



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 17, 2023 – 06:01 AM EDT

PDB ID : 2D3S
Title : Crystal Structure of basic winged bean lectin with Tn-antigen
Authors : Kulkarni, K.A.; Sinha, S.; Katiyar, S.; Surolia, A.; Vijayan, M.; Suguna, K.
Deposited on : 2005-10-01
Resolution : 2.35 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)
Xtrriage (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

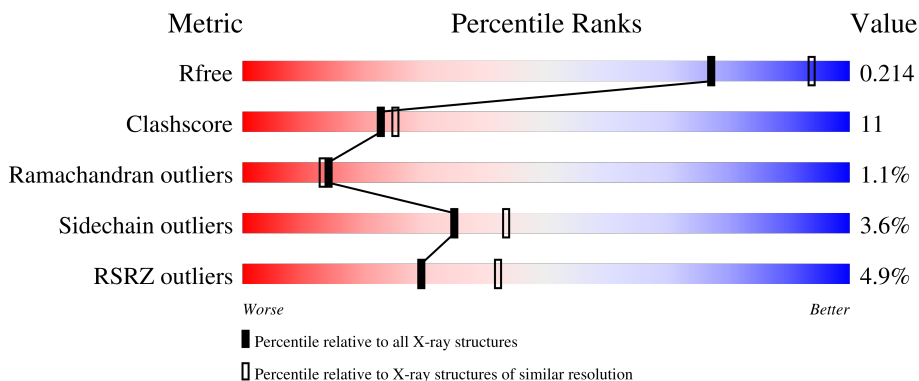
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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.



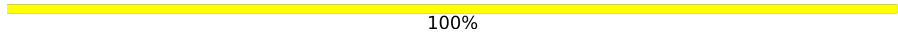

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	242	 4% 75% 19% ..
1	B	242	 2% 78% 19% ..
1	C	242	 5% 74% 21% ..
1	D	242	 8% 69% 26% ..
2	E	3	 33% 67%

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Mol	Chain	Length	Quality of chain
2	G	3	 67% 33%
2	H	3	 100%
2	J	3	 67% 33%
3	F	2	 50% 50%
4	I	4	 25% 75%

2 Entry composition [i](#)

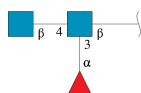
There are 10 unique types of molecules in this entry. The entry contains 7912 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Basic agglutinin.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
1	A	237	Total 1831	C 1179	N 307	O 345	0	0	0
1	B	237	Total 1839	C 1184	N 308	O 347	0	0	0
1	C	237	Total 1831	C 1177	N 307	O 347	0	0	0
1	D	237	Total 1807	C 1166	N 300	O 341	0	0	0

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



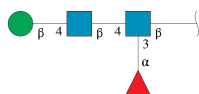
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	E	3	Total 38	C 22	N 2	O 14	0	0	0
2	G	3	Total 38	C 22	N 2	O 14	0	0	0
2	H	3	Total 38	C 22	N 2	O 14	0	0	0
2	J	3	Total 38	C 22	N 2	O 14	0	0	0

- Molecule 3 is an oligosaccharide called alpha-L-fucopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	F	2	24	14	1	9	0	0	0

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
4	I	4	49	28	2	19	0	0	0

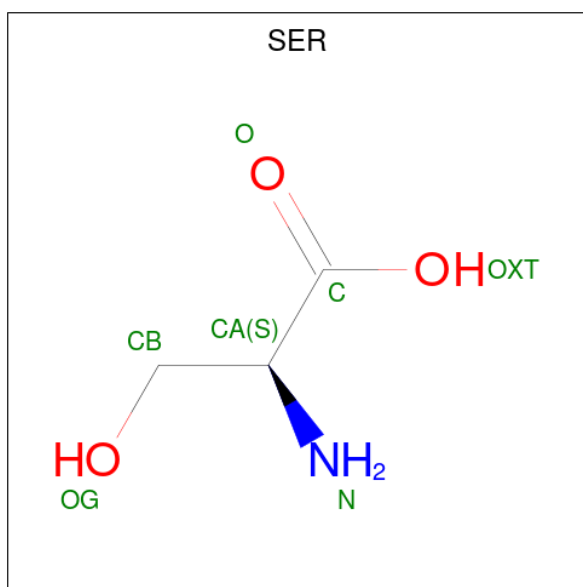
- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 6 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

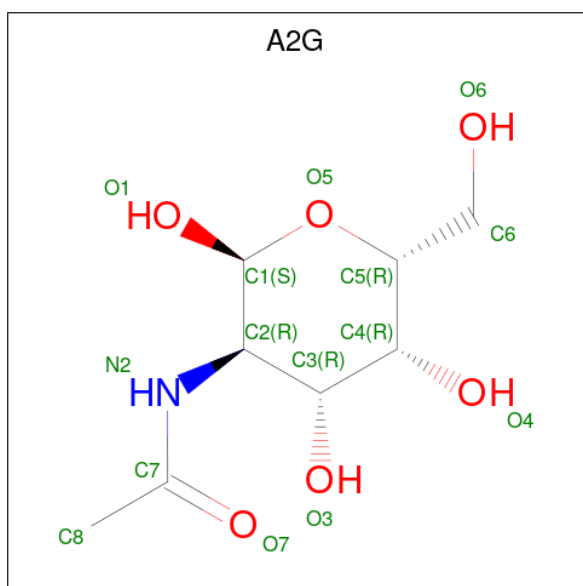
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Mn	0	0
			1	1		
6	B	1	Total	Mn	0	0
			1	1		
6	C	1	Total	Mn	0	0
			1	1		
6	D	1	Total	Mn	0	0
			1	1		

- Molecule 7 is SERINE (three-letter code: SER) (formula: C₃H₇NO₃).



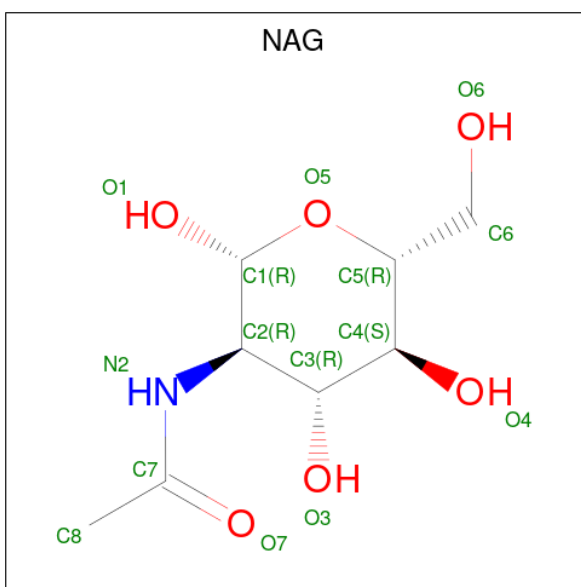
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
7	A	1	Total	C	N	O	0	0
			7	3	1	3		
7	B	1	Total	C	N	O	0	0
			7	3	1	3		
7	C	1	Total	C	N	O	0	0
			7	3	1	3		
7	D	1	Total	C	N	O	0	0
			7	3	1	3		

- Molecule 8 is 2-acetamido-2-deoxy-alpha-D-galactopyranose (three-letter code: A2G) (formula: C₈H₁₅NO₆).



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

- Molecule 9 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	C	1	Total	C	N	O	0	0
			14	8	1	5		

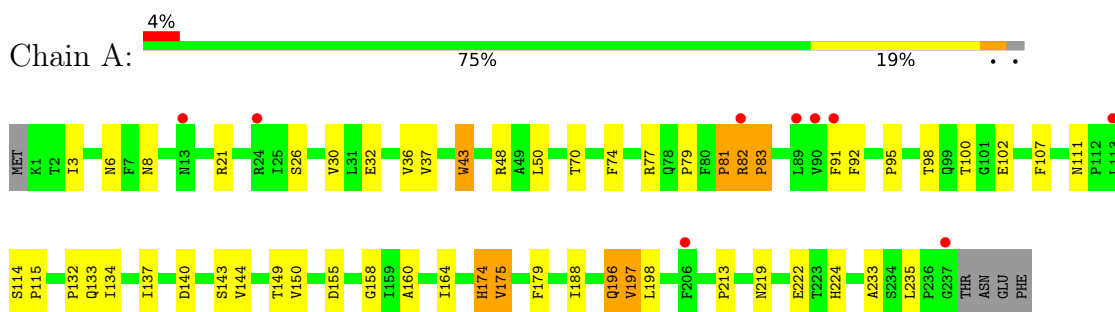
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	86	Total	O	0	0
			86	86		
10	B	83	Total	O	0	0
			83	83		
10	C	53	Total	O	0	0
			53	53		
10	D	51	Total	O	0	0
			51	51		

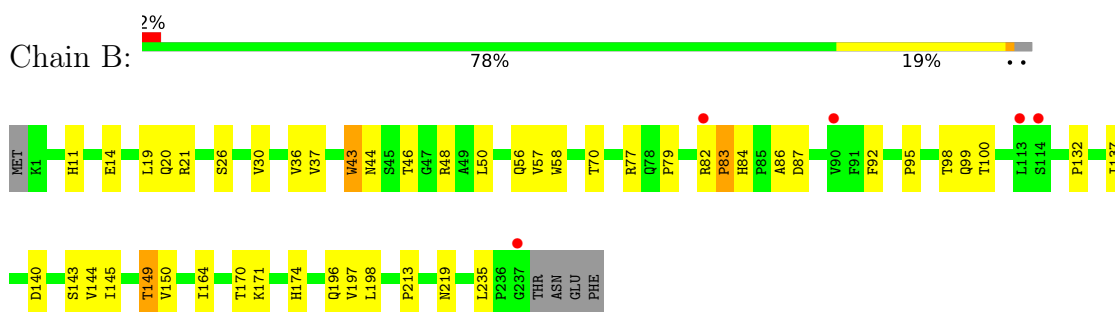
3 Residue-property plots [i](#)

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.

