



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:48 PM UTC

PDB ID : 4COF / pdb_00004cof
Title : Crystal structure of a human gamma-aminobutyric acid receptor, the GABA(A)R-beta3 homopentamer
Authors : Miller, P.S.; Aricescu, A.R.
Deposited on : 2014-01-28
Resolution : 2.97 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)	:	2.0
EDS	:	3.0
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

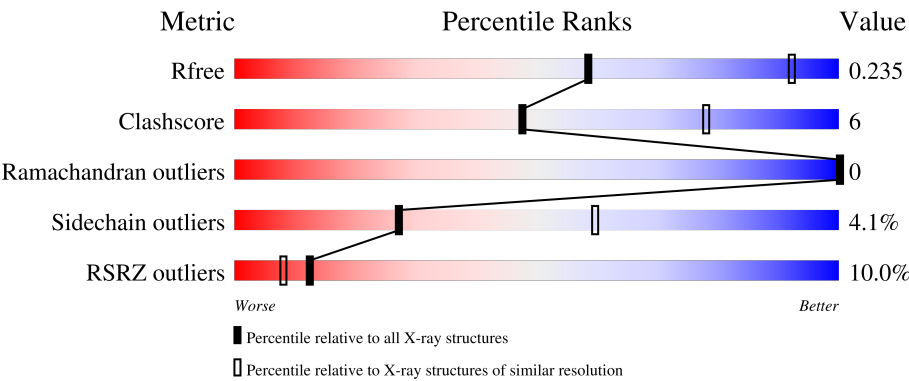
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.97 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.



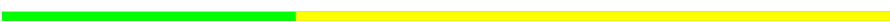
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3580 (3.00-2.96)
Clashscore	190562	3904 (3.00-2.96)
Ramachandran outliers	187476	3761 (3.00-2.96)
Sidechain outliers	187428	3764 (3.00-2.96)
RSRZ outliers	180081	3579 (3.00-2.96)

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	355	11% 80% 15% . .
1	B	355	6% 75% 17% . 7%
1	C	355	10% 75% 17% . 7%
1	D	355	11% 72% 18% . 6%
1	E	355	9% 75% 16% . 7%

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Mol	Chain	Length	Quality of chain
2	F	3	 33% 67%
2	G	3	 33% 67%
2	H	3	 33% 67%
2	I	3	 33% 67%
2	J	3	 100%

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	BEN	A	500	-	X	-	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13972 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 GAMMA-AMINOBUTYRIC ACID RECEPTOR SUBUNIT BETA-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	340	Total	C	N	O	S	0	0	0
			2779	1816	455	492	16			
1	B	331	Total	C	N	O	S	0	0	0
			2714	1779	443	477	15			
1	C	331	Total	C	N	O	S	0	0	0
			2714	1779	443	477	15			
1	D	332	Total	C	N	O	S	0	0	0
			2722	1783	445	479	15			
1	E	331	Total	C	N	O	S	0	0	0
			2714	1779	443	477	15			

There are 105 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLU	-	expression tag	UNP P28472
A	-1	THR	-	expression tag	UNP P28472
A	0	GLY	-	expression tag	UNP P28472
A	449	GLY	-	expression tag	UNP P28472
A	450	ALA	-	expression tag	UNP P28472
A	451	THR	-	expression tag	UNP P28472
A	452	GLU	-	expression tag	UNP P28472
A	453	THR	-	expression tag	UNP P28472
A	454	SER	-	expression tag	UNP P28472
A	455	GLN	-	expression tag	UNP P28472
A	456	VAL	-	expression tag	UNP P28472
A	457	ALA	-	expression tag	UNP P28472
A	458	PRO	-	expression tag	UNP P28472
A	459	ALA	-	expression tag	UNP P28472
A	308	SER	-	linker	UNP P28472
A	309	GLN	-	linker	UNP P28472
A	310	PRO	-	linker	UNP P28472
A	311	ALA	-	linker	UNP P28472

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Chain	Residue	Modelled	Actual	Comment	Reference
A	312	ARG	-	linker	UNP P28472
A	313	ALA	-	linker	UNP P28472
A	314	ALA	-	linker	UNP P28472
B	-2	GLU	-	expression tag	UNP P28472
B	-1	THR	-	expression tag	UNP P28472
B	0	GLY	-	expression tag	UNP P28472
B	449	GLY	-	expression tag	UNP P28472
B	450	ALA	-	expression tag	UNP P28472
B	451	THR	-	expression tag	UNP P28472
B	452	GLU	-	expression tag	UNP P28472
B	453	THR	-	expression tag	UNP P28472
B	454	SER	-	expression tag	UNP P28472
B	455	GLN	-	expression tag	UNP P28472
B	456	VAL	-	expression tag	UNP P28472
B	457	ALA	-	expression tag	UNP P28472
B	458	PRO	-	expression tag	UNP P28472
B	459	ALA	-	expression tag	UNP P28472
B	308	SER	-	linker	UNP P28472
B	309	GLN	-	linker	UNP P28472
B	310	PRO	-	linker	UNP P28472
B	311	ALA	-	linker	UNP P28472
B	312	ARG	-	linker	UNP P28472
B	313	ALA	-	linker	UNP P28472
B	314	ALA	-	linker	UNP P28472
C	-2	GLU	-	expression tag	UNP P28472
C	-1	THR	-	expression tag	UNP P28472
C	0	GLY	-	expression tag	UNP P28472
C	449	GLY	-	expression tag	UNP P28472
C	450	ALA	-	expression tag	UNP P28472
C	451	THR	-	expression tag	UNP P28472
C	452	GLU	-	expression tag	UNP P28472
C	453	THR	-	expression tag	UNP P28472
C	454	SER	-	expression tag	UNP P28472
C	455	GLN	-	expression tag	UNP P28472
C	456	VAL	-	expression tag	UNP P28472
C	457	ALA	-	expression tag	UNP P28472
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C	459	ALA	-	expression tag	UNP P28472
C	308	SER	-	linker	UNP P28472
C	309	GLN	-	linker	UNP P28472
C	310	PRO	-	linker	UNP P28472
C	311	ALA	-	linker	UNP P28472

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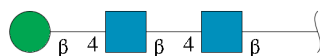
Chain	Residue	Modelled	Actual	Comment	Reference
C	312	ARG	-	linker	UNP P28472
C	313	ALA	-	linker	UNP P28472
C	314	ALA	-	linker	UNP P28472
D	-2	GLU	-	expression tag	UNP P28472
D	-1	THR	-	expression tag	UNP P28472
D	0	GLY	-	expression tag	UNP P28472
D	449	GLY	-	expression tag	UNP P28472
D	450	ALA	-	expression tag	UNP P28472
D	451	THR	-	expression tag	UNP P28472
D	452	GLU	-	expression tag	UNP P28472
D	453	THR	-	expression tag	UNP P28472
D	454	SER	-	expression tag	UNP P28472
D	455	GLN	-	expression tag	UNP P28472
D	456	VAL	-	expression tag	UNP P28472
D	457	ALA	-	expression tag	UNP P28472
D	458	PRO	-	expression tag	UNP P28472
D	459	ALA	-	expression tag	UNP P28472
D	308	SER	-	linker	UNP P28472
D	309	GLN	-	linker	UNP P28472
D	310	PRO	-	linker	UNP P28472
D	311	ALA	-	linker	UNP P28472
D	312	ARG	-	linker	UNP P28472
D	313	ALA	-	linker	UNP P28472
D	314	ALA	-	linker	UNP P28472
E	-2	GLU	-	expression tag	UNP P28472
E	-1	THR	-	expression tag	UNP P28472
E	0	GLY	-	expression tag	UNP P28472
E	449	GLY	-	expression tag	UNP P28472
E	450	ALA	-	expression tag	UNP P28472
E	451	THR	-	expression tag	UNP P28472
E	452	GLU	-	expression tag	UNP P28472
E	453	THR	-	expression tag	UNP P28472
E	454	SER	-	expression tag	UNP P28472
E	455	GLN	-	expression tag	UNP P28472
E	456	VAL	-	expression tag	UNP P28472
E	457	ALA	-	expression tag	UNP P28472
E	458	PRO	-	expression tag	UNP P28472
E	459	ALA	-	expression tag	UNP P28472
E	308	SER	-	linker	UNP P28472
E	309	GLN	-	linker	UNP P28472
E	310	PRO	-	linker	UNP P28472
E	311	ALA	-	linker	UNP P28472

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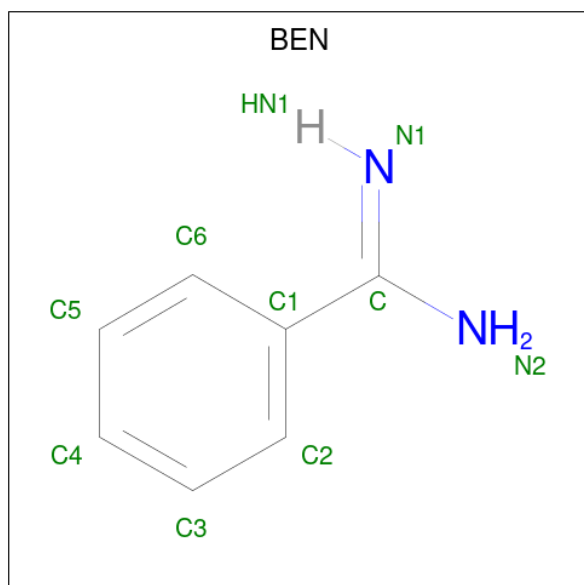
Chain	Residue	Modelled	Actual	Comment	Reference
E	312	ARG	-	linker	UNP P28472
E	313	ALA	-	linker	UNP P28472
E	314	ALA	-	linker	UNP P28472

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



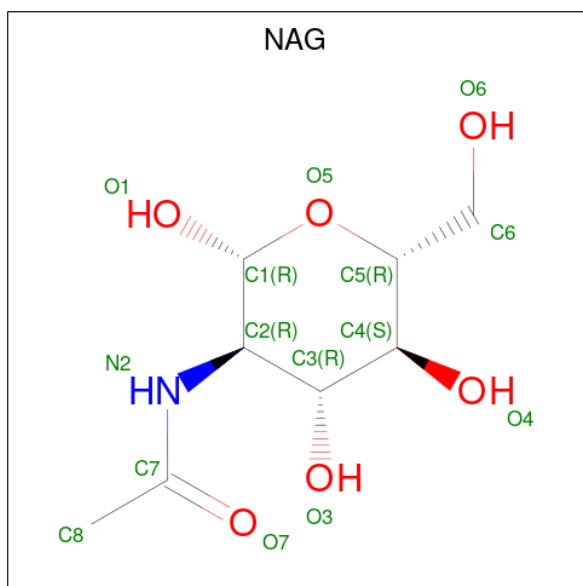
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	F	3	Total	C	N	O	0	0	0
			39	22	2	15			
2	G	3	Total	C	N	O	0	0	0
			39	22	2	15			
2	H	3	Total	C	N	O	0	0	0
			39	22	2	15			
2	I	3	Total	C	N	O	0	0	0
			39	22	2	15			
2	J	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 3 is BENZAMIDINE (CCD ID: BEN) (formula: C₇H₈N₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			9	7	2		
3	B	1	Total	C	N	0	0
			9	7	2		
3	C	1	Total	C	N	0	0
			9	7	2		
3	D	1	Total	C	N	0	0
			9	7	2		
3	E	1	Total	C	N	0	0
			9	7	2		

