



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:42 PM UTC

PDB ID : 6M4E / pdb_00006m4e
Title : Crystal structure of a GH1 beta-glucosidase from *Hamamotoa singularis*
Authors : Uehara, R.; Iwamoto, R.; Aoki, S.; Yoshizawa, T.; Takano, K.; Matsumura, H.; Tanaka, S.-i.
Deposited on : 2020-03-06
Resolution : 2.10 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)
Xtrriage (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

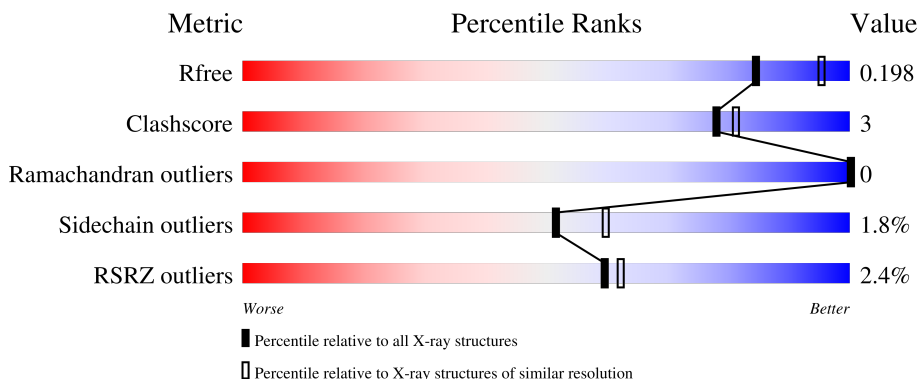
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION




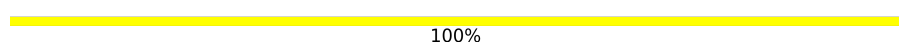
The reported resolution of this entry is 2.10 Å.

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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	578	 2% 88% 6% 6%
2	B	2	 50% 50%
2	C	2	 100%
2	D	2	 100%

2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 4802 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 Beta-galactosidase-like enzyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	542	4288	2755	712	812	9	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

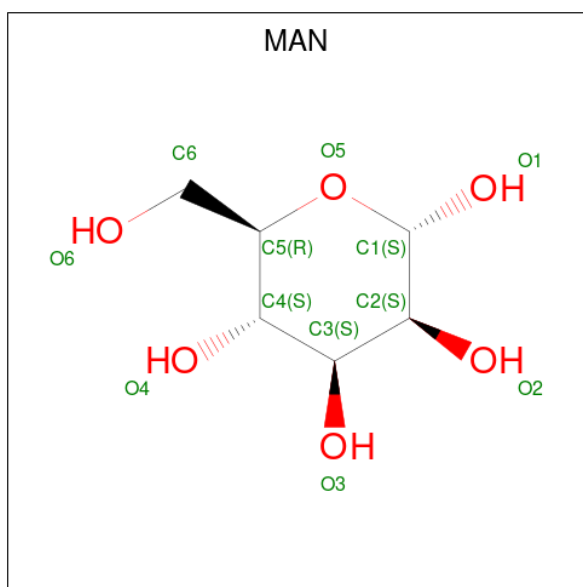
Chain	Residue	Modelled	Actual	Comment	Reference
A	595	HIS	-	expression tag	UNP Q564N5
A	596	HIS	-	expression tag	UNP Q564N5
A	597	HIS	-	expression tag	UNP Q564N5
A	598	HIS	-	expression tag	UNP Q564N5
A	599	HIS	-	expression tag	UNP Q564N5
A	600	HIS	-	expression tag	UNP Q564N5

- 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
2	C	2	28	16	2	10	0	0	0
2	D	2	28	16	2	10	0	0	0

- Molecule 3 is alpha-D-mannopyranose (CCD ID: MAN) (formula: C₆H₁₂O₆).



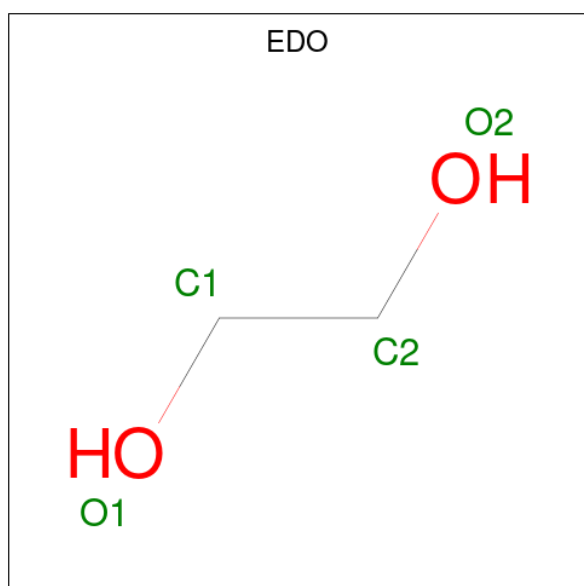
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 11 6 5	0	0
3	A	1	Total C O 11 6 5	0	0
3	A	1	Total C O 11 6 5	0	0

- 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
			Total	C	N	O		
4	A	1	14	8	1	5	0	0

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
5	A	1	4	2	2	0	0
5	A	1	4	2	2	0	0
5	A	1	4	2	2	0	0
5	A	1	4	2	2	0	0
5	A	1	4	2	2	0	0
5	A	1	4	2	2	0	0
5	A	1	4	2	2	0	0
5	A	1	4	2	2	0	0
5	A	1	4	2	2	0	0
5	A	1	4	2	2	0	0
5	A	1	4	2	2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		

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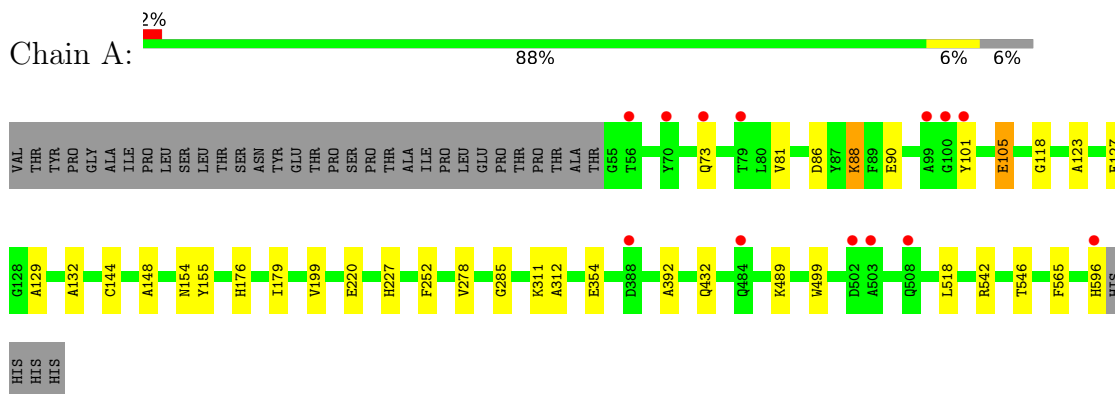
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	171	Total 171	O 171	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.