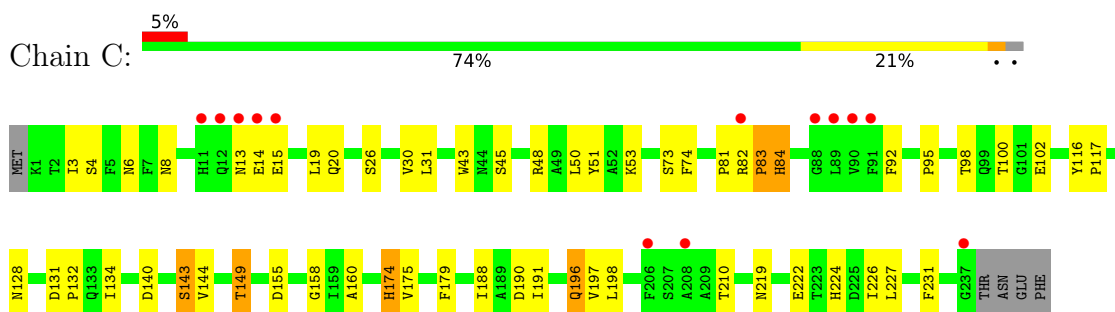
- Molecule 1: Basic agglutinin



- Molecule 1: Basic agglutinin

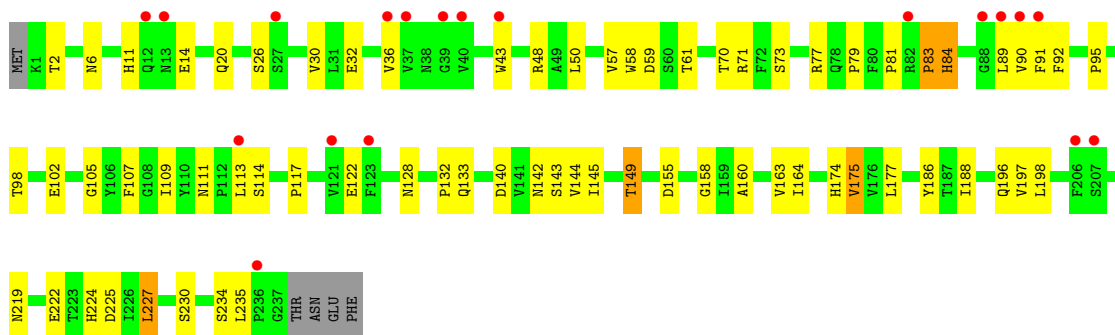


- Molecule 1: Basic agglutinin



- Molecule 1: Basic agglutinin





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

Chain E: 33% 67%



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

Chain G: 67% 33%



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

Chain H: 100%



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

Chain J: 67% 33%

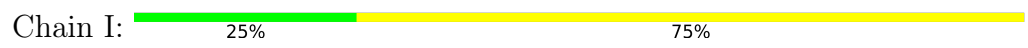


- Molecule 3: alpha-L-fucopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F: 50% 50%



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



MAG1
MAG2
BMA3
FUC4

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	157.87Å 91.90Å 73.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.35 28.56 – 2.34	Depositor EDS
% Data completeness (in resolution range)	97.5 (30.00-2.35) 96.8 (28.56-2.34)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.46 (at 2.34Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.198 , 0.244 0.213 , 0.214	Depositor DCC
R_{free} test set	1802 reflections (4.05%)	wwPDB-VP
Wilson B-factor (Å ²)	39.7	Xtrriage
Anisotropy	0.547	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 46.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7912	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: A2G, FUC, MN, NAG, BMA, CA

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	0.39	0/1884	0.69	1/2580 (0.0%)
1	B	0.40	0/1892	0.70	0/2589
1	C	0.38	0/1884	0.65	0/2579
1	D	0.36	0/1859	0.65	1/2548 (0.0%)
All	All	0.38	0/7519	0.67	2/10296 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	91	PHE	N-CA-C	-5.59	95.90	111.00
1	D	91	PHE	N-CA-C	-5.32	96.64	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1831	0	1777	39	0
1	B	1839	0	1791	36	0
1	C	1831	0	1770	41	0
1	D	1807	0	1741	50	0
2	E	38	0	34	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	G	38	0	34	2	0
2	H	38	0	34	6	0
2	J	38	0	34	0	0
3	F	24	0	22	5	0
4	I	49	0	43	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
6	C	1	0	0	0	0
6	D	1	0	0	0	0
7	A	7	0	3	0	0
7	B	7	0	3	0	0
7	C	7	0	3	0	0
7	D	7	0	3	0	0
8	A	14	0	12	0	0
8	B	14	0	12	0	0
8	C	14	0	12	0	0
8	D	14	0	12	0	0
9	C	14	0	13	0	0
10	A	86	0	0	3	0
10	B	83	0	0	2	0
10	C	53	0	0	1	0
10	D	51	0	0	1	0
All	All	7912	0	7353	172	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