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



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

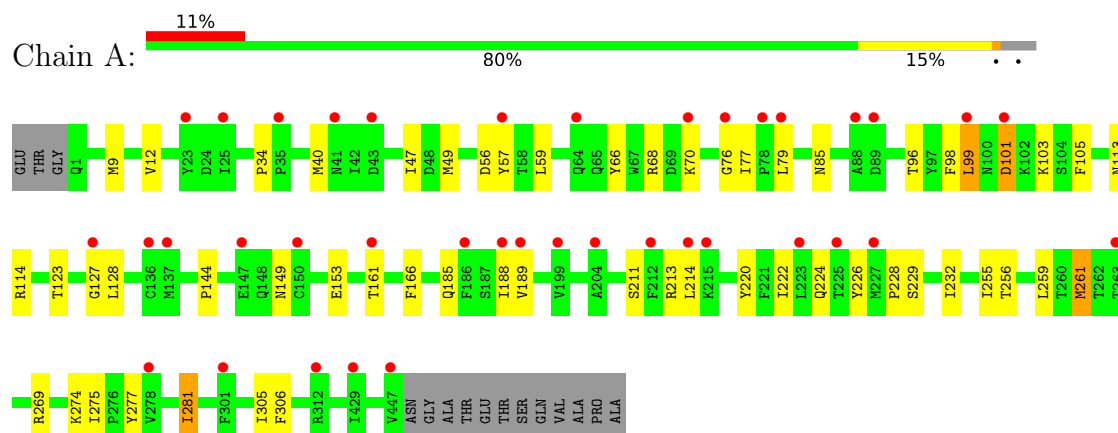
- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total 1	Cl 1	0	0
5	B	1	Total 1	Cl 1	0	0
5	C	1	Total 1	Cl 1	0	0
5	D	1	Total 1	Cl 1	0	0
5	E	1	Total 1	Cl 1	0	0

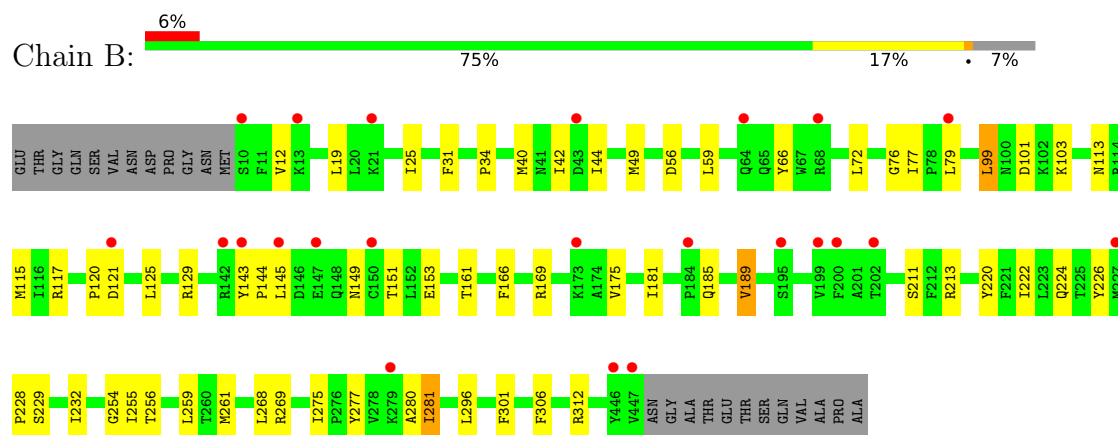
3 Residue-property plots [i](#)

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.

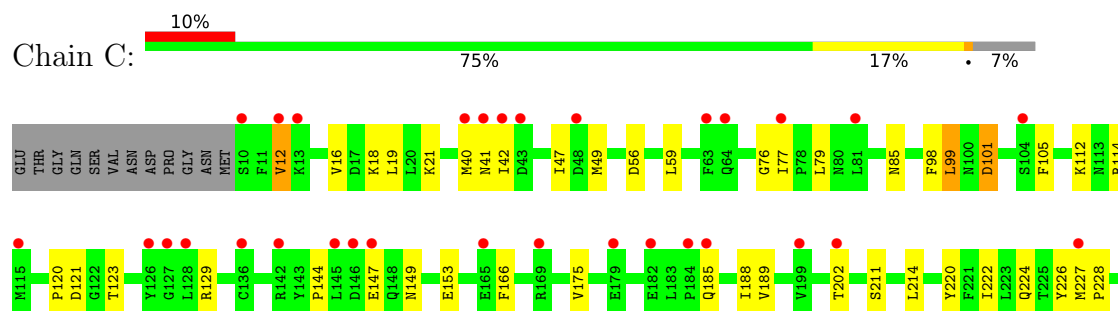
• Molecule 1: GAMMA-AMINOBUTYRIC ACID RECEPTOR SUBUNIT BETA-3



• Molecule 1: GAMMA-AMINOBUTYRIC ACID RECEPTOR SUBUNIT BETA-3

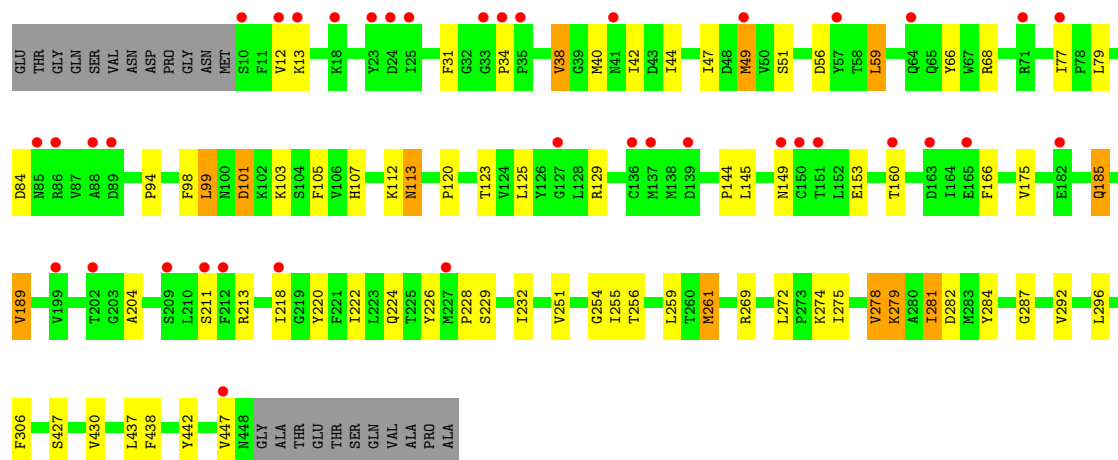
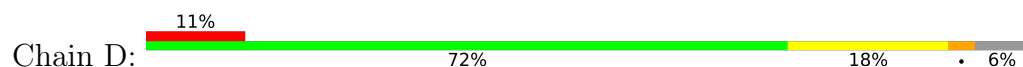


• Molecule 1: GAMMA-AMINOBUTYRIC ACID RECEPTOR SUBUNIT BETA-3

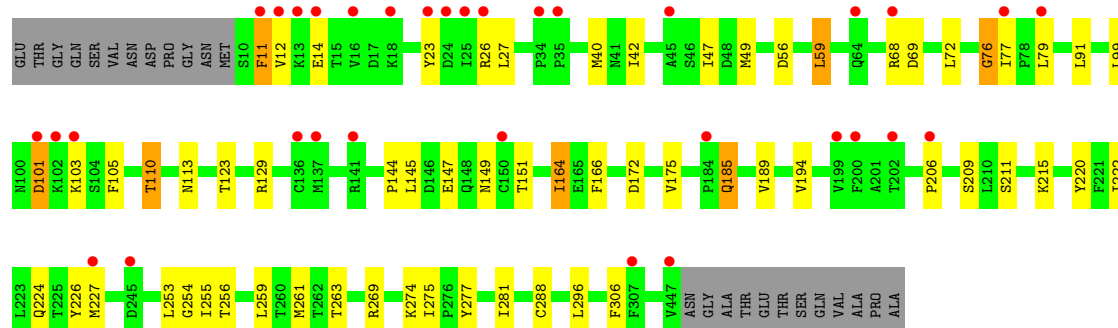




• Molecule 1: GAMMA-AMINOBUTYRIC ACID RECEPTOR SUBUNIT BETA-3



• Molecule 1: GAMMA-AMINOBUTYRIC ACID RECEPTOR SUBUNIT BETA-3



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



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



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

Chain H:  33% 67%

MAG1
MAG2
BMA3

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

Chain I:  33% 67%

MAG1
MAG2
BMA3

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

Chain J:  100%

MAG1
MAG2
BMA3

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	174.10Å 108.90Å 207.44Å 90.00° 107.43° 90.00°	Depositor
Resolution (Å)	40.00 – 2.97 40.00 – 2.97	Depositor EDS
% Data completeness (in resolution range)	99.8 (40.00-2.97) 99.7 (40.00-2.97)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.67 (at 2.96Å)	Xtriage
Refinement program	BUSTER 2.11.2	Depositor
R, R_{free}	0.205 , 0.226 0.215 , 0.235	Depositor DCC
R_{free} test set	3837 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	81.8	Xtriage
Anisotropy	0.684	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 45.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	13972	wwPDB-VP
Average B, all atoms (Å ²)	102.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BEN, NAG, CL, BMA

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	0.87	1/2854 (0.0%)	1.21	9/3886 (0.2%)
1	B	0.88	1/2788 (0.0%)	1.22	12/3796 (0.3%)
1	C	0.88	0/2788	1.24	11/3796 (0.3%)
1	D	0.89	1/2796 (0.0%)	1.25	11/3807 (0.3%)
1	E	0.87	0/2788	1.24	11/3796 (0.3%)
All	All	0.88	3/14014 (0.0%)	1.23	54/19081 (0.3%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	34	PRO	CA-C	5.45	1.54	1.51
1	A	34	PRO	CA-C	5.25	1.54	1.51
1	D	34	PRO	CA-C	5.06	1.54	1.51

