



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 17, 2023 – 05:08 PM EST

PDB ID : 4WSS
Title : The crystal structure of hemagglutinin form A/chicken/New York/14677-13/1998 in complex with LSTa
Authors : Yang, H.; Carney, P.J.; Chang, J.C.; Villanueva, J.M.; Stevens, J.
Deposited on : 2014-10-28
Resolution : 2.80 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

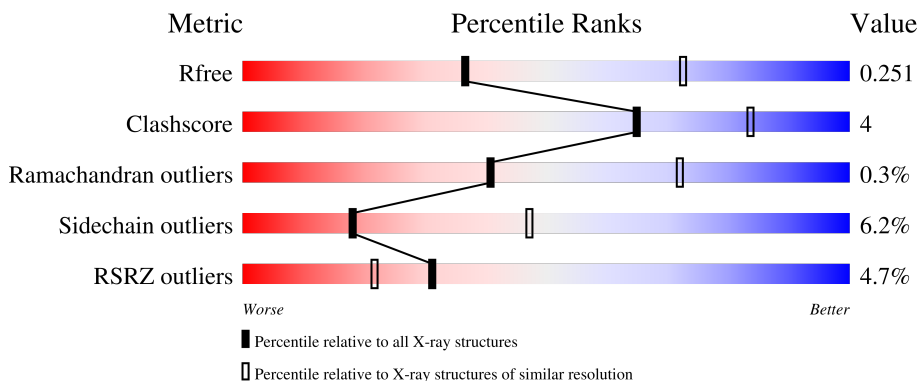
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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	514	 2% 83% 10% • 5%
1	B	514	 2% 80% 12% • 6%
1	C	514	 2% 82% 9% • 6%
1	D	514	 8% 82% 12% • •
1	E	514	 81% 12% • 5%

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Mol	Chain	Length	Quality of chain
1	F	514	 12% 80% 10% 7%
2	G	2	 100%
2	J	2	 100%
3	H	3	 67% 33%
3	I	3	 67% 33%
3	K	3	 100%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	NAG	B	601	X	-	-	-
4	NAG	C	602	X	-	-	-
4	NAG	D	900	X	-	-	-
4	NAG	E	601	X	-	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 23442 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 Hemagglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	490	3895	2446	676	754	19	0	0	0
1	B	481	3821	2399	663	740	19	0	0	0
1	C	482	3832	2405	667	741	19	0	0	0
1	E	489	3886	2441	675	751	19	0	0	0
1	D	493	3909	2458	678	754	19	0	0	0
1	F	479	3810	2391	663	737	19	0	0	0

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	ALA	-	expression tag	UNP Q0A3A5
A	-2	ASP	-	expression tag	UNP Q0A3A5
A	-1	PRO	-	expression tag	UNP Q0A3A5
A	0	GLY	-	expression tag	UNP Q0A3A5
A	317	THR	UNK	conflict	UNP Q0A3A5
A	499	GLU	GLN	conflict	UNP Q0A3A5
A	504	GLY	VAL	conflict	UNP Q0A3A5
A	505	ARG	MET	conflict	UNP Q0A3A5
A	507	VAL	-	expression tag	UNP Q0A3A5
A	508	PRO	-	expression tag	UNP Q0A3A5
A	509	ARG	-	expression tag	UNP Q0A3A5
A	510	GLY	-	expression tag	UNP Q0A3A5
B	-3	ALA	-	expression tag	UNP Q0A3A5
B	-2	ASP	-	expression tag	UNP Q0A3A5
B	-1	PRO	-	expression tag	UNP Q0A3A5
B	0	GLY	-	expression tag	UNP Q0A3A5
B	317	THR	UNK	conflict	UNP Q0A3A5

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Chain	Residue	Modelled	Actual	Comment	Reference
B	499	GLU	GLN	conflict	UNP Q0A3A5
B	504	GLY	VAL	conflict	UNP Q0A3A5
B	505	ARG	MET	conflict	UNP Q0A3A5
B	507	VAL	-	expression tag	UNP Q0A3A5
B	508	PRO	-	expression tag	UNP Q0A3A5
B	509	ARG	-	expression tag	UNP Q0A3A5
B	510	GLY	-	expression tag	UNP Q0A3A5
C	-3	ALA	-	expression tag	UNP Q0A3A5
C	-2	ASP	-	expression tag	UNP Q0A3A5
C	-1	PRO	-	expression tag	UNP Q0A3A5
C	0	GLY	-	expression tag	UNP Q0A3A5
C	317	THR	UNK	conflict	UNP Q0A3A5
C	499	GLU	GLN	conflict	UNP Q0A3A5
C	504	GLY	VAL	conflict	UNP Q0A3A5
C	505	ARG	MET	conflict	UNP Q0A3A5
C	507	VAL	-	expression tag	UNP Q0A3A5
C	508	PRO	-	expression tag	UNP Q0A3A5
C	509	ARG	-	expression tag	UNP Q0A3A5
C	510	GLY	-	expression tag	UNP Q0A3A5
E	-3	ALA	-	expression tag	UNP Q0A3A5
E	-2	ASP	-	expression tag	UNP Q0A3A5
E	-1	PRO	-	expression tag	UNP Q0A3A5
E	0	GLY	-	expression tag	UNP Q0A3A5
E	317	THR	UNK	conflict	UNP Q0A3A5
E	499	GLU	GLN	conflict	UNP Q0A3A5
E	504	GLY	VAL	conflict	UNP Q0A3A5
E	505	ARG	MET	conflict	UNP Q0A3A5
E	507	VAL	-	expression tag	UNP Q0A3A5
E	508	PRO	-	expression tag	UNP Q0A3A5
E	509	ARG	-	expression tag	UNP Q0A3A5
E	510	GLY	-	expression tag	UNP Q0A3A5
D	-3	ALA	-	expression tag	UNP Q0A3A5
D	-2	ASP	-	expression tag	UNP Q0A3A5
D	-1	PRO	-	expression tag	UNP Q0A3A5
D	0	GLY	-	expression tag	UNP Q0A3A5
D	317	THR	UNK	conflict	UNP Q0A3A5
D	499	GLU	GLN	conflict	UNP Q0A3A5
D	504	GLY	VAL	conflict	UNP Q0A3A5
D	505	ARG	MET	conflict	UNP Q0A3A5
D	507	VAL	-	expression tag	UNP Q0A3A5
D	508	PRO	-	expression tag	UNP Q0A3A5
D	509	ARG	-	expression tag	UNP Q0A3A5

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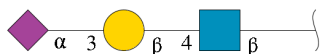
Chain	Residue	Modelled	Actual	Comment	Reference
D	510	GLY	-	expression tag	UNP Q0A3A5
F	-3	ALA	-	expression tag	UNP Q0A3A5
F	-2	ASP	-	expression tag	UNP Q0A3A5
F	-1	PRO	-	expression tag	UNP Q0A3A5
F	0	GLY	-	expression tag	UNP Q0A3A5
F	317	THR	UNK	conflict	UNP Q0A3A5
F	499	GLU	GLN	conflict	UNP Q0A3A5
F	504	GLY	VAL	conflict	UNP Q0A3A5
F	505	ARG	MET	conflict	UNP Q0A3A5
F	507	VAL	-	expression tag	UNP Q0A3A5
F	508	PRO	-	expression tag	UNP Q0A3A5
F	509	ARG	-	expression tag	UNP Q0A3A5
F	510	GLY	-	expression tag	UNP Q0A3A5

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



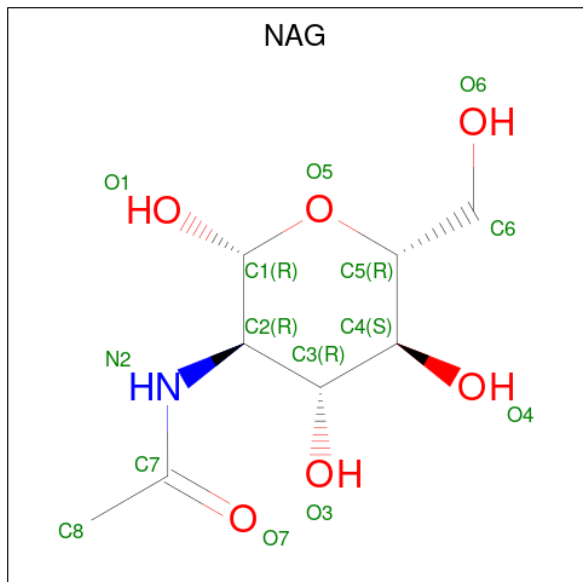
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	J	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	H	3	Total	C	N	O	0	0	0
			45	25	2	18			
3	I	3	Total	C	N	O	0	0	0
			45	25	2	18			
3	K	3	Total	C	N	O	0	0	0
			45	25	2	18			

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

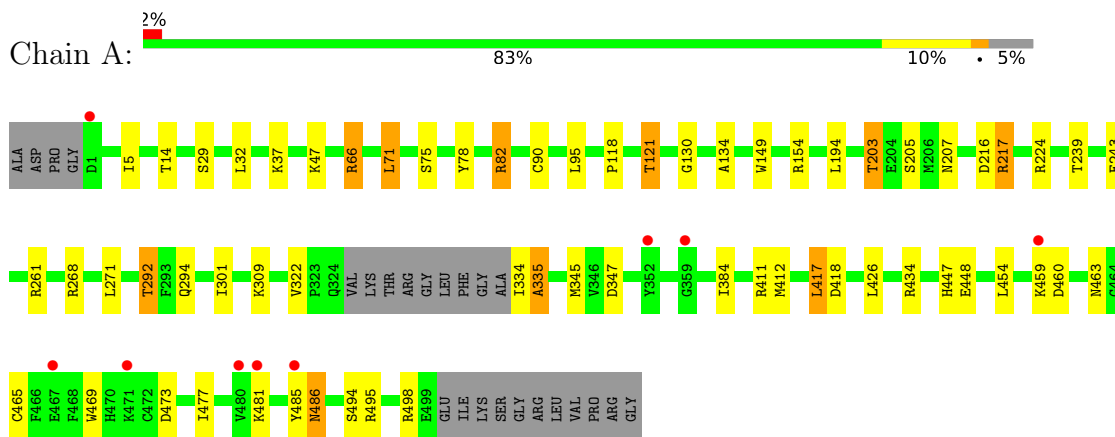


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
4	B	1	Total 14	C 8	N 1	O 5	0	0
4	B	1	Total 14	C 8	N 1	O 5	0	0
4	C	1	Total 14	C 8	N 1	O 5	0	0
4	C	1	Total 14	C 8	N 1	O 5	0	0
4	E	1	Total 14	C 8	N 1	O 5	0	0
4	E	1	Total 14	C 8	N 1	O 5	0	0
4	D	1	Total 14	C 8	N 1	O 5	0	0

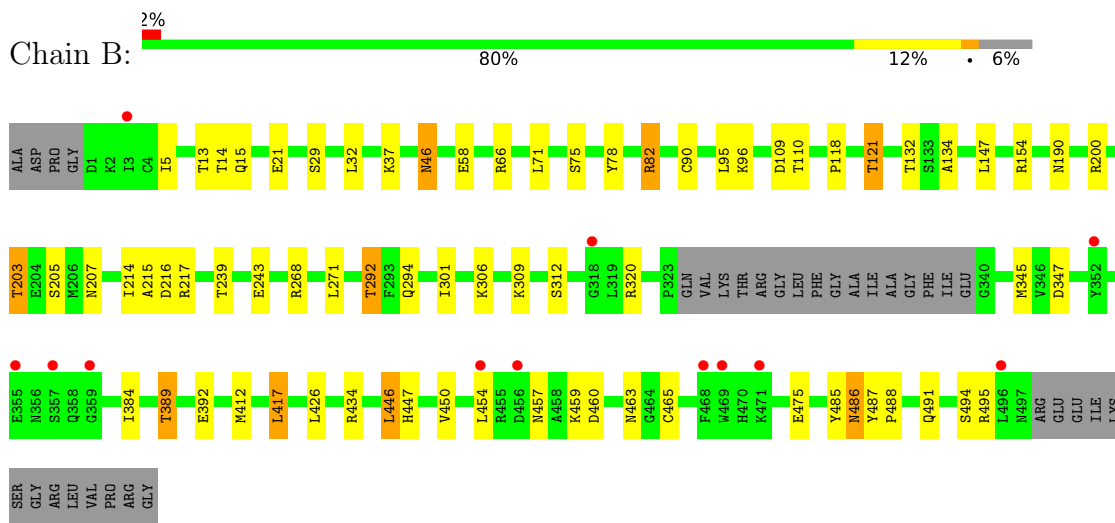
3 Residue-property plots [i](#)

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.

