



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 22, 2020 – 10:10 PM BST

PDB ID : 6Y9L  
Title : Crystal structure of TSWV glycoprotein N ectodomain (sGn)  
Authors : Dessau, M.; Bahat, Y.  
Deposited on : 2020-03-09  
Resolution : 4.10 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 : 2.14.6  
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) : 2.14.6

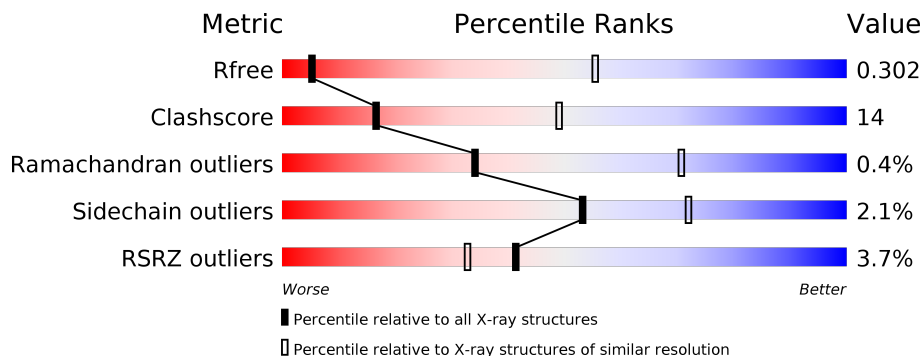
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## X-RAY DIFFRACTION

The reported resolution of this entry is 4.10 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	1193 (4.50-3.70)
Clashscore	141614	1003 (4.44-3.76)
Ramachandran outliers	138981	1005 (4.48-3.72)
Sidechain outliers	138945	1199 (4.50-3.70)
RSRZ outliers	127900	1034 (4.50-3.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on 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	288	 2% 45% 23% 32%
1	B	288	 4% 52% 16% 32%
1	C	288	 2% 47% 20% 33%
1	D	288	 3% 46% 21% 32%
2	E	3	 33% 33% 33%
2	G	3	 33% 67%

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Mol	Chain	Length	Quality of chain
2	J	3	 100%
3	F	2	 50% 50%
3	H	2	 50% 50%
3	I	2	 50% 50%
3	K	2	 50% 50%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	BMA	G	3	-	-	-	X
2	NAG	J	2	-	-	-	X
3	NAG	F	1	-	-	-	X
3	NAG	F	2	-	-	-	X
3	NAG	H	1	-	-	-	X
3	NAG	I	2	-	-	-	X
4	NAG	C	401	-	-	-	X

## 2 Entry composition [i](#)

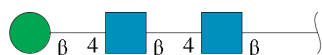
There are 4 unique types of molecules in this entry. The entry contains 6325 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Glycoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	B	196	1522	959	253	301	9	0	0	0
1	D	196	1523	961	254	299	9	0	0	0
1	A	197	1531	965	255	302	9	0	0	0
1	C	194	1506	949	251	297	9	0	0	0

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



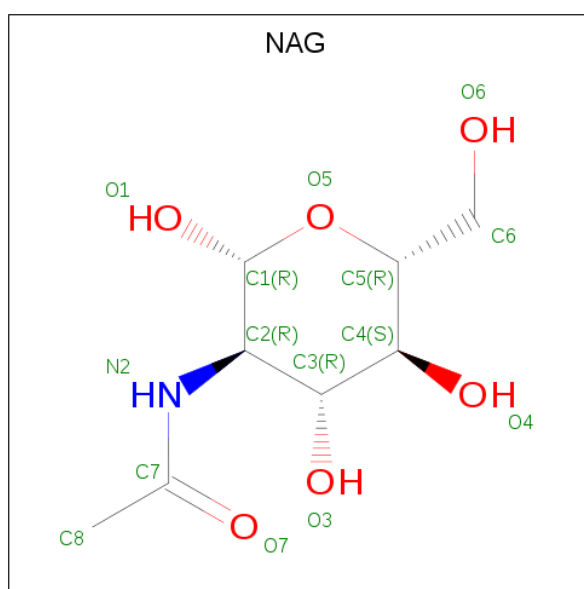
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	E	3	39	22	2	15	0	0	0
2	G	3	39	22	2	15	0	0	0
2	J	3	39	22	2	15	0	0	0

- 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	F	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			
3	K	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:  $C_8H_{15}NO_6$ ).

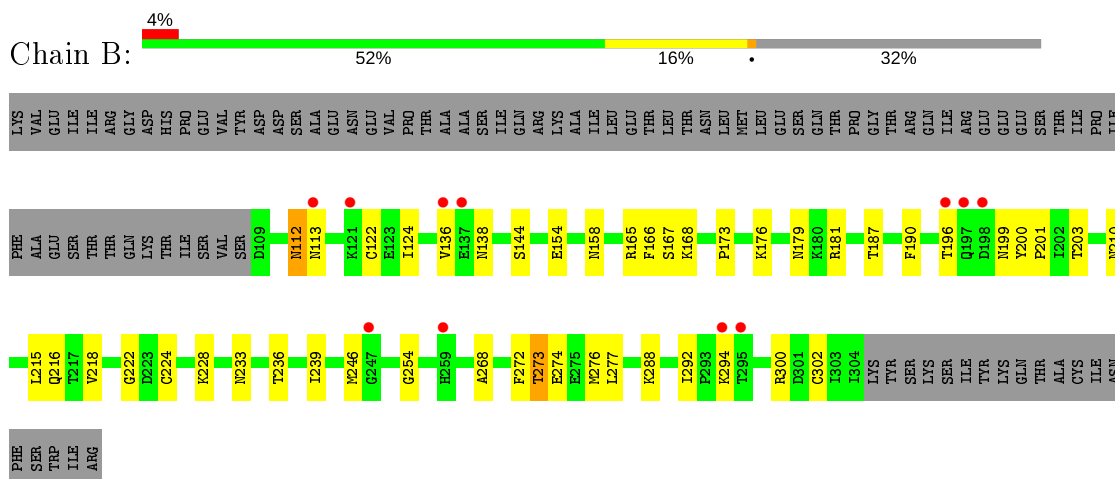


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	C	1	Total	C	N	O	0	0
			14	8	1	5		

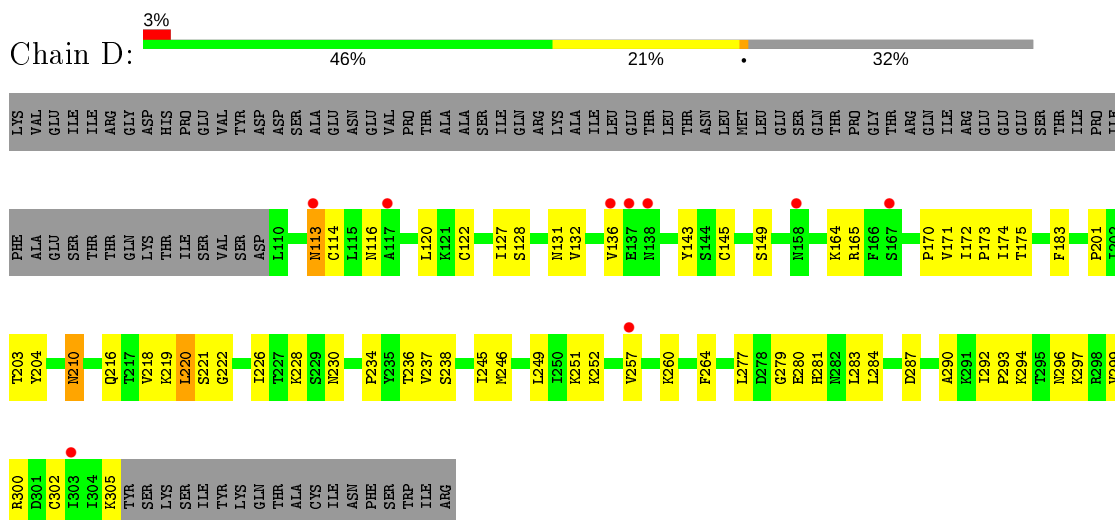
### 3 Residue-property plots [i](#)

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.

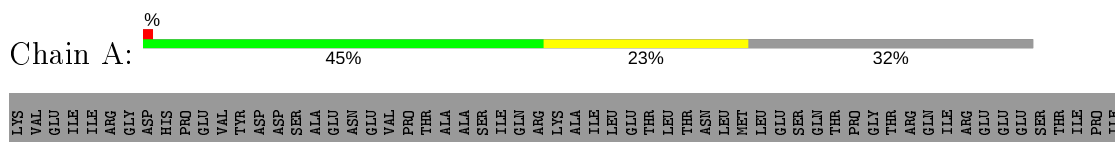
- Molecule 1: Glycoprotein

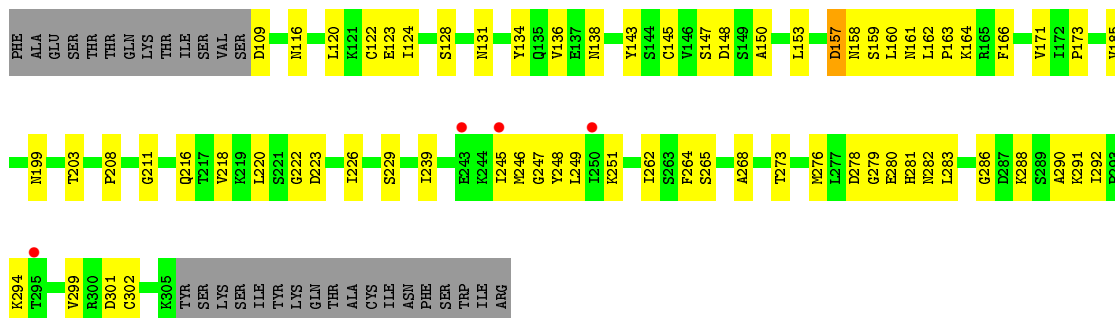


- Molecule 1: Glycoprotein

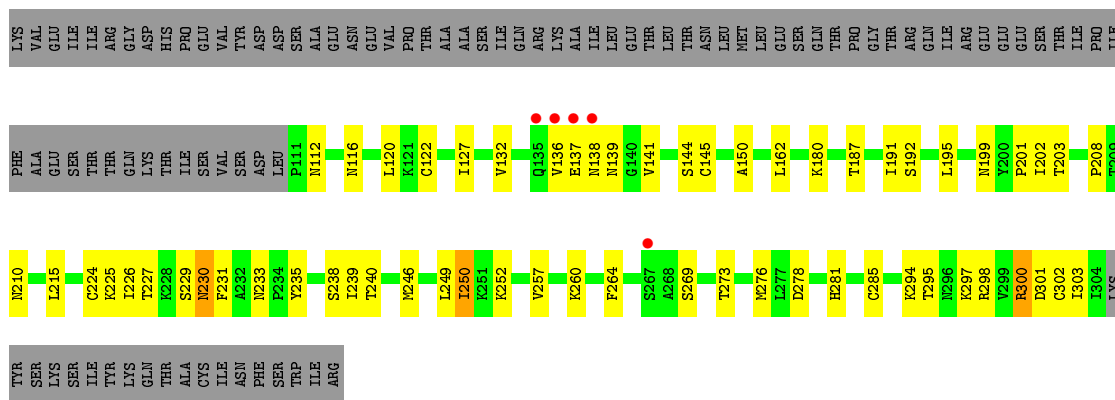


- Molecule 1: Glycoprotein





- Molecule 1: Glycoprotein



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



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



- Molecule 2: beta-D-mannopyranose-(1-4)-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

