



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 26, 2021 – 12:04 PM EDT

PDB ID : 4XQ5  
Title : Human-infecting H10N8 influenza virus retains strong preference for avian-type receptors  
Authors : Tzarum, N.; Zhang, H.; Zhu, X.; Wilson, I.A.  
Deposited on : 2015-01-19  
Resolution : 2.59 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](mailto: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.

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

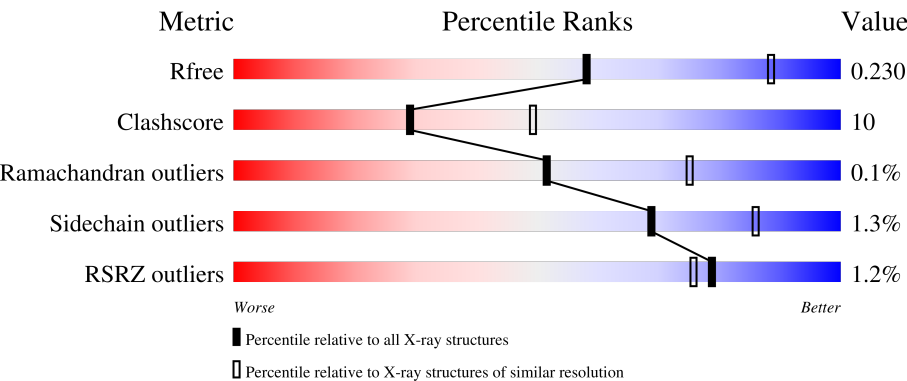
MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.23.1  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.23.1

# 1 Overall quality at a glance i

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

The reported resolution of this entry is 2.59 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 <sub>free</sub>	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	323	<div><div></div><div>81%17%.</div></div>
1	C	323	<div><div>2%</div><div>83%16%.</div></div>
1	E	323	<div><div></div><div>80%19%.</div></div>
2	B	180	<div><div>3%</div><div>71%24%5%</div></div>
2	D	180	<div><div>%</div><div>71%23%6%</div></div>

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Mol	Chain	Length	Quality of chain
2	F	180	<div><div></div><div>3%</div><div>69%</div><div>23%</div><div>6%</div></div>
3	G	2	<div><div></div><div>100%</div></div>
3	H	2	<div><div></div><div>50%</div><div>50%</div></div>
3	I	2	<div><div></div><div>50%</div><div>50%</div></div>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11820 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 HA1 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	318	Total	C	N	O	S	0	0	0
			2437	1506	449	465	17			
1	C	319	Total	C	N	O	S	0	0	0
			2445	1512	450	466	17			
1	E	318	Total	C	N	O	S	0	0	0
			2437	1506	449	465	17			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	7	ALA	-	expression tag	UNP A0A059T4A1
A	8	ASP	-	expression tag	UNP A0A059T4A1
A	9	PRO	-	expression tag	UNP A0A059T4A1
A	10	GLY	-	expression tag	UNP A0A059T4A1
C	7	ALA	-	expression tag	UNP A0A059T4A1
C	8	ASP	-	expression tag	UNP A0A059T4A1
C	9	PRO	-	expression tag	UNP A0A059T4A1
C	10	GLY	-	expression tag	UNP A0A059T4A1
E	7	ALA	-	expression tag	UNP A0A059T4A1
E	8	ASP	-	expression tag	UNP A0A059T4A1
E	9	PRO	-	expression tag	UNP A0A059T4A1
E	10	GLY	-	expression tag	UNP A0A059T4A1

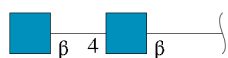
- Molecule 2 is a protein called Hemagglutinin HA2 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	171	Total	C	N	O	S	0	0	0
			1382	854	239	281	8			
2	D	169	Total	C	N	O	S	0	0	0
			1367	846	236	277	8			
2	F	170	Total	C	N	O	S	0	0	0
			1378	852	238	280	8			

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	175	SER	-	expression tag	UNP A0A059T4A1
B	176	GLY	-	expression tag	UNP A0A059T4A1
B	177	ARG	-	expression tag	UNP A0A059T4A1
B	178	LEU	-	expression tag	UNP A0A059T4A1
B	179	VAL	-	expression tag	UNP A0A059T4A1
B	180	PRO	-	expression tag	UNP A0A059T4A1
B	181	ARG	-	expression tag	UNP A0A059T4A1
D	175	SER	-	expression tag	UNP A0A059T4A1
D	176	GLY	-	expression tag	UNP A0A059T4A1
D	177	ARG	-	expression tag	UNP A0A059T4A1
D	178	LEU	-	expression tag	UNP A0A059T4A1
D	179	VAL	-	expression tag	UNP A0A059T4A1
D	180	PRO	-	expression tag	UNP A0A059T4A1
D	181	ARG	-	expression tag	UNP A0A059T4A1
F	175	SER	-	expression tag	UNP A0A059T4A1
F	176	GLY	-	expression tag	UNP A0A059T4A1
F	177	ARG	-	expression tag	UNP A0A059T4A1
F	178	LEU	-	expression tag	UNP A0A059T4A1
F	179	VAL	-	expression tag	UNP A0A059T4A1
F	180	PRO	-	expression tag	UNP A0A059T4A1
F	181	ARG	-	expression tag	UNP A0A059T4A1

- Molecule 3 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
3	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	H	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	I	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



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

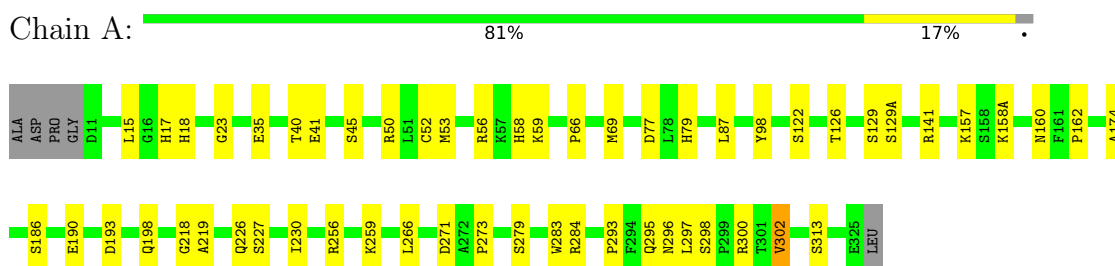
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	77	Total	O	0	0
			77	77		
5	B	26	Total	O	0	0
			26	26		
5	C	76	Total	O	0	0
			76	76		
5	D	15	Total	O	0	0
			15	15		
5	E	53	Total	O	0	0
			53	53		
5	F	15	Total	O	0	0
			15	15		

