



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 16, 2021 – 12:08 PM JST

PDB ID : 7DRV  
Title : Structural basis of SARS-CoV-2-closely-related bat coronavirus RaTG13 to hACE2  
Authors : Liu, K.F.; Pan, X.Q.; Li, L.J.; Feng, Y.; Meng, Y.M.; Zhang, Y.F.; Wu, L.L.; Chen, Q.; Zheng, A.Q.; Song, C.L.; Jia, Y.F.; Niu, S.; Qiao, C.P.; Zhao, X.; Ma, D.L.; Ma, X.P.; Tan, S.G.; Qi, J.X.; Gao, G.F.; Wang, Q.H.  
Deposited on : 2020-12-29  
Resolution : 3.09 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:

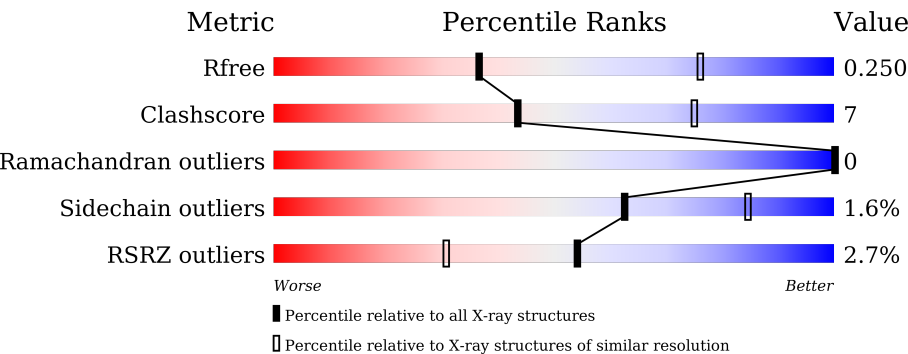
MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.23.1  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : 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 3.09 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	1094 (3.10-3.10)
Clashscore	141614	1184 (3.10-3.10)
Ramachandran outliers	138981	1141 (3.10-3.10)
Sidechain outliers	138945	1141 (3.10-3.10)
RSRZ outliers	127900	1067 (3.10-3.10)

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	596	<div><div></div><div></div><div></div><div></div><div></div></div> <div>87%13%</div>
1	B	596	<div><div></div><div></div><div></div><div></div><div></div></div> <div>%78%21%</div>
2	C	223	<div><div></div><div></div><div></div><div></div><div></div></div> <div>2%74%13%13%</div>
2	D	223	<div><div></div><div></div><div></div><div></div><div></div></div> <div>14%64%23%13%</div>
3	E	2	<div><div></div><div></div><div></div><div></div><div></div></div> <div>100%</div>
3	F	2	<div><div></div><div></div><div></div><div></div><div></div></div> <div>50%50%</div>

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Mol	Chain	Length	Quality of chain
3	G	2	

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
5	ZN	A	703	-	-	-	X

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 12916 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 Angiotensin-converting enzyme 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	596	Total	C	N	O	S	0	0	0
			4856	3108	803	916	29			
1	B	596	Total	C	N	O	S	0	0	0
			4851	3103	803	916	29			

- Molecule 2 is a protein called Spike glycoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	195	Total	C	N	O	S	0	0	0
			1532	984	253	287	8			
2	D	195	Total	C	N	O	S	0	0	0
			1521	978	249	286	8			

- 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	E	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	G	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>) (labeled as "Ligand of Interest" by depositor).



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

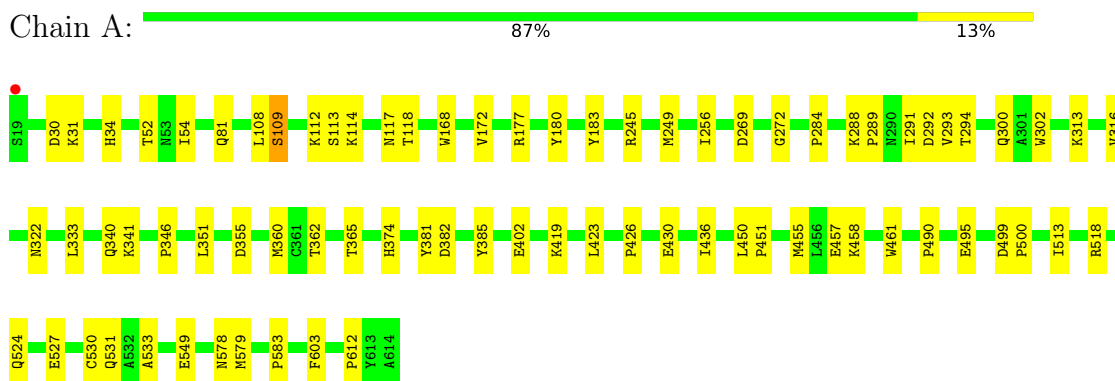
- Molecule 5 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Zn	0	0
			1	1		
5	B	1	Total	Zn	0	0
			1	1		

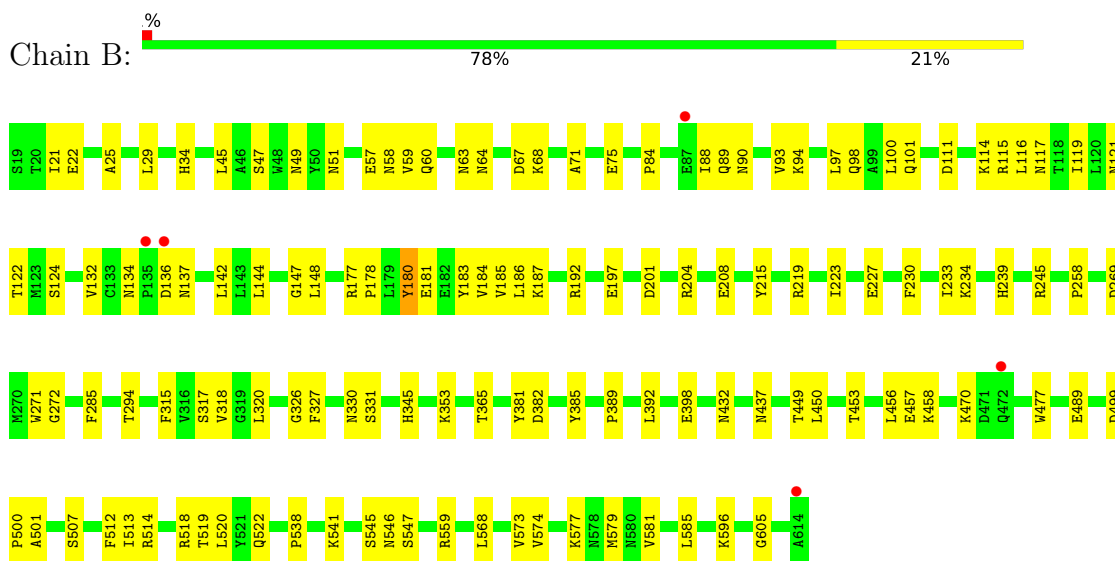
### 3 Residue-property plots

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.

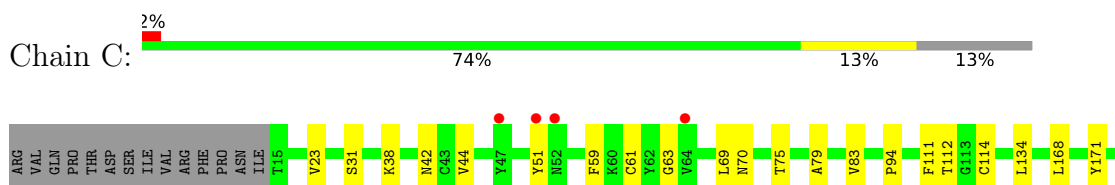
- Molecule 1: Angiotensin-converting enzyme 2

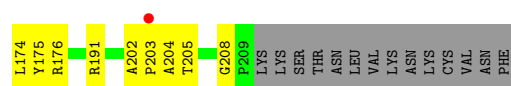


- Molecule 1: Angiotensin-converting enzyme 2

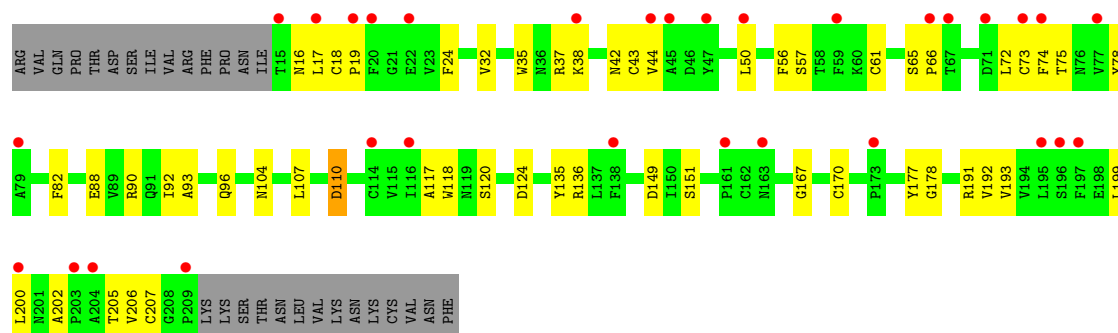


- Molecule 2: Spike glycoprotein





• Molecule 2: Spike glycoprotein



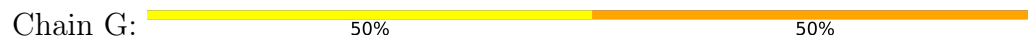
• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



• 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, $\alpha$ , $\beta$ , $\gamma$	81.28Å 122.88Å 110.74Å 90.00° 95.07° 90.00°	Depositor
Resolution (Å)	48.94 – 3.09 48.94 – 3.09	Depositor EDS
% Data completeness (in resolution range)	99.4 (48.94-3.09) 99.4 (48.94-3.09)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.15 (at 3.07Å)	Xtriage
Refinement program	PHENIX 1.16_3549	Depositor
R, $R_{free}$	0.226 , 0.250 0.226 , 0.250	Depositor DCC
$R_{free}$ test set	1934 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	95.0	Xtriage
Anisotropy	0.100	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.26 , 50.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	12916	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	118.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.50% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, 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	0.23	0/4993	0.38	0/6785
1	B	0.23	0/4988	0.39	0/6779
2	C	0.24	0/1577	0.43	0/2148
2	D	0.26	0/1566	0.46	0/2137
All	All	0.24	0/13124	0.40	0/17849

There are no bond length outliers.

There are no bond angle outliers.