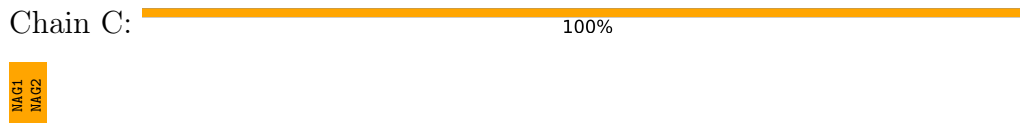
- Molecule 1: Beta-galactosidase-like enzyme



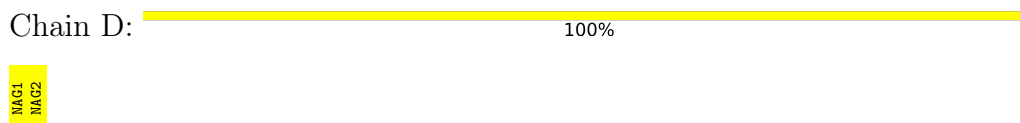
- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



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



- 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 2	Depositor
Cell constants a, b, c, α , β , γ	160.79Å 172.83Å 67.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.24 – 2.10 39.24 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (39.24-2.10) 99.9 (39.24-2.10)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.32 (at 2.10Å)	Xtrriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.174 , 0.197 0.177 , 0.198	Depositor DCC
R_{free} test set	2752 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	47.4	Xtrriage
Anisotropy	0.242	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 49.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4802	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.96% 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: EDO, MAN, 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	0.58	1/4422 (0.0%)	0.56	1/6034 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	546	THR	C-O	-5.01	1.17	1.23

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	499	TRP	N-CA-C	5.18	117.89	109.24

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4288	0	4024	24	0
2	B	28	0	25	1	0
2	C	28	0	25	2	0
2	D	28	0	25	0	0
3	A	33	0	30	1	0
4	A	14	0	13	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	212	0	318	12	0
6	A	171	0	0	1	0
All	All	4802	0	4460	25	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 3.

All (25) 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:86:ASP:O	1:A:88:LYS:HE3	2.01	0.61
1:A:285:GLY:HA2	2:B:1:NAG:O5	2.02	0.58
1:A:105:GLU:O	1:A:105:GLU:HG3	2.01	0.58
1:A:90:GLU:HG3	5:A:734:EDO:O2	2.06	0.55
1:A:542:ARG:HH12	5:A:707:EDO:H22	1.70	0.55
1:A:432:GLN:HE22	2:C:1:NAG:H3	1.71	0.54
1:A:542:ARG:NH1	5:A:707:EDO:H22	2.25	0.52
1:A:118:GLY:HA2	1:A:179:ILE:HG23	1.90	0.52
1:A:81:VAL:HA	3:A:703:MAN:H2	1.92	0.50
1:A:154:ASN:HB2	5:A:724:EDO:H21	1.93	0.50
1:A:596:HIS:H	5:A:716:EDO:H11	1.76	0.50
1:A:432:GLN:OE1	2:C:2:NAG:H83	2.13	0.49
1:A:220:GLU:OE1	5:A:728:EDO:O1	2.30	0.48
1:A:144:CYS:O	1:A:148:ALA:HB2	2.17	0.45
5:A:710:EDO:H22	5:A:726:EDO:O2	2.18	0.44
1:A:392:ALA:HA	5:A:711:EDO:H11	2.01	0.43
1:A:252:PHE:CD2	1:A:312:ALA:HB1	2.54	0.43
1:A:101:TYR:CE1	5:A:733:EDO:H12	2.55	0.42
1:A:155:TYR:CE2	5:A:724:EDO:H22	2.53	0.42
1:A:227:HIS:HD2	6:A:853:HOH:O	2.02	0.42
1:A:518:LEU:HD23	1:A:518:LEU:HA	1.89	0.41
1:A:278:VAL:HG21	5:A:709:EDO:H22	2.01	0.41
1:A:123:ALA:O	1:A:127:GLU:HB2	2.21	0.41
1:A:129:ALA:HB1	1:A:132:ALA:HB3	2.03	0.41
1:A:176:HIS:HE1	5:A:740:EDO:H11	1.86	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	540/578 (93%)	522 (97%)	18 (3%)	0	100 100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	443/475 (93%)	435 (98%)	8 (2%)	51 60

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

Mol	Chain	Res	Type
1	A	73	GLN
1	A	88	LYS
1	A	105	GLU
1	A	199	VAL
1	A	311	LYS
1	A	354	GLU
1	A	489	LYS
1	A	565	PHE

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	73	GLN
1	A	154	ASN
1	A	175	GLN
1	A	176	HIS
1	A	200	ASN
1	A	242	GLN
1	A	283	ASN
1	A	370	GLN
1	A	432	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](#)

6 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	1.35	2 (14%)	17,19,21	1.74	3 (17%)
2	NAG	B	2	2	14,14,15	0.99	0	17,19,21	1.76	3 (17%)
2	NAG	C	1	1,2	14,14,15	1.41	1 (7%)	17,19,21	1.80	4 (23%)
2	NAG	C	2	2	14,14,15	1.08	1 (7%)	17,19,21	1.74	3 (17%)
2	NAG	D	1	1,2	14,14,15	1.05	0	17,19,21	1.38	1 (5%)
2	NAG	D	2	2	14,14,15	0.85	0	17,19,21	1.29	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	-	0/6/23/26	0/1/1/1
2	NAG	C	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
2	NAG	D	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	D	2	2	-	1/6/23/26	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1	NAG	C2-N2	-3.08	1.41	1.46
2	C	2	NAG	C2-N2	-2.15	1.42	1.46
2	B	1	NAG	C2-N2	-2.10	1.42	1.46
2	B	1	NAG	O5-C1	-2.01	1.40	1.43