All (172) 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:C:83:PRO:HG2	1:C:219:ASN:HB2	1.52	0.92
1:D:197:VAL:HG23	1:D:198:LEU:HG	1.53	0.90
1:B:83:PRO:HG2	1:B:219:ASN:HB2	1.55	0.87
3:F:1:NAG:H3	3:F:2:FUC:H5	1.58	0.86
2:H:1:NAG:H4	2:H:3:NAG:HN2	1.40	0.84
1:D:57:VAL:HG23	1:D:58:TRP:HD1	1.46	0.81
1:C:6:ASN:HD21	1:C:8:ASN:ND2	1.84	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:132:PRO:HG3	1:B:149:THR:HG21	1.69	0.74
2:H:1:NAG:H4	2:H:3:NAG:N2	2.03	0.73
1:C:197:VAL:HG23	1:C:198:LEU:HG	1.70	0.73
1:D:132:PRO:HG3	1:D:149:THR:HG21	1.70	0.72
1:B:145:ILE:HB	1:C:196:GLN:HE22	1.54	0.72
1:D:83:PRO:HG2	1:D:219:ASN:CB	2.20	0.71
3:F:1:NAG:O4	3:F:2:FUC:H3	1.91	0.71
1:C:132:PRO:HG3	1:C:149:THR:HG21	1.75	0.68
1:C:50:LEU:HD11	1:C:92:PHE:CZ	2.29	0.67
1:B:48:ARG:HD2	1:B:100:THR:OG1	1.95	0.67
2:H:2:FUC:H5	2:H:3:NAG:C1	2.24	0.67
1:D:175:VAL:HG13	1:D:188:ILE:HG23	1.76	0.66
1:C:48:ARG:HD2	1:C:100:THR:OG1	1.96	0.65
1:D:57:VAL:HG23	1:D:58:TRP:CD1	2.32	0.64
1:A:132:PRO:HG3	1:A:149:THR:HG21	1.78	0.64
1:D:83:PRO:HG2	1:D:219:ASN:HB2	1.80	0.63
1:D:155:ASP:OD2	1:D:158:GLY:HA3	1.99	0.62
1:A:197:VAL:CG2	1:A:198:LEU:HG	2.29	0.61
1:D:59:ASP:OD2	1:D:61:THR:HB	2.00	0.61
1:C:6:ASN:HD21	1:C:8:ASN:HD21	1.45	0.61
1:C:140:ASP:HB3	1:C:143:SER:O	2.00	0.61
1:D:32:GLU:OE2	1:D:36:VAL:HG11	2.02	0.60
1:B:140:ASP:HB3	1:B:143:SER:O	2.02	0.59
1:D:11:HIS:O	1:D:14:GLU:HB2	2.04	0.58
1:B:197:VAL:HG23	1:B:198:LEU:HG	1.85	0.58
1:B:82:ARG:HB2	10:B:647:HOH:O	2.04	0.57
1:C:50:LEU:HD11	1:C:92:PHE:HZ	1.67	0.57
1:A:26:SER:OG	1:A:30:VAL:HG22	2.05	0.57
1:D:113:LEU:HD22	1:D:113:LEU:H	1.69	0.57
1:B:143:SER:OG	1:C:196:GLN:NE2	2.38	0.56
1:C:26:SER:OG	1:C:30:VAL:HG22	2.05	0.56
1:D:26:SER:OG	1:D:30:VAL:HG22	2.04	0.56
1:D:71:ARG:HG2	1:D:163:VAL:HG22	1.87	0.56
1:A:83:PRO:HG2	1:A:219:ASN:HB2	1.87	0.56
1:B:50:LEU:HD11	1:B:92:PHE:CZ	2.41	0.56
1:A:36:VAL:O	1:A:36:VAL:HG13	2.05	0.56
1:D:2:THR:HG22	1:D:234:SER:CB	2.36	0.56
1:D:2:THR:HG22	1:D:234:SER:HB3	1.86	0.56
1:A:82:ARG:HG3	10:A:645:HOH:O	2.05	0.55
1:B:11:HIS:O	1:B:14:GLU:HG2	2.07	0.55
1:D:50:LEU:HD11	1:D:92:PHE:CZ	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:197:VAL:HG23	1:A:198:LEU:HG	1.90	0.54
1:A:6:ASN:HD21	1:A:8:ASN:ND2	2.05	0.54
1:A:175:VAL:HG13	1:A:188:ILE:HG23	1.90	0.54
1:C:31:LEU:HB3	1:C:226:ILE:HB	1.89	0.53
1:B:83:PRO:HG2	1:B:219:ASN:CB	2.33	0.53
1:D:90:VAL:HG21	1:D:109:ILE:HD13	1.91	0.53
3:F:1:NAG:H3	3:F:2:FUC:C5	2.33	0.53
1:B:144:VAL:HG12	1:B:144:VAL:O	2.09	0.53
1:D:140:ASP:HB3	1:D:143:SER:O	2.08	0.53
1:C:222:GLU:OE2	1:C:224:HIS:NE2	2.39	0.53
1:D:175:VAL:HG13	1:D:188:ILE:CG2	2.39	0.53
1:D:177:LEU:HB3	1:D:186:TYR:HB2	1.90	0.53
1:A:160:ALA:HB2	1:A:179:PHE:CE1	2.44	0.52
1:D:50:LEU:HD11	1:D:92:PHE:HZ	1.74	0.52
1:D:77:ARG:O	1:D:79:PRO:HD3	2.10	0.52
1:C:14:GLU:O	1:C:15:GLU:HB2	2.09	0.52
1:C:155:ASP:OD2	1:C:158:GLY:HA3	2.09	0.52
1:A:70:THR:HG22	1:A:164:ILE:HB	1.91	0.51
1:A:81:PRO:HG2	1:A:82:ARG:H	1.75	0.51
1:C:83:PRO:HG2	1:C:219:ASN:CB	2.33	0.51
1:B:77:ARG:O	1:B:79:PRO:HD3	2.11	0.51
1:B:70:THR:HG22	1:B:164:ILE:HB	1.93	0.51
1:A:82:ARG:HB2	10:A:655:HOH:O	2.11	0.50
1:C:95:PRO:HD3	1:C:117:PRO:O	2.12	0.50
1:B:95:PRO:O	1:B:98:THR:HG23	2.11	0.50
1:D:105:GLY:HA3	1:D:128:ASN:HD21	1.77	0.50
1:D:83:PRO:HG2	1:D:219:ASN:HB3	1.94	0.50
1:D:36:VAL:HG13	1:D:36:VAL:O	2.12	0.50
1:D:84:HIS:O	1:D:84:HIS:CD2	2.65	0.49
1:B:56:GLN:HG2	1:B:235:LEU:HD12	1.94	0.49
1:C:84:HIS:HD2	1:C:84:HIS:O	1.94	0.49
1:D:95:PRO:HD3	1:D:117:PRO:O	2.13	0.49
1:C:174:HIS:N	1:C:174:HIS:CD2	2.80	0.49
1:C:73:SER:HA	1:C:160:ALA:O	2.11	0.49
1:D:6:ASN:HD22	1:D:230:SER:HB3	1.78	0.48
1:A:83:PRO:HG2	1:A:219:ASN:CB	2.43	0.48
1:A:144:VAL:HG12	1:A:144:VAL:O	2.12	0.48
1:B:37:VAL:O	1:B:37:VAL:HG23	2.13	0.48
1:C:84:HIS:O	1:C:84:HIS:CD2	2.67	0.48
1:C:128:ASN:HB2	1:C:131:ASP:OD2	2.13	0.48
1:D:32:GLU:HG2	1:D:225:ASP:OD1	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:73:SER:HA	1:D:160:ALA:O	2.13	0.48
1:A:50:LEU:HD11	1:A:92:PHE:CZ	2.49	0.47
1:B:84:HIS:O	1:B:84:HIS:HD2	1.97	0.47
1:A:155:ASP:OD2	1:A:158:GLY:HA3	2.15	0.47
1:C:20:GLN:OE1	1:C:48:ARG:HD3	2.14	0.47
1:D:84:HIS:O	1:D:84:HIS:HD2	1.97	0.47
1:A:137:ILE:HB	1:A:150:VAL:HG12	1.96	0.47
1:B:83:PRO:HG3	2:G:1:NAG:C7	2.45	0.47
1:C:4:SER:HA	1:C:231:PHE:O	2.15	0.47
1:C:19:LEU:N	1:C:19:LEU:HD12	2.29	0.47
1:C:174:HIS:CD2	1:C:174:HIS:H	2.33	0.46
1:B:84:HIS:O	1:B:84:HIS:CD2	2.68	0.46
1:C:45:SER:HB3	1:C:210:THR:OG1	2.16	0.46
1:A:50:LEU:HD22	1:A:50:LEU:N	2.31	0.46
1:B:20:GLN:OE1	1:B:48:ARG:HD3	2.15	0.46
1:B:170:THR:O	1:B:171:LYS:HB2	2.16	0.46
1:C:51:TYR:CE2	1:C:53:LYS:HB2	2.52	0.45
1:D:144:VAL:HG13	10:D:704:HOH:O	2.15	0.45
1:A:144:VAL:O	1:A:144:VAL:CG1	2.64	0.45
1:C:95:PRO:O	1:C:98:THR:HG23	2.16	0.45
1:A:43:TRP:CZ3	1:A:213:PRO:HA	2.51	0.45
1:D:122:GLU:OE1	1:D:140:ASP:OD2	2.35	0.45
1:A:174:HIS:CD2	1:A:174:HIS:N	2.85	0.45
1:B:26:SER:OG	1:B:30:VAL:HG22	2.16	0.45
1:A:213:PRO:HG3	3:F:1:NAG:H82	1.98	0.45
1:D:144:VAL:O	1:D:144:VAL:CG1	2.65	0.45
1:A:222:GLU:OE2	1:A:224:HIS:NE2	2.44	0.45
1:B:144:VAL:O	1:B:144:VAL:CG1	2.65	0.45
1:B:83:PRO:CG	1:B:219:ASN:HD22	2.29	0.44
1:C:175:VAL:HG22	1:C:188:ILE:HG22	1.99	0.44
1:A:77:ARG:O	1:A:79:PRO:HD3	2.18	0.44
1:B:83:PRO:HG3	2:G:1:NAG:C8	2.47	0.44
1:B:99:GLN:HB2	10:B:641:HOH:O	2.17	0.44
1:C:227:LEU:N	1:C:227:LEU:HD12	2.33	0.44
1:A:134:ILE:O	1:A:134:ILE:HD12	2.18	0.44
10:C:640:HOH:O	1:D:174:HIS:HB2	2.18	0.44
2:H:3:NAG:O3	2:H:3:NAG:C7	2.66	0.44
1:C:3:ILE:O	1:C:3:ILE:HG23	2.18	0.44
1:A:37:VAL:O	1:A:37:VAL:HG23	2.18	0.43
1:B:56:GLN:HG2	1:B:235:LEU:CD1	2.48	0.43
1:D:227:LEU:CD1	1:D:227:LEU:N	2.81	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:160:ALA:HB2	1:C:179:PHE:CE1	2.53	0.43
1:A:32:GLU:OE2	1:A:36:VAL:HG11	2.19	0.43
1:A:213:PRO:CG	3:F:1:NAG:H82	2.49	0.43
1:B:21:ARG:HB3	1:B:46:THR:O	2.19	0.43
1:D:70:THR:HG22	1:D:164:ILE:HB	2.01	0.43
1:A:95:PRO:O	1:A:98:THR:HG23	2.19	0.43
1:C:74:PHE:C	1:C:74:PHE:CD1	2.92	0.43
1:D:89:LEU:C	1:D:89:LEU:HD12	2.38	0.43
1:D:20:GLN:OE1	1:D:48:ARG:NH1	2.52	0.43
1:B:83:PRO:HG2	1:B:219:ASN:HD22	1.83	0.42
1:B:44:ASN:HB2	1:B:213:PRO:HG3	2.01	0.42
1:B:137:ILE:HB	1:B:150:VAL:HG12	2.00	0.42
1:B:86:ALA:HA	1:B:87:ASP:HA	1.84	0.42
1:C:134:ILE:HD12	1:C:134:ILE:O	2.19	0.42
1:A:74:PHE:C	1:A:74:PHE:CD1	2.92	0.42
1:C:81:PRO:HG2	1:C:82:ARG:H	1.84	0.42
1:D:117:PRO:HA	1:D:142:ASN:OD1	2.20	0.42
1:D:113:LEU:HD22	1:D:113:LEU:N	2.35	0.42
1:A:114:SER:N	1:A:115:PRO:HD3	2.35	0.41
1:A:196:GLN:OE1	1:D:145:ILE:HB	2.20	0.41
1:D:95:PRO:O	1:D:98:THR:HG23	2.20	0.41
1:A:3:ILE:HG22	1:A:233:ALA:HB3	2.02	0.41
1:A:134:ILE:HD12	1:A:134:ILE:C	2.41	0.41
1:A:48:ARG:HD2	1:A:100:THR:OG1	2.21	0.41
2:H:1:NAG:C4	2:H:3:NAG:HN2	2.22	0.41
1:A:140:ASP:HB3	1:A:143:SER:O	2.20	0.41
1:A:21:ARG:HD3	10:A:639:HOH:O	2.20	0.41
1:A:111:ASN:ND2	1:A:114:SER:OG	2.54	0.41
1:C:134:ILE:HD12	1:C:134:ILE:C	2.41	0.41
1:D:79:PRO:O	1:D:81:PRO:HD3	2.20	0.41
2:H:1:NAG:C4	2:H:3:NAG:N2	2.80	0.41
1:B:43:TRP:CZ3	1:B:213:PRO:HA	2.56	0.40
1:B:57:VAL:HG23	1:B:58:TRP:HD1	1.86	0.40
1:B:82:ARG:HA	1:B:82:ARG:HD3	1.81	0.40
1:C:190:ASP:OD1	1:C:191:ILE:N	2.54	0.40
1:D:111:ASN:ND2	1:D:114:SER:H	2.19	0.40
1:C:144:VAL:O	1:C:144:VAL:CG1	2.69	0.40
1:D:222:GLU:OE2	1:D:224:HIS:NE2	2.52	0.40
1:C:98:THR:HB	1:C:116:TYR:CE2	2.55	0.40
1:D:90:VAL:HG12	1:D:122:GLU:HA	2.02	0.40
1:D:105:GLY:HA3	1:D:128:ASN:ND2	2.37	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:111:ASN:ND2	1:D:114:SER:OG	2.55	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	235/242 (97%)	220 (94%)	11 (5%)	4 (2%)	9	7
1	B	235/242 (97%)	224 (95%)	10 (4%)	1 (0%)	34	38
1	C	235/242 (97%)	219 (93%)	14 (6%)	2 (1%)	17	17
1	D	235/242 (97%)	218 (93%)	14 (6%)	3 (1%)	12	10
All	All	940/968 (97%)	881 (94%)	49 (5%)	10 (1%)	14	13

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	83	PRO
1	C	83	PRO
1	D	83	PRO
1	D	107	PHE
1	B	83	PRO
1	C	13	ASN
1	A	107	PHE
1	A	82	ARG
1	A	81	PRO
1	D	84	HIS

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	202/211 (96%)	194 (96%)	8 (4%)	31	39
1	B	204/211 (97%)	198 (97%)	6 (3%)	42	52
1	C	202/211 (96%)	195 (96%)	7 (4%)	36	44
1	D	196/211 (93%)	188 (96%)	8 (4%)	30	37
All	All	804/844 (95%)	775 (96%)	29 (4%)	35	43

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	TRP
1	A	102	GLU
1	A	133	GLN
1	A	174	HIS
1	A	175	VAL
1	A	196	GLN
1	A	197	VAL
1	A	235	LEU
1	B	19	LEU
1	B	36	VAL
1	B	43	TRP
1	B	149	THR
1	B	174	HIS
1	B	196	GLN
1	C	43	TRP
1	C	84	HIS
1	C	102	GLU
1	C	143	SER
1	C	149	THR
1	C	174	HIS
1	C	196	GLN
1	D	43	TRP
1	D	102	GLU
1	D	133	GLN

