



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 6, 2020 – 08:47 AM BST

PDB ID : 1V3B  
Title : Structure of the hemagglutinin-neuraminidase from human parainfluenza virus type III  
Authors : Lawrence, M.C.; Borg, N.A.; Streltsov, V.A.; Pilling, P.A.; Epa, V.C.; Varghese, J.N.; McKimm-Breschkin, J.L.; Colman, P.M.  
Deposited on : 2003-10-30  
Resolution : 2.00 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	:	2.13.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13.1

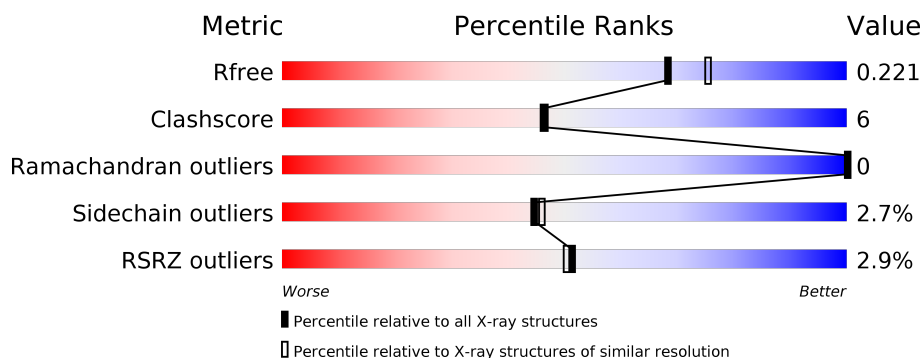
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	431	<div> <div>3%</div> <div> <div></div> <div>86%</div> <div>13%</div> </div> <div></div> </div>
1	B	431	<div> <div>3%</div> <div> <div></div> <div>85%</div> <div>14%</div> </div> <div></div> </div>
2	C	2	<div> <div></div> <div> <div>50%</div> <div>50%</div> </div> <div></div> </div>
2	D	2	<div> <div></div> <div>100%</div> <div></div> </div>

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 7781 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-neuraminidase glycoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	431	Total	C	N	O	S	0	2	0
			3391	2149	581	641	20			
1	B	431	Total	C	N	O	S	0	2	0
			3391	2149	581	641	20			

- Molecule 2 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
2	C	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	D	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 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
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Ca	0	0
			1	1		
4	A	1	Total	Ca	0	0
			1	1		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	O	S	0	0
			5	4	1		
5	B	1	Total	O	S	0	0
			5	4	1		

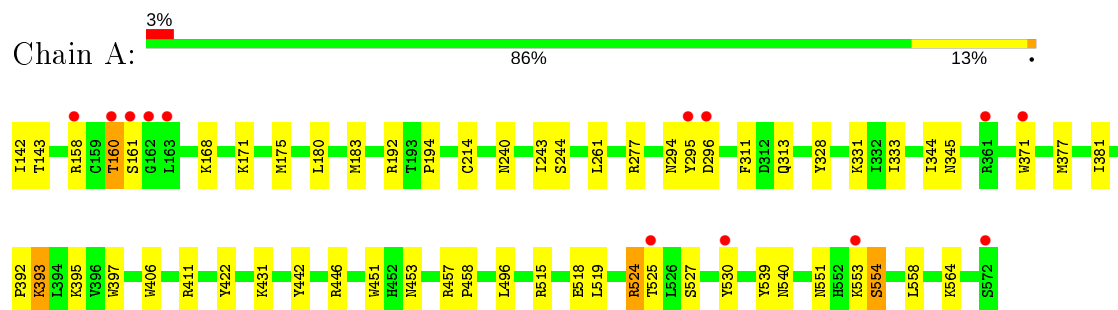
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	425	Total	O	0	0
			425	425		
6	B	464	Total	O	0	0
			464	464		

