wwPDB X-ray Structure Validation Summary Report (i)

Oct 5, 2023 – 08:12 PM EDT

PDB ID	:	8F8W
Title	:	Crystal structure of Nb.X0 bound to the afucosylated human IgG1 fragment crystal form I
Authors	:	Goldgur, Y.; Ravetch, J.; Gupta, A.; Kao, K.; Oren, D.
Deposited on	:	2022-11-22
Resolution	:	2.71 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) 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 (i)) were used in the production of this report:

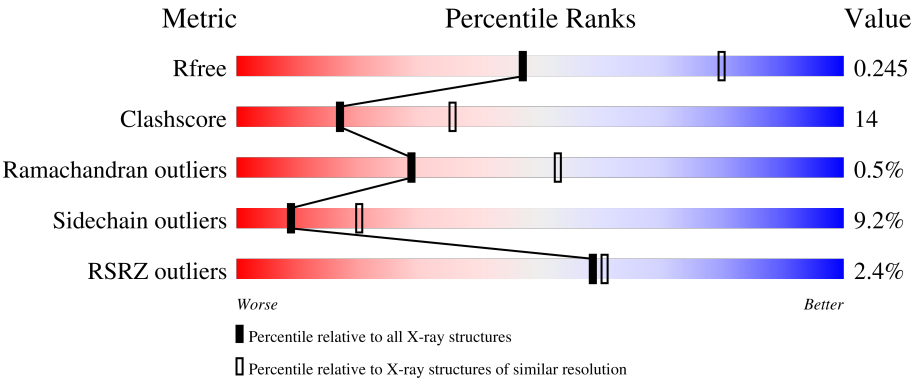
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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 2.71 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	3359 (2.74-2.70)
Clashscore	141614	3686 (2.74-2.70)
Ramachandran outliers	138981	3622 (2.74-2.70)
Sidechain outliers	138945	3623 (2.74-2.70)
RSRZ outliers	127900	3276 (2.74-2.70)

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	224	68% 20% • • 8%
1	B	224	2% 67% 21% • • 8%
1	E	224	% 61% 28% • 8%
1	F	224	3% 55% 29% 6% • 7%
2	C	120	% 67% 29% • •

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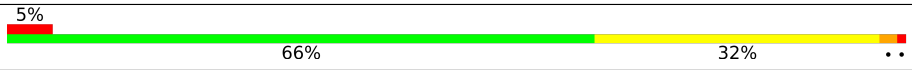
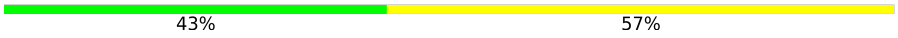
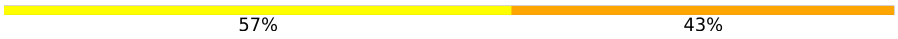
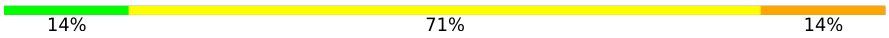
Mol	Chain	Length	Quality of chain
2	D	120	
2	G	120	
2	H	120	
3	I	7	
3	J	7	
3	K	7	
3	L	7	

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 10703 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 afucosylated IgG1 fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	207	Total	C	N	O	S	0	0	0
			1661	1057	282	315	7			
1	B	207	Total	C	N	O	S	0	0	0
			1661	1057	282	315	7			
1	E	206	Total	C	N	O	S	0	0	0
			1657	1055	281	314	7			
1	F	208	Total	C	N	O	S	0	0	0
			1657	1055	283	312	7			

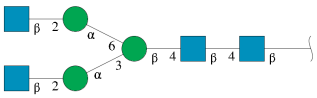
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	382	ARG	GLU	conflict	UNP Q6MZV7
B	382	ARG	GLU	conflict	UNP Q6MZV7
E	382	ARG	GLU	conflict	UNP Q6MZV7
F	382	ARG	GLU	conflict	UNP Q6MZV7

- Molecule 2 is a protein called Nb.X0.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	120	Total	C	N	O	S	0	0	0
			923	578	162	179	4			
2	D	120	Total	C	N	O	S	0	0	0
			923	578	162	179	4			
2	G	120	Total	C	N	O	S	0	0	0
			923	578	162	179	4			
2	H	120	Total	C	N	O	S	0	0	0
			923	578	162	179	4			

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-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
3	I	7	Total C N O 89 50 4 35	0	0	0
3	J	7	Total C N O 89 50 4 35	0	0	0
3	K	7	Total C N O 89 50 4 35	0	0	0
3	L	7	Total C N O 89 50 4 35	0	0	0

- Molecule 4 is water.

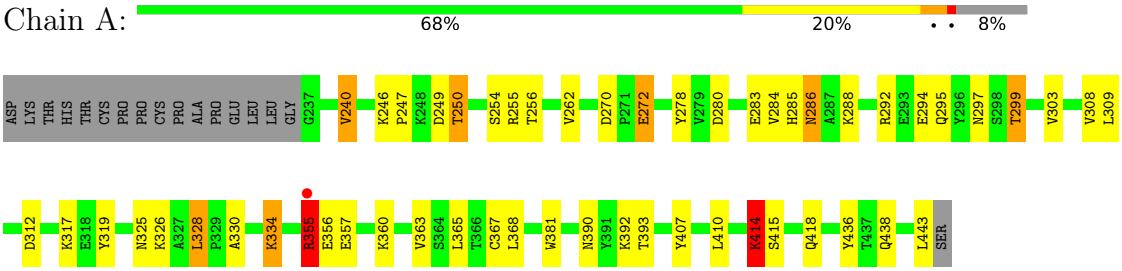
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	7	Total O 7 7	0	0
4	B	5	Total O 5 5	0	0
4	C	1	Total O 1 1	0	0
4	E	5	Total O 5 5	0	0
4	F	1	Total O 1 1	0	0

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.

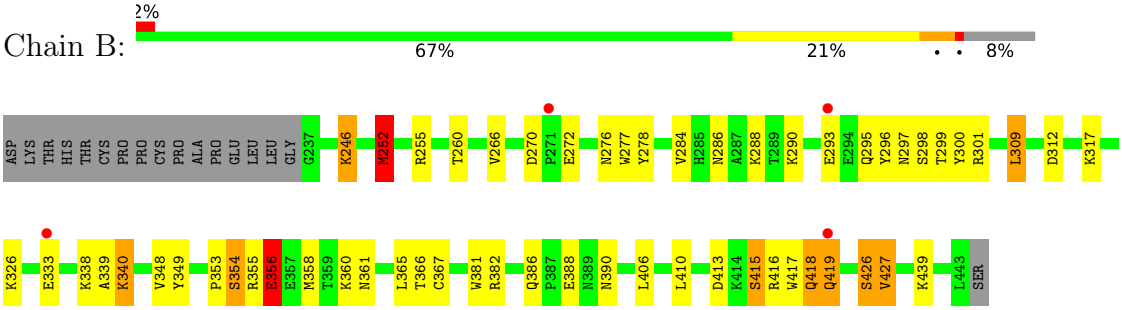
• Molecule 1: afucosylated IgG1 fragment

Chain A:



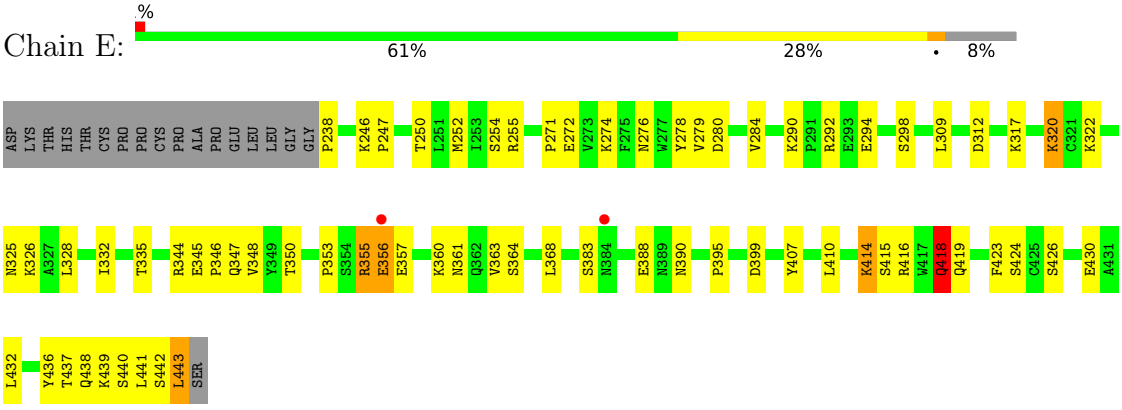
• Molecule 1: afucosylated IgG1 fragment

Chain B:



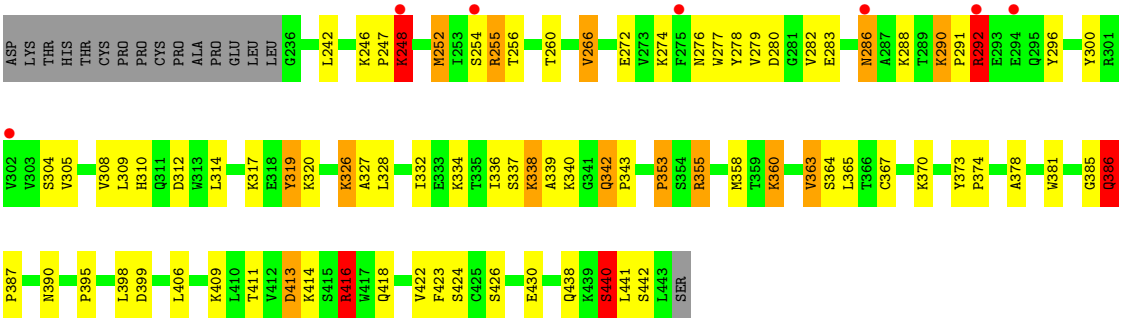
• Molecule 1: afucosylated IgG1 fragment

Chain E:



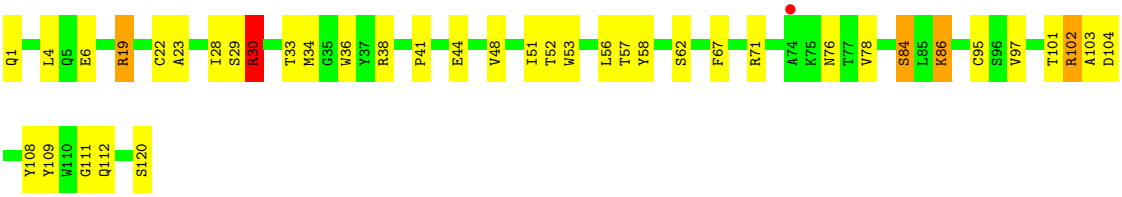
• Molecule 1: afucosylated IgG1 fragment

Chain F:



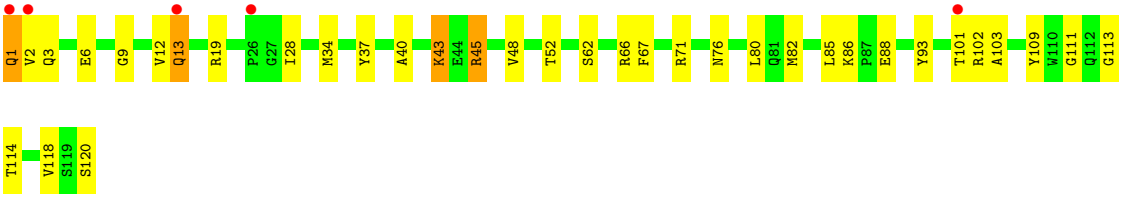
- Molecule 2: Nb.X0

Chain C:



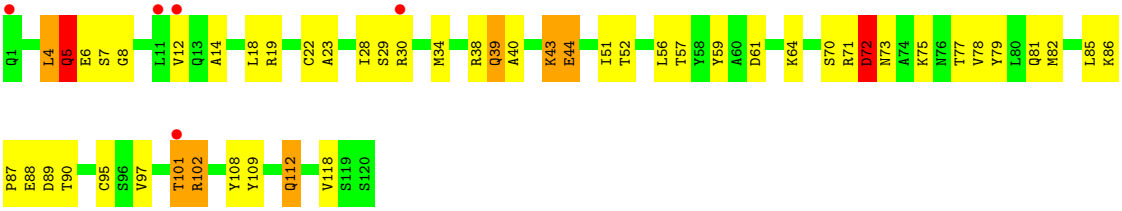
- Molecule 2: Nb.X0

Chain D:



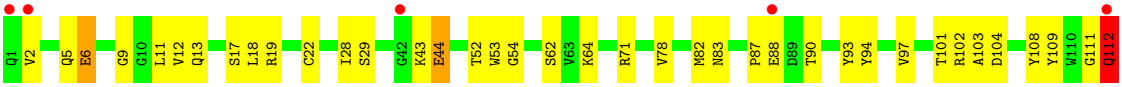
- Molecule 2: Nb.X0

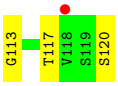
Chain G:



- Molecule 2: Nb.X0

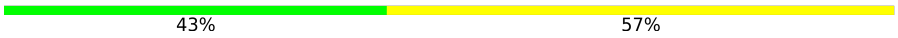
Chain H:





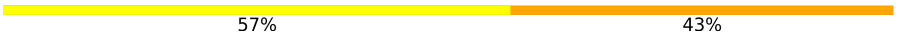
• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:



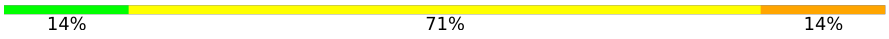
• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:



• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-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 K:



• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-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 L:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α, β, γ	170.64Å 170.64Å 126.42Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.90 – 2.71 45.90 – 2.71	Depositor EDS
% Data completeness (in resolution range)	99.7 (45.90-2.71) 99.7 (45.90-2.71)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.58 (at 2.73Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.208 , 0.241 0.211 , 0.245	Depositor DCC
R_{free} test set	2013 reflections (3.55%)	wwPDB-VP
Wilson B-factor (Å^2)	76.1	Xtriage
Anisotropy	0.474	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.29 , 51.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.037 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	10703	wwPDB-VP
Average B, all atoms (Å^2)	93.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.01% of the height of the origin peak. No significant pseudotranslation is detected.*