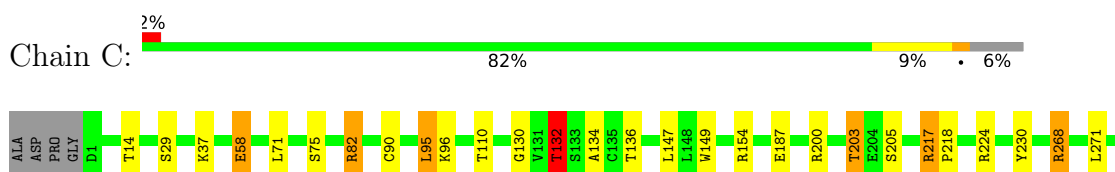
- Molecule 1: Hemagglutinin

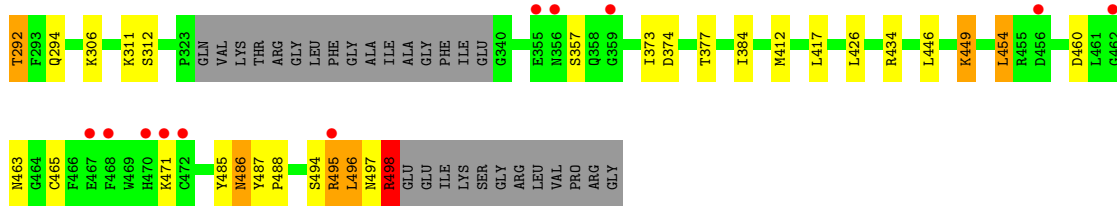


- Molecule 1: Hemagglutinin



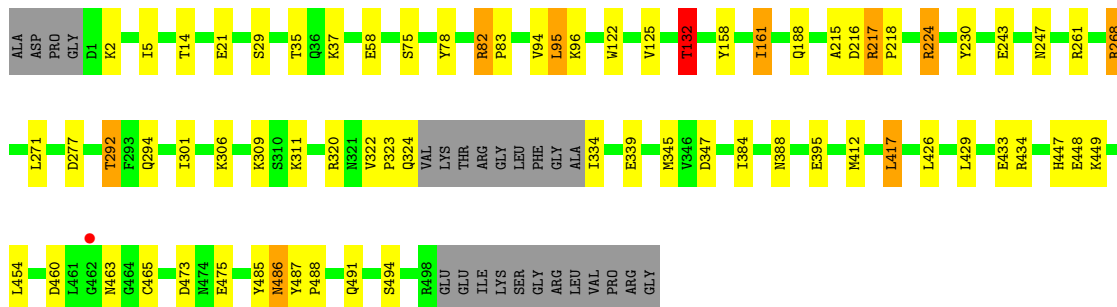
- Molecule 1: Hemagglutinin





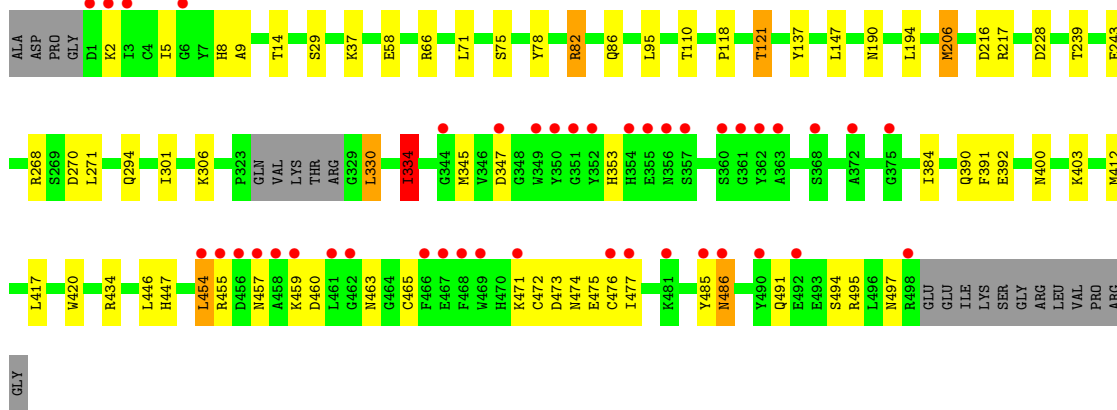
● Molecule 1: Hemagglutinin

Chain E: • 5%



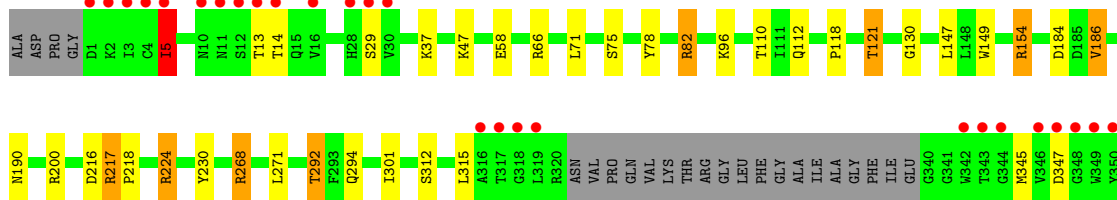
● Molecule 1: Hemagglutinin

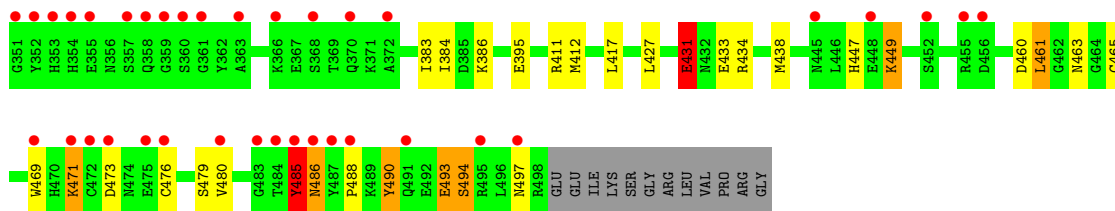
Chain D: • •



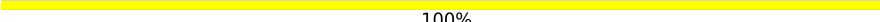
● Molecule 1: Hemagglutinin

Chain F: • • • 7%



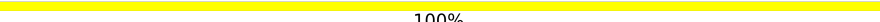


- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:  100%



MAG1
MAG2

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

Chain J:  100%


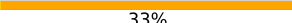
MAG1
MAG2

- Molecule 3: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H:  67%  33%

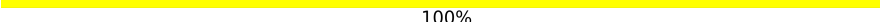
MAG1
GAL2
SIA3

- Molecule 3: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  67%  33%

MAG1
GAL2
SIA3

- Molecule 3: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K:  100%

MAG1
GAL2
SIA3

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	254.97Å 134.78Å 122.62Å 90.00° 112.20° 90.00°	Depositor
Resolution (Å)	45.20 – 2.80 45.20 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.5 (45.20-2.80) 98.5 (45.20-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.63 (at 2.81Å)	Xtrriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.220 , 0.254 0.221 , 0.251	Depositor DCC
R_{free} test set	4650 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	52.3	Xtrriage
Anisotropy	0.359	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 18.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	23442	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.90% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: SIA, GAL, NAG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.64	0/3983	0.77	5/5397 (0.1%)
1	B	0.61	0/3908	0.76	5/5297 (0.1%)
1	C	0.62	0/3919	0.77	4/5311 (0.1%)
1	D	0.64	0/3998	0.80	7/5417 (0.1%)
1	E	0.64	0/3974	0.79	6/5385 (0.1%)
1	F	0.64	0/3896	0.81	7/5278 (0.1%)
All	All	0.63	0/23678	0.79	34/32085 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	1
1	E	0	1
All	All	0	2

There are no bond length outliers.

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	461	LEU	CA-CB-CG	9.70	137.62	115.30
1	E	277	ASP	CB-CG-OD2	8.01	125.51	118.30
1	F	485	TYR	N-CA-C	7.71	131.82	111.00
1	F	431	GLU	CB-CA-C	7.60	125.60	110.40
1	C	496	LEU	CA-CB-CG	-6.86	99.53	115.30
1	D	495	ARG	NE-CZ-NH1	6.77	123.69	120.30
1	A	473	ASP	CB-CG-OD2	6.50	124.15	118.30
1	D	206	MET	CA-CB-CG	6.26	123.94	113.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	112	GLN	CB-CA-C	-6.17	98.07	110.40
1	E	261	ARG	NE-CZ-NH1	6.14	123.37	120.30
1	A	243	GLU	CB-CA-C	-5.83	98.73	110.40
1	F	5	ILE	CB-CA-C	5.82	123.23	111.60
1	B	446	LEU	CB-CG-CD2	5.81	120.88	111.00
1	A	498	ARG	NE-CZ-NH1	5.78	123.19	120.30
1	F	473	ASP	CB-CG-OD2	-5.73	113.14	118.30
1	B	66	ARG	CG-CD-NE	5.66	123.68	111.80
1	D	417	LEU	CB-CG-CD2	5.66	120.62	111.00
1	C	132	THR	CB-CA-C	-5.59	96.51	111.60
1	A	194	LEU	CB-CG-CD2	-5.55	101.57	111.00
1	D	446	LEU	CB-CG-CD2	5.49	120.33	111.00
1	D	243	GLU	CB-CA-C	-5.46	99.48	110.40
1	B	46	ASN	CB-CA-C	-5.45	99.50	110.40
1	D	384	ILE	CA-CB-CG1	5.43	121.32	111.00
1	B	216	ASP	CB-CG-OD1	5.36	123.12	118.30
1	C	498	ARG	NE-CZ-NH1	5.33	122.96	120.30
1	C	446	LEU	CB-CG-CD2	5.31	120.03	111.00
1	D	228	ASP	CB-CG-OD2	-5.22	113.60	118.30
1	E	132	THR	CB-CA-C	-5.21	97.53	111.60
1	E	161	ILE	CB-CA-C	-5.21	101.18	111.60
1	E	473	ASP	CB-CG-OD1	5.20	122.98	118.30
1	E	434	ARG	NE-CZ-NH1	5.17	122.89	120.30
1	B	109	ASP	N-CA-CB	5.13	119.84	110.60
1	F	66	ARG	NE-CZ-NH2	-5.06	117.77	120.30
1	A	473	ASP	CB-CG-OD1	-5.04	113.77	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	334	ILE	Peptide
1	E	323	PRO	Peptide

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	3895	0	3753	33	0
1	B	3821	0	3680	29	0
1	C	3832	0	3693	34	0
1	D	3909	0	3770	32	2
1	E	3886	0	3746	37	0
1	F	3810	0	3672	40	0
2	G	28	0	25	0	0
2	J	28	0	25	0	0
3	H	45	0	38	4	0
3	I	45	0	38	1	0
3	K	45	0	38	0	0
4	B	28	0	26	1	0
4	C	28	0	26	1	0
4	D	14	0	13	0	0
4	E	28	0	26	2	0
All	All	23442	0	22569	187	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 4.