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Mol	Chain	Res	Type
1	D	149	THR
1	D	175	VAL
1	D	196	GLN
1	D	227	LEU
1	D	235	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	12	GLN
1	A	78	GLN
1	A	84	HIS
1	A	97	ASN
1	A	111	ASN
1	A	133	GLN
1	A	174	HIS
1	A	203	ASN
1	B	9	GLN
1	B	64	ASN
1	B	78	GLN
1	B	84	HIS
1	B	99	GLN
1	B	111	ASN
1	B	133	GLN
1	B	203	ASN
1	C	8	ASN
1	C	38	ASN
1	C	78	GLN
1	C	84	HIS
1	C	97	ASN
1	C	133	GLN
1	C	174	HIS
1	C	196	GLN
1	C	203	ASN
1	D	6	ASN
1	D	12	GLN
1	D	64	ASN
1	D	78	GLN
1	D	84	HIS
1	D	111	ASN
1	D	128	ASN

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Mol	Chain	Res	Type
1	D	133	GLN
1	D	174	HIS
1	D	203	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](#)

18 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	E	1	2,1	14,14,15	0.57	0	17,19,21	0.72	1 (5%)
2	FUC	E	2	2	10,10,11	0.62	0	14,14,16	0.36	0
2	NAG	E	3	2	14,14,15	0.54	0	17,19,21	0.66	1 (5%)
3	NAG	F	1	3,1	14,14,15	0.71	0	17,19,21	0.77	1 (5%)
3	FUC	F	2	3	10,10,11	0.51	0	14,14,16	0.64	0
2	NAG	G	1	2,1	14,14,15	0.58	0	17,19,21	0.69	1 (5%)
2	FUC	G	2	2	10,10,11	0.50	0	14,14,16	0.34	0
2	NAG	G	3	2	14,14,15	0.54	0	17,19,21	0.61	0
2	NAG	H	1	2,1	14,14,15	0.66	0	17,19,21	0.68	0
2	FUC	H	2	2	10,10,11	0.50	0	14,14,16	0.33	0
2	NAG	H	3	2	14,14,15	0.70	0	17,19,21	0.67	0
4	NAG	I	1	4,1	14,14,15	0.57	0	17,19,21	0.73	1 (5%)
4	NAG	I	2	4	14,14,15	0.61	0	17,19,21	1.02	1 (5%)
4	BMA	I	3	4	11,11,12	0.42	0	15,15,17	0.88	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	FUC	I	4	4	10,10,11	0.53	0	14,14,16	0.31	0
2	NAG	J	1	2,1	14,14,15	0.58	0	17,19,21	0.69	0
2	FUC	J	2	2	10,10,11	0.54	0	14,14,16	0.30	0
2	NAG	J	3	2	14,14,15	0.53	0	17,19,21	0.74	1 (5%)

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

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

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	I	3	BMA	C1-O5-C5	2.77	115.95	112.19
2	J	3	NAG	C2-N2-C7	-2.20	119.78	122.90
4	I	1	NAG	C2-N2-C7	-2.19	119.79	122.90
3	F	1	NAG	C2-N2-C7	-2.19	119.79	122.90
2	E	1	NAG	C2-N2-C7	-2.14	119.85	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	I	2	NAG	C2-N2-C7	-2.08	119.94	122.90
2	G	1	NAG	C2-N2-C7	-2.01	120.04	122.90
2	E	3	NAG	C2-N2-C7	-2.01	120.04	122.90

There are no chirality outliers.

All (28) torsion outliers are listed below:

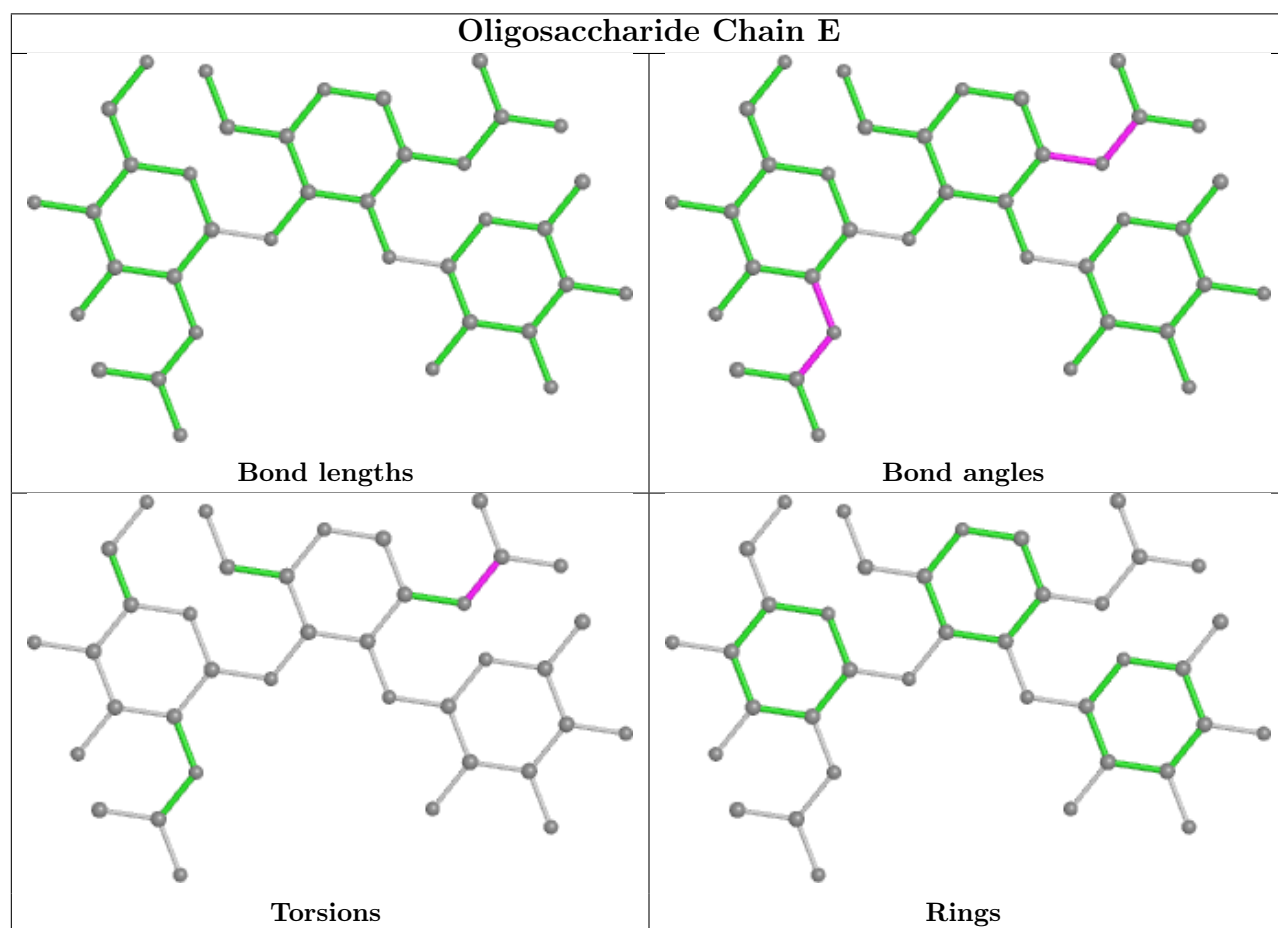
Mol	Chain	Res	Type	Atoms
2	H	3	NAG	C3-C2-N2-C7
2	H	3	NAG	C8-C7-N2-C2
2	H	3	NAG	O7-C7-N2-C2
2	J	3	NAG	C8-C7-N2-C2
2	J	3	NAG	O7-C7-N2-C2
4	I	1	NAG	C8-C7-N2-C2
4	I	1	NAG	O7-C7-N2-C2
2	H	3	NAG	C4-C5-C6-O6
2	J	1	NAG	C8-C7-N2-C2
2	J	1	NAG	O7-C7-N2-C2
2	H	3	NAG	O5-C5-C6-O6
2	G	1	NAG	C8-C7-N2-C2
2	G	1	NAG	O7-C7-N2-C2
3	F	1	NAG	C8-C7-N2-C2
2	J	3	NAG	O5-C5-C6-O6
2	H	1	NAG	C4-C5-C6-O6
2	J	1	NAG	O5-C5-C6-O6
2	G	3	NAG	O5-C5-C6-O6
2	J	1	NAG	C4-C5-C6-O6
2	E	1	NAG	C8-C7-N2-C2
3	F	1	NAG	O7-C7-N2-C2
2	H	1	NAG	O5-C5-C6-O6
2	J	3	NAG	C4-C5-C6-O6
2	G	3	NAG	C4-C5-C6-O6
2	E	1	NAG	O7-C7-N2-C2
4	I	3	BMA	O5-C5-C6-O6
3	F	1	NAG	C4-C5-C6-O6
3	F	1	NAG	O5-C5-C6-O6