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	11	PHE	CA-CB-CG	7.69	121.49	113.80
1	C	166	PHE	CA-CB-CG	7.04	120.84	113.80
1	C	306	PHE	CA-CB-CG	7.02	120.82	113.80
1	E	166	PHE	CA-CB-CG	6.84	120.64	113.80
1	E	101	ASP	CA-CB-CG	6.81	119.41	112.60
1	D	166	PHE	CA-CB-CG	6.78	120.58	113.80
1	D	306	PHE	CA-CB-CG	6.73	120.53	113.80
1	A	113	ASN	CA-CB-CG	6.61	119.21	112.60
1	A	306	PHE	CA-CB-CG	6.54	120.34	113.80
1	D	47	ILE	CA-C-N	6.46	128.94	120.28
1	D	47	ILE	C-N-CA	6.46	128.94	120.28
1	B	306	PHE	CA-CB-CG	6.45	120.25	113.80
1	D	120	PRO	CA-C-N	6.44	128.91	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	120	PRO	C-N-CA	6.44	128.91	120.28
1	D	113	ASN	CA-CB-CG	6.44	119.04	112.60
1	E	306	PHE	CA-CB-CG	6.41	120.21	113.80
1	C	120	PRO	CA-C-N	6.34	129.40	120.28
1	C	120	PRO	C-N-CA	6.34	129.40	120.28
1	B	120	PRO	CA-C-N	6.29	129.56	120.38
1	B	120	PRO	C-N-CA	6.29	129.56	120.38
1	E	172	ASP	CA-CB-CG	6.27	118.87	112.60
1	B	166	PHE	CA-CB-CG	6.18	119.98	113.80
1	A	166	PHE	CA-CB-CG	6.17	119.97	113.80
1	C	121	ASP	CA-CB-CG	6.05	118.65	112.60
1	A	101	ASP	CA-CB-CG	5.99	118.59	112.60
1	D	101	ASP	CA-CB-CG	5.90	118.50	112.60
1	C	149	ASN	CA-CB-CG	5.85	118.45	112.60
1	C	101	ASP	CA-CB-CG	5.83	118.43	112.60
1	A	149	ASN	CA-CB-CG	5.78	118.38	112.60
1	B	101	ASP	CA-CB-CG	5.56	118.16	112.60
1	E	149	ASN	CA-CB-CG	5.55	118.16	112.60
1	B	113	ASN	CA-CB-CG	5.54	118.14	112.60
1	E	113	ASN	CA-CB-CG	5.52	118.12	112.60
1	B	121	ASP	CA-CB-CG	5.50	118.11	112.60
1	B	149	ASN	CA-CB-CG	5.50	118.10	112.60
1	C	47	ILE	CA-C-N	5.47	128.16	120.28
1	C	47	ILE	C-N-CA	5.47	128.16	120.28
1	D	149	ASN	CA-CB-CG	5.47	118.07	112.60
1	B	56	ASP	CA-CB-CG	5.39	117.99	112.60
1	C	76	GLY	N-CA-C	5.37	118.98	112.49
1	A	47	ILE	CA-C-N	5.34	127.97	120.28
1	A	47	ILE	C-N-CA	5.34	127.97	120.28
1	B	25	ILE	CA-C-N	5.31	127.93	120.28
1	B	25	ILE	C-N-CA	5.31	127.93	120.28
1	A	56	ASP	CA-CB-CG	5.22	117.82	112.60
1	D	447	VAL	N-CA-C	-5.21	105.74	113.39
1	D	56	ASP	CA-CB-CG	5.21	117.81	112.60
1	C	56	ASP	CA-CB-CG	5.20	117.80	112.60
1	E	56	ASP	CA-CB-CG	5.20	117.80	112.60
1	A	76	GLY	N-CA-C	5.18	118.76	112.49
1	B	76	GLY	N-CA-C	5.16	118.73	112.49
1	E	76	GLY	N-CA-C	5.12	118.69	112.49
1	E	14	GLU	CA-C-N	5.01	126.95	120.44
1	E	14	GLU	C-N-CA	5.01	126.95	120.44

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	2779	0	2769	36	0
1	B	2714	0	2710	38	0
1	C	2714	0	2710	34	0
1	D	2722	0	2716	51	0
1	E	2714	0	2710	43	0
2	F	39	0	34	0	0
2	G	39	0	34	0	0
2	H	39	0	34	0	0
2	I	39	0	34	0	0
2	J	39	0	34	0	0
3	A	9	0	7	0	0
3	B	9	0	7	0	0
3	C	9	0	7	0	0
3	D	9	0	7	0	0
3	E	9	0	7	0	0
4	A	28	0	26	0	0
4	B	14	0	13	0	0
4	C	14	0	13	0	0
4	D	14	0	13	0	0
4	E	14	0	13	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
5	E	1	0	0	0	0
All	All	13972	0	13898	156	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 (156) 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:E:68:ARG:HG2	1:E:123:THR:HG22	1.41	1.01
1:D:144:PRO:HD3	1:D:281:ILE:HG12	1.52	0.90
1:A:85:ASN:HB2	1:A:114:ARG:HG3	1.61	0.83
1:D:66:TYR:CE2	1:D:125:LEU:HD13	2.18	0.78
1:A:256:THR:HG23	1:B:259:LEU:HD11	1.65	0.78
1:B:66:TYR:CE2	1:B:125:LEU:HD13	2.19	0.78
1:D:256:THR:HG23	1:E:259:LEU:HD11	1.67	0.76
1:A:259:LEU:HD11	1:E:256:THR:HG23	1.66	0.76
1:D:44:ILE:HD12	1:D:59:LEU:HD11	1.70	0.74
1:D:129:ARG:HH22	1:E:101:ASP:HB3	1.53	0.73
1:C:256:THR:HG23	1:D:259:LEU:HD11	1.70	0.71
1:B:228:PRO:O	1:B:232:ILE:HG12	1.91	0.70
1:B:256:THR:HG23	1:C:259:LEU:HD11	1.72	0.70
1:D:66:TYR:CZ	1:D:125:LEU:HD13	2.28	0.68
1:A:224:GLN:HG2	1:B:269:ARG:HD2	1.76	0.65
1:D:256:THR:HG21	1:E:255:ILE:HG23	1.79	0.65
1:D:272:LEU:HD13	1:D:279:LYS:HE3	1.77	0.65
1:E:151:THR:HG22	1:E:211:SER:HB3	1.80	0.64
1:C:256:THR:HG21	1:D:255:ILE:HG23	1.82	0.62
1:C:254:GLY:HA3	1:C:296:LEU:HD13	1.82	0.60
1:A:259:LEU:HD11	1:E:256:THR:CG2	2.32	0.60
1:A:256:THR:HG21	1:B:255:ILE:HG23	1.84	0.59
1:D:144:PRO:HD2	1:D:145:LEU:HD12	1.84	0.58
1:D:224:GLN:HG2	1:E:269:ARG:HD2	1.85	0.58
1:A:269:ARG:HD2	1:E:224:GLN:HG2	1.84	0.58
1:E:144:PRO:HD3	1:E:281:ILE:HB	1.85	0.57
1:A:57:TYR:HE2	1:A:59:LEU:HD23	1.69	0.57
1:C:256:THR:HG21	1:D:255:ILE:CG2	2.34	0.57
1:D:275:ILE:HD11	1:D:279:LYS:HE2	1.87	0.56
1:B:66:TYR:CZ	1:B:125:LEU:HD13	2.41	0.56
1:A:96:THR:O	1:E:110:THR:HB	2.05	0.56
1:E:11:PHE:HZ	1:E:76:GLY:HA3	1.71	0.56
1:D:256:THR:HG21	1:E:255:ILE:CG2	2.36	0.56
1:B:256:THR:HG21	1:C:255:ILE:HG23	1.88	0.55
1:A:114:ARG:HB3	1:A:128:LEU:HD23	1.88	0.55
1:D:229:SER:HA	1:D:261:MET:HE1	1.90	0.54
1:E:11:PHE:CZ	1:E:76:GLY:HA3	2.42	0.54
1:D:254:GLY:HA3	1:D:296:LEU:HD13	1.89	0.54
1:E:164:ILE:HG23	1:E:206:PRO:HG3	1.89	0.54
1:C:16:VAL:O	1:C:19:LEU:HB2	2.07	0.53
1:B:254:GLY:HA3	1:B:296:LEU:HD13	1.90	0.53
1:E:72:LEU:HD21	1:E:91:LEU:HD22	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:SER:HA	1:A:261:MET:HE1	1.90	0.53
1:A:256:THR:CG2	1:B:259:LEU:HD11	2.34	0.53
1:D:256:THR:CG2	1:E:259:LEU:HD11	2.37	0.53
1:C:144:PRO:HD3	1:C:281:ILE:HB	1.91	0.52
1:E:254:GLY:HA3	1:E:296:LEU:HD13	1.91	0.52
1:A:144:PRO:HD3	1:A:281:ILE:HB	1.92	0.52
1:D:272:LEU:HD13	1:D:279:LYS:CE	2.41	0.51
1:E:220:TYR:CZ	1:E:224:GLN:HG3	2.45	0.51
1:A:220:TYR:CZ	1:A:224:GLN:HG3	2.45	0.51
1:A:256:THR:HG21	1:B:255:ILE:CG2	2.40	0.51
1:D:220:TYR:CZ	1:D:224:GLN:HG3	2.46	0.51
1:B:144:PRO:HD2	1:B:145:LEU:HD12	1.92	0.51
1:B:229:SER:HA	1:B:261:MET:HE1	1.91	0.51
1:B:256:THR:HG21	1:C:255:ILE:CG2	2.41	0.50
1:B:220:TYR:CZ	1:B:224:GLN:HG3	2.46	0.50
1:A:66:TYR:CZ	1:A:123:THR:HG21	2.46	0.50
1:B:129:ARG:HH22	1:C:101:ASP:HB3	1.76	0.50
1:C:228:PRO:O	1:C:232:ILE:HG12	2.12	0.50
1:A:255:ILE:HG23	1:E:256:THR:HG21	1.92	0.50
1:B:224:GLN:HG2	1:C:269:ARG:HD2	1.94	0.50
1:D:228:PRO:O	1:D:232:ILE:HG12	2.12	0.49
1:A:255:ILE:CG2	1:E:256:THR:HG21	2.42	0.49
1:E:261:MET:HE3	1:E:288:CYS:HB3	1.95	0.49
1:B:66:TYR:CD1	1:B:169:ARG:HD3	2.48	0.49
1:C:220:TYR:CZ	1:C:224:GLN:HG3	2.47	0.49
1:C:98:PHE:HB2	1:C:101:ASP:HB2	1.93	0.49
1:A:274:LYS:HD2	1:E:185:GLN:NE2	2.28	0.49
1:A:228:PRO:O	1:A:232:ILE:HG12	2.13	0.48
1:D:68:ARG:HA	1:D:123:THR:HA	1.94	0.48
1:C:256:THR:CG2	1:D:259:LEU:HD11	2.40	0.48
1:D:189:VAL:HG12	1:D:213:ARG:HB3	1.95	0.48
1:D:145:LEU:HG	1:D:218:ILE:HD13	1.95	0.48
1:A:68:ARG:HD2	1:A:70:LYS:NZ	2.29	0.48
1:B:117:ARG:NH2	1:C:202:THR:HB	2.29	0.48
1:A:68:ARG:HA	1:A:123:THR:HA	1.95	0.47
1:A:101:ASP:HB3	1:E:129:ARG:HH22	1.79	0.47
1:D:98:PHE:HB2	1:D:101:ASP:HB2	1.97	0.47
1:A:98:PHE:HB2	1:A:101:ASP:HB2	1.96	0.47
1:B:189:VAL:HG12	1:B:213:ARG:HB3	1.96	0.47
1:B:268:LEU:HD21	1:B:281:ILE:HG21	1.96	0.46
1:C:129:ARG:HH22	1:D:101:ASP:HB3	1.79	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:188:ILE:HD13	1:A:214:LEU:HD23	1.98	0.46
1:C:42:ILE:HB	1:C:175:VAL:HG22	1.97	0.46
1:C:99:LEU:HB3	1:C:153:GLU:HB2	1.97	0.46
1:B:275:ILE:HD12	1:B:277:TYR:CE1	2.51	0.46
1:E:144:PRO:HD2	1:E:145:LEU:HD12	1.98	0.46
1:D:77:ILE:HG22	1:D:79:LEU:H	1.81	0.46
1:C:18:LYS:HD3	1:C:21:LYS:HD2	1.97	0.46
1:D:185:GLN:NE2	1:E:274:LYS:HD2	2.30	0.46
1:D:42:ILE:HB	1:D:175:VAL:HG22	1.97	0.45
1:C:105:PHE:CZ	1:D:103:LYS:HD2	2.51	0.45
1:E:275:ILE:HD12	1:E:277:TYR:CE1	2.51	0.45
1:C:272:LEU:HB3	1:C:273:PRO:HD2	1.99	0.45
1:D:278:VAL:HG23	1:D:282:ASP:HB3	1.99	0.45
1:E:23:TYR:CE2	1:E:72:LEU:HD11	2.52	0.45
1:E:77:ILE:HG22	1:E:79:LEU:H	1.82	0.45
1:A:103:LYS:HD2	1:E:105:PHE:CZ	2.53	0.44
1:E:42:ILE:HB	1:E:175:VAL:HG22	1.98	0.44
1:C:222:ILE:HA	1:C:226:TYR:HB2	1.98	0.44
1:E:47:ILE:HG12	1:E:59:LEU:HD12	2.00	0.44
1:C:188:ILE:HD13	1:C:214:LEU:HD23	1.98	0.44
1:D:84:ASP:HB3	1:E:26:ARG:HA	1.98	0.44
1:E:222:ILE:HA	1:E:226:TYR:HB2	1.98	0.44
1:D:107:HIS:HB2	1:D:113:ASN:OD1	2.18	0.44
1:A:105:PHE:CZ	1:B:103:LYS:HD2	2.53	0.44
1:A:222:ILE:HA	1:A:226:TYR:HB2	2.00	0.44
1:D:13:LYS:HG2	1:E:27:LEU:HD22	2.00	0.44
1:A:99:LEU:HB3	1:A:153:GLU:HB2	2.00	0.44
1:C:275:ILE:HD12	1:C:277:TYR:CE1	2.52	0.44
1:B:77:ILE:HG22	1:B:79:LEU:H	1.82	0.44
1:B:99:LEU:HB3	1:B:153:GLU:HB2	2.00	0.44
1:D:105:PHE:CZ	1:E:103:LYS:HD2	2.53	0.44
1:D:272:LEU:HD13	1:D:279:LYS:HZ2	1.82	0.44
1:E:68:ARG:HA	1:E:123:THR:HA	1.99	0.44
1:D:222:ILE:HA	1:D:226:TYR:HB2	2.00	0.43
1:A:9:MET:HG2	1:B:31:PHE:HB3	1.99	0.43
1:B:256:THR:CG2	1:C:259:LEU:HD11	2.43	0.43
1:A:68:ARG:HH11	1:A:70:LYS:HE2	1.83	0.43
1:A:275:ILE:HD12	1:A:277:TYR:CE1	2.54	0.43
1:A:114:ARG:HA	1:A:127:GLY:O	2.18	0.43
1:D:129:ARG:NH2	1:E:101:ASP:HB3	2.27	0.43
1:C:12:VAL:HG11	1:D:31:PHE:HE2	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:77:ILE:HG22	1:A:79:LEU:H	1.83	0.42
1:D:287:GLY:HA3	1:D:438:PHE:CZ	2.54	0.42
1:B:42:ILE:HB	1:B:175:VAL:HG22	2.01	0.42
1:D:427:SER:HA	1:D:430:VAL:HG12	2.01	0.42
1:D:99:LEU:HB3	1:D:153:GLU:HB2	2.00	0.42
1:B:222:ILE:HA	1:B:226:TYR:HB2	2.00	0.42
1:C:220:TYR:CZ	1:D:274:LYS:HG2	2.54	0.42
1:D:49:MET:HE2	1:D:51:SER:HB2	2.01	0.42
1:D:160:THR:HG23	1:D:204:ALA:O	2.20	0.42
1:E:69:ASP:HB3	1:E:72:LEU:HD13	2.01	0.42
1:E:194:VAL:HB	1:E:209:SER:HB3	2.02	0.42
1:B:151:THR:HG22	1:B:211:SER:HB3	2.02	0.42
1:C:77:ILE:HG22	1:C:79:LEU:H	1.83	0.42
1:E:69:ASP:CG	1:E:72:LEU:HD13	2.45	0.42
1:B:44:ILE:HG22	1:B:181:ILE:HD11	2.01	0.41
1:B:220:TYR:CZ	1:C:274:LYS:HG2	2.55	0.41
1:C:85:ASN:HB2	1:C:114:ARG:HB3	2.02	0.41
1:A:255:ILE:HD11	1:E:253:LEU:HG	2.03	0.41
1:B:19:LEU:HD22	1:B:72:LEU:HA	2.02	0.41
1:D:284:TYR:HB2	1:D:442:TYR:CE1	2.56	0.41
1:D:38:VAL:HG11	1:D:94:PRO:HG2	2.03	0.41
1:B:301:PHE:HZ	1:B:312:ARG:HH21	1.67	0.41
1:C:227:MET:HE2	1:C:227:MET:HB3	1.96	0.41
1:C:265:ASN:HD21	1:C:286:MET:HE3	1.85	0.41
1:E:147:GLU:HG3	1:E:215:LYS:HD2	2.02	0.41
1:B:115:MET:HE2	1:B:115:MET:HB3	1.89	0.41
1:B:268:LEU:HD21	1:B:281:ILE:CG2	2.52	0.41
1:B:143:TYR:O	1:B:280:ALA:HB3	2.22	0.40
1:D:437:LEU:HD12	1:D:437:LEU:HA	1.95	0.40
1:A:12:VAL:HG11	1:B:31:PHE:CE2	2.56	0.40
1:C:249:ALA:HB1	1:D:251:VAL:HG21	2.03	0.40
1:D:272:LEU:HD13	1:D:279:LYS:NZ	2.37	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	338/355 (95%)	329 (97%)	9 (3%)	0	100	100
1	B	329/355 (93%)	318 (97%)	11 (3%)	0	100	100
1	C	329/355 (93%)	319 (97%)	10 (3%)	0	100	100
1	D	330/355 (93%)	319 (97%)	11 (3%)	0	100	100
1	E	329/355 (93%)	319 (97%)	10 (3%)	0	100	100
All	All	1655/1775 (93%)	1604 (97%)	51 (3%)	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	304/314 (97%)	293 (96%)	11 (4%)	31	63
1	B	296/314 (94%)	287 (97%)	9 (3%)	36	67
1	C	296/314 (94%)	282 (95%)	14 (5%)	23	56
1	D	297/314 (95%)	281 (95%)	16 (5%)	20	51
1	E	296/314 (94%)	285 (96%)	11 (4%)	30	62
All	All	1489/1570 (95%)	1428 (96%)	61 (4%)	27	60