All (187) 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:D:457:ASN:ND2	1:D:485:TYR:OH	1.98	0.96
1:B:457:ASN:ND2	1:B:485:TYR:OH	1.98	0.96
1:D:473:ASP:O	1:D:476:CYS:SG	2.24	0.96
1:E:2:LYS:NZ	1:E:334:ILE:O	1.98	0.95
1:F:469:TRP:CD1	1:F:494:SER:OG	2.20	0.95
1:A:207:ASN:OD1	1:C:217:ARG:NH1	2.00	0.94
1:F:476:CYS:SG	1:F:485:TYR:OH	2.29	0.89
1:E:268:ARG:HB3	1:E:268:ARG:NH1	1.87	0.88
1:D:476:CYS:SG	1:D:477:ILE:N	2.46	0.88
1:A:463:ASN:O	1:A:463:ASN:OD1	1.99	0.79
1:C:311:LYS:O	4:C:602:NAG:H82	1.83	0.79
1:E:188:GLN:HE22	1:E:247:ASN:HD21	1.33	0.77
1:E:122:TRP:HE3	1:E:161:ILE:CG2	1.99	0.75
1:D:474:ASN:HA	1:D:477:ILE:HG12	1.67	0.74
1:F:490:TYR:O	1:F:493:GLU:OE1	2.05	0.73
1:B:118:PRO:O	1:B:121:THR:HG23	1.90	0.72
1:D:118:PRO:O	1:D:121:THR:HG23	1.90	0.72
1:A:118:PRO:O	1:A:121:THR:HG23	1.90	0.71
1:F:118:PRO:O	1:F:121:THR:HG23	1.90	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:268:ARG:HB3	1:E:268:ARG:HH11	1.54	0.71
1:E:311:LYS:HB3	4:E:601:NAG:H82	1.72	0.70
1:B:203:THR:HG22	1:B:205:SER:H	1.57	0.69
1:D:334:ILE:HD11	1:D:353:HIS:CE1	2.27	0.69
1:A:203:THR:HG22	1:A:205:SER:H	1.58	0.69
1:D:334:ILE:HD11	1:D:353:HIS:NE2	2.09	0.67
1:C:203:THR:HG22	1:C:205:SER:H	1.61	0.65
1:E:122:TRP:CE3	1:E:161:ILE:CG2	2.80	0.64
1:F:427:LEU:O	1:F:431:GLU:HG3	1.97	0.64
1:C:373:ILE:O	1:C:377:THR:HG22	1.98	0.64
1:D:306:LYS:HD3	1:D:390:GLN:HA	1.80	0.64
1:B:446:LEU:O	1:B:450:VAL:HG12	1.96	0.64
1:F:493:GLU:H	1:F:493:GLU:CD	2.01	0.64
1:F:480:VAL:HG22	1:F:485:TYR:CE2	2.34	0.62
1:B:154:ARG:NH1	1:B:190:ASN:OD1	2.32	0.62
1:F:82:ARG:NH2	1:F:271:LEU:O	2.32	0.62
1:C:132:THR:CG2	3:H:3:SIA:O1B	2.49	0.61
1:E:82:ARG:NH2	1:E:271:LEU:O	2.33	0.61
1:A:216:ASP:N	1:B:243:GLU:OE1	2.26	0.60
1:F:383:ILE:HD11	1:F:431:GLU:OE2	2.01	0.60
1:D:306:LYS:HE3	1:D:391:PHE:H	1.66	0.60
1:E:5:ILE:HG13	1:E:447:HIS:HA	1.84	0.60
1:A:82:ARG:NH2	1:A:271:LEU:O	2.34	0.60
1:B:5:ILE:HG13	1:B:447:HIS:HA	1.85	0.59
1:E:388:ASN:O	1:E:388:ASN:OD1	2.21	0.59
1:D:5:ILE:HG13	1:D:447:HIS:HA	1.84	0.59
1:D:82:ARG:NH2	1:D:271:LEU:O	2.35	0.59
1:A:66:ARG:CZ	1:A:66:ARG:HB3	2.33	0.59
1:E:96:LYS:HG3	1:E:230:TYR:CE1	2.38	0.59
1:A:207:ASN:CG	1:C:217:ARG:NH1	2.57	0.58
1:A:5:ILE:HG13	1:A:447:HIS:HA	1.85	0.58
1:B:82:ARG:NH2	1:B:271:LEU:O	2.37	0.58
1:C:82:ARG:NH2	1:C:271:LEU:O	2.35	0.58
1:A:459:LYS:HE2	1:A:469:TRP:CZ2	2.40	0.57
1:D:400:ASN:O	1:D:403:LYS:HE3	2.03	0.57
1:C:96:LYS:HG3	1:C:230:TYR:CE1	2.40	0.56
1:D:334:ILE:HD11	1:D:353:HIS:CD2	2.39	0.56
1:B:15:GLN:OE1	4:B:601:NAG:H82	2.06	0.55
1:C:357:SER:OG	1:C:471:LYS:HE2	2.07	0.55
1:E:14:THR:HG21	1:E:29:SER:OG	2.07	0.55
1:A:418:ASP:HB3	1:B:389:THR:HG21	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:2:LYS:NZ	1:D:334:ILE:O	2.39	0.55
1:F:471:LYS:HD3	1:F:471:LYS:H	1.72	0.55
1:F:315:LEU:HD11	1:F:431:GLU:CD	2.28	0.54
1:A:14:THR:HG21	1:A:29:SER:OG	2.08	0.54
1:E:460:ASP:O	1:E:465:CYS:O	2.26	0.54
1:E:83:PRO:HA	1:E:268:ARG:HH12	1.73	0.53
1:F:14:THR:HG21	1:F:29:SER:OG	2.08	0.53
1:B:14:THR:HG21	1:B:29:SER:OG	2.08	0.53
1:F:75:SER:HB2	1:F:110:THR:HG22	1.91	0.53
1:F:154:ARG:NH2	1:F:190:ASN:OD1	2.34	0.53
1:E:433:GLU:OE2	1:F:434:ARG:NH2	2.42	0.53
1:C:460:ASP:O	1:C:465:CYS:O	2.27	0.52
1:F:96:LYS:HG3	1:F:230:TYR:CE1	2.44	0.52
1:C:14:THR:HG21	1:C:29:SER:OG	2.09	0.52
1:B:460:ASP:O	1:B:465:CYS:O	2.26	0.52
1:D:460:ASP:O	1:D:465:CYS:O	2.27	0.52
1:F:292:THR:HG21	1:F:384:ILE:HG12	1.92	0.52
1:D:14:THR:HG21	1:D:29:SER:OG	2.10	0.51
1:F:469:TRP:CG	1:F:494:SER:OG	2.48	0.51
1:A:460:ASP:O	1:A:465:CYS:O	2.27	0.51
1:F:383:ILE:CD1	1:F:431:GLU:OE2	2.59	0.51
1:D:75:SER:HB2	1:D:110:THR:HG22	1.92	0.51
1:C:132:THR:HG23	3:H:3:SIA:O1B	2.11	0.50
1:C:75:SER:HB2	1:C:110:THR:HG22	1.92	0.50
1:D:8:HIS:CD2	1:D:9:ALA:O	2.65	0.50
1:B:292:THR:HG21	1:B:384:ILE:HG12	1.93	0.50
1:F:431:GLU:OE1	1:F:431:GLU:C	2.50	0.50
1:A:292:THR:HG21	1:A:384:ILE:HG12	1.92	0.50
1:A:334:ILE:O	1:A:335:ALA:HB3	2.12	0.50
1:C:292:THR:HG21	1:C:384:ILE:HG12	1.93	0.49
1:E:454:LEU:HD23	1:E:485:TYR:CE2	2.48	0.49
1:E:215:ALA:HA	1:F:200:ARG:HD3	1.94	0.49
1:D:472:CYS:C	1:D:476:CYS:SG	2.91	0.49
1:F:493:GLU:OE1	1:F:493:GLU:N	2.45	0.49
1:F:460:ASP:O	1:F:465:CYS:O	2.30	0.49
1:E:292:THR:HG21	1:E:384:ILE:HG12	1.95	0.48
1:C:132:THR:HG22	3:H:3:SIA:O1B	2.14	0.48
1:E:125:VAL:HG11	1:E:161:ILE:HD11	1.95	0.48
1:B:75:SER:HB2	1:B:110:THR:HG22	1.95	0.48
1:A:434:ARG:HH22	1:C:434:ARG:CG	2.27	0.48
1:A:477:ILE:HG23	1:A:481:LYS:HZ2	1.79	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:217:ARG:HB2	1:C:218:PRO:HD2	1.97	0.47
1:E:158:TYR:CE2	1:E:161:ILE:HD12	2.50	0.47
1:A:454:LEU:HD23	1:A:485:TYR:CE2	2.50	0.47
1:B:454:LEU:HD23	1:B:485:TYR:CE2	2.49	0.47
1:C:187:GLU:OE2	3:H:3:SIA:H92	2.15	0.47
1:E:243:GLU:OE1	1:D:216:ASP:N	2.42	0.47
1:F:216:ASP:HB2	1:F:224:ARG:HH11	1.79	0.47
1:E:216:ASP:HB2	1:E:224:ARG:HH11	1.79	0.46
1:D:8:HIS:CD2	1:D:9:ALA:N	2.84	0.46
1:B:215:ALA:HB2	1:C:200:ARG:HD3	1.98	0.46
1:A:71:LEU:HD12	1:D:270:ASP:HB2	1.98	0.45
1:B:345:MET:O	1:B:347:ASP:N	2.48	0.45
1:E:21:GLU:OE1	1:E:320:ARG:NH2	2.48	0.45
1:B:214:ILE:O	1:C:200:ARG:HD3	2.16	0.45
1:F:345:MET:O	1:F:347:ASP:N	2.48	0.45
1:D:330:LEU:HD22	1:D:330:LEU:H	1.81	0.45
1:D:345:MET:O	1:D:347:ASP:N	2.48	0.45
1:F:268:ARG:HE	1:F:268:ARG:HB3	1.63	0.45
1:A:411:ARG:HD2	1:B:392:GLU:O	2.17	0.45
1:C:449:LYS:HD3	1:C:449:LYS:HA	1.85	0.45
1:E:78:TYR:CD2	1:E:301:ILE:HD11	2.52	0.45
1:A:463:ASN:OD1	1:A:463:ASN:C	2.54	0.45
1:C:495:ARG:HA	1:C:498:ARG:NH1	2.32	0.45
1:A:78:TYR:CD2	1:A:301:ILE:HD11	2.52	0.45
1:D:78:TYR:CD2	1:D:301:ILE:HD11	2.52	0.44
1:F:217:ARG:HB2	1:F:218:PRO:HD2	1.99	0.44
1:A:130:GLY:HA3	1:A:149:TRP:HB3	1.99	0.44
1:E:58:GLU:HG2	1:E:95:LEU:HD12	1.98	0.44
1:B:78:TYR:CD2	1:B:301:ILE:HD11	2.52	0.44
1:C:374:ASP:HA	1:C:377:THR:HG22	2.00	0.44
1:F:493:GLU:CD	1:F:493:GLU:N	2.69	0.44
1:D:474:ASN:O	1:D:477:ILE:HG13	2.18	0.44
1:F:449:LYS:HA	1:F:449:LYS:HD3	1.72	0.44
1:F:485:TYR:O	1:F:486:ASN:CB	2.66	0.44
1:A:434:ARG:CG	1:B:434:ARG:HH22	2.31	0.44
1:C:454:LEU:HD12	1:C:485:TYR:CE2	2.52	0.44
1:A:309:LYS:HE3	1:A:417:LEU:HD13	2.00	0.43
1:D:454:LEU:HD12	1:D:485:TYR:CE2	2.53	0.43
1:F:479:SER:OG	1:F:485:TYR:CE1	2.71	0.43
1:F:485:TYR:O	1:F:486:ASN:HB3	2.18	0.43
1:B:21:GLU:OE1	1:B:320:ARG:NH2	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:184:ASP:OD1	1:F:186:VAL:HG13	2.18	0.43
1:A:345:MET:O	1:A:347:ASP:N	2.50	0.43
1:A:459:LYS:HE2	1:A:469:TRP:CH2	2.54	0.43
1:F:78:TYR:CD2	1:F:301:ILE:HD11	2.53	0.43
1:A:434:ARG:HH22	1:C:434:ARG:HG2	1.82	0.43
1:E:429:LEU:HD11	1:F:386:LYS:HE3	2.00	0.43
1:D:434:ARG:NH1	1:F:433:GLU:OE1	2.52	0.43
1:A:217:ARG:NH1	1:B:207:ASN:OD1	2.44	0.42
1:F:47:LYS:HA	1:F:47:LYS:HD3	1.90	0.42
1:E:35:THR:HG21	4:E:601:NAG:O7	2.19	0.42
1:E:306:LYS:HB2	1:E:306:LYS:HE2	1.92	0.42
1:E:485:TYR:O	1:E:486:ASN:CB	2.68	0.42
1:E:345:MET:O	1:E:347:ASP:N	2.49	0.42
1:C:58:GLU:HG2	1:C:95:LEU:HD12	2.00	0.41
1:C:485:TYR:O	1:C:486:ASN:CB	2.67	0.41
1:E:158:TYR:CE2	1:E:161:ILE:CD1	3.03	0.41
1:E:217:ARG:HB2	1:E:218:PRO:HD2	2.02	0.41
1:F:130:GLY:HA3	1:F:149:TRP:HB3	2.02	0.41
1:A:477:ILE:HG23	1:A:481:LYS:NZ	2.35	0.41
1:A:90:CYS:HB2	1:A:134:ALA:O	2.21	0.41
1:C:90:CYS:HB2	1:C:134:ALA:O	2.21	0.41
1:A:216:ASP:OD2	1:A:224:ARG:NH2	2.52	0.41
1:D:306:LYS:HE2	1:D:420:TRP:CG	2.55	0.41
1:A:485:TYR:O	1:A:486:ASN:CB	2.69	0.41
1:E:324:GLN:HB2	1:E:339:GLU:O	2.19	0.41
1:D:334:ILE:HG21	1:D:334:ILE:HD13	1.85	0.41
1:B:200:ARG:HB2	1:B:243:GLU:HG3	2.02	0.41
1:E:485:TYR:O	1:E:486:ASN:HB3	2.21	0.41
1:F:5:ILE:HG21	1:F:447:HIS:HA	2.03	0.41
1:B:306:LYS:HB2	1:B:306:LYS:HE2	1.91	0.41
1:B:309:LYS:HE3	1:B:417:LEU:HD13	2.02	0.41
1:B:485:TYR:O	1:B:486:ASN:CB	2.68	0.41
1:C:306:LYS:HB2	1:C:306:LYS:HE2	1.92	0.41
1:E:132:THR:CG2	3:I:3:SIA:O1A	2.69	0.41
1:C:485:TYR:O	1:C:486:ASN:HB3	2.20	0.41
1:E:487:TYR:N	1:E:488:PRO:HD2	2.36	0.41
1:D:485:TYR:O	1:D:486:ASN:CB	2.68	0.41
1:B:487:TYR:N	1:B:488:PRO:HD2	2.36	0.40
1:C:268:ARG:HE	1:C:268:ARG:HB3	1.60	0.40
1:C:374:ASP:HA	1:C:377:THR:CG2	2.51	0.40
1:C:487:TYR:N	1:C:488:PRO:HD2	2.37	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:392:GLU:O	1:F:411:ARG:HD2	2.21	0.40
1:B:90:CYS:HB2	1:B:134:ALA:O	2.21	0.40
1:C:130:GLY:HA3	1:C:149:TRP:HB3	2.03	0.40
1:E:309:LYS:HE3	1:E:417:LEU:HD13	2.02	0.40