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	4856	0	4621	45	0
1	B	4851	0	4611	70	0
2	C	1532	0	1430	18	0
2	D	1521	0	1405	37	0
3	E	28	0	25	0	0
3	F	28	0	25	1	0
3	G	28	0	25	2	0
4	A	42	0	39	1	0
4	B	14	0	13	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	C	14	0	13	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
All	All	12916	0	12207	168	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 (168) 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:A:374:HIS:NE2	1:A:402:GLU:OE2	2.13	0.81
1:B:201:ASP:HA	1:B:204:ARG:HE	1.46	0.81
1:A:457:GLU:HG2	1:A:513:ILE:HB	1.66	0.77
1:A:30:ASP:O	1:A:34:HIS:ND1	2.25	0.69
2:D:17:LEU:HA	2:D:44:VAL:HB	1.75	0.68
2:D:24:PHE:HE1	2:D:193:VAL:HG11	1.60	0.65
1:B:520:LEU:HD22	1:B:579:MET:HE2	1.78	0.64
1:B:477:TRP:NE1	1:B:499:ASP:OD2	2.31	0.63
1:A:245:ARG:NH1	1:A:603:PHE:O	2.33	0.62
1:B:144:LEU:HA	1:B:148:LEU:HB2	1.82	0.61
1:B:121:ASN:HA	1:B:124:SER:HB3	1.82	0.61
2:D:32:VAL:HG13	2:D:104:ASN:HD22	1.66	0.61
2:D:19:PRO:O	2:D:38:LYS:NZ	2.34	0.60
1:B:326:GLY:O	1:B:330:ASN:ND2	2.33	0.60
1:A:419:LYS:NZ	1:A:430:GLU:OE2	2.33	0.60
1:A:288:LYS:HG3	1:A:289:PRO:HD2	1.83	0.59
1:B:389:PRO:HG2	1:B:392:LEU:HB2	1.85	0.59
2:D:72:LEU:HD22	2:D:73:CYS:H	1.67	0.58
1:B:208:GLU:OE2	1:B:219:ARG:HD2	2.03	0.58
3:G:1:NAG:H62	3:G:2:NAG:HN2	1.68	0.58
2:D:16:ASN:ND2	2:D:42:ASN:O	2.28	0.57
1:B:233:ILE:HD13	1:B:450:LEU:HD13	1.85	0.57
1:A:351:LEU:HD13	1:A:355:ASP:OD2	2.05	0.57
1:A:52:THR:O	1:A:340:GLN:NE2	2.38	0.57
1:A:419:LYS:HD3	1:A:426:PRO:HA	1.86	0.57
1:A:524:GLN:HG2	1:A:583:PRO:HG2	1.87	0.56
1:B:581:VAL:HG12	1:B:585:LEU:HD13	1.88	0.55
1:B:204:ARG:HD2	1:B:223:ILE:HD11	1.87	0.55
1:B:573:VAL:HG13	1:B:574:VAL:HG13	1.90	0.54
1:A:291:ILE:HG13	1:A:292:ASP:H	1.71	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:45:LEU:O	1:B:49:ASN:ND2	2.30	0.54
1:B:90:ASN:HB3	1:B:93:VAL:HG22	1.89	0.54
2:C:75:THR:HA	2:C:204:ALA:HA	1.90	0.54
1:B:111:ASP:HA	1:B:114:LYS:HB2	1.90	0.54
1:B:21:ILE:HG21	1:B:84:PRO:HD2	1.90	0.54
1:B:227:GLU:OE1	1:B:458:LYS:NZ	2.33	0.54
1:B:245:ARG:NH1	1:B:605:GLY:O	2.28	0.54
1:B:285:PHE:HB2	1:B:437:ASN:HD21	1.72	0.54
2:D:75:THR:O	2:D:205:THR:OG1	2.20	0.53
2:D:120:SER:OG	2:D:124:ASP:OD1	2.26	0.53
1:B:353:LYS:NZ	2:D:178:GLY:O	2.27	0.53
2:C:94:PRO:HG3	2:C:111:PHE:HB3	1.90	0.53
1:A:81:GLN:HE22	3:F:1:NAG:H82	1.74	0.53
1:A:333:LEU:O	1:A:362:THR:OG1	2.25	0.52
2:C:69:LEU:HD23	2:C:69:LEU:O	2.09	0.52
1:A:31:LYS:NZ	2:C:175:TYR:OH	2.37	0.52
1:A:294:THR:HG23	1:A:365:THR:HA	1.92	0.52
1:A:269:ASP:OD2	1:A:272:GLY:N	2.42	0.52
1:B:63:ASN:O	1:B:67:ASP:N	2.34	0.52
1:B:538:PRO:HD2	1:B:541:LYS:HD3	1.92	0.52
1:B:185:VAL:HG23	1:B:186:LEU:HD12	1.91	0.52
1:B:457:GLU:OE1	1:B:513:ILE:N	2.43	0.51
1:B:116:LEU:HD21	1:B:187:LYS:HE2	1.92	0.51
1:B:192:ARG:NE	1:B:197:GLU:O	2.43	0.51
1:A:293:VAL:HG21	1:A:423:LEU:HB3	1.92	0.50
1:B:34:HIS:ND1	2:D:135:TYR:OH	2.44	0.50
3:G:1:NAG:H62	3:G:2:NAG:N2	2.26	0.50
1:A:313:LYS:HA	1:A:316:VAL:HG22	1.93	0.50
1:B:47:SER:O	1:B:51:ASN:ND2	2.37	0.49
2:D:57:SER:N	2:D:117:ALA:O	2.45	0.49
2:C:134:LEU:HD12	2:C:174:LEU:HB3	1.94	0.49
1:B:116:LEU:HD11	1:B:187:LYS:HG2	1.95	0.49
2:C:23:VAL:HG22	2:C:38:LYS:HE2	1.95	0.49
2:C:63:GLY:HA3	2:C:112:THR:HA	1.94	0.49
2:C:202:ALA:HB1	2:C:203:PRO:HD2	1.95	0.49
1:A:177:ARG:NH2	1:A:495:GLU:O	2.46	0.49
2:C:83:VAL:HG22	2:C:191:ARG:HG2	1.95	0.49
2:D:93:ALA:HB3	2:D:96:GLN:HG3	1.95	0.49
1:A:113:SER:O	1:A:117:ASN:ND2	2.34	0.48
1:B:223:ILE:H	1:B:223:ILE:HD12	1.78	0.48
1:A:578:ASN:OD1	1:A:579:MET:N	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:43:CYS:H	2:D:206:VAL:HG22	1.78	0.48
1:B:545:SER:HB2	4:B:802:NAG:H82	1.96	0.47
1:A:249:MET:HG2	1:A:256:ILE:HB	1.96	0.47
1:A:490:PRO:HA	1:A:612:PRO:HG2	1.97	0.47
2:C:23:VAL:HG11	2:C:79:ALA:HB1	1.96	0.47
1:A:168:TRP:O	1:A:172:VAL:HG22	2.14	0.47
1:B:184:VAL:HG11	1:B:470:LYS:HZ1	1.78	0.47
2:D:18:CYS:HA	2:D:43:CYS:HB2	1.97	0.47
1:A:300:GLN:HB3	1:A:302:TRP:CD1	2.50	0.47
1:A:530:CYS:HA	1:A:533:ALA:HB3	1.96	0.47
2:C:63:GLY:HA3	2:C:112:THR:HG23	1.96	0.47
1:B:499:ASP:N	1:B:500:PRO:HD2	2.30	0.47
2:D:200:LEU:HB2	2:D:202:ALA:HB2	1.96	0.47
1:B:132:VAL:HG21	1:B:148:LEU:HD11	1.97	0.47
1:B:453:THR:HG22	1:B:512:PHE:HE1	1.80	0.47
2:D:120:SER:HB3	2:D:191:ARG:HG2	1.97	0.47
1:A:402:GLU:HB3	1:A:518:ARG:HG3	1.97	0.46
2:C:61:CYS:HA	2:C:114:CYS:HA	1.97	0.46
1:A:54:ILE:HB	1:A:341:LYS:HG3	1.98	0.46
1:A:180:TYR:HA	1:A:183:TYR:HB3	1.98	0.46
1:B:177:ARG:HB3	1:B:178:PRO:HD3	1.97	0.46
2:D:65:SER:HB3	2:D:66:PRO:HD2	1.98	0.45
1:B:134:ASN:HB3	1:B:137:ASN:O	2.16	0.45
2:D:56:PHE:HA	2:D:118:TRP:HB3	1.97	0.45
1:A:109:SER:OG	1:A:112:LYS:HG2	2.16	0.45
2:D:149:ASP:OD2	2:D:151:SER:HB3	2.17	0.45
1:B:449:THR:O	1:B:453:THR:HG23	2.17	0.45
2:D:19:PRO:HG2	2:D:38:LYS:HZ3	1.82	0.45
1:B:215:TYR:CZ	1:B:568:LEU:HD13	2.51	0.44
1:B:239:HIS:CE1	1:B:596:LYS:HG2	2.52	0.44
1:B:294:THR:HG23	1:B:365:THR:HA	1.99	0.44
1:B:432:ASN:OD1	1:B:432:ASN:N	2.49	0.44
1:A:31:LYS:HD2	2:C:171:TYR:CD2	2.52	0.44
2:D:35:TRP:CZ3	2:D:37:ARG:HB2	2.52	0.44
2:D:104:ASN:HD21	2:D:136:ARG:H	1.64	0.44
1:B:134:ASN:OD1	1:B:136:ASP:N	2.51	0.44
1:A:499:ASP:N	1:A:500:PRO:HD2	2.33	0.44
1:B:269:ASP:OD1	1:B:272:GLY:N	2.49	0.44
2:D:61:CYS:SG	2:D:66:PRO:HD3	2.58	0.44
2:C:70:ASN:OD1	2:C:70:ASN:N	2.51	0.43
1:A:114:LYS:NZ	1:A:118:THR:OG1	2.50	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:315:PHE:HD1	1:B:320:LEU:HD12	1.83	0.43
1:B:382:ASP:HA	1:B:385:TYR:CZ	2.53	0.43
1:B:64:ASN:O	1:B:68:LYS:HG2	2.18	0.43
1:B:317:SER:HB2	1:B:546:ASN:H	1.84	0.43
1:A:382:ASP:HA	1:A:385:TYR:CZ	2.53	0.43
1:B:88:ILE:HD13	1:B:94:LYS:HD3	2.00	0.43
2:D:24:PHE:CE1	2:D:193:VAL:HG11	2.47	0.43
2:D:74:PHE:HB2	2:D:199:LEU:HB3	2.01	0.43
1:A:322:ASN:HD22	4:A:702:NAG:C7	2.32	0.43
1:B:59:VAL:HG13	1:B:60:GLN:N	2.34	0.43
1:B:177:ARG:O	1:B:181:GLU:HG3	2.19	0.43
1:B:398:GLU:HB2	1:B:514:ARG:HG2	2.01	0.43
2:D:88:GLU:OE1	2:D:177:TYR:OH	2.23	0.42
2:C:44:VAL:HG13	2:C:208:GLY:HA2	2.01	0.42
2:C:168:LEU:HD12	2:C:168:LEU:H	1.83	0.42
2:D:110:ASP:OD1	2:D:110:ASP:N	2.52	0.42
2:D:167:GLY:H	2:D:170:CYS:HB2	1.84	0.42
1:B:119:ILE:HA	1:B:122:THR:HG1	1.84	0.42
1:B:142:LEU:HG	1:B:147:GLY:HA3	2.02	0.42
1:B:547:SER:O	1:B:547:SER:OG	2.36	0.42
1:A:346:PRO:HG3	1:A:360:MET:HG3	2.02	0.42
2:C:42:ASN:H	2:C:205:THR:HB	1.83	0.42
2:D:92:ILE:HG23	2:D:107:LEU:HD11	2.00	0.42
1:A:527:GLU:O	1:A:531:GLN:HB2	2.19	0.42
1:B:230:PHE:O	1:B:234:LYS:HG3	2.19	0.42
1:B:245:ARG:HD3	1:B:258:PRO:HA	2.01	0.42
1:B:501:ALA:O	1:B:507:SER:HB3	2.19	0.42
1:A:54:ILE:HD12	1:A:341:LYS:HE2	2.01	0.42
1:A:341:LYS:HE3	1:A:341:LYS:HB2	1.82	0.42
2:D:37:ARG:HE	2:D:78:TYR:HB3	1.85	0.42
1:A:458:LYS:HA	1:A:461:TRP:HB2	2.02	0.42
1:B:98:GLN:HA	1:B:101:GLN:HB2	2.02	0.42
2:D:136:ARG:NH2	2:D:149:ASP:OD2	2.53	0.41
1:B:71:ALA:O	1:B:75:GLU:HG2	2.20	0.41
2:D:35:TRP:HZ3	2:D:37:ARG:HB2	1.85	0.41
1:A:549:GLU:OE1	1:A:549:GLU:N	2.50	0.41
1:B:84:PRO:O	1:B:88:ILE:HD11	2.20	0.41
1:A:450:LEU:HB2	1:A:451:PRO:HD3	2.02	0.41
1:B:22:GLU:OE2	1:B:89:GLN:HG3	2.21	0.41
1:B:327:PHE:O	1:B:331:SER:OG	2.31	0.41
1:B:519:THR:O	1:B:522:GLN:HG2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:LEU:HG	1:A:112:LYS:HG3	2.02	0.41
1:A:423:LEU:HD23	1:A:423:LEU:HA	1.90	0.41
2:D:43:CYS:O	2:D:206:VAL:HG13	2.21	0.41
2:D:82:PHE:CE1	2:D:192:VAL:HG11	2.55	0.41
2:C:31:SER:HB2	2:C:134:LEU:O	2.21	0.40
1:B:59:VAL:HG13	1:B:60:GLN:H	1.85	0.40
1:A:284:PRO:HG2	1:A:436:ILE:HG22	2.04	0.40
1:B:180:TYR:HA	1:B:183:TYR:HB3	2.04	0.40
1:B:318:VAL:HG23	1:B:320:LEU:HG	2.02	0.40
1:B:93:VAL:O	1:B:97:LEU:HG	2.21	0.40
1:B:25:ALA:O	1:B:29:LEU:HG	2.21	0.40
1:B:215:TYR:CD1	1:B:577:LYS:HD3	2.57	0.40
2:D:24:PHE:HE2	2:D:50:LEU:HG	1.87	0.40
2:D:32:VAL:HB	2:D:82:PHE:CD1	2.57	0.40
2:D:117:ALA:HA	2:D:191:ARG:O	2.22	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	594/596 (100%)	583 (98%)	11 (2%)	0	100	100
1	B	594/596 (100%)	580 (98%)	14 (2%)	0	100	100
2	C	193/223 (86%)	178 (92%)	15 (8%)	0	100	100
2	D	193/223 (86%)	177 (92%)	16 (8%)	0	100	100
All	All	1574/1638 (96%)	1518 (96%)	56 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	524/526 (100%)	521 (99%)	3 (1%)	86	94
1	B	523/526 (99%)	510 (98%)	13 (2%)	47	75
2	C	163/195 (84%)	160 (98%)	3 (2%)	59	82
2	D	160/195 (82%)	157 (98%)	3 (2%)	57	81
All	All	1370/1442 (95%)	1348 (98%)	22 (2%)	62	84

