



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 8, 2020 – 07:22 PM BST

PDB ID : 5E2Z
Title : Crystal structure of H5 hemagglutinin Q226L mutant from the influenza virus A/duck/Egypt/10185SS/2010 (H5N1) with LSTa
Authors : Zhu, X.; Wilson, I.A.
Deposited on : 2015-10-01
Resolution : 2.62 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 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 : 2.13.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

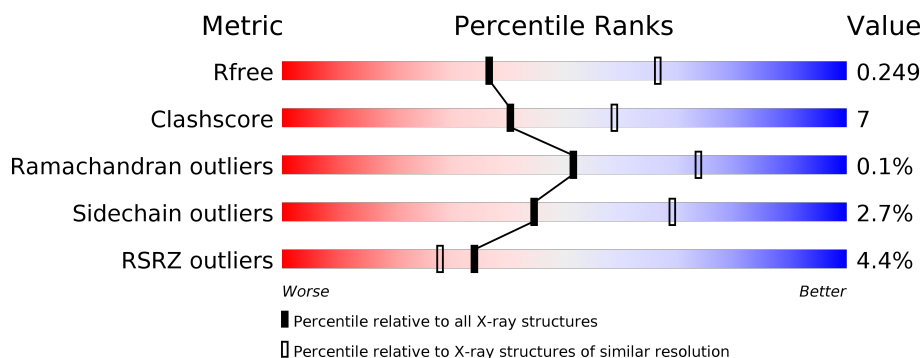
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.62 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	3797 (2.64-2.60)
Clashscore	141614	4168 (2.64-2.60)
Ramachandran outliers	138981	4093 (2.64-2.60)
Sidechain outliers	138945	4093 (2.64-2.60)
RSRZ outliers	127900	3731 (2.64-2.60)

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 on 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	333	
1	C	333	
1	E	333	
2	B	180	
2	D	180	
2	F	180	

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Mol	Chain	Length	Quality of chain
3	G	3	 33% 67%
3	K	3	 33% 67%
4	H	2	 50% 50%
4	L	2	 100%
5	I	4	 25% 25% 50%
6	J	3	 67% 33%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 12369 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
1	C	323	Total	C	N	O	S	0	0	0
			2558	1611	443	490	14			
1	A	323	Total	C	N	O	S	0	0	0
			2558	1611	443	490	14			
1	E	323	Total	C	N	O	S	0	0	0
			2558	1611	443	490	14			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	7	ALA	-	expression tag	UNP G8IPF0
C	8	ASP	-	expression tag	UNP G8IPF0
C	9	PRO	-	expression tag	UNP G8IPF0
C	10	GLY	-	expression tag	UNP G8IPF0
C	226	LEU	GLN	engineered mutation	UNP G8IPF0
A	7	ALA	-	expression tag	UNP G8IPF0
A	8	ASP	-	expression tag	UNP G8IPF0
A	9	PRO	-	expression tag	UNP G8IPF0
A	10	GLY	-	expression tag	UNP G8IPF0
A	226	LEU	GLN	engineered mutation	UNP G8IPF0
E	7	ALA	-	expression tag	UNP G8IPF0
E	8	ASP	-	expression tag	UNP G8IPF0
E	9	PRO	-	expression tag	UNP G8IPF0
E	10	GLY	-	expression tag	UNP G8IPF0
E	226	LEU	GLN	engineered mutation	UNP G8IPF0

- Molecule 2 is a protein called Hemagglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	175	Total	C	N	O	S	0	0	0
			1418	880	248	282	8			

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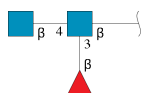
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	175	Total	C	N	O	S	0	0	0
			1418	880	248	282	8			
2	F	175	Total	C	N	O	S	0	0	0
			1418	880	248	282	8			

There are 18 discrepancies between the modelled and reference sequences:

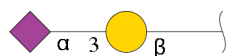
Chain	Residue	Modelled	Actual	Comment	Reference
D	1	GLY	-	expression tag	UNP G8IPF0
D	176	ARG	-	expression tag	UNP G8IPF0
D	177	LEU	-	expression tag	UNP G8IPF0
D	178	VAL	-	expression tag	UNP G8IPF0
D	179	PRO	-	expression tag	UNP G8IPF0
D	180	ARG	-	expression tag	UNP G8IPF0
B	1	GLY	-	expression tag	UNP G8IPF0
B	176	ARG	-	expression tag	UNP G8IPF0
B	177	LEU	-	expression tag	UNP G8IPF0
B	178	VAL	-	expression tag	UNP G8IPF0
B	179	PRO	-	expression tag	UNP G8IPF0
B	180	ARG	-	expression tag	UNP G8IPF0
F	1	GLY	-	expression tag	UNP G8IPF0
F	176	ARG	-	expression tag	UNP G8IPF0
F	177	LEU	-	expression tag	UNP G8IPF0
F	178	VAL	-	expression tag	UNP G8IPF0
F	179	PRO	-	expression tag	UNP G8IPF0
F	180	ARG	-	expression tag	UNP G8IPF0

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



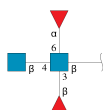
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	G	3	Total	C	N	O	0	0	0
			38	22	2	14			
3	K	3	Total	C	N	O	0	0	0
			38	22	2	14			

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



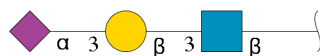
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	H	2	Total	C	N	O	0	0	0
			31	17	1	13			
4	L	2	Total	C	N	O	0	0	0
			31	17	1	13			

- Molecule 5 is an oligosaccharide called beta-L-fucopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)][alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	I	4	Total	C	N	O	0	0	0
			48	28	2	18			

- Molecule 6 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	J	3	Total	C	N	O	0	0	0
			46	25	2	19			

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

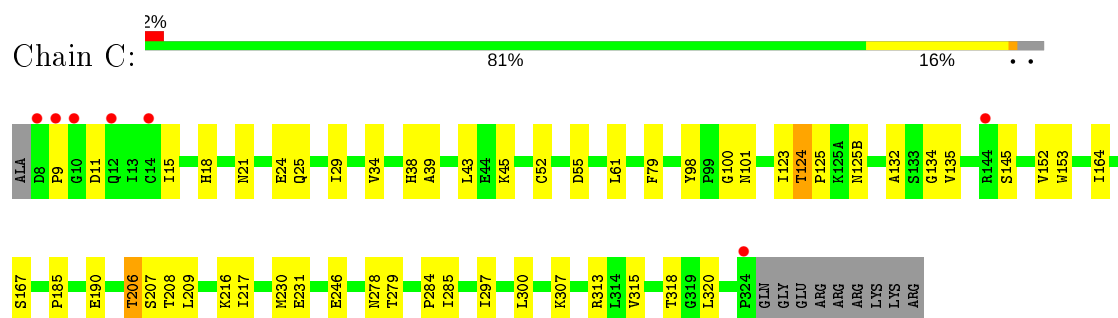
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	C	47	Total	O	0	0
			47	47		
8	D	7	Total	O	0	0
			7	7		
8	A	54	Total	O	0	0
			54	54		
8	B	8	Total	O	0	0
			8	8		
8	E	43	Total	O	0	0
			43	43		
8	F	8	Total	O	0	0
			8	8		

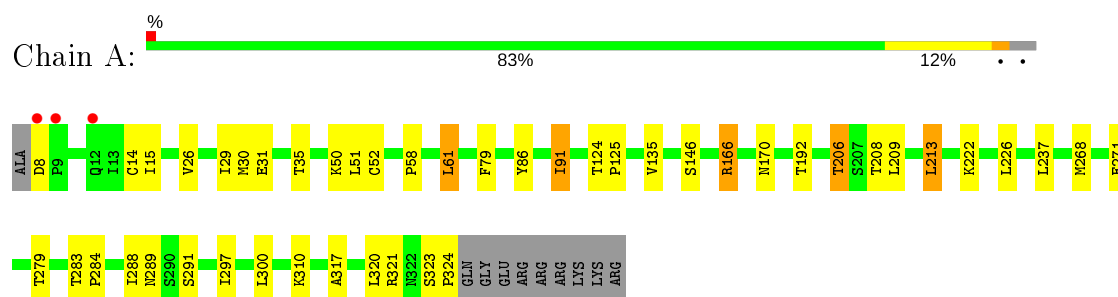
3 Residue-property plots [i](#)

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.

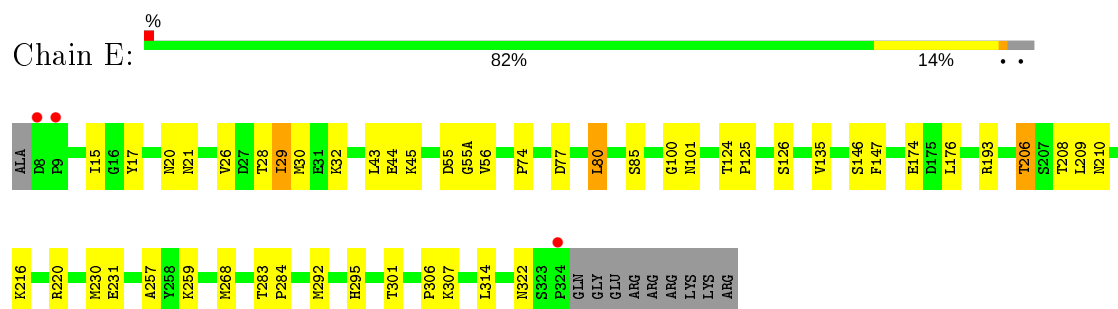
• Molecule 1: Hemagglutinin



• Molecule 1: Hemagglutinin

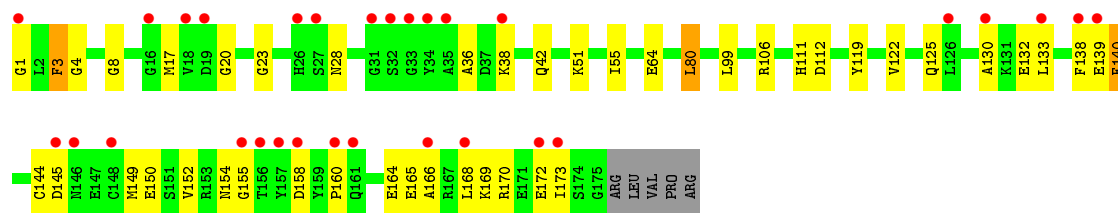


• Molecule 1: Hemagglutinin

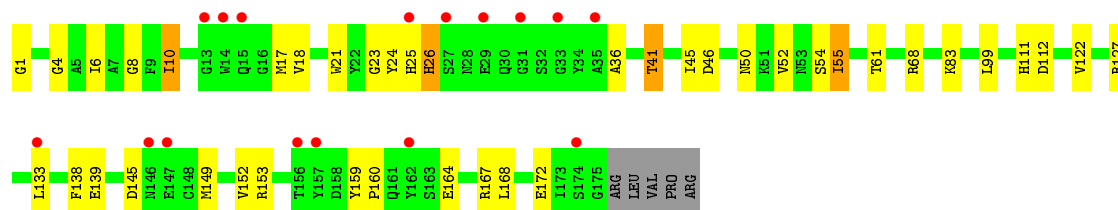
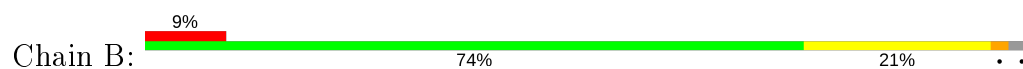