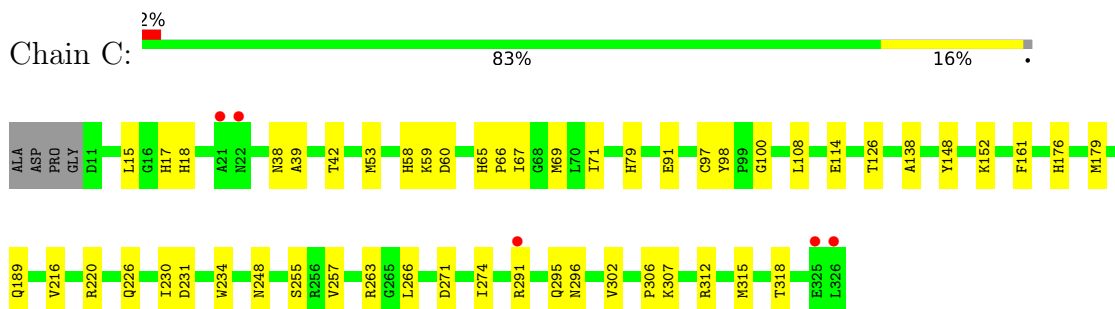
### 3 Residue-property plots

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

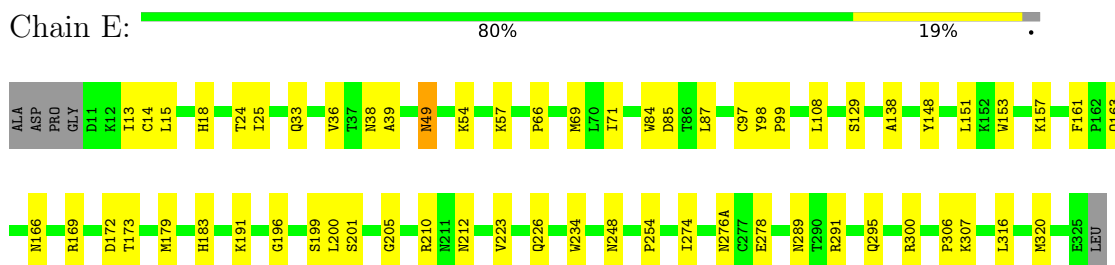
- Molecule 1: Hemagglutinin HA1 chain



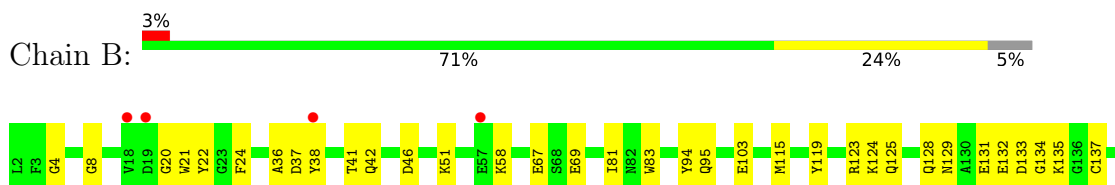
- Molecule 1: Hemagglutinin HA1 chain



- Molecule 1: Hemagglutinin HA1 chain

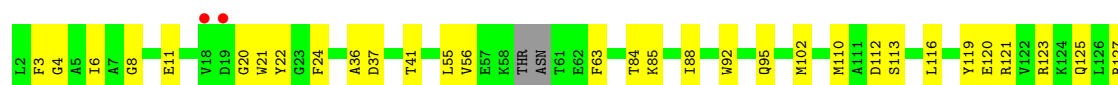


- Molecule 2: Hemagglutinin HA2 chain





• Molecule 2: Hemagglutinin HA2 chain



• Molecule 2: Hemagglutinin HA2 chain



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



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



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





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	64.07Å 242.99Å 70.38Å 90.00° 111.26° 90.00°	Depositor
Resolution (Å)	48.06 – 2.59 48.06 – 2.59	Depositor EDS
% Data completeness (in resolution range)	97.6 (48.06-2.59) 94.1 (48.06-2.59)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.53 (at 2.58Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, $R_{free}$	0.191 , 0.230 0.194 , 0.230	Depositor DCC
$R_{free}$ test set	3010 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	26.9	Xtriage
Anisotropy	0.364	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 37.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	11820	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: 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	0.69	1/2486 (0.0%)	0.76	0/3368
1	C	0.60	0/2494	0.68	0/3379
1	E	0.63	0/2486	0.69	0/3368
2	B	0.56	0/1407	0.66	0/1899
2	D	0.56	0/1391	0.65	0/1875
2	F	0.53	0/1402	0.71	2/1891 (0.1%)
All	All	0.61	1/11666 (0.0%)	0.70	2/15780 (0.0%)

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	E	0	1
2	F	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	190	GLU	CG-CD	6.41	1.61	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	144	CYS	CA-CB-SG	5.75	124.34	114.00
2	F	31	GLY	N-CA-C	5.13	125.94	113.10

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	84	TRP	Peptide
2	F	143	ALA	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	2437	0	2385	47	0
1	C	2445	0	2396	44	0
1	E	2437	0	2386	44	0
2	B	1382	0	1279	36	0
2	D	1367	0	1265	40	0
2	F	1378	0	1274	48	0
3	G	28	0	25	0	0
3	H	28	0	25	0	0
3	I	28	0	25	1	0
4	A	14	0	13	0	0
4	C	14	0	13	0	0
5	A	77	0	0	3	0
5	B	26	0	0	2	0
5	C	76	0	0	6	0
5	D	15	0	0	0	0
5	E	53	0	0	3	0
5	F	15	0	0	4	0
All	All	11820	0	11086	218	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 10.

