wwPDB X-ray Structure Validation Summary Report ⓘ

Dec 22, 2022 – 06:32 pm GMT

PDB ID	:	8BET
Title	:	Structure of D188A-fructofuranosidase from Rhodotorula dairenesis in complex with sucrose
Authors	:	Jimenez-Ortega, E.; Sanz-Aparicio, J.
Deposited on	:	2022-10-21
Resolution	:	2.38 Å(reported)

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

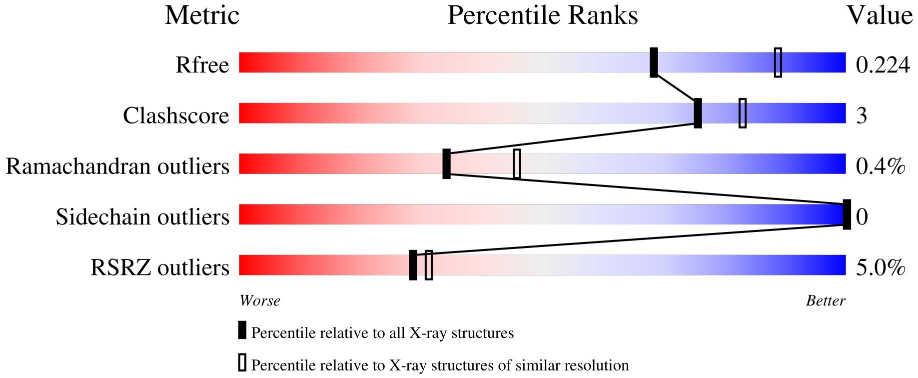
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.31.3
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.3

1 Overall quality at a glance

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

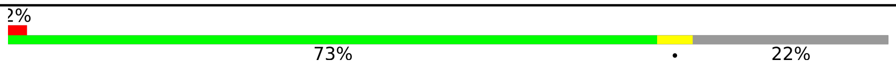
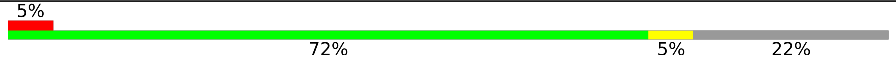
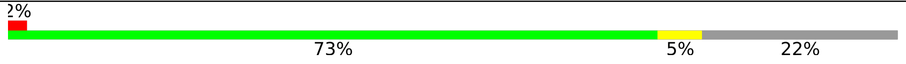
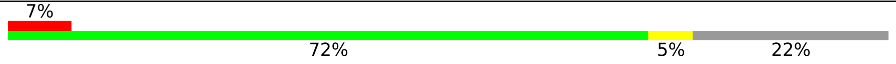
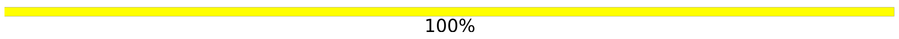
The reported resolution of this entry is 2.38 Å.

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}	130704	5509 (2.40-2.36)
Clashscore	141614	6082 (2.40-2.36)
Ramachandran outliers	138981	5973 (2.40-2.36)
Sidechain outliers	138945	5975 (2.40-2.36)
RSRZ outliers	127900	5397 (2.40-2.36)

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 <=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	675	
1	B	675	
1	C	675	
1	D	675	
2	E	2	

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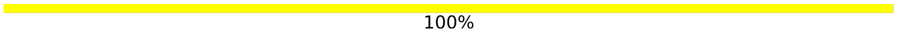
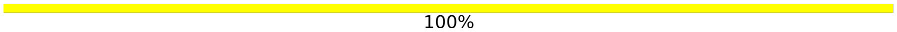
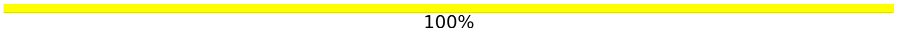
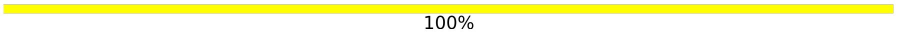
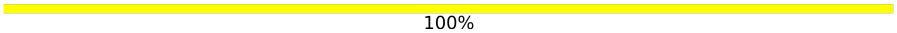
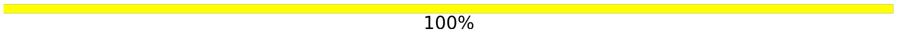
Mol	Chain	Length	Quality of chain
2	H	2	 100%
2	K	2	 100%
3	F	3	 100%
3	I	3	 67% 33%
3	J	3	 100%
4	G	6	 33% 67%
5	L	2	 100%
5	M	2	 100%
5	N	2	 50% 50%
5	O	2	 50% 50%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	MAN	B	702	-	-	-	X
6	MAN	C	703	-	-	-	X
6	MAN	D	701	-	-	-	X
7	NAG	B	705	-	-	-	X

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 17723 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.

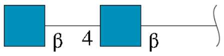
- Molecule 1 is a protein called Beta-fructofuranosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	525	Total 4032	C 2569	N 656	O 796	S 11	0	0	0
1	B	526	Total 4038	C 2572	N 657	O 798	S 11	0	0	0
1	C	528	Total 4050	C 2578	N 659	O 802	S 11	0	0	0
1	D	524	Total 4026	C 2566	N 655	O 794	S 11	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

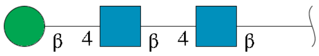
Chain	Residue	Modelled	Actual	Comment	Reference
A	188	ALA	ASP	engineered mutation	UNP A0A856TAI5
B	188	ALA	ASP	engineered mutation	UNP A0A856TAI5
C	188	ALA	ASP	engineered mutation	UNP A0A856TAI5
D	188	ALA	ASP	engineered mutation	UNP A0A856TAI5

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



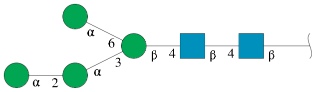
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	2	Total 28	C 16	N 2	O 10	0	0	0
2	H	2	Total 28	C 16	N 2	O 10	0	0	0
2	K	2	Total 28	C 16	N 2	O 10	0	0	0

- Molecule 3 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.



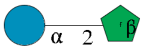
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
3	F	3	Total C N O 39 22 2 15	0	0	0
3	I	3	Total C N O 39 22 2 15	0	0	0
3	J	3	Total C N O 39 22 2 15	0	0	0

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



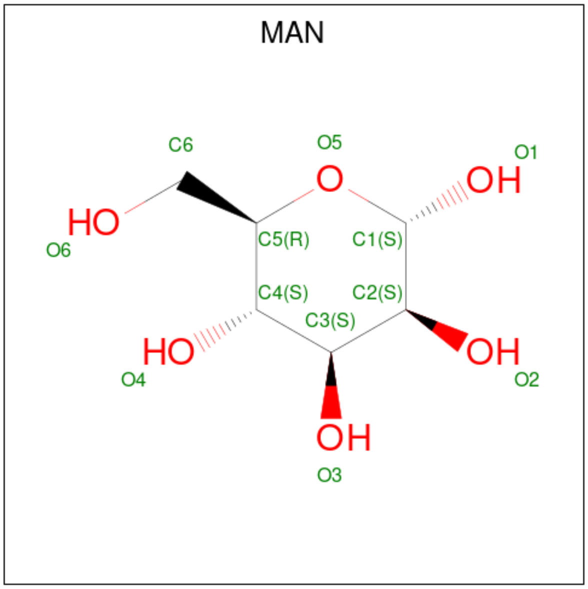
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
4	G	6	Total C N O 72 40 2 30	0	0	0

- Molecule 5 is an oligosaccharide called beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
5	L	2	Total C O 23 12 11	0	0	0
5	M	2	Total C O 23 12 11	0	0	0
5	N	2	Total C O 23 12 11	0	0	0
5	O	2	Total C O 23 12 11	0	0	0

- Molecule 6 is alpha-D-mannopyranose (three-letter code: MAN) (formula: $C_6H_{12}O_6$).



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

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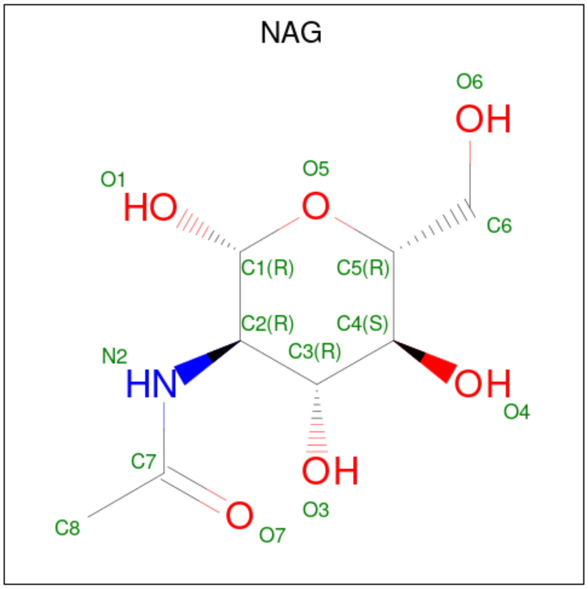
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	D	1	Total C O 11 6 5	0	0

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

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

- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	200	Total O 200 200	0	0
8	B	176	Total O 176 176	0	0
8	C	166	Total O 166 166	0	0

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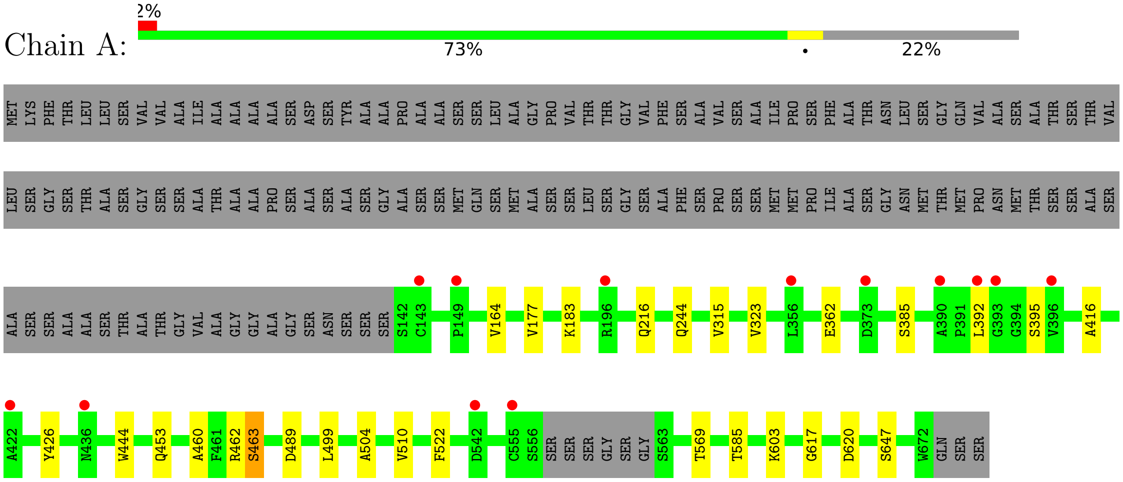
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	D	152	Total O 152 152	0	0