All (61) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	MET
1	A	49	MET
1	A	99	LEU
1	A	161	THR
1	A	185	GLN
1	A	189	VAL

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Mol	Chain	Res	Type
1	A	211	SER
1	A	213	ARG
1	A	261	MET
1	A	281	ILE
1	A	305	ILE
1	B	12	VAL
1	B	40	MET
1	B	49	MET
1	B	59	LEU
1	B	99	LEU
1	B	161	THR
1	B	185	GLN
1	B	189	VAL
1	B	281	ILE
1	C	12	VAL
1	C	40	MET
1	C	41	ASN
1	C	49	MET
1	C	59	LEU
1	C	99	LEU
1	C	112	LYS
1	C	123	THR
1	C	147	GLU
1	C	185	GLN
1	C	189	VAL
1	C	211	SER
1	C	242	ILE
1	C	440	LEU
1	D	12	VAL
1	D	38	VAL
1	D	40	MET
1	D	49	MET
1	D	59	LEU
1	D	99	LEU
1	D	112	LYS
1	D	185	GLN
1	D	189	VAL
1	D	211	SER
1	D	261	MET
1	D	269	ARG
1	D	278	VAL
1	D	279	LYS

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Mol	Chain	Res	Type
1	D	281	ILE
1	D	292	VAL
1	E	12	VAL
1	E	40	MET
1	E	49	MET
1	E	59	LEU
1	E	99	LEU
1	E	110	THR
1	E	164	ILE
1	E	185	GLN
1	E	189	VAL
1	E	227	MET
1	E	263	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	ASN
1	A	191	HIS
1	A	265	ASN
1	B	54	ASN
1	B	65	GLN
1	B	85	ASN
1	B	265	ASN
1	C	41	ASN
1	C	65	GLN
1	C	265	ASN
1	D	65	GLN
1	E	265	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 ⓘ

15 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	F	1	1,2	14,14,15	0.56	0	17,19,21	1.29	2 (11%)
2	NAG	F	2	2	14,14,15	0.57	0	17,19,21	1.47	3 (17%)
2	BMA	F	3	2	11,11,12	0.42	0	15,15,17	0.67	0
2	NAG	G	1	1,2	14,14,15	0.51	0	17,19,21	1.49	3 (17%)
2	NAG	G	2	2	14,14,15	0.64	0	17,19,21	1.35	2 (11%)
2	BMA	G	3	2	11,11,12	0.40	0	15,15,17	0.64	0
2	NAG	H	1	1,2	14,14,15	0.57	0	17,19,21	1.57	3 (17%)
2	NAG	H	2	2	14,14,15	0.65	0	17,19,21	1.50	3 (17%)
2	BMA	H	3	2	11,11,12	0.40	0	15,15,17	0.57	0
2	NAG	I	1	1,2	14,14,15	0.56	0	17,19,21	1.48	4 (23%)
2	NAG	I	2	2	14,14,15	0.61	0	17,19,21	1.47	3 (17%)
2	BMA	I	3	2	11,11,12	0.41	0	15,15,17	0.67	0
2	NAG	J	1	1,2	14,14,15	0.52	0	17,19,21	1.34	3 (17%)
2	NAG	J	2	2	14,14,15	0.61	0	17,19,21	1.49	2 (11%)
2	BMA	J	3	2	11,11,12	0.43	0	15,15,17	0.77	1 (6%)

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	F	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	F	2	2	-	2/6/23/26	0/1/1/1
2	BMA	F	3	2	-	0/2/19/22	0/1/1/1
2	NAG	G	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	G	2	2	-	2/6/23/26	0/1/1/1
2	BMA	G	3	2	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	H	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	H	2	2	-	2/6/23/26	0/1/1/1
2	BMA	H	3	2	-	0/2/19/22	0/1/1/1
2	NAG	I	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	I	2	2	-	2/6/23/26	0/1/1/1
2	BMA	I	3	2	-	0/2/19/22	0/1/1/1
2	NAG	J	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	J	2	2	-	2/6/23/26	0/1/1/1
2	BMA	J	3	2	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	2	NAG	C1-O5-C5	4.34	118.00	112.19
2	I	2	NAG	C1-O5-C5	4.05	117.61	112.19
2	F	2	NAG	C1-O5-C5	4.04	117.61	112.19
2	H	2	NAG	C1-O5-C5	3.77	117.23	112.19
2	G	2	NAG	C1-O5-C5	3.68	117.12	112.19
2	H	1	NAG	C1-O5-C5	3.58	116.99	112.19
2	G	1	NAG	C1-O5-C5	3.38	116.72	112.19
2	I	1	NAG	C1-O5-C5	3.22	116.50	112.19
2	F	1	NAG	C1-O5-C5	3.20	116.47	112.19
2	J	1	NAG	C1-O5-C5	3.14	116.39	112.19
2	H	1	NAG	C4-C3-C2	3.13	115.61	111.02
2	I	1	NAG	C4-C3-C2	3.01	115.42	111.02
2	G	1	NAG	C2-N2-C7	2.94	126.84	122.90
2	G	1	NAG	C4-C3-C2	2.81	115.14	111.02
2	H	1	NAG	C2-N2-C7	2.67	126.48	122.90
2	H	2	NAG	C4-C3-C2	2.60	114.83	111.02
2	F	2	NAG	C4-C3-C2	2.46	114.62	111.02
2	J	3	BMA	C1-O5-C5	2.40	115.40	112.19
2	F	1	NAG	C4-C3-C2	2.40	114.53	111.02
2	J	1	NAG	C4-C3-C2	2.39	114.52	111.02
2	I	2	NAG	C4-C3-C2	2.35	114.47	111.02
2	I	1	NAG	C2-N2-C7	2.35	126.05	122.90
2	F	2	NAG	O5-C1-C2	-2.35	107.65	111.29
2	G	2	NAG	C4-C3-C2	2.24	114.30	111.02
2	J	2	NAG	C4-C3-C2	2.19	114.22	111.02
2	J	1	NAG	C2-N2-C7	2.11	125.72	122.90
2	H	2	NAG	O5-C1-C2	-2.10	108.05	111.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	2	NAG	O5-C1-C2	-2.07	108.08	111.29
2	I	1	NAG	O5-C5-C6	2.02	111.59	107.66

There are no chirality outliers.