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	NAG	O5-C1-C2	-5.22	103.21	111.29
2	B	1	NAG	C1-O5-C5	4.96	118.83	112.19
2	C	1	NAG	C1-O5-C5	4.84	118.67	112.19
2	C	2	NAG	O5-C1-C2	-4.61	104.15	111.29
2	D	2	NAG	O5-C1-C2	-4.19	104.80	111.29
2	D	1	NAG	O5-C1-C2	-3.56	105.78	111.29
2	B	2	NAG	C1-O5-C5	3.54	116.94	112.19
2	C	1	NAG	C3-C4-C5	-3.01	104.77	110.23
2	C	2	NAG	O3-C3-C4	-2.94	103.44	110.38
2	B	1	NAG	C2-N2-C7	2.31	125.99	122.90
2	B	1	NAG	C1-C2-N2	-2.25	106.88	110.43
2	C	1	NAG	C2-N2-C7	-2.18	119.98	122.90
2	C	1	NAG	O4-C4-C3	-2.14	105.33	110.38
2	B	2	NAG	O4-C4-C3	-2.11	105.39	110.38
2	C	2	NAG	C1-O5-C5	2.05	114.94	112.19

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	2	NAG	C8-C7-N2-C2

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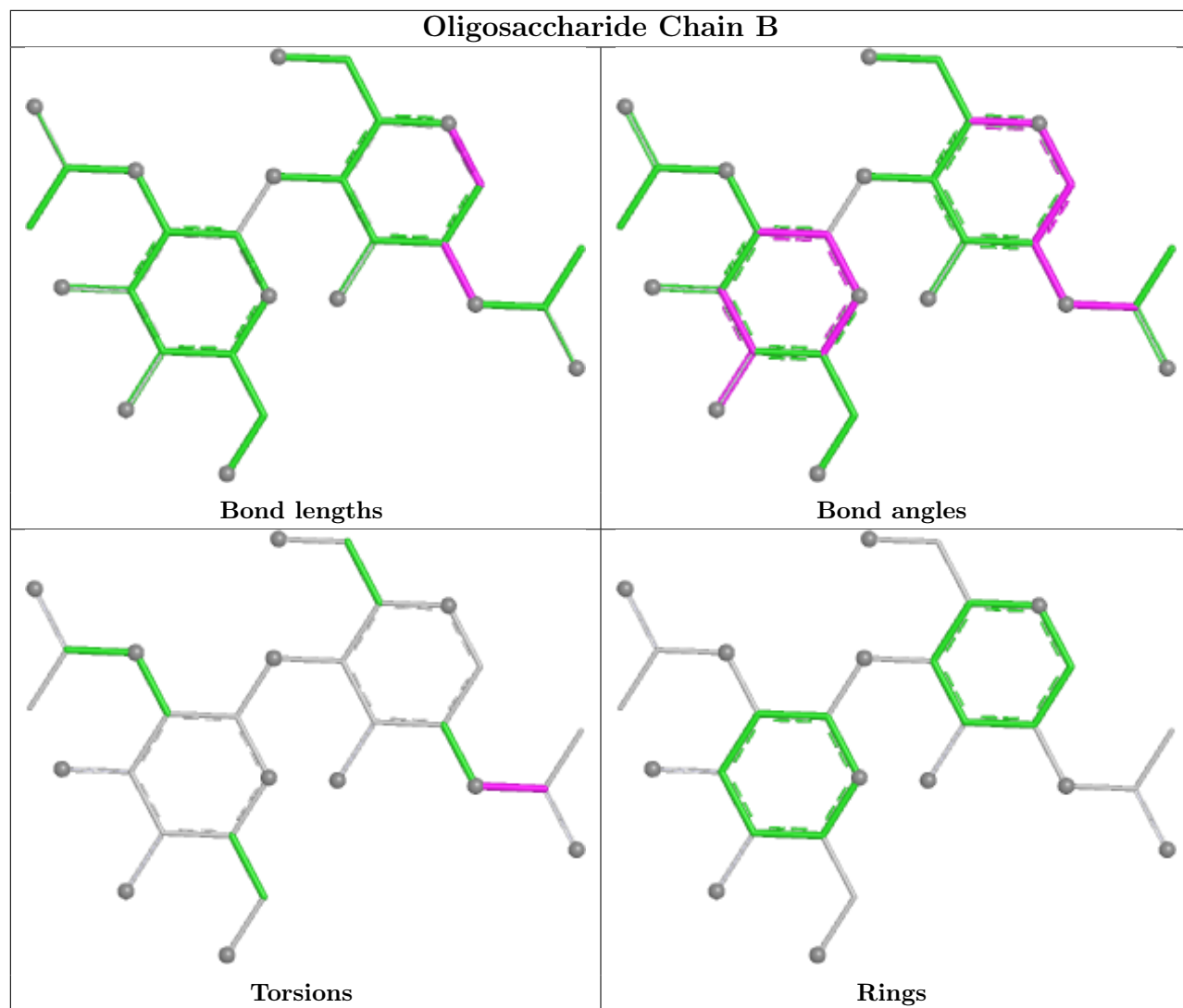
Mol	Chain	Res	Type	Atoms
2	C	2	NAG	O7-C7-N2-C2
2	D	1	NAG	C8-C7-N2-C2
2	D	1	NAG	O7-C7-N2-C2
2	B	1	NAG	C8-C7-N2-C2
2	B	1	NAG	O7-C7-N2-C2
2	C	1	NAG	C8-C7-N2-C2
2	D	2	NAG	O5-C5-C6-O6
2	C	1	NAG	O7-C7-N2-C2