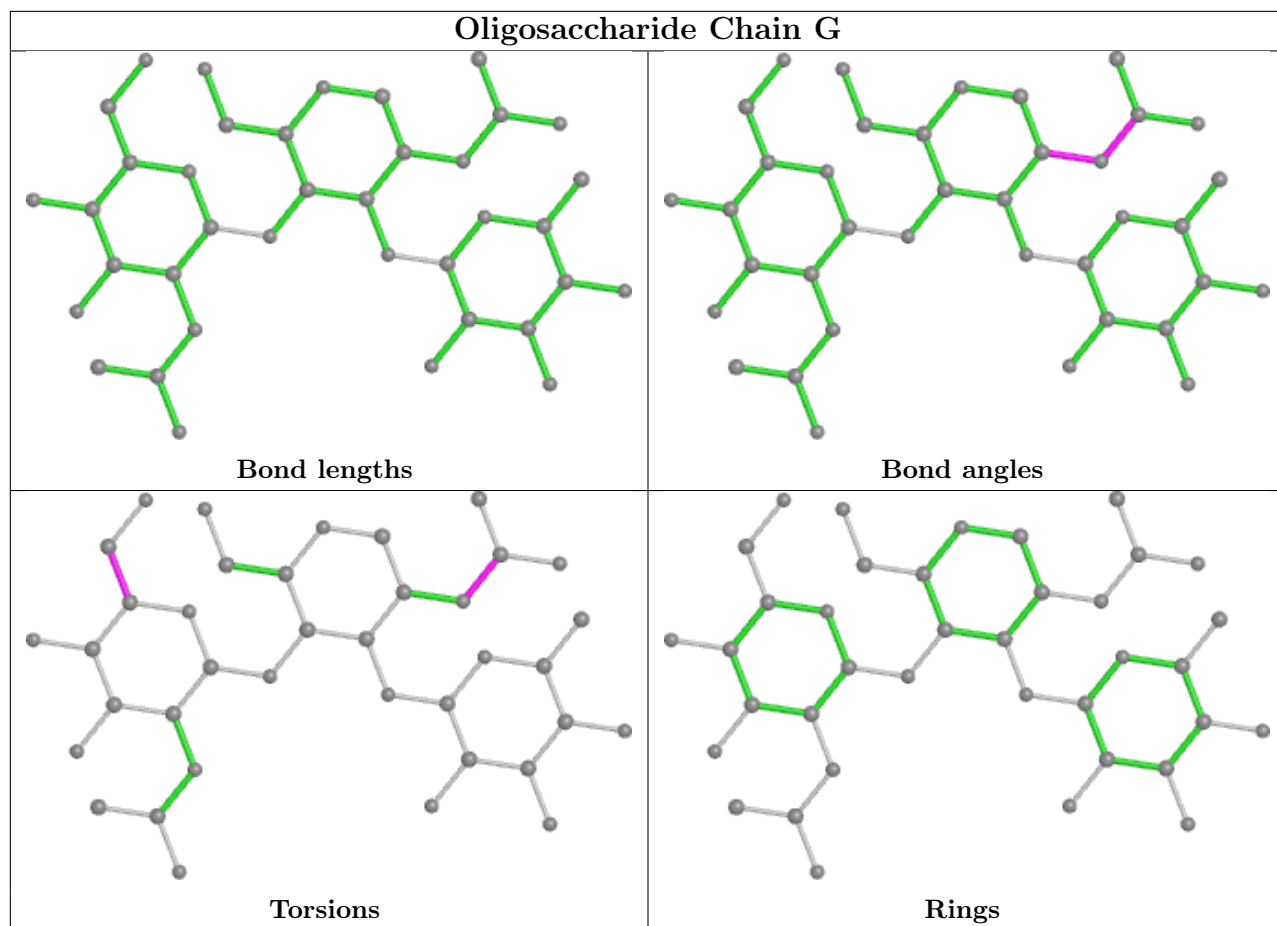
There are no ring outliers.

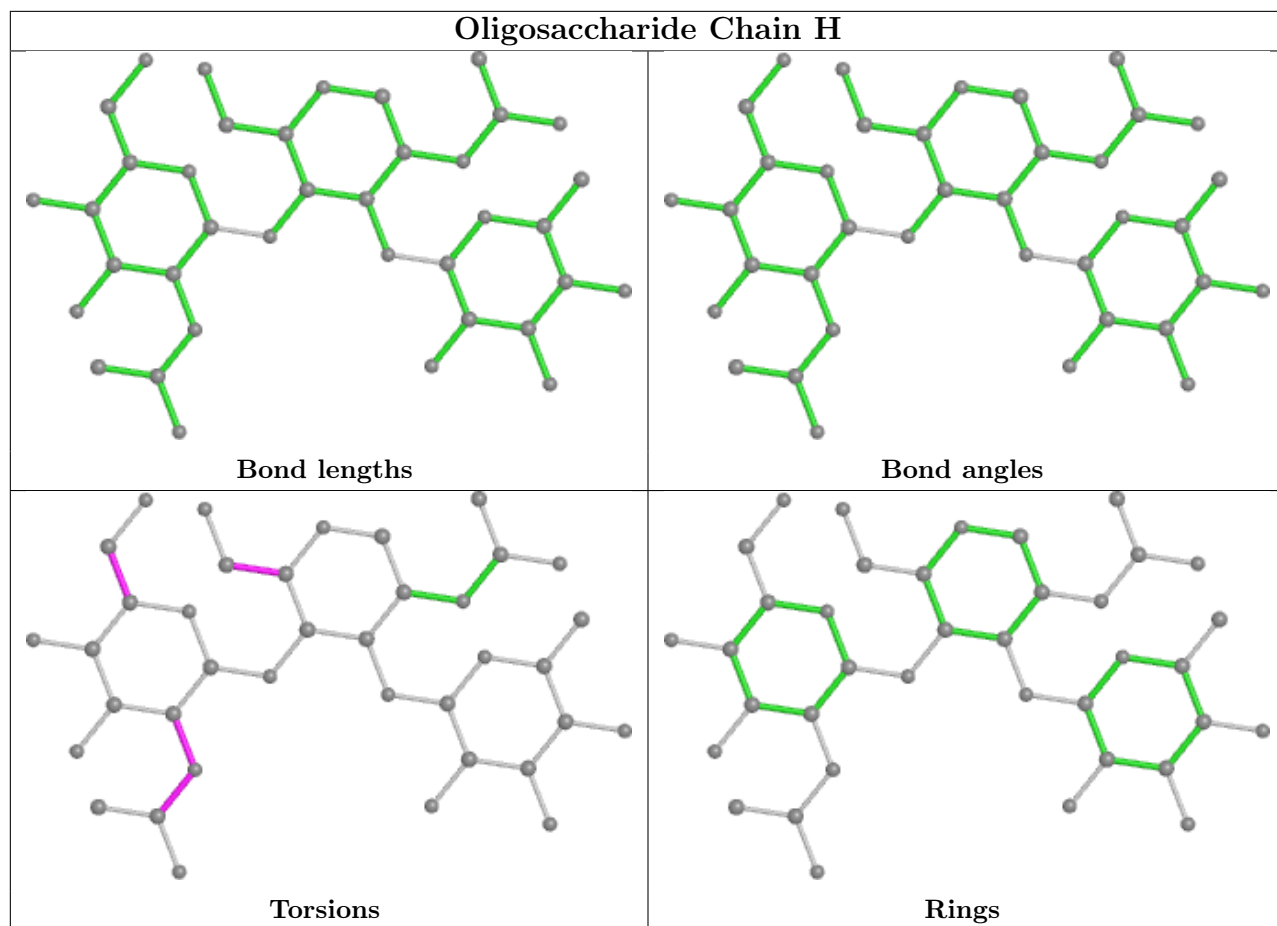
6 monomers are involved in 13 short contacts:

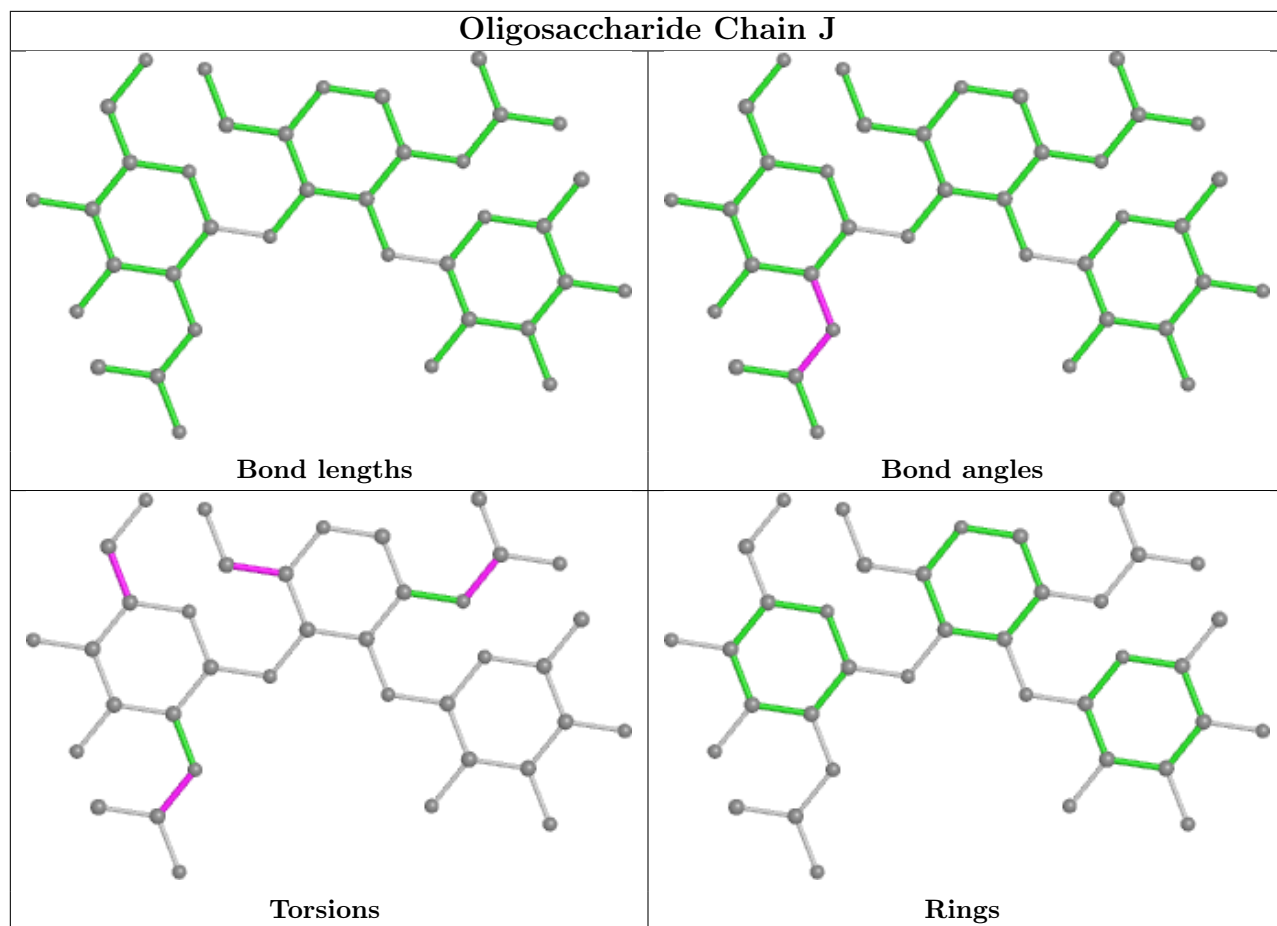
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	H	1	NAG	4	0
2	H	3	NAG	6	0
3	F	1	NAG	5	0
3	F	2	FUC	3	0
2	G	1	NAG	2	0
2	H	2	FUC	1	0