All (218) 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:E:49:ASN:OD1	5:E:738:HOH:O	1.56	1.19
2:F:29:ALA:O	2:F:30:GLN:NE2	1.95	0.99
2:F:134:GLY:O	2:F:135:LYS:HD2	1.65	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:119:TYR:OH	2:D:132:GLU:OE2	1.87	0.92
2:F:8:GLY:O	5:F:205:HOH:O	1.95	0.85
1:A:271:ASP:OD1	1:A:284:ARG:NH1	2.11	0.83
1:C:291:ARG:HD2	2:D:56:VAL:HB	1.63	0.81
2:F:125:GLN:HE22	2:F:155:ASN:HA	1.46	0.81
1:A:41:GLU:OE2	1:A:313:SER:OG	2.00	0.78
1:A:295:GLN:NE2	1:A:298:SER:H	1.81	0.78
1:C:53:MET:HE2	1:C:58:HIS:HB3	1.65	0.77
2:F:165:GLU:O	2:F:169:ASN:ND2	2.21	0.73
1:E:278:GLU:O	1:E:289:ASN:ND2	2.19	0.72
2:B:133:ASP:OD1	2:B:134:GLY:N	2.21	0.72
1:A:18:HIS:CD2	2:B:21:TRP:HA	2.24	0.71
1:C:18:HIS:CD2	2:D:21:TRP:HA	2.26	0.71
1:A:129:SER:O	1:A:157:LYS:HE2	1.91	0.71
2:D:120:GLU:OE1	2:D:123:ARG:NH2	2.24	0.70
2:D:133:ASP:OD1	2:D:137:CYS:HB2	1.91	0.70
1:E:295:GLN:HG2	1:E:306:PRO:HG2	1.74	0.70
1:A:77:ASP:O	5:A:529:HOH:O	2.10	0.69
1:A:300:ARG:HH11	2:B:67:GLU:HB3	1.56	0.69
2:B:95:GLN:HE21	2:F:95:GLN:HE22	1.40	0.69
2:F:125:GLN:NE2	2:F:155:ASN:OD1	2.26	0.69
2:F:144:CYS:HA	2:F:145:ASP:HB2	1.74	0.68
1:A:18:HIS:HD2	2:B:20:GLY:O	1.77	0.68
1:E:98:TYR:CE2	1:E:226:GLN:HG2	2.30	0.67
2:F:44:ALA:HA	2:F:110:MET:HE1	1.76	0.66
1:A:284:ARG:HG2	1:A:284:ARG:HH11	1.61	0.66
2:D:121:ARG:HD3	2:D:155:ASN:HD21	1.60	0.66
2:B:119:TYR:OH	2:B:132:GLU:OE1	2.07	0.65
1:C:189:GLN:NE2	5:C:647:HOH:O	2.31	0.64
2:F:44:ALA:HA	2:F:110:MET:CE	2.28	0.64
1:C:98:TYR:CE2	1:C:226:GLN:HG2	2.33	0.64
2:B:133:ASP:OD2	2:B:137:CYS:HB2	1.98	0.63
1:A:266:LEU:HD11	1:A:302:VAL:HG13	1.81	0.63
1:A:295:GLN:HE21	1:A:297:LEU:H	1.47	0.63
2:F:4:GLY:O	2:F:8:GLY:HA3	1.98	0.62
1:A:50:ARG:NH2	1:A:273:PRO:HG2	2.13	0.62
2:B:37:ASP:O	2:B:41:THR:HG23	1.99	0.62
2:F:134:GLY:C	2:F:135:LYS:HD2	2.20	0.62
1:C:291:ARG:CD	2:D:56:VAL:HB	2.28	0.62
1:C:231:ASP:OD2	1:E:210:ARG:NH1	2.33	0.61
1:A:295:GLN:HE21	1:A:297:LEU:N	1.99	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:117:ASN:ND2	5:F:212:HOH:O	2.34	0.60
2:D:121:ARG:HD3	2:D:155:ASN:ND2	2.16	0.60
1:C:38:ASN:HB3	1:C:318:THR:HG23	1.84	0.60
1:E:191:LYS:NZ	1:E:196:GLY:O	2.34	0.59
2:F:21:TRP:H	2:F:41:THR:HG21	1.67	0.59
1:C:315:MET:O	5:C:626:HOH:O	2.17	0.59
2:D:127:ARG:HD3	2:D:159:HIS:CE1	2.38	0.59
2:F:133:ASP:OD1	2:F:137:CYS:O	2.20	0.58
1:C:53:MET:HE2	1:C:58:HIS:CB	2.32	0.58
1:E:300:ARG:HH11	1:E:300:ARG:HB2	1.69	0.58
1:E:179:MET:HG2	1:E:234:TRP:HB3	1.84	0.58
2:F:28:ASN:HD22	2:F:145:ASP:H	1.51	0.58
1:E:99:PRO:HB3	1:E:223:VAL:HB	1.86	0.57
1:A:295:GLN:HE22	1:A:298:SER:H	1.49	0.57
1:C:296:ASN:ND2	5:C:648:HOH:O	2.36	0.57
2:D:125:GLN:OE1	2:D:155:ASN:HA	2.04	0.57
1:C:295:GLN:HG2	1:C:306:PRO:HG2	1.87	0.57
2:B:148:CYS:O	2:B:151:SER:HB3	2.05	0.57
2:F:28:ASN:OD1	2:F:31:GLY:HA3	2.05	0.56
1:A:283:TRP:CE3	1:A:298:SER:HB2	2.40	0.56
2:D:37:ASP:O	2:D:41:THR:HG23	2.06	0.56
2:F:145:ASP:HB3	2:F:148:CYS:HB3	1.88	0.56
1:C:100:GLY:HA3	1:C:230:ILE:O	2.06	0.56
1:A:53:MET:HE2	1:A:58:HIS:CB	2.36	0.56
2:B:58:LYS:HE3	5:F:210:HOH:O	2.04	0.56
2:B:141:TYR:O	2:B:166:ALA:HA	2.06	0.56
1:C:266:LEU:HD11	1:C:302:VAL:HG13	1.88	0.56
2:D:21:TRP:H	2:D:41:THR:HG21	1.71	0.56
1:C:60:ASP:HB2	1:C:274:ILE:HD11	1.88	0.56
1:A:284:ARG:HH11	1:A:284:ARG:CG	2.19	0.55
1:A:53:MET:CE	1:A:58:HIS:HB3	2.37	0.55
1:A:293:PRO:O	5:A:543:HOH:O	2.18	0.55
2:D:20:GLY:HA3	2:D:36:ALA:HB1	1.89	0.55
2:F:99:LEU:O	2:F:103:GLU:HB2	2.07	0.55
1:C:271:ASP:HB2	5:C:663:HOH:O	2.05	0.55
2:D:4:GLY:O	2:D:8:GLY:HA3	2.06	0.55
2:B:131:GLU:HG3	2:B:141:TYR:CE1	2.42	0.55
1:C:216:VAL:HG11	1:E:212:ASN:HB2	1.87	0.55
1:A:53:MET:HE1	1:A:56:ARG:HD2	1.89	0.54
1:C:15:LEU:HD13	2:D:119:TYR:HA	1.88	0.54
2:B:51:LYS:HD3	2:B:103:GLU:HG2	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:220:ARG:HG2	1:E:205:GLY:HA3	1.90	0.54
2:F:21:TRP:H	2:F:41:THR:CG2	2.21	0.54
1:A:53:MET:HE2	1:A:58:HIS:HB3	1.90	0.53
1:A:98:TYR:CE2	1:A:226:GLN:HG2	2.43	0.53
2:B:95:GLN:HE21	2:F:95:GLN:NE2	2.06	0.53
1:C:59:LYS:HE3	1:C:79:HIS:ND1	2.23	0.53
1:A:174:ALA:HB3	1:A:259:LYS:HE3	1.91	0.53
1:E:18:HIS:CD2	2:F:21:TRP:HA	2.44	0.53
1:A:59:LYS:HE3	1:A:79:HIS:ND1	2.24	0.53
1:E:71:ILE:HD13	1:E:179:MET:SD	2.48	0.53
1:E:161:PHE:HB3	1:E:248:ASN:O	2.08	0.53
1:A:17:HIS:HB3	2:B:115:MET:CE	2.39	0.53
1:A:271:ASP:CG	1:A:284:ARG:NH1	2.63	0.52
1:E:36:VAL:HG12	1:E:38:ASN:H	1.74	0.52
1:A:53:MET:HE1	1:A:56:ARG:CD	2.41	0.51
1:E:166:ASN:ND2	5:E:716:HOH:O	2.44	0.51
2:F:144:CYS:HB2	2:F:145:ASP:HB3	1.92	0.51
1:E:13:ILE:HG22	2:F:138:PHE:HB2	1.92	0.51
2:B:22:TYR:H	2:B:41:THR:HG22	1.77	0.50
2:D:131:GLU:OE2	2:F:128:GLN:HG3	2.11	0.50
1:A:17:HIS:HB3	2:B:115:MET:HE2	1.94	0.49
1:C:161:PHE:HB3	1:C:248:ASN:O	2.12	0.49
1:C:66:PRO:HA	1:C:69:MET:CE	2.43	0.49
1:C:71:ILE:O	1:C:148:TYR:HB3	2.12	0.49
1:E:71:ILE:HG23	1:E:148:TYR:CD1	2.48	0.49
2:D:167:LEU:O	2:D:171:LEU:HG	2.12	0.48
1:E:320:MET:HE1	2:F:21:TRP:HB3	1.94	0.48
1:E:172:ASP:OD1	1:E:173:THR:N	2.44	0.48