Chain F:  50% 50%

  
MAG1  
MAG2

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

Chain H:  50% 50%


  
MAG1  
MAG2

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

Chain I:  50% 50%

  
MAG1  
MAG2

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

Chain K:  50% 50%

  
MAG1  
MAG2



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	165.99Å 215.88Å 146.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.02 – 4.10 49.02 – 4.10	Depositor EDS
% Data completeness (in resolution range)	98.3 (49.02-4.10) 98.6 (49.02-4.10)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.97 (at 4.14Å)	Xtrriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, $R_{free}$	0.277 , 0.302 0.277 , 0.302	Depositor DCC
$R_{free}$ test set	1033 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	288.8	Xtrriage
Anisotropy	0.268	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 387.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.39$ , $\langle L^2 \rangle = 0.23$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	6325	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	331.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.43% 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: BMA, 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	0.27	0/1559	0.49	0/2108
1	B	0.28	0/1550	0.49	0/2097
1	C	0.27	0/1534	0.48	0/2074
1	D	0.25	0/1551	0.46	0/2097
All	All	0.27	0/6194	0.48	0/8376

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	A	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	157	ASP	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	1531	0	1532	49	0
1	B	1522	0	1519	31	0
1	C	1506	0	1505	61	0
1	D	1523	0	1528	45	0
2	E	39	0	34	1	0
2	G	39	0	34	0	0
2	J	39	0	34	1	0
3	F	28	0	25	1	0
3	H	28	0	25	1	0
3	I	28	0	25	0	0
3	K	28	0	25	1	0
4	C	14	0	13	0	0
All	All	6325	0	6299	179	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 14.