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	109	SER
1	A	381	TYR
1	A	455	MET
2	C	51	TYR
2	C	59	PHE
2	C	176	ARG
1	B	57	GLU
1	B	58	ASN
1	B	100	LEU
1	B	115	ARG
1	B	117	ASN
1	B	180	TYR
1	B	271	TRP
1	B	345	HIS
1	B	381	TYR
1	B	456	LEU
1	B	489	GLU
1	B	518	ARG
1	B	559	ARG
2	D	90	ARG
2	D	110	ASP
2	D	207	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	D	104	ASN

### 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 ⓘ

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	E	1	1,3	14,14,15	0.28	0	17,19,21	0.47	0
3	NAG	E	2	3	14,14,15	0.29	0	17,19,21	0.39	0
3	NAG	F	1	1,3	14,14,15	0.31	0	17,19,21	0.66	1 (5%)
3	NAG	F	2	3	14,14,15	0.28	0	17,19,21	0.63	0
3	NAG	G	1	1,3	14,14,15	0.28	0	17,19,21	0.65	1 (5%)
3	NAG	G	2	3	14,14,15	0.30	0	17,19,21	0.46	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	E	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	E	2	3	-	0/6/23/26	0/1/1/1
3	NAG	F	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	F	2	3	-	0/6/23/26	0/1/1/1
3	NAG	G	1	1,3	-	0/6/23/26	0/1/1/1

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

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	1	NAG	C1-O5-C5	2.24	115.22	112.19
3	G	1	NAG	C1-O5-C5	2.19	115.17	112.19

There are no chirality outliers.

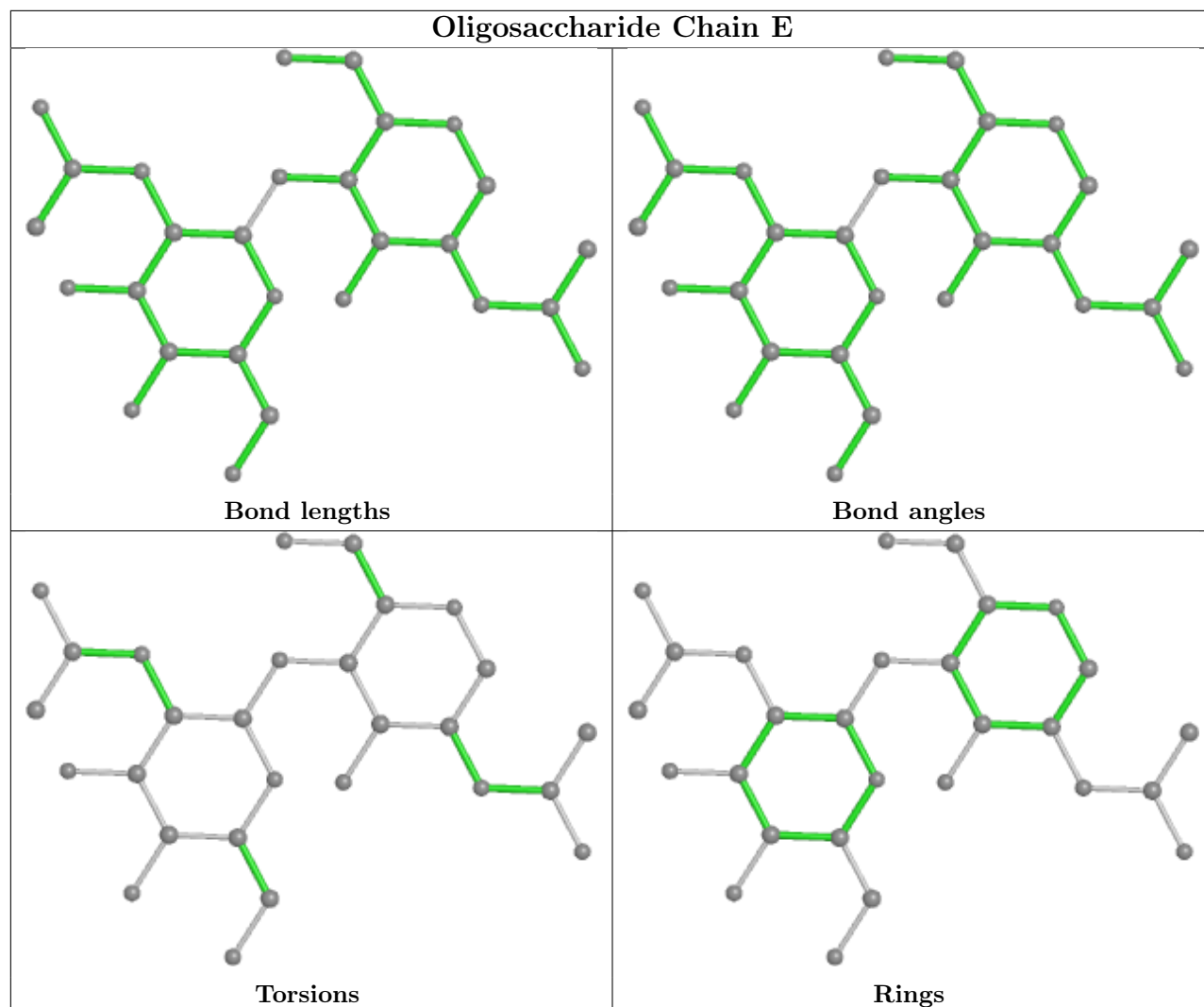
There are no torsion outliers.