1:A:284:ARG:NH1	1:A:284:ARG:CG	2.76	0.48
2:B:21:TRP:H	2:B:41:THR:HG21	1.79	0.48
2:D:95:GLN:NE2	2:F:95:GLN:HE22	2.12	0.47
2:B:46:ASP:OD1	5:B:219:HOH:O	2.20	0.47
1:E:69:MET:HE1	1:E:87:LEU:HD21	1.96	0.47
2:B:94:TYR:HE1	2:D:95:GLN:OE1	1.98	0.47
1:C:271:ASP:N	5:C:663:HOH:O	2.32	0.47
2:B:38:TYR:CZ	2:B:42:GLN:HG3	2.50	0.47
2:D:21:TRP:H	2:D:41:THR:CG2	2.28	0.47
2:D:21:TRP:HB2	2:D:41:THR:HB	1.97	0.47
2:D:139:GLU:OE1	2:F:127:ARG:NH2	2.27	0.47
2:F:89:THR:HA	5:F:203:HOH:O	2.14	0.47
2:D:131:GLU:OE2	2:F:128:GLN:CG	2.62	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:291:ARG:NH1	2:F:56:VAL:O	2.48	0.47
1:C:65:HIS:CE1	1:C:67:ILE:HG12	2.49	0.47
1:E:97:CYS:HB2	1:E:138:ALA:O	2.15	0.47
1:C:69:MET:HG2	1:C:79:HIS:O	2.15	0.46
2:D:6:ILE:HG13	2:D:112:ASP:HA	1.96	0.46
1:E:25:ILE:HG23	1:E:33:GLN:HA	1.97	0.46
2:F:125:GLN:NE2	2:F:155:ASN:HA	2.24	0.46
1:A:15:LEU:HD21	2:B:24:PHE:CE1	2.50	0.46
1:A:50:ARG:CZ	1:A:273:PRO:HG2	2.46	0.46
1:C:152:LYS:HE3	1:C:255:SER:HB3	1.95	0.46
1:C:91:GLU:O	1:C:91:GLU:HG3	2.16	0.46
1:E:108:LEU:HD22	1:E:234:TRP:CD1	2.50	0.46
1:A:186:SER:HA	1:A:218:GLY:O	2.15	0.46
1:A:129(A):SER:HB2	1:A:162:PRO:HG3	1.98	0.46
1:C:71:ILE:HD12	1:C:148:TYR:CE1	2.51	0.46
1:E:25:ILE:CG2	1:E:33:GLN:HA	2.46	0.45
1:E:57:LYS:N	1:E:85:ASP:OD2	2.41	0.45
1:C:114:GLU:O	1:C:263:ARG:HD3	2.17	0.45
2:D:110:MET:O	2:D:113:SER:HB3	2.16	0.45
1:E:15:LEU:HD22	2:F:119:TYR:HA	1.98	0.45
2:B:167:LEU:HD23	2:B:167:LEU:HA	1.80	0.45
1:E:24:THR:HG21	1:E:39:ALA:HB3	1.98	0.45
1:A:158(A):LYS:O	1:A:193:ASP:O	2.34	0.45
1:C:38:ASN:HB3	1:C:318:THR:CG2	2.47	0.45
2:F:159:HIS:CG	2:F:160:SER:N	2.85	0.45
2:B:83:TRP:CD1	2:D:63:PHE:HD1	2.35	0.45
2:D:95:GLN:HE21	2:F:95:GLN:NE2	2.15	0.45
1:E:153:TRP:HZ2	1:E:183:HIS:CE1	2.34	0.45
1:A:160:ASN:OD1	1:A:198:GLN:HG2	2.17	0.45
1:C:312:ARG:HA	5:C:648:HOH:O	2.17	0.45
2:D:11:GLU:H	2:D:11:GLU:HG2	1.53	0.44
1:A:69:MET:HE1	1:A:87:LEU:HD21	2.00	0.44
1:C:15:LEU:HD21	2:D:24:PHE:CE1	2.53	0.44
1:C:18:HIS:HD2	2:D:20:GLY:O	1.99	0.44
2:D:134:GLY:HA2	2:F:124:LYS:HD3	1.99	0.44
1:A:66:PRO:HA	1:A:69:MET:HE3	2.00	0.44
2:B:131:GLU:OE2	2:B:170:ARG:NE	2.50	0.44
2:F:37:ASP:O	2:F:41:THR:HG23	2.18	0.44
2:F:144:CYS:HB2	2:F:148:CYS:HB3	1.18	0.44
2:F:128:GLN:O	2:F:170:ARG:NH1	2.48	0.44
2:B:145:ASP:OD1	2:B:146:ASP:N	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:58:LYS:O	5:B:210:HOH:O	2.21	0.43
2:B:67:GLU:HG3	2:B:81:ILE:HG21	1.99	0.43
2:D:3:PHE:HB3	2:D:116:LEU:HD22	2.00	0.43
1:A:23:GLY:HA3	1:A:35:GLU:OE2	2.19	0.43
1:C:179:MET:HG2	1:C:234:TRP:HB3	2.00	0.43
1:C:307:LYS:HG3	2:D:92:TRP:CE2	2.53	0.43
1:E:163:GLN:NE2	5:E:750:HOH:O	2.51	0.43
1:A:271:ASP:CG	1:A:284:ARG:HH11	2.22	0.43
1:C:42:THR:HG22	2:D:55:LEU:HD21	2.00	0.43
1:E:18:HIS:HB3	2:F:17:MET:HE3	2.01	0.43
1:C:291:ARG:CG	2:D:56:VAL:HB	2.49	0.43
2:D:22:TYR:H	2:D:41:THR:CG2	2.31	0.43
1:E:151:LEU:HD23	1:E:254:PRO:HA	2.00	0.43
1:E:316:LEU:HD11	2:F:103:GLU:HG2	2.00	0.43
1:A:52:CYS:HB2	1:A:279:SER:HB2	2.01	0.43
1:E:169:ARG:NH1	3:I:1:NAG:O5	2.52	0.43
2:F:123:ARG:HA	2:F:138:PHE:CE2	2.53	0.43
1:A:122:SER:HA	1:A:256:ARG:HG2	2.01	0.42
1:C:67:ILE:HD12	1:C:108:LEU:HD23	2.01	0.42
1:E:66:PRO:HA	1:E:69:MET:HE3	2.01	0.42
1:E:307:LYS:HG3	2:F:92:TRP:CE2	2.53	0.42
1:A:45:SER:HA	1:A:296:ASN:OD1	2.19	0.42
1:A:53:MET:HE2	1:A:58:HIS:HB2	2.01	0.42
2:B:83:TRP:CH2	2:D:85:LYS:HG2	2.55	0.42
1:C:291:ARG:HG3	2:D:56:VAL:HB	2.01	0.42
1:C:39:ALA:N	1:C:318:THR:HG22	2.35	0.42
1:E:199:SER:O	1:E:248:ASN:ND2	2.52	0.42
2:B:129:ASN:N	2:B:129:ASN:OD1	2.52	0.41
1:C:17:HIS:NE2	2:D:6:ILE:HG23	2.35	0.41
1:E:201:SER:H	1:E:248:ASN:HB2	1.85	0.41
2:B:125:GLN:OE1	2:B:155:ASN:HA	2.20	0.41
1:A:98:TYR:CD1	1:A:230:ILE:HG13	2.55	0.41
1:C:97:CYS:HB2	1:C:138:ALA:O	2.20	0.41
1:C:176:HIS:HB3	1:C:257:VAL:CG1	2.50	0.41
2:B:128:GLN:O	2:B:170:ARG:NH1	2.40	0.41
1:E:54:LYS:HD2	1:E:276(A):ASN:O	2.20	0.41
2:F:157:TYR:CE1	2:F:159:HIS:HA	2.56	0.41
1:A:219:ALA:O	5:A:560:HOH:O	2.21	0.41
1:E:71:ILE:HD12	1:E:71:ILE:HA	1.74	0.41
1:A:284:ARG:HG3	2:B:69:GLU:HG2	2.03	0.41
2:B:123:ARG:HH21	2:B:124:LYS:HG2	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:84:THR:O	2:D:88:ILE:HG12	2.21	0.41
2:F:47:GLN:HB2	2:F:110:MET:HE3	2.03	0.41
1:E:14:CYS:HA	2:F:137:CYS:HA	2.03	0.41
1:E:191:LYS:HE2	1:E:200:LEU:HG	2.02	0.40
2:F:145:ASP:HB3	2:F:148:CYS:CB	2.50	0.40
1:A:141:ARG:HH11	1:A:141:ARG:HD2	1.70	0.40
1:E:129:SER:O	1:E:157:LYS:HD2	2.21	0.40
2:F:28:ASN:OD1	2:F:31:GLY:CA	2.69	0.40
2:B:4:GLY:O	2:B:8:GLY:HA3	2.21	0.40
2:B:20:GLY:HA3	2:B:36:ALA:HB1	2.04	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	316/323 (98%)	305 (96%)	11 (4%)	0	100	100
1	C	317/323 (98%)	307 (97%)	10 (3%)	0	100	100
1	E	316/323 (98%)	306 (97%)	10 (3%)	0	100	100
2	B	169/180 (94%)	160 (95%)	9 (5%)	0	100	100
2	D	165/180 (92%)	161 (98%)	4 (2%)	0	100	100
2	F	166/180 (92%)	154 (93%)	11 (7%)	1 (1%)	25	47
All	All	1449/1509 (96%)	1393 (96%)	55 (4%)	1 (0%)	51	75