[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: MAN, NAG, 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	0.63	3/1707 (0.2%)	0.83	3/2324 (0.1%)
1	B	0.73	3/1707 (0.2%)	1.02	10/2324 (0.4%)
1	E	0.81	7/1703 (0.4%)	1.00	16/2318 (0.7%)
1	F	0.75	6/1703 (0.4%)	1.38	27/2319 (1.2%)
2	C	0.58	0/945	1.01	6/1280 (0.5%)
2	D	0.62	2/945 (0.2%)	1.06	7/1280 (0.5%)
2	G	0.89	6/945 (0.6%)	1.30	16/1280 (1.2%)
2	H	0.59	0/945	1.55	5/1280 (0.4%)
All	All	0.72	27/10600 (0.3%)	1.14	90/14405 (0.6%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	E	0	3
1	F	0	6
2	C	0	1
2	D	0	1
2	G	0	4
2	H	0	1
All	All	0	17

The worst 5 of 27 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	356	GLU	CD-OE2	14.44	1.41	1.25
1	E	272	GLU	CD-OE2	11.29	1.38	1.25
1	B	356	GLU	CG-CD	10.00	1.67	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	356	GLU	CD-OE2	9.76	1.36	1.25
2	G	5	GLN	CG-CD	-9.55	1.29	1.51

The worst 5 of 90 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	H	44	GLU	OE1-CD-OE2	-37.08	78.80	123.30
2	H	44	GLU	CG-CD-OE1	20.90	160.09	118.30
2	H	44	GLU	CG-CD-OE2	-18.40	81.50	118.30
1	F	416	ARG	CG-CD-NE	-16.84	76.44	111.80
1	F	292	ARG	CA-CB-CG	15.83	148.23	113.40

There are no chirality outliers.

5 of 17 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	356	GLU	Sidechain
2	C	30	ARG	Sidechain
2	D	45	ARG	Sidechain
1	E	271	PRO	Peptide
1	E	355	ARG	Peptide

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	1661	0	1632	31	1
1	B	1661	0	1632	38	1
1	E	1657	0	1630	40	1
1	F	1657	0	1627	68	1
2	C	923	0	884	26	0
2	D	923	0	884	22	0
2	G	923	0	884	51	0
2	H	923	0	884	28	0
3	I	89	0	76	1	0
3	J	89	0	76	4	0
3	K	89	0	76	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	L	89	0	76	0	0
4	A	7	0	0	1	0
4	B	5	0	0	0	0
4	C	1	0	0	0	0
4	E	5	0	0	1	0
4	F	1	0	0	0	0
All	All	10703	0	10361	295	2

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 14.

The worst 5 of 295 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:F:360:LYS:CE	1:F:360:LYS:NZ	1.69	1.56
1:F:290:LYS:CE	1:F:290:LYS:NZ	1.69	1.52
2:G:5:GLN:HE22	2:G:23:ALA:N	1.51	1.06
1:F:360:LYS:NZ	1:F:360:LYS:CD	2.19	1.05
2:G:57:THR:HG1	2:G:59:TYR:HE1	1.08	1.01

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:436:TYR:OH	1:B:426:SER:OG[5_554]	1.93	0.27
1:E:436:TYR:OH	1:F:426:SER:OG[5_554]	1.99	0.21

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	205/224 (92%)	200 (98%)	5 (2%)	0	100	100
1	B	205/224 (92%)	199 (97%)	6 (3%)	0	100	100
1	E	204/224 (91%)	200 (98%)	4 (2%)	0	100	100
1	F	206/224 (92%)	198 (96%)	7 (3%)	1 (0%)	29	53
2	C	118/120 (98%)	114 (97%)	3 (2%)	1 (1%)	19	41
2	D	118/120 (98%)	114 (97%)	3 (2%)	1 (1%)	19	41
2	G	118/120 (98%)	114 (97%)	2 (2%)	2 (2%)	9	21
2	H	118/120 (98%)	117 (99%)	0	1 (1%)	19	41
All	All	1292/1376 (94%)	1256 (97%)	30 (2%)	6 (0%)	29	53

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	102	ARG
2	D	102	ARG
2	G	5	GLN
2	H	102	ARG
1	F	292	ARG

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	193/208 (93%)	178 (92%)	15 (8%)	12	28
1	B	193/208 (93%)	170 (88%)	23 (12%)	5	11
1	E	193/208 (93%)	181 (94%)	12 (6%)	18	39
1	F	191/208 (92%)	170 (89%)	21 (11%)	6	14
2	C	95/95 (100%)	85 (90%)	10 (10%)	7	15
2	D	95/95 (100%)	89 (94%)	6 (6%)	18	38
2	G	95/95 (100%)	87 (92%)	8 (8%)	11	24
2	H	95/95 (100%)	84 (88%)	11 (12%)	5	12

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1150/1212 (95%)	1044 (91%)	106 (9%)	9 20

5 of 106 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	326	LYS
1	F	286	ASN
2	H	43	LYS
1	E	356	GLU
1	E	443	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 12 such sidechains are listed below:

Mol	Chain	Res	Type
2	G	76	ASN
2	G	81	GLN
2	H	112	GLN
2	H	39	GLN
1	F	390	ASN

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)

28 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	$\#\|Z\|>2$
3	NAG	I	1	3,1	14,14,15	0.33	0	17,19,21	0.56	0
3	NAG	I	2	3	14,14,15	0.55	0	17,19,21	0.66	0
3	BMA	I	3	3	11,11,12	1.77	3 (27%)	15,15,17	1.52	3 (20%)
3	MAN	I	4	3	11,11,12	1.42	3 (27%)	15,15,17	1.45	2 (13%)
3	NAG	I	5	3	14,14,15	0.47	0	17,19,21	0.57	0
3	MAN	I	6	3	11,11,12	0.93	0	15,15,17	1.43	2 (13%)
3	NAG	I	7	3	14,14,15	0.42	0	17,19,21	0.57	0
3	NAG	J	1	3,1	14,14,15	0.71	0	17,19,21	0.78	1 (5%)
3	NAG	J	2	3	14,14,15	0.70	0	17,19,21	0.87	1 (5%)
3	BMA	J	3	3	11,11,12	1.00	0	15,15,17	1.30	1 (6%)
3	MAN	J	4	3	11,11,12	0.99	0	15,15,17	1.46	2 (13%)
3	NAG	J	5	3	14,14,15	0.65	1 (7%)	17,19,21	0.76	1 (5%)
3	MAN	J	6	3	11,11,12	0.94	0	15,15,17	1.30	3 (20%)
3	NAG	J	7	3	14,14,15	0.28	0	17,19,21	0.58	0
3	NAG	K	1	3,1	14,14,15	0.49	0	17,19,21	0.58	0
3	NAG	K	2	3	14,14,15	0.70	1 (7%)	17,19,21	0.73	1 (5%)
3	BMA	K	3	3	11,11,12	1.15	2 (18%)	15,15,17	1.44	1 (6%)
3	MAN	K	4	3	11,11,12	1.31	1 (9%)	15,15,17	1.22	1 (6%)
3	NAG	K	5	3	14,14,15	0.45	0	17,19,21	0.59	0
3	MAN	K	6	3	11,11,12	1.20	1 (9%)	15,15,17	1.57	2 (13%)
3	NAG	K	7	3	14,14,15	0.46	0	17,19,21	0.56	0
3	NAG	L	1	3,1	14,14,15	0.38	0	17,19,21	0.79	0
3	NAG	L	2	3	14,14,15	0.35	0	17,19,21	0.74	0
3	BMA	L	3	3	11,11,12	1.33	1 (9%)	15,15,17	1.06	1 (6%)
3	MAN	L	4	3	11,11,12	1.69	2 (18%)	15,15,17	1.40	2 (13%)
3	NAG	L	5	3	14,14,15	0.67	0	17,19,21	0.50	0
3	MAN	L	6	3	11,11,12	1.61	2 (18%)	15,15,17	1.52	2 (13%)
3	NAG	L	7	3	14,14,15	0.43	0	17,19,21	0.52	0