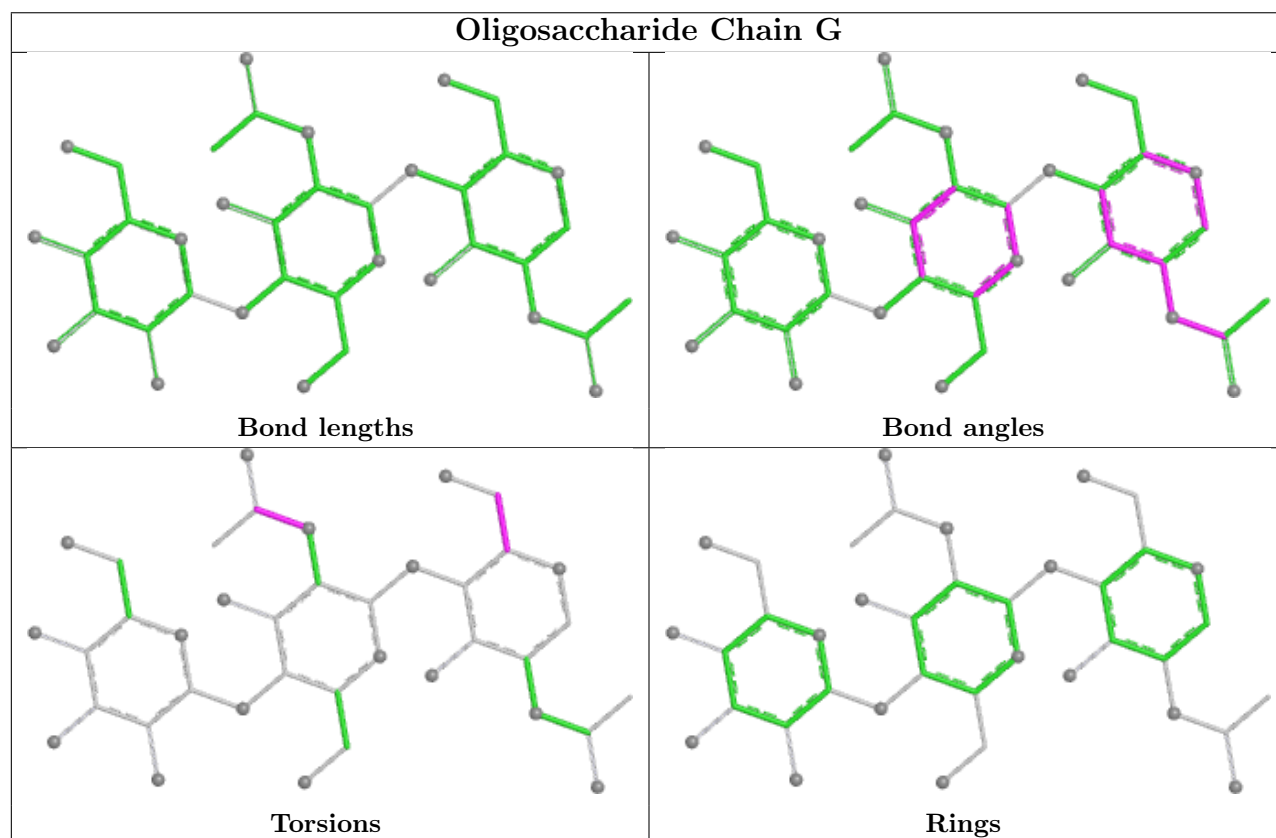
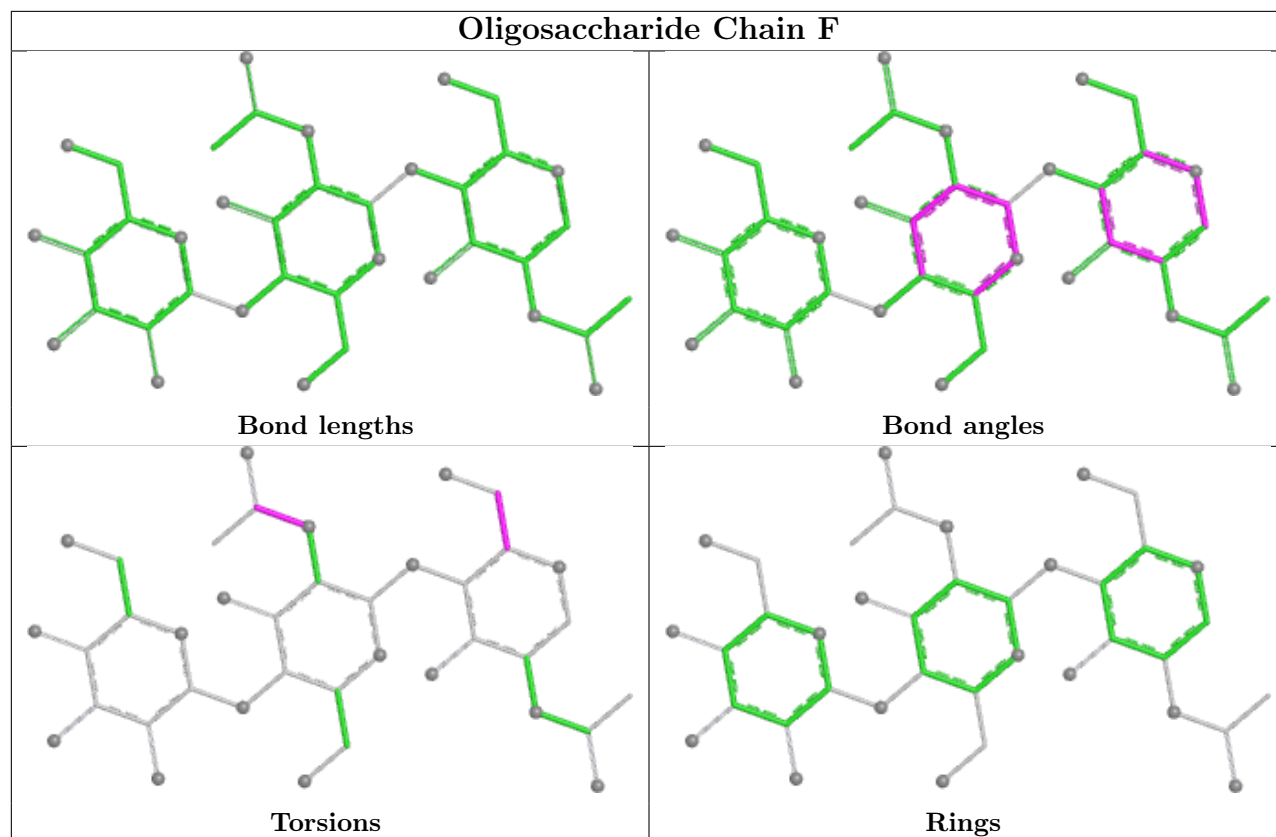
All (15) torsion outliers are listed below:

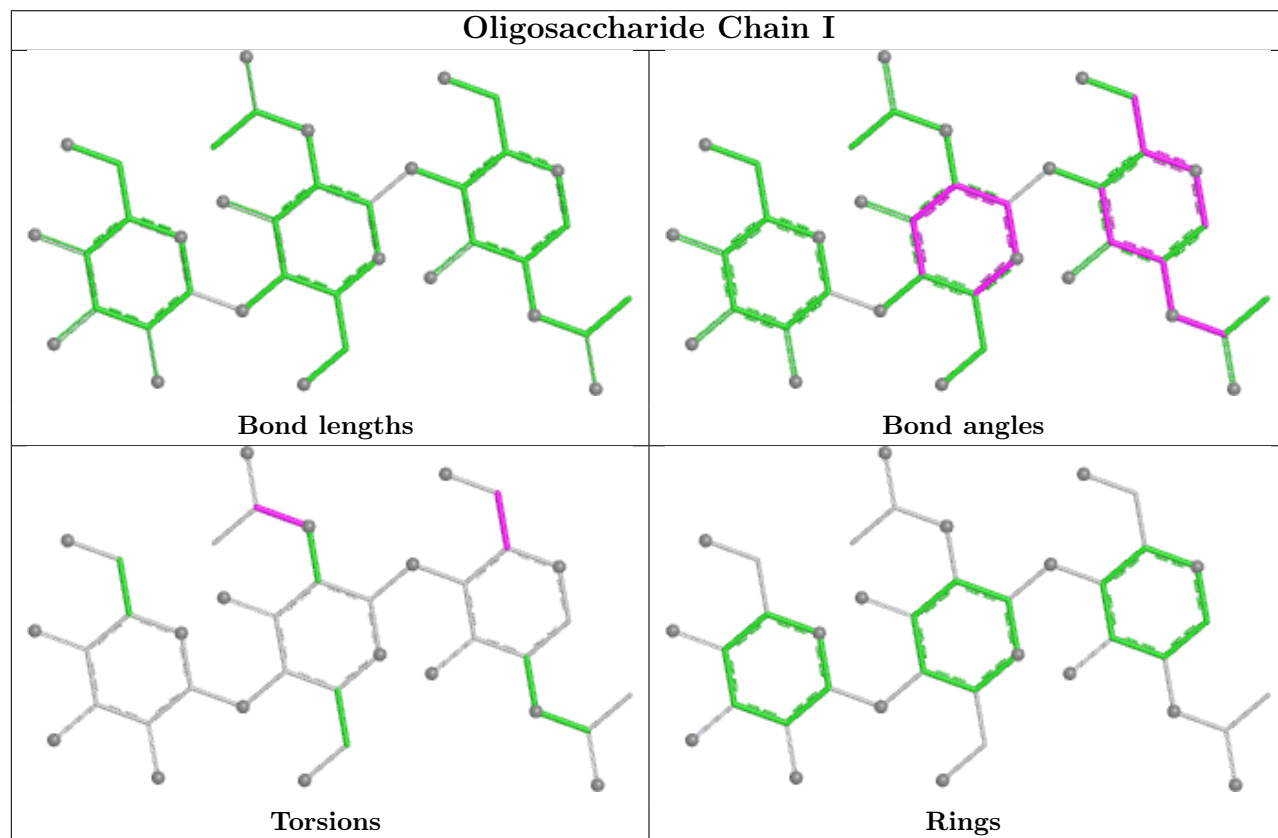
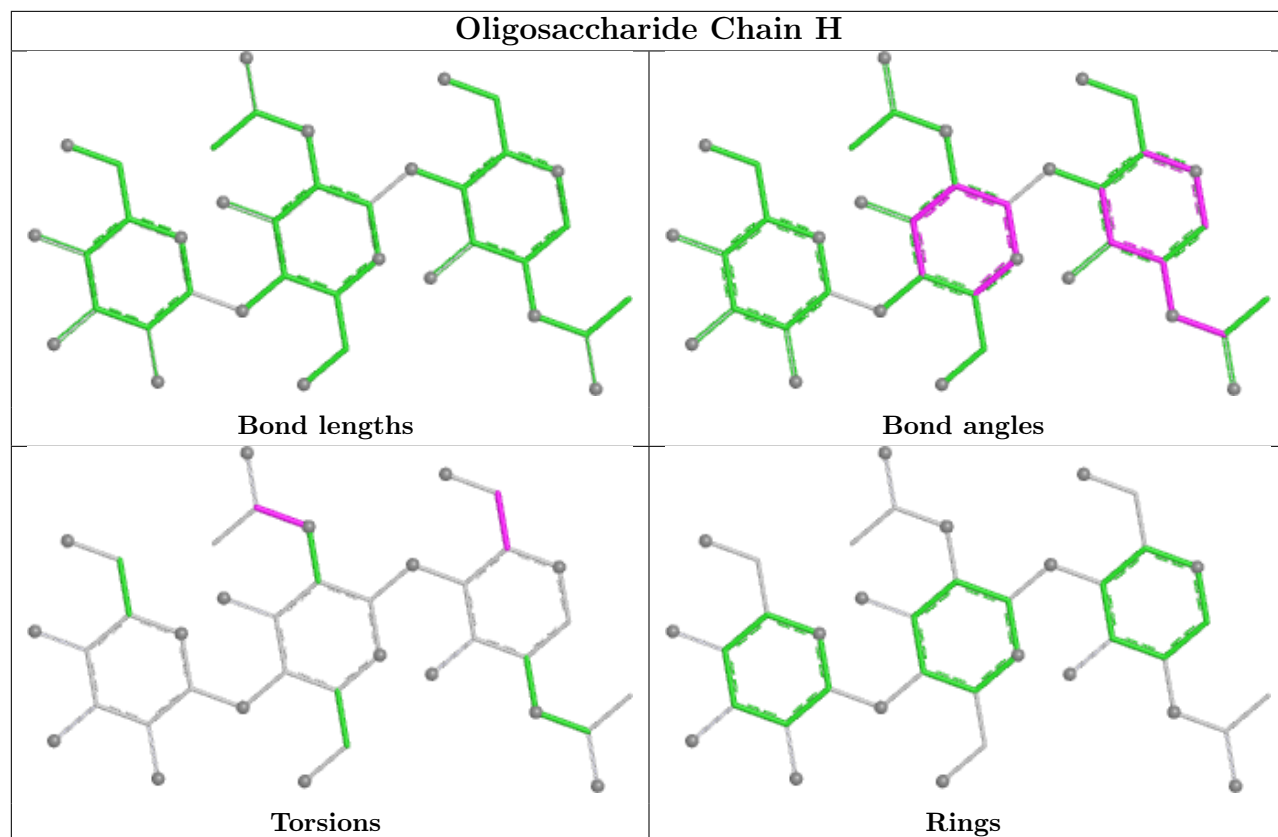
Mol	Chain	Res	Type	Atoms
2	F	2	NAG	C8-C7-N2-C2
2	G	2	NAG	C8-C7-N2-C2
2	H	2	NAG	C8-C7-N2-C2
2	H	2	NAG	O7-C7-N2-C2
2	I	2	NAG	C8-C7-N2-C2
2	I	2	NAG	O7-C7-N2-C2
2	J	2	NAG	C8-C7-N2-C2
2	F	2	NAG	O7-C7-N2-C2
2	G	2	NAG	O7-C7-N2-C2
2	J	2	NAG	O7-C7-N2-C2
2	I	1	NAG	C4-C5-C6-O6
2	H	1	NAG	C4-C5-C6-O6
2	F	1	NAG	C4-C5-C6-O6
2	G	1	NAG	C4-C5-C6-O6
2	J	1	NAG	C4-C5-C6-O6

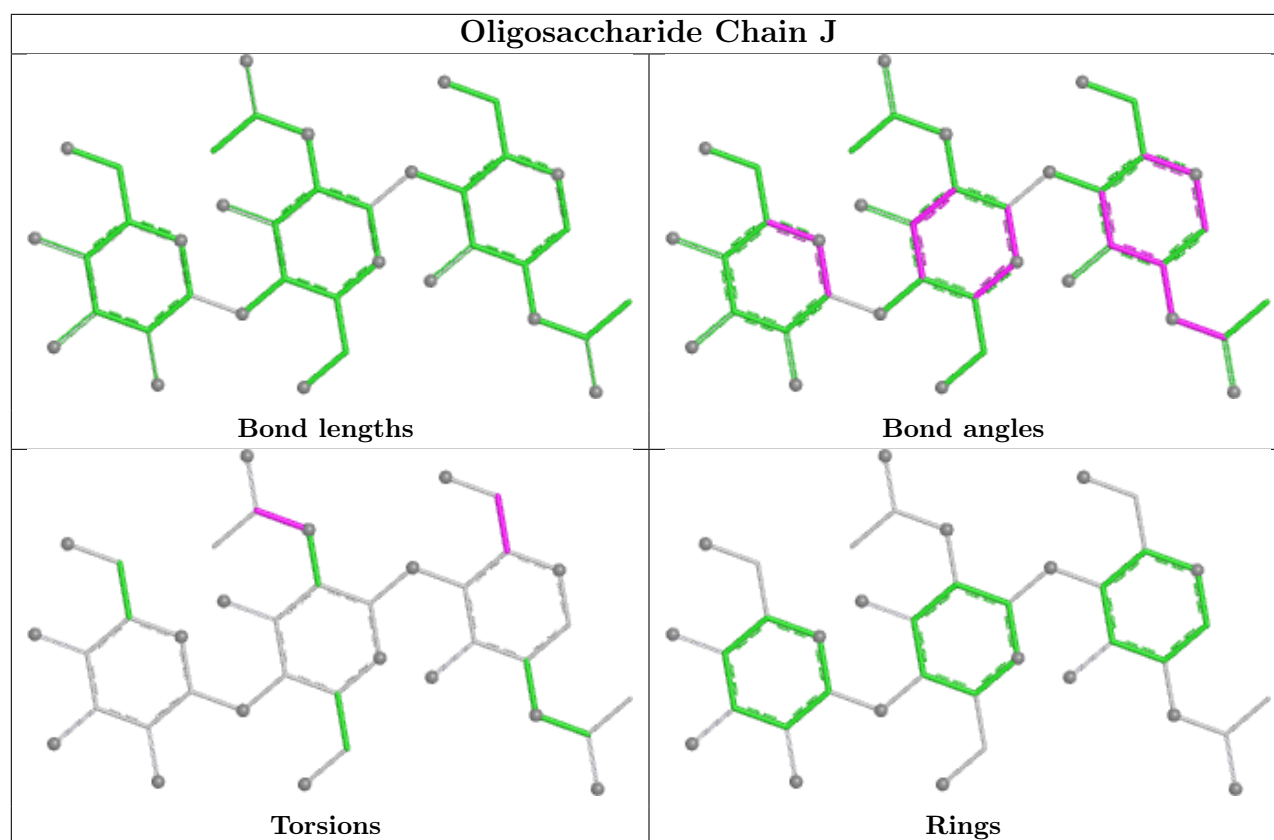
There are no ring outliers.

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](#)

Of 16 ligands modelled in this entry, 5 are monoatomic - leaving 11 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	$\# Z > 2$
3	BEN	C	500	-	9,9,9	2.32	4 (44%)	7,11,11	1.23	0
4	NAG	B	2000	1	14,14,15	0.54	0	17,19,21	1.58	3 (17%)
4	NAG	C	2000	1	14,14,15	0.59	0	17,19,21	1.40	2 (11%)
3	BEN	D	500	-	9,9,9	2.20	4 (44%)	7,11,11	1.16	0
3	BEN	B	500	-	9,9,9	2.56	5 (55%)	7,11,11	1.19	0
4	NAG	E	2000	1	14,14,15	0.63	0	17,19,21	1.48	3 (17%)
3	BEN	E	500	-	9,9,9	2.16	3 (33%)	7,11,11	1.31	1 (14%)
4	NAG	A	1000	1	14,14,15	0.60	0	17,19,21	1.68	4 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	A	2000	1	14,14,15	0.60	0	17,19,21	1.71	4 (23%)
4	NAG	D	2000	1	14,14,15	0.56	0	17,19,21	1.50	2 (11%)
3	BEN	A	500	-	9,9,9	2.59	5 (55%)	7,11,11	1.50	1 (14%)

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	BEN	C	500	-	-	4/4/4/4	0/1/1/1
4	NAG	B	2000	1	-	3/6/23/26	0/1/1/1
4	NAG	C	2000	1	-	4/6/23/26	0/1/1/1
3	BEN	D	500	-	-	4/4/4/4	0/1/1/1
3	BEN	B	500	-	-	4/4/4/4	0/1/1/1
4	NAG	E	2000	1	-	4/6/23/26	0/1/1/1
3	BEN	E	500	-	-	4/4/4/4	0/1/1/1
4	NAG	A	1000	1	-	5/6/23/26	0/1/1/1
4	NAG	A	2000	1	-	4/6/23/26	0/1/1/1
4	NAG	D	2000	1	-	4/6/23/26	0/1/1/1
3	BEN	A	500	-	-	4/4/4/4	0/1/1/1

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	500	BEN	C6-C1	4.50	1.46	1.39
3	B	500	BEN	C2-C1	4.43	1.46	1.39
3	D	500	BEN	C2-C1	4.03	1.45	1.39
3	C	500	BEN	C2-C1	3.98	1.45	1.39
3	A	500	BEN	C2-C1	3.81	1.45	1.39
3	B	500	BEN	C6-C1	3.80	1.45	1.39
3	E	500	BEN	C6-C1	3.46	1.44	1.39
3	E	500	BEN	C2-C1	3.43	1.44	1.39
3	D	500	BEN	C6-C1	2.93	1.43	1.39
3	B	500	BEN	C-N2	2.76	1.41	1.33
3	C	500	BEN	C6-C1	2.72	1.43	1.39
3	C	500	BEN	C-N2	2.54	1.40	1.33
3	C	500	BEN	C5-C6	2.44	1.43	1.38
3	D	500	BEN	C-N2	2.38	1.40	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	500	BEN	C5-C6	2.34	1.42	1.38
3	B	500	BEN	C5-C4	2.25	1.43	1.38
3	A	500	BEN	C1-C	2.20	1.51	1.47
3	D	500	BEN	C5-C6	2.13	1.42	1.38
3	A	500	BEN	C4-C3	2.06	1.42	1.38
3	E	500	BEN	C-N2	2.06	1.39	1.33
3	B	500	BEN	C5-C6	2.02	1.42	1.38