All (179) 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:B:124:ILE:HB	1:B:158:ASN:HB3	1.63	0.81
1:B:273:THR:HG22	1:B:276:MET:HB2	1.64	0.81
1:A:223:ASP:H	1:A:246:MET:HE1	1.46	0.80
1:C:144:SER:HB2	1:C:202:ILE:HG22	1.66	0.75
1:C:273:THR:HB	1:C:276:MET:HG2	1.69	0.74
1:B:112:ASN:HB2	3:F:1:NAG:H81	1.68	0.73
1:B:173:PRO:HA	1:B:218:VAL:HA	1.72	0.71
1:D:145:CYS:SG	1:D:203:THR:OG1	2.47	0.70
1:B:122:CYS:HB3	1:B:136:VAL:HG12	1.72	0.70
1:A:150:ALA:HA	1:A:153:LEU:HD21	1.74	0.69
1:D:234:PRO:HD3	1:C:295:THR:OG1	1.96	0.66
1:A:245:ILE:HA	1:A:265:SER:HA	1.76	0.66
1:D:128:SER:OG	1:D:131:ASN:OD1	2.13	0.66
1:C:224:CYS:HB3	1:C:239:ILE:HD11	1.78	0.66
1:A:185:VAL:HG11	1:A:220:LEU:HG	1.79	0.65
1:D:173:PRO:HA	1:D:218:VAL:HA	1.77	0.65
1:D:113:ASN:HA	1:D:116:ASN:HB2	1.79	0.65
1:A:279:GLY:N	1:A:294:LYS:HB2	2.12	0.65
1:C:191:ILE:HG13	1:C:202:ILE:HD11	1.77	0.65
1:A:145:CYS:SG	1:A:203:THR:OG1	2.56	0.64
1:C:239:ILE:HG23	1:C:264:PHE:CZ	2.33	0.64
1:C:238:SER:HA	1:C:269:SER:HA	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:201:PRO:HA	1:C:210:ASN:HA	1.80	0.63
1:A:124:ILE:HD13	1:A:134:TYR:HB3	1.80	0.63
1:A:124:ILE:HG13	1:A:159:SER:HB3	1.81	0.63
1:A:220:LEU:HB3	1:A:286:GLY:HA2	1.81	0.63
1:D:127:ILE:HG12	1:D:132:VAL:HG23	1.81	0.63
1:D:296:ASN:HB3	1:D:299:VAL:HG23	1.80	0.63
1:A:278:ASP:HA	1:A:294:LYS:HD2	1.82	0.62
1:A:120:LEU:HD13	1:A:138:ASN:HB2	1.81	0.61
1:A:122:CYS:O	1:A:157:ASP:N	2.32	0.61
1:C:145:CYS:SG	1:C:203:THR:OG1	2.58	0.61
1:D:246:MET:HB3	1:D:264:PHE:HD2	1.66	0.61
1:B:144:SER:HB2	1:B:200:TYR:HE2	1.66	0.61
2:J:1:NAG:H61	2:J:2:NAG:HN2	1.66	0.61
1:B:187:THR:HB	1:B:215:LEU:HD11	1.83	0.59
1:D:252:LYS:HB3	1:D:281:HIS:HA	1.85	0.58
1:C:122:CYS:HB3	1:C:136:VAL:HG12	1.85	0.57
1:D:283:LEU:HB3	1:D:290:ALA:HB3	1.86	0.57
1:D:113:ASN:HB2	1:D:120:LEU:HD11	1.85	0.57
1:B:239:ILE:HG23	1:B:268:ALA:HB3	1.87	0.57
1:A:281:HIS:HB3	1:A:292:ILE:HD11	1.87	0.56
1:C:257:VAL:HG23	1:C:260:LYS:HD2	1.88	0.56
1:D:172:ILE:HG12	1:D:221:SER:HB2	1.88	0.56
1:D:174:ILE:HD13	1:D:219:LYS:HB2	1.88	0.56
1:A:109:ASP:OD1	1:A:109:ASP:N	2.40	0.55
1:B:272:PHE:HE1	1:B:292:ILE:HG21	1.71	0.55
1:C:225:LYS:HB2	1:C:240:THR:CG2	2.37	0.55
1:C:252:LYS:HD3	1:C:281:HIS:CE1	2.42	0.55
1:D:136:VAL:HG22	1:D:143:TYR:HB2	1.89	0.55
1:C:139:ASN:H	1:C:141:VAL:HG22	1.72	0.55
1:A:173:PRO:HA	1:A:218:VAL:HA	1.89	0.55
1:D:293:PRO:HD3	1:C:231:PHE:CE1	2.42	0.54
1:B:176:LYS:HB2	1:B:179:ASN:HB2	1.88	0.54
3:H:1:NAG:H3	3:H:1:NAG:H83	1.90	0.54
1:A:222:GLY:HA3	1:A:246:MET:HE1	1.89	0.54
1:D:219:LYS:NZ	1:D:220:LEU:O	2.35	0.54
1:C:235:TYR:HE2	1:C:300:ARG:HD2	1.71	0.54
1:C:246:MET:HB3	1:C:264:PHE:HB3	1.88	0.54
1:A:249:LEU:HD21	1:A:251:LYS:HB2	1.90	0.54
1:A:160:LEU:HG	1:A:161:ASN:N	2.23	0.53
1:B:113:ASN:HD21	1:B:138:ASN:HB2	1.74	0.53
1:C:278:ASP:OD2	1:C:281:HIS:ND1	2.41	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:274:GLU:HG3	1:B:300:ARG:HE	1.74	0.53
1:D:183:PHE:HA	1:D:284:LEU:HD21	1.91	0.53
1:C:138:ASN:H	1:C:141:VAL:HG23	1.74	0.53
1:C:150:ALA:HB1	1:C:162:LEU:HD11	1.91	0.52
1:A:301:ASP:OD1	1:A:302:CYS:N	2.42	0.52
1:C:225:LYS:HB2	1:C:240:THR:HG22	1.91	0.52
3:K:1:NAG:N2	3:K:1:NAG:O4	2.41	0.52
1:C:137:GLU:HG3	1:C:141:VAL:N	2.25	0.52
1:C:120:LEU:CD2	1:C:138:ASN:HA	2.39	0.52
1:A:122:CYS:HB3	1:A:136:VAL:HG12	1.92	0.52
1:C:252:LYS:HB3	1:C:281:HIS:CD2	2.44	0.52
1:A:147:SER:OG	1:A:148:ASP:N	2.42	0.52
1:B:165:ARG:NH2	1:B:167:SER:HA	2.25	0.52
1:C:203:THR:HA	1:C:208:PRO:HA	1.92	0.51
1:D:201:PRO:HA	1:D:210:ASN:HA	1.92	0.51
1:D:257:VAL:HG12	1:D:260:LYS:HE2	1.93	0.51
1:B:233:ASN:HB3	1:B:236:THR:HB	1.92	0.51
1:D:283:LEU:N	1:D:290:ALA:O	2.41	0.51
1:C:199:ASN:N	1:C:199:ASN:OD1	2.43	0.51
1:A:123:GLU:HA	1:A:157:ASP:O	2.10	0.50
1:A:226:ILE:H	1:A:288:LYS:NZ	2.09	0.50
1:C:230:ASN:OD1	1:C:230:ASN:N	2.45	0.50
1:D:164:LYS:HG3	1:D:165:ARG:H	1.76	0.50
1:C:239:ILE:HG12	1:C:285:CYS:SG	2.51	0.50
1:D:302:CYS:SG	1:C:302:CYS:HB2	2.52	0.50
1:B:222:GLY:HA3	1:B:246:MET:HE1	1.94	0.50
1:A:128:SER:OG	1:A:131:ASN:O	2.30	0.50
1:A:216:GLN:HG2	1:C:192:SER:HA	1.93	0.50
1:B:203:THR:O	1:B:203:THR:OG1	2.29	0.50
1:D:305:LYS:HG3	1:C:298:ARG:NH2	2.26	0.50
1:C:187:THR:HB	1:C:215:LEU:HD11	1.93	0.49
1:C:273:THR:HB	1:C:276:MET:CG	2.41	0.49
1:A:148:ASP:N	1:A:148:ASP:OD1	2.45	0.49
1:A:203:THR:HA	1:A:208:PRO:HA	1.93	0.49
1:D:230:ASN:HB3	1:D:237:VAL:HG12	1.93	0.49
1:A:273:THR:H	1:A:276:MET:HB2	1.76	0.49
1:A:223:ASP:N	1:A:246:MET:HE1	2.23	0.49
1:C:226:ILE:HA	1:C:239:ILE:HD12	1.94	0.49
1:C:249:LEU:HD12	1:C:250:ILE:H	1.78	0.49
1:C:246:MET:HB3	1:C:264:PHE:CD2	2.48	0.49
1:D:280:GLU:N	1:D:280:GLU:OE1	2.46	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:137:GLU:HA	1:C:141:VAL:O	2.13	0.48
1:A:171:VAL:HG12	1:A:220:LEU:HA	1.94	0.48
1:C:238:SER:OG	1:C:239:ILE:N	2.46	0.48
1:A:223:ASP:H	1:A:246:MET:CE	2.20	0.48
1:C:297:LYS:HE2	1:C:301:ASP:OD2	2.13	0.48
1:A:248:TYR:HE2	1:A:264:PHE:HB3	1.79	0.48
1:C:233:ASN:C	1:C:235:TYR:H	2.15	0.48
1:D:149:SER:O	1:D:165:ARG:NH2	2.47	0.48
1:A:239:ILE:HB	1:A:268:ALA:HB3	1.96	0.48
1:B:190:PHE:HE1	1:D:175:THR:HG23	1.79	0.48
1:A:128:SER:OG	1:A:131:ASN:OD1	2.20	0.47
1:D:293:PRO:HD3	1:C:231:PHE:HE1	1.78	0.47
1:C:252:LYS:HB3	1:C:281:HIS:CG	2.49	0.47
1:D:222:GLY:O	1:D:287:ASP:HB2	2.15	0.47
1:D:230:ASN:ND2	1:D:237:VAL:HA	2.30	0.47
1:B:113:ASN:ND2	1:B:138:ASN:HB2	2.31	0.46
1:B:224:CYS:SG	1:B:246:MET:HE3	2.56	0.46
1:B:224:CYS:O	1:B:288:LYS:HD2	2.15	0.46
1:B:274:GLU:HG3	1:B:300:ARG:HH11	1.81	0.46
1:C:276:MET:C	1:C:278:ASP:H	2.18	0.46
1:A:136:VAL:HG22	1:A:143:TYR:HB2	1.98	0.46
1:A:278:ASP:HA	1:A:294:LYS:HB2	1.98	0.45
1:A:199:ASN:HD22	1:A:211:GLY:HA2	1.82	0.45
1:B:181:ARG:HH21	1:B:254:GLY:HA3	1.81	0.45
1:B:166:PHE:O	1:B:168:LYS:HG3	2.17	0.45
1:C:227:THR:HG23	1:C:239:ILE:HA	1.99	0.45
1:A:163:PRO:HG2	1:A:166:PHE:CE1	2.50	0.45
1:C:233:ASN:C	1:C:235:TYR:N	2.70	0.45
1:B:201:PRO:HA	1:B:210:ASN:HA	1.99	0.45
1:C:235:TYR:CE2	1:C:300:ARG:HD2	2.51	0.45
1:B:176:LYS:HD3	1:B:179:ASN:ND2	2.33	0.44
1:C:294:LYS:O	1:C:294:LYS:HG3	2.18	0.44
1:C:127:ILE:HG12	1:C:132:VAL:HG23	1.98	0.43
1:D:226:ILE:HG22	1:D:228:LYS:HG2	2.00	0.43
1:C:138:ASN:H	1:C:141:VAL:CG2	2.30	0.43
1:A:278:ASP:OD1	1:A:281:HIS:HD2	2.01	0.43
1:A:282:ASN:OD1	1:A:291:LYS:HD2	2.18	0.43
1:D:131:ASN:ND2	1:D:204:TYR:OH	2.47	0.43
1:D:171:VAL:HG13	1:D:218:VAL:HB	2.00	0.43
1:C:273:THR:HB	1:C:276:MET:HB2	2.00	0.43
1:D:277:LEU:HA	1:D:292:ILE:HG21	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:278:ASP:OD1	1:A:281:HIS:CD2	2.71	0.43
1:B:228:LYS:C	1:A:229:SER:HB3	2.38	0.43
1:B:277:LEU:O	1:B:294:LYS:HA	2.18	0.43
1:B:302:CYS:SG	1:A:299:VAL:HG13	2.59	0.43
1:D:183:PHE:HB3	1:D:219:LYS:HG3	2.00	0.43
1:D:170:PRO:HD3	1:D:245:ILE:HG23	2.00	0.43
1:D:251:LYS:HB3	1:D:251:LYS:HE3	1.85	0.42
1:B:199:ASN:HB2	2:E:1:NAG:O7	2.20	0.42
1:D:297:LYS:HG2	1:D:300:ARG:HD2	2.01	0.42
1:B:216:GLN:OE1	1:D:216:GLN:NE2	2.52	0.42
1:A:280:GLU:OE1	1:A:280:GLU:N	2.37	0.42
1:B:154:GLU:N	1:B:154:GLU:OE2	2.48	0.42
1:C:139:ASN:N	1:C:141:VAL:HG22	2.33	0.42
1:D:230:ASN:HD22	1:D:238:SER:H	1.67	0.42
1:C:227:THR:C	1:C:229:SER:H	2.22	0.42
1:C:276:MET:HA	1:C:278:ASP:OD1	2.20	0.42
1:C:112:ASN:O	1:C:112:ASN:OD1	2.38	0.42
1:A:248:TYR:HD2	1:A:264:PHE:CD2	2.38	0.42
1:C:246:MET:HB3	1:C:264:PHE:HD2	1.85	0.42
1:A:247:GLY:HA2	1:A:262:ILE:O	2.20	0.41
1:A:251:LYS:HE3	1:A:251:LYS:HB3	1.91	0.41
1:C:249:LEU:HD21	1:C:257:VAL:HG21	2.01	0.41
1:A:239:ILE:O	1:A:268:ALA:N	2.53	0.41
1:D:249:LEU:HD13	1:D:260:LYS:O	2.21	0.41
1:C:137:GLU:HG3	1:C:141:VAL:H	1.84	0.41
1:C:252:LYS:HB3	1:C:281:HIS:CE1	2.56	0.41
1:C:195:LEU:HD22	1:C:195:LEU:H	1.86	0.40
1:C:180:LYS:HE2	1:C:180:LYS:HB3	1.79	0.40
1:D:279:GLY:N	1:D:294:LYS:HB2	2.36	0.40
1:A:283:LEU:HB3	1:A:290:ALA:HB3	2.02	0.40
1:C:273:THR:HB	1:C:276:MET:CB	2.52	0.40
1:D:252:LYS:HD2	1:D:281:HIS:CD2	2.57	0.40
1:D:279:GLY:H	1:D:294:LYS:HD3	1.86	0.40
1:D:251:LYS:HD3	1:D:257:VAL:HG21	2.03	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	195/288 (68%)	181 (93%)	13 (7%)	1 (0%)	29	67
1	B	194/288 (67%)	166 (86%)	28 (14%)	0	100	100
1	C	192/288 (67%)	172 (90%)	19 (10%)	1 (0%)	29	67
1	D	194/288 (67%)	170 (88%)	23 (12%)	1 (0%)	29	67
All	All	775/1152 (67%)	689 (89%)	83 (11%)	3 (0%)	34	71

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	158	ASN
1	C	116	ASN
1	D	113	ASN

### 5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	181/264 (69%)	178 (98%)	3 (2%)	60	78
1	B	180/264 (68%)	177 (98%)	3 (2%)	60	78
1	C	178/264 (67%)	174 (98%)	4 (2%)	52	71
1	D	180/264 (68%)	175 (97%)	5 (3%)	43	65
All	All	719/1056 (68%)	704 (98%)	15 (2%)	53	72

All (15) residues with a non-rotameric sidechain are listed below:



Mol	Chain	Res	Type
1	B	112	ASN
1	B	196	THR
1	B	273	THR
1	D	114	CYS
1	D	122	CYS
1	D	210	ASN
1	D	220	LEU
1	D	236	THR
1	A	116	ASN
1	A	162	LEU
1	A	164	LYS
1	C	230	ASN
1	C	250	ILE
1	C	300	ARG
1	C	303	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	216	GLN
1	C	112	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](#)

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
2	NAG	E	1	1,2	14,14,15	1.00	1 (7%)	17,19,21	1.11	2 (11%)
2	NAG	E	2	2	14,14,15	0.87	1 (7%)	17,19,21	0.52	0
2	BMA	E	3	2	11,11,12	0.55	0	15,15,17	0.72	0
3	NAG	F	1	1,3	14,14,15	0.26	0	17,19,21	0.77	1 (5%)
3	NAG	F	2	3	14,14,15	0.32	0	17,19,21	0.43	0
2	NAG	G	1	1,2	14,14,15	0.51	0	17,19,21	0.59	0
2	NAG	G	2	2	14,14,15	0.41	0	17,19,21	1.34	2 (11%)
2	BMA	G	3	2	11,11,12	0.65	0	15,15,17	1.02	1 (6%)
3	NAG	H	1	1,3	14,14,15	0.56	0	17,19,21	1.41	3 (17%)
3	NAG	H	2	3	14,14,15	0.35	0	17,19,21	0.54	0
3	NAG	I	1	1,3	14,14,15	1.46	1 (7%)	17,19,21	1.41	3 (17%)
3	NAG	I	2	3	14,14,15	0.24	0	17,19,21	0.47	0
2	NAG	J	1	1,2	14,14,15	0.43	0	17,19,21	0.67	0
2	NAG	J	2	2	14,14,15	0.23	0	17,19,21	0.52	0
2	BMA	J	3	2	11,11,12	0.82	1 (9%)	15,15,17	0.80	0
3	NAG	K	1	1,3	14,14,15	0.52	0	17,19,21	1.05	2 (11%)
3	NAG	K	2	3	14,14,15	0.38	0	17,19,21	0.83	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	1,2	-	4/6/23/26	0/1/1/1
2	NAG	E	2	2	-	4/6/23/26	0/1/1/1
2	BMA	E	3	2	-	2/2/19/22	0/1/1/1
3	NAG	F	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1
2	NAG	G	1	1,2	-	3/6/23/26	0/1/1/1
2	NAG	G	2	2	-	0/6/23/26	0/1/1/1
2	BMA	G	3	2	-	1/2/19/22	0/1/1/1
3	NAG	H	1	1,3	-	5/6/23/26	0/1/1/1
3	NAG	H	2	3	-	2/6/23/26	0/1/1/1
3	NAG	I	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	I	2	3	-	2/6/23/26	0/1/1/1
2	NAG	J	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	J	2	2	-	0/6/23/26	0/1/1/1

