1:A:469:VAL:HG22	1.84	0.58
1:A:318:ARG:HD3	1:A:333:LYS:HD3	1.85	0.57
1:A:269:ASN:HD22	1:A:320:ALA:H	1.51	0.57
1:A:578:VAL:HG12	1:A:580:ILE:HG13	1.86	0.57
1:B:435:ARG:HD3	1:B:479:TRP:HB2	1.87	0.57
1:B:236:LYS:HA	1:B:245:ALA:HA	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:448:LEU:HB2	1:A:516:ARG:HH12	1.71	0.56
1:A:319:LEU:HD22	1:A:332:GLN:HG2	1.87	0.56
1:A:220:PRO:HB2	1:A:577:LEU:HD21	1.88	0.56
1:B:583:THR:HG23	1:B:584:GLY:H	1.71	0.56
1:A:514:ILE:HD11	1:A:520:VAL:HG12	1.87	0.56
1:A:239:SER:O	1:A:243:GLY:N	2.36	0.55
1:A:402:ASN:H	1:A:405:SER:HB3	1.71	0.55
1:A:337:ILE:HD13	1:A:426:LEU:HD11	1.87	0.55
1:B:257:ASP:HB3	1:B:568:LEU:HD21	1.89	0.55
1:B:339:LYS:HB2	1:B:424:ILE:O	2.06	0.55
1:A:602:SER:OG	1:A:603:GLU:OE1	2.25	0.55
1:B:338:THR:H	1:B:424:ILE:HG21	1.70	0.55
1:A:528:SER:OG	1:A:531:THR:HB	2.07	0.55
1:A:394:ASN:HD22	1:A:502:VAL:CG2	2.18	0.54
1:B:450:GLN:HB3	1:B:516:ARG:NH2	2.21	0.54
1:A:462:THR:HG22	1:A:487:ARG:HE	1.71	0.54
1:B:203:ILE:HD13	1:B:594:ALA:HB2	1.90	0.54
1:B:360:ASP:CB	1:B:418:LEU:HB2	2.28	0.54
1:B:244:ILE:HD12	1:B:244:ILE:H	1.73	0.54
1:A:271:TRP:CH2	1:A:273:PRO:HB3	2.43	0.54
1:A:537:PHE:HB2	1:A:551:LEU:HD11	1.90	0.53
1:B:451:PRO:HB2	1:B:469:VAL:HG21	1.90	0.53
1:A:345:TYR:CD1	1:A:370:LEU:HB2	2.43	0.53
1:B:509:ASN:HD22	1:B:561:THR:HG23	1.73	0.52
1:B:221:LEU:O	1:B:231:TYR:HA	2.10	0.52
1:B:484:VAL:HG23	1:B:485:ILE:HG12	1.92	0.52
1:B:509:ASN:ND2	1:B:561:THR:HG23	2.25	0.52
1:B:524:VAL:CG1	1:B:561:THR:HG21	2.39	0.51
1:A:462:THR:HG22	1:A:487:ARG:HH21	1.76	0.51
1:B:188:ILE:CD1	1:B:514:ILE:HG21	2.39	0.51
1:B:218:THR:HG21	1:B:236:LYS:HD2	1.92	0.51
1:B:256:LEU:O	1:B:264:SER:N	2.43	0.51
1:B:402:ASN:H	1:B:405:SER:HB3	1.74	0.51
1:B:300:VAL:O	3:B:702:NAG:H83	2.10	0.51
1:B:345:TYR:CD1	1:B:370:LEU:HB2	2.45	0.51
1:B:602:SER:OG	1:B:603:GLU:OE1	2.29	0.51
1:B:351:TYR:CE2	1:B:441:PRO:HD3	2.47	0.50
1:B:290:PHE:CE2	1:B:318:ARG:HD2	2.47	0.50
1:A:468:ASP:OD1	1:A:519:TRP:HH2	1.94	0.50
1:A:508:TYR:HB3	1:A:560:LYS:HE3	1.93	0.50
1:A:402:ASN:HB2	1:A:405:SER:N	2.27	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:463:MET:HE2	1:B:483:SER:HA	1.94	0.50
1:A:451:PRO:HB2	1:A:469:VAL:HG21	1.92	0.49
1:A:394:ASN:ND2	1:A:502:VAL:HG23	2.19	0.49
1:B:271:TRP:CH2	1:B:273:PRO:HB3	2.47	0.49
1:B:352:GLY:HA3	1:B:442:SER:O	2.12	0.49
1:B:239:SER:O	1:B:243:GLY:N	2.45	0.49
1:A:433:ASP:O	1:A:436:LEU:HD23	2.13	0.49
1:A:457:SER:O	1:A:487:ARG:HD3	2.12	0.49
1:A:355:GLY:HA2	1:A:444:ILE:HG23	1.95	0.49
1:A:524:VAL:CG1	1:A:561:THR:HG21	2.43	0.48
1:B:371:PRO:HA	1:B:407:TYR:HA	1.94	0.48
1:A:578:VAL:CG1	1:A:580:ILE:HG13	2.42	0.48
1:A:288:GLU:O	1:A:289:ASP:HB2	2.13	0.48
1:B:293:THR:HG22	1:B:317:ILE:HB	1.96	0.48
1:A:201:ARG:NH1	1:A:262:VAL:HG21	2.29	0.48
1:A:221:LEU:O	1:A:231:TYR:HA	2.13	0.48
1:A:541:LYS:HB2	1:A:541:LYS:HE3	1.57	0.48
1:A:419:SER:O	1:A:420:LEU:C	2.55	0.48
1:B:440:SER:HB2	1:B:458:TYR:HD2	1.78	0.48
1:B:355:GLY:HA2	1:B:444:ILE:HG23	1.96	0.48
1:A:528:SER:HB3	1:A:531:THR:O	2.15	0.47
1:B:233:HIS:CE1	1:B:249:ILE:HG12	2.49	0.47
1:B:288:GLU:O	1:B:289:ASP:HB2	2.13	0.47
1:A:553:GLU:H	1:A:556:THR:HG21	1.79	0.47
1:A:293:THR:CG2	1:A:317:ILE:HB	2.45	0.47
1:A:462:THR:HG22	1:A:487:ARG:NE	2.29	0.47
1:B:235:GLU:O	1:B:246:LYS:N	2.40	0.47
1:B:412:GLY:HA3	1:B:429:ILE:O	2.15	0.47
1:B:318:ARG:O	1:B:318:ARG:HG3	2.13	0.47
1:B:399:MET:HE1	1:B:436:LEU:HG	1.96	0.47
1:A:233:HIS:O	1:A:248:ARG:HD3	2.15	0.47
1:A:318:ARG:HH11	1:A:333:LYS:HB2	1.80	0.47
1:B:252:VAL:HG13	1:B:319:LEU:HD12	1.97	0.46
1:A:529:ASN:O	1:A:530:GLN:HB2	2.14	0.46
1:B:375:PHE:CG	1:B:408:ILE:HD13	2.50	0.46
1:B:287:HIS:HB3	1:B:292:TYR:HE2	1.79	0.46
1:A:602:SER:HG	1:A:603:GLU:H	1.64	0.46
1:B:486:SER:HB3	1:B:495:ARG:CG	2.42	0.46
1:A:318:ARG:CG	1:A:333:LYS:HB3	2.46	0.46