• Molecule 2: Hemagglutinin

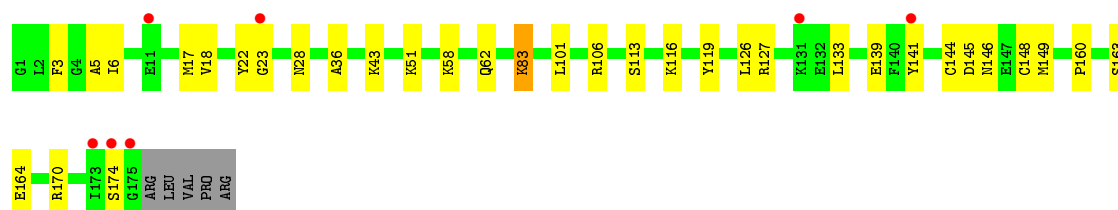
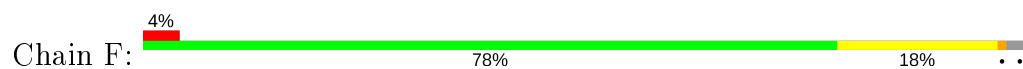




• Molecule 2: Hemagglutinin



• Molecule 2: Hemagglutinin



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



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



• Molecule 4: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose



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

Chain L:  100%

GAL1
S1A2

- Molecule 5: beta-L-fucopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)]|alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  25% 25% 50%

MAG1
FUL2
MAG3
FUC4

- Molecule 6: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  67% 33%

MAG1
GAL2
S1A3

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	73.06Å 234.15Å 72.95Å 90.00° 115.49° 90.00°	Depositor
Resolution (Å)	50.00 – 2.62 48.42 – 2.62	Depositor EDS
% Data completeness (in resolution range)	86.8 (50.00-2.62) 86.8 (48.42-2.62)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.89 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309	Depositor
R, R_{free}	0.202 , 0.248 0.203 , 0.249	Depositor DCC
R_{free} test set	2919 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	54.6	Xtriage
Anisotropy	0.303	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 53.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.027 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12369	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.65% 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: SIA, GAL, NAG, FUL, FUC

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	0.57	0/2619	0.68	1/3559 (0.0%)
1	C	0.55	0/2619	0.66	0/3559
1	E	0.49	0/2619	0.61	0/3559
2	B	0.45	0/1445	0.56	0/1942
2	D	0.46	0/1445	0.60	1/1942 (0.1%)
2	F	0.44	0/1445	0.55	0/1942
All	All	0.51	0/12192	0.62	2/16503 (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	213	LEU	CA-CB-CG	5.50	127.94	115.30
2	D	80	LEU	CA-CB-CG	5.34	127.58	115.30

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	2558	0	2492	32	0
1	C	2558	0	2492	36	0
1	E	2558	0	2492	35	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1418	0	1322	32	0
2	D	1418	0	1322	37	0
2	F	1418	0	1322	22	0
3	G	38	0	34	0	0
3	K	38	0	34	0	0
4	H	31	0	26	1	0
4	L	31	0	26	0	0
5	I	48	0	43	3	0
6	J	46	0	40	3	0
7	A	14	0	13	0	0
7	C	14	0	13	0	0
7	E	14	0	13	0	0
8	A	54	0	0	3	0
8	B	8	0	0	0	0
8	C	47	0	0	0	0
8	D	7	0	0	0	0
8	E	43	0	0	1	0
8	F	8	0	0	0	0
All	All	12369	0	11684	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 7.