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

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

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	I	5	3	-	4/6/23/26	0/1/1/1
3	MAN	I	6	3	-	0/2/19/22	0/1/1/1
3	NAG	I	7	3	-	0/6/23/26	0/1/1/1
3	NAG	J	1	3,1	-	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	-	0/2/19/22	0/1/1/1
3	MAN	J	4	3	-	2/2/19/22	0/1/1/1
3	NAG	J	5	3	-	3/6/23/26	0/1/1/1
3	MAN	J	6	3	-	0/2/19/22	0/1/1/1
3	NAG	J	7	3	-	4/6/23/26	0/1/1/1
3	NAG	K	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	K	2	3	-	0/6/23/26	0/1/1/1
3	BMA	K	3	3	-	0/2/19/22	0/1/1/1
3	MAN	K	4	3	-	2/2/19/22	0/1/1/1
3	NAG	K	5	3	-	2/6/23/26	0/1/1/1
3	MAN	K	6	3	-	0/2/19/22	0/1/1/1
3	NAG	K	7	3	-	2/6/23/26	0/1/1/1
3	NAG	L	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	L	2	3	-	0/6/23/26	0/1/1/1
3	BMA	L	3	3	-	0/2/19/22	0/1/1/1
3	MAN	L	4	3	-	2/2/19/22	0/1/1/1
3	NAG	L	5	3	-	4/6/23/26	0/1/1/1
3	MAN	L	6	3	-	0/2/19/22	0/1/1/1
3	NAG	L	7	3	-	4/6/23/26	0/1/1/1

The worst 5 of 17 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	L	4	MAN	O5-C1	-4.75	1.36	1.43
3	L	6	MAN	O5-C1	-4.45	1.36	1.43
3	K	4	MAN	O5-C1	-3.50	1.38	1.43
3	I	4	MAN	O5-C1	-3.18	1.38	1.43
3	I	3	BMA	C4-C5	2.90	1.59	1.53

The worst 5 of 26 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	K	3	BMA	O2-C2-C3	-4.40	101.32	110.14
3	I	6	MAN	O2-C2-C3	-4.16	101.81	110.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	J	4	MAN	O2-C2-C3	-4.08	101.95	110.14
3	I	3	BMA	O2-C2-C3	-3.91	102.30	110.14
3	K	6	MAN	C1-C2-C3	-3.66	105.17	109.67

There are no chirality outliers.

5 of 35 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	K	4	MAN	O5-C5-C6-O6
3	J	4	MAN	O5-C5-C6-O6
3	L	4	MAN	O5-C5-C6-O6
3	I	4	MAN	O5-C5-C6-O6
3	L	7	NAG	O5-C5-C6-O6

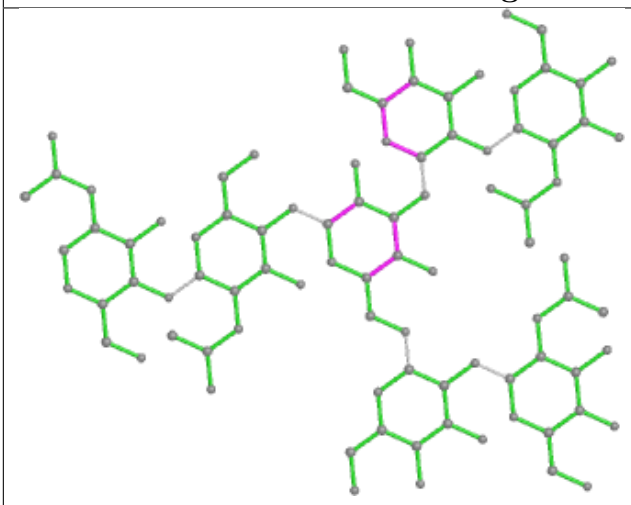
There are no ring outliers.

8 monomers are involved in 7 short contacts:

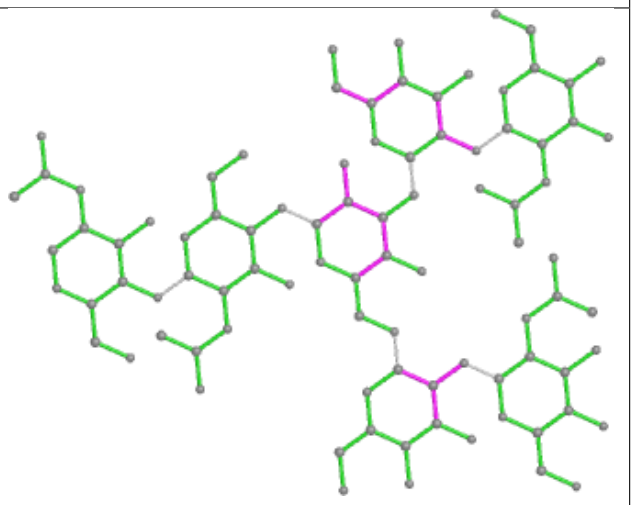
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	I	1	NAG	1	0
3	J	1	NAG	2	0
3	J	7	NAG	1	0
3	K	7	NAG	1	0
3	J	4	MAN	1	0
3	K	4	MAN	1	0
3	K	5	NAG	1	0
3	J	5	NAG	1	0