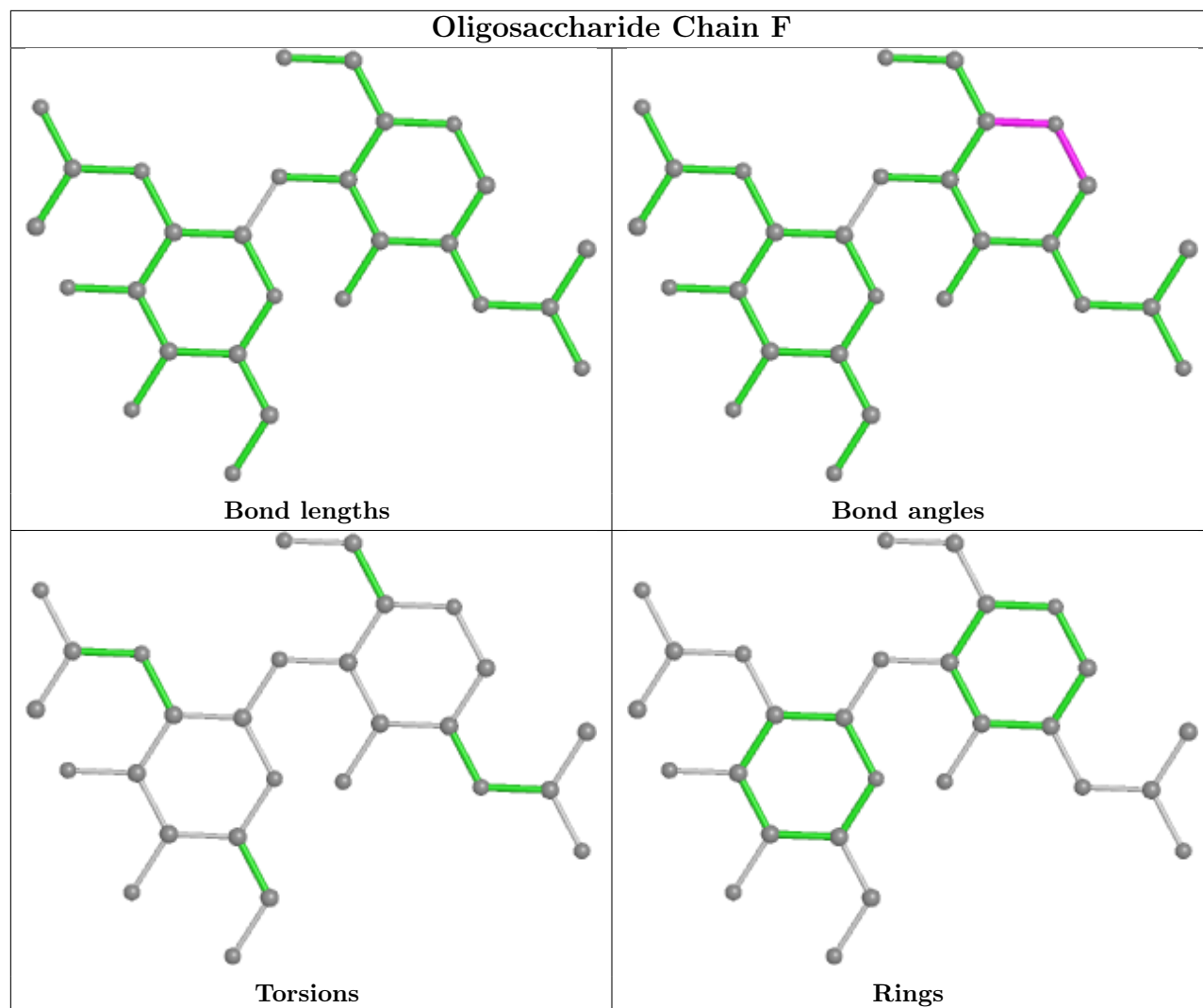
There are no ring outliers.

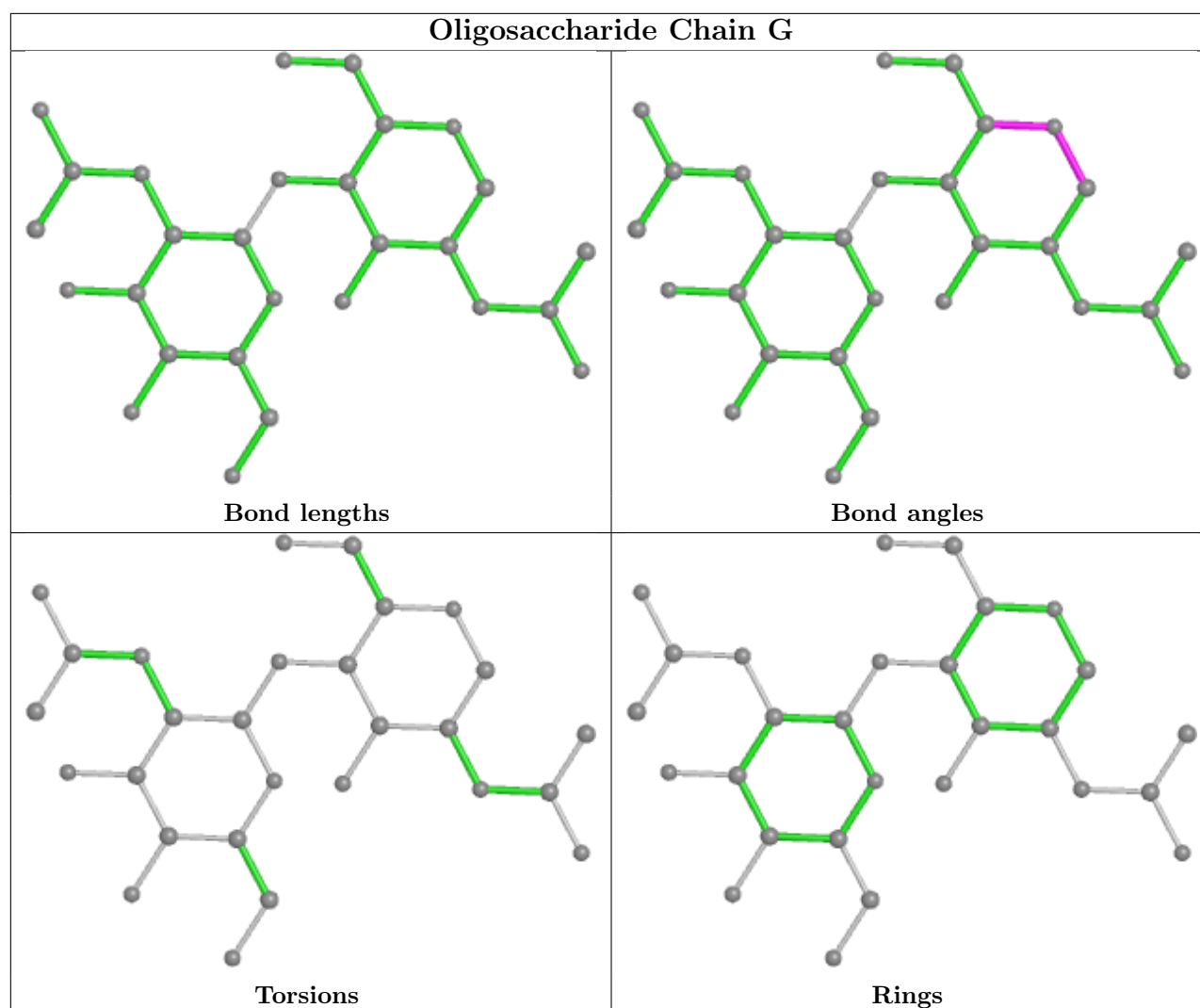
3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	1	NAG	1	0
3	G	2	NAG	2	0
3	G	1	NAG	2	0

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







## 5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
4	NAG	C	301	2	14,14,15	0.24	0	17,19,21	0.47	0
4	NAG	A	702	1	14,14,15	0.32	0	17,19,21	0.38	0
4	NAG	A	704	1	14,14,15	0.17	0	17,19,21	0.42	0
4	NAG	A	701	1	14,14,15	0.30	0	17,19,21	0.51	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	NAG	B	802	1	14,14,15	0.29	0	17,19,21	0.50	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	301	2	-	1/6/23/26	0/1/1/1
4	NAG	A	702	1	-	2/6/23/26	0/1/1/1
4	NAG	A	704	1	-	2/6/23/26	0/1/1/1
4	NAG	A	701	1	-	2/6/23/26	0/1/1/1
4	NAG	B	802	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	702	NAG	O5-C5-C6-O6
4	A	704	NAG	O5-C5-C6-O6
4	A	701	NAG	O5-C5-C6-O6
4	A	702	NAG	C4-C5-C6-O6
4	A	704	NAG	C4-C5-C6-O6
4	A	701	NAG	C4-C5-C6-O6
4	C	301	NAG	O5-C5-C6-O6

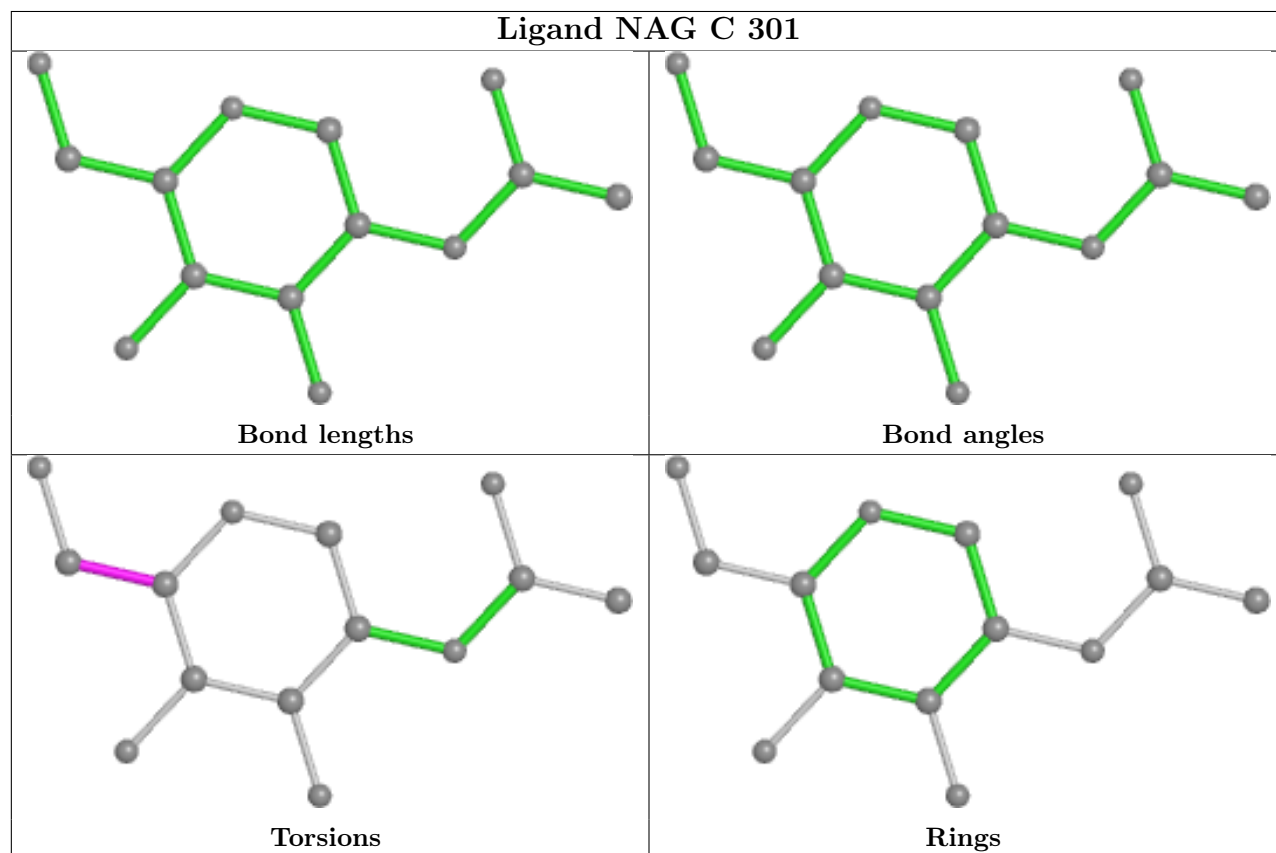
There are no ring outliers.