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

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	I	1	NAG	O5-C1	-5.29	1.35	1.43
2	E	1	NAG	O5-C1	-3.66	1.37	1.43
2	E	2	NAG	O5-C1	-3.19	1.38	1.43
2	J	3	BMA	C1-C2	2.06	1.56	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	1	NAG	C2-N2-C7	4.28	129.00	122.90
2	G	2	NAG	O4-C4-C5	3.67	118.42	109.30
3	I	1	NAG	C4-C3-C2	3.67	116.39	111.02
3	I	1	NAG	C3-C4-C5	3.59	116.65	110.24
2	G	2	NAG	C1-O5-C5	3.42	116.83	112.19
2	E	1	NAG	O4-C4-C5	-3.12	101.55	109.30
2	G	3	BMA	C1-O5-C5	2.92	116.14	112.19
3	K	1	NAG	C1-O5-C5	2.75	115.92	112.19
3	H	1	NAG	C1-O5-C5	2.43	115.49	112.19
2	E	1	NAG	C3-C4-C5	2.35	114.43	110.24
3	I	1	NAG	C1-O5-C5	-2.29	109.09	112.19
3	H	1	NAG	C1-C2-N2	2.19	114.23	110.49
3	K	2	NAG	C1-C2-N2	2.15	114.16	110.49
3	F	1	NAG	C1-O5-C5	2.12	115.07	112.19
3	K	1	NAG	O4-C4-C3	-2.08	105.55	110.35

There are no chirality outliers.

All (37) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	1	NAG	C4-C5-C6-O6
2	E	2	NAG	O5-C5-C6-O6
2	E	1	NAG	O5-C5-C6-O6
3	H	1	NAG	O5-C5-C6-O6
2	E	2	NAG	C4-C5-C6-O6
3	H	1	NAG	C4-C5-C6-O6

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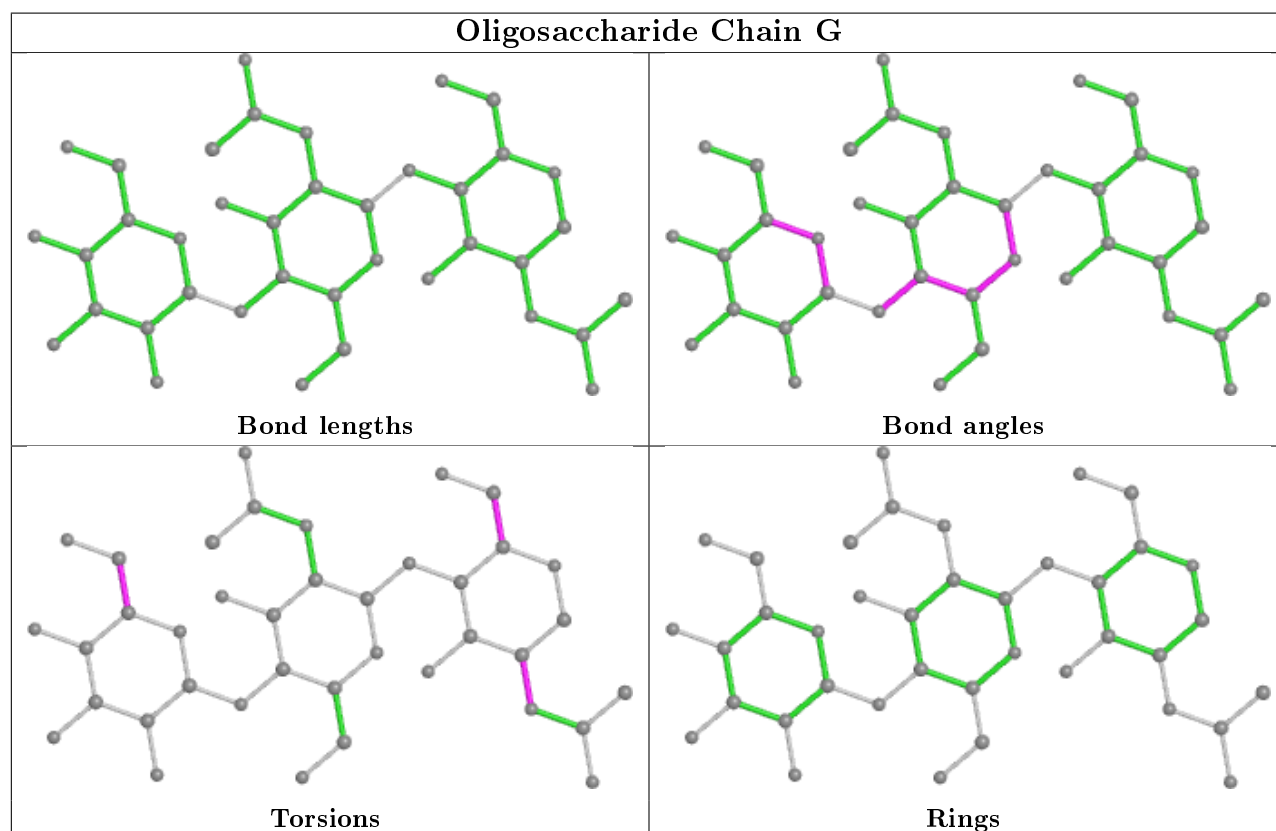
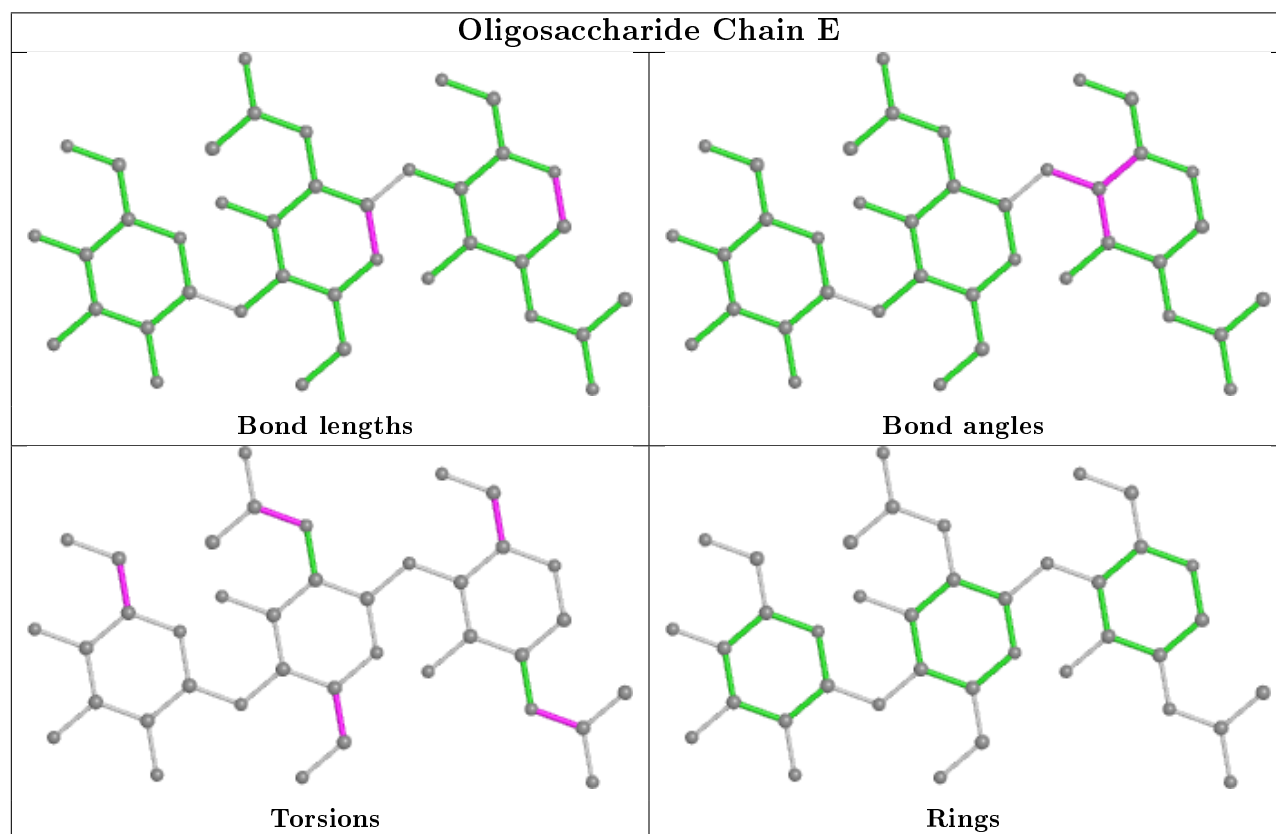
Mol	Chain	Res	Type	Atoms
2	E	1	NAG	C8-C7-N2-C2
2	E	1	NAG	O7-C7-N2-C2
2	E	2	NAG	C8-C7-N2-C2
2	E	2	NAG	O7-C7-N2-C2
3	H	2	NAG	C8-C7-N2-C2
3	H	2	NAG	O7-C7-N2-C2
3	K	1	NAG	C8-C7-N2-C2
3	K	1	NAG	O7-C7-N2-C2
3	H	1	NAG	C8-C7-N2-C2
3	H	1	NAG	O7-C7-N2-C2
2	G	1	NAG	O5-C5-C6-O6
2	G	1	NAG	C4-C5-C6-O6
3	K	2	NAG	C1-C2-N2-C7
3	I	2	NAG	O5-C5-C6-O6
2	J	1	NAG	C1-C2-N2-C7
3	I	1	NAG	C1-C2-N2-C7
2	E	3	BMA	C4-C5-C6-O6
3	F	2	NAG	C1-C2-N2-C7
2	J	3	BMA	O5-C5-C6-O6
2	G	3	BMA	O5-C5-C6-O6
3	K	2	NAG	C4-C5-C6-O6
3	K	2	NAG	O5-C5-C6-O6
2	E	3	BMA	O5-C5-C6-O6
2	J	1	NAG	C3-C2-N2-C7
3	I	2	NAG	C4-C5-C6-O6
2	G	1	NAG	C1-C2-N2-C7
3	F	1	NAG	C3-C2-N2-C7
3	F	2	NAG	C3-C2-N2-C7
3	H	1	NAG	C3-C2-N2-C7
3	I	1	NAG	C3-C2-N2-C7
3	F	1	NAG	C1-C2-N2-C7