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:D:137:TYR:OH	1:D:137:TYR:OH[2_555]	1.91	0.29
1:D:137:TYR:CZ	1:D:137:TYR:OH[2_555]	2.19	0.01

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	486/514 (95%)	465 (96%)	19 (4%)	2 (0%)	34	66
1	B	477/514 (93%)	457 (96%)	19 (4%)	1 (0%)	47	78
1	C	478/514 (93%)	457 (96%)	19 (4%)	2 (0%)	34	66
1	D	489/514 (95%)	468 (96%)	19 (4%)	2 (0%)	34	66
1	E	485/514 (94%)	464 (96%)	20 (4%)	1 (0%)	47	78
1	F	475/514 (92%)	451 (95%)	22 (5%)	2 (0%)	34	66
All	All	2890/3084 (94%)	2762 (96%)	118 (4%)	10 (0%)	41	72

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	486	ASN
1	A	486	ASN
1	B	486	ASN

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Mol	Chain	Res	Type
1	C	486	ASN
1	D	486	ASN
1	D	497	ASN
1	F	486	ASN
1	F	497	ASN
1	A	335	ALA
1	C	497	ASN

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	426/443 (96%)	402 (94%)	24 (6%)	21	51
1	B	419/443 (95%)	390 (93%)	29 (7%)	15	41
1	C	420/443 (95%)	394 (94%)	26 (6%)	18	47
1	D	426/443 (96%)	399 (94%)	27 (6%)	18	46
1	E	425/443 (96%)	403 (95%)	22 (5%)	23	55
1	F	417/443 (94%)	387 (93%)	30 (7%)	14	38
All	All	2533/2658 (95%)	2375 (94%)	158 (6%)	18	47

All (158) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	LEU
1	A	37	LYS
1	A	47	LYS
1	A	66	ARG
1	A	71	LEU
1	A	75	SER
1	A	82	ARG
1	A	95	LEU
1	A	121	THR
1	A	154	ARG
1	A	203	THR

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Mol	Chain	Res	Type
1	A	217	ARG
1	A	239	THR
1	A	261	ARG
1	A	268	ARG
1	A	292	THR
1	A	294	GLN
1	A	322	VAL
1	A	412	MET
1	A	417	LEU
1	A	426	LEU
1	A	448	GLU
1	A	494	SER
1	A	495	ARG
1	B	13	THR
1	B	32	LEU
1	B	37	LYS
1	B	46	ASN
1	B	58	GLU
1	B	71	LEU
1	B	82	ARG
1	B	95	LEU
1	B	96	LYS
1	B	121	THR
1	B	132	THR
1	B	147	LEU
1	B	203	THR
1	B	217	ARG
1	B	239	THR
1	B	268	ARG
1	B	292	THR
1	B	294	GLN
1	B	312	SER
1	B	389	THR
1	B	412	MET
1	B	417	LEU
1	B	426	LEU
1	B	459	LYS
1	B	463	ASN
1	B	475	GLU
1	B	491	GLN
1	B	494	SER
1	B	495	ARG

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Mol	Chain	Res	Type
1	C	37	LYS
1	C	58	GLU
1	C	71	LEU
1	C	82	ARG
1	C	95	LEU
1	C	132	THR
1	C	136	THR
1	C	147	LEU
1	C	154	ARG
1	C	203	THR
1	C	217	ARG
1	C	224	ARG
1	C	268	ARG
1	C	292	THR
1	C	294	GLN
1	C	312	SER
1	C	412	MET
1	C	417	LEU
1	C	426	LEU
1	C	449	LYS
1	C	454	LEU
1	C	463	ASN
1	C	494	SER
1	C	495	ARG
1	C	496	LEU
1	C	498	ARG
1	E	37	LYS
1	E	75	SER
1	E	82	ARG
1	E	94	VAL
1	E	95	LEU
1	E	132	THR
1	E	217	ARG
1	E	224	ARG
1	E	268	ARG
1	E	292	THR
1	E	294	GLN
1	E	322	VAL
1	E	395	GLU
1	E	412	MET
1	E	417	LEU
1	E	426	LEU

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Mol	Chain	Res	Type
1	E	448	GLU
1	E	449	LYS
1	E	463	ASN
1	E	475	GLU
1	E	491	GLN
1	E	494	SER
1	D	37	LYS
1	D	58	GLU
1	D	66	ARG
1	D	71	LEU
1	D	82	ARG
1	D	86	GLN
1	D	95	LEU
1	D	121	THR
1	D	147	LEU
1	D	190	ASN
1	D	194	LEU
1	D	206	MET
1	D	217	ARG
1	D	239	THR
1	D	268	ARG
1	D	294	GLN
1	D	330	LEU
1	D	334	ILE
1	D	412	MET
1	D	454	LEU
1	D	455	ARG
1	D	459	LYS
1	D	463	ASN
1	D	471	LYS
1	D	475	GLU
1	D	491	GLN
1	D	494	SER
1	F	5	ILE
1	F	13	THR
1	F	37	LYS
1	F	58	GLU
1	F	71	LEU
1	F	82	ARG
1	F	121	THR
1	F	147	LEU
1	F	154	ARG

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Mol	Chain	Res	Type
1	F	186	VAL
1	F	217	ARG
1	F	224	ARG
1	F	268	ARG
1	F	292	THR
1	F	294	GLN
1	F	312	SER
1	F	395	GLU
1	F	412	MET
1	F	417	LEU
1	F	431	GLU
1	F	438	MET
1	F	449	LYS
1	F	461	LEU
1	F	463	ASN
1	F	471	LYS
1	F	485	TYR
1	F	488	PRO
1	F	490	TYR
1	F	493	GLU
1	F	494	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	GLN
1	A	112	GLN
1	A	453	GLN
1	A	463	ASN
1	B	112	GLN
1	B	453	GLN
1	B	457	ASN
1	E	86	GLN
1	E	112	GLN
1	E	247	ASN
1	E	453	GLN
1	E	491	GLN
1	D	8	HIS
1	D	457	ASN
1	F	8	HIS
1	F	171	GLN
1	F	378	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](#)

17 non-standard protein/DNA/RNA residues 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	C	601	1	14,14,15	0.69	0	17,19,21	1.41	2 (11%)
4	NAG	B	602	1	14,14,15	1.05	1 (7%)	17,19,21	1.97	6 (35%)
3	NAG	I	1	3	14,14,15	0.86	1 (7%)	17,19,21	1.99	5 (29%)
4	NAG	B	601	1	14,14,15	0.43	0	17,19,21	1.67	4 (23%)
2	NAG	J	1	2,1	14,14,15	0.57	0	17,19,21	1.62	3 (17%)
4	NAG	E	601	1	14,14,15	1.52	2 (14%)	17,19,21	3.19	12 (70%)
4	NAG	D	900	1	14,14,15	0.84	0	17,19,21	1.60	3 (17%)
4	NAG	E	602	1	14,14,15	0.56	0	17,19,21	1.87	2 (11%)
2	NAG	G	2	2	14,14,15	0.54	0	17,19,21	1.31	3 (17%)
4	NAG	C	602	1	14,14,15	1.06	1 (7%)	17,19,21	2.13	6 (35%)
3	NAG	H	1	3	14,14,15	0.72	0	17,19,21	1.37	2 (11%)
3	NAG	K	1	3	14,14,15	0.90	1 (7%)	17,19,21	2.07	7 (41%)
2	NAG	G	1	2,1	14,14,15	0.86	1 (7%)	17,19,21	1.33	2 (11%)
3	GAL	I	2	3	11,11,12	0.89	0	15,15,17	2.13	5 (33%)
3	GAL	K	2	3	11,11,12	0.79	0	15,15,17	2.45	6 (40%)
3	GAL	H	2	3	11,11,12	0.88	0	15,15,17	1.94	3 (20%)
2	NAG	J	2	2	14,14,15	0.66	0	17,19,21	1.25	2 (11%)

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	C	601	1	-	0/6/23/26	0/1/1/1
4	NAG	B	602	1	-	0/6/23/26	0/1/1/1
3	NAG	I	1	3	-	2/6/23/26	0/1/1/1
4	NAG	B	601	1	1/1/5/7	1/6/23/26	0/1/1/1
4	NAG	D	900	1	1/1/5/7	2/6/23/26	0/1/1/1
4	NAG	E	601	1	1/1/5/7	2/6/23/26	0/1/1/1
2	NAG	J	1	2,1	-	0/6/23/26	0/1/1/1
4	NAG	E	602	1	-	0/6/23/26	0/1/1/1
4	NAG	C	602	1	1/1/5/7	2/6/23/26	0/1/1/1
2	NAG	G	2	2	-	0/6/23/26	0/1/1/1
3	NAG	H	1	3	-	0/6/23/26	0/1/1/1
3	NAG	K	1	3	-	2/6/23/26	0/1/1/1
2	NAG	G	1	2,1	-	1/6/23/26	0/1/1/1
3	GAL	I	2	3	-	0/2/19/22	0/1/1/1
3	GAL	K	2	3	-	2/2/19/22	0/1/1/1
3	GAL	H	2	3	-	0/2/19/22	0/1/1/1
2	NAG	J	2	2	-	0/6/23/26	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	601	NAG	C1-C2	4.53	1.59	1.52
4	B	602	NAG	C1-C2	2.65	1.56	1.52
3	I	1	NAG	C1-C2	2.49	1.56	1.52
2	G	1	NAG	C1-C2	2.38	1.55	1.52
4	E	601	NAG	C3-C2	2.28	1.57	1.52
4	C	602	NAG	C1-C2	2.20	1.55	1.52
3	K	1	NAG	C3-C2	2.16	1.57	1.52

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	602	NAG	C1-O5-C5	5.85	120.12	112.19
4	E	601	NAG	O5-C5-C6	5.64	116.05	107.20
4	E	601	NAG	C3-C4-C5	-5.13	101.08	110.24
4	E	601	NAG	O5-C5-C4	-5.09	98.44	110.83
3	K	2	GAL	O3-C3-C2	-4.71	100.97	109.99
3	H	2	GAL	O5-C1-C2	-4.52	103.79	110.77
3	K	2	GAL	O5-C1-C2	-4.39	104.00	110.77
4	C	602	NAG	O5-C5-C4	-4.33	100.28	110.83
3	K	2	GAL	O2-C2-C3	-4.29	101.55	110.14
3	I	1	NAG	O5-C5-C6	4.26	113.88	107.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	1	NAG	O5-C1-C2	-4.12	104.78	111.29
4	E	601	NAG	C1-O5-C5	4.05	117.68	112.19
3	I	1	NAG	O5-C1-C2	3.97	117.56	111.29
3	K	1	NAG	C4-C3-C2	3.93	116.78	111.02
3	I	2	GAL	O5-C1-C2	-3.93	104.70	110.77
4	B	601	NAG	O5-C5-C6	3.82	113.19	107.20
4	D	900	NAG	O5-C1-C2	-3.75	105.37	111.29
3	H	2	GAL	O3-C3-C2	-3.67	102.97	109.99
3	I	2	GAL	O5-C5-C6	3.66	112.94	107.20
3	I	2	GAL	O3-C3-C2	-3.55	103.20	109.99
4	E	601	NAG	C6-C5-C4	3.55	121.31	113.00
4	C	602	NAG	O5-C1-C2	-3.51	105.75	111.29
4	B	602	NAG	C4-C3-C2	3.45	116.08	111.02
4	C	602	NAG	C8-C7-N2	3.40	121.85	116.10
4	C	602	NAG	C1-C2-N2	3.22	115.99	110.49
4	E	601	NAG	O4-C4-C5	3.22	117.29	109.30
4	B	602	NAG	C2-N2-C7	3.17	127.42	122.90
2	J	1	NAG	C1-C2-N2	3.15	115.87	110.49
4	E	601	NAG	C1-C2-N2	3.13	115.83	110.49
4	B	601	NAG	C1-O5-C5	3.13	116.43	112.19
4	C	601	NAG	C1-O5-C5	3.10	116.39	112.19
3	H	1	NAG	O4-C4-C3	-3.09	103.20	110.35
4	C	602	NAG	O7-C7-C8	-3.04	116.42	122.06
4	B	602	NAG	C1-C2-N2	2.91	115.46	110.49
4	E	601	NAG	O7-C7-C8	-2.90	116.66	122.06
3	K	2	GAL	O2-C2-C1	2.89	115.06	109.15
4	B	602	NAG	O7-C7-C8	-2.88	116.71	122.06
3	K	2	GAL	O4-C4-C5	2.85	116.39	109.30
3	K	1	NAG	C1-O5-C5	2.85	116.05	112.19
4	D	900	NAG	C4-C3-C2	2.76	115.07	111.02
3	K	1	NAG	O5-C1-C2	2.76	115.65	111.29
4	E	601	NAG	O4-C4-C3	2.74	116.68	110.35
4	E	601	NAG	O3-C3-C2	2.73	115.11	109.47
3	I	1	NAG	C1-O5-C5	2.69	115.83	112.19
2	G	1	NAG	C4-C3-C2	2.62	114.86	111.02
4	C	602	NAG	C6-C5-C4	2.62	119.13	113.00
3	I	2	GAL	C1-C2-C3	-2.61	106.45	109.67
3	K	1	NAG	O3-C3-C2	2.60	114.85	109.47
4	B	602	NAG	O5-C5-C6	2.55	111.20	107.20
3	K	1	NAG	O7-C7-C8	-2.54	117.34	122.06
2	G	1	NAG	C1-C2-N2	2.54	114.83	110.49
2	G	2	NAG	C3-C4-C5	-2.53	105.73	110.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	602	NAG	O5-C1-C2	2.47	115.19	111.29
2	J	2	NAG	C1-C2-N2	2.45	114.68	110.49
3	H	1	NAG	C4-C3-C2	2.37	114.49	111.02
4	E	601	NAG	O3-C3-C4	2.35	115.79	110.35
4	B	601	NAG	O7-C7-C8	-2.33	117.73	122.06
3	H	2	GAL	O2-C2-C3	-2.32	105.48	110.14
2	G	2	NAG	O4-C4-C5	2.31	115.03	109.30
3	K	2	GAL	O5-C5-C6	2.30	110.80	107.20
3	I	2	GAL	O3-C3-C4	-2.27	105.09	110.35
2	J	2	NAG	C3-C4-C5	-2.27	106.19	110.24
3	I	1	NAG	O4-C4-C5	2.26	114.90	109.30
3	K	1	NAG	O4-C4-C5	2.24	114.87	109.30
4	B	601	NAG	C6-C5-C4	-2.24	107.75	113.00
2	G	2	NAG	O5-C5-C6	2.22	110.69	107.20
3	I	1	NAG	O3-C3-C2	2.22	114.05	109.47
4	C	601	NAG	O3-C3-C2	2.21	114.03	109.47
4	B	602	NAG	O4-C4-C3	2.18	115.38	110.35
2	J	1	NAG	O7-C7-C8	-2.16	118.05	122.06
3	K	1	NAG	O4-C4-C3	-2.15	105.37	110.35
4	D	900	NAG	O5-C5-C6	-2.07	103.96	107.20
4	E	601	NAG	C8-C7-N2	2.02	119.52	116.10