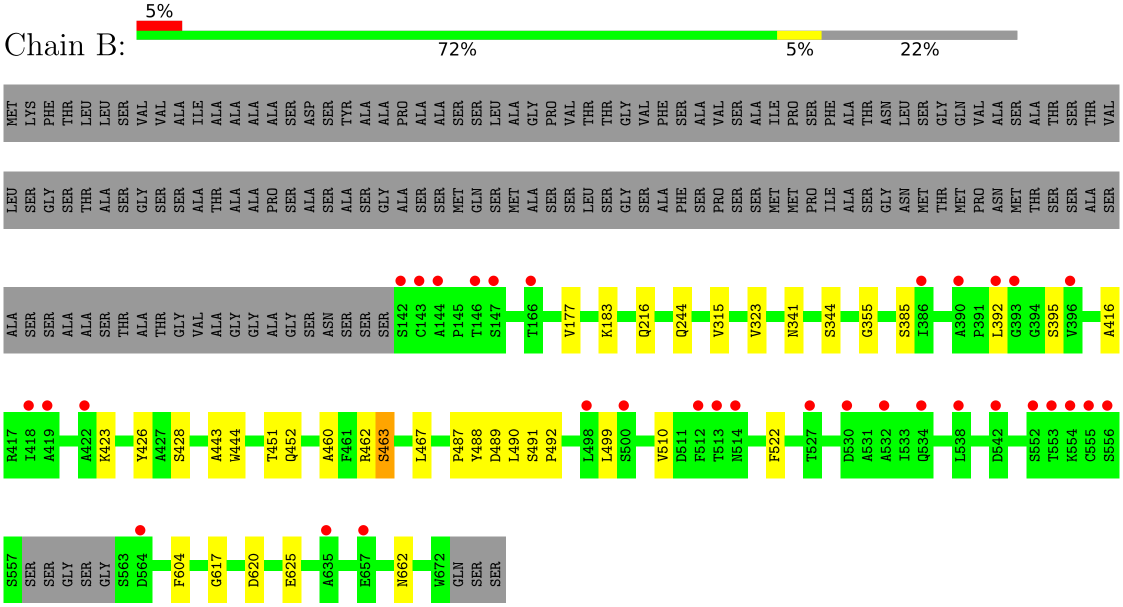
3 Residue-property plots

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

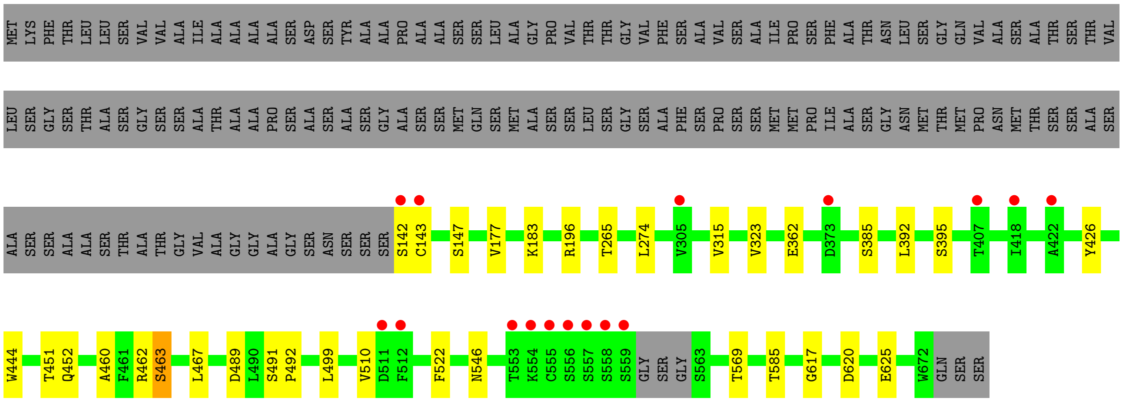
- Molecule 1: Beta-fructofuranosidase



- Molecule 1: Beta-fructofuranosidase

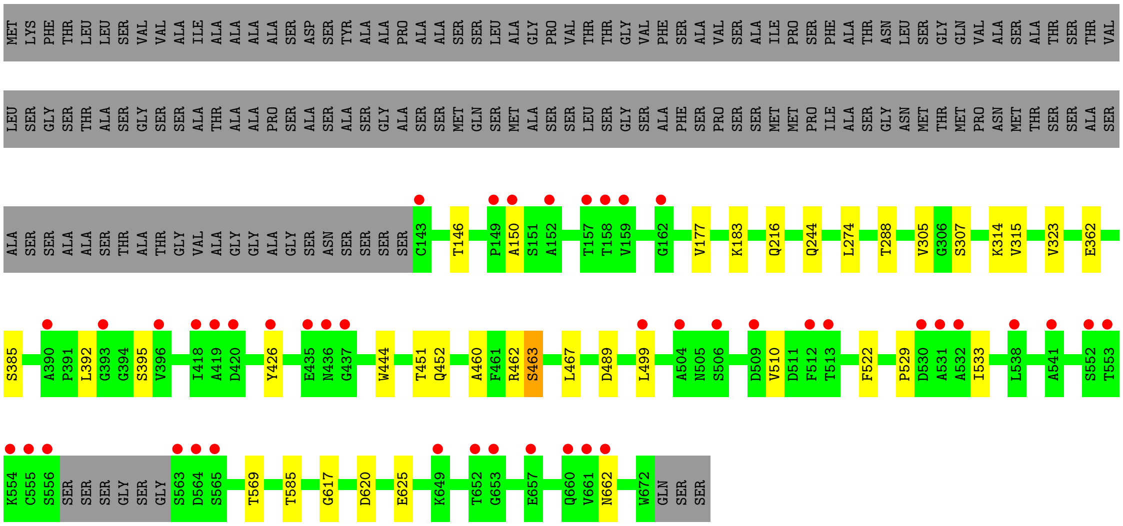


- Molecule 1: Beta-fructofuranosidase



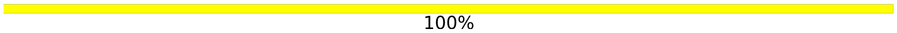
- Molecule 1: Beta-fructofuranosidase

Chain D:



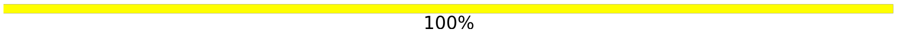
- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:



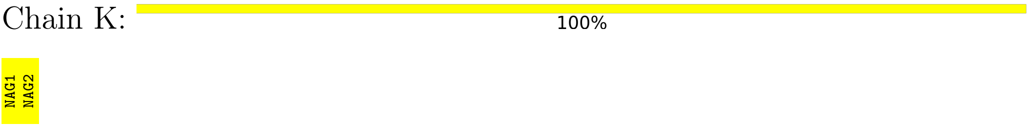
- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H:



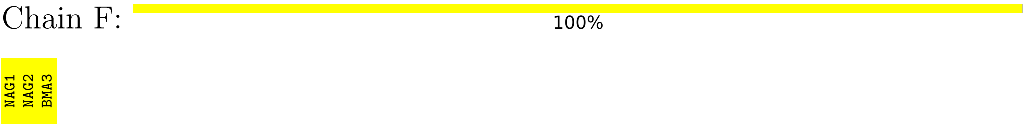
- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K:



- Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:

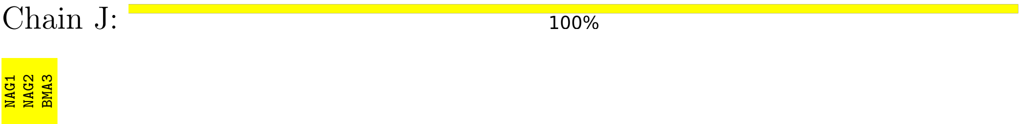


- Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:

- Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:

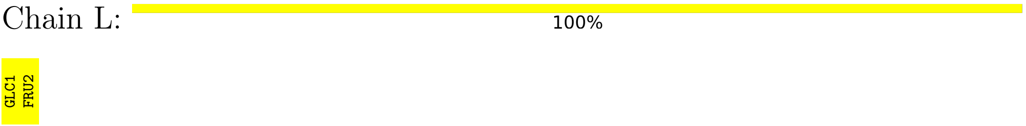


- Molecule 4: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:

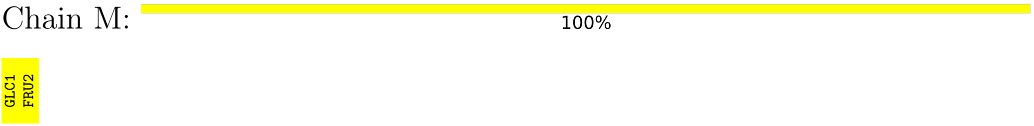
- Molecule 5: beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose

Chain L:



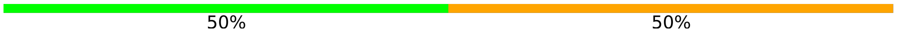
- Molecule 5: beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose

Chain M:



- Molecule 5: beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose

Chain N:



- Molecule 5: beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose

Chain O:

4 Data and refinement statistics ⓘ

Property	**Value**	**Source**
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	80.34Å 112.81Å 139.08Å 90.00° 104.74° 90.00°	Depositor
Resolution (Å)	45.69 – 2.38 45.65 – 2.38	Depositor EDS
% Data completeness (in resolution range)	99.8 (45.69-2.38) 99.8 (45.65-2.38)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.31 (at 2.37Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.184 , 0.222 0.186 , 0.224	Depositor DCC
R_{free} test set	4728 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å^2)	36.8	Xtriage
Anisotropy	0.184	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 27.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.021 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	17723	wwPDB-VP
Average B, all atoms (Å^2)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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, GLC, FRU, BMA, MAN

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	0.65	0/4160	0.75	0/5721
1	B	0.65	0/4166	0.75	0/5729
1	C	0.66	0/4178	0.75	0/5745
1	D	0.64	0/4154	0.74	0/5713
All	All	0.65	0/16658	0.75	0/22908

There are no bond length outliers.

There are no bond angle outliers.

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	4032	0	3751	21	0
1	B	4038	0	3757	29	0
1	C	4050	0	3765	25	0
1	D	4026	0	3747	27	0
2	E	28	0	25	0	0
2	H	28	0	25	0	0
2	K	28	0	25	0	0
3	F	39	0	34	0	0
3	I	39	0	34	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	J	39	0	34	0	0
4	G	72	0	61	0	0
5	L	23	0	21	1	0
5	M	23	0	21	0	0
5	N	23	0	21	1	0
5	O	23	0	21	1	0
6	A	44	0	40	0	0
6	B	33	0	30	0	0
6	C	44	0	40	2	0
6	D	33	0	30	2	0
7	A	98	0	91	2	0
7	B	98	0	91	3	0
7	C	84	0	78	1	0
7	D	84	0	78	4	0
8	A	200	0	0	5	0
8	B	176	0	0	0	0
8	C	166	0	0	0	0
8	D	152	0	0	1	0
All	All	17723	0	15820	102	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.

The worst 5 of 102 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:487:PRO:HB2	1:B:490:LEU:HD11	1.14	1.12
1:B:487:PRO:HB2	1:B:490:LEU:CD1	1.96	0.95
1:C:467:LEU:HD12	1:C:625:GLU:OE1	1.80	0.80
1:B:467:LEU:HD12	1:B:625:GLU:OE1	1.81	0.80
1:D:467:LEU:HD12	1:D:625:GLU:OE1	1.81	0.79

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	521/675 (77%)	501 (96%)	18 (4%)	2 (0%)	34	46
1	B	522/675 (77%)	501 (96%)	19 (4%)	2 (0%)	34	46
1	C	524/675 (78%)	502 (96%)	20 (4%)	2 (0%)	34	46
1	D	520/675 (77%)	503 (97%)	15 (3%)	2 (0%)	34	46
All	All	2087/2700 (77%)	2007 (96%)	72 (3%)	8 (0%)	34	46

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	463	SER
1	C	463	SER
1	A	463	SER
1	A	489	ASP
1	B	489	ASP

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	441/545 (81%)	441 (100%)	0	100	100
1	B	442/545 (81%)	442 (100%)	0	100	100
1	C	444/545 (82%)	444 (100%)	0	100	100
1	D	440/545 (81%)	440 (100%)	0	100	100
All	All	1767/2180 (81%)	1767 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	518	ASN

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Mol	Chain	Res	Type
1	D	664	GLN
1	D	666	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 ⓘ

29 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	E	1	2,1	14,14,15	0.79	0	17,19,21	0.95	1 (5%)
2	NAG	E	2	2	14,14,15	0.60	0	17,19,21	1.19	2 (11%)
3	NAG	F	1	1,3	14,14,15	0.64	0	17,19,21	1.11	1 (5%)
3	NAG	F	2	3	14,14,15	0.64	0	17,19,21	1.03	1 (5%)
3	BMA	F	3	3	11,11,12	0.71	0	15,15,17	1.09	1 (6%)
4	NAG	G	1	4,1	14,14,15	0.51	0	17,19,21	1.00	0
4	NAG	G	2	4	14,14,15	0.36	0	17,19,21	0.94	0
4	BMA	G	3	4	11,11,12	0.61	0	15,15,17	0.99	1 (6%)
4	MAN	G	4	4	11,11,12	0.50	0	15,15,17	1.09	1 (6%)
4	MAN	G	5	4	11,11,12	0.58	0	15,15,17	1.24	2 (13%)
4	MAN	G	6	4	11,11,12	0.84	0	15,15,17	1.55	3 (20%)
2	NAG	H	1	2,1	14,14,15	0.76	0	17,19,21	1.09	2 (11%)
2	NAG	H	2	2	14,14,15	0.81	0	17,19,21	1.40	1 (5%)
3	NAG	I	1	1,3	14,14,15	0.51	0	17,19,21	1.04	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\lvert Z\rvert > 2$	Counts	RMSZ	$\#\lvert Z\rvert > 2$
3	NAG	I	2	3	14,14,15	0.56	0	17,19,21	1.29	2 (11%)
3	BMA	I	3	3	11,11,12	0.62	0	15,15,17	1.12	1 (6%)
3	NAG	J	1	1,3	14,14,15	0.49	0	17,19,21	1.28	2 (11%)
3	NAG	J	2	3	14,14,15	0.54	0	17,19,21	1.12	2 (11%)
3	BMA	J	3	3	11,11,12	0.64	0	15,15,17	0.91	1 (6%)
2	NAG	K	1	2,1	14,14,15	0.57	0	17,19,21	1.20	1 (5%)
2	NAG	K	2	2	14,14,15	0.49	0	17,19,21	1.01	1 (5%)
5	GLC	L	1	5	11,11,12	0.73	0	15,15,17	1.04	0
5	FRU	L	2	5	11,12,12	0.53	0	10,18,18	0.89	1 (10%)
5	GLC	M	1	5	11,11,12	0.69	0	15,15,17	1.57	5 (33%)
5	FRU	M	2	5	11,12,12	0.75	1 (9%)	10,18,18	0.77	0
5	GLC	N	1	5	11,11,12	0.81	1 (9%)	15,15,17	0.99	2 (13%)
5	FRU	N	2	5	11,12,12	0.81	0	10,18,18	0.70	0
5	GLC	O	1	5	11,11,12	0.54	0	15,15,17	1.93	6 (40%)
5	FRU	O	2	5	11,12,12	0.62	0	10,18,18	0.97	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	E	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	E	2	2	-	0/6/23/26	0/1/1/1
3	NAG	F	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1
3	BMA	F	3	3	-	2/2/19/22	0/1/1/1
4	NAG	G	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	G	2	4	-	1/6/23/26	0/1/1/1
4	BMA	G	3	4	-	2/2/19/22	0/1/1/1
4	MAN	G	4	4	-	1/2/19/22	0/1/1/1
4	MAN	G	5	4	-	2/2/19/22	0/1/1/1
4	MAN	G	6	4	-	0/2/19/22	0/1/1/1
2	NAG	H	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	H	2	2	-	0/6/23/26	0/1/1/1
3	NAG	I	1	1,3	-	4/6/23/26	0/1/1/1
3	NAG	I	2	3	-	0/6/23/26	0/1/1/1
3	BMA	I	3	3	-	2/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	J	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	J	2	3	-	2/6/23/26	0/1/1/1
3	BMA	J	3	3	-	2/2/19/22	0/1/1/1
2	NAG	K	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	K	2	2	-	0/6/23/26	0/1/1/1
5	GLC	L	1	5	-	0/2/19/22	0/1/1/1
5	FRU	L	2	5	-	1/5/24/24	0/1/1/1
5	GLC	M	1	5	-	2/2/19/22	0/1/1/1
5	FRU	M	2	5	-	3/5/24/24	0/1/1/1
5	GLC	N	1	5	-	0/2/19/22	0/1/1/1
5	FRU	N	2	5	-	3/5/24/24	0/1/1/1
5	GLC	O	1	5	-	2/2/19/22	0/1/1/1
5	FRU	O	2	5	-	1/5/24/24	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	N	1	GLC	O5-C1	-2.07	1.40	1.43
5	M	2	FRU	O5-C2	-2.05	1.40	1.43

The worst 5 of 42 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	H	2	NAG	C1-O5-C5	4.33	118.05	112.19
5	O	1	GLC	O3-C3-C4	-3.47	102.32	110.35
3	J	1	NAG	O5-C1-C2	-3.34	106.01	111.29
2	K	2	NAG	C1-O5-C5	3.26	116.61	112.19
3	I	2	NAG	C1-O5-C5	3.14	116.44	112.19

There are no chirality outliers.