1:A:435:ARG:HD3	1:A:479:TRP:CB	2.45	0.46
1:A:448:LEU:HB2	1:A:516:ARG:NH1	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:495:ARG:O	1:B:496:PHE:HB2	2.16	0.46
1:B:581:TYR:CZ	1:B:588:ILE:HG12	2.51	0.46
1:A:287:HIS:HB3	1:A:292:TYR:CE2	2.51	0.45
1:A:263:PRO:HG3	1:A:568:LEU:HD22	1.97	0.45
1:A:351:TYR:CE2	1:A:441:PRO:HD3	2.51	0.45
1:B:313:SER:O	1:B:314:LEU:HD23	2.16	0.45
1:B:440:SER:HB2	1:B:458:TYR:CD2	2.50	0.45
1:A:463:MET:HE2	1:A:483:SER:HA	1.99	0.45
1:A:526:LEU:HD11	1:A:558:ALA:O	2.17	0.45
1:B:201:ARG:HD2	1:B:262:VAL:HG11	1.98	0.45
1:A:218:THR:HG21	1:A:236:LYS:HD2	1.99	0.45
1:B:437:THR:HG21	1:B:465:LYS:HD2	1.99	0.45
1:B:508:TYR:HB3	1:B:560:LYS:HE3	1.98	0.45
1:A:318:ARG:CD	1:A:333:LYS:HB3	2.47	0.44
1:B:431:ILE:HG13	1:B:475:LEU:HG	1.99	0.44
1:B:450:GLN:HB3	1:B:516:ARG:CZ	2.47	0.44
1:B:488:PRO:HG3	1:B:526:LEU:HD12	1.99	0.44
1:A:584:GLY:O	1:A:585:ASP:HB2	2.18	0.44
1:A:188:ILE:HD11	1:A:514:ILE:HD12	1.99	0.44
1:B:433:ASP:OD1	1:B:433:ASP:N	2.51	0.44
1:B:526:LEU:HD11	1:B:558:ALA:O	2.17	0.44
1:A:514:ILE:HD11	1:A:539:VAL:HG13	2.00	0.44
1:A:435:ARG:HD3	1:A:479:TRP:HB2	1.99	0.44
1:A:510:ASP:OD1	1:A:562:ILE:HA	2.17	0.44
1:B:290:PHE:CD2	1:B:318:ARG:HD2	2.53	0.44
1:A:287:HIS:HB3	1:A:292:TYR:HE2	1.82	0.44
1:A:293:THR:HG22	1:A:317:ILE:HB	1.99	0.44
1:B:223:ALA:HB2	1:B:354:SER:HB3	2.00	0.43
1:B:397:LEU:HD22	1:B:403:SER:HB2	2.00	0.43
1:B:553:GLU:H	1:B:556:THR:CG2	2.31	0.43
1:B:297:VAL:HB	1:B:313:SER:HB3	2.00	0.43
1:A:316:LEU:HD22	1:A:335:ILE:HD12	2.00	0.43
1:A:490:GLN:O	1:A:530:GLN:HG2	2.17	0.43
1:A:534:ASN:O	1:A:536:VAL:HG23	2.18	0.43
1:A:252:VAL:HG22	1:A:269:ASN:HB3	2.01	0.43
1:A:345:TYR:CE1	1:A:370:LEU:HB2	2.53	0.43
1:B:360:ASP:O	1:B:418:LEU:HD12	2.18	0.43
1:A:258:ARG:HD3	1:A:258:ARG:N	2.34	0.42
1:A:322:ARG:HD3	1:B:258:ARG:HH12	1.84	0.42
1:A:418:LEU:O	1:A:419:SER:C	2.62	0.42
1:B:402:ASN:HB2	1:B:405:SER:N	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:494:PRO:HB3	1:A:530:GLN:HG3	2.00	0.42
1:A:371:PRO:HA	1:A:407:TYR:HA	2.01	0.42
1:A:462:THR:HG22	1:A:487:ARG:NH2	2.34	0.42
1:B:252:VAL:HG22	1:B:269:ASN:HB3	2.02	0.42
1:A:440:SER:HB2	1:A:458:TYR:HD2	1.84	0.42
1:B:348:VAL:HG12	1:B:368:GLY:HA2	2.02	0.42
1:A:203:ILE:HD13	1:A:594:ALA:HB2	2.02	0.42
1:A:509:ASN:HD22	1:A:561:THR:HG23	1.85	0.42
1:A:512:PHE:CD2	1:A:564:ASP:HA	2.54	0.42
1:B:566:PHE:CE2	1:B:575:ILE:HG12	2.54	0.42
1:B:582:ASP:OD1	1:B:583:THR:HG22	2.20	0.42
1:A:361:THR:HG21	1:A:415:LYS:HE3	2.02	0.41
1:A:435:ARG:O	1:A:465:LYS:NZ	2.47	0.41
1:A:574:CYS:O	1:A:575:ILE:HD13	2.19	0.41
1:B:583:THR:HG23	1:B:584:GLY:N	2.34	0.41
1:A:318:ARG:HH11	1:A:333:LYS:CB	2.33	0.41
1:B:433:ASP:O	1:B:436:LEU:HB2	2.20	0.41
1:A:190:LEU:CD2	1:A:520:VAL:HG21	2.46	0.41
1:A:399:MET:HB3	1:A:408:ILE:HB	2.02	0.41
1:A:337:ILE:HD13	1:A:426:LEU:CD1	2.51	0.41
1:B:263:PRO:HG3	1:B:568:LEU:HD22	2.02	0.41
1:B:356:ILE:HD12	1:B:446:ASN:HD21	1.86	0.41
1:A:349:MET:HE3	1:A:350:PRO:O	2.21	0.41
1:B:360:ASP:HB3	1:B:418:LEU:CB	2.30	0.41
1:B:370:LEU:O	1:B:408:ILE:HG12	2.21	0.41
1:B:435:ARG:HD3	1:B:479:TRP:CG	2.55	0.41
1:B:509:ASN:HD22	1:B:561:THR:CG2	2.34	0.41
1:A:491:SER:HA	1:A:530:GLN:NE2	2.36	0.41
1:B:233:HIS:O	1:B:248:ARG:HD3	2.21	0.41
1:A:314:LEU:HD23	1:A:314:LEU:HA	1.79	0.40
1:A:322:ARG:HA	1:A:323:PRO:HD3	1.94	0.40
1:B:486:SER:CB	1:B:495:ARG:HG3	2.45	0.40
1:B:290:PHE:HE2	1:B:330:TYR:O	2.05	0.40
1:B:426:LEU:HD23	1:B:426:LEU:HA	1.81	0.40
1:A:464:ILE:HA	1:A:464:ILE:HD12	1.86	0.40
1:A:495:ARG:NH1	1:A:527:ASN:OD1	2.51	0.40
1:B:399:MET:CE	1:B:436:LEU:HG	2.51	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	394/416 (95%)	370 (94%)	22 (6%)	2 (0%)	24	55
1	B	397/416 (95%)	368 (93%)	28 (7%)	1 (0%)	36	65
All	All	791/832 (95%)	738 (93%)	50 (6%)	3 (0%)	30	60