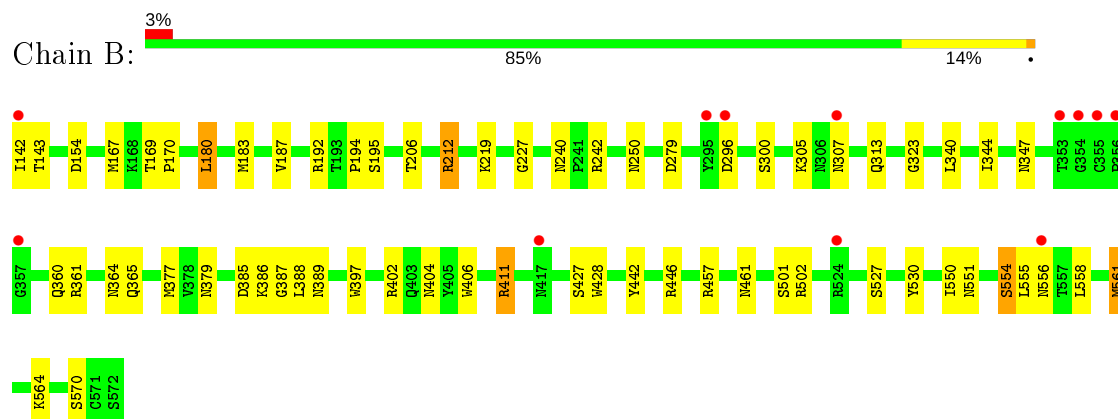
### 3 Residue-property plots [i](#)

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

- Molecule 1: hemagglutinin-neuraminidase glycoprotein



- Molecule 1: hemagglutinin-neuraminidase glycoprotein



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



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



MAG1  
MAG2

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 6 <sub>2</sub> 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	219.14Å 219.14Å 109.38Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	26.53 – 2.00 26.53 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (26.53-2.00) 100.0 (26.53-2.00)	Depositor EDS
$R_{merge}$	0.17	Depositor
$R_{sym}$	0.17	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.39 (at 1.99Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.182 , 0.215 0.187 , 0.221	Depositor DCC
$R_{free}$ test set	5234 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	29.2	Xtriage
Anisotropy	0.139	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 60.3	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	0.96	EDS
Total number of atoms	7781	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.07% 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: CA, NAG, SO4

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	0.82	1/3482 (0.0%)	0.91	5/4748 (0.1%)
1	B	0.81	1/3482 (0.0%)	0.91	2/4748 (0.0%)
All	All	0.81	2/6964 (0.0%)	0.91	7/9496 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	422	TYR	CE2-CZ	5.79	1.46	1.38
1	B	411	ARG	CG-CD	5.18	1.65	1.51

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	515	ARG	NE-CZ-NH2	-8.52	116.04	120.30
1	A	515	ARG	NE-CZ-NH1	6.01	123.30	120.30
1	B	411	ARG	NE-CZ-NH2	-5.13	117.74	120.30
1	B	279	ASP	CB-CG-OD1	5.10	122.89	118.30
1	A	411	ARG	NE-CZ-NH2	-5.10	117.75	120.30
1	A	519	LEU	CA-CB-CG	5.05	126.91	115.30
1	A	515	ARG	CB-CG-CD	-5.00	98.59	111.60

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	3391	0	3354	43	0
1	B	3391	0	3353	46	1
2	C	28	0	25	1	0
2	D	28	0	25	0	0
3	A	14	0	13	0	0
3	B	28	0	26	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	B	10	0	0	0	0
6	A	425	0	0	7	0
6	B	464	0	0	10	1
All	All	7781	0	6796	86	2

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 6.