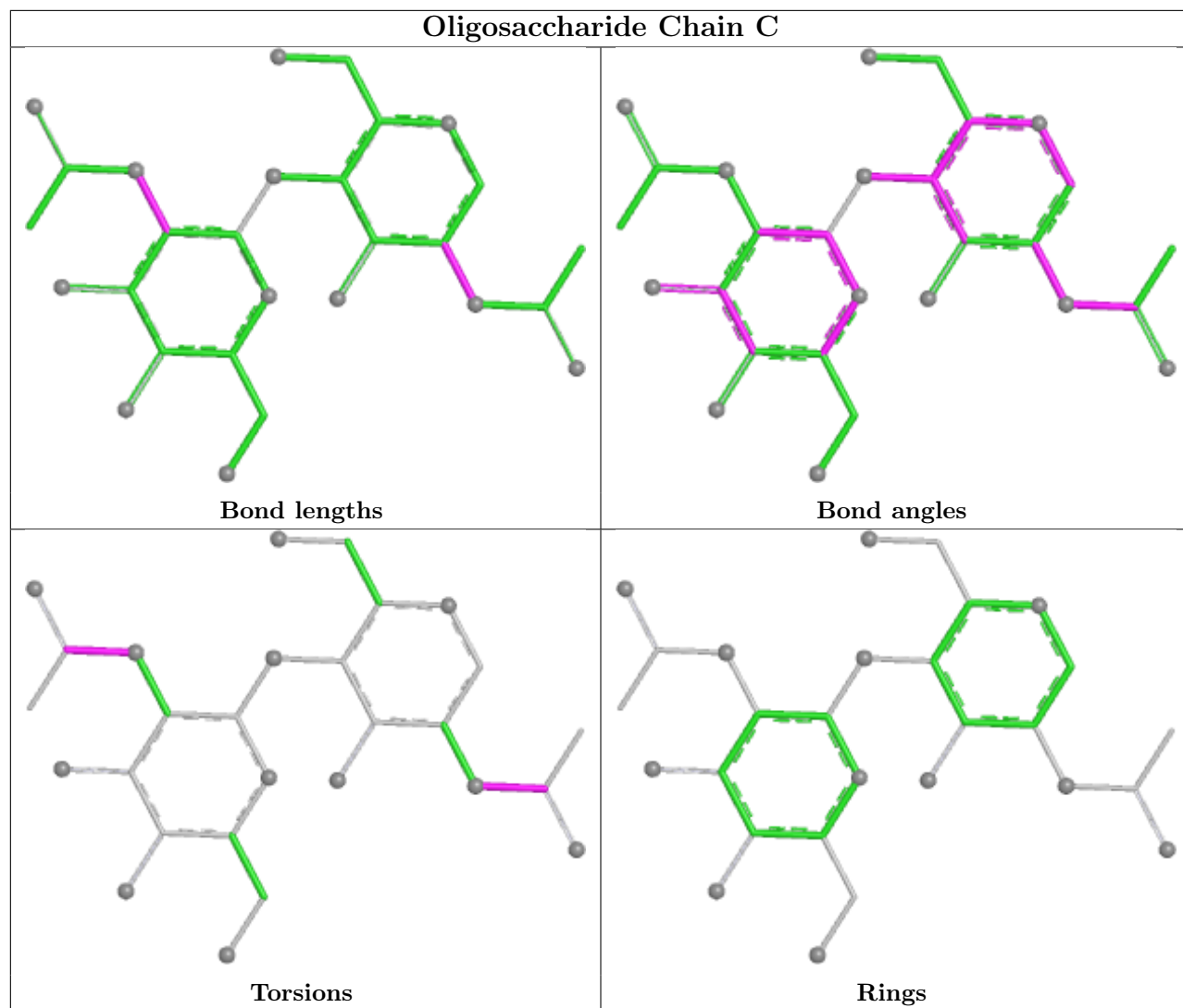
There are no ring outliers.

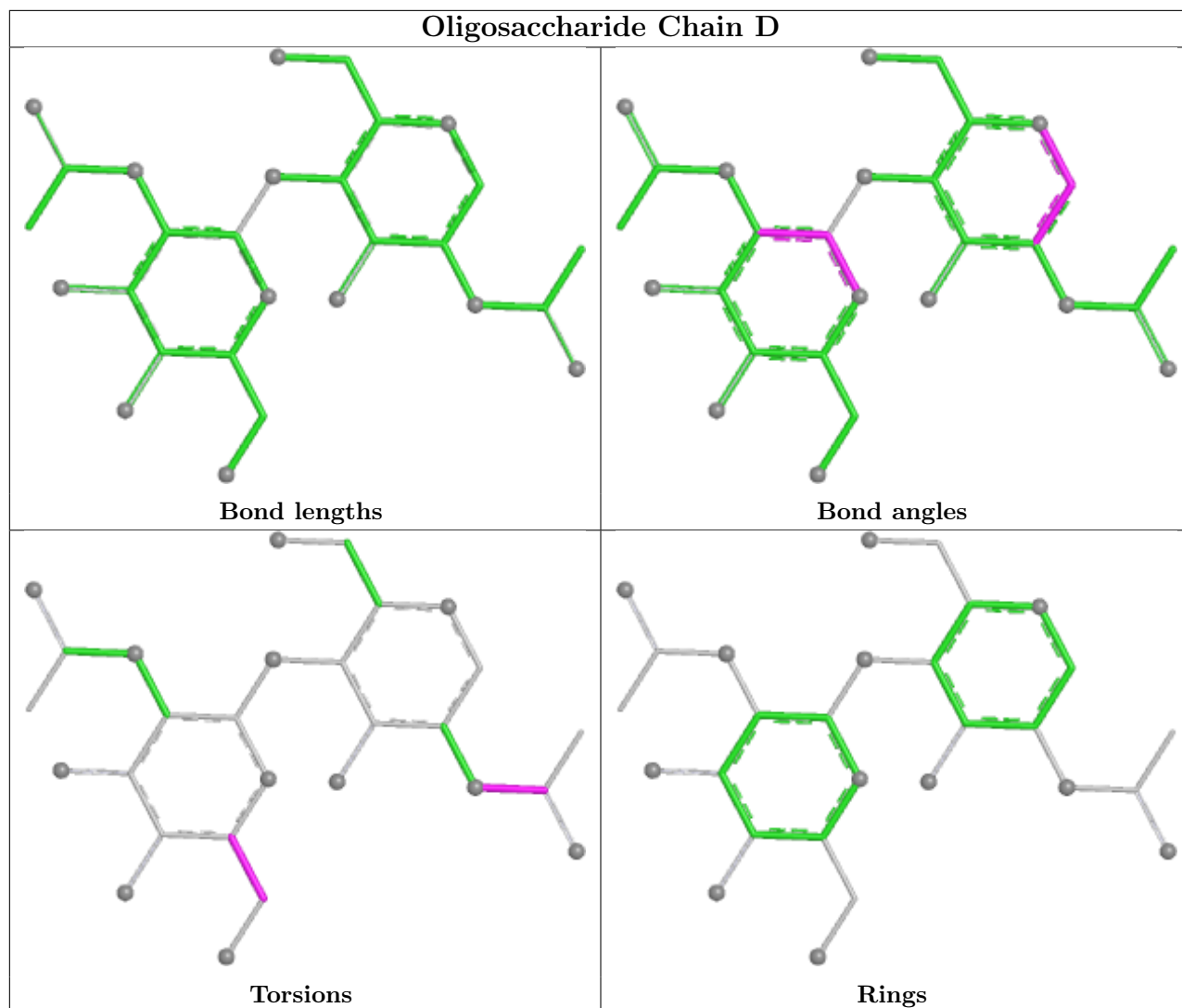
3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	NAG	1	0
2	B	1	NAG	1	0
2	C	2	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 oligosaccharide.