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	585	ASP
1	B	353	PRO
1	A	353	PRO

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	360/373 (96%)	354 (98%)	6 (2%)	53	71
1	B	360/373 (96%)	358 (99%)	2 (1%)	78	81
All	All	720/746 (96%)	712 (99%)	8 (1%)	65	76

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	291	TYR
1	A	430	GLU
1	A	436	LEU

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Mol	Chain	Res	Type
1	A	541	LYS
1	A	542	ASP
1	A	587	VAL
1	B	249	ILE
1	B	291	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	226	ASN
1	A	233	HIS
1	A	247	GLN
1	A	394	ASN
1	A	402	ASN
1	B	233	HIS
1	B	394	ASN
1	B	427	GLN
1	B	548	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](#)

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	C	1	2,1	14,14,15	0.73	1 (7%)	17,19,21	1.79	3 (17%)
2	NAG	C	2	2	14,14,15	0.83	1 (7%)	17,19,21	0.90	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	C	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	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	NAG	C1-C2	-2.59	1.48	1.52
2	C	1	NAG	O5-C1	2.26	1.47	1.43

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	NAG	C1-O5-C5	4.77	118.57	112.19
2	C	1	NAG	O4-C4-C3	-3.97	101.02	110.38
2	C	1	NAG	O4-C4-C5	2.59	115.69	109.32
2	C	2	NAG	C3-C4-C5	2.35	114.50	110.23

There are no chirality outliers.

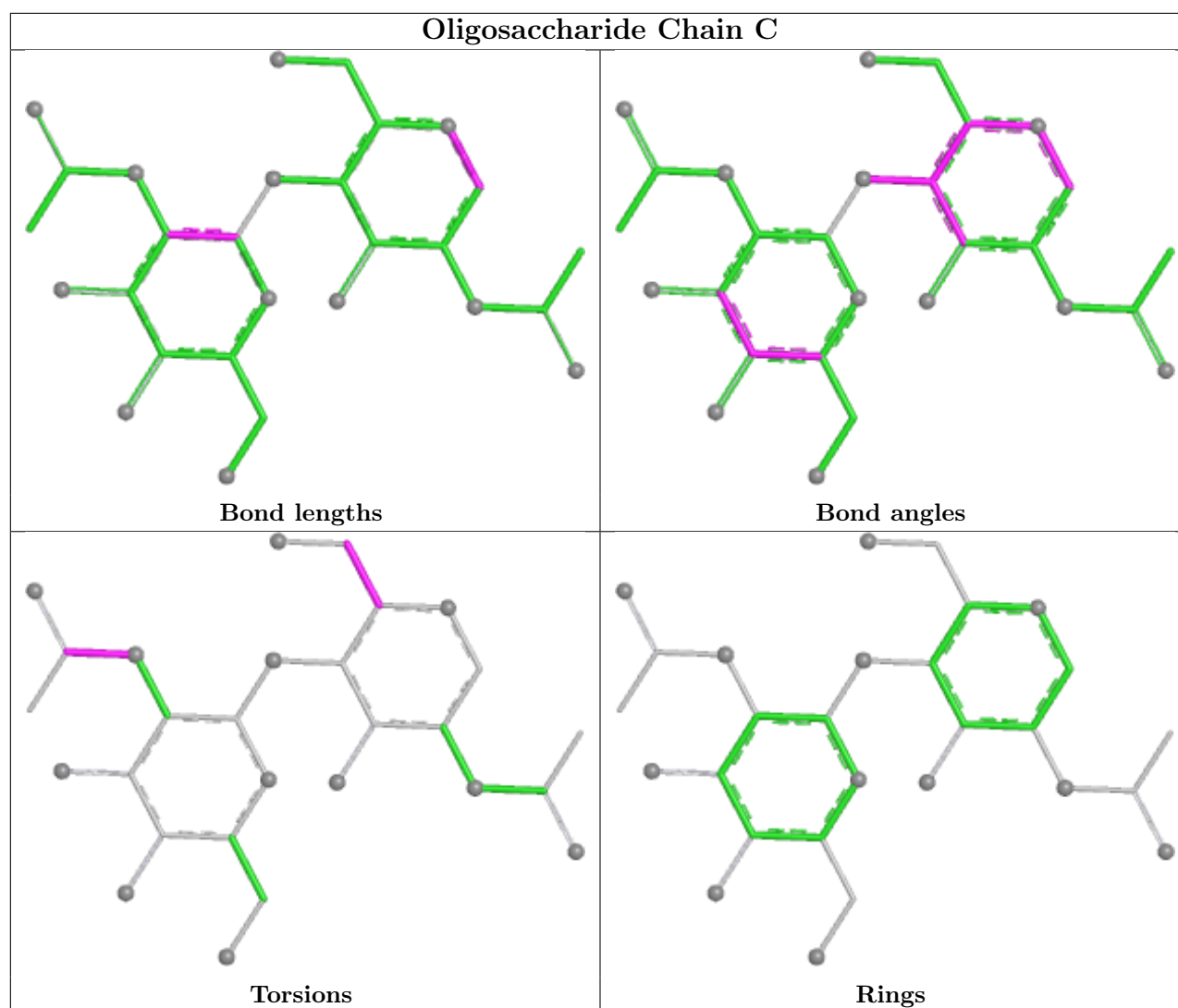
All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	1	NAG	O5-C5-C6-O6
2	C	1	NAG	C4-C5-C6-O6
2	C	2	NAG	C8-C7-N2-C2
2	C	2	NAG	O7-C7-N2-C2

There are no ring outliers.