All (86) 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:180:LEU:HD22	1:B:180:LEU:HD13	1.56	0.88
1:A:142:ILE:HG23	1:A:143:THR:HG23	1.59	0.83
1:B:195:SER:HB2	6:B:5438:HOH:O	1.82	0.78
1:A:277:ARG:HD2	1:A:371:TRP:CE2	2.20	0.76
1:A:524:ARG:H	1:A:524:ARG:HD2	1.50	0.76
1:A:392:PRO:C	1:A:393:LYS:HD2	2.07	0.74
1:B:365:GLN:HG2	6:B:5645:HOH:O	1.85	0.74
1:B:570:SER:HB2	6:B:5571:HOH:O	1.87	0.73
1:A:392:PRO:O	1:A:393:LYS:HD2	1.91	0.71
1:B:361:ARG:O	1:B:365:GLN:HG3	1.90	0.71
1:A:142:ILE:HA	6:A:3807:HOH:O	1.91	0.70
1:B:167:MET:HG2	6:B:5633:HOH:O	1.94	0.68
1:B:360:GLN:HE21	1:B:364:ASN:HD21	1.44	0.66
1:A:524:ARG:CD	1:A:524:ARG:H	2.09	0.65
1:A:277:ARG:HD2	1:A:371:TRP:CZ2	2.35	0.62
1:B:555:LEU:O	1:B:556:ASN:HB3	2.00	0.61
1:B:219:LYS:HD2	1:B:250:ASN:HD22	1.65	0.60
1:B:323:GLY:HA3	1:B:411:ARG:HB3	1.84	0.60
1:A:554:SER:OG	6:A:3516:HOH:O	2.16	0.60
1:A:168:LYS:HG3	6:A:3854:HOH:O	2.01	0.59
1:A:539:TYR:O	1:A:540:ASN:HB2	2.03	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:MET:CE	1:B:183:MET:HE3	2.34	0.57
1:B:340:LEU:HD13	1:B:344:ILE:HD12	1.85	0.57
1:A:527:SER:HB2	1:A:551:ASN:HB3	1.87	0.56
1:A:331:LYS:HE2	6:A:3726:HOH:O	2.06	0.56
1:B:360:GLN:HE21	1:B:364:ASN:ND2	2.03	0.56
1:A:525:THR:O	1:A:553:LYS:NZ	2.32	0.56
1:B:427:SER:HB3	6:B:5322:HOH:O	2.06	0.55
6:A:3516:HOH:O	1:B:554:SER:OG	2.19	0.53
6:A:3786:HOH:O	2:C:1:NAG:H81	2.09	0.53
1:A:539:TYR:O	1:A:540:ASN:CB	2.59	0.51
1:B:219:LYS:CD	1:B:250:ASN:HD22	2.22	0.51
1:B:386:LYS:HE3	1:B:387:GLY:O	2.10	0.51
1:A:160:THR:HG23	1:A:161:SER:H	1.75	0.51
1:A:158:ARG:HD2	6:A:3883:HOH:O	2.11	0.50
1:B:227:GLY:HA2	1:B:242:ARG:HG2	1.93	0.50
1:B:550:ILE:HG13	1:B:561:MET:HE3	1.94	0.50
1:A:377:MET:HG3	1:A:406:TRP:CZ2	2.47	0.50
1:A:451:TRP:CH2	1:A:453:ASN:HB2	2.48	0.49
1:B:570:SER:OG	6:B:5633:HOH:O	2.20	0.49
1:A:277:ARG:HD2	1:A:371:TRP:NE1	2.27	0.48
1:A:551:ASN:HA	1:A:558:LEU:HD23	1.95	0.48
1:B:340:LEU:CD1	1:B:344:ILE:HD12	2.42	0.48
1:B:502:ARG:HD2	6:B:5695:HOH:O	2.13	0.48
1:A:183:MET:CE	1:B:183:MET:CE	2.92	0.48
1:A:192:ARG:HH11	1:A:214:CYS:HB3	1.78	0.48
1:B:212:ARG:HG2	1:B:212:ARG:HH11	1.79	0.47
1:B:194:PRO:O	1:B:530[B]:TYR:CZ	2.68	0.47
1:A:183:MET:HE3	1:B:183:MET:HE3	1.95	0.47
1:A:328:TYR:HB3	1:A:333:ILE:HD12	1.97	0.46
1:B:195:SER:HB3	1:B:206:THR:HG22	1.96	0.46
1:B:555:LEU:O	1:B:556:ASN:CB	2.60	0.46
1:A:296:ASP:OD1	1:A:296:ASP:C	2.53	0.45
1:A:171:LYS:HE3	1:A:518:GLU:O	2.16	0.45
1:B:142:ILE:HG23	1:B:143:THR:HG23	1.98	0.45
1:B:169:THR:HA	1:B:170:PRO:C	2.36	0.45
1:B:313:GLN:OE1	1:B:344:ILE:HD11	2.16	0.45
1:A:431:LYS:NZ	1:A:453:ASN:OD1	2.48	0.45
1:B:402:ARG:NH1	6:B:5643:HOH:O	2.44	0.45
1:A:333:ILE:CG2	1:A:381:ILE:HG23	2.48	0.44
1:B:187:VAL:HG12	6:B:5251:HOH:O	2.18	0.44
1:B:461:ASN:ND2	6:B:5510:HOH:O	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:305:LYS:HB2	1:B:307:ASN:ND2	2.32	0.44
1:A:333:ILE:HG23	1:A:381:ILE:HG23	2.00	0.43
1:B:377:MET:HG3	1:B:406:TRP:CZ2	2.53	0.43
1:A:261:LEU:HD23	1:A:261:LEU:C	2.38	0.43
1:B:183:MET:HB3	1:B:183:MET:HE3	1.88	0.43
1:A:395:LYS:HB2	1:A:395:LYS:HE3	1.89	0.43
1:B:388:LEU:HD23	1:B:389:ASN:ND2	2.33	0.43
1:A:194:PRO:O	1:A:530[B]:TYR:CZ	2.72	0.42
1:A:175:MET:HE3	1:A:564:LYS:HB3	2.01	0.42
1:B:388:LEU:HD23	1:B:389:ASN:HD21	1.85	0.42
1:B:347:ASN:HA	1:B:364:ASN:HD21	1.84	0.42
1:B:379:ASN:HB2	1:B:404:ASN:ND2	2.35	0.41
1:B:527:SER:HB2	1:B:551:ASN:HB3	2.01	0.41
1:B:427:SER:OG	1:B:428:TRP:N	2.52	0.41
1:A:243:ILE:HG12	1:A:244:SER:N	2.35	0.41
1:B:397:TRP:CD2	1:B:442:TYR:HB3	2.56	0.41
1:A:458:PRO:HD3	1:A:496:LEU:O	2.21	0.41
1:A:524:ARG:N	1:A:524:ARG:HD2	2.28	0.41
1:A:311:PHE:HB3	1:A:313:GLN:O	2.21	0.41
1:A:344:ILE:HG22	1:A:345:ASN:N	2.36	0.41
1:B:194:PRO:O	1:B:530[B]:TYR:CE2	2.74	0.41
1:A:397:TRP:CD2	1:A:442:TYR:HB3	2.56	0.41
1:B:192:ARG:CD	1:B:558:LEU:HD12	2.52	0.40
1:A:393:LYS:N	1:A:393:LYS:HD2	2.34	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B:5383:HOH:O	6:B:5383:HOH:O[7_555]	2.05	0.15
1:B:296:ASP:OD2	1:B:296:ASP:OD2[10_665]	2.13	0.07