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:206:THR:HG22	1:C:208:THR:H	1.38	0.85
1:E:206:THR:HG22	1:E:209:LEU:H	1.43	0.83
1:E:28:THR:HG22	1:E:30:MET:H	1.45	0.81
2:D:125:GLN:HE22	2:D:155:GLY:HA2	1.47	0.78
1:A:222:LYS:NZ	6:J:1:NAG:H82	2.04	0.73
1:E:174:GLU:HG3	1:E:259:LYS:HB3	1.72	0.72
1:E:20:ASN:OD1	1:E:322:ASN:ND2	2.25	0.70
2:D:140:PHE:H	2:D:140:PHE:HD1	1.37	0.69
1:E:307:LYS:HE2	2:F:62:GLN:HB3	1.73	0.69
1:A:206:THR:HG22	1:A:209:LEU:H	1.58	0.68
1:E:21:ASN:ND2	8:E:1102:HOH:O	2.29	0.65
1:A:206:THR:HG22	1:A:208:THR:H	1.62	0.65
5:I:3:NAG:H5	5:I:4:FUC:H63	1.77	0.65
2:D:130:ALA:HA	2:D:140:PHE:HA	1.77	0.65
2:D:28:ASN:ND2	2:D:144:CYS:O	2.30	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:29:ILE:HD11	2:F:51:LYS:HE2	1.79	0.64
2:B:17:MET:HE1	2:B:36:ALA:HA	1.80	0.63
1:A:26:VAL:HG21	1:A:317:ALA:HB2	1.79	0.63
2:B:167:ARG:HD2	2:F:174:SER:HA	1.81	0.63
2:B:24:TYR:CZ	2:B:153:ARG:HG2	2.36	0.61
2:B:26:HIS:HD2	2:B:149:MET:CE	2.13	0.61
1:E:206:THR:HG22	1:E:208:THR:H	1.65	0.61
2:D:1:GLY:HA3	2:D:3:PHE:CZ	2.36	0.60
2:D:3:PHE:HD1	2:D:4:GLY:H	1.46	0.60
2:F:145:ASP:OD1	2:F:148:CYS:N	2.33	0.60
1:A:289:ASN:ND2	8:A:1103:HOH:O	2.35	0.60
1:C:15:ILE:HD13	2:D:119:TYR:HA	1.86	0.58
1:A:222:LYS:HZ2	6:J:1:NAG:H82	1.67	0.58
2:B:6:ILE:HD13	2:B:112:ASP:HA	1.86	0.57
1:C:43:LEU:O	1:C:45:LYS:NZ	2.37	0.57
2:D:132:GLU:HG2	2:D:138:PHE:HE1	1.70	0.57
2:D:168:LEU:O	2:D:172:GLU:HG3	2.06	0.56
2:B:168:LEU:O	2:B:172:GLU:HG3	2.06	0.54
1:C:206:THR:HB	1:C:209:LEU:HB3	1.89	0.54
2:D:150:GLU:O	2:D:154:ASN:ND2	2.40	0.54
2:D:133:LEU:HA	2:F:127:ARG:HH12	1.72	0.54
2:B:46:ASP:O	2:B:50:ASN:ND2	2.28	0.54
1:C:185:PRO:HG2	1:C:217:ILE:HG12	1.88	0.54
2:F:164:GLU:CD	2:F:164:GLU:H	2.11	0.54
1:C:313:ARG:HH21	1:C:315:VAL:HG11	1.73	0.54
1:E:206:THR:HG22	1:E:209:LEU:N	2.20	0.53
2:D:3:PHE:HD1	2:D:4:GLY:N	2.06	0.53
1:A:58:PRO:HB3	1:A:86:TYR:CE1	2.44	0.53
2:D:3:PHE:HB2	2:D:112:ASP:OD1	2.08	0.53
1:C:15:ILE:HD11	2:D:122:VAL:HG21	1.90	0.53
1:C:52:CYS:HB2	1:C:279:THR:HG22	1.91	0.52
2:D:106:ARG:NH2	2:F:106:ARG:HH21	2.06	0.52
2:B:21:TRP:H	2:B:41:THR:HG21	1.73	0.52
1:C:285:ILE:HD11	1:C:297:ILE:HG23	1.91	0.52
1:A:15:ILE:O	2:B:10:ILE:HD12	2.09	0.52
2:D:51:LYS:HG3	1:A:29:ILE:HD12	1.92	0.52
2:F:133:LEU:HD11	2:F:139:GLU:HB2	1.91	0.51
1:C:98:TYR:CD1	1:C:230:MET:HG3	2.46	0.51
1:A:206:THR:HB	1:A:209:LEU:HB3	1.91	0.51
1:E:206:THR:CG2	1:E:208:THR:H	2.23	0.51
2:D:51:LYS:HE2	1:A:29:ILE:HD11	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:149:MET:O	2:B:152:VAL:HG22	2.11	0.51
2:D:169:LYS:O	2:D:173:ILE:HD12	2.11	0.50
1:A:310:LYS:NZ	8:A:1106:HOH:O	2.44	0.50
1:E:43:LEU:O	1:E:45:LYS:NZ	2.44	0.50
1:C:164:ILE:O	1:C:246:GLU:HA	2.11	0.50
2:D:64:GLU:O	2:B:83:LYS:NZ	2.33	0.50
2:B:133:LEU:HD11	2:B:139:GLU:HB2	1.94	0.50
2:B:41:THR:O	2:B:45:ILE:HG12	2.12	0.49
1:C:135:VAL:HG13	1:C:145:SER:HB3	1.94	0.49
1:E:101:ASN:OD1	1:E:231:GLU:HG3	2.11	0.49
1:E:206:THR:HB	1:E:209:LEU:HB3	1.94	0.49
2:D:149:MET:O	2:D:152:VAL:HG22	2.13	0.49
1:C:134:GLY:HA3	1:C:153:TRP:HB3	1.93	0.49
2:D:17:MET:HE3	2:D:23:GLY:HA3	1.95	0.49
1:E:268:MET:HG2	1:E:284:PRO:HG3	1.94	0.49
1:E:32:LYS:HD2	1:E:32:LYS:N	2.28	0.49
1:E:77:ASP:O	1:E:80:LEU:HD12	2.13	0.49
2:D:17:MET:HE3	2:D:36:ALA:HA	1.93	0.48
1:C:100:GLY:HA3	1:C:230:MET:O	2.12	0.48
1:C:216:LYS:NZ	1:E:210:ASN:O	2.42	0.48
2:D:158:ASP:OD1	2:D:160:PRO:HD2	2.14	0.48
2:D:125:GLN:NE2	2:D:155:GLY:HA2	2.24	0.48
2:B:21:TRP:HB2	2:B:41:THR:HG22	1.94	0.48
1:A:288:ILE:HD11	1:A:297:ILE:HD13	1.96	0.48
2:B:145:ASP:O	2:B:149:MET:HG2	2.14	0.47
1:A:14:CYS:HB2	2:B:25:HIS:HB3	1.97	0.47
2:B:127:ARG:NH2	2:F:133:LEU:O	2.47	0.47
2:D:38:LYS:O	2:D:42:GLN:N	2.43	0.47
1:E:44:GLU:HB2	1:E:292:MET:HG3	1.95	0.47
2:B:54:SER:OG	1:E:32:LYS:NZ	2.35	0.47
2:F:3:PHE:CE2	2:F:113:SER:HB2	2.50	0.47
1:E:56:VAL:HB	1:E:85:SER:HB3	1.96	0.47
2:F:17:MET:HE3	2:F:23:GLY:HA3	1.97	0.47
2:B:26:HIS:HD2	2:B:149:MET:HE3	1.80	0.46
1:A:170:ASN:HB2	1:A:237:LEU:HD23	1.98	0.46
1:A:29:ILE:HG13	1:A:30:MET:N	2.30	0.46
1:C:206:THR:CG2	1:C:207:SER:N	2.78	0.46
1:A:61:LEU:HD22	1:A:61:LEU:H	1.81	0.46
2:B:52:VAL:O	2:B:55:ILE:HG22	2.16	0.46
2:F:28:ASN:ND2	2:F:146:ASN:OD1	2.30	0.46
1:A:31:GLU:OE1	1:A:321:ARG:NH2	2.43	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:320:LEU:HB3	2:D:111:HIS:CD2	2.52	0.45
2:B:159:TYR:HB3	2:B:160:PRO:HD3	1.98	0.45
1:C:284:PRO:HD3	1:C:300:LEU:O	2.16	0.45
2:D:130:ALA:HB2	2:D:140:PHE:HB3	1.97	0.45
1:A:61:LEU:HA	1:A:79:PHE:CZ	2.52	0.45
1:C:307:LYS:HA	1:C:307:LYS:HD3	1.76	0.45
1:A:320:LEU:HB3	2:B:111:HIS:CD2	2.52	0.45
1:C:123:ILE:HG13	1:C:124:THR:HG22	1.99	0.45
2:B:1:GLY:HA3	2:B:112:ASP:OD2	2.17	0.45
1:A:166:ARG:HG3	8:A:1120:HOH:O	2.17	0.44
1:A:14:CYS:O	2:B:24:TYR:HA	2.18	0.44
1:E:295:HIS:HD2	1:E:306:PRO:HB2	1.82	0.44
5:I:3:NAG:H5	5:I:4:FUC:C6	2.46	0.44
1:E:74:PRO:HG3	1:E:147:PHE:O	2.16	0.44
2:D:55:ILE:HG23	2:D:99:LEU:HD21	2.00	0.44
1:E:100:GLY:HA3	1:E:230:MET:O	2.18	0.44
1:E:135:VAL:HG22	1:E:146:SER:HA	2.00	0.44
1:E:216:LYS:O	1:E:220:ARG:NH2	2.51	0.44
2:F:141:TYR:CE2	2:F:170:ARG:HD3	2.52	0.44
2:B:26:HIS:CD2	2:B:149:MET:CE	2.98	0.44
2:F:160:PRO:HA	2:F:163:SER:HB2	1.99	0.44
1:A:135:VAL:HG22	1:A:146:SER:HA	1.99	0.44
1:C:61:LEU:HA	1:C:79:PHE:CZ	2.53	0.44
1:A:124:THR:HA	1:A:125:PRO:HD3	1.88	0.43
2:D:140:PHE:CE2	2:D:144:CYS:HB2	2.53	0.43
1:C:190:GLU:OE2	4:H:2:SIA:O9	2.31	0.43
1:C:21:ASN:OD1	1:C:21:ASN:N	2.49	0.43
1:A:323:SER:HA	1:A:324:PRO:HD3	1.90	0.43
2:D:164:GLU:HG2	2:D:165:GLU:N	2.33	0.43
1:E:55:ASP:HB3	1:E:55(A):GLY:H	1.68	0.43
1:A:50:LYS:HB3	1:A:50:LYS:HE2	1.75	0.43
1:C:29:ILE:HD12	2:F:51:LYS:HG3	2.01	0.43
2:B:145:ASP:OD1	2:B:145:ASP:N	2.52	0.43
2:B:17:MET:HE1	2:B:23:GLY:HA3	2.01	0.43
2:D:4:GLY:O	2:D:8:GLY:HA3	2.19	0.43
2:B:4:GLY:O	2:B:8:GLY:HA3	2.19	0.43
1:C:125:PRO:HB2	1:C:125(B):ASN:OD1	2.18	0.43
1:C:9:PRO:HA	2:D:139:GLU:OE2	2.19	0.43
1:A:222:LYS:HZ3	6:J:1:NAG:H82	1.80	0.42
1:A:91:ILE:HG13	1:A:271:GLU:OE2	2.19	0.42
2:B:55:ILE:HD11	2:B:99:LEU:HG	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:132:ALA:HB1	1:C:152:VAL:HG11	2.01	0.42
1:E:176:LEU:HD22	1:E:257:ALA:HB1	2.00	0.42
1:E:268:MET:HE3	1:E:284:PRO:HA	2.01	0.42
1:E:283:THR:HG22	1:E:301:THR:HG22	2.01	0.42
1:C:101:ASN:OD1	1:C:231:GLU:HG3	2.19	0.42
1:E:15:ILE:HD13	2:F:119:TYR:HA	2.01	0.42
5:I:2:FUL:H5	5:I:3:NAG:O3	2.19	0.42
1:A:222:LYS:HA	1:A:226:LEU:O	2.19	0.42
2:D:164:GLU:H	2:D:164:GLU:CD	2.22	0.42
2:F:17:MET:HE3	2:F:36:ALA:HA	2.01	0.42
1:A:52:CYS:HB2	1:A:279:THR:HG22	2.01	0.42
1:C:24:GLU:OE1	1:C:39:ALA:HB3	2.19	0.42
1:E:124:THR:HA	1:E:125:PRO:HD2	1.88	0.42
1:E:28:THR:CG2	1:E:29:ILE:N	2.82	0.42
1:A:268:MET:HE3	1:A:284:PRO:HA	2.01	0.42
2:B:122:VAL:HG23	2:B:138:PHE:HE2	1.85	0.42
2:D:3:PHE:CD1	2:D:4:GLY:N	2.83	0.42
1:E:125:PRO:HG2	1:E:126:SER:HB2	2.02	0.42
1:C:55:ASP:O	1:C:278:ASN:OD1	2.37	0.41
1:C:124:THR:HA	1:C:125:PRO:HD2	1.92	0.41
2:B:68:ARG:NH2	2:F:83:LYS:HD3	2.34	0.41
1:C:25:GLN:HA	1:C:34:VAL:O	2.20	0.41
1:E:314:LEU:HA	1:E:314:LEU:HD23	1.92	0.41
1:A:284:PRO:HD3	1:A:300:LEU:O	2.20	0.41
2:D:166:ALA:O	2:D:170:ARG:HG2	2.21	0.41
2:F:145:ASP:O	2:F:149:MET:HG2	2.20	0.41
2:F:5:ALA:HB2	2:F:116:LYS:HB2	2.02	0.40
1:C:11:ASP:HB2	2:D:140:PHE:CE1	2.57	0.40
2:F:58:LYS:HA	2:F:58:LYS:HD3	1.88	0.40
1:C:38:HIS:C	1:C:318:THR:HG22	2.42	0.40
1:E:17:TYR:CZ	2:F:6:ILE:HG23	2.57	0.40
1:E:268:MET:CE	1:E:284:PRO:HA	2.52	0.40
2:B:164:GLU:H	2:B:164:GLU:CD	2.24	0.40
1:C:18:HIS:HB2	2:D:20:GLY:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	321/333 (96%)	309 (96%)	12 (4%)	0	100	100
1	C	321/333 (96%)	308 (96%)	13 (4%)	0	100	100
1	E	321/333 (96%)	309 (96%)	12 (4%)	0	100	100
2	B	173/180 (96%)	160 (92%)	13 (8%)	0	100	100
2	D	173/180 (96%)	160 (92%)	12 (7%)	1 (1%)	25	45
2	F	173/180 (96%)	158 (91%)	15 (9%)	0	100	100
All	All	1482/1539 (96%)	1404 (95%)	77 (5%)	1 (0%)	51	74

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	145	ASP

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	290/298 (97%)	279 (96%)	11 (4%)	33	57
1	C	290/298 (97%)	287 (99%)	3 (1%)	76	89
1	E	290/298 (97%)	285 (98%)	5 (2%)	60	80
2	B	149/154 (97%)	143 (96%)	6 (4%)	31	55
2	D	149/154 (97%)	146 (98%)	3 (2%)	55	77
2	F	149/154 (97%)	142 (95%)	7 (5%)	26	49