No monomer is involved in short contacts.

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](#)

4 ligands 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	$\# Z > 2$
3	NAG	A	702	1	14,14,15	0.73	1 (7%)	17,19,21	0.61	0
3	NAG	A	701	1	14,14,15	1.02	1 (7%)	17,19,21	1.12	1 (5%)
3	NAG	B	702	1	14,14,15	0.42	0	17,19,21	0.70	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	B	701	1	14,14,15	1.01	1 (7%)	17,19,21	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
3	NAG	A	702	1	-	4/6/23/26	0/1/1/1
3	NAG	A	701	1	-	2/6/23/26	0/1/1/1
3	NAG	B	702	1	-	2/6/23/26	0/1/1/1
3	NAG	B	701	1	-	2/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	701	NAG	O5-C1	3.53	1.49	1.43
3	A	701	NAG	O5-C1	3.41	1.49	1.43
3	A	702	NAG	O5-C1	2.29	1.47	1.43

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	701	NAG	C2-N2-C7	2.61	126.40	122.90

There are no chirality outliers.

All (10) torsion outliers are listed below:

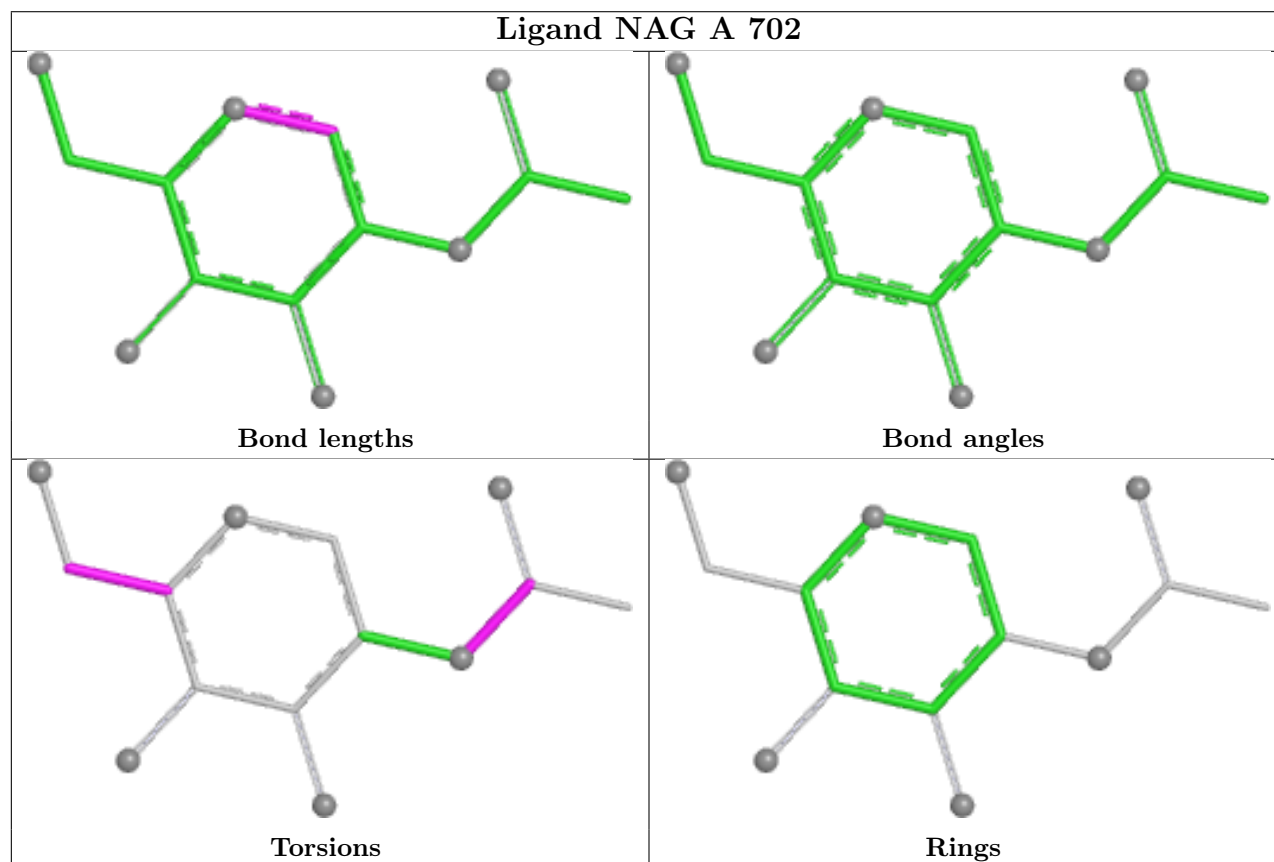
Mol	Chain	Res	Type	Atoms
3	A	702	NAG	C4-C5-C6-O6
3	B	701	NAG	O5-C5-C6-O6
3	A	702	NAG	O5-C5-C6-O6
3	B	701	NAG	C4-C5-C6-O6
3	A	702	NAG	C8-C7-N2-C2
3	A	702	NAG	O7-C7-N2-C2
3	B	702	NAG	C8-C7-N2-C2
3	B	702	NAG	O7-C7-N2-C2
3	A	701	NAG	C4-C5-C6-O6
3	A	701	NAG	C3-C2-N2-C7