## 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	431/431 (100%)	410 (95%)	21 (5%)	0	100	100
1	B	431/431 (100%)	413 (96%)	18 (4%)	0	100	100
All	All	862/862 (100%)	823 (96%)	39 (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	392/390 (100%)	383 (98%)	9 (2%)	50	53
1	B	392/390 (100%)	380 (97%)	12 (3%)	40	40
All	All	784/780 (100%)	763 (97%)	21 (3%)	44	46

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	160	THR
1	A	240	ASN
1	A	294	ASN
1	A	295	TYR
1	A	393	LYS
1	A	446	ARG
1	A	457	ARG
1	A	524	ARG
1	A	554	SER
1	B	154	ASP
1	B	180	LEU
1	B	212	ARG
1	B	240	ASN
1	B	300	SER
1	B	385	ASP

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Mol	Chain	Res	Type
1	B	446	ARG
1	B	457	ARG
1	B	501	SER
1	B	554	SER
1	B	561	MET
1	B	564	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	240	ASN
1	A	262	ASN
1	A	294	ASN
1	A	347	ASN
1	A	504	ASN
1	A	556	ASN
1	B	240	ASN
1	B	250	ASN
1	B	262	ASN
1	B	307	ASN
1	B	347	ASN
1	B	364	ASN
1	B	389	ASN
1	B	499	GLN
1	B	504	ASN
1	B	556	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](#)