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1317/1356 (97%)	1282 (97%)	35 (3%)	44 69

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	124	THR
1	C	167	SER
1	C	206	THR
2	D	3	PHE
2	D	80	LEU
2	D	140	PHE
1	A	8	ASP
1	A	35	THR
1	A	51	LEU
1	A	61	LEU
1	A	91	ILE
1	A	166	ARG
1	A	192	THR
1	A	206	THR
1	A	213	LEU
1	A	283	THR
1	A	291	SER
2	B	10	ILE
2	B	18	VAL
2	B	26	HIS
2	B	41	THR
2	B	55	ILE
2	B	61	THR
1	E	26	VAL
1	E	29	ILE
1	E	80	LEU
1	E	193	ARG
1	E	206	THR
2	F	18	VAL
2	F	22	TYR
2	F	43	LYS
2	F	83	LYS
2	F	101	LEU
2	F	126	LEU
2	F	144	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such

sidechains are listed below:

Mol	Chain	Res	Type
2	D	25	HIS
2	D	125	GLN
1	A	47	HIS
2	B	25	HIS
2	B	26	HIS
2	B	125	GLN
2	F	142	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

17 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NAG	G	1	1,3	14,14,15	0.70	1 (7%)	17,19,21	1.00	0
3	FUL	G	2	3	10,10,11	1.58	2 (20%)	14,14,16	1.54	4 (28%)
3	NAG	G	3	3	14,14,15	0.45	0	17,19,21	0.42	0
4	GAL	H	1	4	11,11,12	1.85	3 (27%)	15,15,17	1.08	1 (6%)
4	SIA	H	2	4	17,20,21	0.60	0	21,28,31	1.51	2 (9%)
5	NAG	I	1	1,5	14,14,15	0.48	0	17,19,21	0.65	0
5	FUL	I	2	5	10,10,11	1.35	1 (10%)	14,14,16	1.41	2 (14%)
5	NAG	I	3	5	14,14,15	0.58	0	17,19,21	0.46	0
5	FUC	I	4	5	10,10,11	1.51	3 (30%)	14,14,16	2.00	4 (28%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NAG	J	1	6	15,15,15	0.57	1 (6%)	21,21,21	0.78	0
6	GAL	J	2	6	11,11,12	1.35	2 (18%)	15,15,17	1.87	2 (13%)
6	SIA	J	3	6	17,20,21	0.78	0	21,28,31	1.24	2 (9%)
3	NAG	K	1	1,3	14,14,15	0.70	1 (7%)	17,19,21	0.69	0
3	FUL	K	2	3	10,10,11	1.20	1 (10%)	14,14,16	0.83	0
3	NAG	K	3	3	14,14,15	0.31	0	17,19,21	0.38	0
4	GAL	L	1	4	11,11,12	1.83	5 (45%)	15,15,17	1.22	1 (6%)
4	SIA	L	2	4	17,20,21	0.67	0	21,28,31	1.38	3 (14%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	G	1	1,3	-	1/6/23/26	0/1/1/1
3	FUL	G	2	3	-	-	0/1/1/1
3	NAG	G	3	3	-	2/6/23/26	0/1/1/1
4	GAL	H	1	4	-	0/2/19/22	0/1/1/1
4	SIA	H	2	4	-	1/14/34/38	0/1/1/1
5	NAG	I	1	1,5	-	2/6/23/26	0/1/1/1
5	FUL	I	2	5	-	-	0/1/1/1
5	NAG	I	3	5	-	0/6/23/26	0/1/1/1
5	FUC	I	4	5	-	-	0/1/1/1
6	NAG	J	1	6	-	0/6/26/26	0/1/1/1
6	GAL	J	2	6	-	0/2/19/22	0/1/1/1
6	SIA	J	3	6	-	0/14/34/38	0/1/1/1
3	NAG	K	1	1,3	-	2/6/23/26	0/1/1/1
3	FUL	K	2	3	-	-	0/1/1/1
3	NAG	K	3	3	-	2/6/23/26	0/1/1/1
4	GAL	L	1	4	-	0/2/19/22	0/1/1/1
4	SIA	L	2	4	-	0/14/34/38	0/1/1/1

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	1	GAL	C4-C3	3.57	1.61	1.52
5	I	4	FUC	C1-C2	3.32	1.59	1.52
3	G	2	FUL	C1-C2	2.85	1.58	1.52
4	L	1	GAL	O3-C3	2.85	1.49	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	1	GAL	C2-C3	2.83	1.56	1.52
3	G	2	FUL	O5-C5	2.80	1.49	1.43
5	I	2	FUL	C4-C3	2.79	1.59	1.52
4	L	1	GAL	C4-C5	2.69	1.58	1.53
5	I	4	FUC	O5-C5	2.59	1.49	1.43
6	J	2	GAL	C1-C2	2.49	1.57	1.52
4	H	1	GAL	O3-C3	2.37	1.48	1.43
3	G	1	NAG	O5-C1	-2.33	1.40	1.43
4	L	1	GAL	C2-C3	2.29	1.55	1.52
3	K	1	NAG	O5-C1	-2.23	1.40	1.43
4	L	1	GAL	O5-C5	2.17	1.47	1.43
4	L	1	GAL	C1-C2	2.12	1.57	1.52
6	J	2	GAL	C2-C3	2.08	1.55	1.52
3	K	2	FUL	O5-C5	2.07	1.48	1.43
6	J	1	NAG	O5-C1	2.05	1.48	1.42
5	I	4	FUC	O5-C1	2.04	1.47	1.43

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	2	SIA	C6-O6-C2	5.32	122.73	111.34
6	J	2	GAL	C1-C2-C3	5.08	115.91	109.67
5	I	4	FUC	C1-O5-C5	3.98	121.79	112.78
6	J	2	GAL	C1-O5-C5	3.93	117.52	112.19
5	I	4	FUC	O5-C1-C2	3.74	116.55	110.77
5	I	4	FUC	C1-C2-C3	3.49	113.95	109.67
5	I	2	FUL	C1-C2-C3	-3.49	105.38	109.67
4	L	2	SIA	C6-O6-C2	3.43	118.67	111.34
4	H	2	SIA	C4-C3-C2	3.19	115.53	109.81
4	H	1	GAL	C1-C2-C3	2.96	113.30	109.67
3	G	2	FUL	C2-C3-C4	-2.91	105.85	110.89
5	I	2	FUL	O2-C2-C1	2.88	115.05	109.15
6	J	3	SIA	C4-C3-C2	2.80	114.83	109.81
4	L	2	SIA	C4-C3-C2	2.78	114.78	109.81
6	J	3	SIA	C6-O6-C2	2.72	117.16	111.34
3	G	2	FUL	O2-C2-C1	2.57	114.42	109.15
3	G	2	FUL	O5-C1-C2	2.32	114.36	110.77
4	L	1	GAL	C1-O5-C5	2.28	115.28	112.19
5	I	4	FUC	O5-C5-C4	2.22	113.51	109.52
4	L	2	SIA	C3-C2-C1	-2.20	107.11	111.93
3	G	2	FUL	C3-C4-C5	-2.07	106.55	109.77

There are no chirality outliers.

All (10) torsion outliers are listed below:

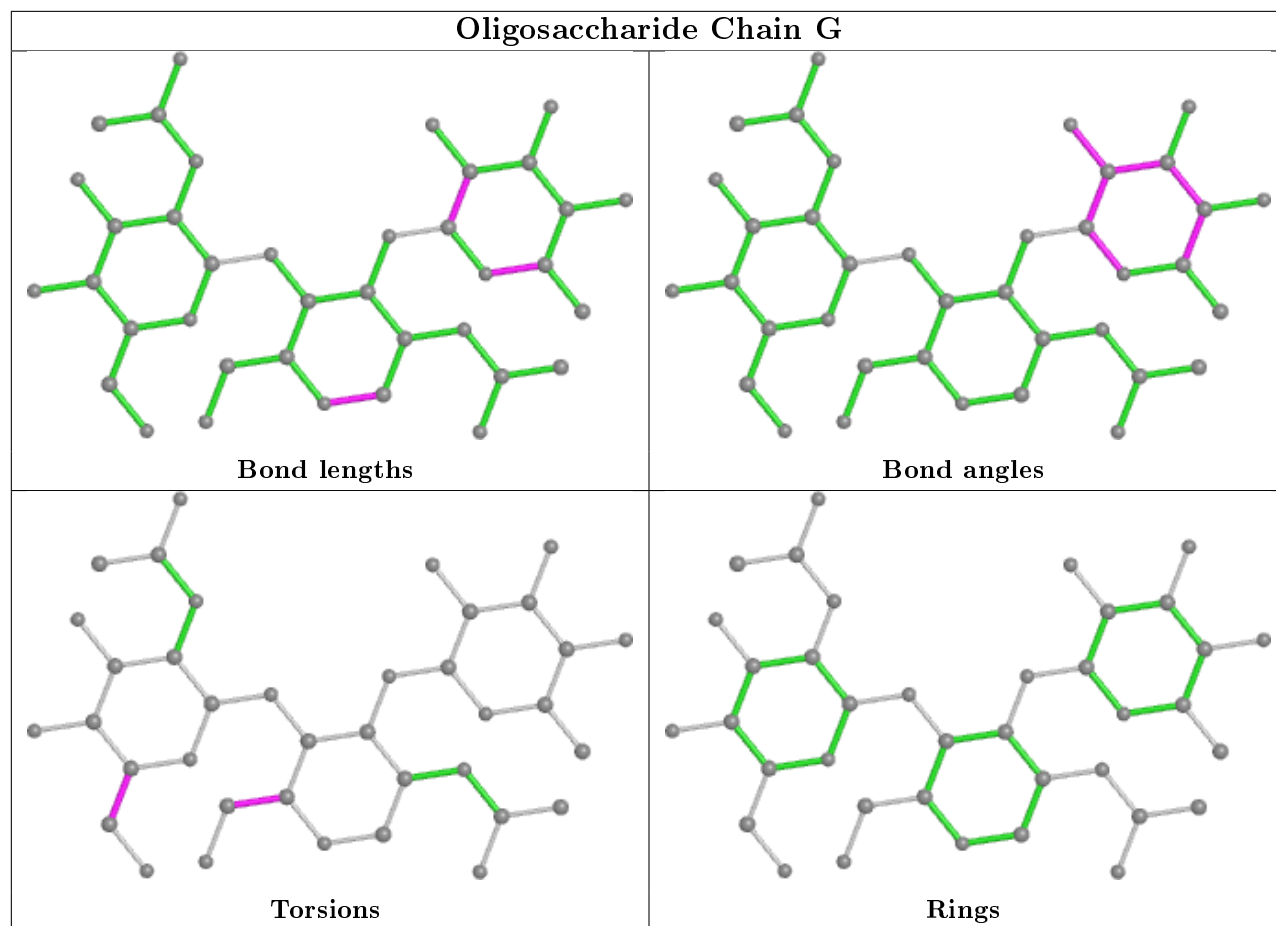
Mol	Chain	Res	Type	Atoms
3	K	3	NAG	C4-C5-C6-O6
3	G	3	NAG	C4-C5-C6-O6
3	K	3	NAG	O5-C5-C6-O6
5	I	1	NAG	O5-C5-C6-O6
5	I	1	NAG	C4-C5-C6-O6
3	K	1	NAG	C4-C5-C6-O6
3	G	3	NAG	O5-C5-C6-O6
3	K	1	NAG	O5-C5-C6-O6
4	H	2	SIA	C6-C7-C8-O8
3	G	1	NAG	C4-C5-C6-O6