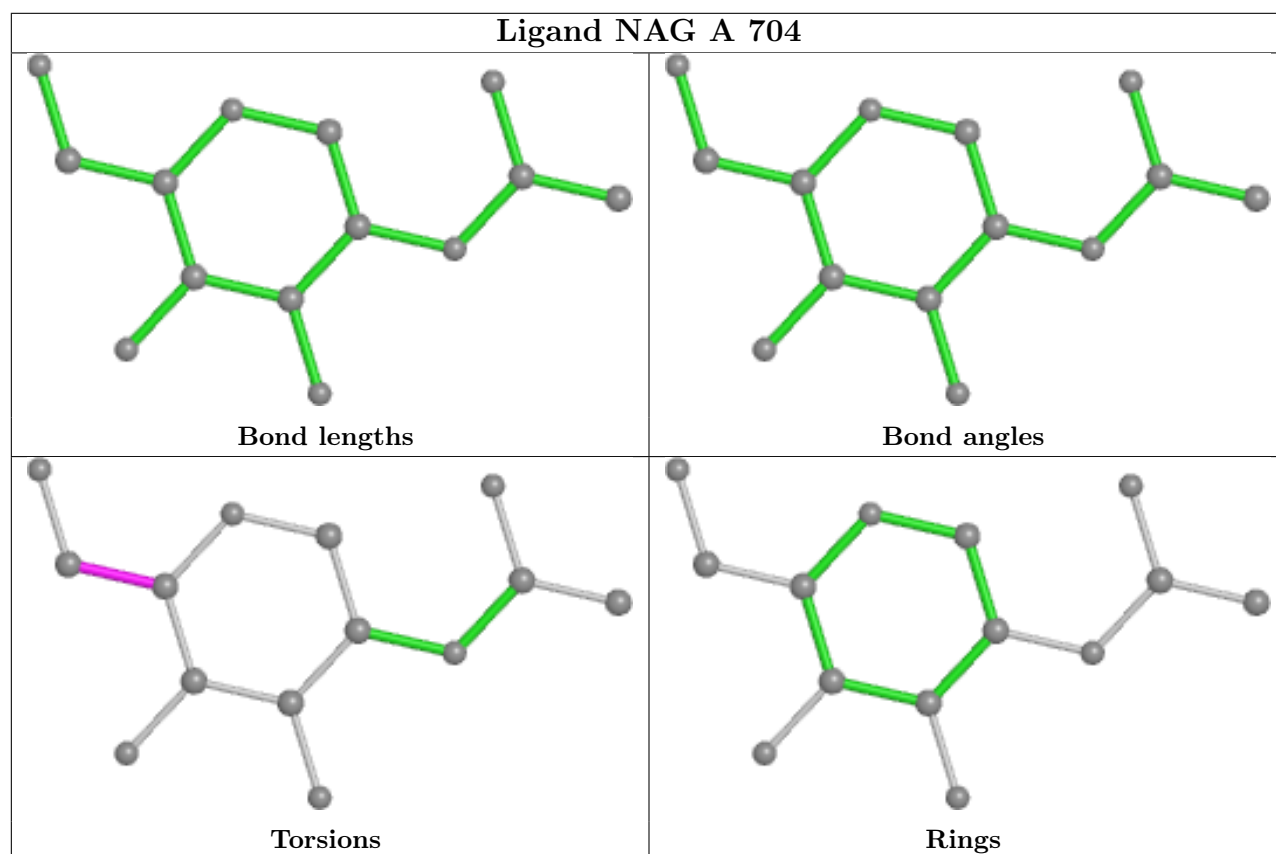
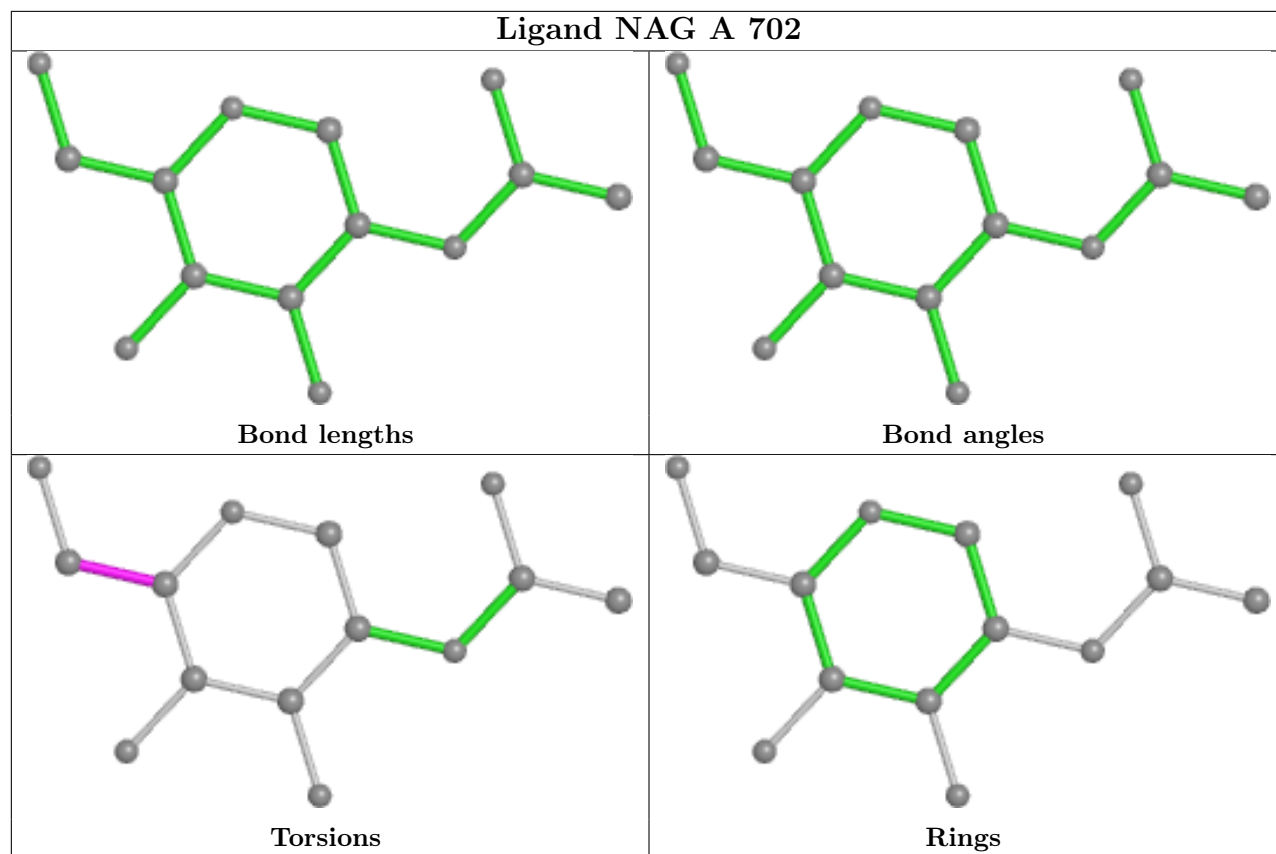
2 monomers are involved in 2 short contacts:

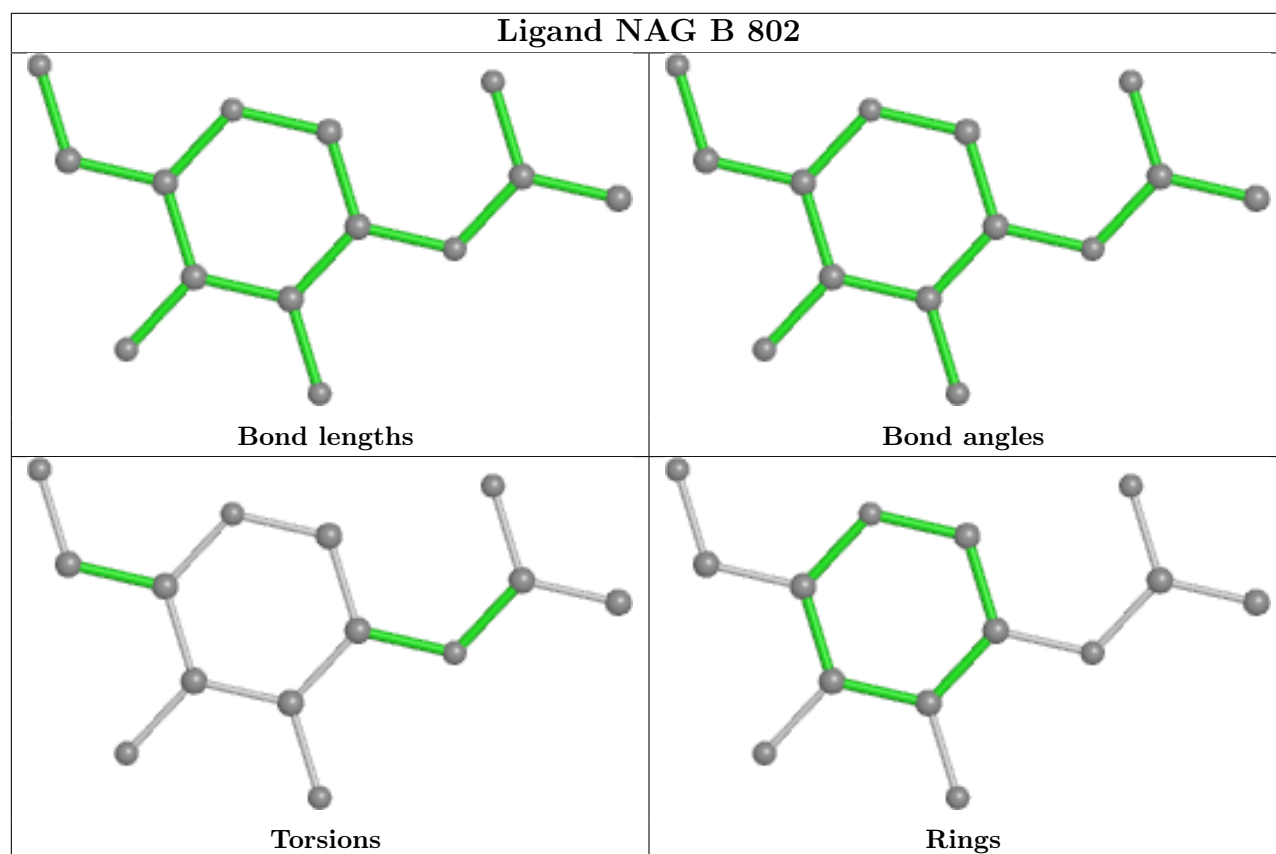
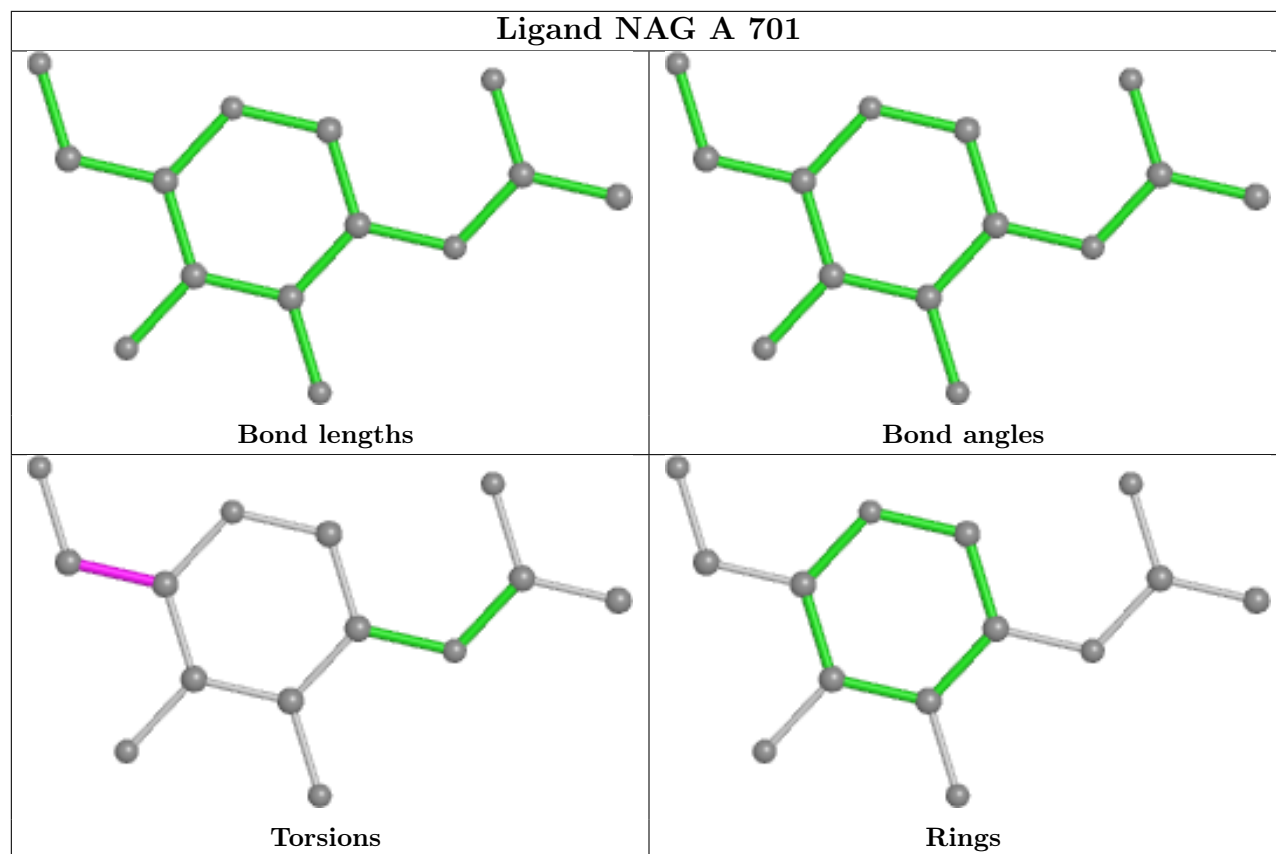
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	702	NAG	1	0
4	B	802	NAG	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier.

Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









## 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, 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	596/596 (100%)	-0.15	1 (0%) 95 90	60, 98, 139, 213	0
1	B	596/596 (100%)	-0.06	5 (0%) 86 72	60, 115, 194, 255	0
2	C	195/223 (87%)	-0.10	5 (2%) 56 33	56, 94, 163, 243	0
2	D	195/223 (87%)	0.61	31 (15%) 1 1	103, 166, 265, 304	0
All	All	1582/1638 (96%)	-0.02	42 (2%) 54 31	56, 108, 202, 304	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	77	VAL	8.1
2	D	197	PHE	5.7
2	D	45	ALA	5.2
2	D	20	PHE	4.5
2	D	195	LEU	4.4
2	D	74	PHE	4.4
2	D	38	LYS	4.3
2	D	204	ALA	4.2
2	D	114	CYS	4.0
1	B	472	GLN	3.8
2	D	44	VAL	3.8
2	D	196	SER	3.7
2	D	79	ALA	3.5
2	D	138	PHE	3.2
1	B	87	GLU	3.1
2	D	66	PRO	3.1
2	D	17	LEU	3.0
2	D	67	THR	2.9
2	D	116	ILE	2.8
2	C	52	ASN	2.6
2	D	203	PRO	2.6

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Mol	Chain	Res	Type	RSRZ
2	C	51	TYR	2.5
2	D	22	GLU	2.5
2	D	19	PRO	2.5
2	D	59	PHE	2.5
2	D	200	LEU	2.5
2	C	47	TYR	2.4
2	D	50	LEU	2.4
2	D	163	ASN	2.4
2	C	64	VAL	2.3
1	B	136	ASP	2.3
2	D	15	THR	2.3
2	D	161	PRO	2.3
2	D	47	TYR	2.2
2	C	203	PRO	2.2
2	D	209	PRO	2.2
1	B	614	ALA	2.2
1	B	135	PRO	2.1
2	D	173	PRO	2.1
1	A	19	SER	2.1
2	D	73	CYS	2.1
2	D	71	ASP	2.1

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

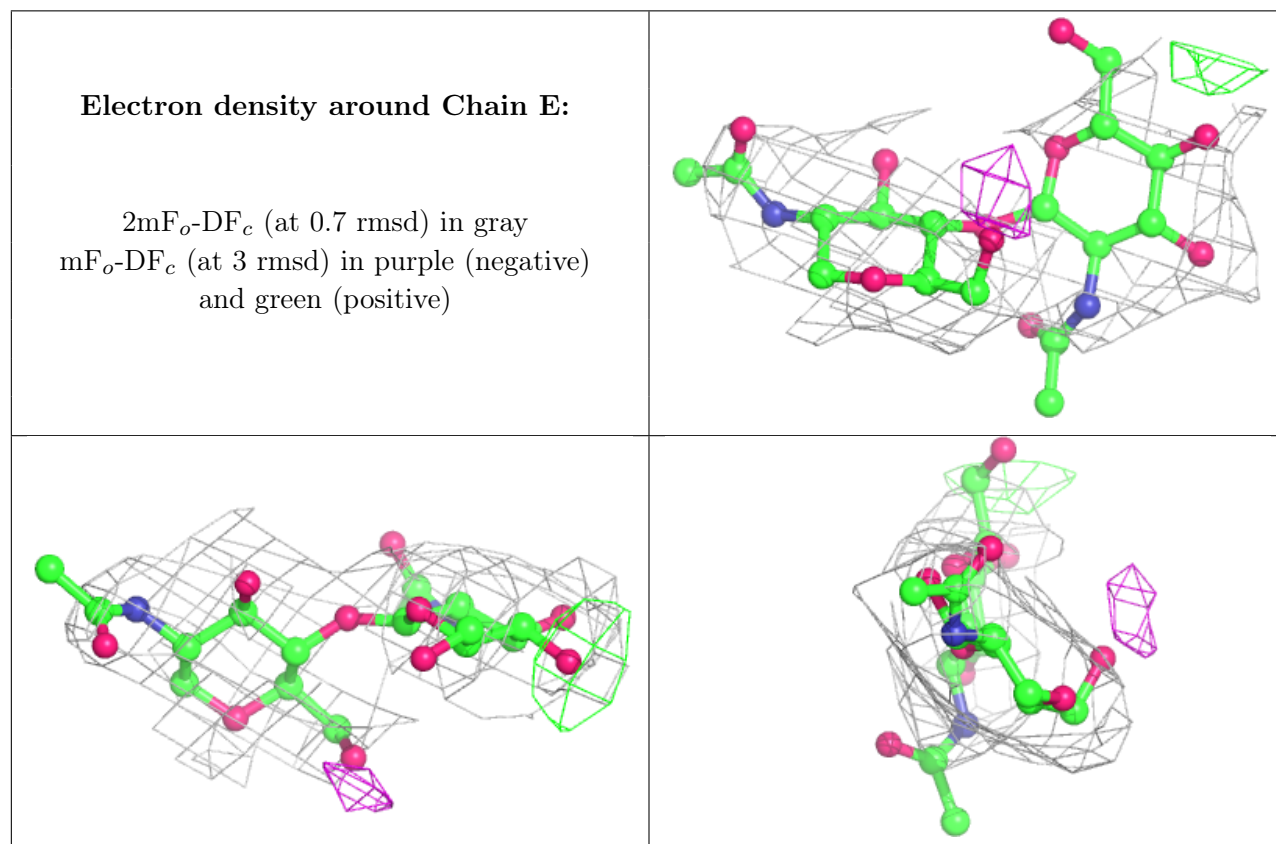
There are no non-standard protein/DNA/RNA residues in this entry.