4 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	C	1	1,2	14,14,15	0.66	0	17,19,21	1.15	3 (17%)
2	NAG	C	2	2	14,14,15	0.57	0	17,19,21	0.80	1 (5%)
2	NAG	D	1	1,2	14,14,15	0.77	0	17,19,21	1.02	2 (11%)
2	NAG	D	2	2	14,14,15	0.91	1 (7%)	17,19,21	0.75	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
2	NAG	C	1	1,2	-	4/6/23/26	0/1/1/1
2	NAG	C	2	2	-	3/6/23/26	0/1/1/1
2	NAG	D	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	D	2	2	-	3/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	2	NAG	C1-C2	2.51	1.56	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	NAG	C2-N2-C7	-3.09	118.50	122.90
2	D	1	NAG	C2-N2-C7	-2.69	119.07	122.90
2	C	1	NAG	C4-C3-C2	-2.39	107.51	111.02
2	D	1	NAG	O5-C1-C2	-2.09	107.99	111.29
2	C	1	NAG	O5-C1-C2	-2.06	108.03	111.29
2	C	2	NAG	C2-N2-C7	-2.04	119.99	122.90

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	2	NAG	C3-C2-N2-C7
2	C	1	NAG	O5-C5-C6-O6
2	C	1	NAG	C8-C7-N2-C2
2	C	1	NAG	C4-C5-C6-O6
2	D	2	NAG	C8-C7-N2-C2
2	D	2	NAG	O7-C7-N2-C2
2	C	2	NAG	C8-C7-N2-C2
2	C	1	NAG	O7-C7-N2-C2
2	D	1	NAG	O5-C5-C6-O6
2	C	2	NAG	O7-C7-N2-C2
2	D	1	NAG	C4-C5-C6-O6
2	C	2	NAG	O5-C5-C6-O6

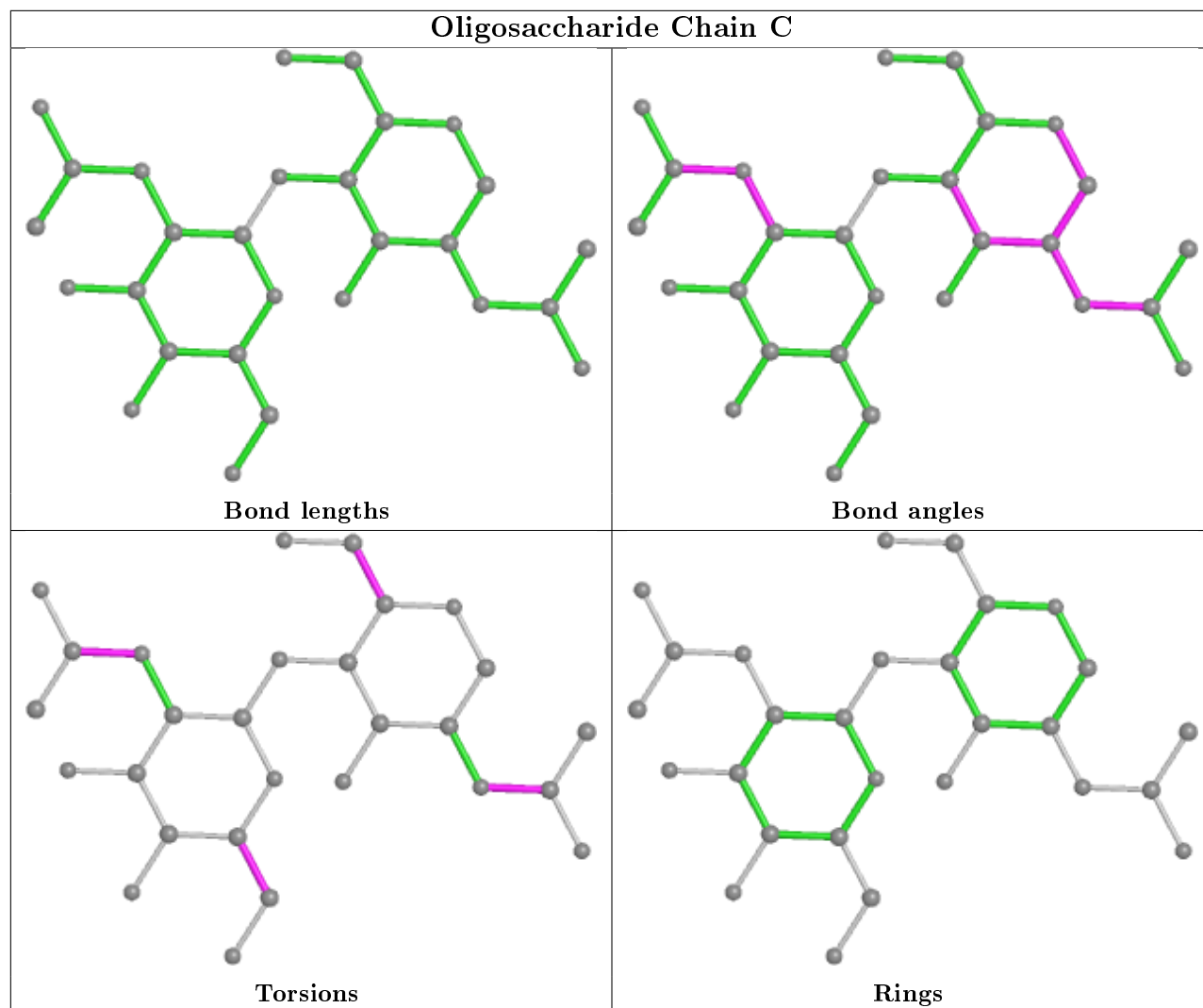
There are no ring outliers.