5 of 36 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	M	2	FRU	O1-C1-C2-C3
5	M	2	FRU	O1-C1-C2-O2
5	M	2	FRU	O1-C1-C2-O5
5	N	2	FRU	O1-C1-C2-O2
5	N	2	FRU	O1-C1-C2-O5

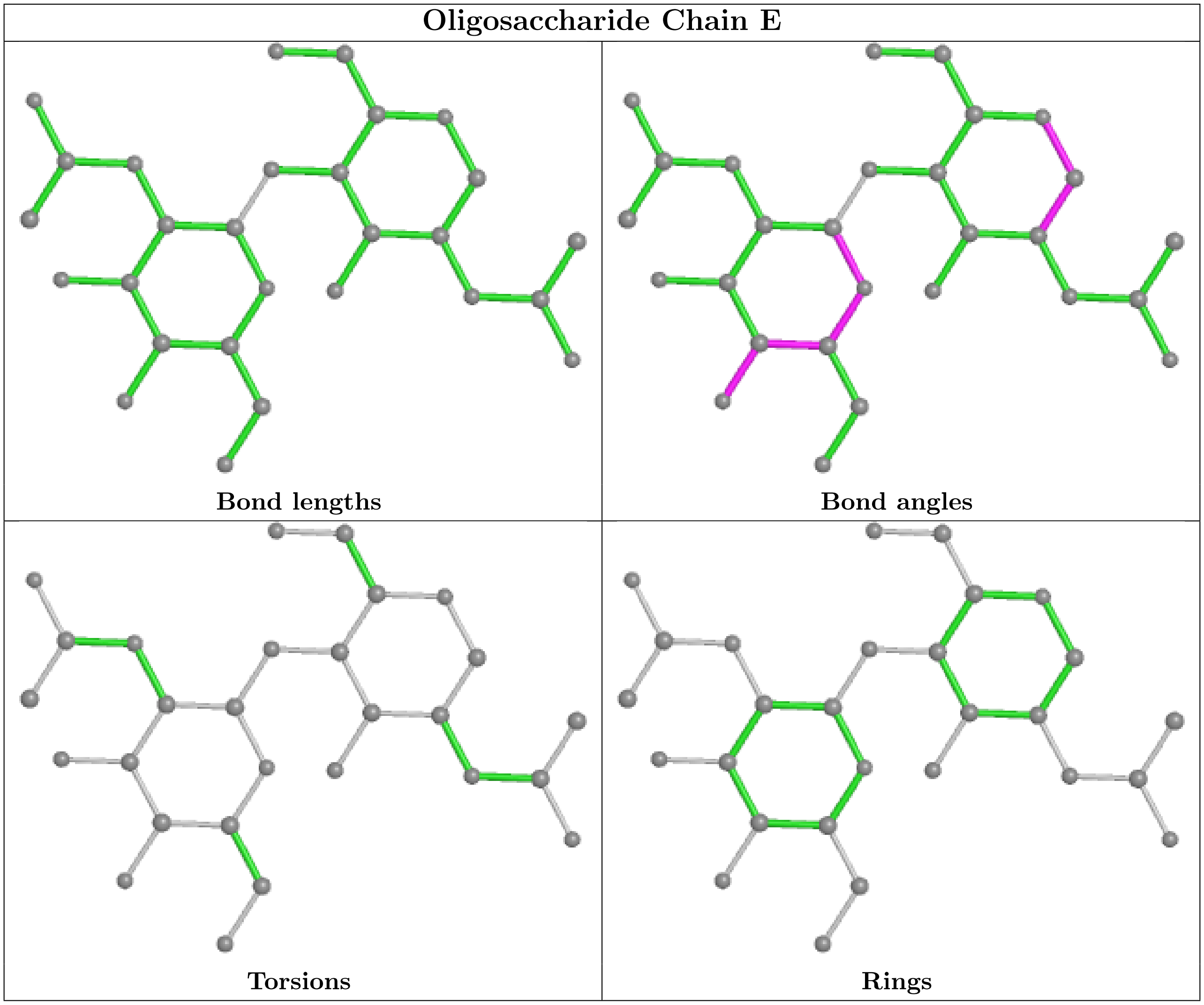
There are no ring outliers.

4 monomers are involved in 5 short contacts:

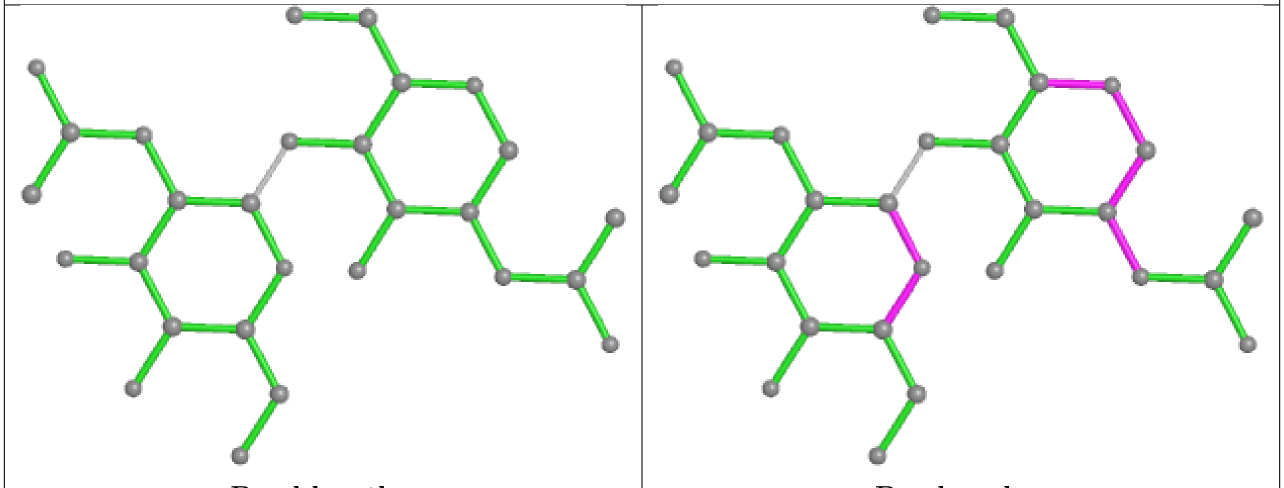
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	I	1	NAG	2	0
5	L	1	GLC	1	0
5	O	1	GLC	1	0
5	N	1	GLC	1	0