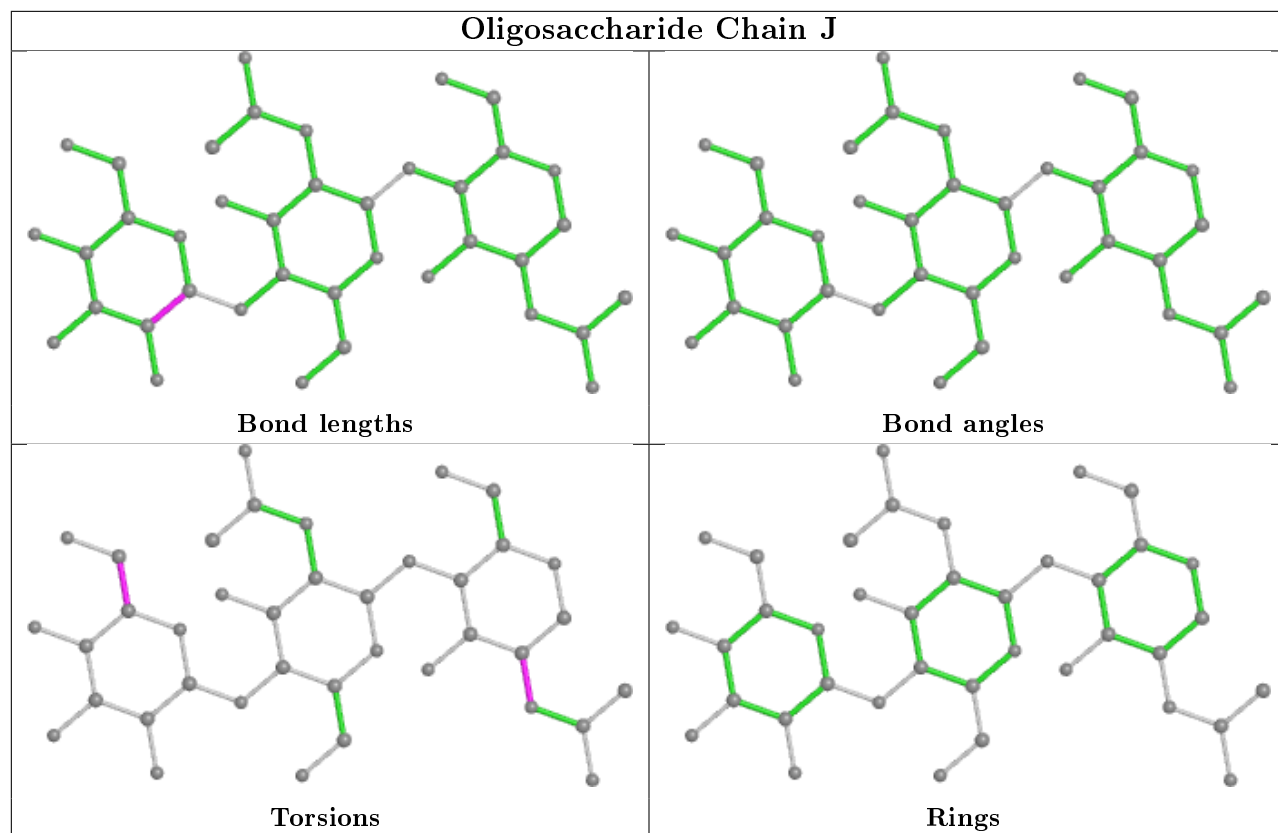
There are no ring outliers.

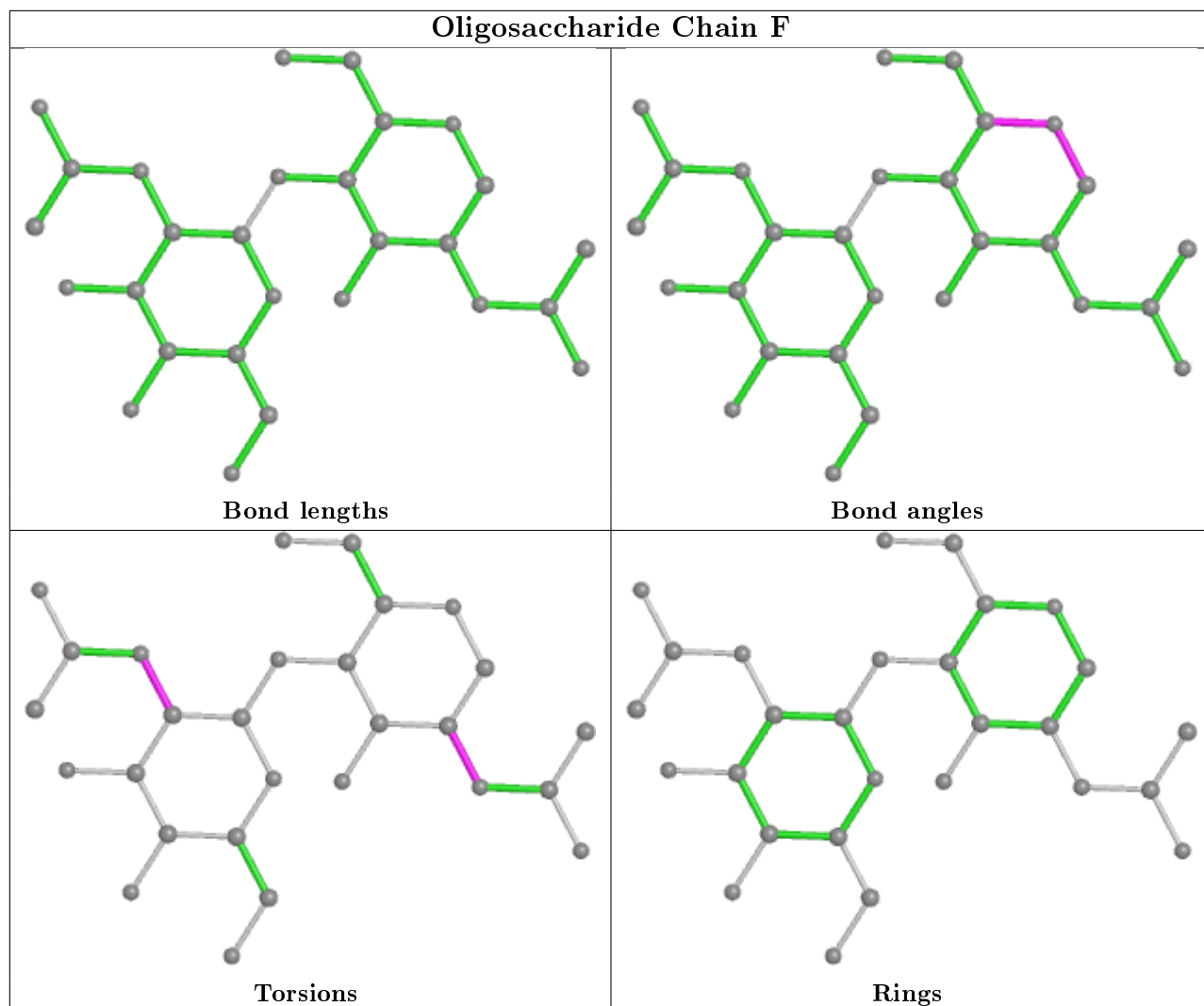
6 monomers are involved in 5 short contacts:

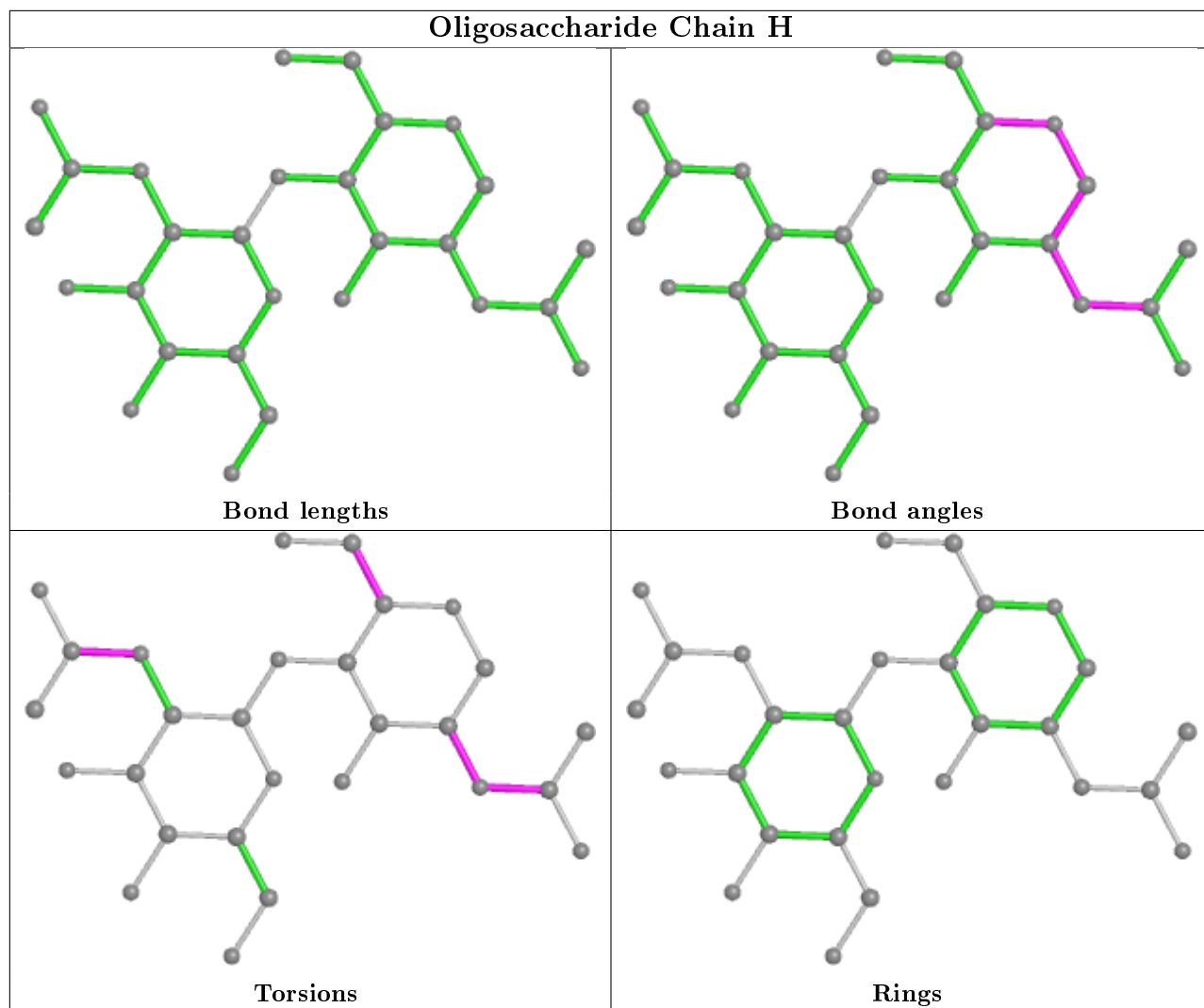
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	1	NAG	1	0
2	E	1	NAG	1	0
2	J	2	NAG	1	0
3	K	1	NAG	1	0
3	H	1	NAG	1	0
2	J	1	NAG	1	0