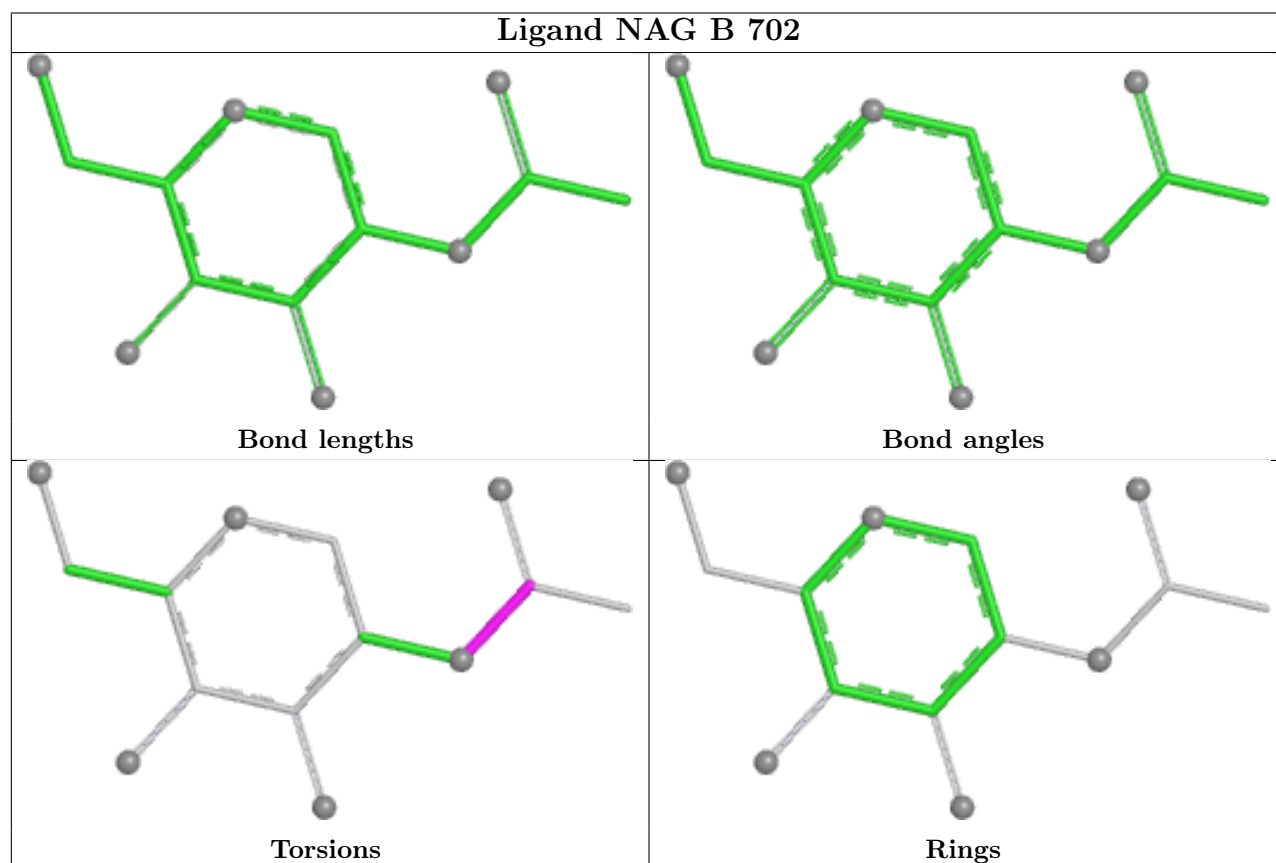
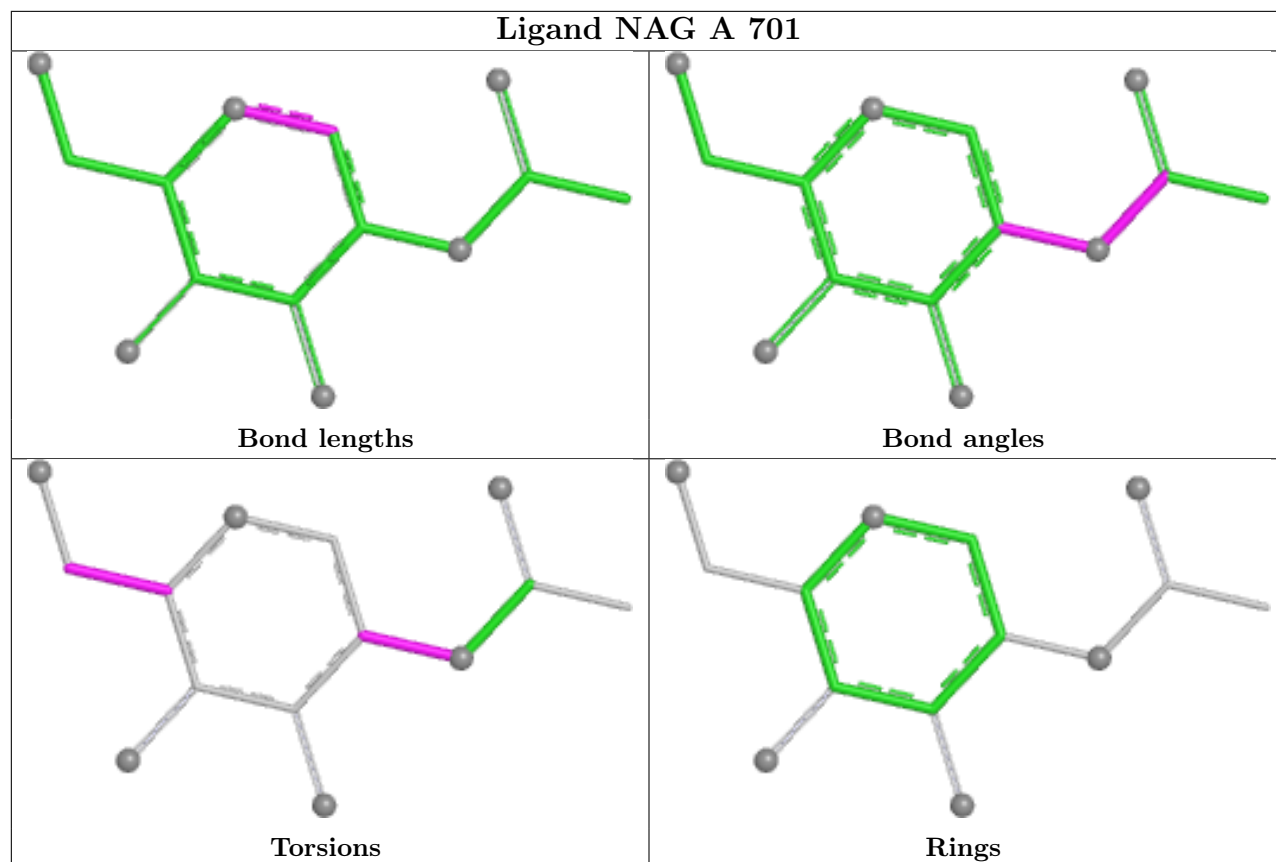
There are no ring outliers.