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

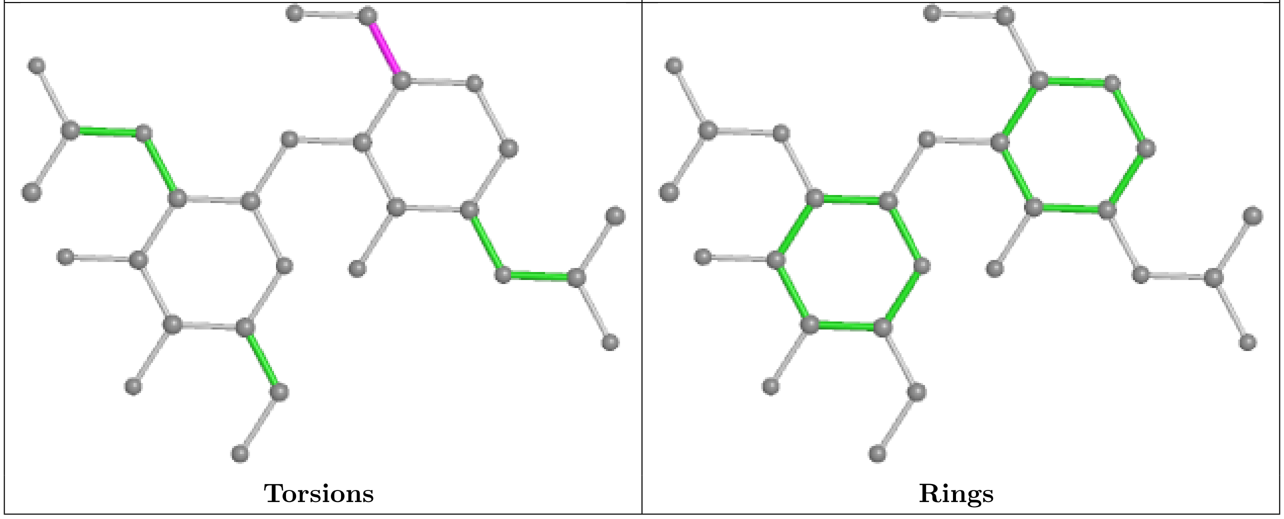
Oligosaccharide Chain E



Oligosaccharide Chain H

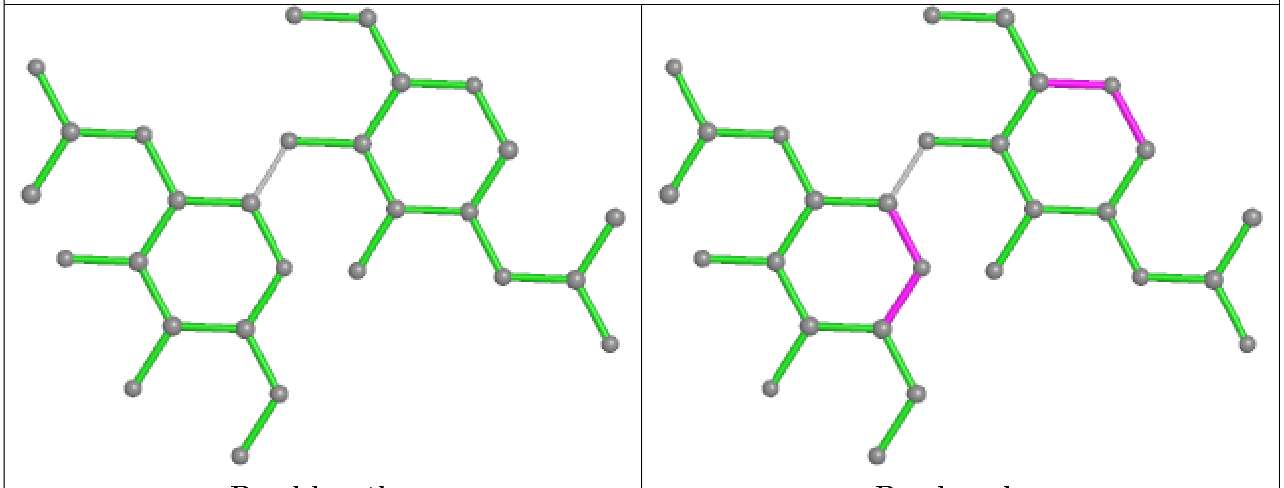


Bond lengths **Bond angles**

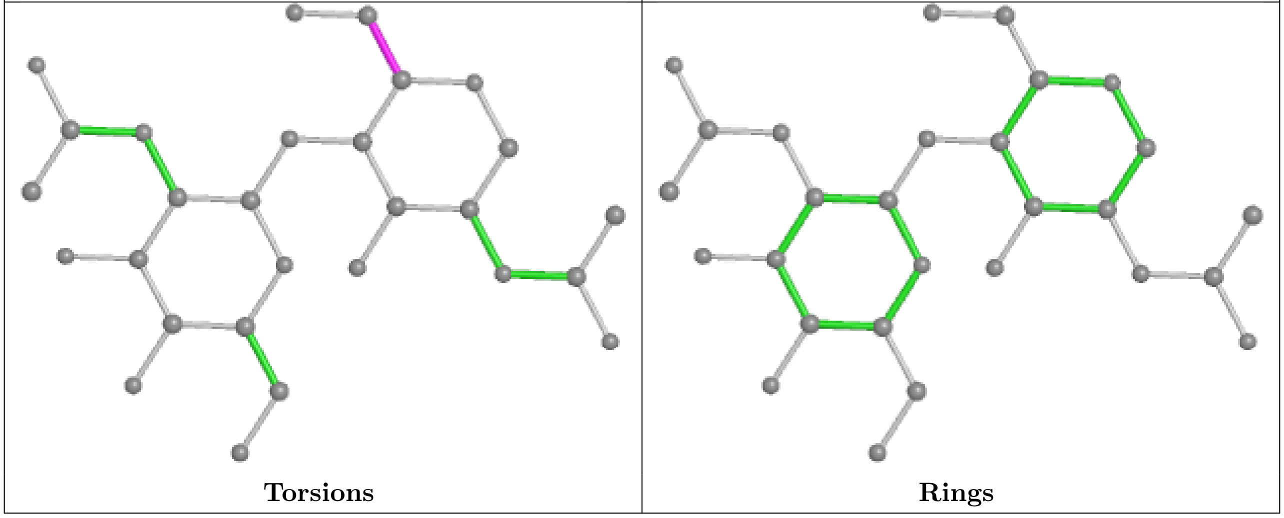


Torsions **Rings**

Oligosaccharide Chain K

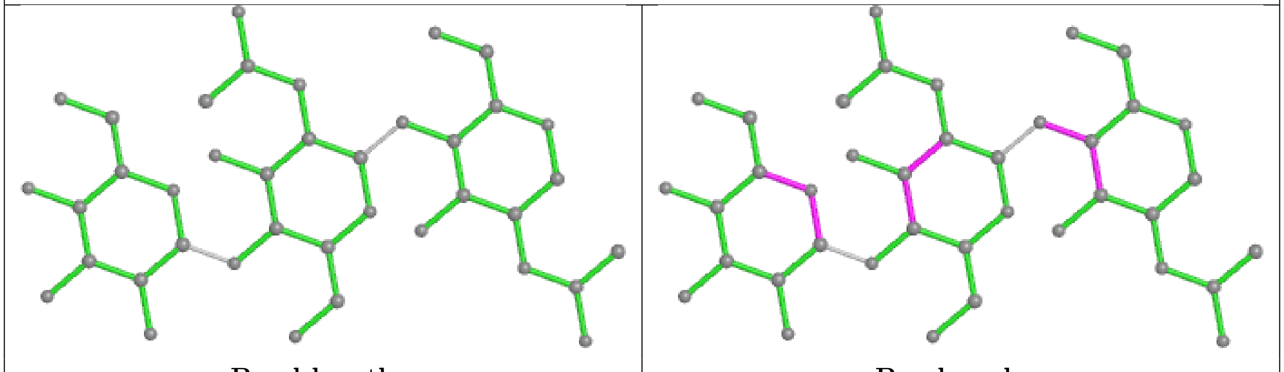


Bond lengths **Bond angles**

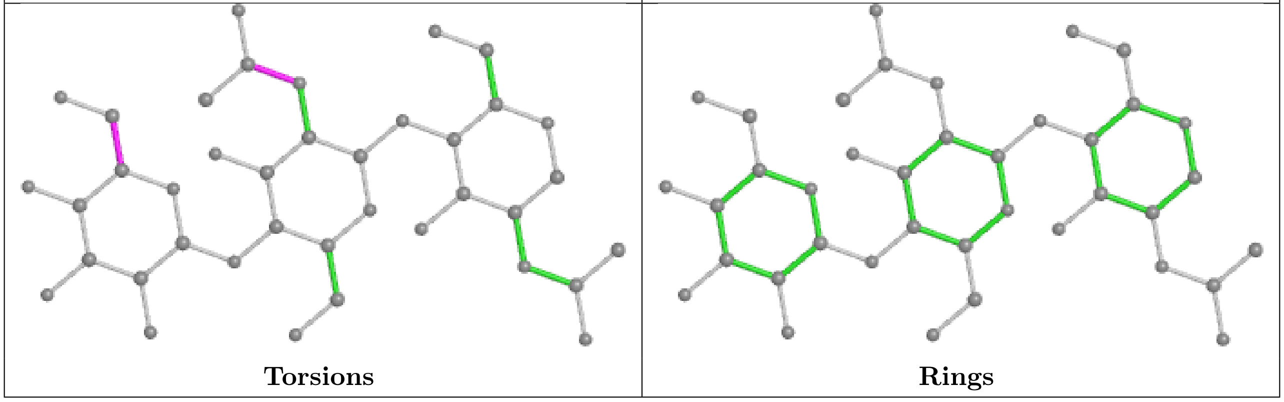


Torsions **Rings**

Oligosaccharide Chain F

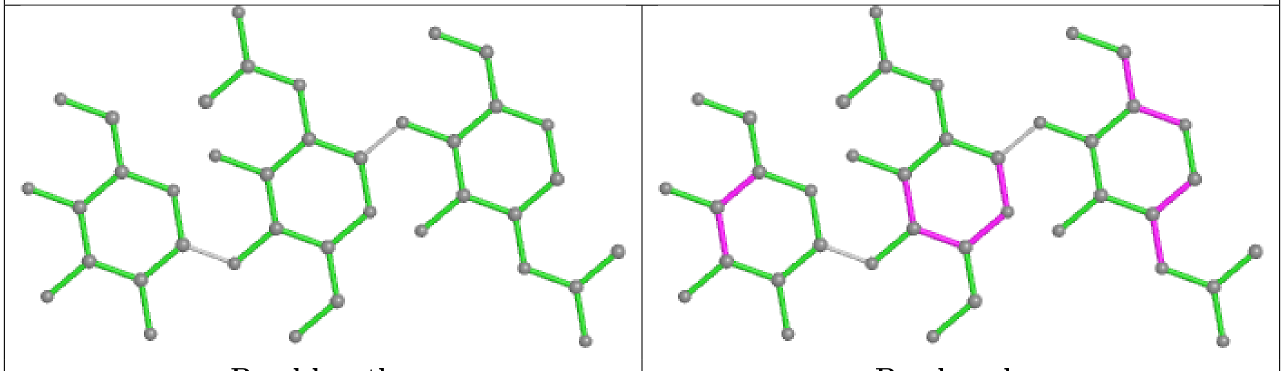


Bond lengths Bond angles

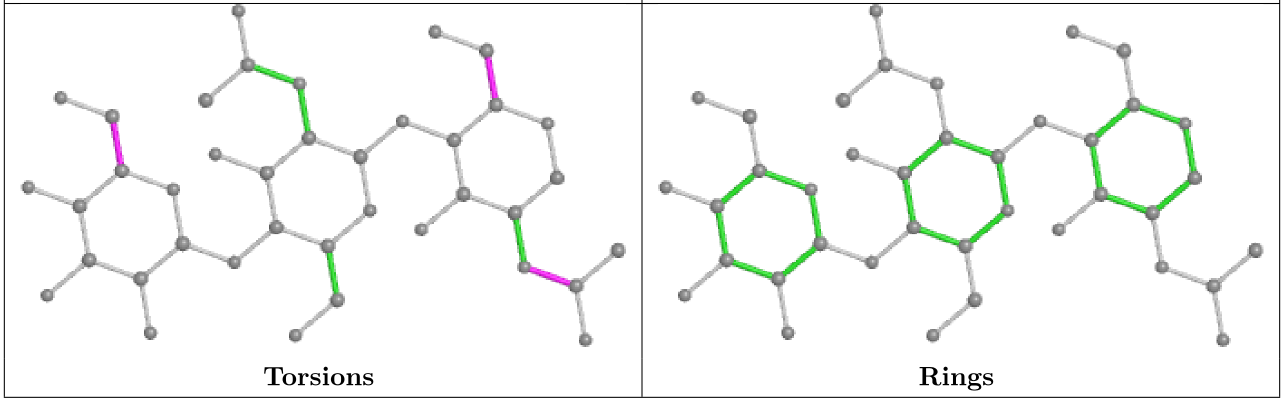


Torsions Rings

Oligosaccharide Chain I

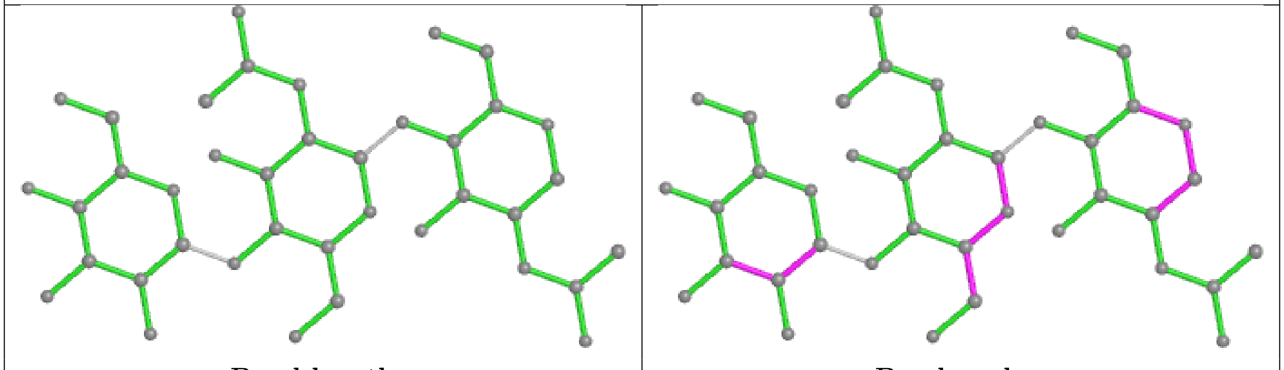