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

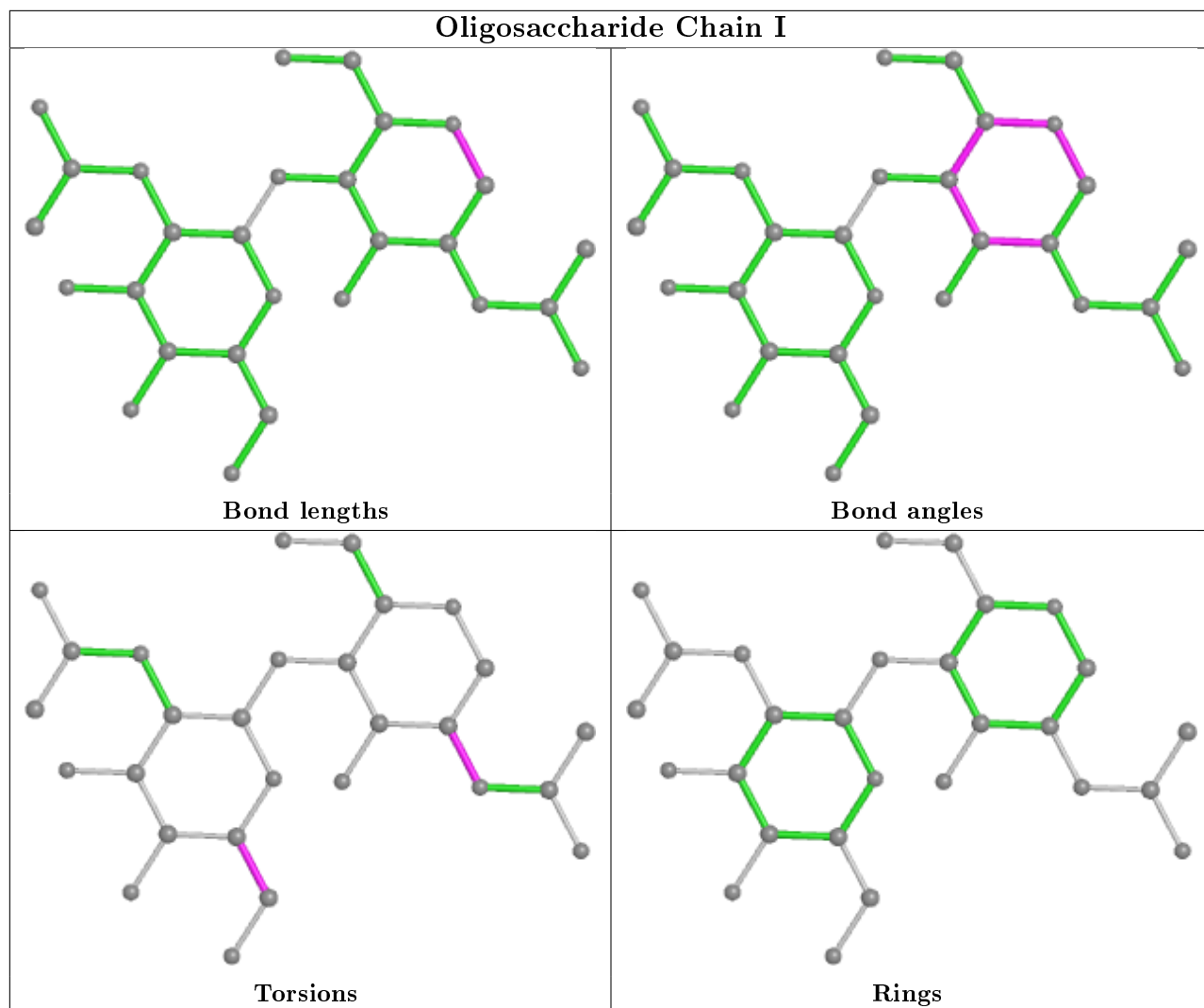


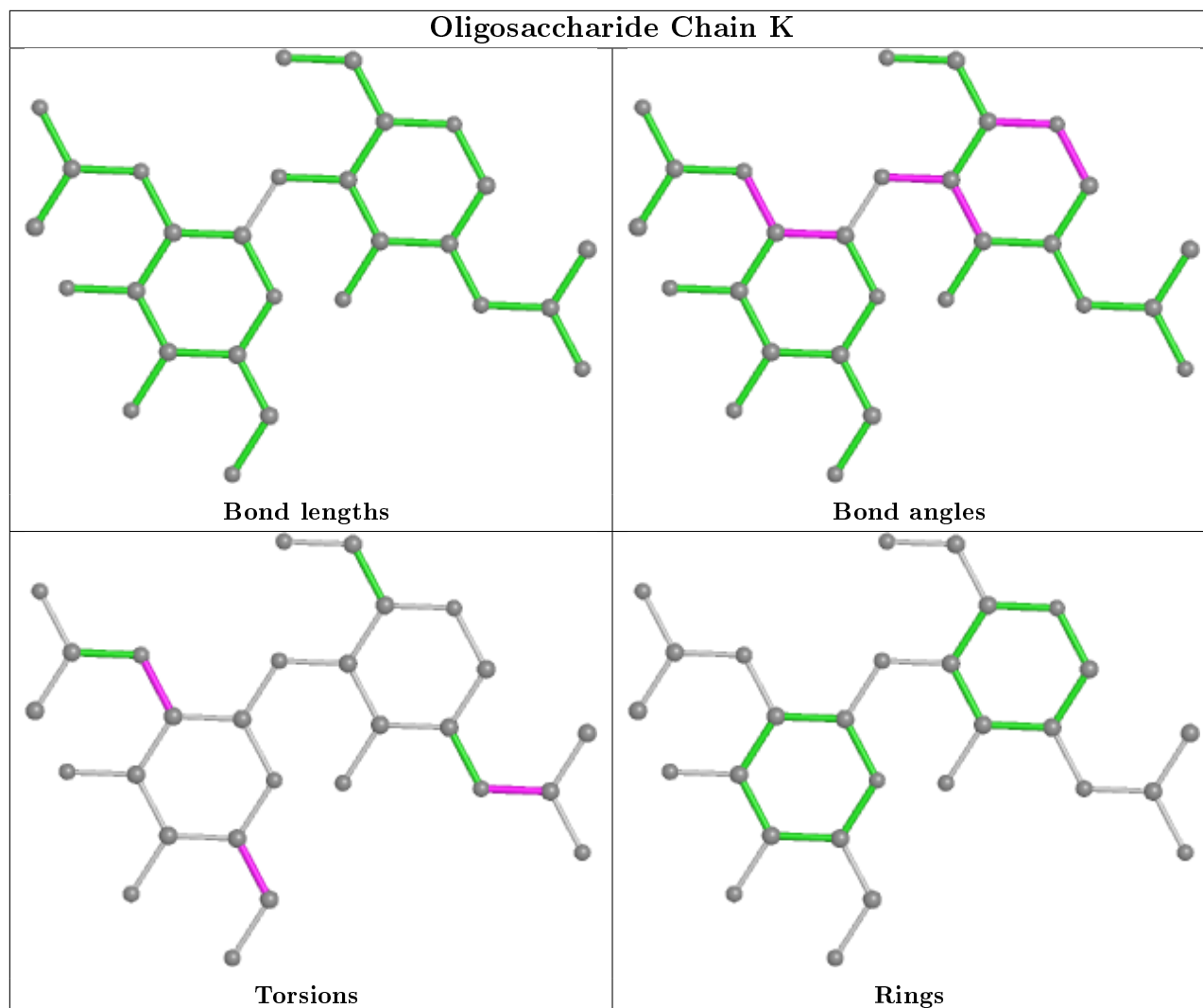












## 5.6 Ligand geometry [i](#)

1 ligand is 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	401	1	14,14,15	0.39	0	17,19,21	0.64	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
4	NAG	C	401	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	401	NAG	C1-O5-C5	2.25	115.24	112.19

There are no chirality outliers.

There are no torsion outliers.