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

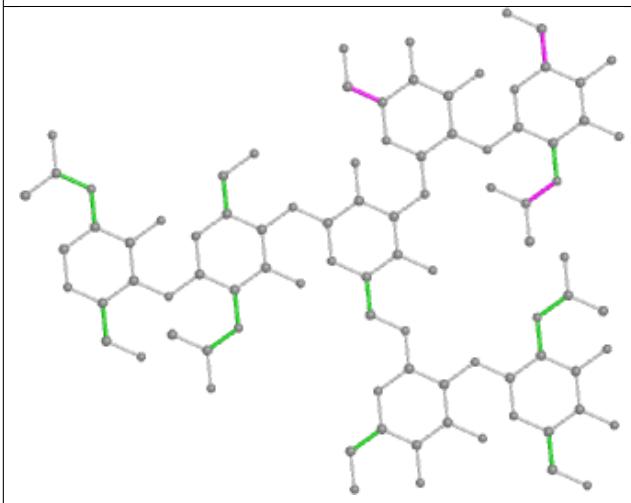
Oligosaccharide Chain I



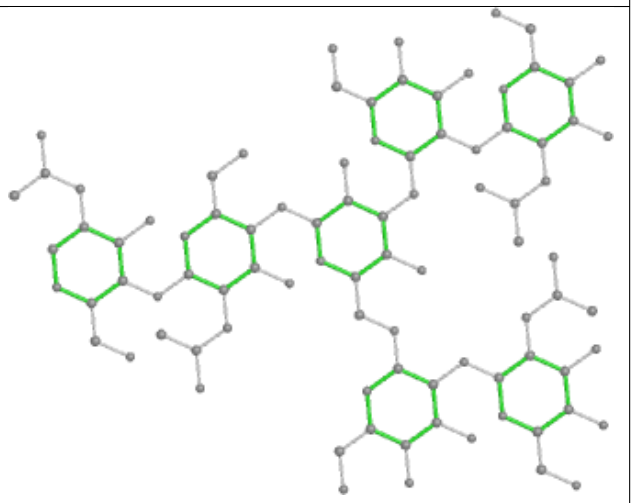
Bond lengths



Bond angles

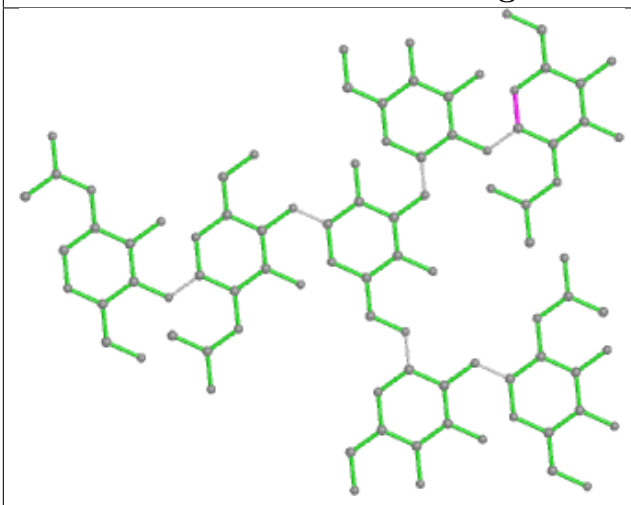


Torsions

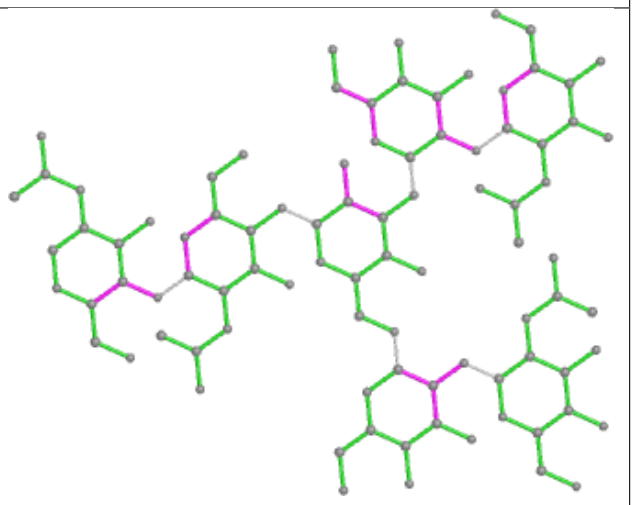


Rings

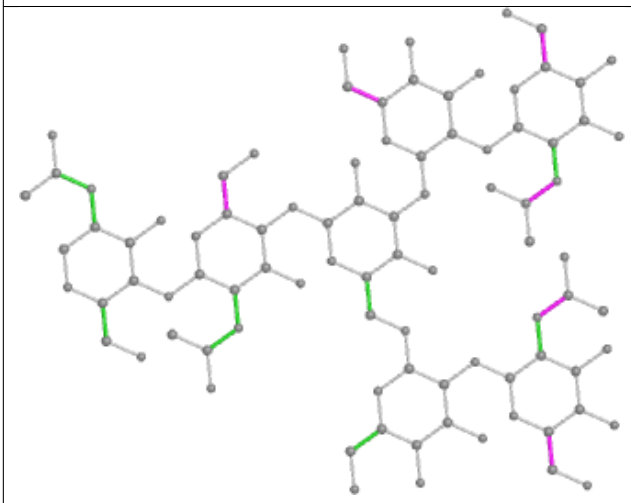
Oligosaccharide Chain J



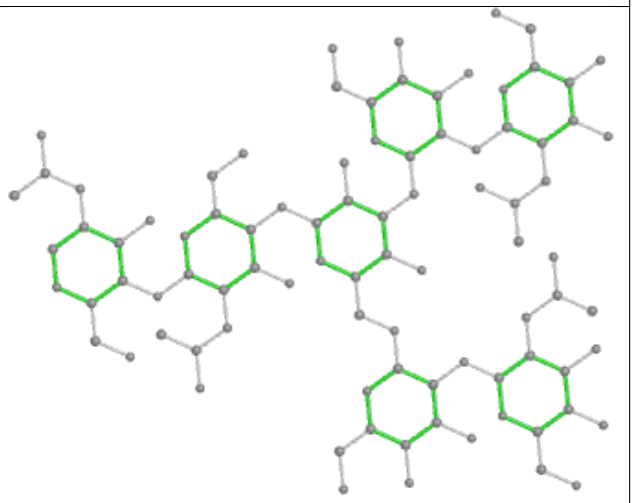
Bond lengths



Bond angles

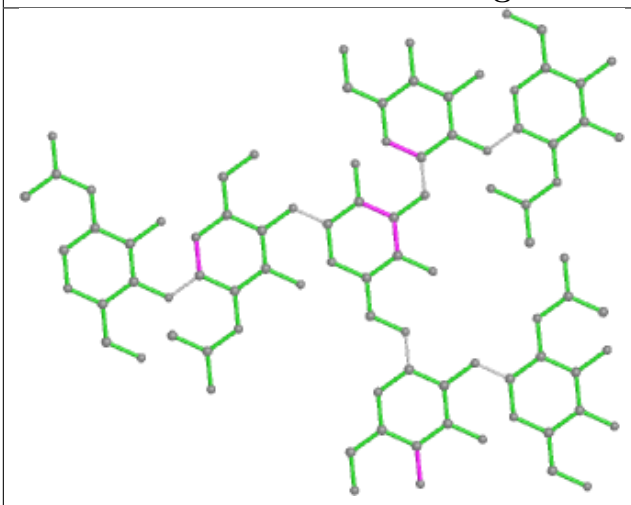


Torsions

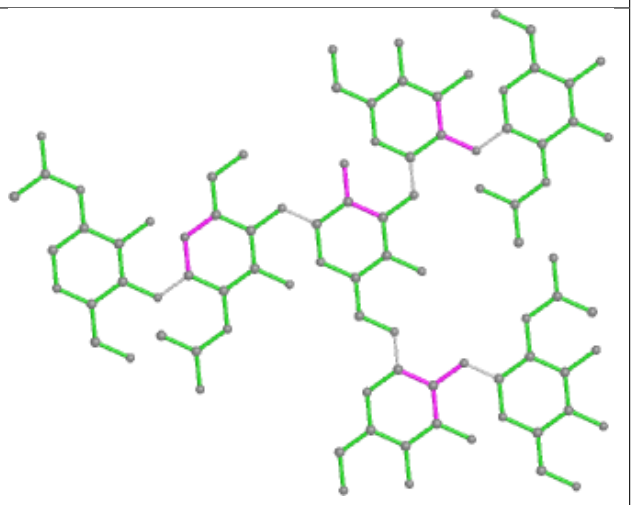


Rings

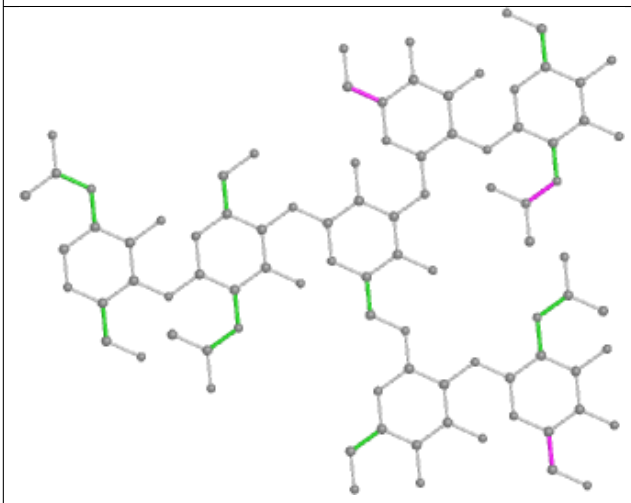
Oligosaccharide Chain K



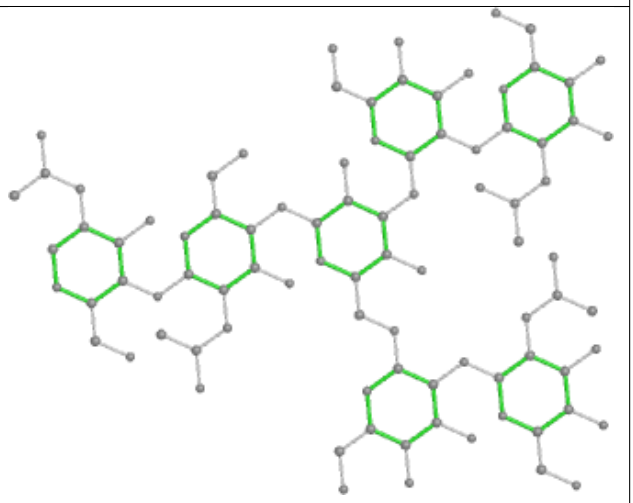
Bond lengths



Bond angles

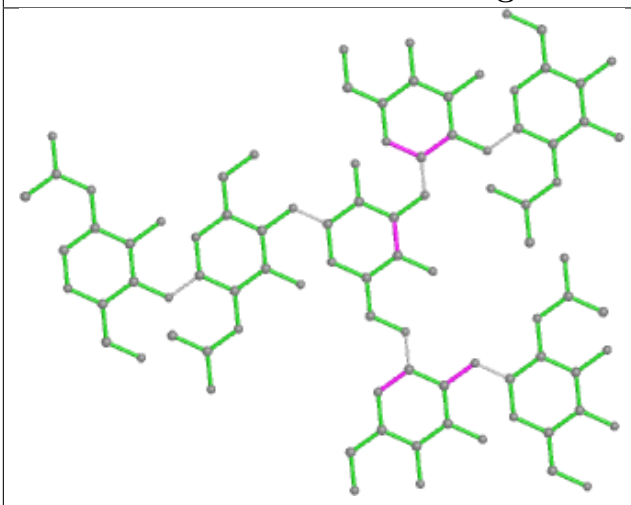


Torsions

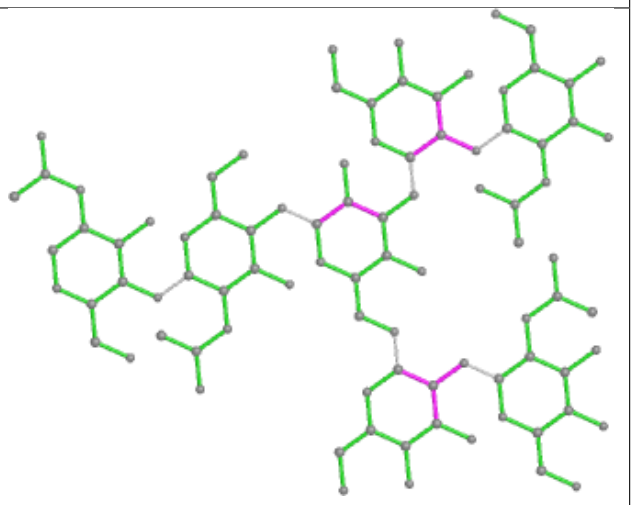


Rings