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	B	601	NAG	C1
4	C	602	NAG	C1
4	E	601	NAG	C1
4	D	900	NAG	C1

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	K	1	NAG	C4-C5-C6-O6
4	E	601	NAG	O5-C5-C6-O6
4	C	602	NAG	O5-C5-C6-O6
3	K	1	NAG	O5-C5-C6-O6
4	D	900	NAG	O5-C5-C6-O6
4	C	602	NAG	C4-C5-C6-O6
3	K	2	GAL	C4-C5-C6-O6
4	D	900	NAG	C4-C5-C6-O6
4	E	601	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
3	K	2	GAL	O5-C5-C6-O6
3	I	1	NAG	C4-C5-C6-O6
2	G	1	NAG	O5-C5-C6-O6
3	I	1	NAG	O5-C5-C6-O6
4	B	601	NAG	O5-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	601	NAG	1	0
4	E	601	NAG	2	0
4	C	602	NAG	1	0

5.5 Carbohydrates [i](#)

13 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	G	1	2,1	14,14,15	0.86	1 (7%)	17,19,21	1.33	2 (11%)
2	NAG	G	2	2	14,14,15	0.54	0	17,19,21	1.31	3 (17%)
3	NAG	H	1	3	14,14,15	0.72	0	17,19,21	1.37	2 (11%)
3	GAL	H	2	3	11,11,12	0.88	0	15,15,17	1.94	3 (20%)
3	SIA	H	3	3	20,20,21	0.96	2 (10%)	24,28,31	1.28	4 (16%)
3	NAG	I	1	3	14,14,15	0.86	1 (7%)	17,19,21	1.99	5 (29%)
3	GAL	I	2	3	11,11,12	0.89	0	15,15,17	2.13	5 (33%)
3	SIA	I	3	3	20,20,21	1.09	2 (10%)	24,28,31	1.79	6 (25%)
2	NAG	J	1	2,1	14,14,15	0.57	0	17,19,21	1.62	3 (17%)
2	NAG	J	2	2	14,14,15	0.66	0	17,19,21	1.25	2 (11%)
3	NAG	K	1	3	14,14,15	0.90	1 (7%)	17,19,21	2.07	7 (41%)
3	GAL	K	2	3	11,11,12	0.79	0	15,15,17	2.45	6 (40%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SIA	K	3	3	20,20,21	0.85	0	24,28,31	1.42	5 (20%)

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	G	1	2,1	-	1/6/23/26	0/1/1/1
2	NAG	G	2	2	-	0/6/23/26	0/1/1/1
3	NAG	H	1	3	-	0/6/23/26	0/1/1/1
3	GAL	H	2	3	-	0/2/19/22	0/1/1/1
3	SIA	H	3	3	-	2/18/34/38	0/1/1/1
3	NAG	I	1	3	-	2/6/23/26	0/1/1/1
3	GAL	I	2	3	-	0/2/19/22	0/1/1/1
3	SIA	I	3	3	-	1/18/34/38	0/1/1/1
2	NAG	J	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	J	2	2	-	0/6/23/26	0/1/1/1
3	NAG	K	1	3	-	2/6/23/26	0/1/1/1
3	GAL	K	2	3	-	2/2/19/22	0/1/1/1
3	SIA	K	3	3	-	0/18/34/38	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	3	SIA	C2-C1	-2.49	1.50	1.52
3	I	1	NAG	C1-C2	2.49	1.56	1.52
3	I	3	SIA	C5-N5	-2.47	1.41	1.45
2	G	1	NAG	C1-C2	2.38	1.55	1.52
3	K	1	NAG	C3-C2	2.16	1.57	1.52
3	I	3	SIA	O1B-C1	-2.16	1.23	1.30
3	H	3	SIA	O1A-C1	2.02	1.28	1.22

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	K	2	GAL	O3-C3-C2	-4.71	100.97	109.99
3	H	2	GAL	O5-C1-C2	-4.52	103.79	110.77
3	K	2	GAL	O5-C1-C2	-4.39	104.00	110.77
3	K	2	GAL	O2-C2-C3	-4.29	101.55	110.14
3	I	1	NAG	O5-C5-C6	4.26	113.88	107.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	1	NAG	O5-C1-C2	-4.12	104.78	111.29
3	I	3	SIA	C3-C4-C5	-4.08	106.53	111.46
3	I	1	NAG	O5-C1-C2	3.97	117.56	111.29
3	K	1	NAG	C4-C3-C2	3.93	116.78	111.02
3	I	2	GAL	O5-C1-C2	-3.93	104.70	110.77
3	H	2	GAL	O3-C3-C2	-3.67	102.97	109.99
3	I	2	GAL	O5-C5-C6	3.66	112.94	107.20
3	I	2	GAL	O3-C3-C2	-3.55	103.20	109.99
2	J	1	NAG	C1-C2-N2	3.15	115.87	110.49
3	H	1	NAG	O4-C4-C3	-3.09	103.20	110.35
3	I	3	SIA	O4-C4-C3	3.03	117.47	109.94
3	K	2	GAL	O2-C2-C1	2.89	115.06	109.15
3	I	3	SIA	C9-C8-C7	-2.89	106.15	112.41
3	K	2	GAL	O4-C4-C5	2.85	116.39	109.30
3	K	1	NAG	C1-O5-C5	2.85	116.05	112.19
3	H	3	SIA	C6-O6-C2	2.84	117.41	111.34
3	K	1	NAG	O5-C1-C2	2.76	115.65	111.29
3	K	3	SIA	C6-C5-N5	-2.72	106.39	110.91
3	K	3	SIA	O1B-C1-C2	2.70	120.74	113.03
3	I	1	NAG	C1-O5-C5	2.69	115.83	112.19
2	G	1	NAG	C4-C3-C2	2.62	114.86	111.02
3	I	3	SIA	O1B-C1-C2	2.62	120.51	113.03
3	I	2	GAL	C1-C2-C3	-2.61	106.45	109.67
3	K	1	NAG	O3-C3-C2	2.60	114.85	109.47
3	H	3	SIA	C9-C8-C7	-2.59	106.79	112.41
3	K	1	NAG	O7-C7-C8	-2.54	117.34	122.06
2	G	1	NAG	C1-C2-N2	2.54	114.83	110.49
2	G	2	NAG	C3-C4-C5	-2.53	105.73	110.24
3	I	3	SIA	O6-C2-C3	-2.50	107.02	110.46
3	K	3	SIA	C6-O6-C2	2.48	116.66	111.34
2	J	2	NAG	C1-C2-N2	2.45	114.68	110.49
3	H	3	SIA	O6-C2-C1	-2.42	102.95	107.70
3	H	1	NAG	C4-C3-C2	2.37	114.49	111.02
3	H	2	GAL	O2-C2-C3	-2.32	105.48	110.14
2	G	2	NAG	O4-C4-C5	2.31	115.03	109.30
3	K	2	GAL	O5-C5-C6	2.30	110.80	107.20
3	I	2	GAL	O3-C3-C4	-2.27	105.09	110.35
2	J	2	NAG	C3-C4-C5	-2.27	106.19	110.24
3	I	1	NAG	O4-C4-C5	2.26	114.90	109.30
3	H	3	SIA	O6-C2-C3	2.25	113.56	110.46
3	K	1	NAG	O4-C4-C5	2.24	114.87	109.30
2	G	2	NAG	O5-C5-C6	2.22	110.69	107.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	K	3	SIA	O1A-C1-C2	-2.22	117.32	122.57
3	I	1	NAG	O3-C3-C2	2.22	114.05	109.47
2	J	1	NAG	O7-C7-C8	-2.16	118.05	122.06
3	K	1	NAG	O4-C4-C3	-2.15	105.37	110.35
3	K	3	SIA	O9-C9-C8	-2.13	106.43	111.07
3	I	3	SIA	C4-C5-N5	-2.05	106.32	110.38

There are no chirality outliers.

All (10) torsion outliers are listed below:

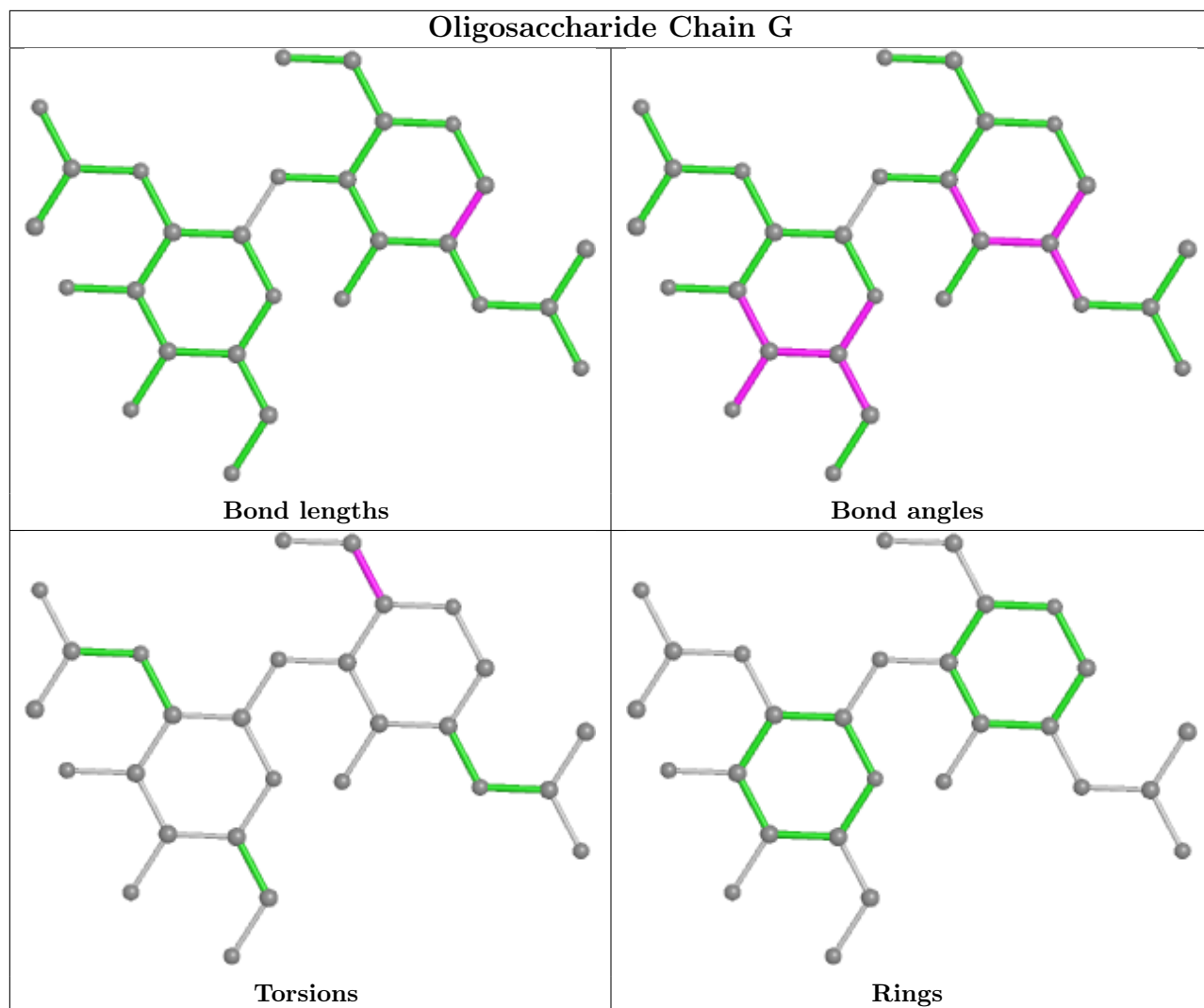
Mol	Chain	Res	Type	Atoms
3	K	1	NAG	C4-C5-C6-O6
3	K	1	NAG	O5-C5-C6-O6
3	K	2	GAL	C4-C5-C6-O6
3	K	2	GAL	O5-C5-C6-O6
3	I	1	NAG	C4-C5-C6-O6
3	H	3	SIA	O8-C8-C9-O9
2	G	1	NAG	O5-C5-C6-O6
3	I	1	NAG	O5-C5-C6-O6
3	I	3	SIA	O1A-C1-C2-O6
3	H	3	SIA	C7-C8-C9-O9