5.6 Ligand geometry [i](#)

57 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$
4	NAG	A	704	1	14,14,15	1.17	1 (7%)	17,19,21	1.99	6 (35%)
5	EDO	A	706	-	3,3,3	0.28	0	2,2,2	0.63	0
5	EDO	A	708	-	3,3,3	0.39	0	2,2,2	0.55	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	EDO	A	740	-	3,3,3	0.45	0	2,2,2	0.51	0
5	EDO	A	738	-	3,3,3	0.49	0	2,2,2	0.48	0
5	EDO	A	709	-	3,3,3	0.41	0	2,2,2	0.50	0
5	EDO	A	717	-	3,3,3	0.45	0	2,2,2	0.87	0
5	EDO	A	722	-	3,3,3	0.38	0	2,2,2	0.60	0
5	EDO	A	716	-	3,3,3	0.43	0	2,2,2	0.48	0
5	EDO	A	752	-	3,3,3	0.30	0	2,2,2	0.30	0
3	MAN	A	703	-	11,11,12	0.25	0	15,15,17	0.52	0
5	EDO	A	736	-	3,3,3	0.44	0	2,2,2	0.37	0
5	EDO	A	718	-	3,3,3	0.46	0	2,2,2	0.35	0
3	MAN	A	702	1	11,11,12	0.25	0	15,15,17	0.53	0
5	EDO	A	726	-	3,3,3	0.45	0	2,2,2	0.64	0
5	EDO	A	725	-	3,3,3	0.43	0	2,2,2	0.60	0
5	EDO	A	729	-	3,3,3	0.60	0	2,2,2	0.26	0
5	EDO	A	719	-	3,3,3	0.35	0	2,2,2	0.69	0
5	EDO	A	732	-	3,3,3	0.47	0	2,2,2	0.24	0
5	EDO	A	742	-	3,3,3	0.46	0	2,2,2	0.42	0
5	EDO	A	743	-	3,3,3	0.43	0	2,2,2	0.45	0
5	EDO	A	754	-	3,3,3	0.51	0	2,2,2	0.33	0
5	EDO	A	737	-	3,3,3	0.22	0	2,2,2	0.18	0
5	EDO	A	727	-	3,3,3	0.50	0	2,2,2	0.40	0
5	EDO	A	713	-	3,3,3	0.51	0	2,2,2	0.40	0
5	EDO	A	714	-	3,3,3	0.42	0	2,2,2	0.51	0
5	EDO	A	733	-	3,3,3	0.47	0	2,2,2	0.64	0
5	EDO	A	735	-	3,3,3	0.46	0	2,2,2	0.41	0
5	EDO	A	705	-	3,3,3	0.58	0	2,2,2	0.17	0
3	MAN	A	701	1	11,11,12	0.84	0	15,15,17	1.27	2 (13%)
5	EDO	A	756	-	3,3,3	0.47	0	2,2,2	0.39	0
5	EDO	A	747	-	3,3,3	0.48	0	2,2,2	0.23	0
5	EDO	A	757	-	3,3,3	0.43	0	2,2,2	0.67	0
5	EDO	A	731	-	3,3,3	0.49	0	2,2,2	0.40	0
5	EDO	A	723	-	3,3,3	0.48	0	2,2,2	0.61	0
5	EDO	A	707	-	3,3,3	0.64	0	2,2,2	0.10	0
5	EDO	A	755	-	3,3,3	0.51	0	2,2,2	0.36	0
5	EDO	A	744	-	3,3,3	0.50	0	2,2,2	0.24	0
5	EDO	A	745	-	3,3,3	0.52	0	2,2,2	0.26	0
5	EDO	A	746	-	3,3,3	0.41	0	2,2,2	0.58	0
5	EDO	A	741	-	3,3,3	0.48	0	2,2,2	0.27	0
5	EDO	A	712	-	3,3,3	0.46	0	2,2,2	0.41	0
5	EDO	A	715	-	3,3,3	0.38	0	2,2,2	0.65	0
5	EDO	A	753	-	3,3,3	0.46	0	2,2,2	0.15	0
5	EDO	A	749	-	3,3,3	0.45	0	2,2,2	0.25	0
5	EDO	A	748	-	3,3,3	0.47	0	2,2,2	0.45	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	EDO	A	730	-	3,3,3	0.24	0	2,2,2	1.01	0
5	EDO	A	728	-	3,3,3	0.40	0	2,2,2	0.75	0
5	EDO	A	721	-	3,3,3	0.48	0	2,2,2	0.39	0
5	EDO	A	739	-	3,3,3	0.43	0	2,2,2	0.38	0
5	EDO	A	724	-	3,3,3	0.52	0	2,2,2	0.29	0
5	EDO	A	751	-	3,3,3	0.47	0	2,2,2	0.33	0
5	EDO	A	734	-	3,3,3	0.31	0	2,2,2	0.68	0
5	EDO	A	720	-	3,3,3	0.40	0	2,2,2	0.35	0
5	EDO	A	750	-	3,3,3	0.49	0	2,2,2	0.42	0
5	EDO	A	711	-	3,3,3	0.17	0	2,2,2	0.90	0
5	EDO	A	710	-	3,3,3	0.52	0	2,2,2	0.51	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
4	NAG	A	704	1	-	2/6/23/26	0/1/1/1
5	EDO	A	706	-	-	1/1/1/1	-
5	EDO	A	708	-	-	0/1/1/1	-
5	EDO	A	740	-	-	1/1/1/1	-
5	EDO	A	738	-	-	1/1/1/1	-
5	EDO	A	709	-	-	0/1/1/1	-
5	EDO	A	717	-	-	1/1/1/1	-
5	EDO	A	722	-	-	0/1/1/1	-
5	EDO	A	716	-	-	0/1/1/1	-
5	EDO	A	752	-	-	1/1/1/1	-
3	MAN	A	703	-	-	0/2/19/22	0/1/1/1
5	EDO	A	736	-	-	0/1/1/1	-
5	EDO	A	718	-	-	0/1/1/1	-
3	MAN	A	702	1	-	1/2/19/22	0/1/1/1
5	EDO	A	726	-	-	1/1/1/1	-
5	EDO	A	725	-	-	1/1/1/1	-
5	EDO	A	729	-	-	0/1/1/1	-
5	EDO	A	719	-	-	1/1/1/1	-
5	EDO	A	732	-	-	1/1/1/1	-
5	EDO	A	742	-	-	1/1/1/1	-
5	EDO	A	743	-	-	0/1/1/1	-
5	EDO	A	754	-	-	0/1/1/1	-
5	EDO	A	737	-	-	0/1/1/1	-
5	EDO	A	727	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	A	713	-	-	1/1/1/1	-
5	EDO	A	714	-	-	1/1/1/1	-
5	EDO	A	733	-	-	1/1/1/1	-
5	EDO	A	735	-	-	1/1/1/1	-
5	EDO	A	705	-	-	0/1/1/1	-
3	MAN	A	701	1	-	2/2/19/22	0/1/1/1
5	EDO	A	756	-	-	0/1/1/1	-
5	EDO	A	747	-	-	0/1/1/1	-
5	EDO	A	757	-	-	1/1/1/1	-
5	EDO	A	731	-	-	0/1/1/1	-
5	EDO	A	723	-	-	0/1/1/1	-
5	EDO	A	707	-	-	1/1/1/1	-
5	EDO	A	755	-	-	1/1/1/1	-
5	EDO	A	744	-	-	0/1/1/1	-
5	EDO	A	745	-	-	0/1/1/1	-
5	EDO	A	746	-	-	0/1/1/1	-
5	EDO	A	741	-	-	0/1/1/1	-
5	EDO	A	712	-	-	1/1/1/1	-
5	EDO	A	715	-	-	1/1/1/1	-
5	EDO	A	753	-	-	0/1/1/1	-
5	EDO	A	749	-	-	0/1/1/1	-
5	EDO	A	748	-	-	0/1/1/1	-
5	EDO	A	730	-	-	1/1/1/1	-
5	EDO	A	728	-	-	1/1/1/1	-
5	EDO	A	721	-	-	0/1/1/1	-
5	EDO	A	739	-	-	0/1/1/1	-
5	EDO	A	724	-	-	1/1/1/1	-
5	EDO	A	751	-	-	0/1/1/1	-
5	EDO	A	734	-	-	1/1/1/1	-
5	EDO	A	720	-	-	0/1/1/1	-
5	EDO	A	750	-	-	0/1/1/1	-
5	EDO	A	711	-	-	0/1/1/1	-
5	EDO	A	710	-	-	0/1/1/1	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	704	NAG	C2-N2	-2.25	1.42	1.46