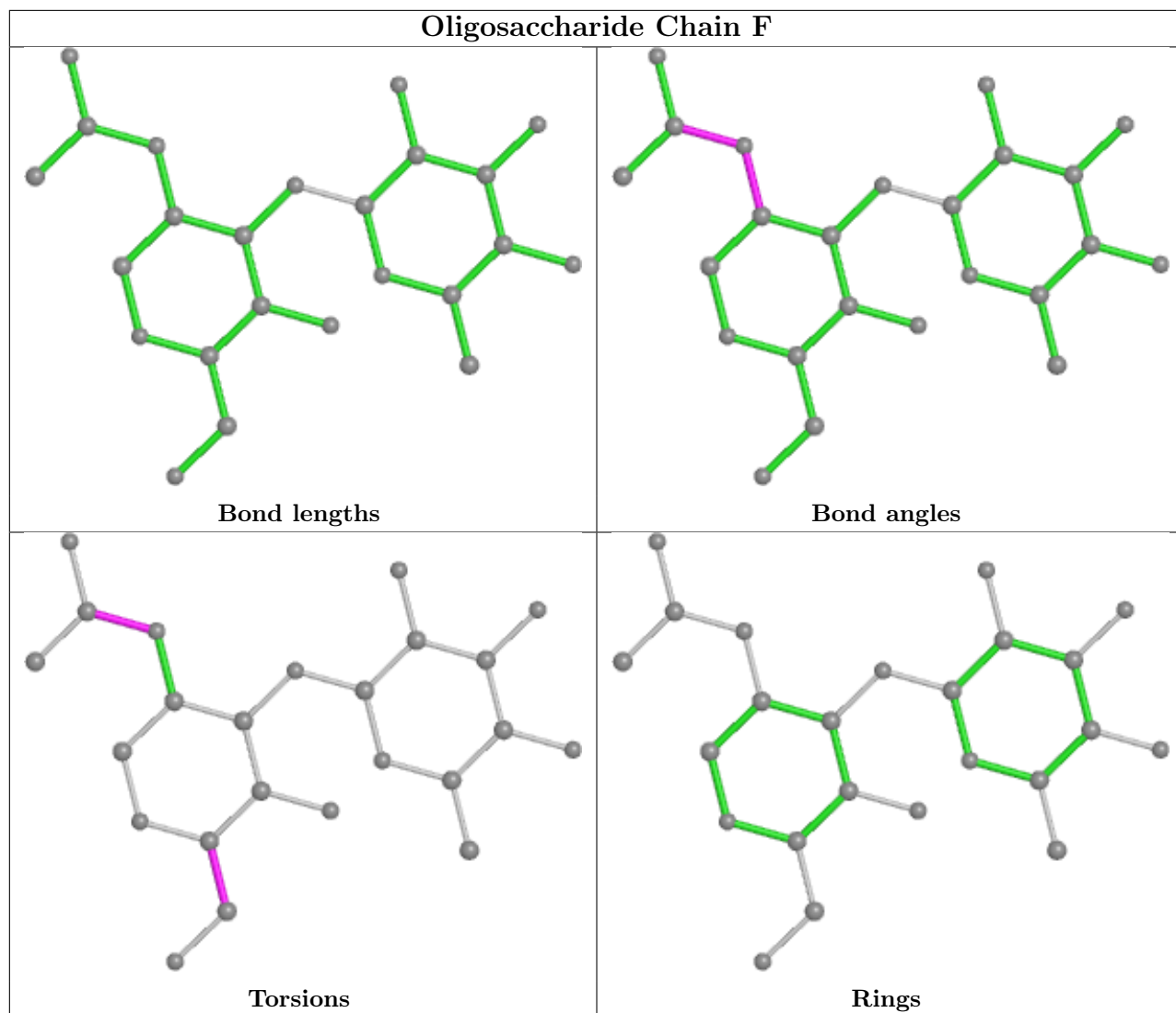
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

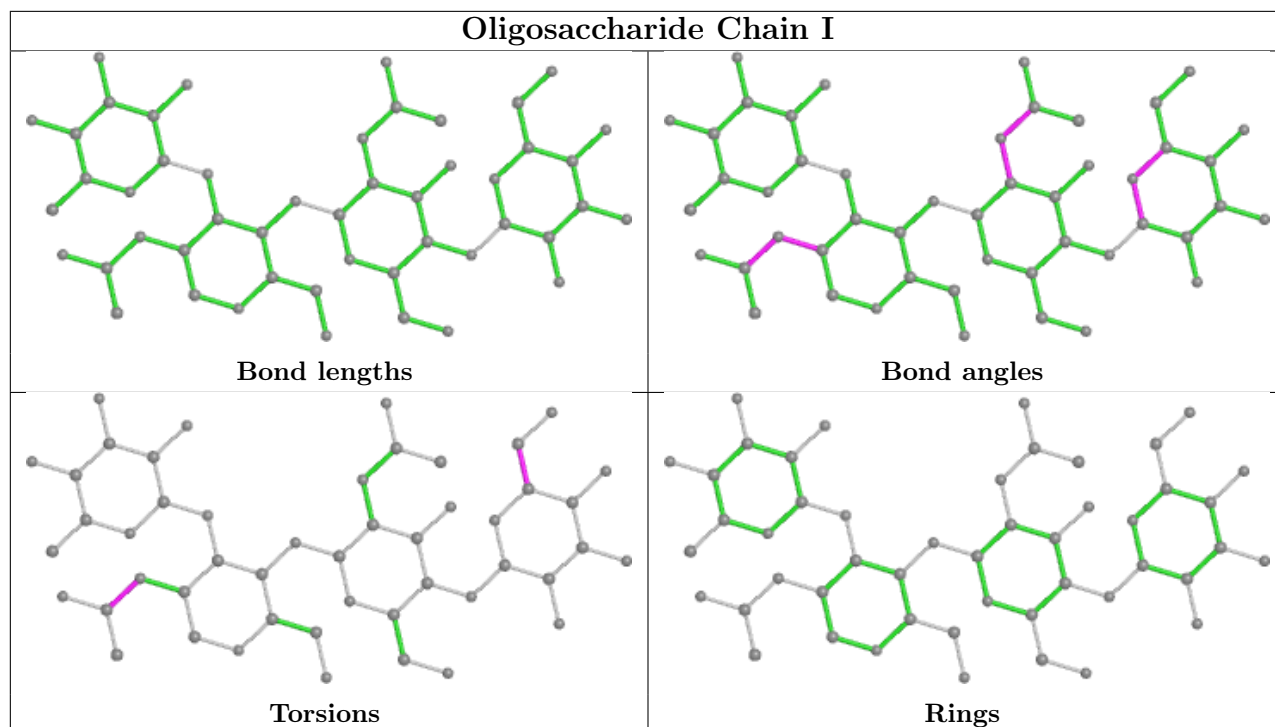












5.6 Ligand geometry [i](#)

Of 17 ligands modelled in this entry, 8 are monoatomic - leaving 9 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
7	SER	C	508	8	5,6,6	0.99	1 (20%)	5,7,7	1.38	2 (40%)
8	A2G	A	509	7	14,14,15	0.64	0	17,19,21	0.79	0
8	A2G	B	510	7	14,14,15	0.69	0	17,19,21	0.79	0
7	SER	B	509	8	5,6,6	0.95	1 (20%)	5,7,7	1.37	2 (40%)
9	NAG	C	505	1	14,14,15	0.69	0	17,19,21	0.69	0
7	SER	A	508	8	5,6,6	0.92	0	5,7,7	1.37	2 (40%)
8	A2G	C	509	7	14,14,15	0.84	1 (7%)	17,19,21	0.79	0
8	A2G	D	607	7	14,14,15	0.82	1 (7%)	17,19,21	0.78	0
7	SER	D	606	8	5,6,6	1.01	1 (20%)	5,7,7	1.36	2 (40%)

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
7	SER	C	508	8	-	0/6/6/6	-
8	A2G	A	509	7	-	0/6/23/26	0/1/1/1
8	A2G	B	510	7	-	0/6/23/26	0/1/1/1
7	SER	B	509	8	-	0/6/6/6	-
9	NAG	C	505	1	-	2/6/23/26	0/1/1/1
7	SER	A	508	8	-	2/6/6/6	-
8	A2G	C	509	7	-	0/6/23/26	0/1/1/1
8	A2G	D	607	7	-	0/6/23/26	0/1/1/1
7	SER	D	606	8	-	3/6/6/6	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	C	509	A2G	C1-C2	2.38	1.55	1.52
8	D	607	A2G	C1-C2	2.14	1.55	1.52
7	D	606	SER	OXT-C	-2.07	1.23	1.30
7	C	508	SER	OXT-C	-2.05	1.23	1.30
7	B	509	SER	OXT-C	-2.01	1.24	1.30

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	509	SER	OXT-C-O	-2.21	119.08	124.09
7	C	508	SER	OXT-C-O	-2.19	119.12	124.09
7	A	508	SER	OXT-C-O	-2.19	119.13	124.09
7	D	606	SER	OXT-C-O	-2.15	119.21	124.09
7	C	508	SER	OXT-C-CA	2.09	120.52	113.38
7	B	509	SER	OXT-C-CA	2.07	120.45	113.38
7	A	508	SER	OXT-C-CA	2.07	120.44	113.38
7	D	606	SER	OXT-C-CA	2.07	120.43	113.38

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	D	606	SER	N-CA-CB-OG
9	C	505	NAG	O5-C5-C6-O6
9	C	505	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
7	D	606	SER	C-CA-CB-OG
7	A	508	SER	OXT-C-CA-CB
7	A	508	SER	O-C-CA-CB
7	D	606	SER	OXT-C-CA-CB

There are no ring outliers.

No monomer is involved in short contacts.