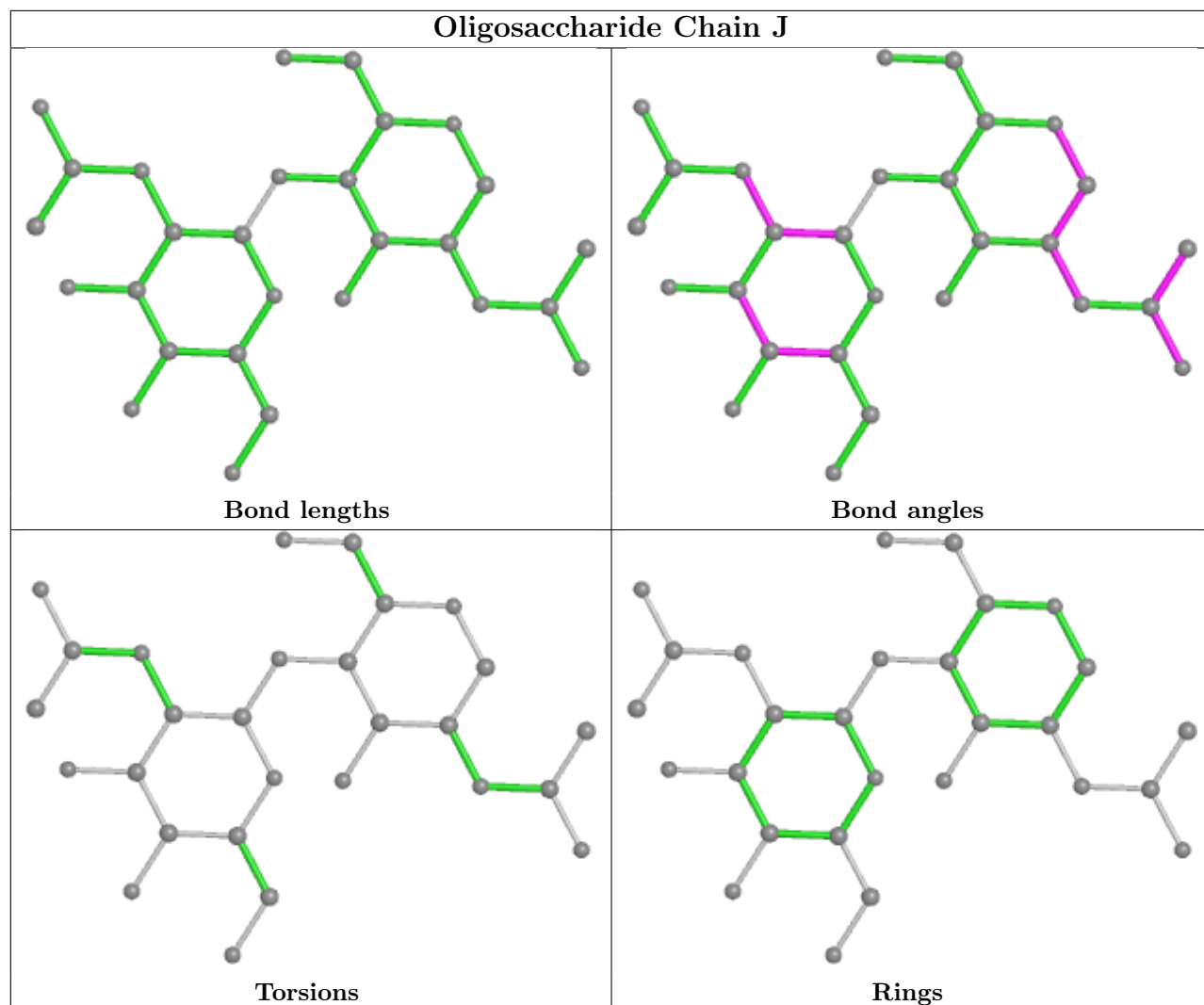
There are no ring outliers.

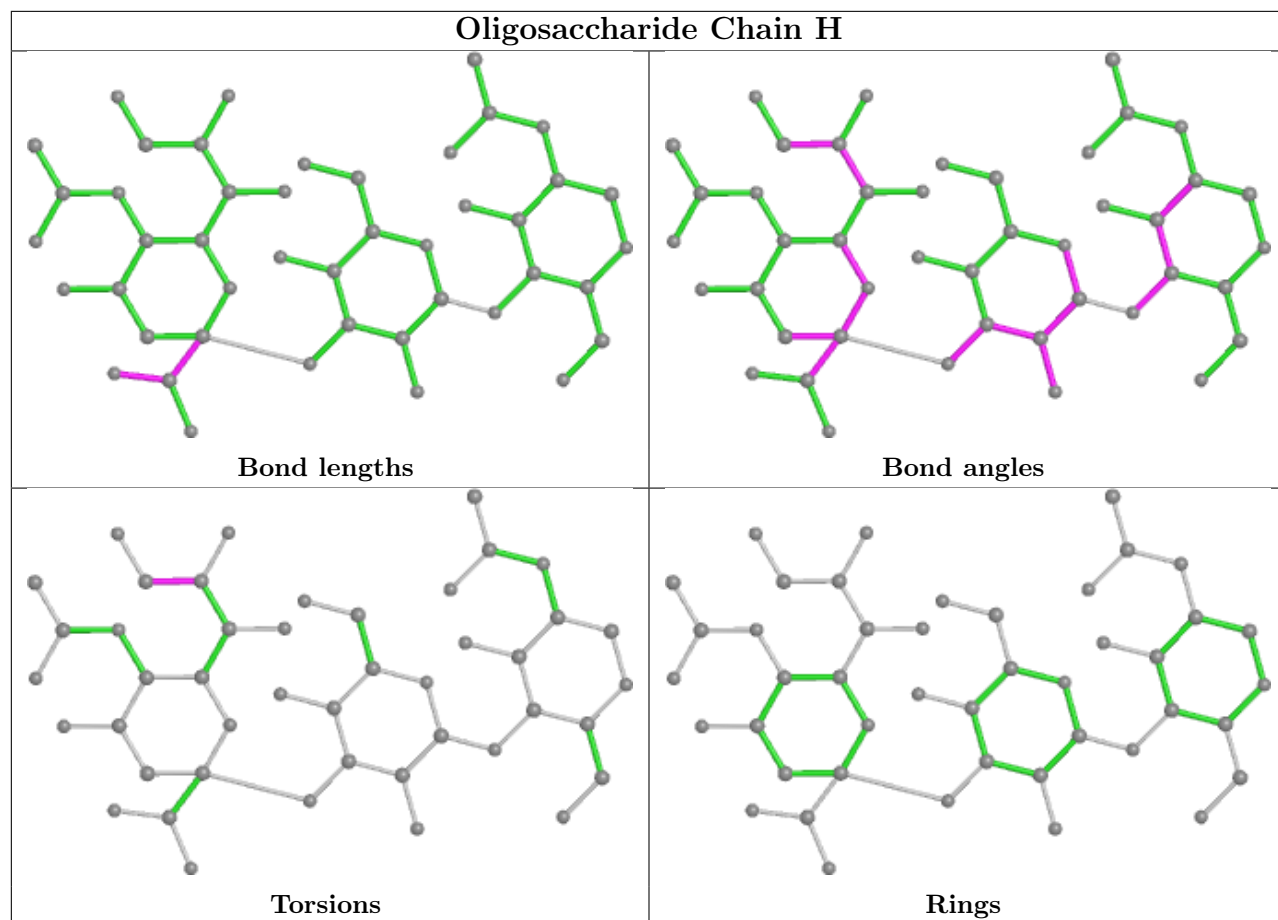
2 monomers are involved in 5 short contacts:

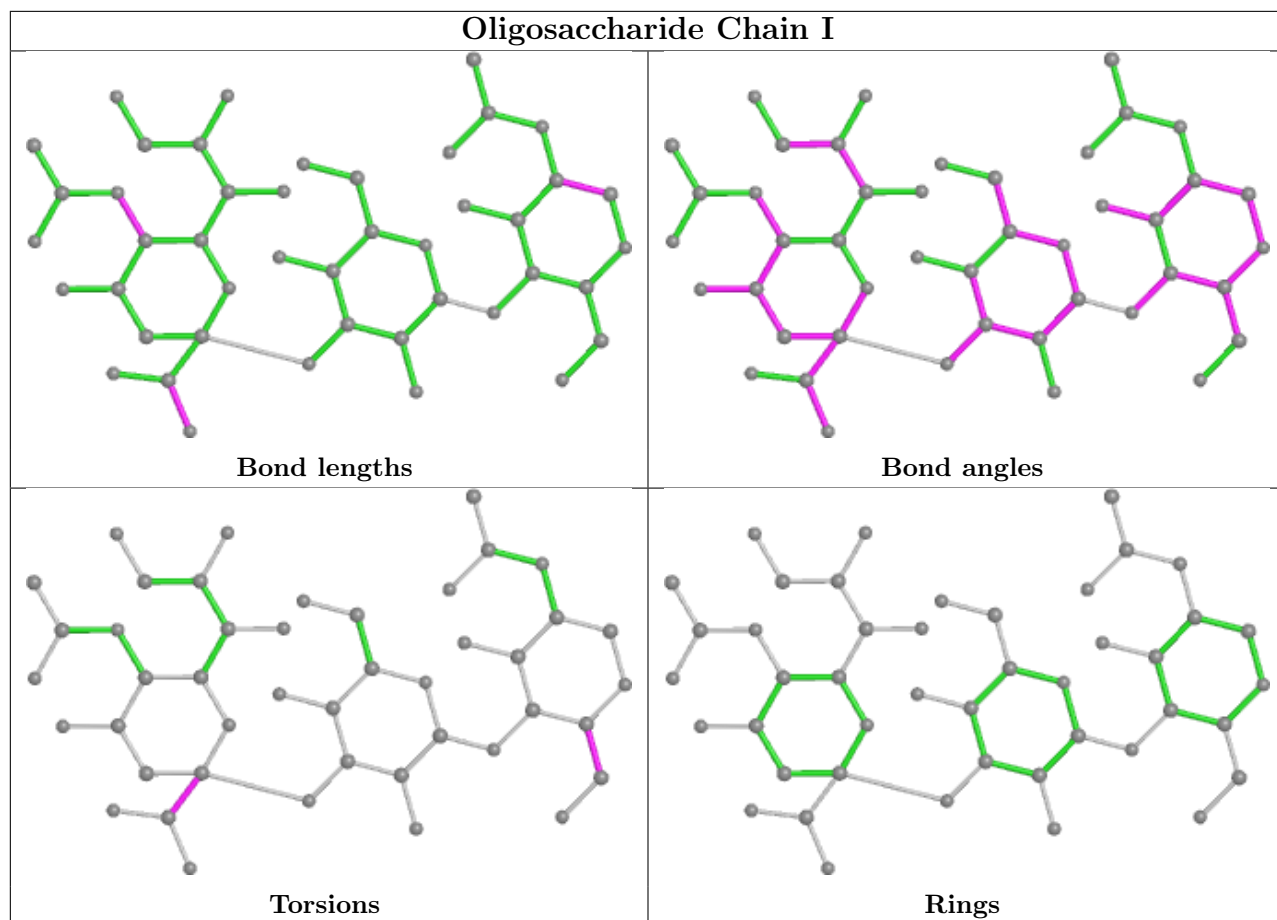
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	I	3	SIA	1	0
3	H	3	SIA	4	0