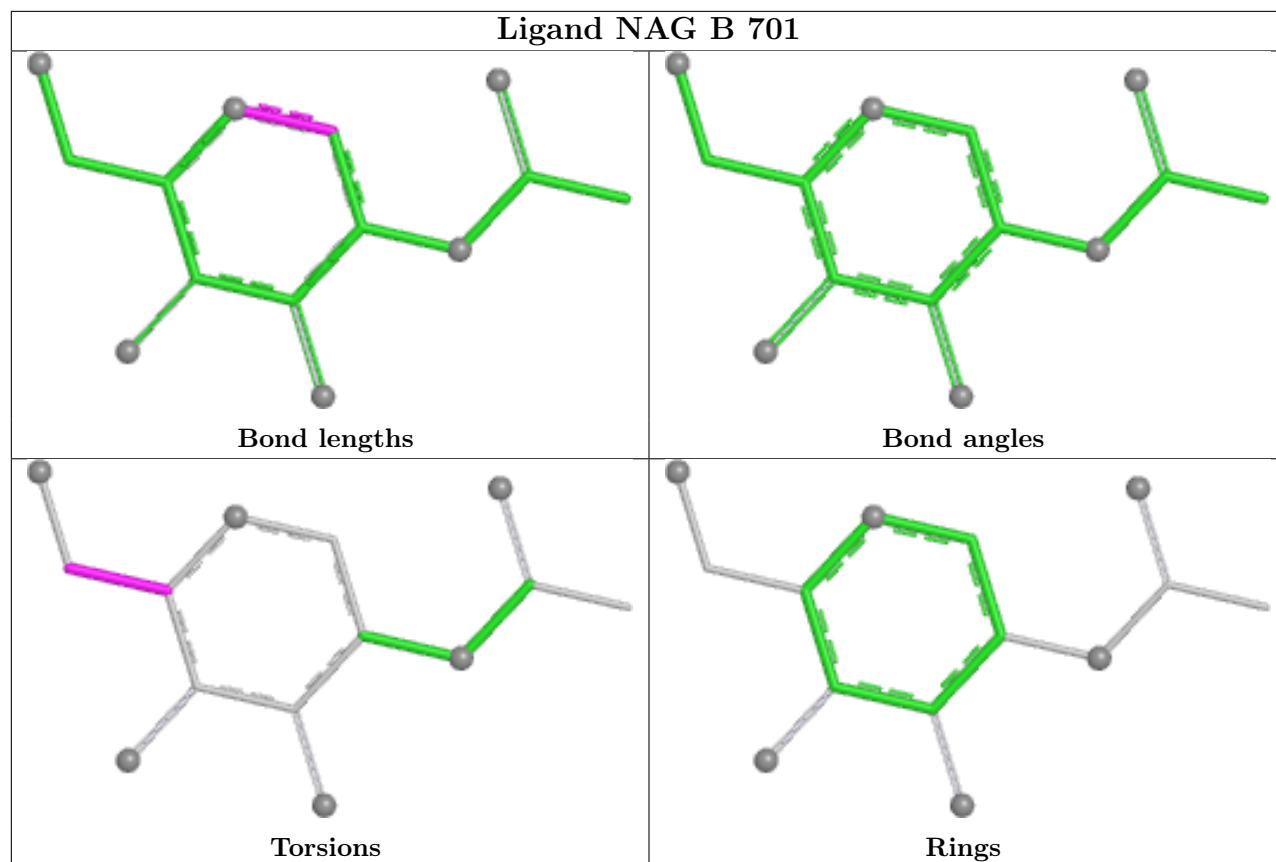
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	702	NAG	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	402/416 (96%)	0.61	28 (6%)	22 15	54, 78, 118, 155	0
1	B	403/416 (96%)	0.58	21 (5%)	33 22	65, 90, 128, 165	0
All	All	805/832 (96%)	0.60	49 (6%)	27 18	54, 84, 125, 165	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	555	ASP	4.5
1	A	533	GLU	4.1
1	A	424	ILE	3.9
1	B	308	THR	3.6
1	B	346	ASP	3.5
1	A	519	TRP	3.5
1	A	470	ASP	3.5
1	A	482	ASN	3.3
1	B	425	ILE	3.2
1	B	295	CYS	3.1
1	A	284	SER	3.0
1	A	532	ALA	3.0
1	B	307	SER	2.7
1	A	589	ARG	2.7
1	A	585	ASP	2.7
1	A	204	SER	2.6
1	B	305	LEU	2.6
1	A	559	GLN	2.6
1	B	204	SER	2.6
1	B	493	CYS	2.5
1	B	492	GLN	2.5
1	A	286	TYR	2.5
1	B	589	ARG	2.5
1	A	406	HIS	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	188	ILE	2.4
1	B	371	PRO	2.4
1	A	265	MET	2.4
1	A	240	CYS	2.4
1	A	588	ILE	2.3
1	B	519	TRP	2.3
1	A	434	ASN	2.3
1	A	534	ASN	2.3
1	B	542	ASP	2.3
1	A	322	ARG	2.2
1	B	265	MET	2.2
1	A	602	SER	2.2
1	B	533	GLU	2.2
1	B	432	ALA	2.2
1	A	289	ASP	2.2
1	B	196	THR	2.2
1	A	542	ASP	2.2
1	B	457	SER	2.1
1	B	188	ILE	2.1
1	A	479	TRP	2.1
1	A	448	LEU	2.1
1	B	407	TYR	2.1
1	B	601	CYS	2.1
1	A	386	HIS	2.0
1	A	579	GLU	2.0