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	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	269/272 (99%)	265 (98%)	4 (2%)	65	83
1	C	270/272 (99%)	269 (100%)	1 (0%)	91	97
1	E	269/272 (99%)	267 (99%)	2 (1%)	84	94
2	B	146/154 (95%)	145 (99%)	1 (1%)	84	94
2	D	144/154 (94%)	142 (99%)	2 (1%)	67	85
2	F	146/154 (95%)	140 (96%)	6 (4%)	30	56
All	All	1244/1278 (97%)	1228 (99%)	16 (1%)	69	86

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	THR
1	A	126	THR
1	A	227	SER
1	A	302	VAL
2	B	135	LYS
1	C	126	THR
2	D	102	MET
2	D	148	CYS
1	E	49	ASN
1	E	274	ILE
2	F	27	GLN
2	F	30	GLN
2	F	72	GLU
2	F	135	LYS
2	F	146	ASP
2	F	164	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	HIS

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Mol	Chain	Res	Type
1	A	295	GLN
1	C	159	GLN
1	C	189	GLN
2	D	95	GLN
2	D	155	ASN
1	E	17	HIS
1	E	163	GLN
1	E	212	ASN
1	E	233	HIS
2	F	95	GLN
2	F	125	GLN
2	F	155	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](#)

6 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	3,1	14,14,15	0.76	1 (7%)	17,19,21	0.89	0
3	NAG	G	2	3	14,14,15	0.44	0	17,19,21	0.92	1 (5%)
3	NAG	H	1	3,1	14,14,15	0.78	1 (7%)	17,19,21	0.79	0
3	NAG	H	2	3	14,14,15	0.23	0	17,19,21	0.53	0
3	NAG	I	1	3,1	14,14,15	0.74	1 (7%)	17,19,21	0.95	1 (5%)
3	NAG	I	2	3	14,14,15	0.34	0	17,19,21	0.78	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	G	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	G	2	3	-	4/6/23/26	0/1/1/1
3	NAG	H	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	H	2	3	-	2/6/23/26	0/1/1/1
3	NAG	I	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	I	2	3	-	0/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	I	1	NAG	O5-C1	-2.46	1.39	1.43
3	G	1	NAG	O5-C1	-2.43	1.39	1.43
3	H	1	NAG	O5-C1	-2.42	1.39	1.43

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	1	NAG	C1-C2-N2	-2.57	106.10	110.49
3	G	2	NAG	C2-N2-C7	-2.43	119.45	122.90

There are no chirality outliers.