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	237/242 (97%)	-0.11	9 (3%) 40 53	24, 36, 57, 71	0
1	B	237/242 (97%)	-0.08	5 (2%) 63 74	21, 34, 63, 75	0
1	C	237/242 (97%)	0.06	13 (5%) 25 36	29, 42, 65, 86	0
1	D	237/242 (97%)	0.30	19 (8%) 12 18	28, 47, 74, 85	0
All	All	948/968 (97%)	0.04	46 (4%) 29 42	21, 40, 68, 86	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	237	GLY	6.6
1	B	82	ARG	5.2
1	D	113	LEU	5.0
1	B	113	LEU	4.9
1	C	82	ARG	4.8
1	C	237	GLY	4.8
1	C	90	VAL	4.7
1	D	82	ARG	4.7
1	A	82	ARG	4.3
1	C	89	LEU	4.1
1	C	13	ASN	4.1
1	C	12	GLN	4.1
1	D	90	VAL	3.9
1	D	37	VAL	3.8
1	D	13	ASN	3.4
1	B	90	VAL	3.3
1	C	15	GLU	3.3
1	A	90	VAL	3.2
1	D	206	PHE	3.0
1	D	207	SER	2.9
1	D	36	VAL	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	123	PHE	2.8
1	D	236	PRO	2.8
1	C	88	GLY	2.7
1	D	89	LEU	2.7
1	A	89	LEU	2.5
1	C	91	PHE	2.5
1	D	39	GLY	2.4
1	A	13	ASN	2.4
1	C	206	PHE	2.4
1	D	43	TRP	2.4
1	D	12	GLN	2.3
1	A	113	LEU	2.3
1	C	11	HIS	2.3
1	D	91	PHE	2.3
1	B	114	SER	2.3
1	D	88	GLY	2.3
1	C	208	ALA	2.3
1	D	40	VAL	2.2
1	C	14	GLU	2.2
1	B	237	GLY	2.2
1	A	91	PHE	2.2
1	D	121	VAL	2.1
1	D	27	SER	2.1
1	A	24	ARG	2.0
1	A	206	PHE	2.0

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

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

6.3 Carbohydrates [i](#)

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

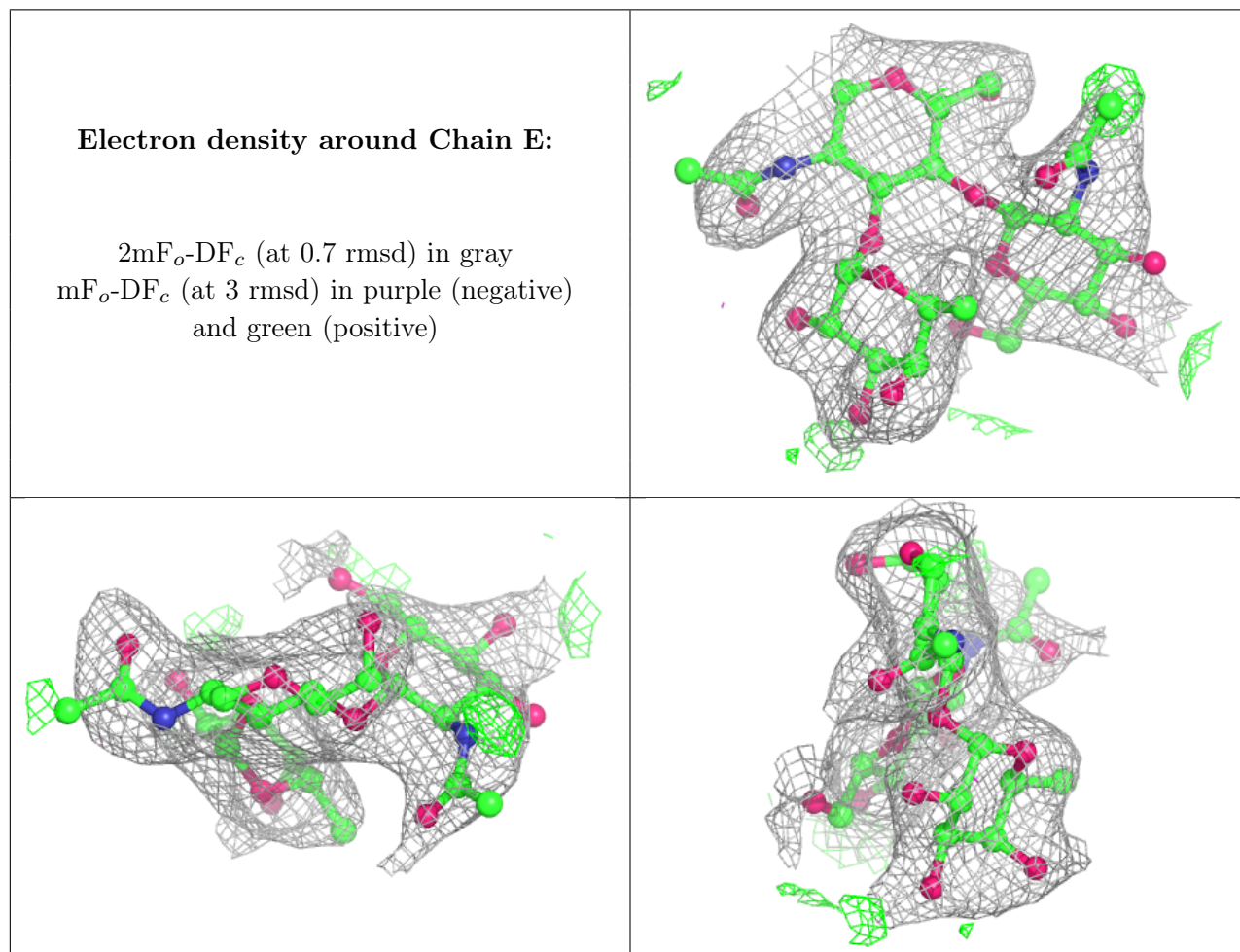
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	FUC	F	2	10/11	0.71	0.39	80,82,83,84	0
4	BMA	I	3	11/12	0.77	0.27	83,85,87,87	0
2	NAG	H	3	14/15	0.79	0.31	78,80,84,85	0

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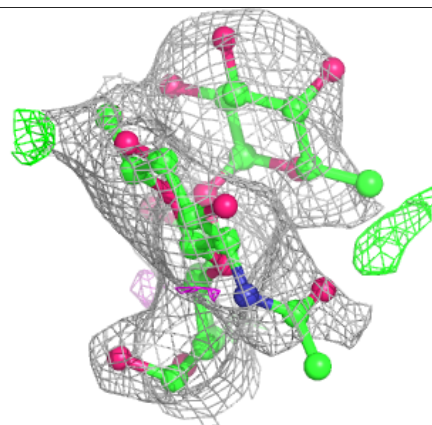
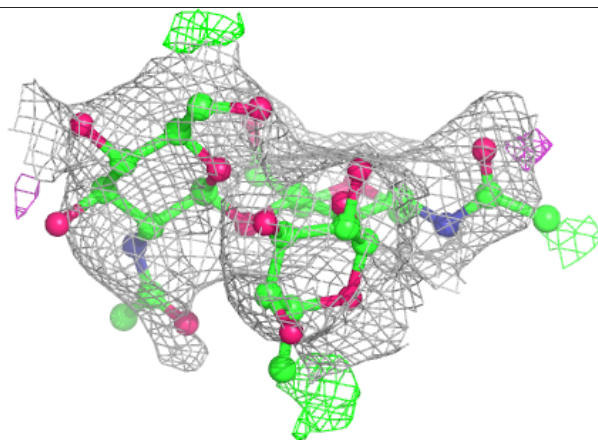
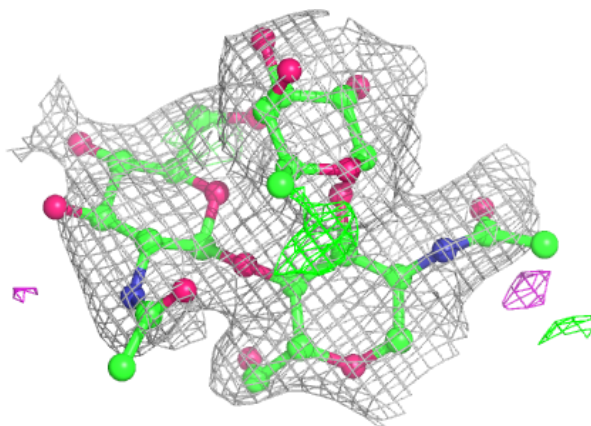
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	FUC	J	2	10/11	0.82	0.40	86,88,89,89	0
3	NAG	F	1	14/15	0.83	0.22	60,64,68,76	0
2	NAG	J	3	14/15	0.85	0.36	88,90,91,91	0
4	NAG	I	2	14/15	0.85	0.19	64,72,76,79	0
2	FUC	G	2	10/11	0.85	0.24	77,79,80,80	0
2	NAG	G	3	14/15	0.86	0.26	75,77,79,79	0
2	NAG	E	3	14/15	0.86	0.27	76,79,82,83	0
2	NAG	H	1	14/15	0.89	0.21	49,60,67,73	0
2	NAG	G	1	14/15	0.89	0.17	60,66,71,73	0
4	NAG	I	1	14/15	0.90	0.17	58,60,66,67	0
2	NAG	J	1	14/15	0.91	0.14	75,79,83,84	0
2	NAG	E	1	14/15	0.91	0.11	59,61,64,71	0
4	FUC	I	4	10/11	0.91	0.25	70,71,72,75	0
2	FUC	E	2	10/11	0.92	0.14	66,68,68,69	0
2	FUC	H	2	10/11	0.92	0.21	58,62,63,63	0

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



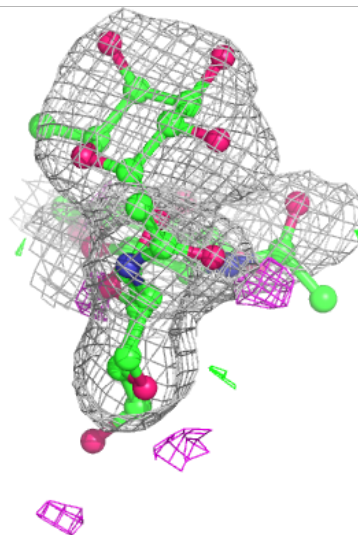
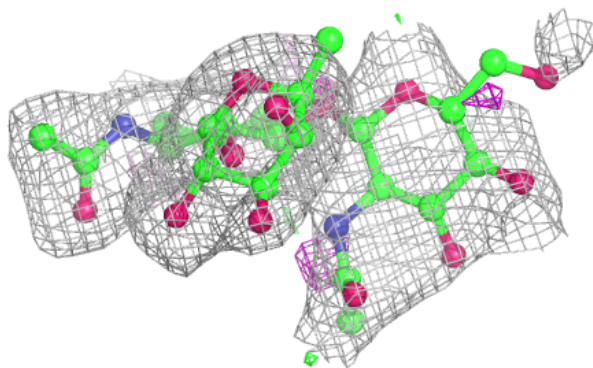
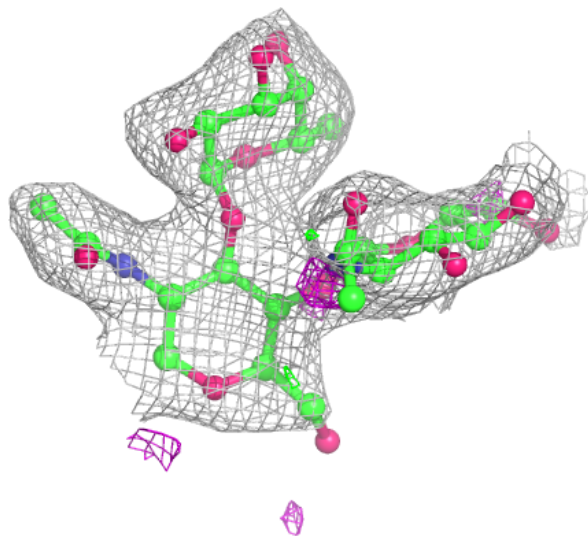
Electron density around Chain G:

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



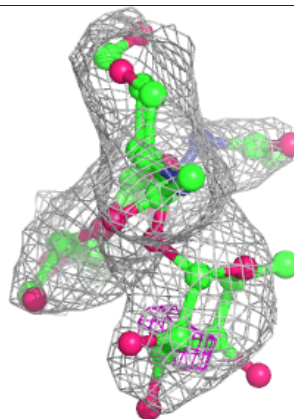
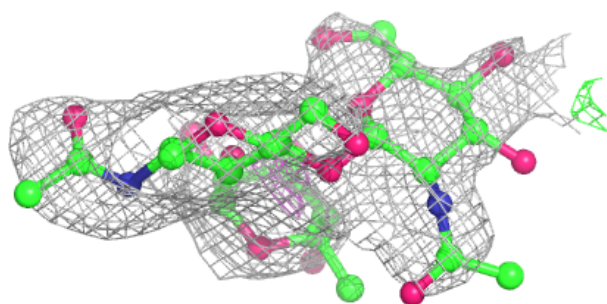
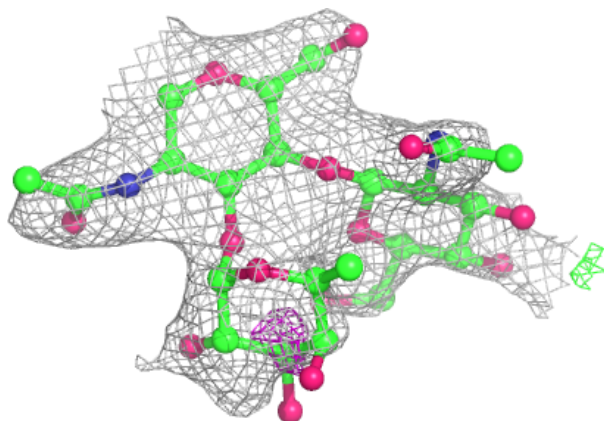
Electron density around Chain H:

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



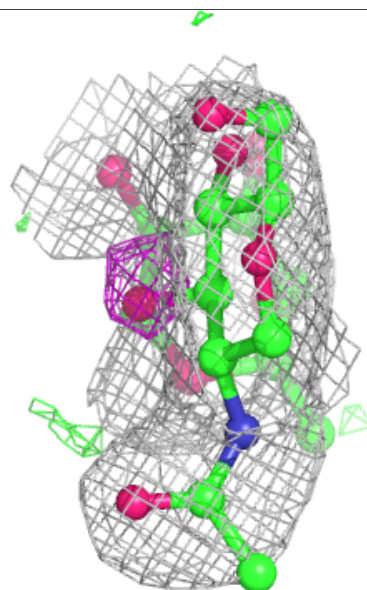
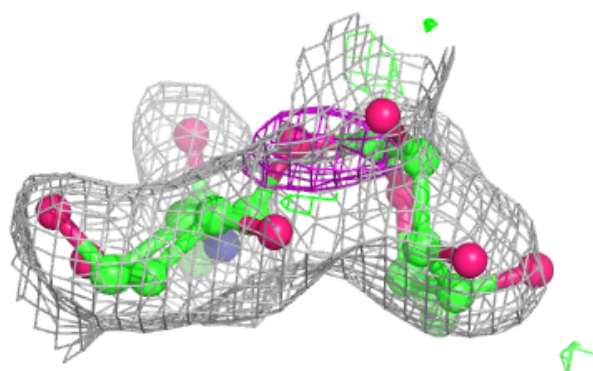
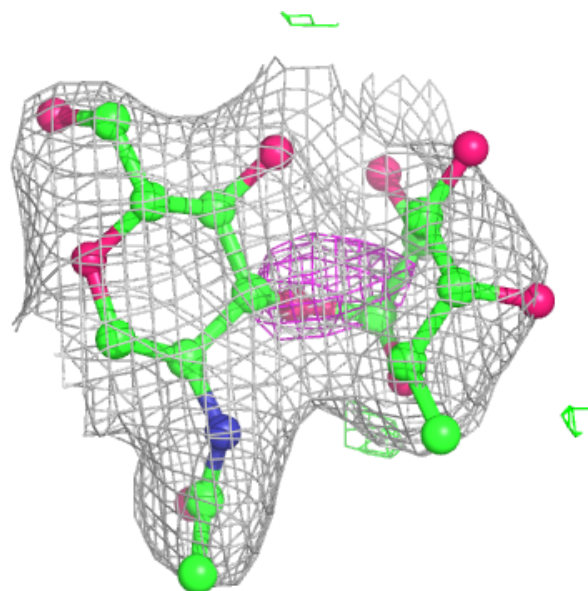
Electron density around Chain J:

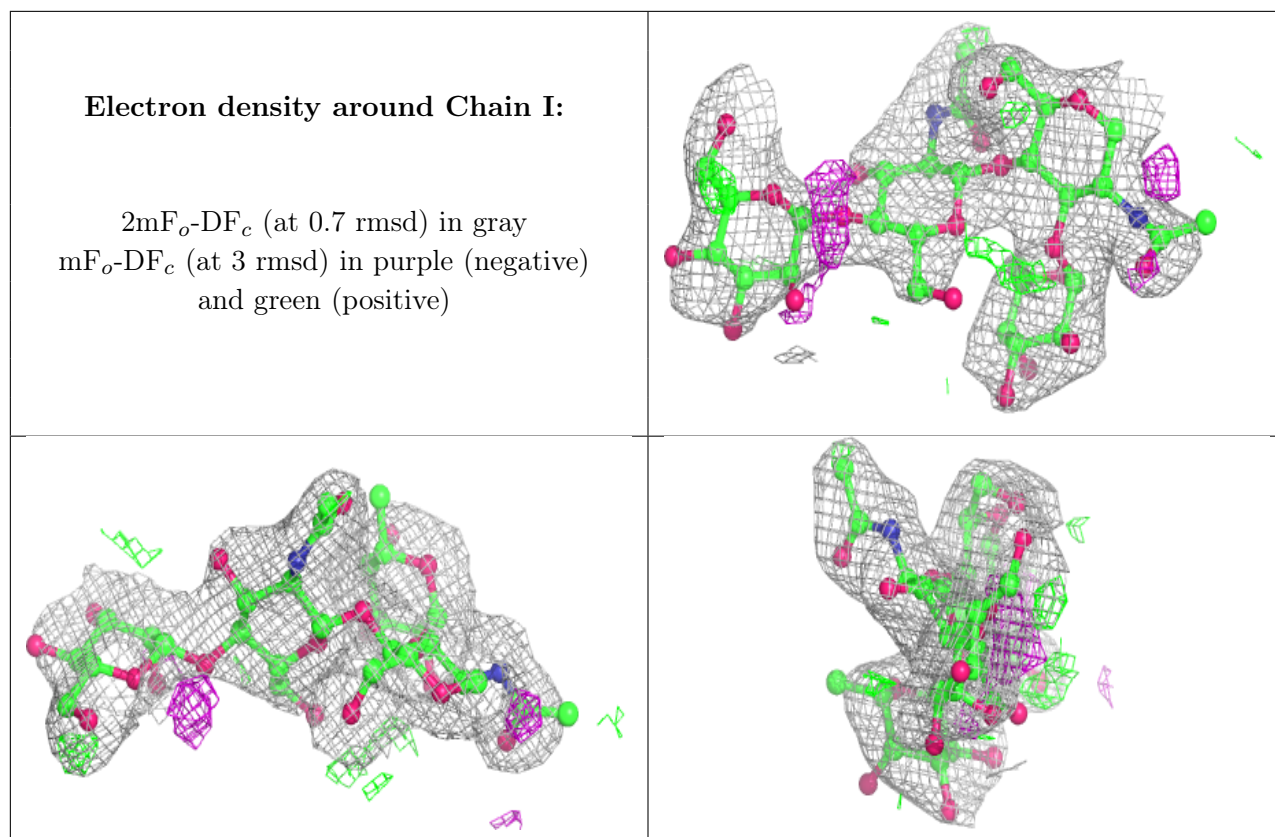
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain F:

$2mF_o-DF_c$ (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
9	NAG	C	505	14/15	0.80	0.25	63,65,68,70	0
7	SER	D	606	7/7	0.84	0.23	55,63,66,67	0
7	SER	C	508	7/7	0.88	0.13	53,58,60,60	0
7	SER	A	508	7/7	0.89	0.17	50,55,57,58	0
7	SER	B	509	7/7	0.92	0.17	43,51,54,55	0
8	A2G	C	509	14/15	0.94	0.10	45,49,51,51	0
8	A2G	D	607	14/15	0.95	0.10	44,49,52,53	0
5	CA	C	506	1/1	0.96	0.15	43,43,43,43	0
8	A2G	B	510	14/15	0.97	0.09	33,36,38,39	0
8	A2G	A	509	14/15	0.97	0.09	40,43,45,46	0
5	CA	D	604	1/1	0.98	0.18	38,38,38,38	0
6	MN	C	507	1/1	0.98	0.06	49,49,49,49	0
6	MN	D	605	1/1	0.98	0.05	49,49,49,49	0
6	MN	B	508	1/1	0.99	0.09	35,35,35,35	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	CA	A	506	1/1	0.99	0.13	34,34,34,34	0
5	CA	B	507	1/1	0.99	0.14	34,34,34,34	0
6	MN	A	507	1/1	0.99	0.06	43,43,43,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.