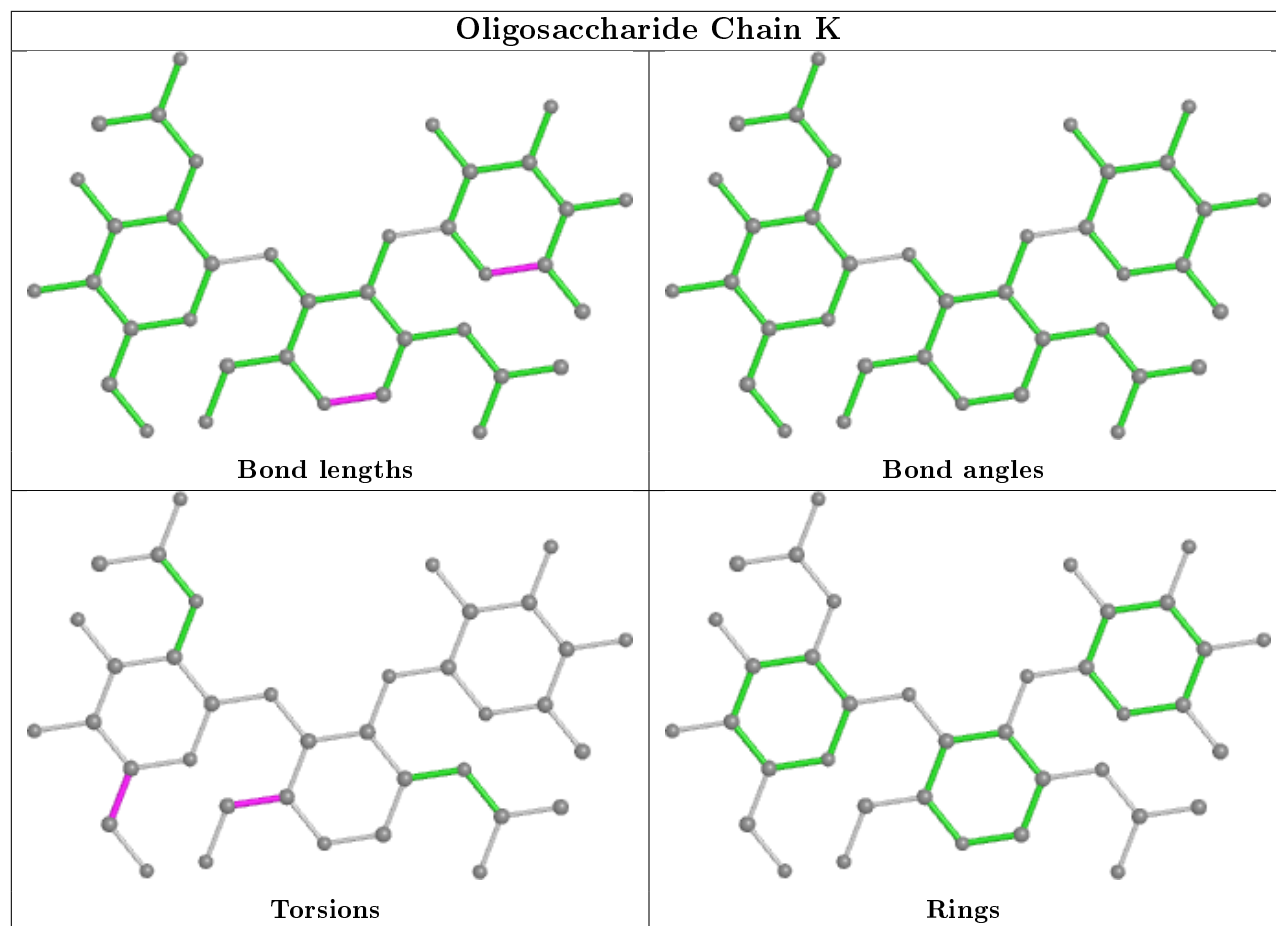
There are no ring outliers.

5 monomers are involved in 7 short contacts:

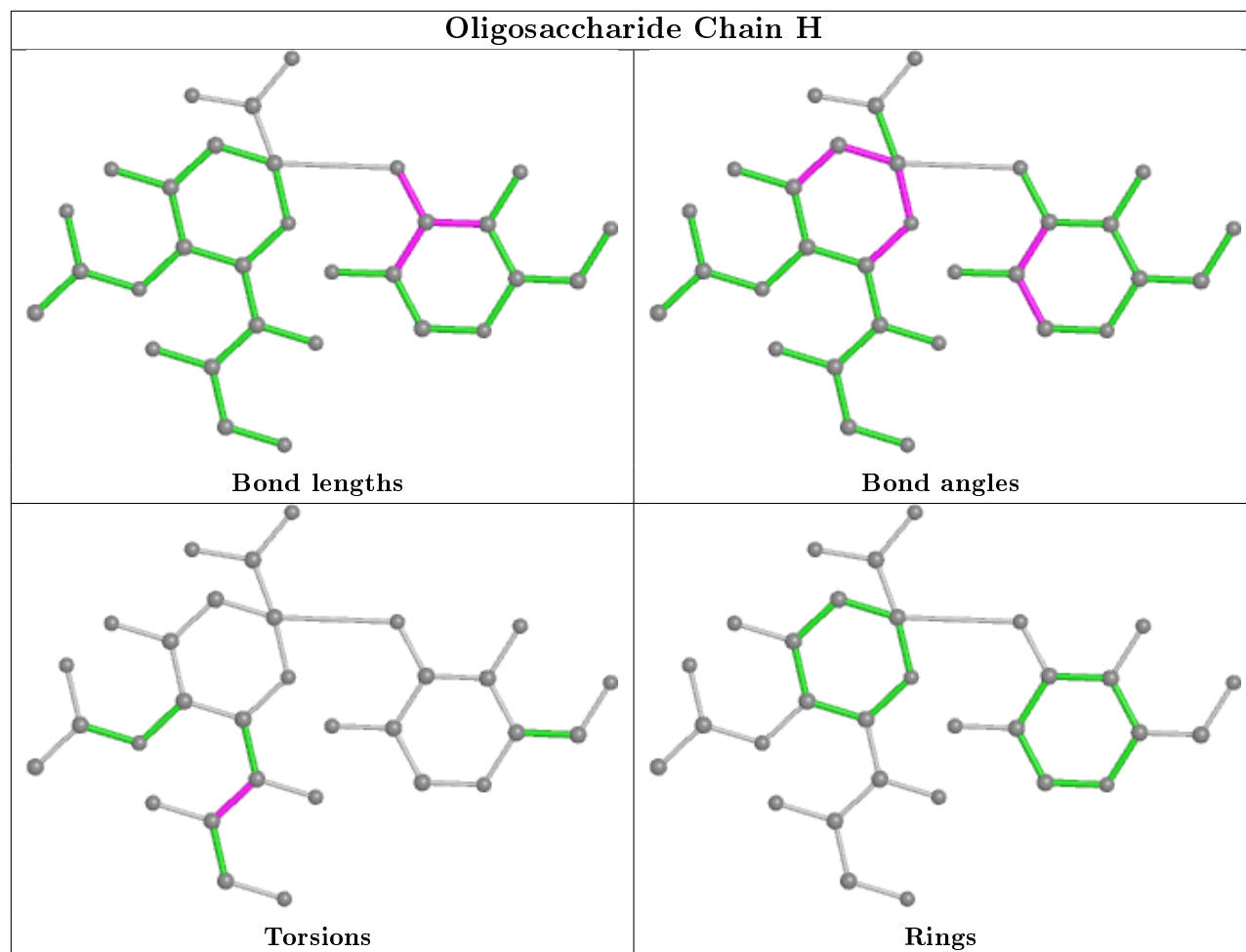
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	J	1	NAG	3	0
5	I	2	FUL	1	0
5	I	4	FUC	2	0
4	H	2	SIA	1	0
5	I	3	NAG	3	0