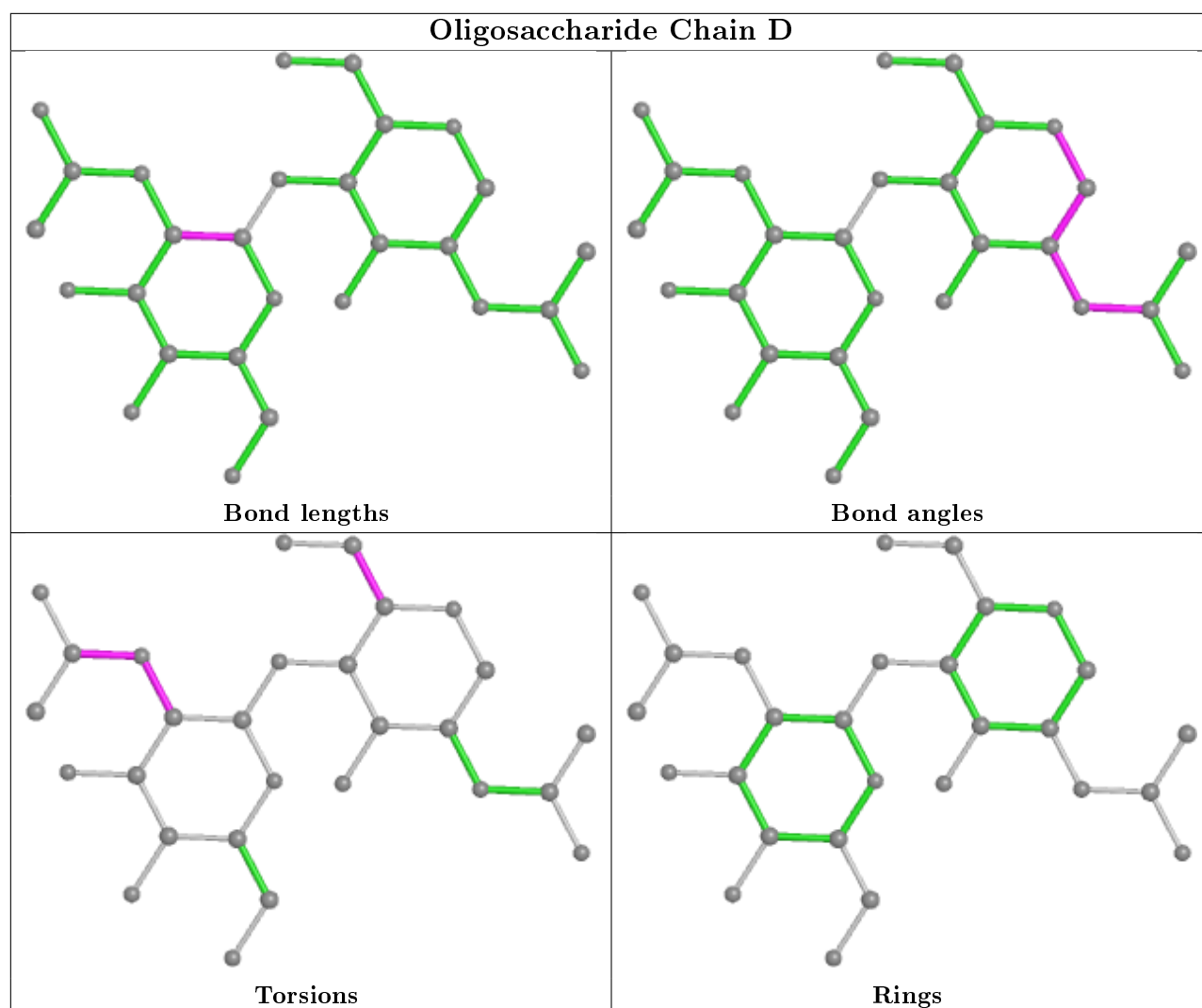
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	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](#)