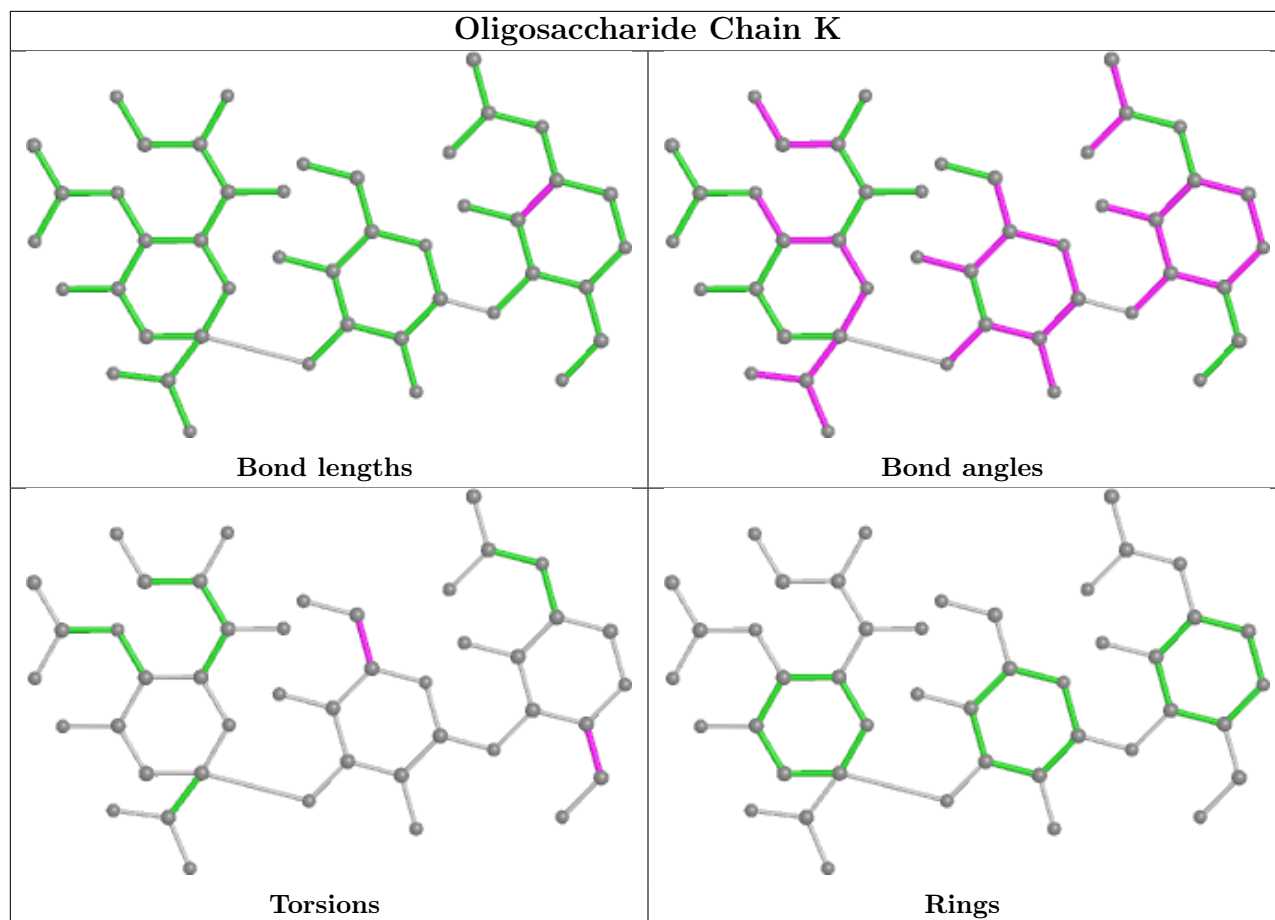
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











5.6 Ligand geometry [i](#)

7 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	C	601	1	14,14,15	0.69	0	17,19,21	1.41	2 (11%)
4	NAG	B	602	1	14,14,15	1.05	1 (7%)	17,19,21	1.97	6 (35%)
4	NAG	B	601	1	14,14,15	0.43	0	17,19,21	1.67	4 (23%)
4	NAG	D	900	1	14,14,15	0.84	0	17,19,21	1.60	3 (17%)
4	NAG	E	601	1	14,14,15	1.52	2 (14%)	17,19,21	3.19	12 (70%)
4	NAG	E	602	1	14,14,15	0.56	0	17,19,21	1.87	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	C	602	1	14,14,15	1.06	1 (7%)	17,19,21	2.13	6 (35%)

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	C	601	1	-	0/6/23/26	0/1/1/1
4	NAG	B	602	1	-	0/6/23/26	0/1/1/1
4	NAG	B	601	1	1/1/5/7	1/6/23/26	0/1/1/1
4	NAG	D	900	1	1/1/5/7	2/6/23/26	0/1/1/1
4	NAG	E	601	1	1/1/5/7	2/6/23/26	0/1/1/1
4	NAG	E	602	1	-	0/6/23/26	0/1/1/1
4	NAG	C	602	1	1/1/5/7	2/6/23/26	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	601	NAG	C1-C2	4.53	1.59	1.52
4	B	602	NAG	C1-C2	2.65	1.56	1.52
4	E	601	NAG	C3-C2	2.28	1.57	1.52
4	C	602	NAG	C1-C2	2.20	1.55	1.52

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	602	NAG	C1-O5-C5	5.85	120.12	112.19
4	E	601	NAG	O5-C5-C6	5.64	116.05	107.20
4	E	601	NAG	C3-C4-C5	-5.13	101.08	110.24
4	E	601	NAG	O5-C5-C4	-5.09	98.44	110.83
4	C	602	NAG	O5-C5-C4	-4.33	100.28	110.83
4	E	601	NAG	C1-O5-C5	4.05	117.68	112.19
4	B	601	NAG	O5-C5-C6	3.82	113.19	107.20
4	D	900	NAG	O5-C1-C2	-3.75	105.37	111.29
4	E	601	NAG	C6-C5-C4	3.55	121.31	113.00
4	C	602	NAG	O5-C1-C2	-3.51	105.75	111.29
4	B	602	NAG	C4-C3-C2	3.45	116.08	111.02
4	C	602	NAG	C8-C7-N2	3.40	121.85	116.10
4	C	602	NAG	C1-C2-N2	3.22	115.99	110.49
4	E	601	NAG	O4-C4-C5	3.22	117.29	109.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	602	NAG	C2-N2-C7	3.17	127.42	122.90
4	E	601	NAG	C1-C2-N2	3.13	115.83	110.49
4	B	601	NAG	C1-O5-C5	3.13	116.43	112.19
4	C	601	NAG	C1-O5-C5	3.10	116.39	112.19
4	C	602	NAG	O7-C7-C8	-3.04	116.42	122.06
4	B	602	NAG	C1-C2-N2	2.91	115.46	110.49
4	E	601	NAG	O7-C7-C8	-2.90	116.66	122.06
4	B	602	NAG	O7-C7-C8	-2.88	116.71	122.06
4	D	900	NAG	C4-C3-C2	2.76	115.07	111.02
4	E	601	NAG	O4-C4-C3	2.74	116.68	110.35
4	E	601	NAG	O3-C3-C2	2.73	115.11	109.47
4	C	602	NAG	C6-C5-C4	2.62	119.13	113.00
4	B	602	NAG	O5-C5-C6	2.55	111.20	107.20
4	E	602	NAG	O5-C1-C2	2.47	115.19	111.29
4	E	601	NAG	O3-C3-C4	2.35	115.79	110.35
4	B	601	NAG	O7-C7-C8	-2.33	117.73	122.06
4	B	601	NAG	C6-C5-C4	-2.24	107.75	113.00
4	C	601	NAG	O3-C3-C2	2.21	114.03	109.47
4	B	602	NAG	O4-C4-C3	2.18	115.38	110.35
4	D	900	NAG	O5-C5-C6	-2.07	103.96	107.20
4	E	601	NAG	C8-C7-N2	2.02	119.52	116.10

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	B	601	NAG	C1
4	C	602	NAG	C1
4	E	601	NAG	C1
4	D	900	NAG	C1

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	E	601	NAG	O5-C5-C6-O6
4	C	602	NAG	O5-C5-C6-O6
4	D	900	NAG	O5-C5-C6-O6
4	C	602	NAG	C4-C5-C6-O6
4	D	900	NAG	C4-C5-C6-O6
4	E	601	NAG	C4-C5-C6-O6
4	B	601	NAG	O5-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	601	NAG	1	0
4	E	601	NAG	2	0
4	C	602	NAG	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

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	490/514 (95%)	-0.17	9 (1%) 68 61	33, 52, 101, 128	0
1	B	481/514 (93%)	-0.12	12 (2%) 57 47	34, 56, 115, 140	0
1	C	482/514 (93%)	-0.20	11 (2%) 60 51	35, 52, 98, 133	0
1	D	493/514 (95%)	0.12	42 (8%) 10 5	32, 55, 160, 224	0
1	E	489/514 (95%)	-0.29	1 (0%) 95 94	36, 54, 76, 99	0
1	F	479/514 (93%)	0.34	62 (12%) 3 2	36, 58, 164, 200	0
All	All	2914/3084 (94%)	-0.05	137 (4%) 31 22	32, 54, 130, 224	0

All (137) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	357	SER	10.3
1	F	359	GLY	9.3
1	F	354	HIS	7.6
1	F	472	CYS	7.1
1	F	452	SER	6.9
1	F	360	SER	6.7
1	D	468	PHE	6.4
1	F	456	ASP	5.6
1	F	353	HIS	5.4
1	F	480	VAL	5.2
1	F	11	ASN	5.1
1	D	469	TRP	5.1
1	F	361	GLY	5.0
1	D	486	ASN	5.0
1	F	352	TYR	4.9
1	D	490	TYR	4.9
1	B	469	TRP	4.8
1	F	366	LYS	4.8
1	F	476	CYS	4.7

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Mol	Chain	Res	Type	RSRZ
1	F	483	GLY	4.7
1	F	488	PRO	4.7
1	D	485	TYR	4.7
1	D	354	HIS	4.7
1	D	352	TYR	4.5
1	D	3	ILE	4.5
1	F	347	ASP	4.5
1	D	456	ASP	4.4
1	F	355	GLU	4.4
1	D	454	LEU	4.3
1	B	359	GLY	4.2
1	F	2	LYS	4.2
1	D	459	LYS	4.2
1	A	481	LYS	4.1
1	F	317	THR	3.9
1	F	351	GLY	3.9
1	F	491	GLN	3.9
1	D	461	LEU	3.7
1	D	351	GLY	3.7
1	D	344	GLY	3.7
1	A	480	VAL	3.7
1	F	485	TYR	3.6
1	F	12	SER	3.6
1	D	357	SER	3.6
1	A	1	ASP	3.5
1	F	348	GLY	3.5
1	F	10	ASN	3.5
1	C	495	ARG	3.5
1	D	360	SER	3.5
1	B	496	LEU	3.4
1	F	16	VAL	3.4
1	D	362	TYR	3.4
1	D	466	PHE	3.4
1	F	342	TRP	3.4
1	B	456	ASP	3.3
1	D	368	SER	3.3
1	A	467	GLU	3.3
1	F	471	LYS	3.2
1	D	455	ARG	3.2
1	E	462	GLY	3.1
1	D	6	GLY	3.1
1	C	468	PHE	3.1

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Mol	Chain	Res	Type	RSRZ
1	F	346	VAL	3.1
1	F	29	SER	3.1
1	C	462	GLY	3.1
1	F	3	ILE	3.1
1	B	355	GLU	3.1
1	F	487	TYR	3.0
1	D	355	GLU	3.0
1	D	361	GLY	3.0
1	B	357	SER	3.0
1	F	358	GLN	3.0
1	C	359	GLY	3.0
1	B	468	PHE	3.0
1	F	13	THR	3.0
1	D	1	ASP	3.0
1	F	484	THR	2.9
1	A	352	TYR	2.9
1	A	485	TYR	2.9
1	F	343	THR	2.9
1	F	495	ARG	2.8
1	C	471	LYS	2.8
1	D	2	LYS	2.8
1	F	318	GLY	2.8
1	C	472	CYS	2.8
1	F	349	TRP	2.7
1	F	497	ASN	2.7
1	F	368	SER	2.7
1	D	363	ALA	2.7
1	F	455	ARG	2.7
1	F	475	GLU	2.7
1	D	458	ALA	2.6
1	F	14	THR	2.6
1	F	363	ALA	2.6
1	D	492	GLU	2.6
1	F	448	GLU	2.6
1	C	355	GLU	2.6
1	F	370	GLN	2.6
1	B	454	LEU	2.6
1	F	350	TYR	2.5
1	F	469	TRP	2.5
1	D	498	ARG	2.5
1	F	4	CYS	2.5
1	D	477	ILE	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	350	TYR	2.5
1	D	471	LYS	2.5
1	D	462	GLY	2.5
1	F	5	ILE	2.5
1	F	30	VAL	2.5
1	F	473	ASP	2.4
1	B	3	ILE	2.4
1	D	476	CYS	2.4
1	D	467	GLU	2.4
1	B	471	LYS	2.3
1	F	28	HIS	2.3
1	D	347	ASP	2.3
1	D	356	ASN	2.3
1	F	486	ASN	2.3
1	D	375	GLY	2.3
1	F	319	LEU	2.3
1	C	467	GLU	2.3
1	C	356	ASN	2.3
1	D	481	LYS	2.2
1	F	372	ALA	2.2
1	A	359	GLY	2.2
1	C	470	HIS	2.2
1	D	372	ALA	2.2
1	A	459	LYS	2.1
1	A	471	LYS	2.1
1	F	344	GLY	2.1
1	D	349	TRP	2.1
1	F	445	ASN	2.1
1	D	457	ASN	2.1
1	C	456	ASP	2.0
1	B	352	TYR	2.0
1	F	316	ALA	2.0
1	B	318	GLY	2.0
1	F	1	ASP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
4	NAG	B	601	14/15	0.74	0.26	106,117,129,130	0
4	NAG	C	601	14/15	0.76	0.25	104,117,132,133	0
4	NAG	D	900	14/15	0.77	0.19	101,115,122,126	0
3	NAG	K	1	14/15	0.78	0.20	87,98,110,110	0
4	NAG	E	601	14/15	0.78	0.25	55,85,93,97	0
4	NAG	E	602	14/15	0.79	0.21	97,114,122,122	0
2	NAG	J	2	14/15	0.82	0.38	88,109,115,115	0
3	NAG	I	1	14/15	0.84	0.17	85,102,118,127	0
4	NAG	B	602	14/15	0.84	0.25	68,75,78,81	0
4	NAG	C	602	14/15	0.85	0.32	57,80,94,100	0
2	NAG	G	2	14/15	0.86	0.29	86,101,109,110	0
3	NAG	H	1	14/15	0.87	0.16	73,94,107,113	0
2	NAG	J	1	14/15	0.90	0.33	82,87,99,100	0
2	NAG	G	1	14/15	0.91	0.20	68,84,92,94	0
3	GAL	K	2	11/12	0.93	0.09	57,65,74,77	0
3	GAL	I	2	11/12	0.94	0.09	59,65,74,76	0
3	GAL	H	2	11/12	0.96	0.08	53,57,65,66	0