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	704	NAG	O4-C4-C5	-4.01	99.46	109.32
4	A	704	NAG	O5-C1-C2	-3.36	106.09	111.29
4	A	704	NAG	O5-C5-C6	3.28	114.04	107.66
3	A	701	MAN	C1-O5-C5	3.27	116.56	112.19
4	A	704	NAG	C1-O5-C5	2.79	115.92	112.19
3	A	701	MAN	O2-C2-C3	-2.57	104.82	110.15
4	A	704	NAG	C8-C7-N2	2.19	119.75	116.12
4	A	704	NAG	O4-C4-C3	-2.17	105.27	110.38

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	704	NAG	C8-C7-N2-C2
4	A	704	NAG	O7-C7-N2-C2
3	A	701	MAN	C4-C5-C6-O6
5	A	714	EDO	O1-C1-C2-O2
5	A	730	EDO	O1-C1-C2-O2
5	A	755	EDO	O1-C1-C2-O2
3	A	702	MAN	O5-C5-C6-O6
3	A	701	MAN	O5-C5-C6-O6
5	A	713	EDO	O1-C1-C2-O2
5	A	717	EDO	O1-C1-C2-O2
5	A	738	EDO	O1-C1-C2-O2
5	A	707	EDO	O1-C1-C2-O2
5	A	734	EDO	O1-C1-C2-O2
5	A	727	EDO	O1-C1-C2-O2
5	A	757	EDO	O1-C1-C2-O2
5	A	706	EDO	O1-C1-C2-O2
5	A	712	EDO	O1-C1-C2-O2
5	A	715	EDO	O1-C1-C2-O2
5	A	724	EDO	O1-C1-C2-O2
5	A	732	EDO	O1-C1-C2-O2
5	A	742	EDO	O1-C1-C2-O2
5	A	752	EDO	O1-C1-C2-O2
5	A	719	EDO	O1-C1-C2-O2
5	A	725	EDO	O1-C1-C2-O2
5	A	733	EDO	O1-C1-C2-O2
5	A	726	EDO	O1-C1-C2-O2
5	A	728	EDO	O1-C1-C2-O2
5	A	740	EDO	O1-C1-C2-O2
5	A	735	EDO	O1-C1-C2-O2

There are no ring outliers.

12 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	740	EDO	1	0
5	A	709	EDO	1	0
5	A	716	EDO	1	0
3	A	703	MAN	1	0
5	A	726	EDO	1	0
5	A	733	EDO	1	0
5	A	707	EDO	2	0
5	A	728	EDO	1	0
5	A	724	EDO	2	0
5	A	734	EDO	1	0
5	A	711	EDO	1	0
5	A	710	EDO	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

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	542/578 (93%)	0.13	13 (2%) 59 62	36, 46, 62, 82	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	388	ASP	4.0
1	A	101	TYR	3.4
1	A	70	TYR	3.3
1	A	99	ALA	3.1
1	A	56	THR	2.8
1	A	508	GLN	2.4
1	A	100	GLY	2.4
1	A	79	THR	2.3
1	A	502	ASP	2.3
1	A	73	GLN	2.2
1	A	484	GLN	2.1
1	A	503	ALA	2.0
1	A	596	HIS	2.0