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	2000	NAG	C1-O5-C5	4.31	117.97	112.19
4	B	2000	NAG	C1-O5-C5	4.15	117.75	112.19
4	D	2000	NAG	C1-O5-C5	3.89	117.40	112.19
3	A	500	BEN	C1-C-N2	3.89	123.91	118.01
4	A	1000	NAG	C2-N2-C7	3.82	128.02	122.90
4	C	2000	NAG	C1-O5-C5	3.56	116.96	112.19
4	A	1000	NAG	C1-O5-C5	3.53	116.92	112.19
4	E	2000	NAG	C4-C3-C2	3.29	115.84	111.02
4	A	1000	NAG	C3-C4-C5	3.07	115.80	110.23
4	B	2000	NAG	C4-C3-C2	3.07	115.52	111.02
4	A	2000	NAG	C4-C3-C2	3.02	115.45	111.02
4	E	2000	NAG	C1-O5-C5	2.76	115.88	112.19
4	D	2000	NAG	C4-C3-C2	2.76	115.06	111.02
4	C	2000	NAG	C4-C3-C2	2.66	114.91	111.02
3	E	500	BEN	C1-C-N2	2.56	121.90	118.01
4	A	1000	NAG	C4-C3-C2	2.51	114.70	111.02
4	A	2000	NAG	C3-C4-C5	2.36	114.50	110.23
4	A	2000	NAG	O7-C7-N2	2.25	125.96	121.98
4	E	2000	NAG	C2-N2-C7	2.24	125.90	122.90
4	B	2000	NAG	C3-C4-C5	2.13	114.10	110.23

There are no chirality outliers.

All (44) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1000	NAG	C3-C2-N2-C7
4	A	1000	NAG	C8-C7-N2-C2
4	A	1000	NAG	O7-C7-N2-C2
4	A	1000	NAG	C4-C5-C6-O6
4	E	2000	NAG	C4-C5-C6-O6
4	A	2000	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
4	D	2000	NAG	C8-C7-N2-C2
4	B	2000	NAG	C8-C7-N2-C2
4	C	2000	NAG	C8-C7-N2-C2
4	D	2000	NAG	O7-C7-N2-C2
4	E	2000	NAG	C8-C7-N2-C2
4	A	2000	NAG	C8-C7-N2-C2
4	B	2000	NAG	O7-C7-N2-C2
4	C	2000	NAG	C4-C5-C6-O6
4	C	2000	NAG	O7-C7-N2-C2
4	E	2000	NAG	O7-C7-N2-C2
4	D	2000	NAG	C4-C5-C6-O6
4	A	1000	NAG	O5-C5-C6-O6
4	A	2000	NAG	O7-C7-N2-C2
4	B	2000	NAG	C4-C5-C6-O6
3	A	500	BEN	N2-C-C1-C2
3	A	500	BEN	N2-C-C1-C6
3	B	500	BEN	N2-C-C1-C2
3	B	500	BEN	N2-C-C1-C6
3	C	500	BEN	N2-C-C1-C2
3	C	500	BEN	N2-C-C1-C6
3	D	500	BEN	N2-C-C1-C2
3	D	500	BEN	N2-C-C1-C6
3	E	500	BEN	N2-C-C1-C2
3	E	500	BEN	N2-C-C1-C6
4	C	2000	NAG	O5-C5-C6-O6
4	A	2000	NAG	O5-C5-C6-O6
4	E	2000	NAG	O5-C5-C6-O6
3	A	500	BEN	N1-C-C1-C2
3	A	500	BEN	N1-C-C1-C6
3	B	500	BEN	N1-C-C1-C2
3	B	500	BEN	N1-C-C1-C6
3	C	500	BEN	N1-C-C1-C2
3	C	500	BEN	N1-C-C1-C6
3	D	500	BEN	N1-C-C1-C2
3	D	500	BEN	N1-C-C1-C6
3	E	500	BEN	N1-C-C1-C2
3	E	500	BEN	N1-C-C1-C6
4	D	2000	NAG	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	340/355 (95%)	0.62	38 (11%) 10 7	68, 99, 127, 158	0
1	B	331/355 (93%)	0.39	23 (6%) 23 14	69, 93, 135, 158	0
1	C	331/355 (93%)	0.58	34 (10%) 12 8	67, 94, 130, 168	0
1	D	332/355 (93%)	0.73	39 (11%) 9 6	70, 102, 139, 191	0
1	E	331/355 (93%)	0.68	33 (9%) 12 8	66, 103, 138, 161	0
All	All	1665/1775 (93%)	0.60	167 (10%) 12 8	66, 99, 136, 191	0

All (167) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	136	CYS	6.3
1	D	150	CYS	5.5
1	C	182	GLU	5.3
1	D	10	SER	5.0
1	D	137	MET	5.0
1	D	151	THR	4.8
1	E	447	VAL	4.8
1	D	34	PRO	4.7
1	C	127	GLY	4.6
1	A	78	PRO	4.4
1	D	88	ALA	4.4
1	B	184	PRO	4.3
1	B	142	ARG	4.3
1	A	447	VAL	4.2
1	E	18	LYS	4.2
1	E	307	PHE	4.1
1	E	11	PHE	4.1
1	A	89	ASP	4.0
1	E	34	PRO	3.9
1	C	447	VAL	3.8

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Mol	Chain	Res	Type	RSRZ
1	A	127	GLY	3.7
1	A	227	MET	3.7
1	D	64	GLN	3.7
1	A	147	GLU	3.6
1	D	89	ASP	3.6
1	C	146	ASP	3.5
1	C	184	PRO	3.5
1	D	139	ASP	3.5
1	D	160	THR	3.5
1	E	150	CYS	3.5
1	B	446	TYR	3.4
1	C	63	PHE	3.4
1	C	43	ASP	3.4
1	E	35	PRO	3.3
1	B	43	ASP	3.3
1	E	77	ILE	3.3
1	E	206	PRO	3.2
1	E	13	LYS	3.2
1	B	68	ARG	3.1
1	E	102	LYS	3.1
1	D	447	VAL	3.0
1	A	25	ILE	3.0
1	B	447	VAL	3.0
1	C	147	GLU	3.0
1	A	70	LYS	3.0
1	E	200	PHE	2.9
1	E	25	ILE	2.9
1	B	64	GLN	2.9
1	C	64	GLN	2.9
1	D	202	THR	2.9
1	A	312	ARG	2.9
1	C	142	ARG	2.9
1	C	199	VAL	2.8
1	D	12	VAL	2.8
1	E	101	ASP	2.8
1	A	188	ILE	2.8
1	A	189	VAL	2.8
1	C	126	TYR	2.8
1	A	99	LEU	2.8
1	B	279	LYS	2.8
1	D	57	TYR	2.7
1	C	446	TYR	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	43	ASP	2.7
1	A	64	GLN	2.7
1	B	199	VAL	2.7
1	D	218	ILE	2.7
1	A	101	ASP	2.7
1	A	150	CYS	2.7
1	D	23	TYR	2.7
1	A	263	THR	2.7
1	D	149	ASN	2.7
1	A	79	LEU	2.6
1	D	209	SER	2.6
1	D	127	GLY	2.6
1	A	212	PHE	2.6
1	A	301	PHE	2.6
1	C	202	THR	2.6
1	D	212	PHE	2.6
1	C	10	SER	2.5
1	C	145	LEU	2.5
1	C	227	MET	2.5
1	B	202	THR	2.5
1	D	165	GLU	2.5
1	E	103	LYS	2.5
1	A	429	ILE	2.5
1	C	77	ILE	2.5
1	B	150	CYS	2.5
1	C	41	ASN	2.5
1	B	227	MET	2.5
1	C	81	LEU	2.5
1	A	161	THR	2.5
1	A	186	PHE	2.5
1	E	184	PRO	2.4
1	A	223	LEU	2.4
1	B	147	GLU	2.4
1	D	18	LYS	2.4
1	D	227	MET	2.4
1	C	307	PHE	2.4
1	C	115	MET	2.4
1	C	12	VAL	2.4
1	B	21	LYS	2.4
1	E	14	GLU	2.4
1	E	45	ALA	2.4
1	E	79	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	13	LYS	2.3
1	E	199	VAL	2.3
1	E	64	GLN	2.3
1	A	76	GLY	2.3
1	D	24	ASP	2.3
1	E	12	VAL	2.3
1	B	195	SER	2.3
1	E	23	TYR	2.3
1	B	200	PHE	2.2
1	A	137	MET	2.2
1	D	49	MET	2.2
1	C	165	GLU	2.2
1	D	41	ASN	2.2
1	B	173	LYS	2.2
1	E	245	ASP	2.2
1	E	202	THR	2.2
1	B	10	SER	2.2
1	D	211	SER	2.2
1	E	227	MET	2.2
1	A	88	ALA	2.2
1	A	41	ASN	2.2
1	B	145	LEU	2.2
1	D	13	LYS	2.2
1	D	85	ASN	2.2
1	C	104	SER	2.2
1	A	214	LEU	2.2
1	C	179	GLU	2.2
1	C	42	ILE	2.2
1	D	71	ARG	2.2
1	E	136	CYS	2.2
1	A	57	TYR	2.2
1	A	215	LYS	2.2
1	B	79	LEU	2.2
1	B	121	ASP	2.2
1	D	163	ASP	2.2
1	A	23	TYR	2.2
1	A	136	CYS	2.2
1	C	185	GLN	2.1
1	E	137	MET	2.1
1	A	225	THR	2.1
1	B	13	LYS	2.1
1	E	24	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	33	GLY	2.1
1	E	68	ARG	2.1
1	D	25	ILE	2.1
1	C	48	ASP	2.1
1	E	26	ARG	2.1
1	C	40	MET	2.1
1	C	169	ARG	2.1
1	C	136	CYS	2.0
1	A	199	VAL	2.0
1	D	199	VAL	2.0
1	E	16	VAL	2.0
1	A	35	PRO	2.0
1	D	35	PRO	2.0
1	C	128	LEU	2.0
1	D	182	GLU	2.0
1	B	143	TYR	2.0
1	A	278	VAL	2.0
1	D	86	ARG	2.0
1	E	141	ARG	2.0
1	D	77	ILE	2.0
1	A	204	ALA	2.0

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

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

6.3 Carbohydrates [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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.