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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	197/288 (68%)	0.18	4 (2%) 65 56	230, 313, 383, 413	0
1	B	196/288 (68%)	0.31	11 (5%) 24 20	250, 323, 380, 563	0
1	C	194/288 (67%)	0.33	5 (2%) 56 45	258, 329, 384, 418	0
1	D	196/288 (68%)	0.39	9 (4%) 32 27	257, 335, 390, 421	0
All	All	783/1152 (67%)	0.30	29 (3%) 41 33	230, 325, 385, 563	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	197	GLN	4.8
1	D	113	ASN	4.3
1	B	196	THR	4.1
1	C	137	GLU	4.1
1	B	113	ASN	3.4
1	C	136	VAL	3.3
1	A	243	GLU	3.0
1	B	295	THR	2.9
1	D	158	ASN	2.9
1	B	198	ASP	2.8
1	D	137	GLU	2.8
1	A	295	THR	2.8
1	B	121	LYS	2.7
1	D	138	ASN	2.7
1	D	117	ALA	2.7
1	D	303	ILE	2.6
1	C	138	ASN	2.5
1	B	247	GLY	2.4
1	B	137	GLU	2.4
1	D	257	VAL	2.4
1	D	136	VAL	2.3

*Continued on next page...*

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Mol	Chain	Res	Type	RSRZ
1	C	267	SER	2.3
1	B	259	HIS	2.2
1	A	245	ILE	2.2
1	B	136	VAL	2.1
1	D	167	SER	2.1
1	B	294	LYS	2.1
1	C	135	GLN	2.0
1	A	250	ILE	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](#)

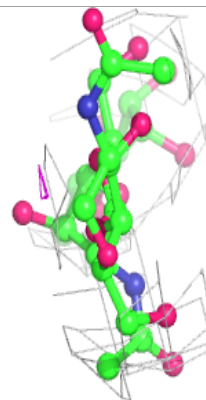
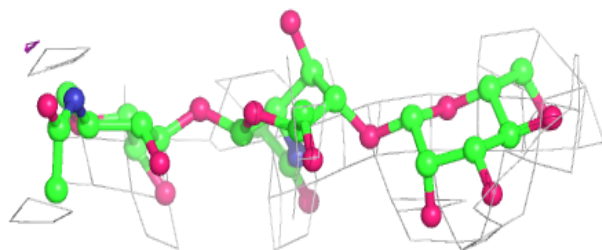
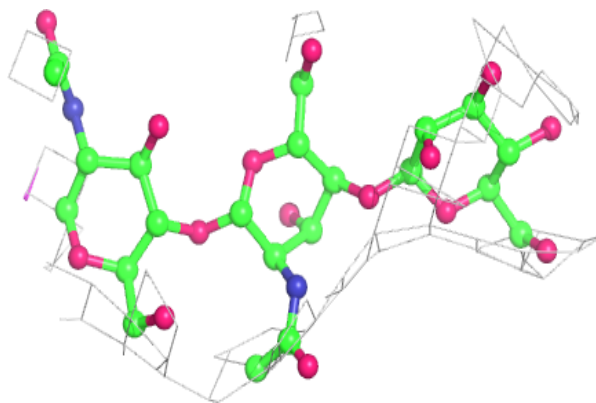
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	BMA	J	3	11/12	0.43	0.36	446,461,470,476	0
2	BMA	G	3	11/12	0.47	0.71	439,484,492,493	0
2	NAG	E	1	14/15	0.54	0.39	356,416,437,465	0
2	BMA	E	3	11/12	0.59	0.25	444,478,490,498	0
3	NAG	I	1	14/15	0.60	0.31	324,399,417,419	0
2	NAG	E	2	14/15	0.61	0.29	409,461,481,491	0
3	NAG	F	1	14/15	0.64	0.60	384,437,464,483	0
2	NAG	J	2	14/15	0.65	0.40	389,448,463,472	0
3	NAG	H	1	14/15	0.69	0.56	358,433,444,445	0
2	NAG	J	1	14/15	0.74	0.36	319,383,412,423	0
3	NAG	I	2	14/15	0.74	0.51	360,389,412,414	0
2	NAG	G	1	14/15	0.77	0.28	429,441,449,453	0
3	NAG	F	2	14/15	0.79	0.83	491,513,525,545	0
3	NAG	K	1	14/15	0.83	0.19	398,398,398,398	0
3	NAG	K	2	14/15	0.88	0.25	417,417,417,417	0
3	NAG	H	2	14/15	0.91	0.28	393,415,427,428	0
2	NAG	G	2	14/15	0.92	0.29	402,439,455,475	0

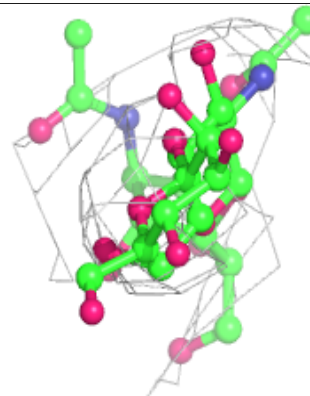
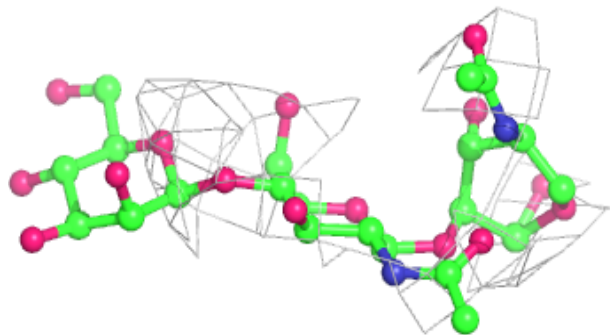
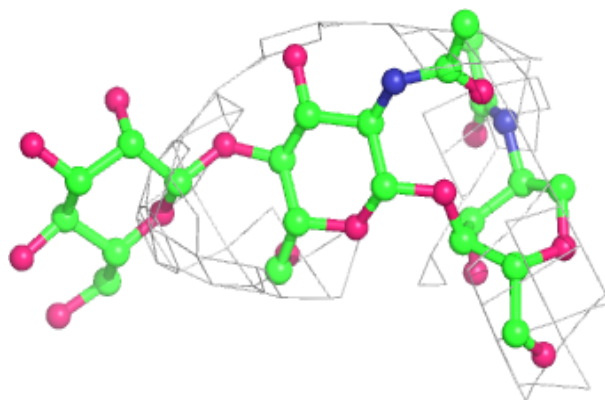
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

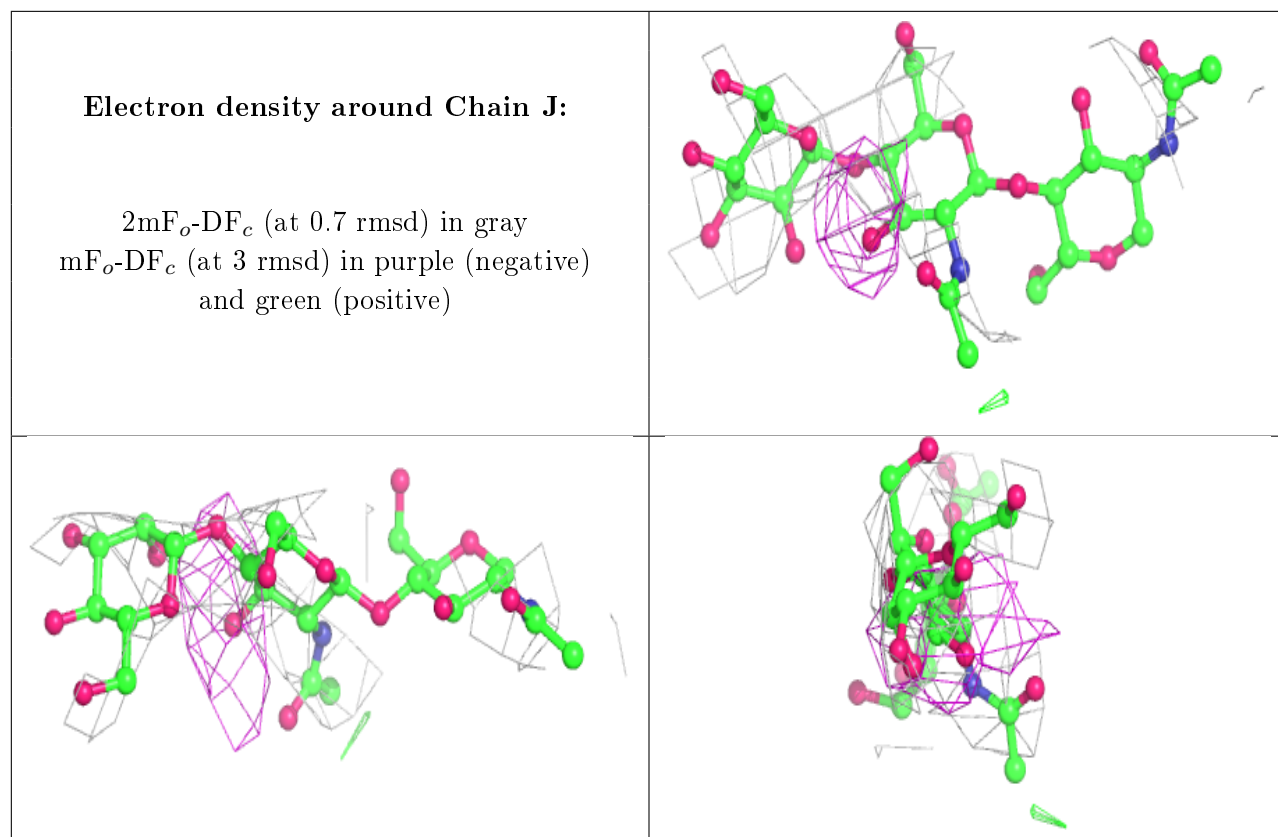
**Electron density around Chain E:**

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

**Electron density around Chain G:**

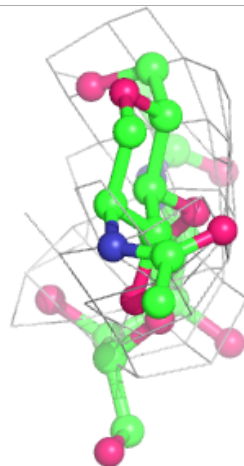
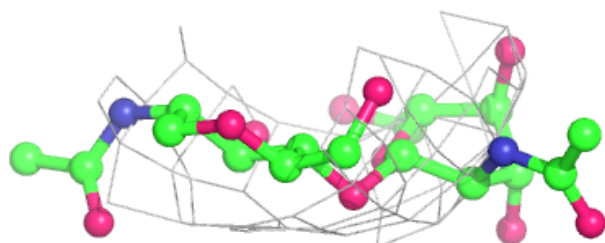
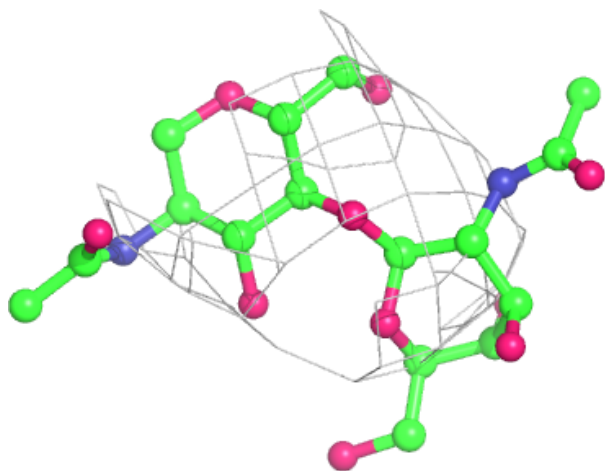
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