## 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.

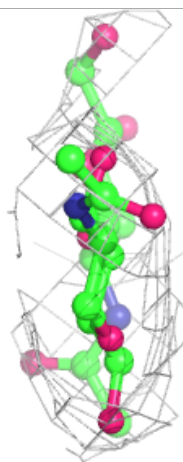
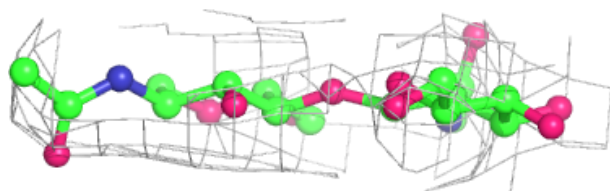
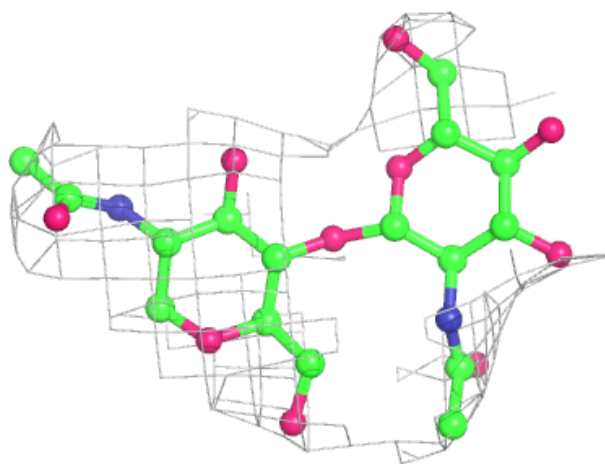
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	NAG	G	2	14/15	0.76	0.23	241,250,256,257	0
3	NAG	G	1	14/15	0.77	0.19	173,186,200,206	0
3	NAG	F	2	14/15	0.78	0.15	171,194,207,213	0
3	NAG	E	2	14/15	0.80	0.28	149,163,175,177	0
3	NAG	E	1	14/15	0.91	0.20	121,129,142,144	0
3	NAG	F	1	14/15	0.93	0.18	122,138,147,158	0

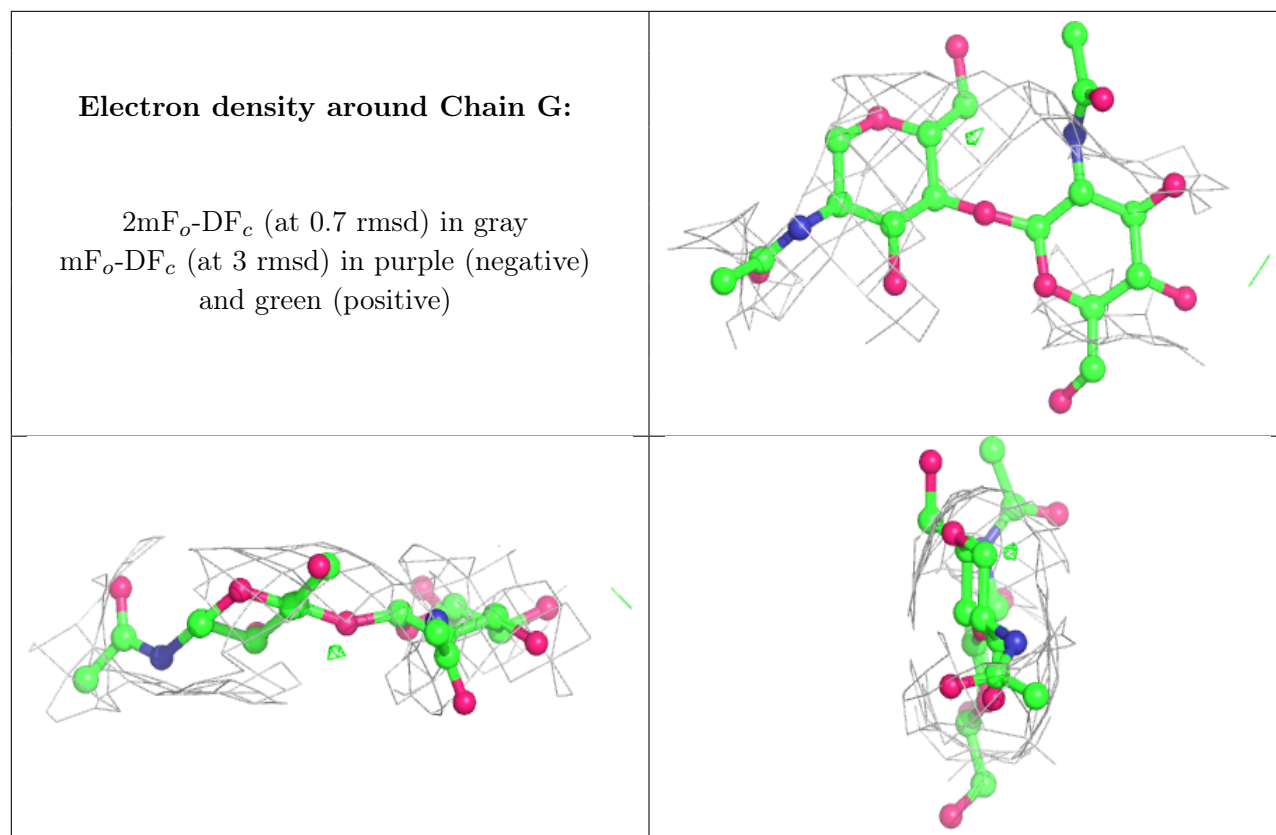
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



**Electron density around Chain F:**

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





## 6.4 Ligands [i](#)

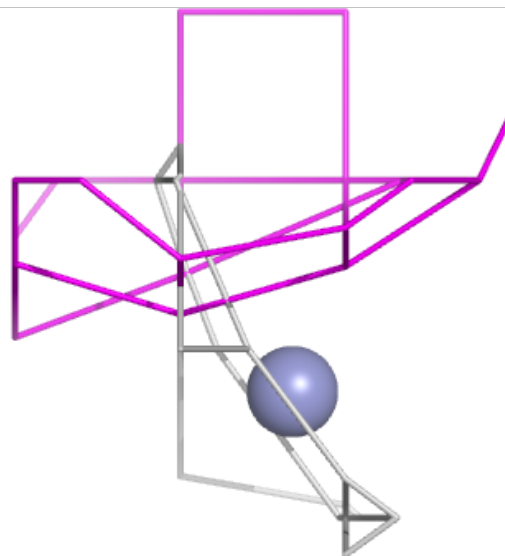
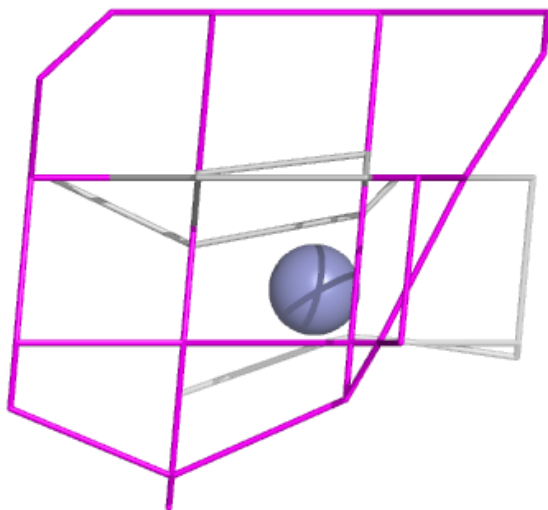
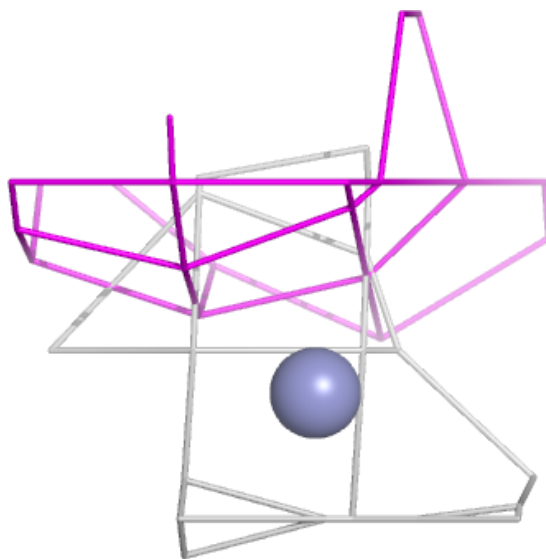
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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
5	ZN	A	703	1/1	0.26	0.61	298,298,298,298	0
4	NAG	A	702	14/15	0.58	0.28	160,182,192,196	0
4	NAG	B	802	14/15	0.69	0.26	122,135,148,150	0
4	NAG	C	301	14/15	0.75	0.21	112,122,129,130	0
4	NAG	A	701	14/15	0.81	0.12	146,154,161,165	0
4	NAG	A	704	14/15	0.83	0.18	134,140,148,149	0
5	ZN	B	801	1/1	0.87	0.40	140,140,140,140	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

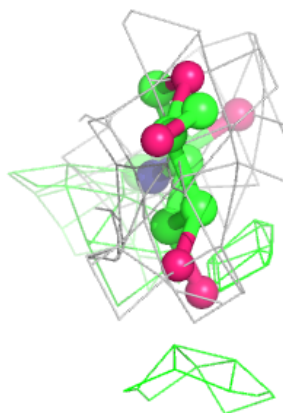
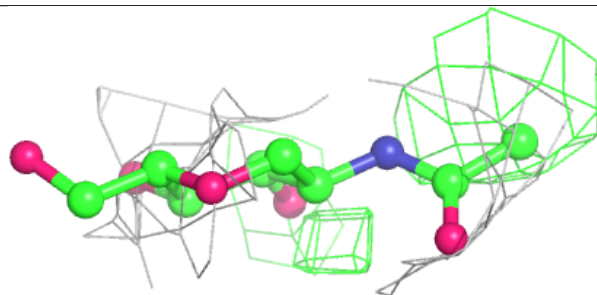
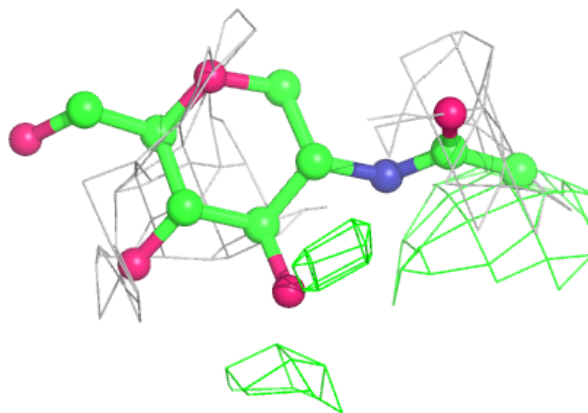
**Electron density around ZN A 703:**