All (10) torsion outliers are listed below:

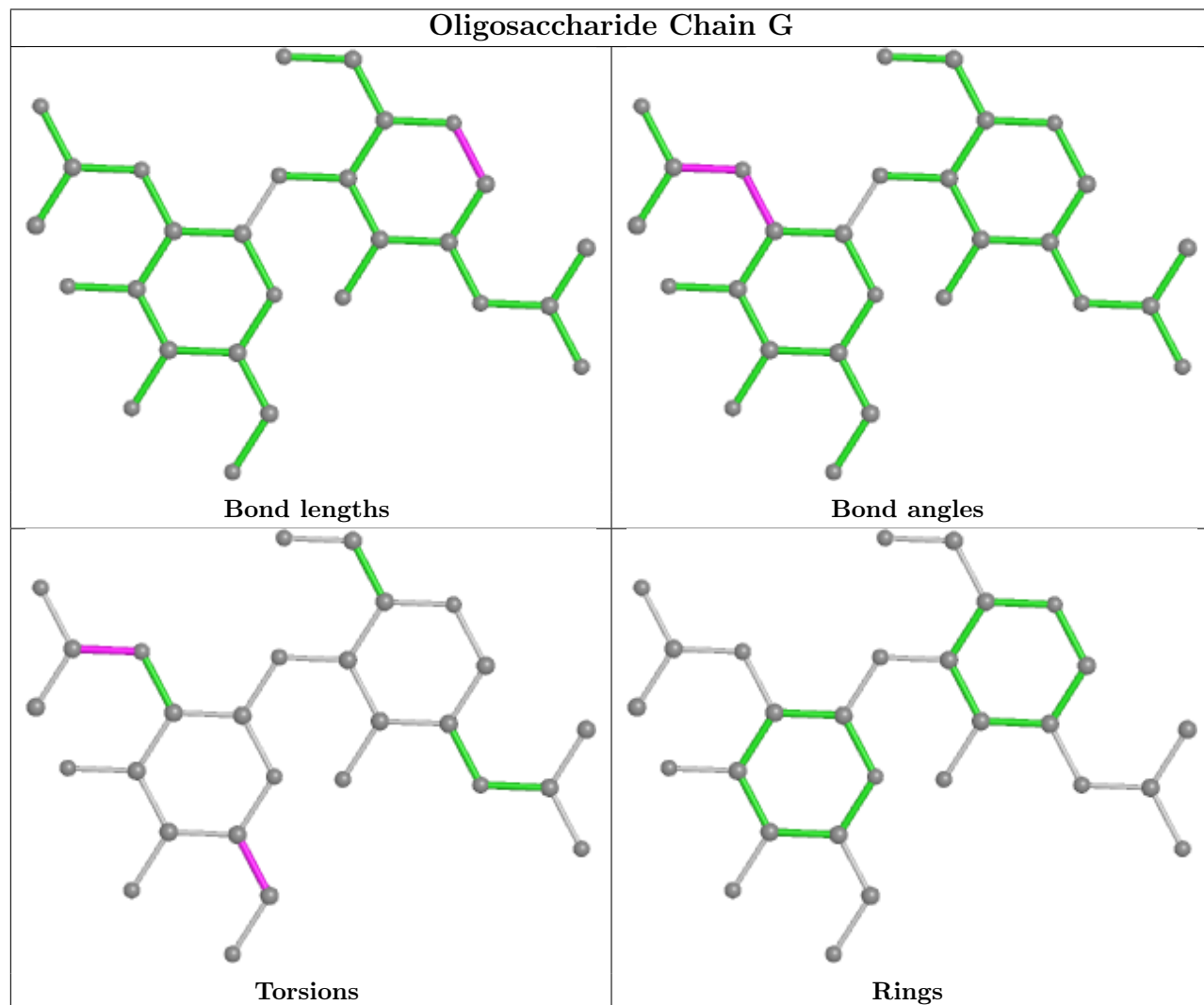
Mol	Chain	Res	Type	Atoms
3	G	2	NAG	O5-C5-C6-O6
3	I	1	NAG	O5-C5-C6-O6
3	G	2	NAG	C4-C5-C6-O6
3	H	2	NAG	O5-C5-C6-O6
3	I	1	NAG	C4-C5-C6-O6
3	H	2	NAG	C4-C5-C6-O6
3	H	1	NAG	C4-C5-C6-O6
3	H	1	NAG	O5-C5-C6-O6
3	G	2	NAG	C8-C7-N2-C2
3	G	2	NAG	O7-C7-N2-C2

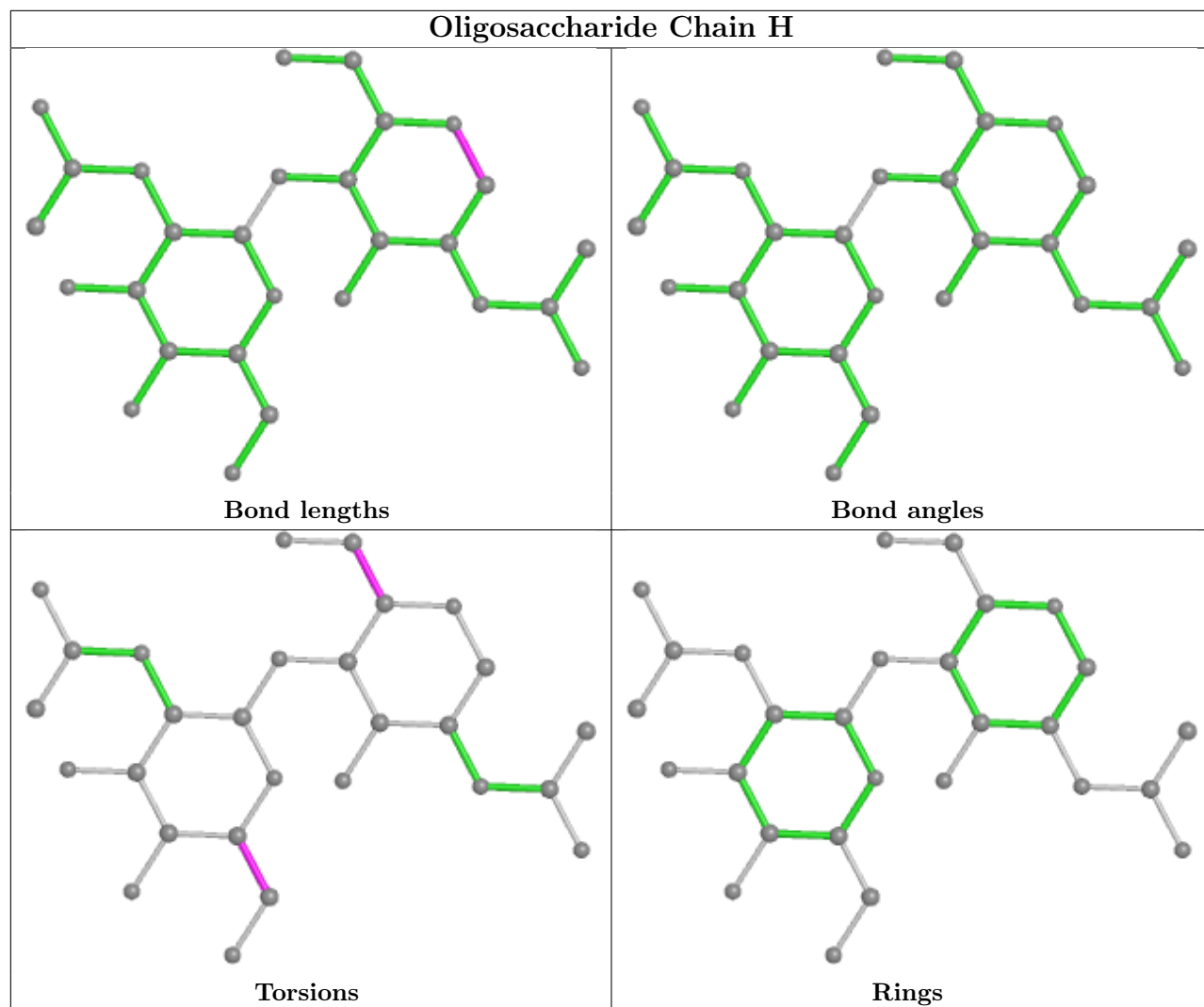
There are no ring outliers.