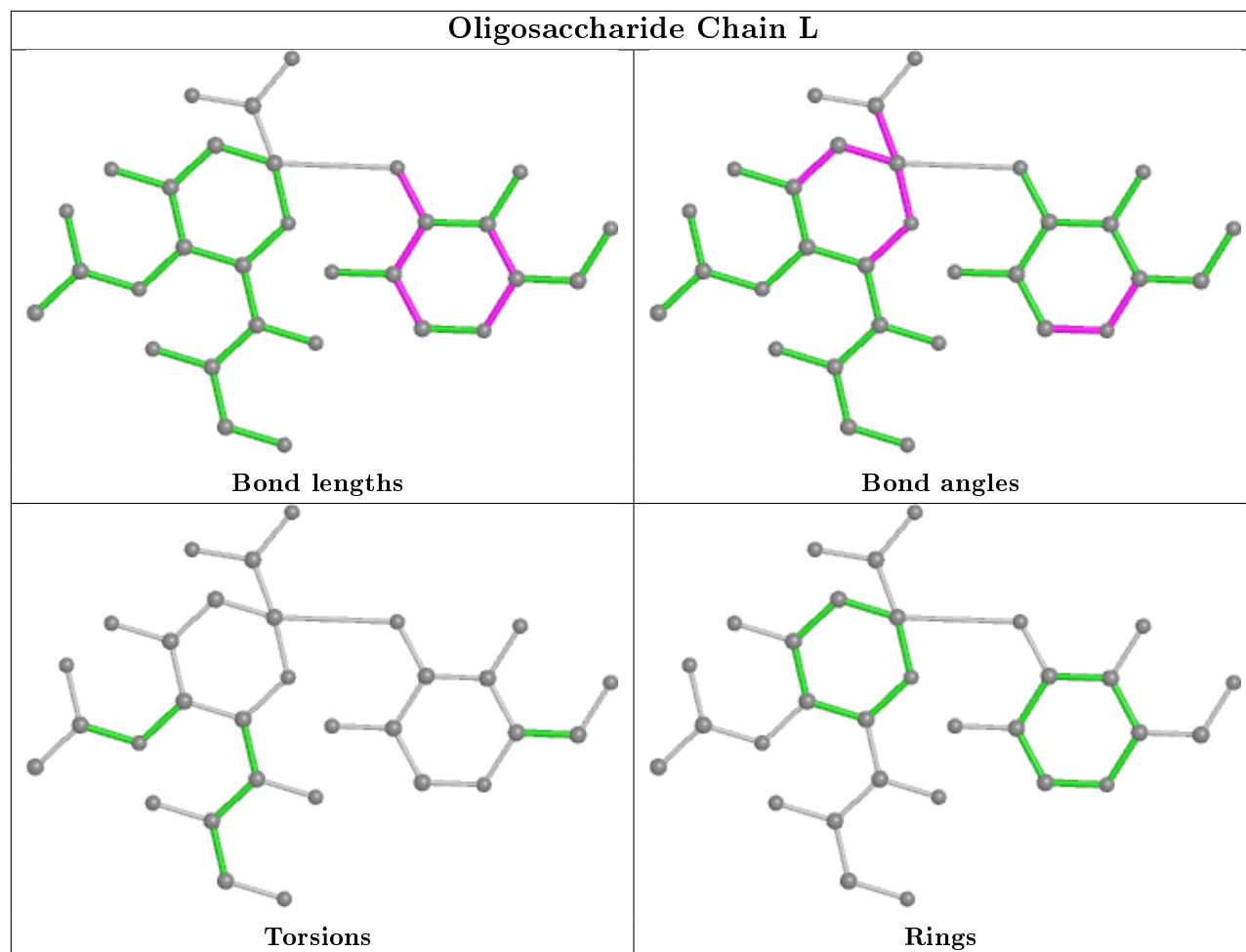
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

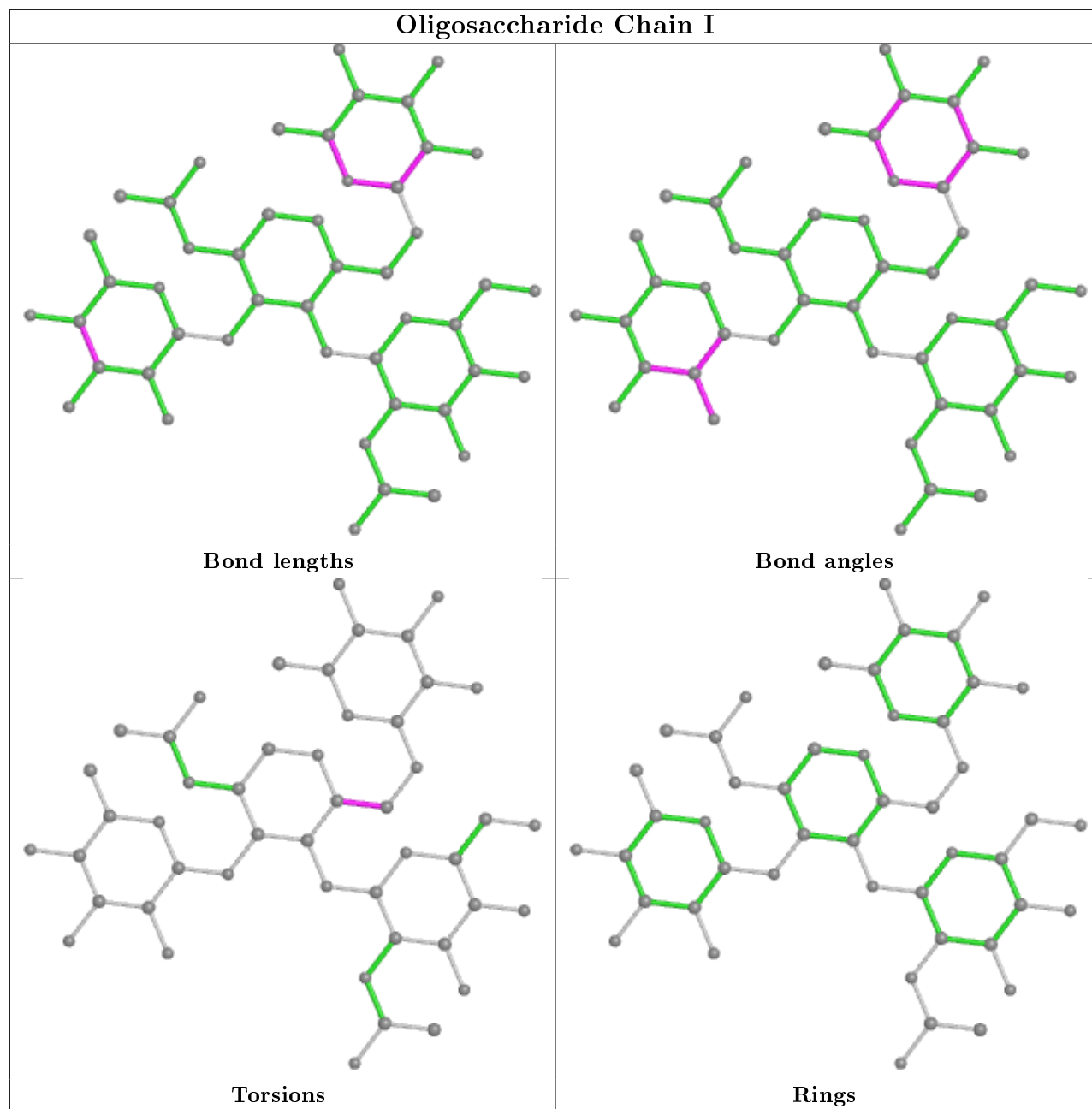


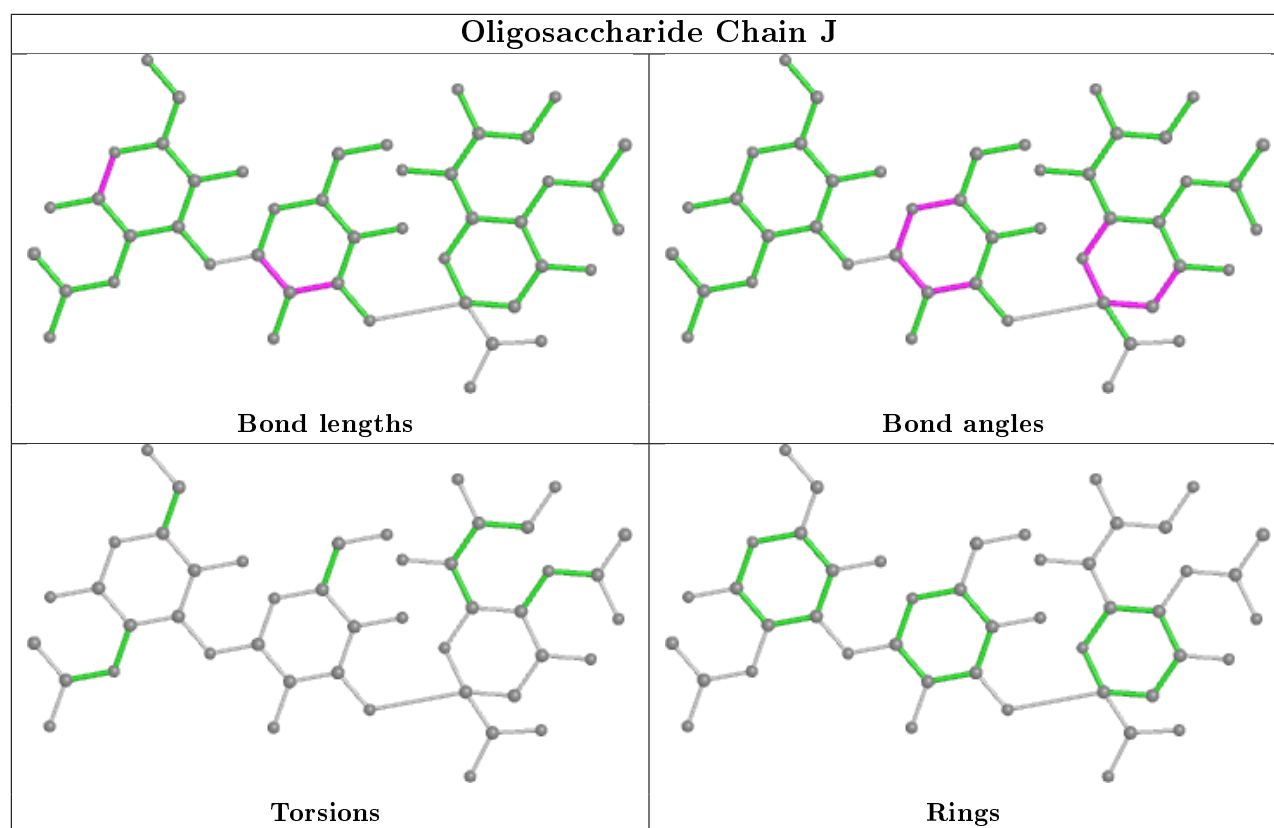


Oligosaccharide Chain H









5.6 Ligand geometry [i](#)

3 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	$\# Z > 2$
7	NAG	A	1001	1	14,14,15	0.78	1 (7%)	17,19,21	0.51	0
7	NAG	C	1001	1	14,14,15	0.38	0	17,19,21	0.47	0
7	NAG	E	1001	1	14,14,15	0.50	0	17,19,21	0.55	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	NAG	A	1001	1	-	2/6/23/26	0/1/1/1
7	NAG	C	1001	1	-	2/6/23/26	0/1/1/1
7	NAG	E	1001	1	-	4/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	1001	NAG	O5-C1	2.12	1.47	1.43

There are no bond angle outliers.