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

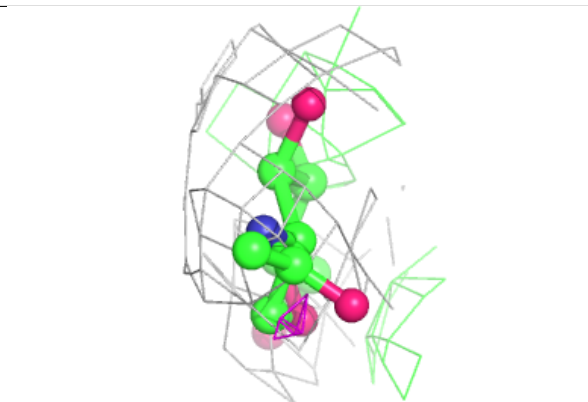
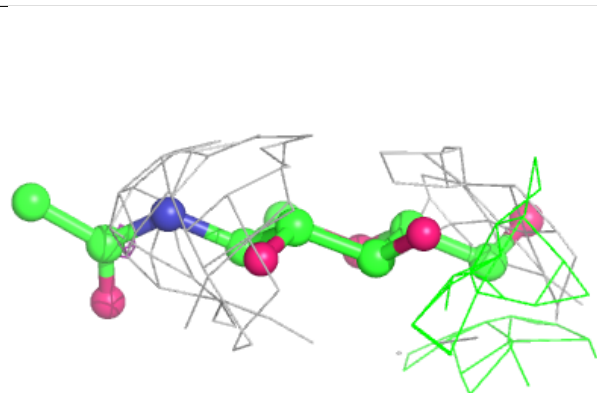
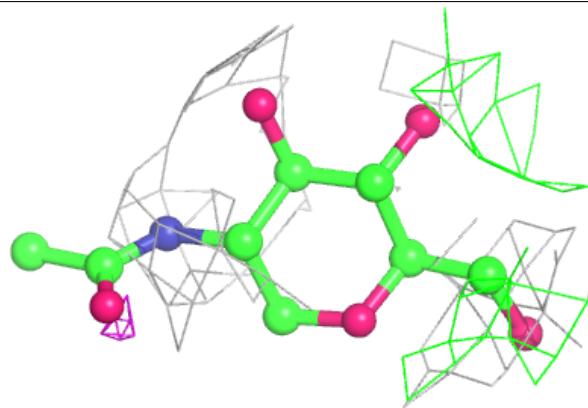


**Electron density around NAG A 702:**

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

**Electron density around NAG B 802:**

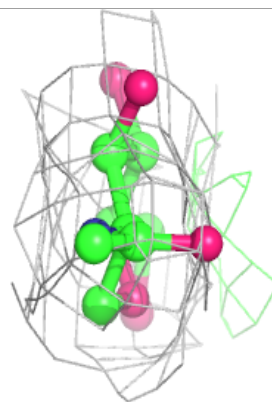
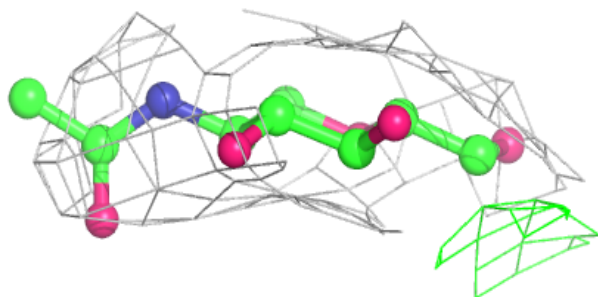
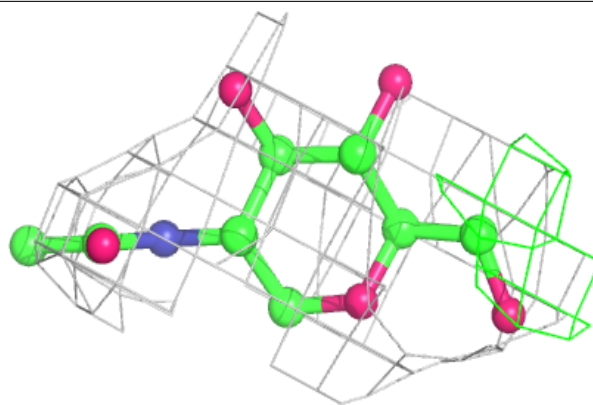
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



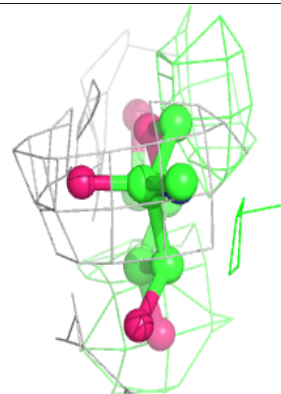
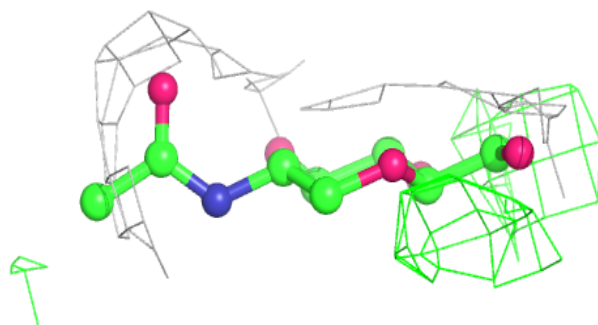
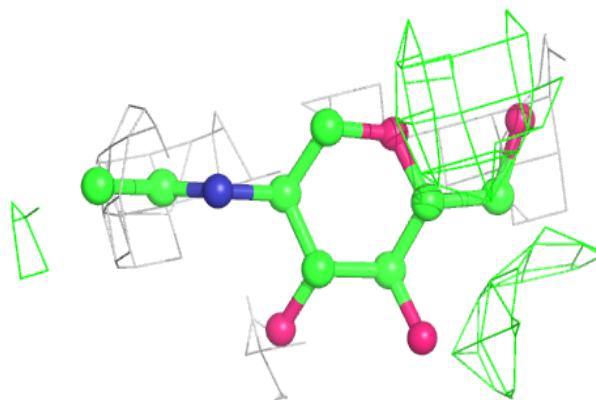


**Electron density around NAG C 301:**

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

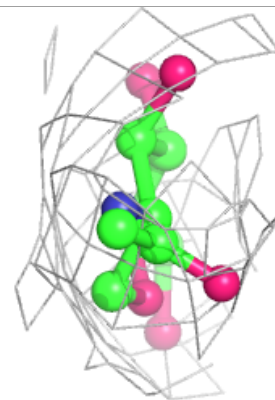
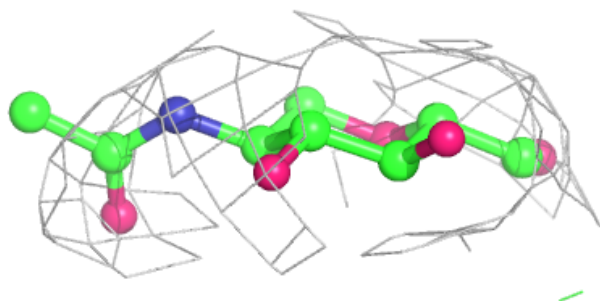
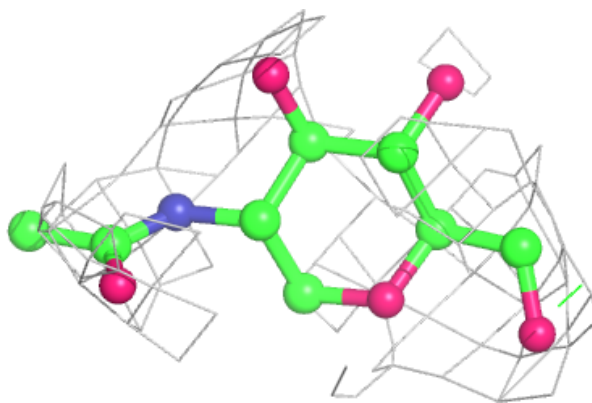
**Electron density around NAG A 701:**

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



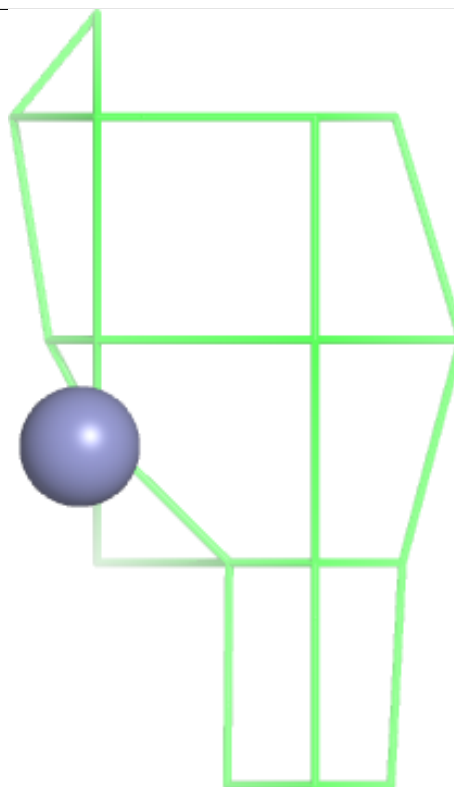
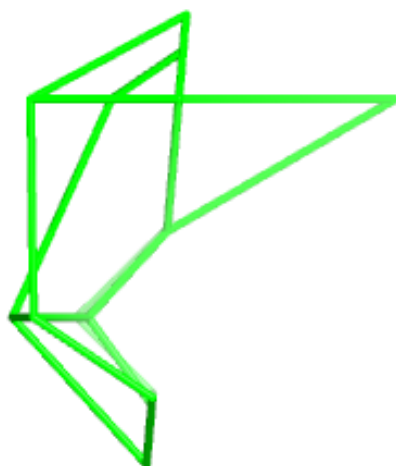
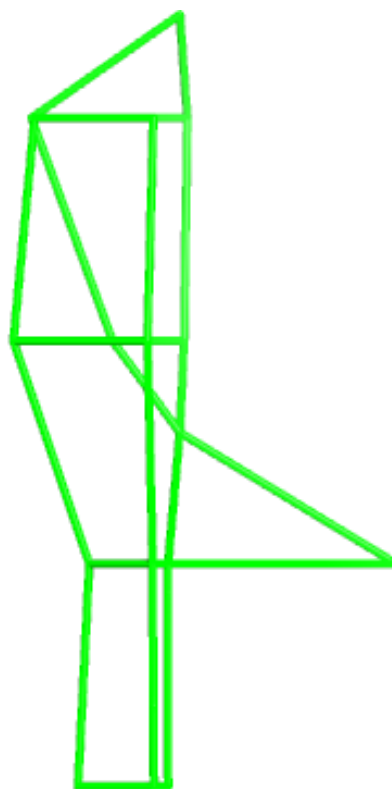
**Electron density around NAG A 704:**

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



**Electron density around ZN B 801:**

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



## 6.5 Other polymers [i](#)

There are no such residues in this entry.