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

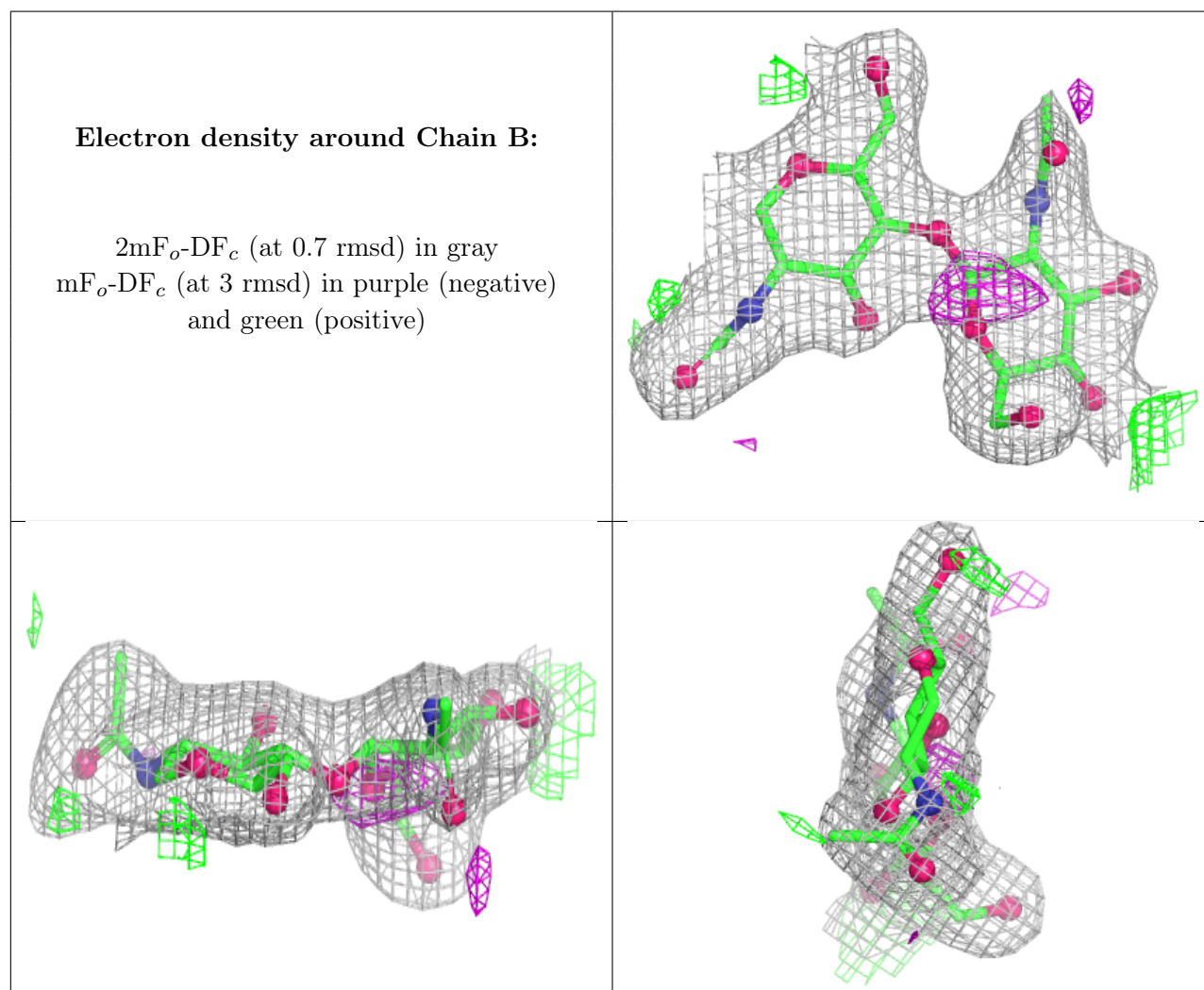
There are no non-standard protein/DNA/RNA residues in this entry.

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.

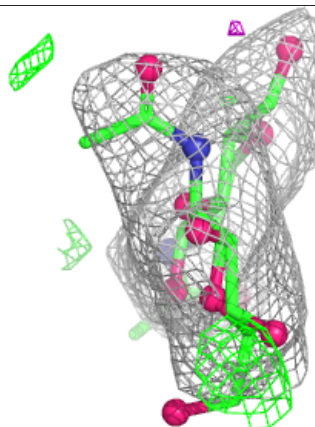
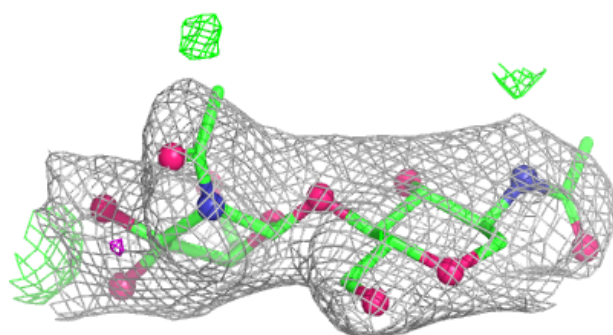
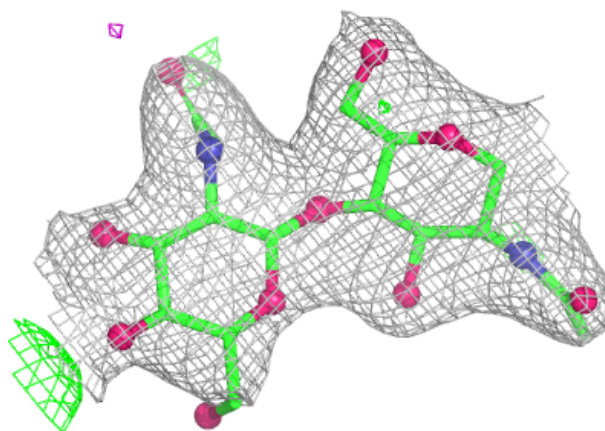
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	NAG	B	2	14/15	0.78	0.16	66,73,80,92	0
2	NAG	C	2	14/15	0.79	0.13	80,94,106,109	0
2	NAG	D	2	14/15	0.79	0.12	89,96,103,103	0
2	NAG	D	1	14/15	0.86	0.11	70,76,85,88	0
2	NAG	C	1	14/15	0.90	0.12	64,72,75,79	0
2	NAG	B	1	14/15	0.91	0.10	54,59,62,65	0

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.