Bond lengths Bond angles



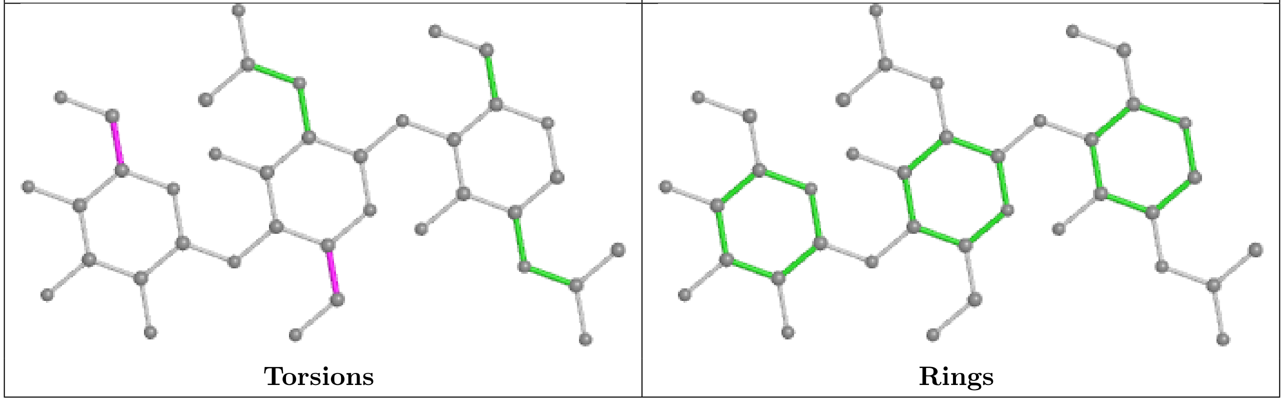
Torsions Rings

Oligosaccharide Chain J



Bond lengths

Bond angles

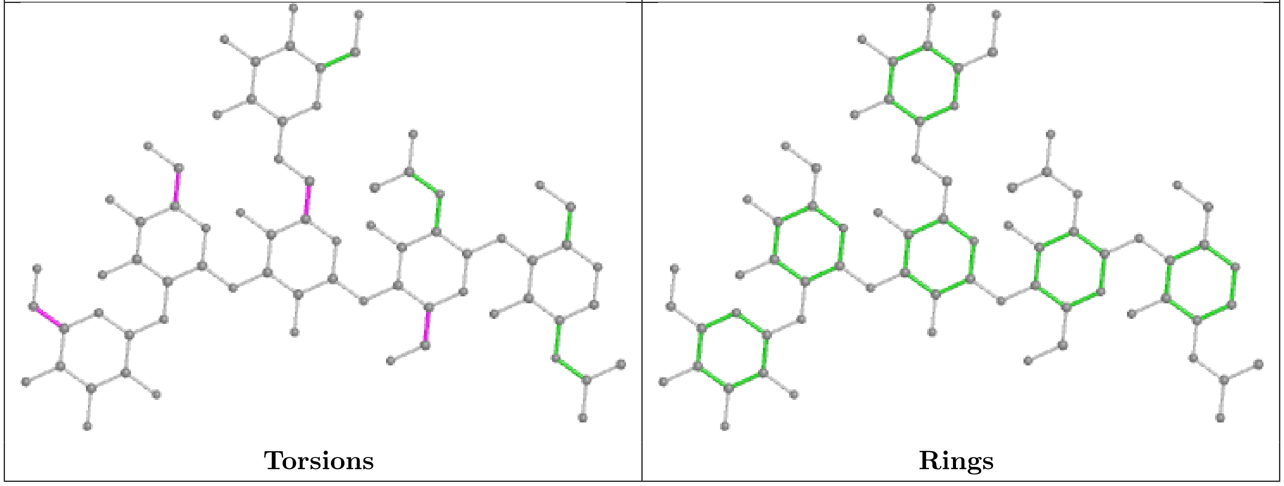


Torsions

Rings

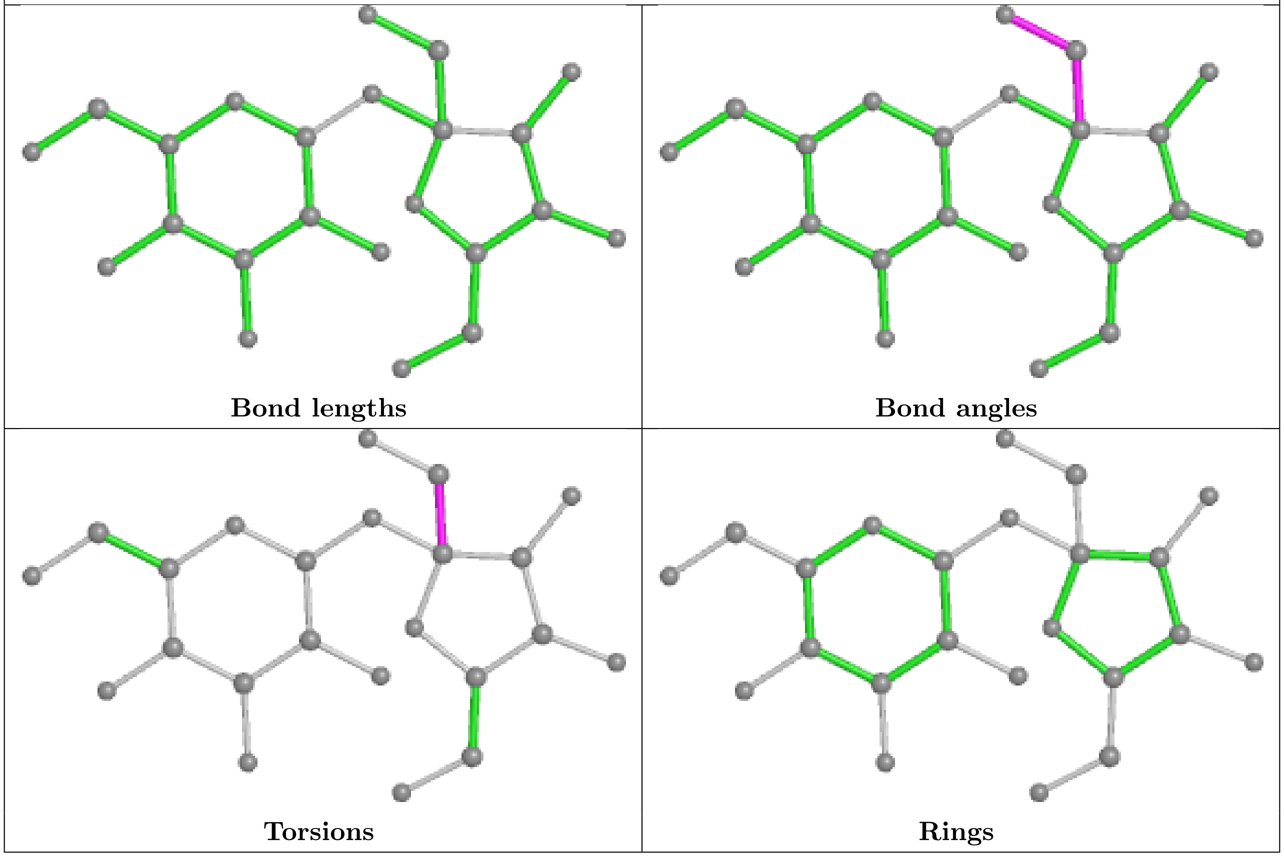
Oligosaccharide Chain G

Bond lengths **Bond angles**



Torsions **Rings**

Oligosaccharide Chain L



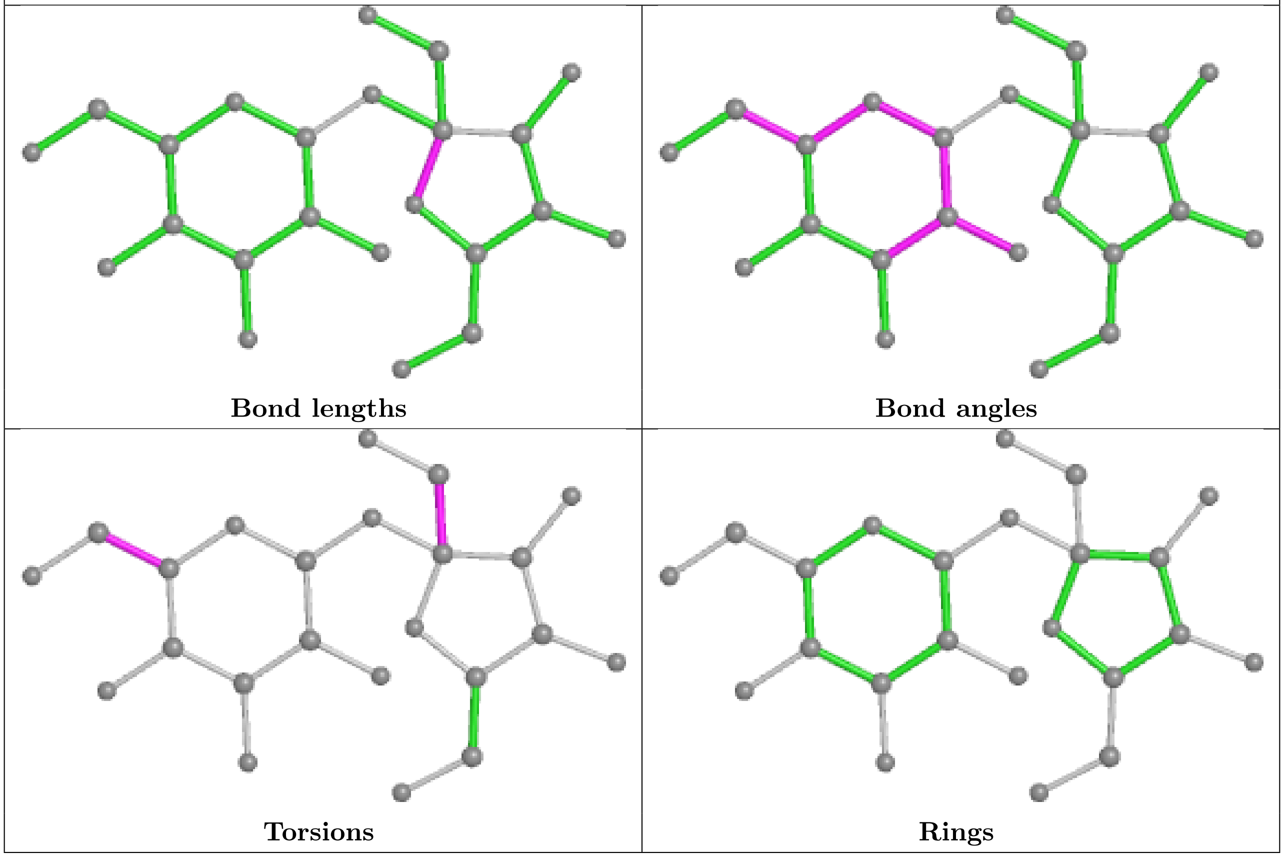
Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain M



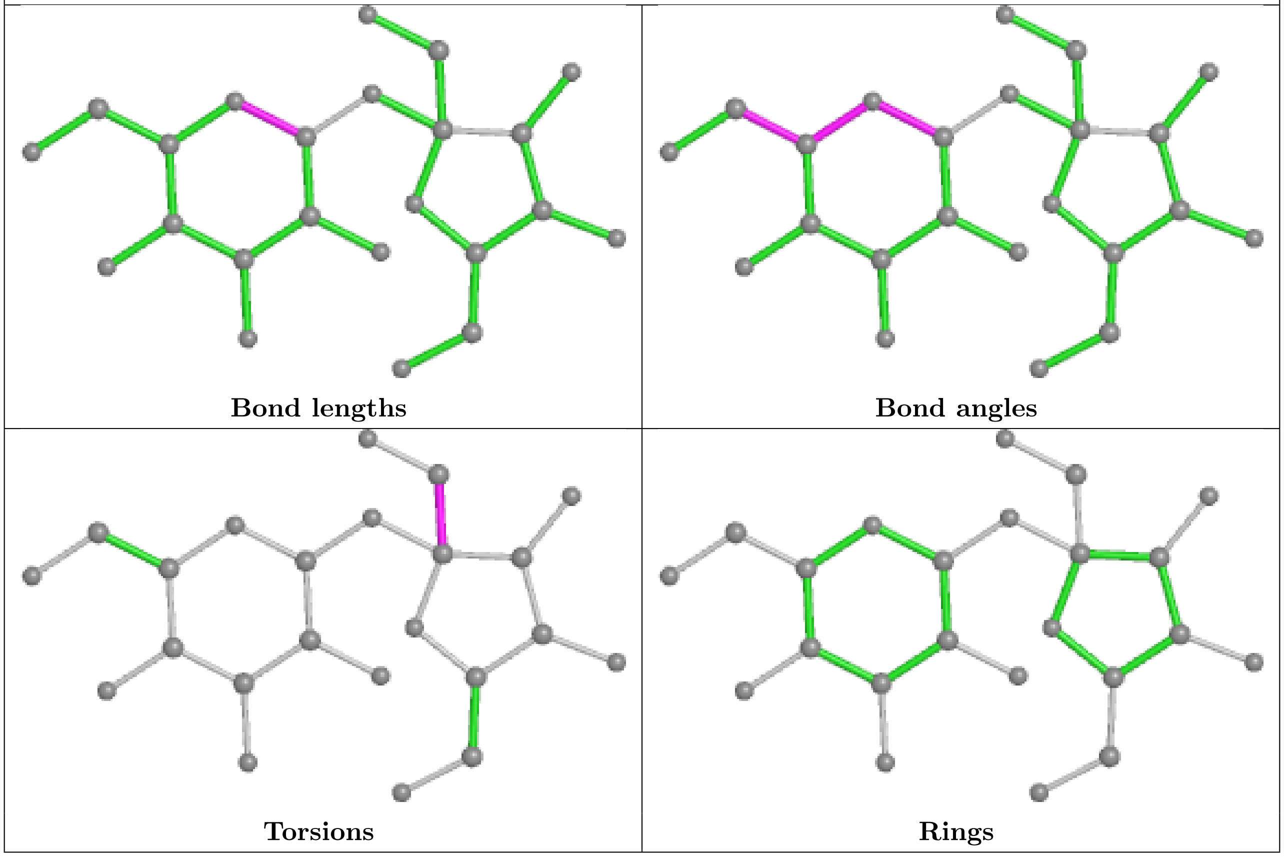
Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain N



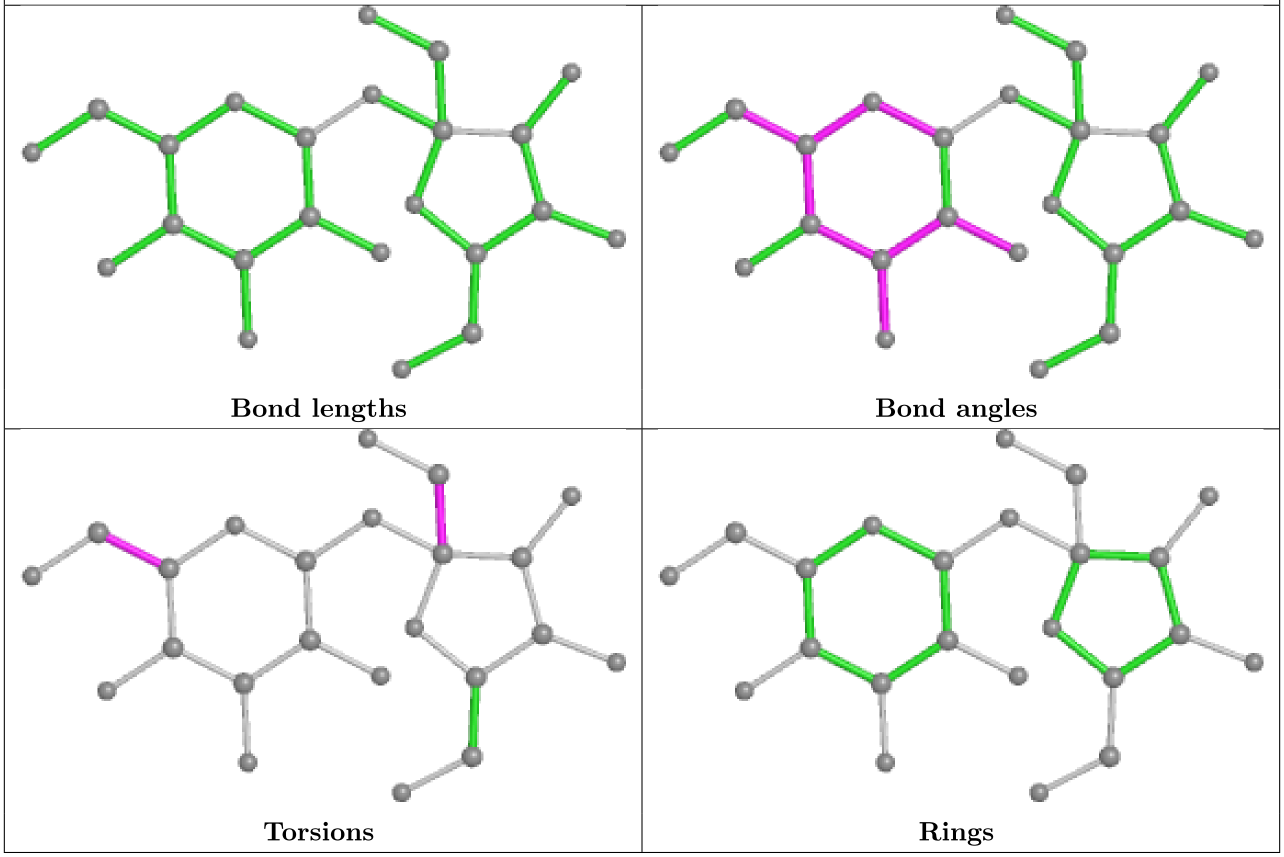
Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain O