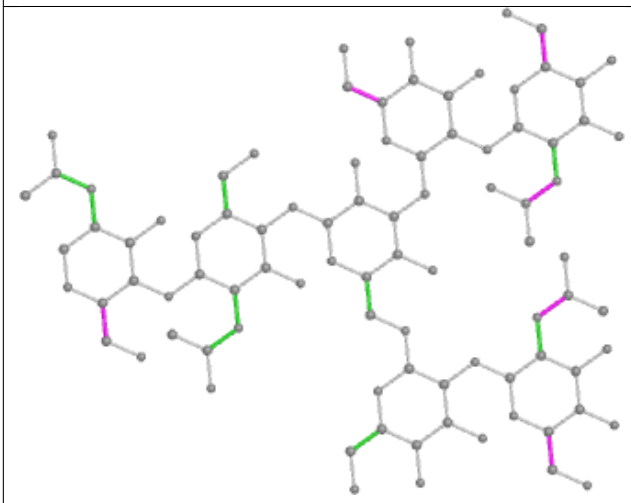
Oligosaccharide Chain L



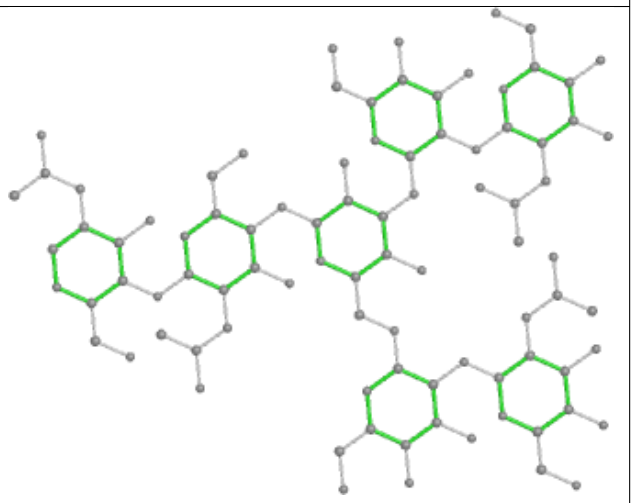
Bond lengths



Bond angles



Torsions



Rings

5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} 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	207/224 (92%)	0.04	1 (0%)	91	92	59, 77, 107, 124	0
1	B	207/224 (92%)	0.05	4 (1%)	66	69	61, 85, 112, 129	0
1	E	206/224 (91%)	0.01	2 (0%)	82	83	61, 84, 128, 146	0
1	F	208/224 (92%)	0.31	7 (3%)	45	45	68, 103, 135, 146	0
2	C	120/120 (100%)	0.02	1 (0%)	86	87	78, 94, 122, 139	0
2	D	120/120 (100%)	0.24	5 (4%)	36	35	69, 95, 131, 144	0
2	G	120/120 (100%)	0.23	5 (4%)	36	35	81, 104, 130, 171	0
2	H	120/120 (100%)	0.35	6 (5%)	28	28	80, 102, 132, 152	0
All	All	1308/1376 (95%)	0.14	31 (2%)	59	60	59, 92, 128, 171	0