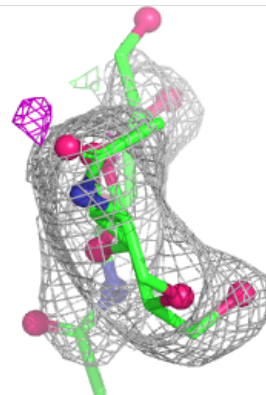
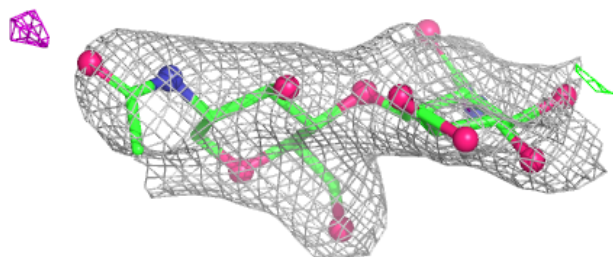
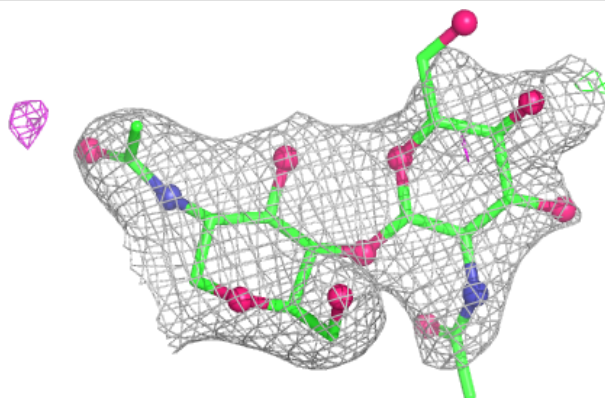


Electron density around Chain C:

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

**Electron density around Chain D:**

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



6.4 Ligands 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	MAN	A	703	11/12	0.66	0.19	63,69,84,85	0
3	MAN	A	702	11/12	0.67	0.14	69,77,87,88	0
5	EDO	A	757	4/4	0.68	0.26	66,68,71,77	0
5	EDO	A	750	4/4	0.69	0.18	67,69,79,80	0
5	EDO	A	739	4/4	0.74	0.24	64,65,69,73	0
5	EDO	A	748	4/4	0.76	0.22	63,64,66,68	0
5	EDO	A	746	4/4	0.77	0.23	64,65,66,82	0
5	EDO	A	752	4/4	0.78	0.19	69,69,72,74	0
5	EDO	A	743	4/4	0.78	0.18	72,73,76,83	0
5	EDO	A	736	4/4	0.79	0.18	60,69,71,75	0
5	EDO	A	707	4/4	0.80	0.20	51,55,65,66	0
5	EDO	A	713	4/4	0.81	0.19	58,60,65,68	0
5	EDO	A	721	4/4	0.82	0.19	65,69,72,74	0
5	EDO	A	740	4/4	0.82	0.20	66,66,72,72	0
5	EDO	A	718	4/4	0.83	0.15	61,65,69,71	0
4	NAG	A	704	14/15	0.83	0.13	50,62,73,83	0
5	EDO	A	723	4/4	0.83	0.19	54,61,66,67	0
5	EDO	A	732	4/4	0.83	0.15	62,68,69,78	0
5	EDO	A	744	4/4	0.83	0.15	69,71,72,75	0
5	EDO	A	738	4/4	0.84	0.17	61,62,67,73	0
5	EDO	A	726	4/4	0.84	0.21	55,58,59,67	0
5	EDO	A	729	4/4	0.84	0.18	58,64,65,68	0
5	EDO	A	756	4/4	0.84	0.33	62,62,64,70	0
5	EDO	A	747	4/4	0.84	0.19	65,66,69,72	0
5	EDO	A	722	4/4	0.85	0.18	60,61,63,68	0
5	EDO	A	706	4/4	0.85	0.19	53,54,56,59	0
5	EDO	A	715	4/4	0.85	0.21	54,56,57,60	0
5	EDO	A	755	4/4	0.85	0.18	66,67,68,75	0
5	EDO	A	727	4/4	0.85	0.17	58,60,65,69	0
5	EDO	A	728	4/4	0.85	0.19	58,60,66,67	0
5	EDO	A	749	4/4	0.86	0.15	65,66,67,68	0
5	EDO	A	741	4/4	0.86	0.14	66,66,69,75	0
5	EDO	A	734	4/4	0.86	0.19	45,50,53,54	0
5	EDO	A	708	4/4	0.87	0.15	60,60,64,66	0
5	EDO	A	724	4/4	0.87	0.15	61,67,69,70	0
5	EDO	A	711	4/4	0.87	0.17	58,66,66,68	0
5	EDO	A	716	4/4	0.87	0.14	72,73,76,78	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	EDO	A	745	4/4	0.88	0.13	64,66,67,74	0
5	EDO	A	725	4/4	0.88	0.14	57,59,60,70	0
5	EDO	A	753	4/4	0.88	0.15	59,60,68,73	0
5	EDO	A	717	4/4	0.88	0.15	52,60,61,62	0
5	EDO	A	710	4/4	0.88	0.14	48,53,54,56	0
5	EDO	A	733	4/4	0.88	0.16	57,60,62,67	0
3	MAN	A	701	11/12	0.89	0.11	56,60,68,76	0
5	EDO	A	742	4/4	0.89	0.23	55,59,60,63	0
5	EDO	A	735	4/4	0.89	0.14	67,70,70,77	0
5	EDO	A	754	4/4	0.91	0.14	64,67,68,70	0
5	EDO	A	751	4/4	0.91	0.19	53,56,65,68	0
5	EDO	A	720	4/4	0.91	0.14	54,56,57,67	0
5	EDO	A	714	4/4	0.91	0.14	62,63,64,76	0
5	EDO	A	719	4/4	0.92	0.12	52,52,52,58	0
5	EDO	A	730	4/4	0.92	0.18	54,58,59,64	0
5	EDO	A	731	4/4	0.92	0.14	56,56,60,64	0
5	EDO	A	709	4/4	0.92	0.14	59,59,61,61	0
5	EDO	A	712	4/4	0.93	0.15	51,54,59,60	0
5	EDO	A	737	4/4	0.93	0.14	55,60,67,69	0
5	EDO	A	705	4/4	0.95	0.14	44,52,52,56	0

6.5 Other polymers [i](#)

There are no such residues in this entry.