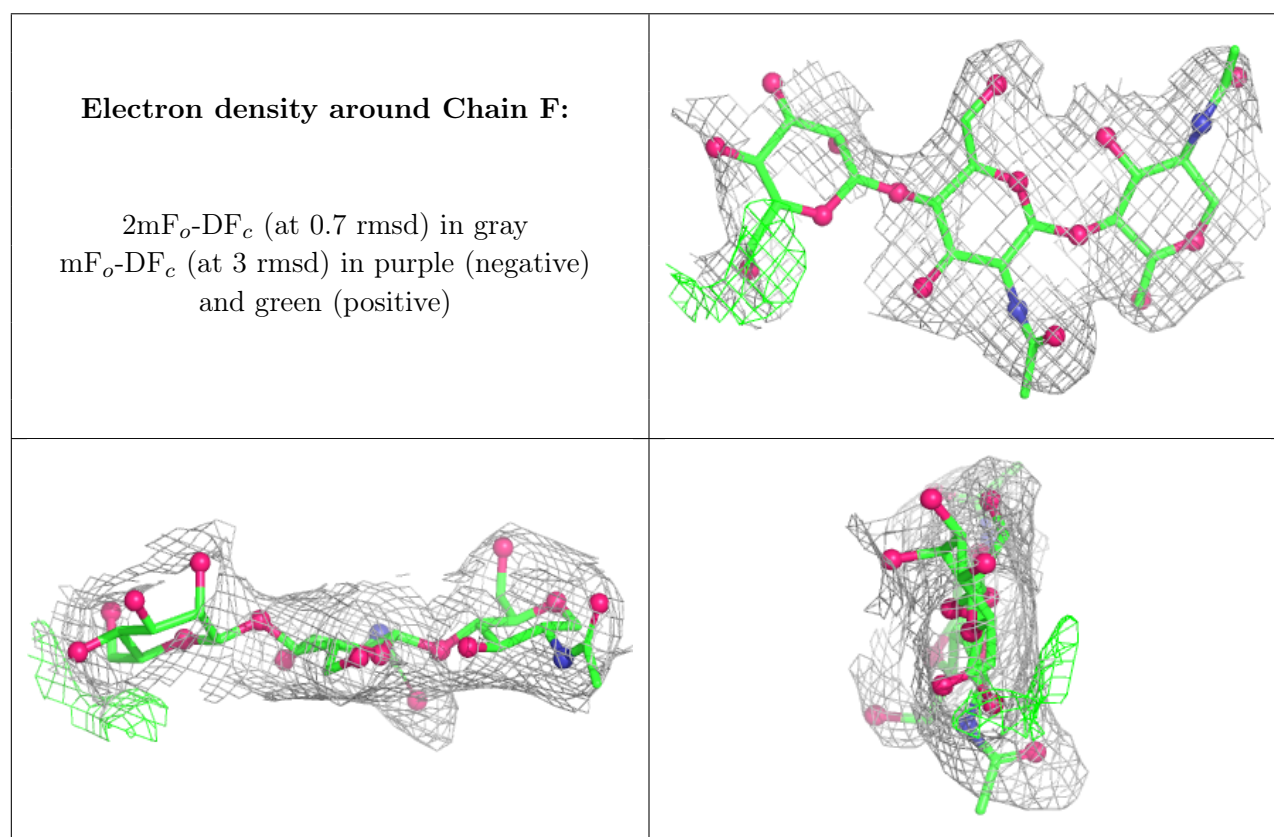
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	BMA	F	3	11/12	0.14	0.17	155,158,160,160	0
2	BMA	J	3	11/12	0.28	0.18	157,160,161,162	0
2	BMA	I	3	11/12	0.34	0.15	159,163,163,164	0
2	BMA	H	3	11/12	0.36	0.14	167,169,171,172	0
2	BMA	G	3	11/12	0.51	0.14	155,157,160,161	0
2	NAG	H	2	14/15	0.73	0.16	134,147,153,161	0
2	NAG	J	2	14/15	0.86	0.13	120,132,140,150	0

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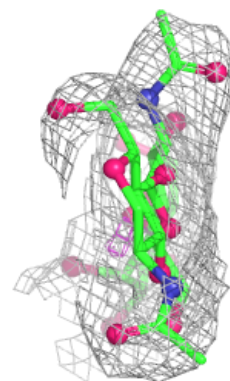
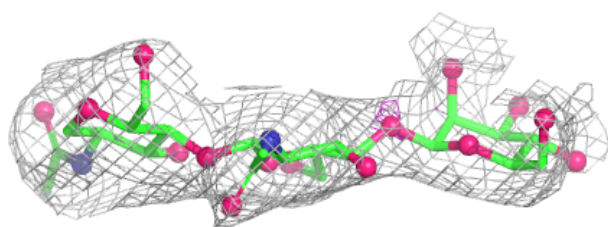
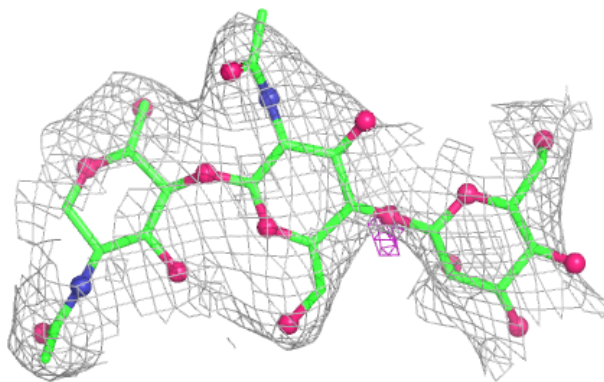
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	I	2	14/15	0.88	0.12	124,132,141,151	0
2	NAG	G	2	14/15	0.89	0.12	115,131,140,149	0
2	NAG	I	1	14/15	0.90	0.12	101,111,117,120	0
2	NAG	J	1	14/15	0.90	0.15	108,114,119,124	0
2	NAG	F	2	14/15	0.91	0.10	120,128,137,148	0
2	NAG	H	1	14/15	0.92	0.09	123,125,130,138	0
2	NAG	G	1	14/15	0.93	0.09	107,112,118,124	0
2	NAG	F	1	14/15	0.94	0.11	113,115,118,121	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.

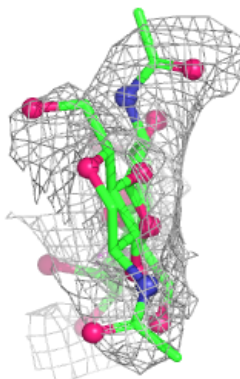
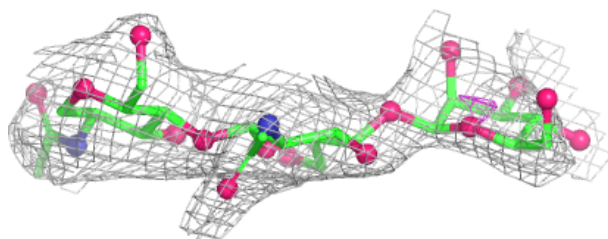
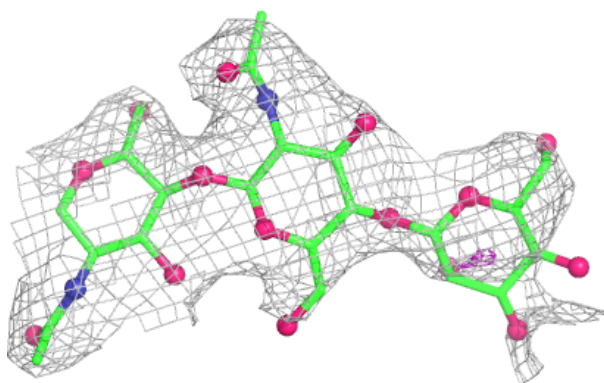


Electron density around Chain G:

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

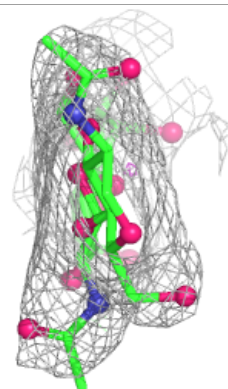
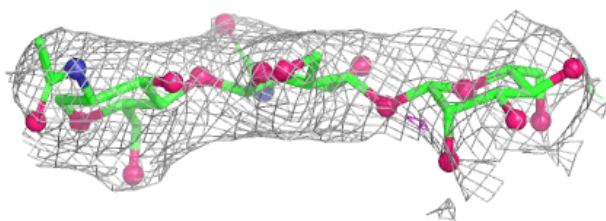
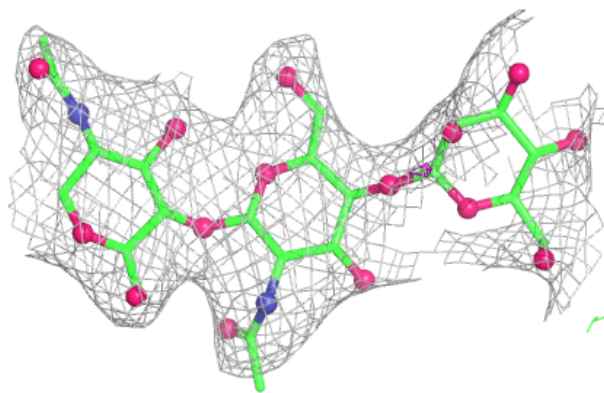
**Electron density around Chain H:**

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

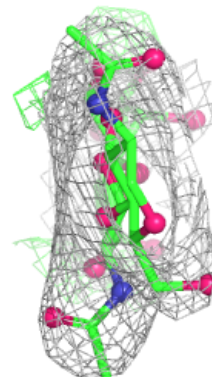
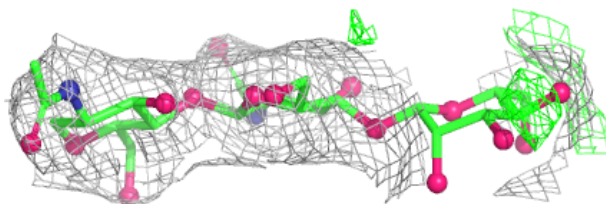
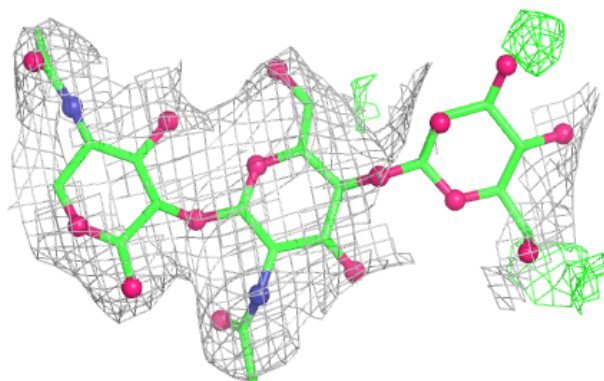


Electron density around Chain I:

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

**Electron density around Chain J:**

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



6.4 Ligands

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
4	NAG	B	2000	14/15	0.61	0.17	152,156,157,157	0
4	NAG	A	1000	14/15	0.67	0.17	144,147,148,149	0
4	NAG	D	2000	14/15	0.70	0.17	158,162,166,166	0
4	NAG	E	2000	14/15	0.74	0.13	136,139,141,142	0
3	BEN	B	500	9/9	0.79	0.35	125,126,126,126	0
3	BEN	A	500	9/9	0.80	0.32	135,136,136,136	0
3	BEN	C	500	9/9	0.80	0.37	129,129,129,129	0
3	BEN	E	500	9/9	0.83	0.35	138,139,139,139	0
4	NAG	C	2000	14/15	0.86	0.13	147,151,154,156	0
5	CL	B	1448	1/1	0.87	0.18	100,100,100,100	0
3	BEN	D	500	9/9	0.88	0.26	117,117,118,118	0
4	NAG	A	2000	14/15	0.89	0.13	114,117,124,124	0
5	CL	C	1448	1/1	0.90	0.15	109,109,109,109	0
5	CL	D	1449	1/1	0.92	0.19	111,111,111,111	0
5	CL	E	1448	1/1	0.93	0.12	100,100,100,100	0
5	CL	A	1448	1/1	0.95	0.14	97,97,97,97	0

6.5 Other polymers

There are no such residues in this entry.