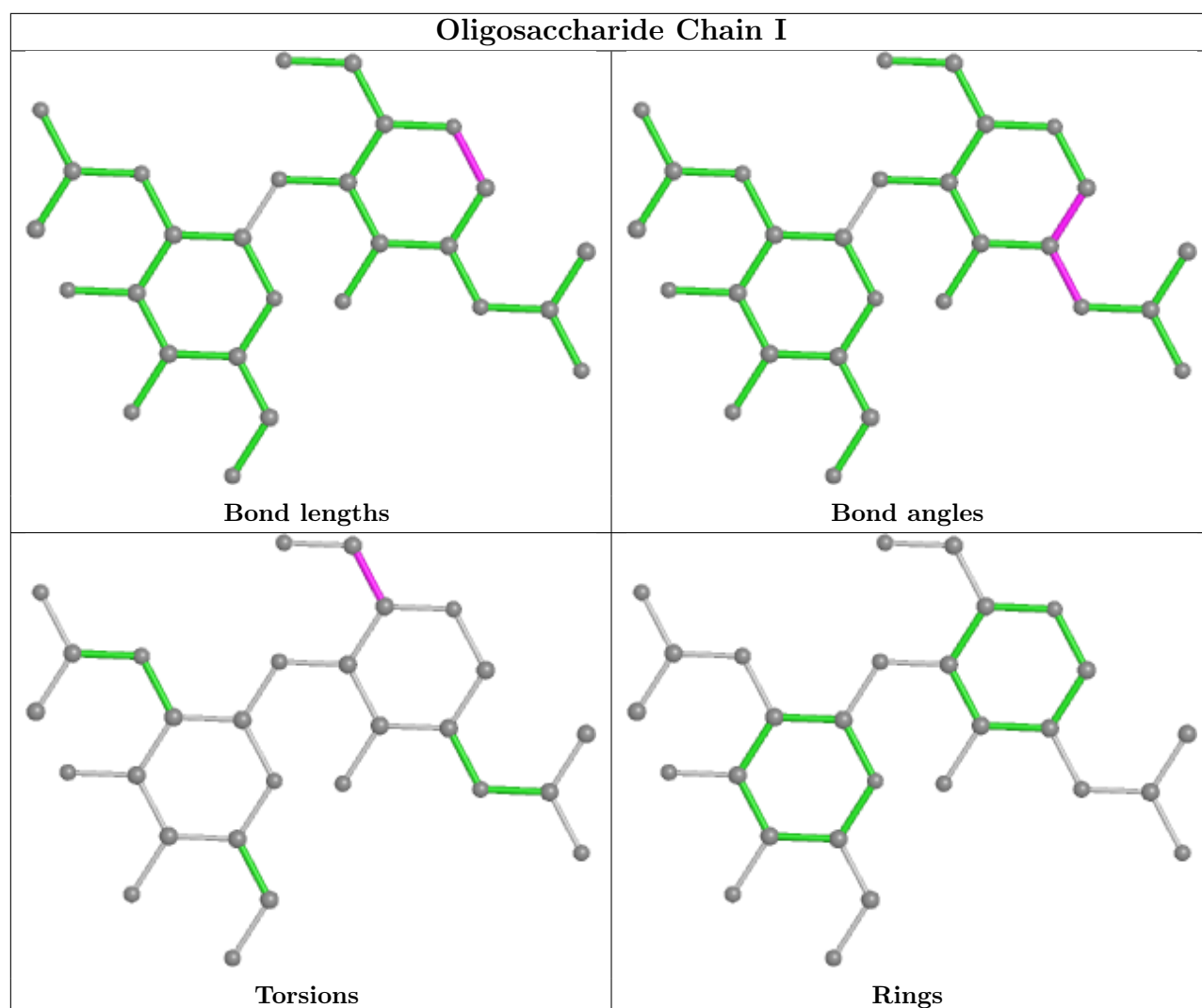
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	I	1	NAG	1	0

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







## 5.6 Ligand geometry ⓘ

2 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	500	1	14,14,15	0.68	1 (7%)	17,19,21	0.80	0
4	NAG	A	400	1	14,14,15	0.65	1 (7%)	17,19,21	0.56	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
4	NAG	C	500	1	-	3/6/23/26	0/1/1/1
4	NAG	A	400	1	-	0/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	500	NAG	O5-C1	-2.36	1.40	1.43
4	A	400	NAG	O5-C1	-2.33	1.40	1.43

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	500	NAG	C8-C7-N2-C2
4	C	500	NAG	O7-C7-N2-C2
4	C	500	NAG	C1-C2-N2-C7

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.



## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	318/323 (98%)	-0.47	0 100 100	11, 21, 42, 63	0
1	C	319/323 (98%)	-0.38	5 (1%) 72 68	14, 24, 45, 69	0
1	E	318/323 (98%)	-0.34	0 100 100	15, 27, 51, 67	0
2	B	171/180 (95%)	0.07	5 (2%) 51 45	18, 38, 56, 69	0
2	D	169/180 (93%)	-0.06	2 (1%) 79 76	17, 40, 55, 69	0
2	F	170/180 (94%)	0.26	5 (2%) 51 45	16, 45, 72, 78	0
All	All	1465/1509 (97%)	-0.23	17 (1%) 79 76	11, 29, 56, 78	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	59	THR	3.5
1	C	21	ALA	3.4
2	F	29	ALA	3.4
1	C	326	LEU	3.0
2	F	23	GLY	3.0
1	C	325	GLU	2.6
2	D	19	ASP	2.6
2	B	57	GLU	2.6
1	C	22	ASN	2.5
2	B	19	ASP	2.4
2	D	18	VAL	2.4
2	B	171	LEU	2.3
2	F	2	LEU	2.2
2	F	38	TYR	2.1
2	B	18	VAL	2.1
2	B	38	TYR	2.1
1	C	291	ARG	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