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	E	1001	NAG	C4-C5-C6-O6
7	C	1001	NAG	C4-C5-C6-O6
7	E	1001	NAG	O5-C5-C6-O6
7	C	1001	NAG	O5-C5-C6-O6
7	A	1001	NAG	C1-C2-N2-C7
7	E	1001	NAG	C1-C2-N2-C7
7	A	1001	NAG	C3-C2-N2-C7
7	E	1001	NAG	C3-C2-N2-C7

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	323/333 (96%)	-0.33	3 (0%) 84 82	18, 40, 92, 124	0
1	C	323/333 (96%)	-0.41	7 (2%) 62 57	22, 42, 77, 128	0
1	E	323/333 (96%)	-0.38	3 (0%) 84 82	27, 51, 76, 119	0
2	B	175/180 (97%)	0.48	16 (9%) 9 6	24, 87, 110, 115	0
2	D	175/180 (97%)	0.74	30 (17%) 1 1	26, 89, 130, 134	0
2	F	175/180 (97%)	0.19	7 (4%) 38 32	23, 77, 106, 122	0
All	All	1494/1539 (97%)	-0.08	66 (4%) 34 28	18, 52, 110, 134	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	175	GLY	7.3
2	D	138	PHE	4.4
2	D	32	SER	4.4
2	B	157	TYR	4.4
2	D	168	LEU	4.3
2	D	33	GLY	4.2
2	F	173	ILE	4.1
2	D	155	GLY	4.0
2	B	35	ALA	3.6
2	F	11	GLU	3.5
1	E	9	PRO	3.5
1	E	8	ASP	3.4
2	B	33	GLY	3.3
1	C	324	PRO	3.3
1	C	9	PRO	3.2
2	B	174	SER	3.2
2	D	31	GLY	3.2
2	F	174	SER	3.2
1	C	14	CYS	3.1

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Mol	Chain	Res	Type	RSRZ
2	D	166	ALA	3.1
2	B	31	GLY	3.0
2	D	139	GLU	3.0
2	B	147	GLU	3.0
2	D	146	ASN	2.9
2	D	27	SER	2.9
2	D	156	THR	2.9
2	D	34	TYR	2.9
2	B	13	GLY	2.9
2	B	14	TRP	2.8
2	D	158	ASP	2.8
2	D	130	ALA	2.8
1	C	12	GLN	2.8
1	C	10	GLY	2.8
2	B	29	GLU	2.8
2	D	18	VAL	2.7
2	D	160	PRO	2.7
2	D	126	LEU	2.7
2	D	35	ALA	2.7
2	B	162	TYR	2.6
1	A	8	ASP	2.6
1	E	324	PRO	2.6
2	B	27	SER	2.5
1	A	12	GLN	2.5
1	C	8	ASP	2.5
2	B	25	HIS	2.5
2	D	157	TYR	2.5
2	F	141	TYR	2.4
2	D	133	LEU	2.4
2	B	133	LEU	2.4
2	D	38	LYS	2.3
2	B	156	THR	2.3
2	D	172	GLU	2.3
2	D	1	GLY	2.3
2	D	161	GLN	2.3
2	B	15	GLN	2.2
2	D	26	HIS	2.2
1	C	144	ARG	2.2
2	D	148	CYS	2.2
2	D	145	ASP	2.1
2	D	173	ILE	2.1
2	F	131	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
2	D	16	GLY	2.1
1	A	9	PRO	2.0
2	D	19	ASP	2.0
2	B	146	ASN	2.0
2	F	23	GLY	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](#)

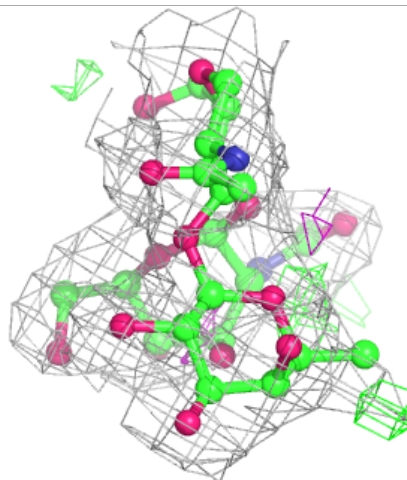
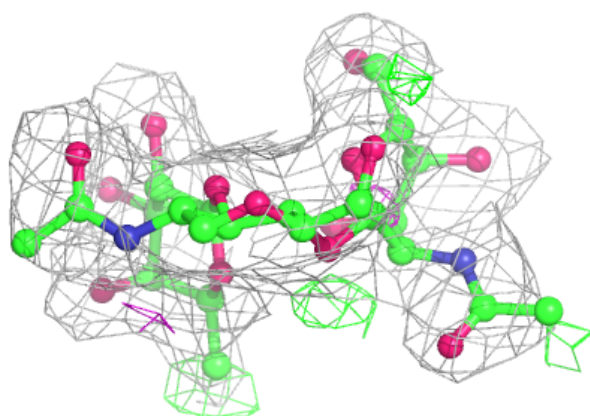
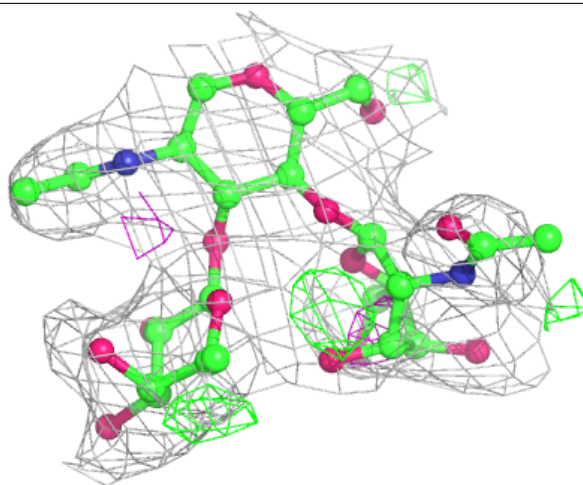
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	NAG	J	1	15/15	0.79	0.26	85,94,101,104	0
4	GAL	H	1	11/12	0.80	0.41	76,80,87,90	0
5	NAG	I	3	14/15	0.82	0.29	78,86,93,94	0
5	FUL	I	2	10/11	0.83	0.29	78,85,87,90	0
3	FUL	K	2	10/11	0.85	0.34	90,96,103,108	0
3	FUL	G	2	10/11	0.86	0.17	72,77,83,86	0
3	NAG	G	3	14/15	0.87	0.24	74,81,87,87	0
4	GAL	L	1	11/12	0.87	0.28	60,74,83,85	0
5	NAG	I	1	14/15	0.89	0.17	52,69,76,81	0
3	NAG	K	3	14/15	0.90	0.29	89,97,103,107	0
4	SIA	H	2	20/21	0.90	0.29	50,68,78,79	0
5	FUC	I	4	10/11	0.91	0.32	59,69,78,81	0
6	GAL	J	2	11/12	0.91	0.15	51,58,70,73	0
3	NAG	K	1	14/15	0.91	0.14	52,72,84,87	0
4	SIA	L	2	20/21	0.92	0.13	42,57,68,70	0
3	NAG	G	1	14/15	0.94	0.14	50,61,71,75	0
6	SIA	J	3	20/21	0.97	0.10	25,36,42,45	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.

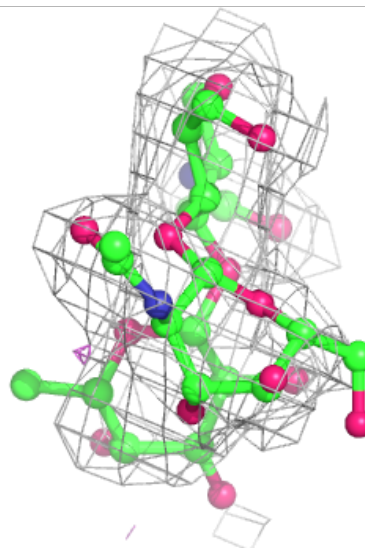
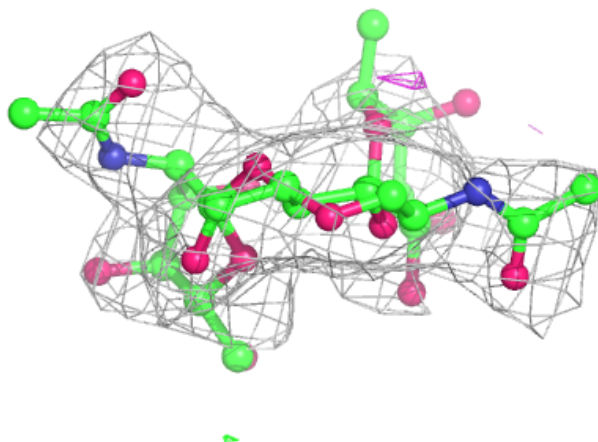
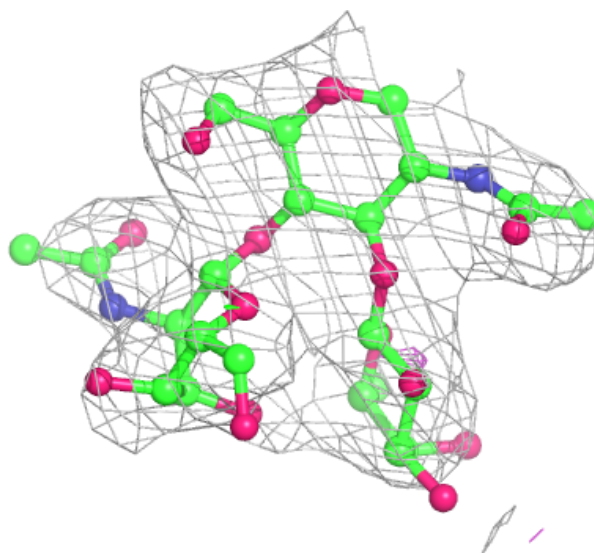
Electron density around Chain G:

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



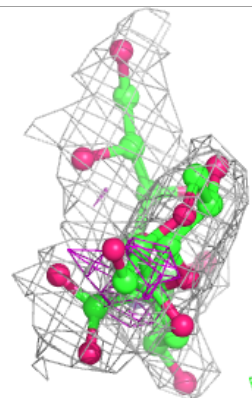
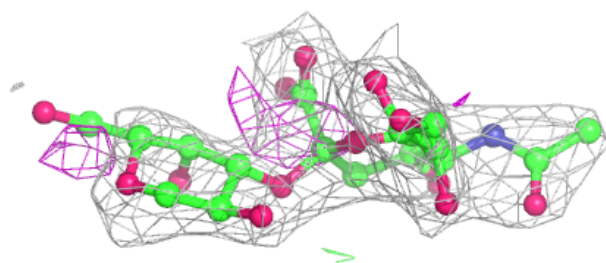
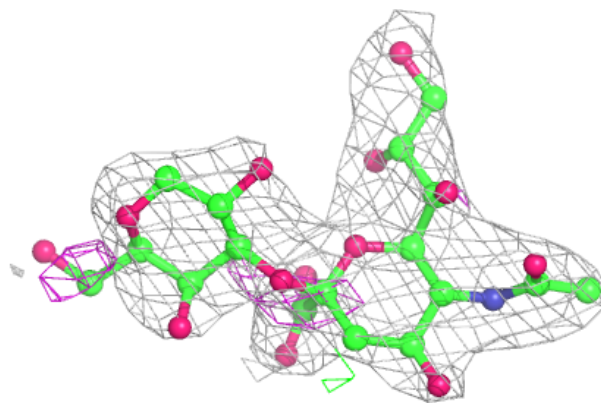
Electron density around Chain K:

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



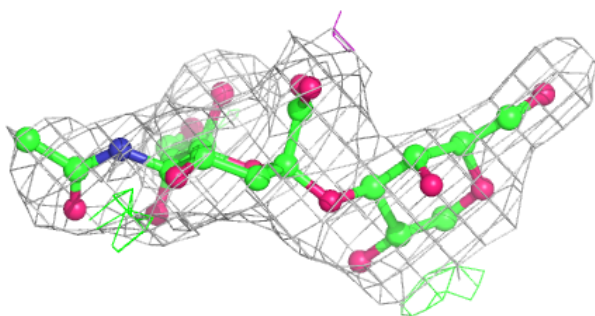
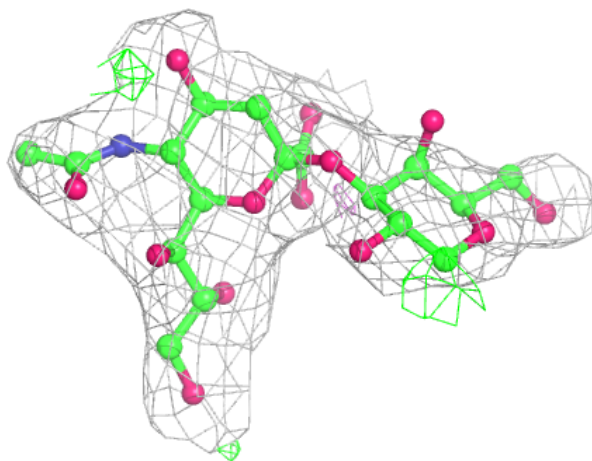
Electron density around Chain H:

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



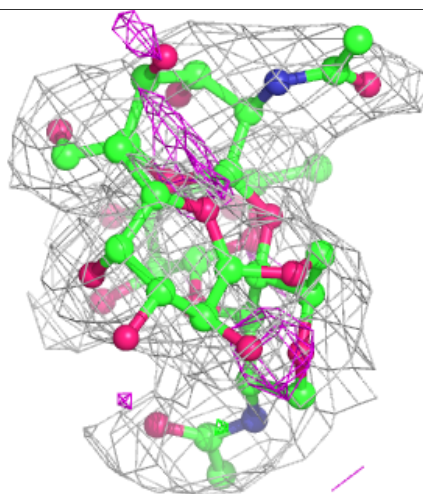
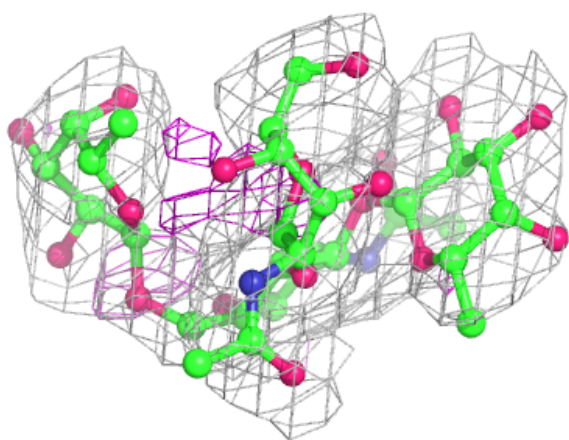
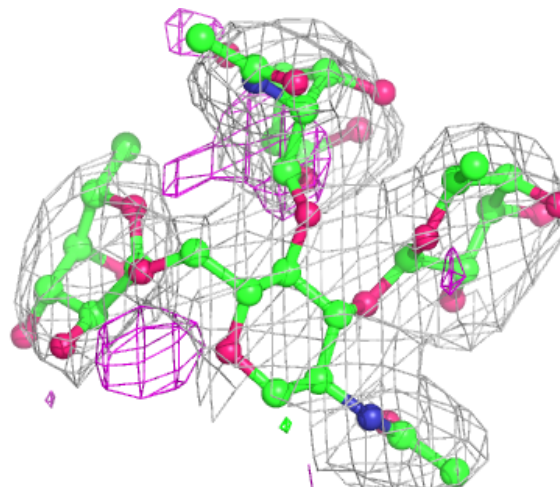
Electron density around Chain L:

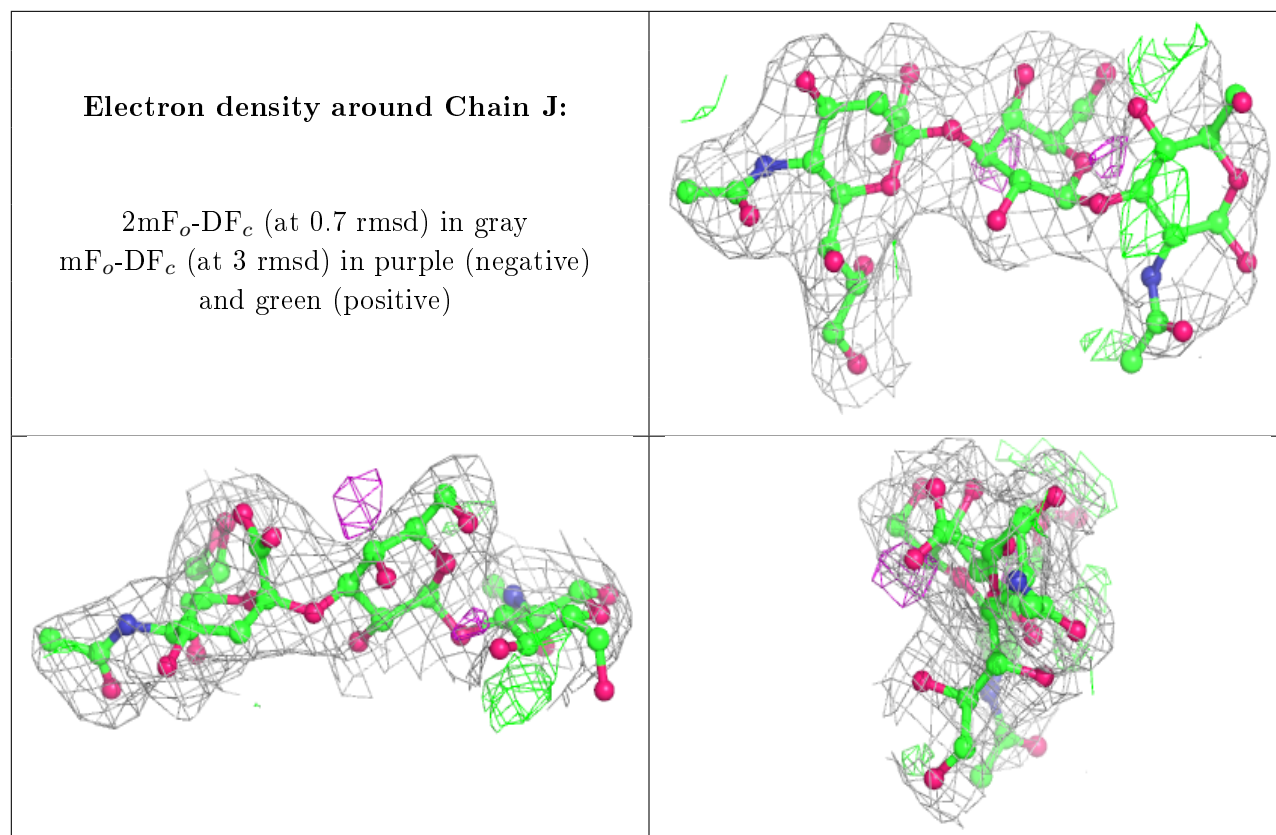
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain I:

$2mF_o-DF_c$ (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
7	NAG	E	1001	14/15	0.84	0.24	95,101,103,105	0
7	NAG	A	1001	14/15	0.85	0.15	90,103,107,107	0
7	NAG	C	1001	14/15	0.88	0.18	83,90,92,94	0

6.5 Other polymers [i](#)

There are no such residues in this entry.