Bond lengths

Bond angles

Torsions

Rings

5.6 Ligand geometry ⓘ

40 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	$\#\|Z\| > 2$
6	MAN	C	703	1	11,11,12	2.60	6 (54%)	15,15,17	1.77	3 (20%)
7	NAG	D	708	1	14,14,15	0.62	0	17,19,21	1.31	2 (11%)
6	MAN	B	702	1	11,11,12	2.28	4 (36%)	15,15,17	1.68	3 (20%)
7	NAG	A	705	1	14,14,15	0.52	0	17,19,21	1.07	1 (5%)
7	NAG	D	707	1	14,14,15	0.39	0	17,19,21	1.41	3 (17%)
7	NAG	C	705	1	14,14,15	0.64	0	17,19,21	1.18	2 (11%)
6	MAN	D	702	1	11,11,12	0.50	0	15,15,17	1.16	1 (6%)
6	MAN	C	702	1	11,11,12	0.52	0	15,15,17	1.91	4 (26%)
7	NAG	B	707	1	14,14,15	0.55	0	17,19,21	1.19	1 (5%)
6	MAN	C	704	1	11,11,12	0.61	0	15,15,17	0.74	0
7	NAG	A	706	1	14,14,15	0.41	0	17,19,21	1.68	2 (11%)
7	NAG	B	705	1	14,14,15	0.95	1 (7%)	17,19,21	1.43	3 (17%)
6	MAN	B	703	1	11,11,12	0.56	0	15,15,17	1.29	1 (6%)
6	MAN	A	702	1	11,11,12	0.75	0	15,15,17	1.30	3 (20%)
7	NAG	B	709	1	14,14,15	0.35	0	17,19,21	0.76	0
7	NAG	A	711	1	14,14,15	0.78	0	17,19,21	1.48	4 (23%)
7	NAG	C	706	1	14,14,15	0.37	0	17,19,21	1.23	1 (5%)
6	MAN	A	703	1	11,11,12	0.87	0	15,15,17	1.48	1 (6%)
7	NAG	A	709	1	14,14,15	0.74	0	17,19,21	1.19	2 (11%)
7	NAG	B	706	1	14,14,15	0.60	0	17,19,21	1.65	4 (23%)
6	MAN	D	701	1	11,11,12	0.74	0	15,15,17	1.43	2 (13%)
7	NAG	B	704	1	14,14,15	0.70	0	17,19,21	1.46	1 (5%)
7	NAG	C	708	1	14,14,15	0.48	0	17,19,21	0.98	0
7	NAG	C	710	1	14,14,15	0.48	0	17,19,21	1.50	3 (17%)
7	NAG	D	704	1	14,14,15	0.60	0	17,19,21	1.34	1 (5%)
7	NAG	C	707	1	14,14,15	0.73	1 (7%)	17,19,21	0.87	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	#|Z| > 2	Counts	RMSZ	#|Z| > 2
6	MAN	B	701	1	11,11,12	0.63	0	15,15,17	1.65	2 (13%)
7	NAG	A	708	1	14,14,15	0.90	1 (7%)	17,19,21	0.94	1 (5%)
7	NAG	B	710	1	14,14,15	0.64	0	17,19,21	1.33	2 (11%)
7	NAG	D	705	1	14,14,15	0.72	0	17,19,21	2.43	6 (35%)
7	NAG	B	708	1	14,14,15	0.83	0	17,19,21	0.97	0
7	NAG	D	706	1	14,14,15	0.48	0	17,19,21	1.09	2 (11%)
6	MAN	A	704	1	11,11,12	0.64	0	15,15,17	1.62	2 (13%)
6	MAN	D	703	1	11,11,12	0.95	0	15,15,17	1.38	2 (13%)
7	NAG	C	709	1	14,14,15	0.67	0	17,19,21	1.25	1 (5%)
7	NAG	A	707	1	14,14,15	0.67	1 (7%)	17,19,21	0.79	0
6	MAN	C	701	1	11,11,12	0.69	0	15,15,17	1.52	3 (20%)
6	MAN	A	701	1	11,11,12	0.63	0	15,15,17	1.02	0
7	NAG	A	710	1	14,14,15	0.64	0	17,19,21	1.12	0
7	NAG	D	709	1	14,14,15	0.52	0	17,19,21	0.94	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
6	MAN	C	703	1	-	1/2/19/22	0/1/1/1
7	NAG	D	708	1	-	2/6/23/26	0/1/1/1
6	MAN	B	702	1	-	2/2/19/22	0/1/1/1
7	NAG	A	705	1	-	2/6/23/26	0/1/1/1
7	NAG	D	707	1	-	2/6/23/26	0/1/1/1
7	NAG	C	705	1	-	4/6/23/26	0/1/1/1
6	MAN	D	702	1	-	1/2/19/22	0/1/1/1
6	MAN	C	702	1	-	2/2/19/22	0/1/1/1
7	NAG	B	707	1	-	2/6/23/26	0/1/1/1
6	MAN	C	704	1	-	0/2/19/22	0/1/1/1
7	NAG	A	706	1	-	2/6/23/26	0/1/1/1
7	NAG	B	705	1	-	4/6/23/26	0/1/1/1
6	MAN	B	703	1	-	2/2/19/22	0/1/1/1
6	MAN	A	702	1	-	2/2/19/22	0/1/1/1
7	NAG	B	709	1	-	0/6/23/26	0/1/1/1
7	NAG	A	711	1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	NAG	C	706	1	-	2/6/23/26	0/1/1/1
6	MAN	A	703	1	-	1/2/19/22	0/1/1/1
7	NAG	A	709	1	-	2/6/23/26	0/1/1/1
7	NAG	B	706	1	-	4/6/23/26	0/1/1/1
6	MAN	D	701	1	-	0/2/19/22	0/1/1/1
7	NAG	B	704	1	-	5/6/23/26	0/1/1/1
7	NAG	C	708	1	-	2/6/23/26	0/1/1/1
7	NAG	C	710	1	-	2/6/23/26	0/1/1/1
7	NAG	D	704	1	-	2/6/23/26	0/1/1/1
7	NAG	C	707	1	-	2/6/23/26	0/1/1/1
6	MAN	B	701	1	-	0/2/19/22	0/1/1/1
7	NAG	A	708	1	-	0/6/23/26	0/1/1/1
7	NAG	B	710	1	-	2/6/23/26	0/1/1/1
7	NAG	D	705	1	-	5/6/23/26	0/1/1/1
7	NAG	B	708	1	-	0/6/23/26	0/1/1/1
7	NAG	D	706	1	-	2/6/23/26	0/1/1/1
6	MAN	A	704	1	-	2/2/19/22	0/1/1/1
6	MAN	D	703	1	-	2/2/19/22	0/1/1/1
7	NAG	C	709	1	-	3/6/23/26	0/1/1/1
7	NAG	A	707	1	-	1/6/23/26	0/1/1/1
6	MAN	C	701	1	-	2/2/19/22	0/1/1/1
6	MAN	A	701	1	-	0/2/19/22	0/1/1/1
7	NAG	A	710	1	-	1/6/23/26	0/1/1/1
7	NAG	D	709	1	-	2/6/23/26	0/1/1/1

The worst 5 of 14 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	C	703	MAN	O3-C3	4.44	1.53	1.43
6	B	702	MAN	O4-C4	4.41	1.53	1.43
6	C	703	MAN	O2-C2	3.90	1.51	1.43
6	B	702	MAN	O5-C1	3.81	1.49	1.43
6	C	703	MAN	O5-C1	3.60	1.49	1.43

The worst 5 of 70 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	D	705	NAG	C2-N2-C7	6.66	132.39	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	706	NAG	O5-C1-C2	-5.33	102.87	111.29
6	C	702	MAN	O5-C1-C2	-5.19	102.77	110.77
7	B	704	NAG	C2-N2-C7	-4.89	115.93	122.90
6	A	704	MAN	O5-C1-C2	-4.72	103.49	110.77

There are no chirality outliers.

5 of 74 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	B	704	NAG	C3-C2-N2-C7
7	B	704	NAG	C8-C7-N2-C2
7	B	704	NAG	O7-C7-N2-C2
7	B	705	NAG	C3-C2-N2-C7
7	B	705	NAG	C8-C7-N2-C2

There are no ring outliers.

12 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	C	705	NAG	1	0
6	D	702	MAN	1	0
6	C	704	MAN	1	0
7	A	706	NAG	1	0
7	A	709	NAG	1	0
7	B	706	NAG	2	0
6	D	701	MAN	1	0
7	D	704	NAG	1	0
7	B	710	NAG	1	0
7	D	705	NAG	2	0
6	C	701	MAN	1	0
7	D	709	NAG	1	0

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	525/675 (77%)	0.06	13 (2%)	57	59	22, 32, 56, 87	0
1	B	526/675 (77%)	0.28	33 (6%)	20	21	22, 35, 72, 105	0
1	C	528/675 (78%)	0.16	16 (3%)	50	53	22, 36, 62, 109	0
1	D	524/675 (77%)	0.46	44 (8%)	11	12	22, 38, 78, 103	0
All	All	2103/2700 (77%)	0.24	106 (5%)	28	31	22, 35, 68, 109	0

The worst 5 of 106 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	555	CYS	6.2
1	C	143	CYS	4.9
1	D	554	LYS	4.8
1	B	513	THR	4.7
1	D	436	ASN	4.2

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
4	MAN	G	5	11/12	0.68	0.40	92,102,108,109	0
3	BMA	I	3	11/12	0.69	0.36	89,103,108,111	0

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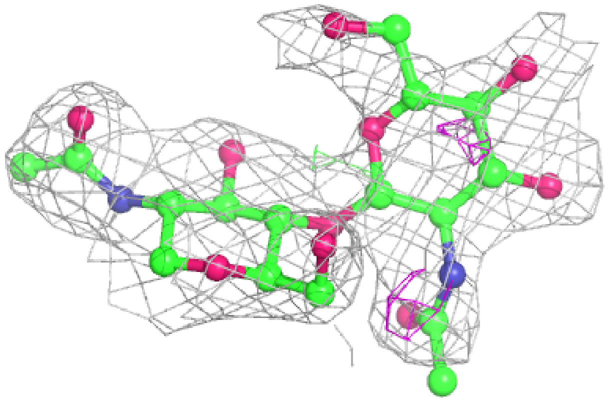
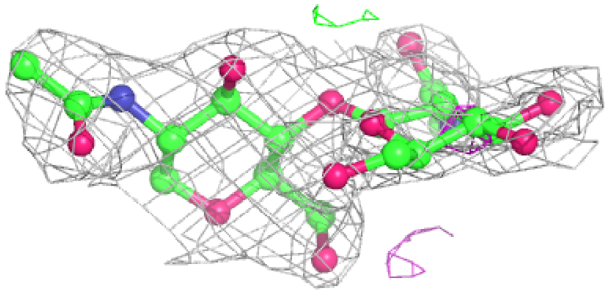
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\AA^2$)	Q<0.9
4	MAN	G	6	11/12	0.71	0.25	27,28,29,31	11
3	BMA	F	3	11/12	0.78	0.29	70,74,80,81	0
2	NAG	H	2	14/15	0.78	0.28	71,78,91,91	0
3	NAG	I	2	14/15	0.80	0.25	69,89,95,97	0
3	NAG	J	2	14/15	0.80	0.21	59,72,80,85	0
2	NAG	E	2	14/15	0.82	0.39	60,67,72,73	0
3	BMA	J	3	11/12	0.82	0.28	85,95,101,101	0
2	NAG	K	2	14/15	0.83	0.32	79,92,102,103	0
3	NAG	F	2	14/15	0.85	0.31	54,67,79,80	0
4	BMA	G	3	11/12	0.86	0.16	66,78,82,84	0
3	NAG	I	1	14/15	0.91	0.20	56,68,72,77	0
4	NAG	G	2	14/15	0.92	0.24	51,61,72,75	0
4	MAN	G	4	11/12	0.92	0.23	84,86,90,96	0
2	NAG	E	1	14/15	0.94	0.14	42,47,50,57	0
5	GLC	N	1	11/12	0.94	0.13	44,48,52,53	0
5	GLC	O	1	11/12	0.94	0.12	34,37,41,43	0
4	NAG	G	1	14/15	0.95	0.18	46,51,55,56	0
3	NAG	J	1	14/15	0.95	0.15	34,38,41,51	0
3	NAG	F	1	14/15	0.95	0.13	35,40,45,47	0
2	NAG	K	1	14/15	0.95	0.11	44,48,60,66	0
2	NAG	H	1	14/15	0.96	0.14	37,42,46,57	0
5	FRU	M	2	12/12	0.96	0.17	27,30,34,44	0
5	GLC	M	1	11/12	0.97	0.12	35,37,41,43	0
5	FRU	N	2	12/12	0.97	0.15	29,35,40,44	0
5	FRU	L	2	12/12	0.97	0.11	28,32,34,37	0
5	FRU	O	2	12/12	0.97	0.12	26,29,33,38	0
5	GLC	L	1	11/12	0.98	0.11	37,39,41,44	0

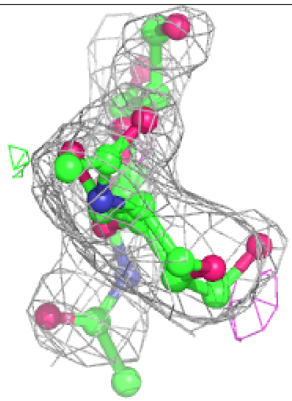
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain E:

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

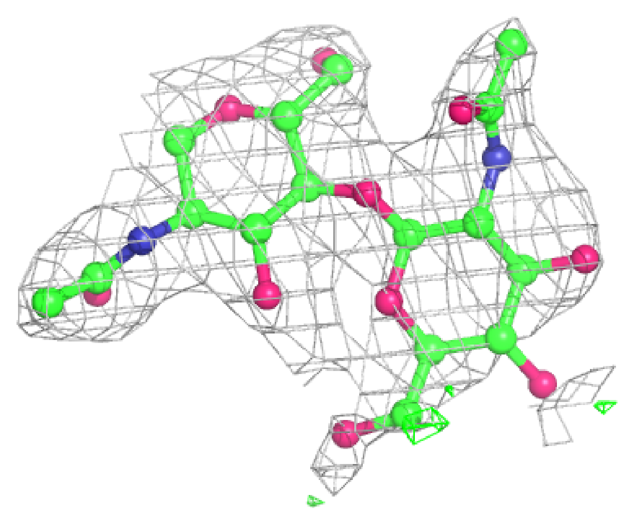


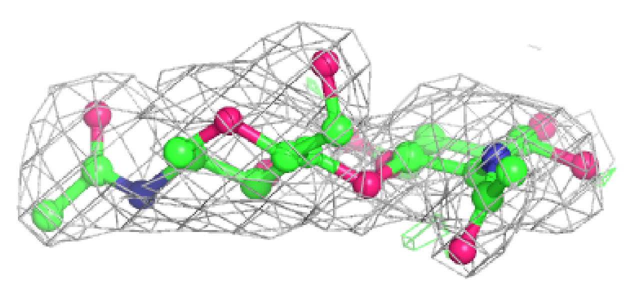


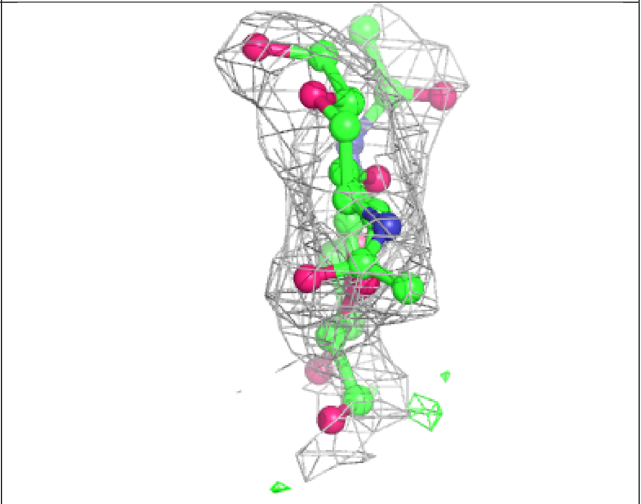


Electron density around Chain H:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
and green (positive)