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

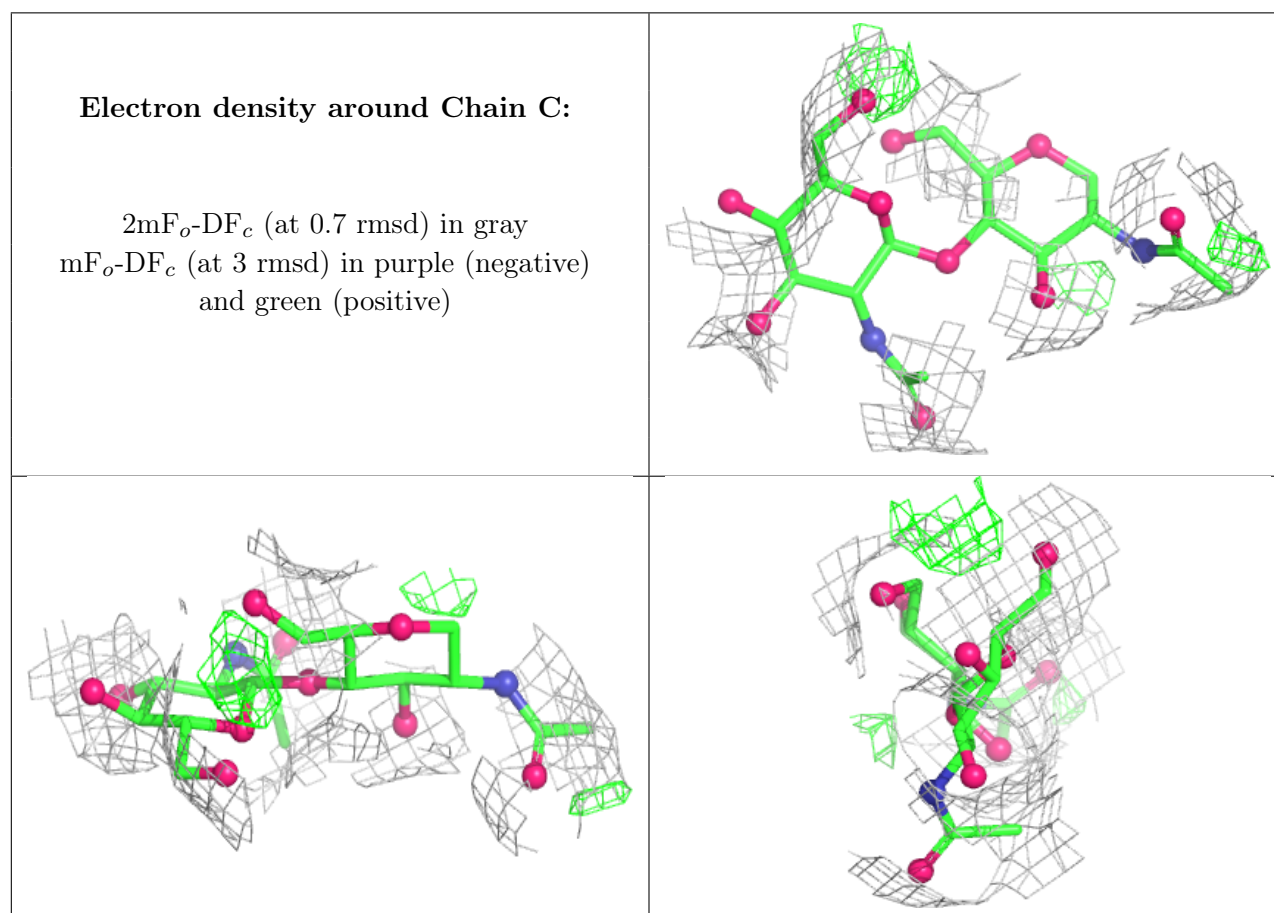
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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(Å ²)	Q<0.9
2	NAG	C	2	14/15	0.40	0.14	140,150,163,177	0
2	NAG	C	1	14/15	0.76	0.10	88,113,131,136	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](#)

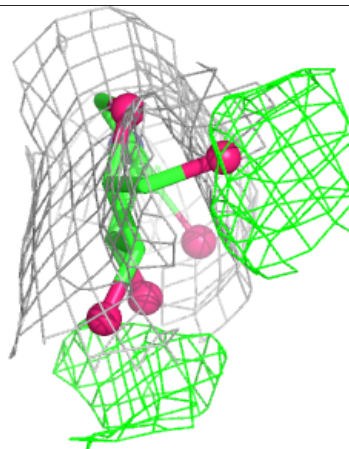
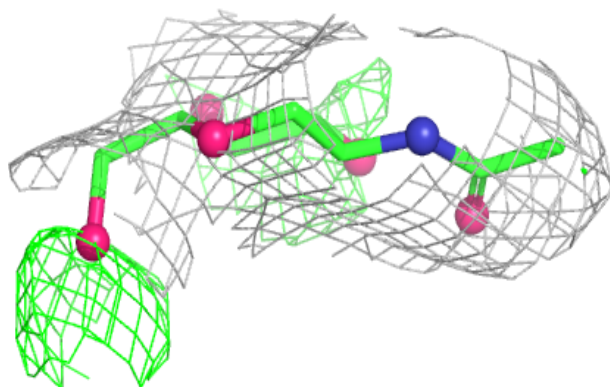
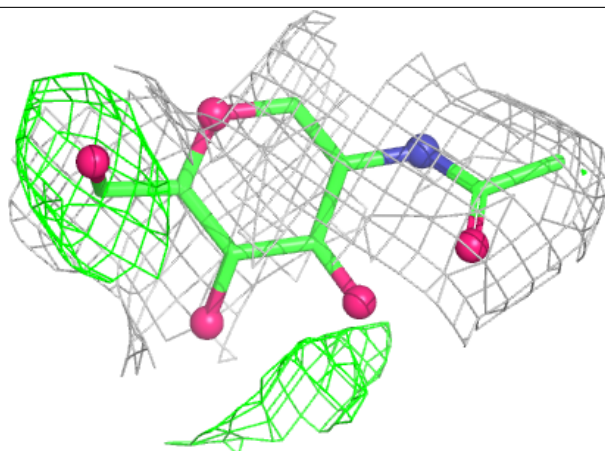
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NAG	A	701	14/15	0.66	0.19	71,90,99,100	0
3	NAG	B	701	14/15	0.66	0.17	85,95,99,103	0
3	NAG	B	702	14/15	0.72	0.12	110,120,136,143	0
3	NAG	A	702	14/15	0.79	0.10	73,90,97,97	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.

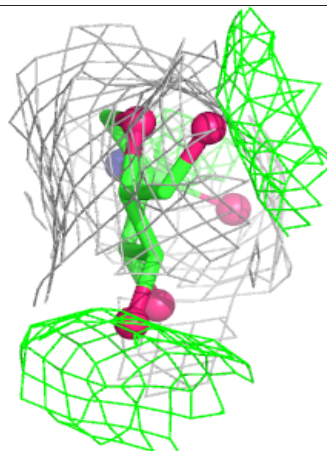
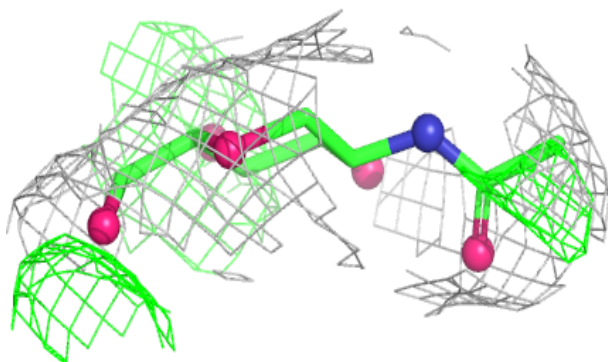
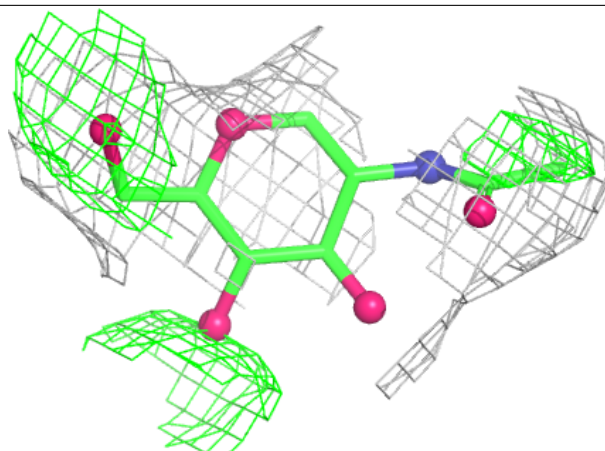
Electron density around NAG A 701:

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

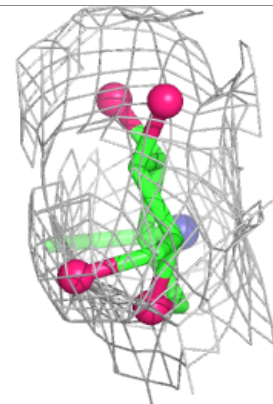
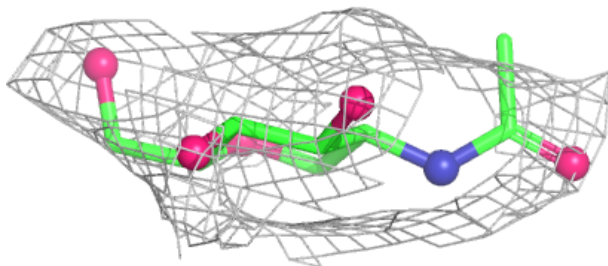
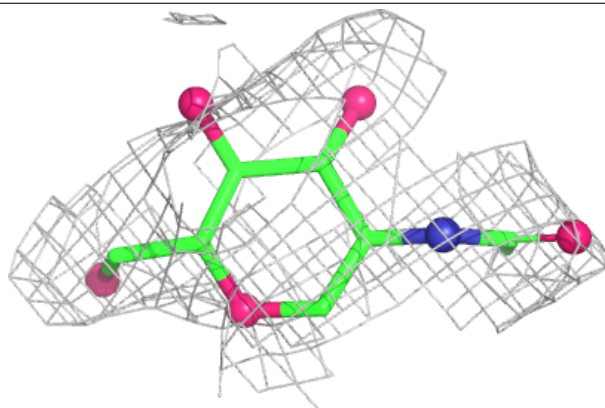


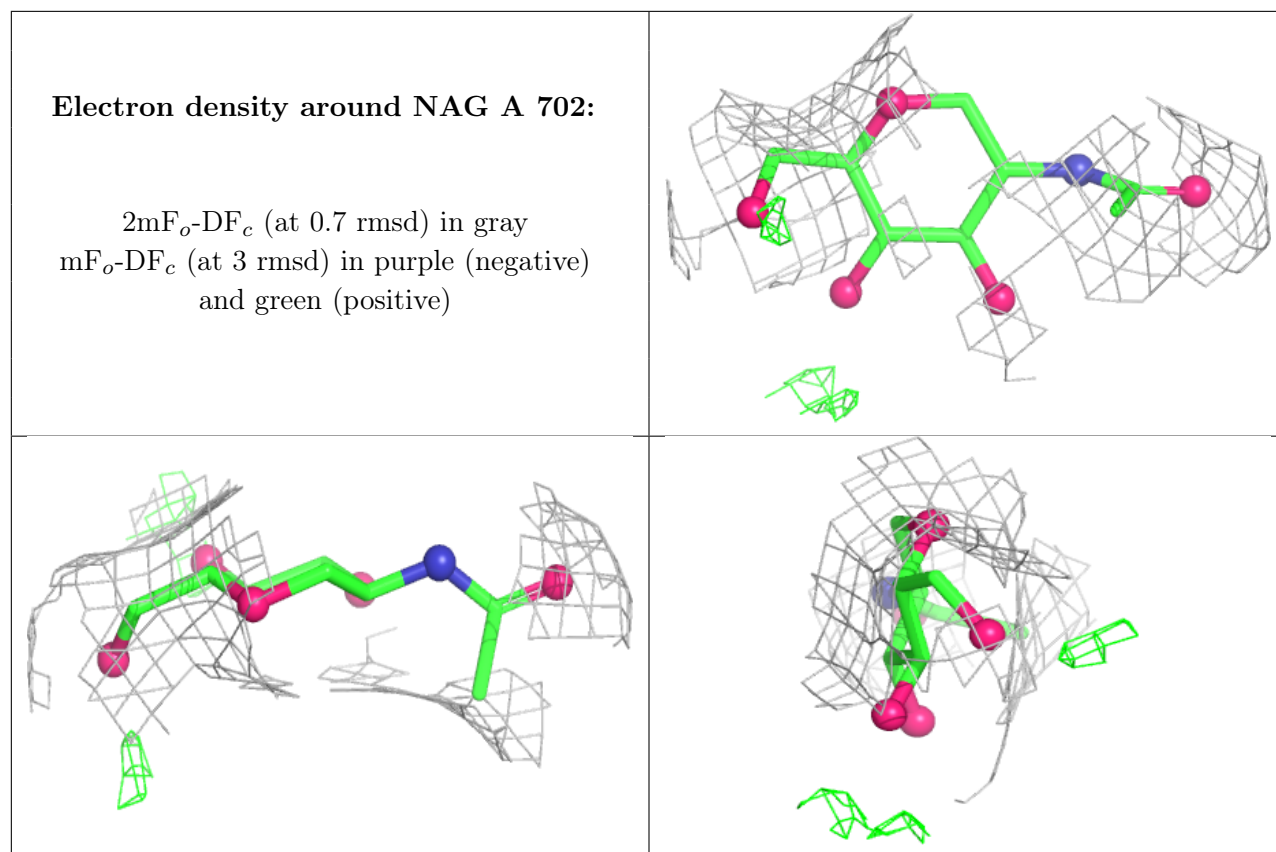
Electron density around NAG B 701:

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

**Electron density around NAG B 702:**

$2mF_o-DF_c$ (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.