The worst 5 of 31 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	1	GLN	11.2
2	H	2	VAL	7.2
2	D	2	VAL	5.5
2	G	1	GLN	5.0
1	F	286	ASN	3.9

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,

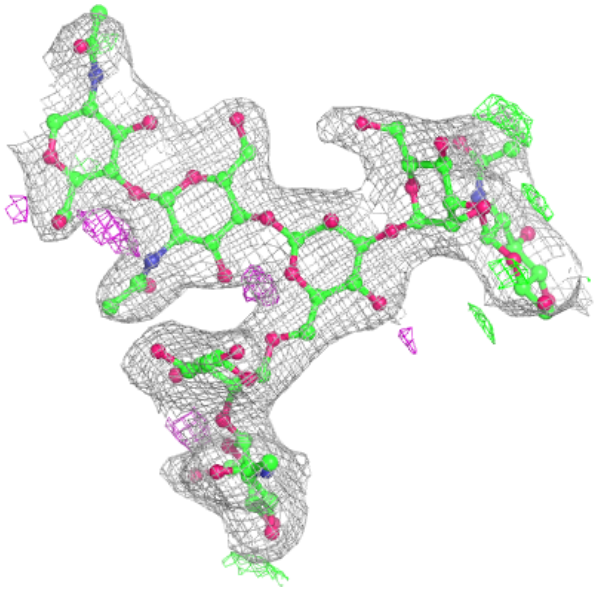
median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

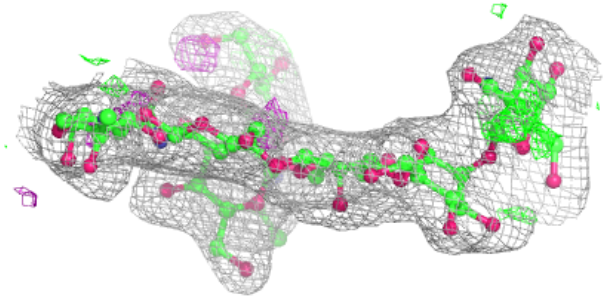
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	NAG	L	5	14/15	0.84	0.22	96,103,109,114	0
3	NAG	J	5	14/15	0.88	0.15	97,107,110,112	0
3	NAG	K	5	14/15	0.88	0.22	80,91,104,104	0
3	NAG	I	5	14/15	0.88	0.20	89,99,107,107	0
3	NAG	J	7	14/15	0.91	0.17	67,75,80,87	0
3	MAN	J	4	11/12	0.93	0.16	95,98,101,105	0
3	NAG	L	7	14/15	0.93	0.14	69,77,81,89	0
3	NAG	I	1	14/15	0.94	0.17	73,77,81,83	0
3	NAG	J	1	14/15	0.94	0.13	90,92,95,96	0
3	MAN	L	4	11/12	0.95	0.16	73,87,92,92	0
3	NAG	I	7	14/15	0.95	0.16	59,66,70,75	0
3	MAN	I	4	11/12	0.95	0.21	81,83,85,88	0
3	BMA	J	3	11/12	0.96	0.15	74,87,90,95	0
3	NAG	K	7	14/15	0.96	0.20	65,70,73,73	0
3	NAG	L	1	14/15	0.96	0.16	72,81,85,86	0
3	NAG	L	2	14/15	0.96	0.15	67,73,75,76	0
3	BMA	L	3	11/12	0.96	0.18	72,75,82,82	0
3	NAG	I	2	14/15	0.96	0.17	57,68,74,76	0
3	BMA	I	3	11/12	0.96	0.19	63,72,77,78	0
3	NAG	J	2	14/15	0.96	0.13	79,86,89,90	0
3	NAG	K	2	14/15	0.97	0.17	57,67,72,72	0
3	MAN	K	4	11/12	0.97	0.21	65,75,80,83	0
3	MAN	J	6	11/12	0.97	0.20	70,74,80,84	0
3	MAN	I	6	11/12	0.97	0.18	62,67,71,71	0
3	MAN	L	6	11/12	0.97	0.17	68,71,74,79	0
3	NAG	K	1	14/15	0.97	0.14	68,72,76,78	0
3	BMA	K	3	11/12	0.98	0.20	59,64,69,70	0
3	MAN	K	6	11/12	0.98	0.21	63,67,71,72	0

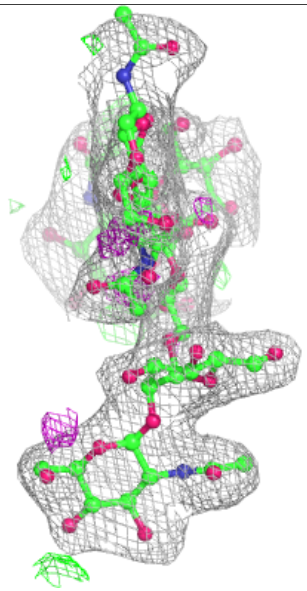
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain I:

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

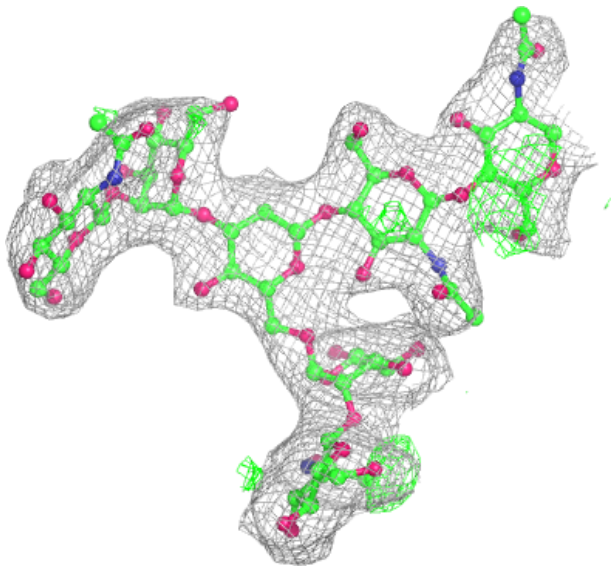


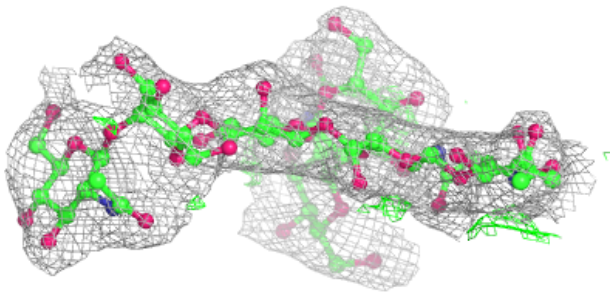


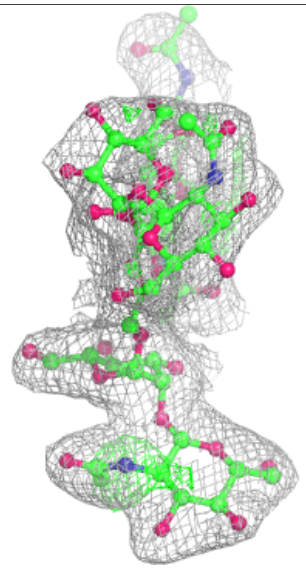


Electron density around Chain J:

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

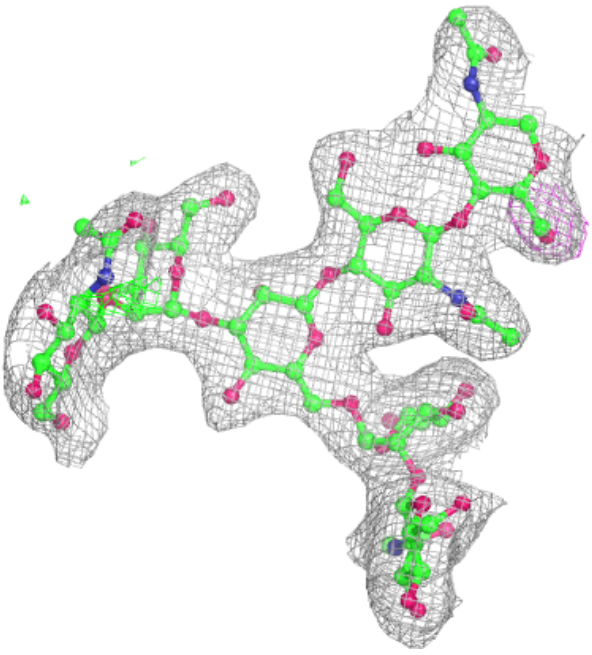


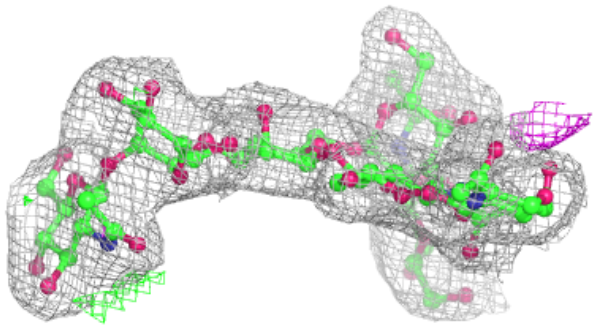


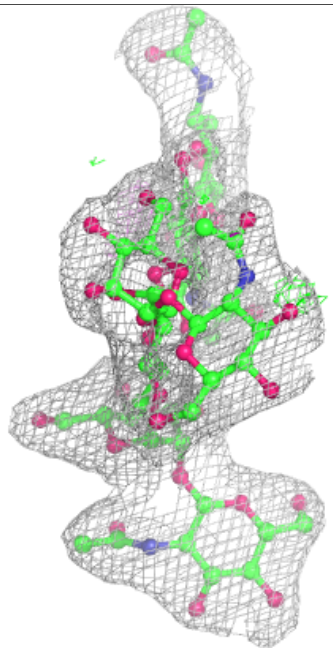


Electron density around Chain K:

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

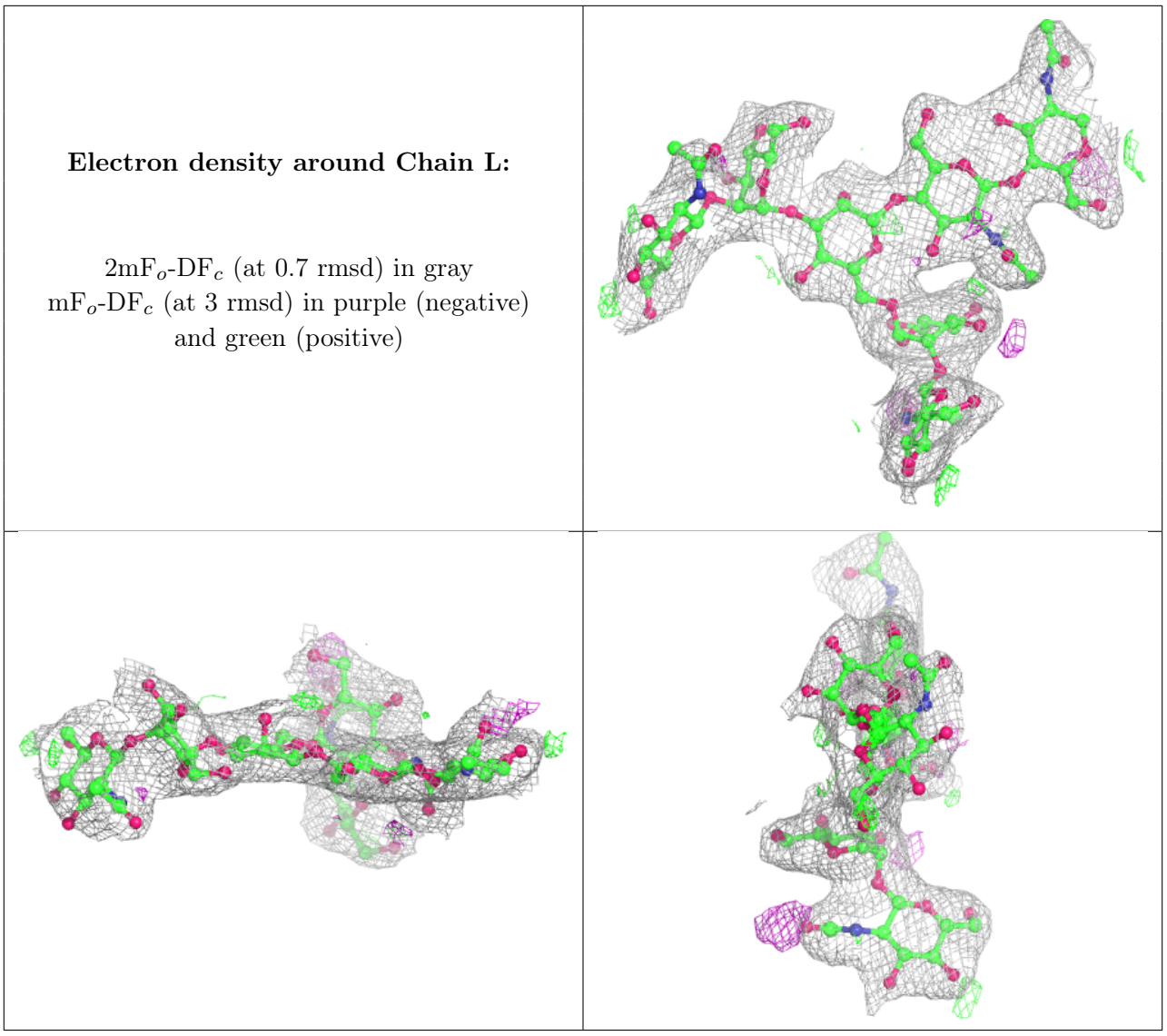






Electron density around Chain L:

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



6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.