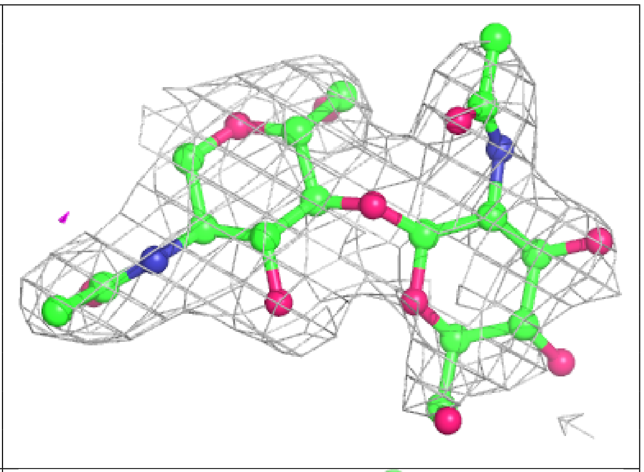


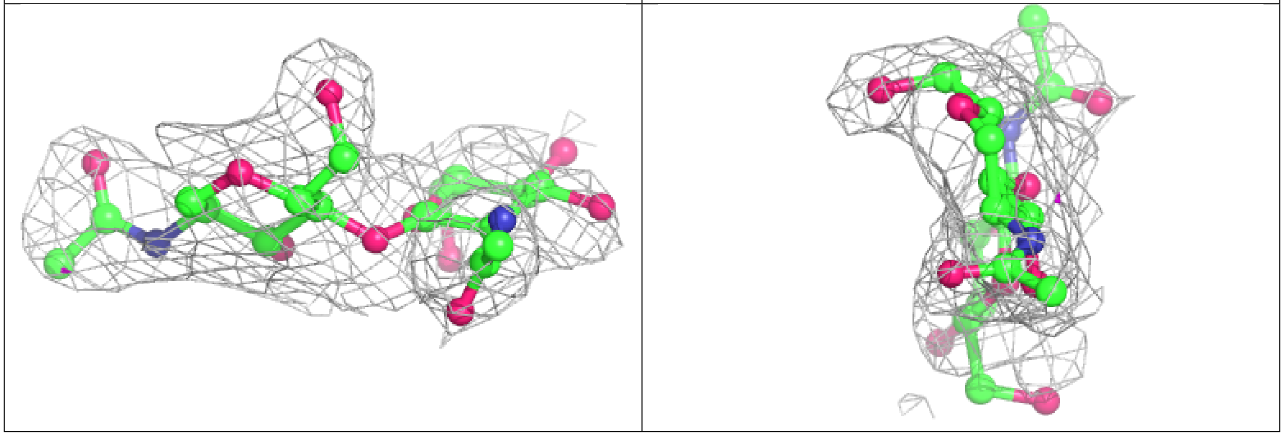




Electron density around Chain K:

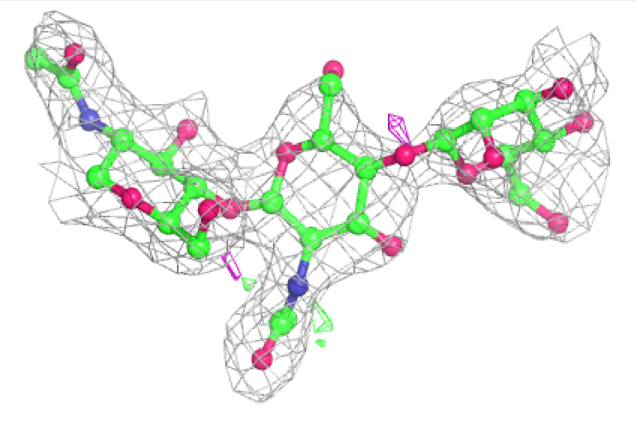
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

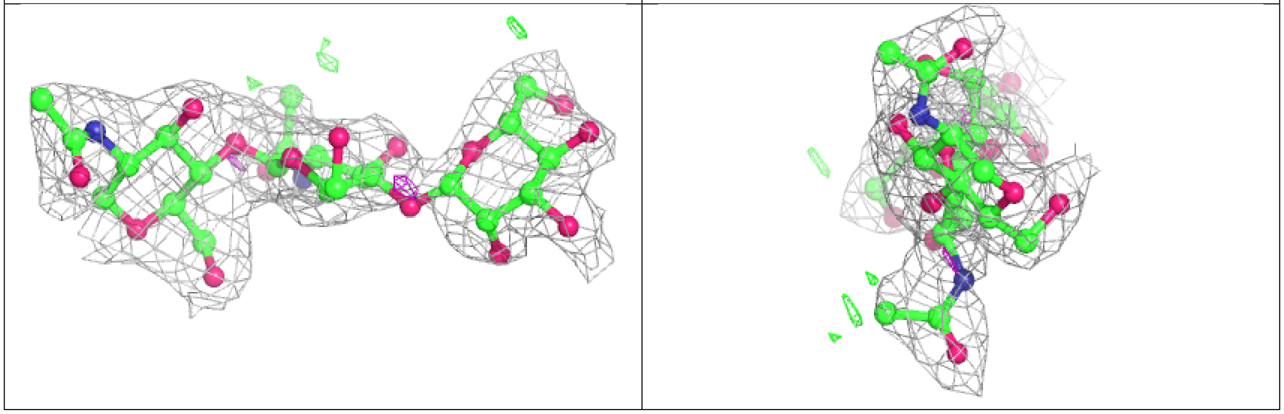




Electron density around Chain F:

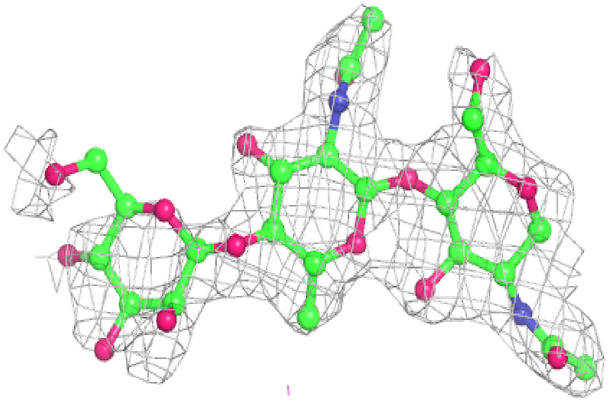
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

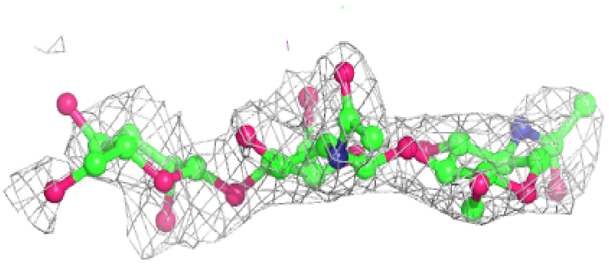


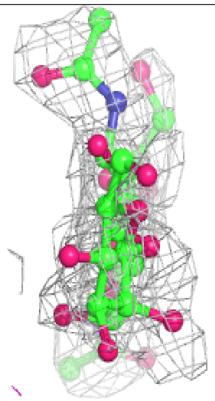


Electron density around Chain I:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
and green (positive)

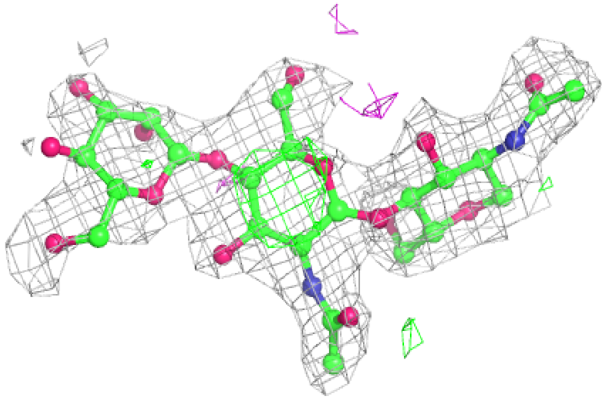


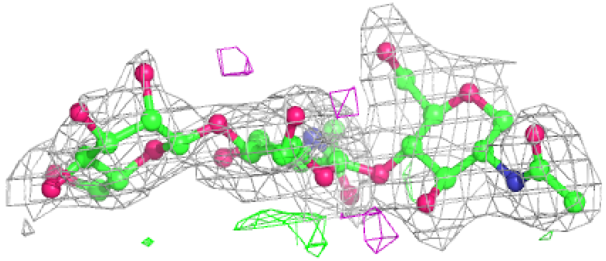


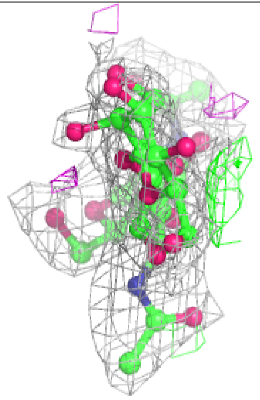


Electron density around Chain J:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
and green (positive)

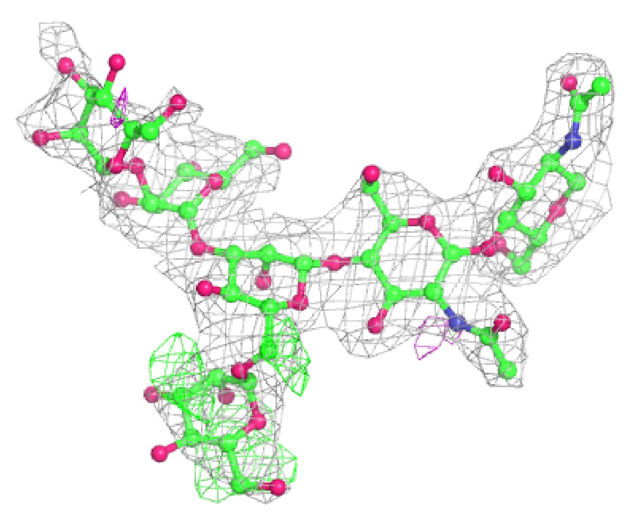


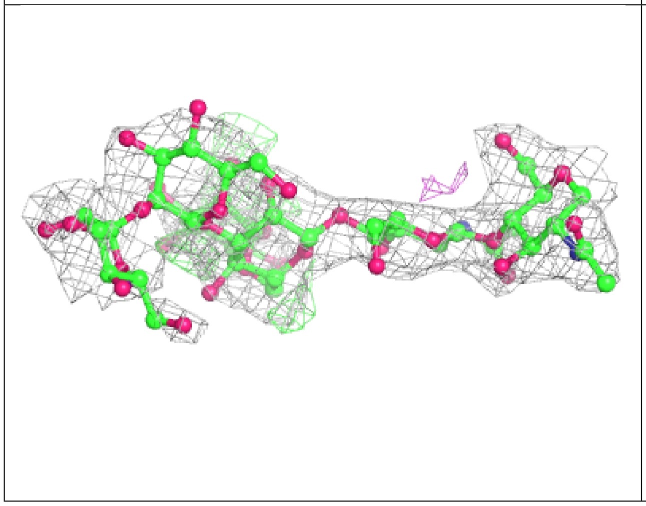


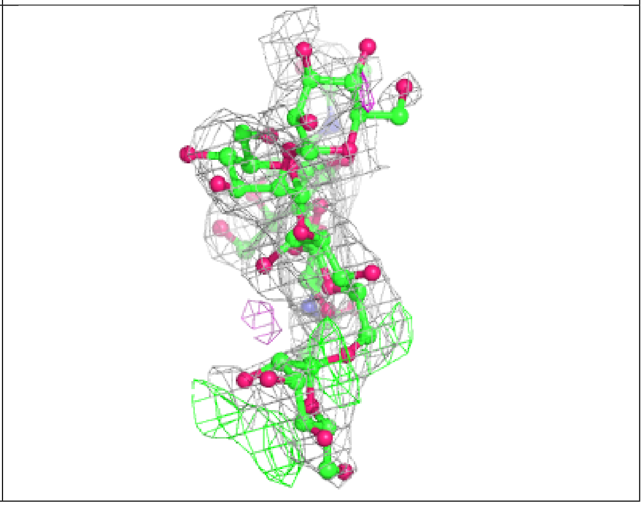


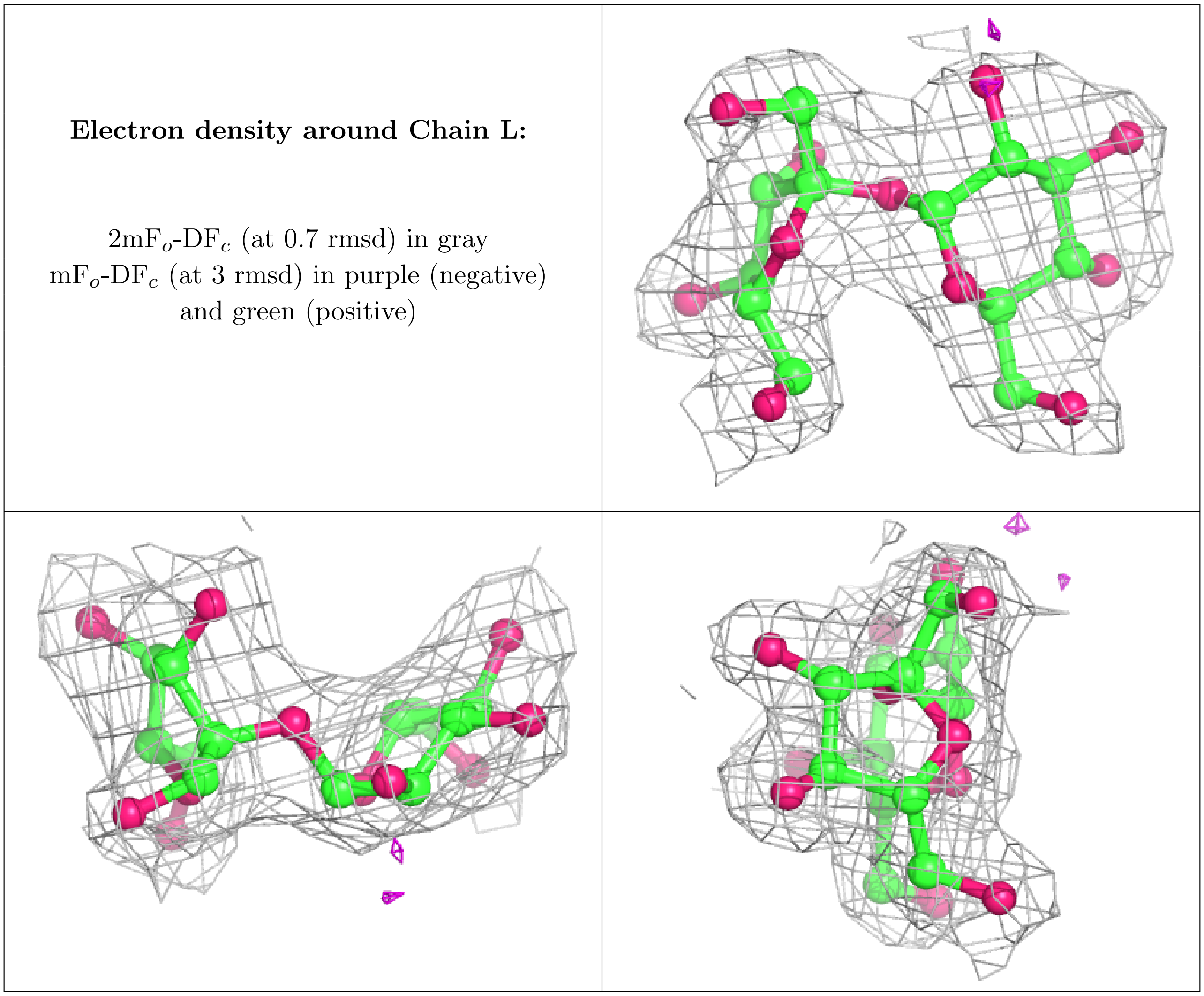
Electron density around Chain G:

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







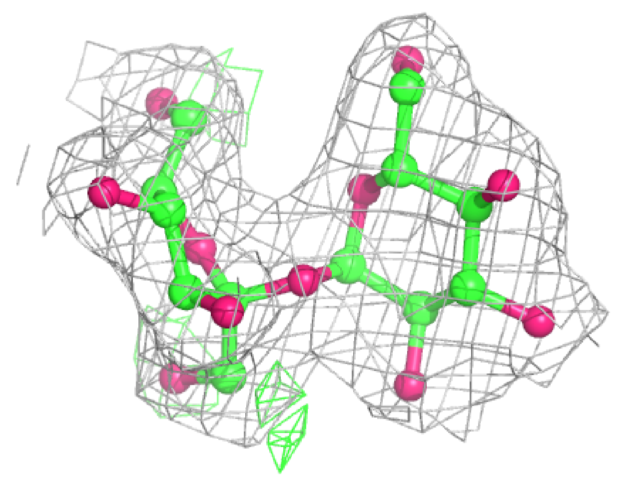


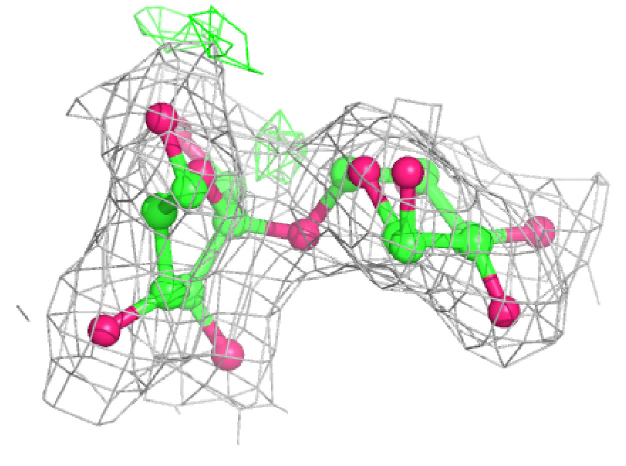
Electron density around Chain L:

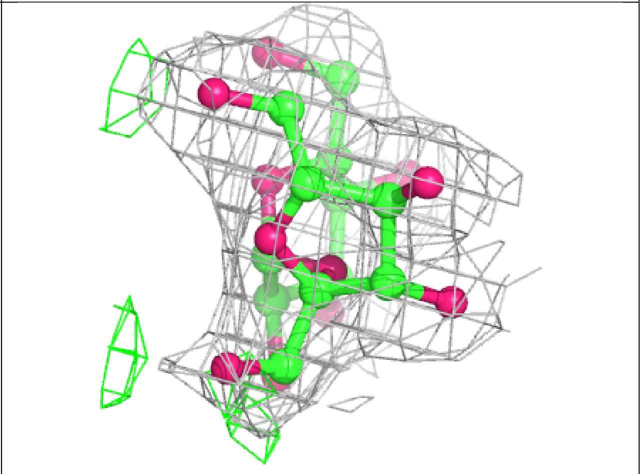
2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
and green (positive)

Electron density around Chain M:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
and green (positive)

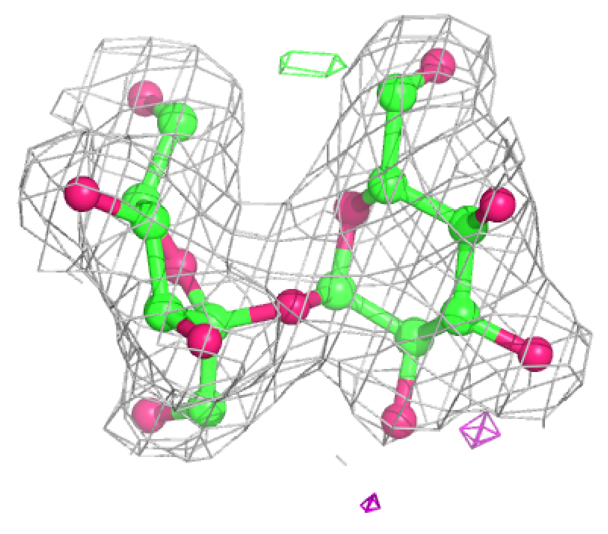


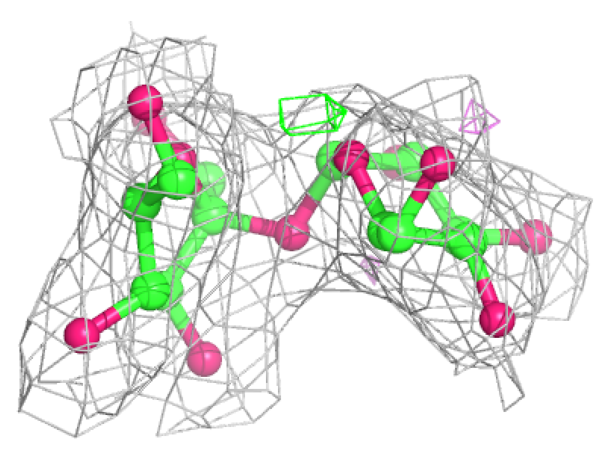


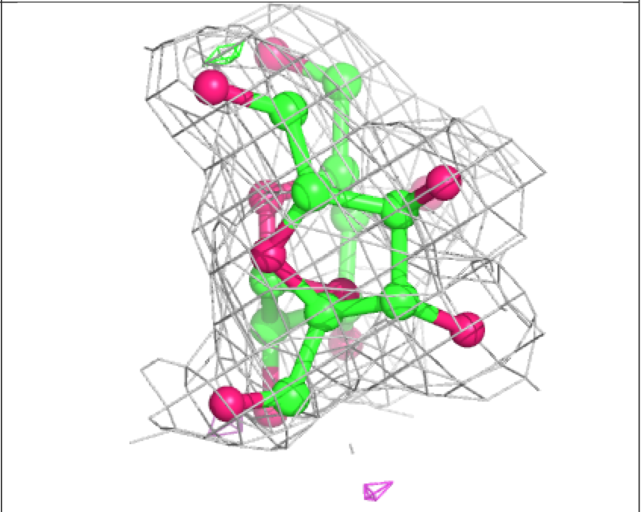


Electron density around Chain N:

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

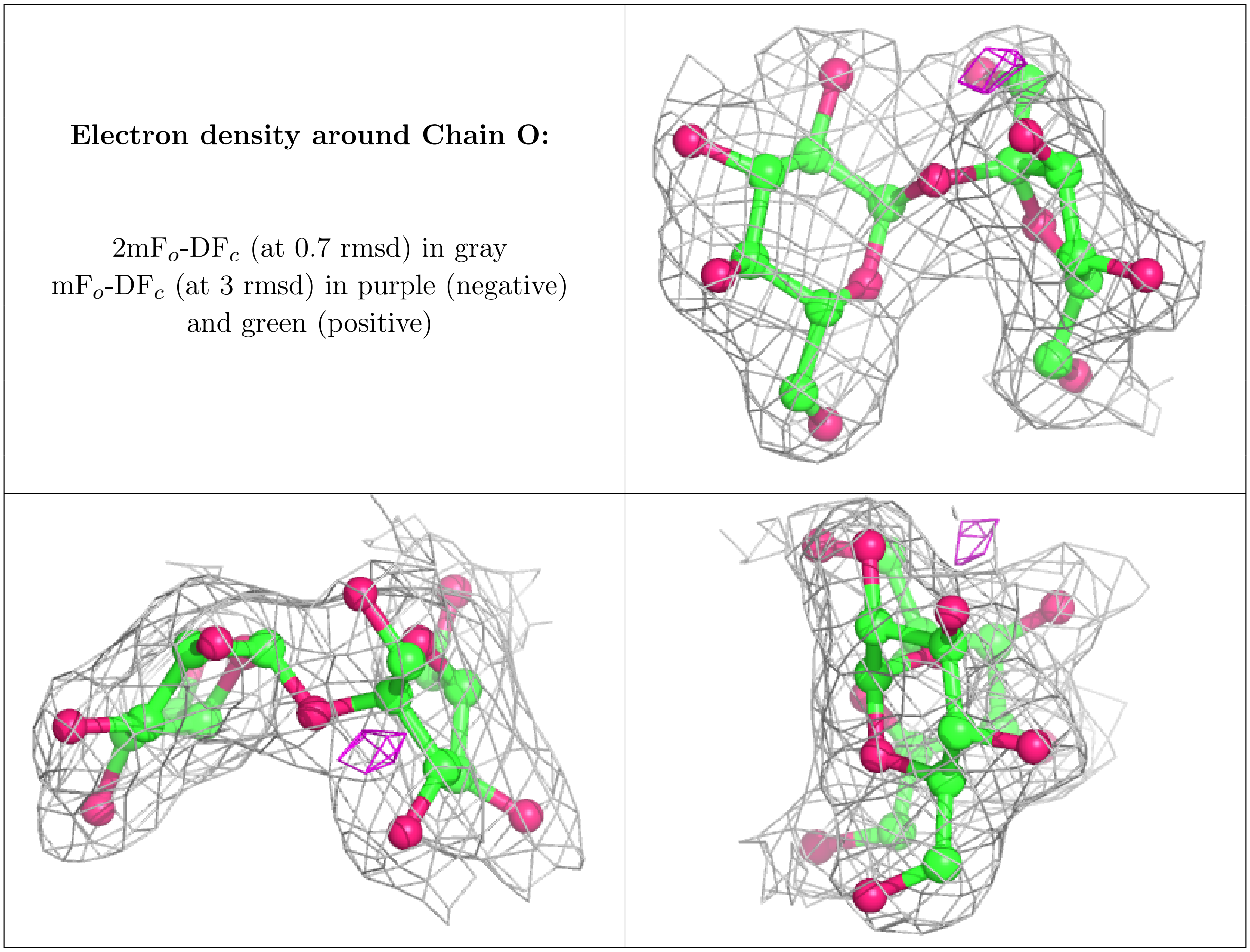






Electron density around Chain O:

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
6	MAN	D	701	11/12	0.69	0.43	88,106,121,121	0
6	MAN	C	703	11/12	0.71	0.43	27,29,31,32	0
6	MAN	B	702	11/12	0.75	0.43	27,29,31,31	0
6	MAN	A	702	11/12	0.76	0.31	82,90,95,96	0
7	NAG	B	705	14/15	0.76	0.45	72,80,83,84	0
6	MAN	D	703	11/12	0.77	0.34	74,79,84,86	0
7	NAG	B	704	14/15	0.80	0.31	59,67,82,84	0
7	NAG	D	708	14/15	0.80	0.35	96,112,121,129	0
6	MAN	C	702	11/12	0.81	0.26	76,85,88,89	0
7	NAG	A	706	14/15	0.82	0.33	66,74,77,78	0
6	MAN	A	704	11/12	0.84	0.25	64,71,78,78	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
7	NAG	D	705	14/15	0.84	0.19	57,63,69,69	0
6	MAN	C	704	11/12	0.84	0.21	56,64,69,71	0
6	MAN	B	703	11/12	0.85	0.21	77,80,89,90	0
7	NAG	B	706	14/15	0.85	0.34	63,74,84,85	0
6	MAN	C	701	11/12	0.86	0.21	71,79,84,85	0
6	MAN	A	703	11/12	0.86	0.23	63,68,75,76	0
7	NAG	D	706	14/15	0.86	0.49	59,70,77,78	0
6	MAN	B	701	11/12	0.86	0.25	79,86,89,89	0
7	NAG	D	709	14/15	0.86	0.22	69,72,78,85	0
7	NAG	A	705	14/15	0.87	0.25	58,65,79,80	0
7	NAG	B	710	14/15	0.88	0.26	66,75,80,80	0
7	NAG	C	709	14/15	0.88	0.18	54,63,71,76	0
6	MAN	A	701	11/12	0.88	0.34	72,79,82,83	0
7	NAG	C	705	14/15	0.89	0.26	62,71,81,84	0
6	MAN	D	702	11/12	0.90	0.45	81,92,99,110	0
7	NAG	B	709	14/15	0.91	0.21	64,71,80,82	0
7	NAG	A	710	14/15	0.91	0.15	51,65,70,75	0
7	NAG	A	707	14/15	0.92	0.15	46,51,55,56	0
7	NAG	D	704	14/15	0.92	0.14	55,59,66,67	0
7	NAG	C	707	14/15	0.93	0.22	51,58,61,65	0
7	NAG	A	711	14/15	0.93	0.16	46,56,64,67	0
7	NAG	C	710	14/15	0.93	0.15	50,52,62,68	0
7	NAG	C	706	14/15	0.93	0.15	55,63,72,72	0
7	NAG	A	709	14/15	0.94	0.16	46,53,56,60	0
7	NAG	A	708	14/15	0.94	0.13	37,40,46,47	0
7	NAG	D	707	14/15	0.95	0.16	43,48,56,58	0
7	NAG	C	708	14/15	0.95	0.17	48,52,57,58	0
7	NAG	B	708	14/15	0.95	0.14	40,43,49,50	0
7	NAG	B	707	14/15	0.97	0.10	30,33,41,42	0

6.5 Other polymers ⓘ

There are no such residues in this entry.