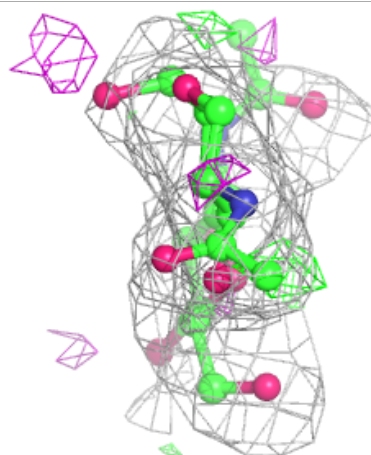
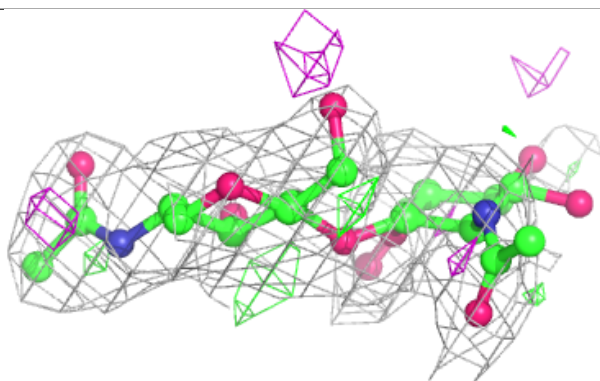
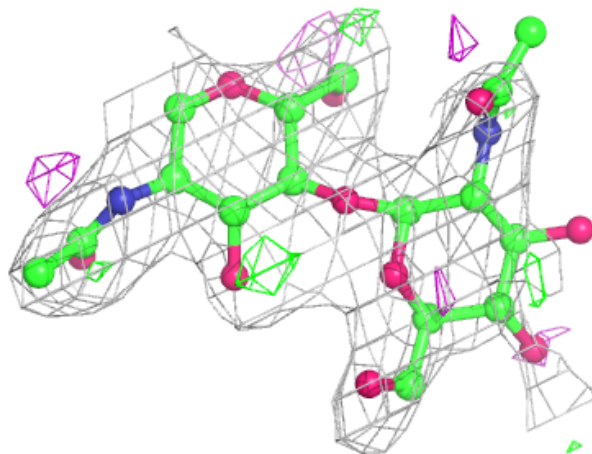
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	NAG	I	2	14/15	0.68	0.36	72,85,93,99	0
3	NAG	G	2	14/15	0.71	0.39	68,76,84,95	0
3	NAG	H	2	14/15	0.79	0.38	67,80,92,95	0
3	NAG	I	1	14/15	0.88	0.19	29,54,67,79	0
3	NAG	H	1	14/15	0.89	0.18	30,50,59,69	0
3	NAG	G	1	14/15	0.89	0.21	27,51,58,67	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.

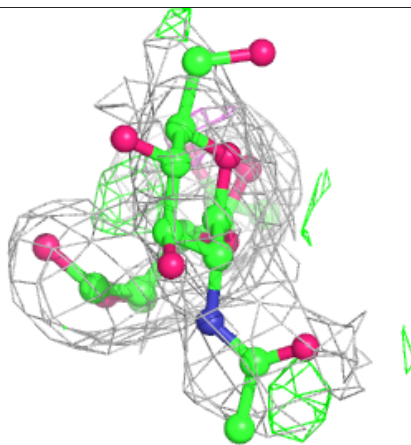
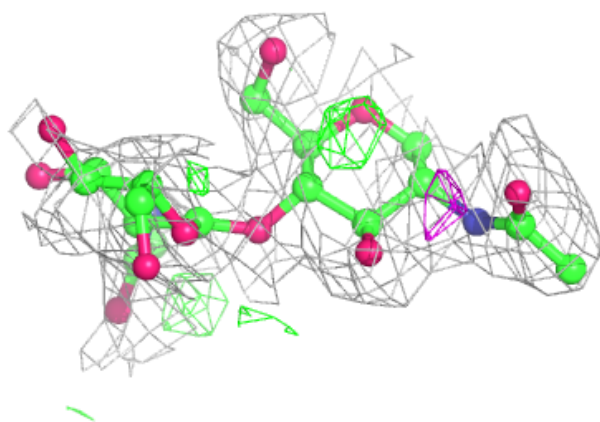
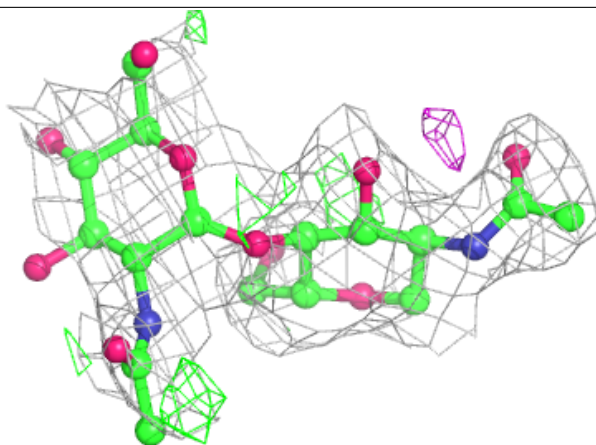
**Electron density around Chain G:**

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

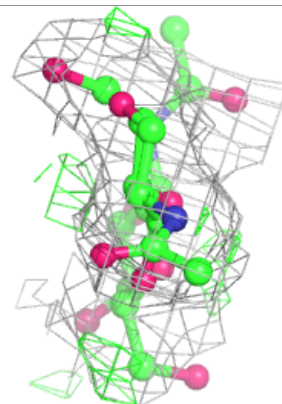
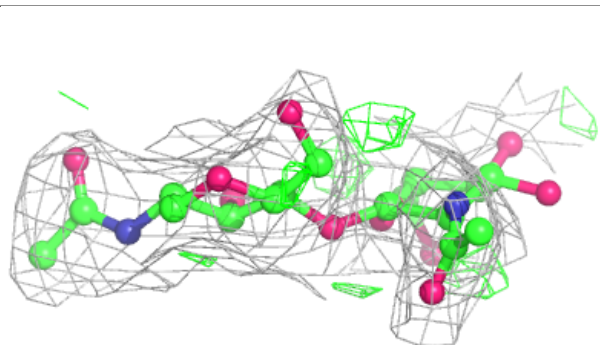
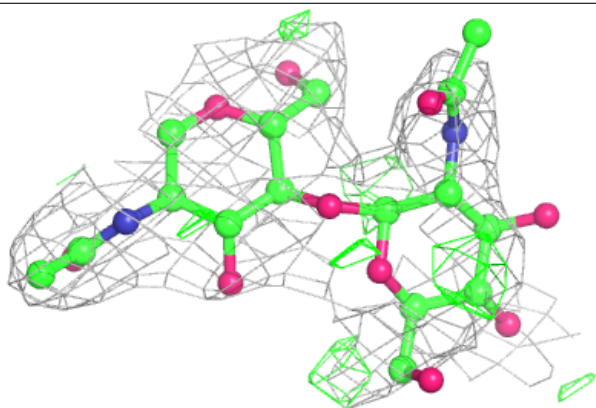


**Electron density around Chain H:**

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

**Electron density around Chain 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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
4	NAG	A	400	14/15	0.79	0.33	52,64,72,76	0
4	NAG	C	500	14/15	0.79	0.34	55,78,82,85	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.