**Electron density around Chain F:**

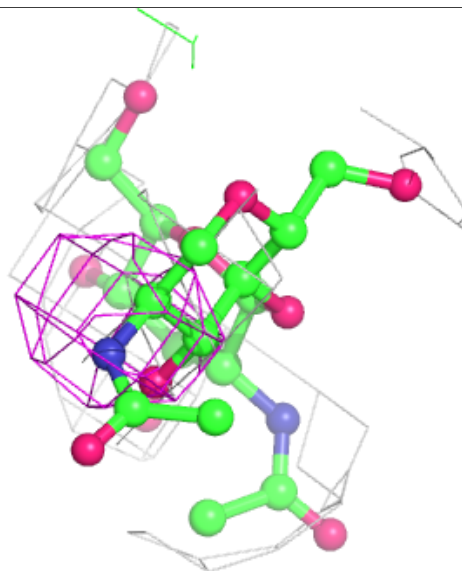
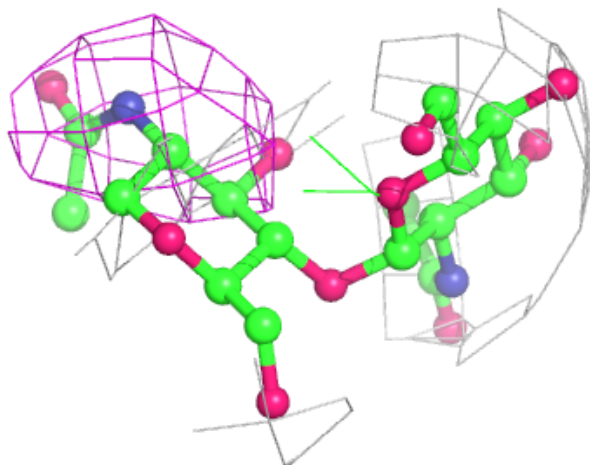
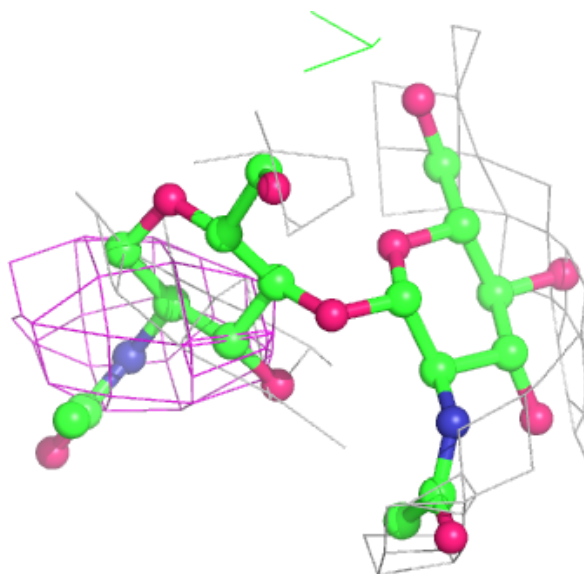
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





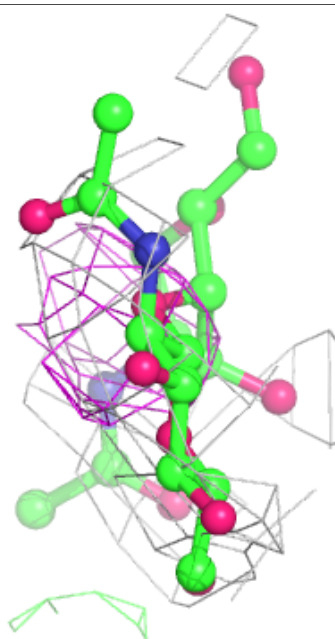
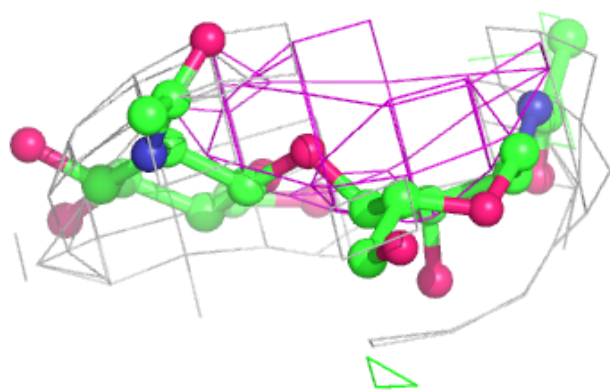
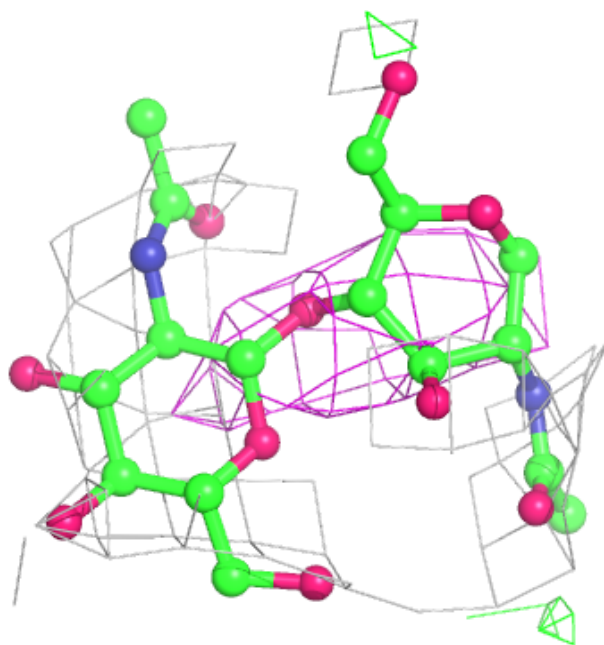
**Electron density around Chain H:**

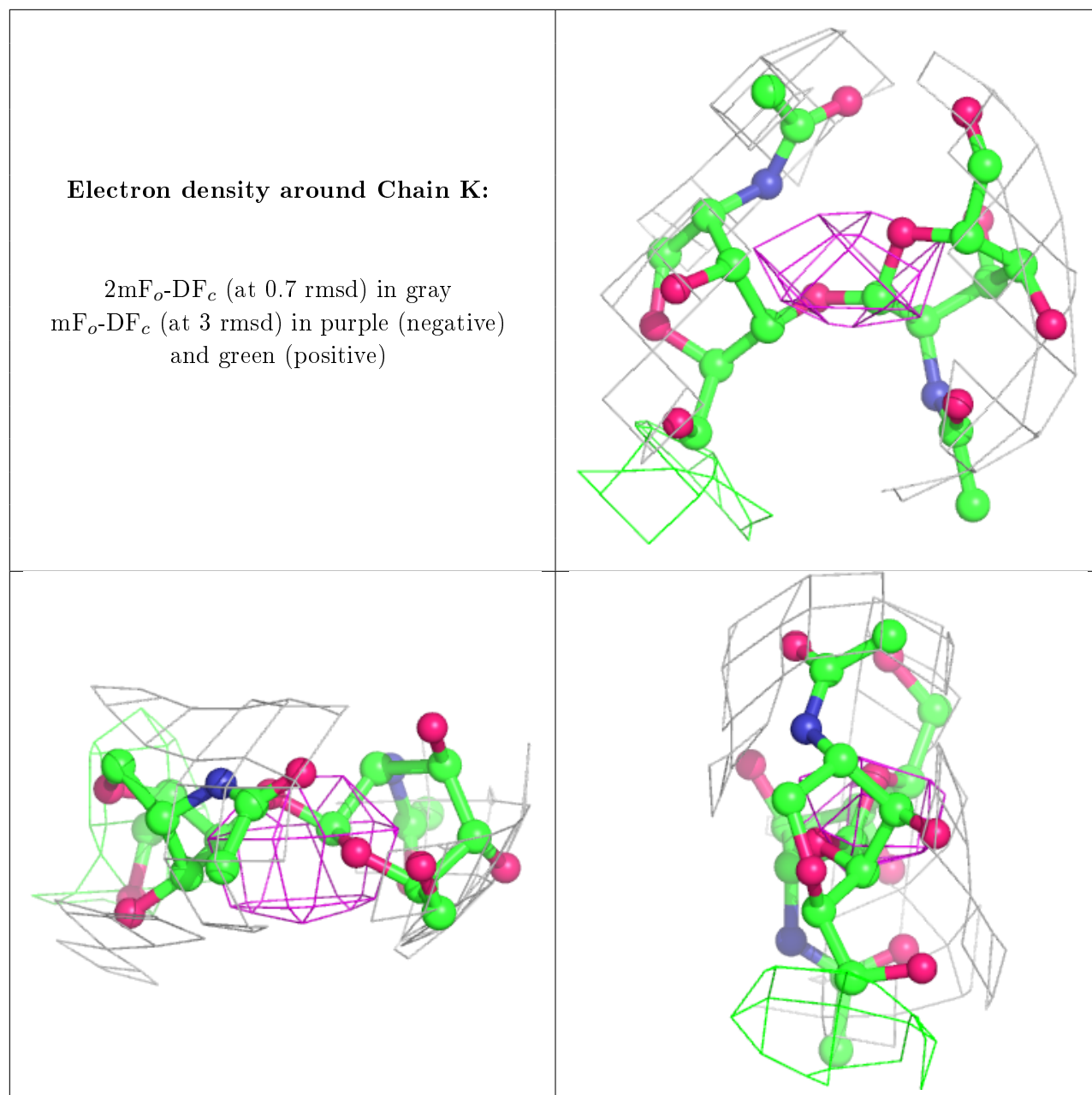
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**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(Å <sup>2</sup> )	Q<0.9
4	NAG	C	401	14/15	0.16	0.79	380,434,447,457	0

## 6.5 Other polymers

There are no such residues in this entry.