6.3 Carbohydrates [i](#)

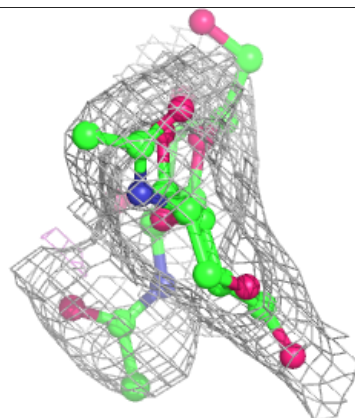
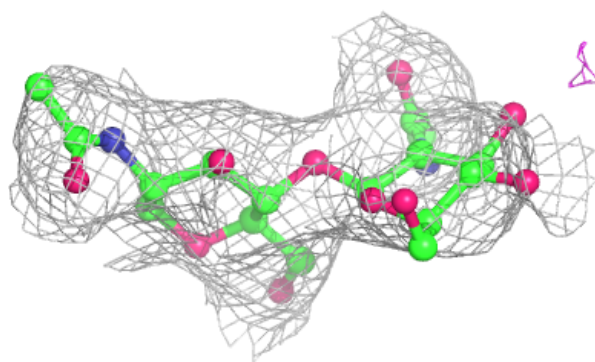
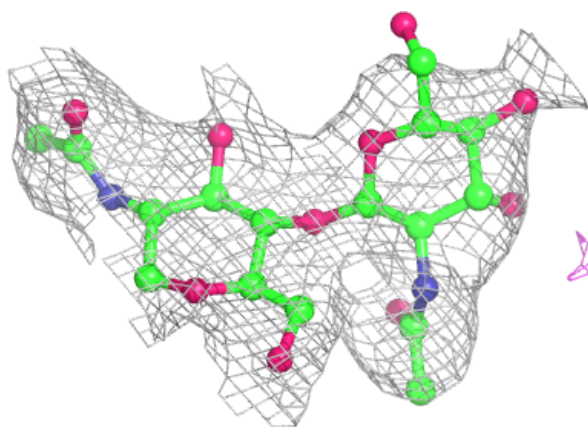
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	NAG	K	1	14/15	0.78	0.20	87,98,110,110	0
2	NAG	J	2	14/15	0.82	0.38	88,109,115,115	0
3	NAG	I	1	14/15	0.84	0.17	85,102,118,127	0
2	NAG	G	2	14/15	0.86	0.29	86,101,109,110	0
3	NAG	H	1	14/15	0.87	0.16	73,94,107,113	0
2	NAG	J	1	14/15	0.90	0.33	82,87,99,100	0
2	NAG	G	1	14/15	0.91	0.20	68,84,92,94	0
3	GAL	K	2	11/12	0.93	0.09	57,65,74,77	0
3	GAL	I	2	11/12	0.94	0.09	59,65,74,76	0
3	GAL	H	2	11/12	0.96	0.08	53,57,65,66	0
3	SIA	H	3	20/21	0.96	0.15	38,45,57,65	0
3	SIA	K	3	20/21	0.96	0.18	49,52,64,67	0
3	SIA	I	3	20/21	0.97	0.12	42,46,53,55	0

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

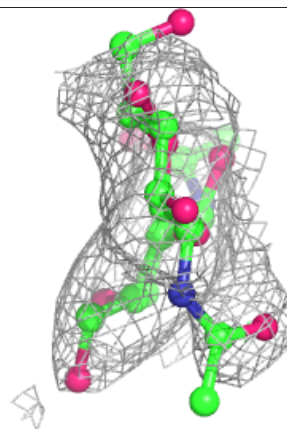
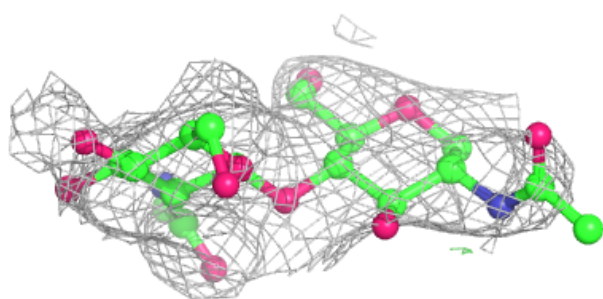
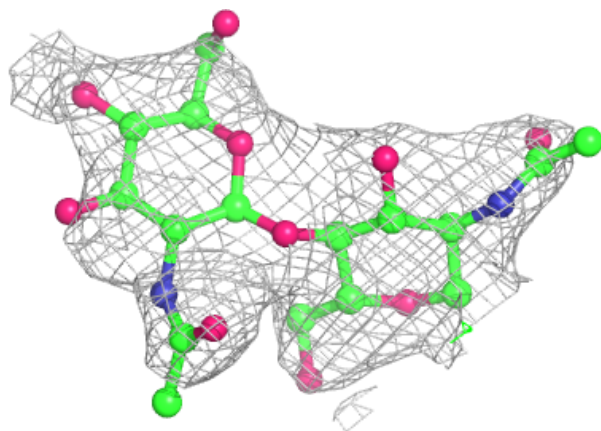
Electron density around Chain G:

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

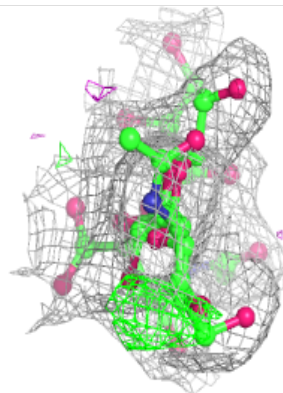
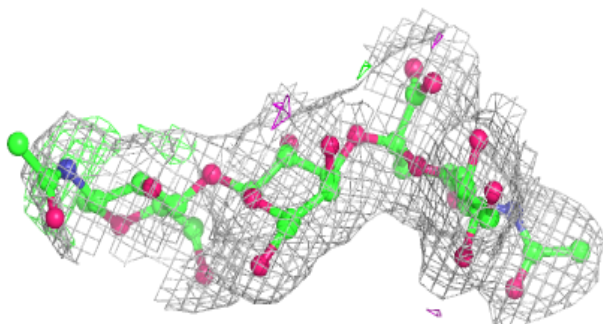
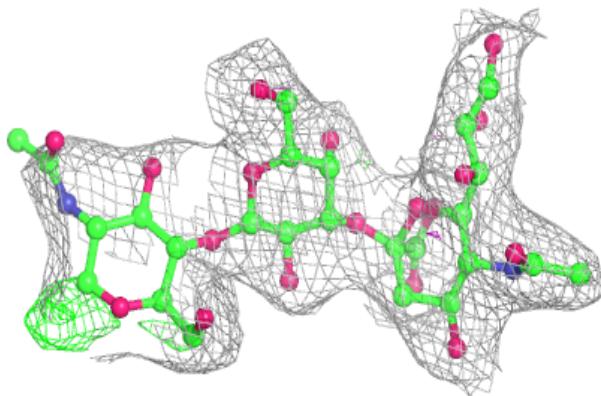


Electron density around Chain J:

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

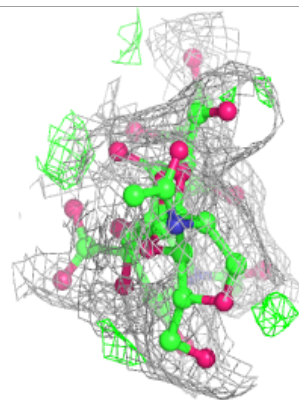
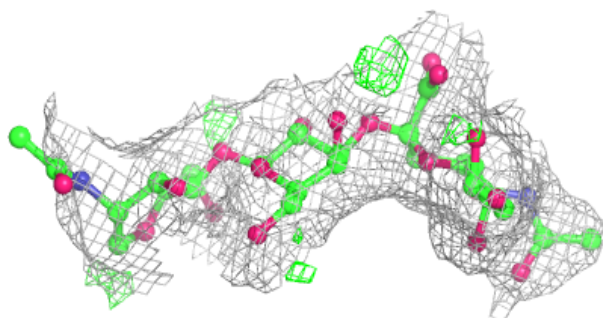
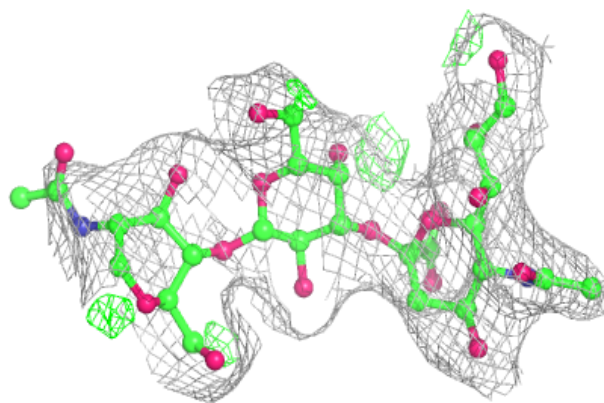
**Electron density around Chain H:**

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

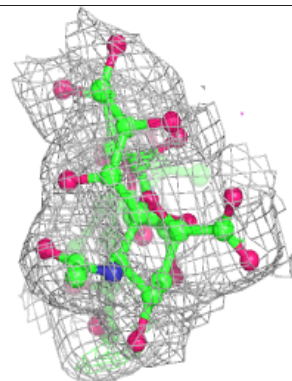
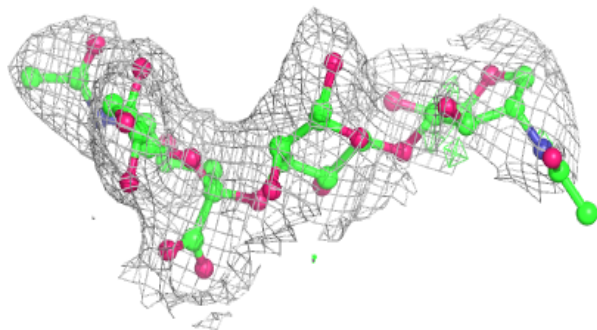
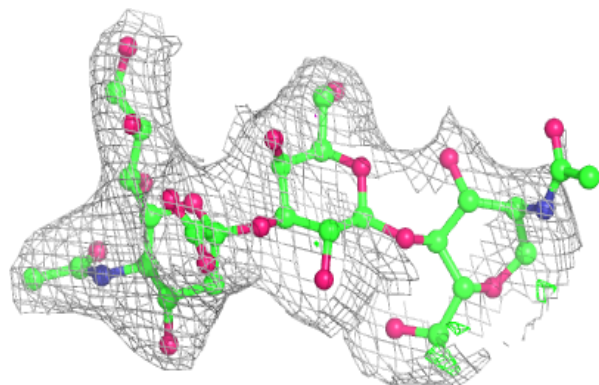


Electron density around Chain I:

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

**Electron density around Chain K:**

$2mF_o-DF_c$ (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
4	NAG	B	601	14/15	0.74	0.26	106,117,129,130	0
4	NAG	C	601	14/15	0.76	0.25	104,117,132,133	0
4	NAG	D	900	14/15	0.77	0.19	101,115,122,126	0
4	NAG	E	601	14/15	0.78	0.25	55,85,93,97	0
4	NAG	E	602	14/15	0.79	0.21	97,114,122,122	0
4	NAG	B	602	14/15	0.84	0.25	68,75,78,81	0
4	NAG	C	602	14/15	0.85	0.32	57,80,94,100	0

6.5 Other polymers [i](#)

There are no such residues in this entry.