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	$\# Z  > 2$
3	NAG	A	3081	1	14,14,15	0.60	0	17,19,21	0.71	0
5	SO4	B	4002	-	4,4,4	0.59	0	6,6,6	0.34	0
5	SO4	B	4001	-	4,4,4	0.27	0	6,6,6	0.25	0
3	NAG	B	3081	1	14,14,15	0.63	0	17,19,21	0.64	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	NAG	B	5231	1	14,14,15	0.90	1 (7%)	17,19,21	0.87	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	A	3081	1	-	3/6/23/26	0/1/1/1
3	NAG	B	5231	1	-	2/6/23/26	0/1/1/1
3	NAG	B	3081	1	-	0/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	5231	NAG	C1-C2	2.41	1.55	1.52

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	5231	NAG	C8-C7-N2-C2
3	B	5231	NAG	O7-C7-N2-C2
3	A	3081	NAG	C8-C7-N2-C2
3	A	3081	NAG	O7-C7-N2-C2
3	A	3081	NAG	C4-C5-C6-O6

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	431/431 (100%)	-0.23	13 (3%)	50	49	20, 32, 57, 82	0
1	B	431/431 (100%)	-0.30	12 (2%)	53	51	21, 31, 54, 71	0
All	All	862/862 (100%)	-0.27	25 (2%)	51	50	20, 32, 56, 82	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	572	SER	7.2
1	A	161	SER	4.4
1	A	371	TRP	3.8
1	A	160	THR	3.7
1	A	295	TYR	3.6
1	A	163	LEU	3.4
1	A	158	ARG	3.3
1	B	142	ILE	3.1
1	B	357	GLY	3.1
1	B	417	ASN	2.8
1	B	295	TYR	2.8
1	B	356	PRO	2.7
1	A	525	THR	2.6
1	B	354	GLY	2.6
1	B	355	CYS	2.5
1	A	553	LYS	2.5
1	B	556	ASN	2.5
1	A	361	ARG	2.5
1	B	296	ASP	2.3
1	B	353	THR	2.3
1	A	530[A]	TYR	2.2
1	A	162	GLY	2.2
1	A	296	ASP	2.2
1	B	307	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	524	ARG	2.2

## 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](#)

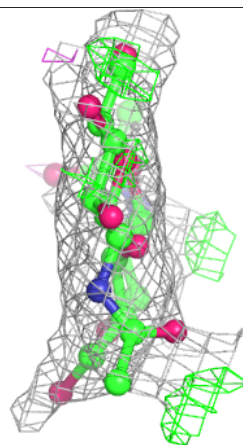
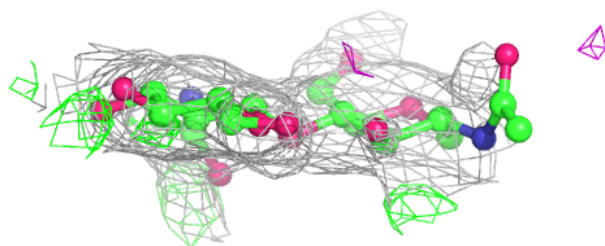
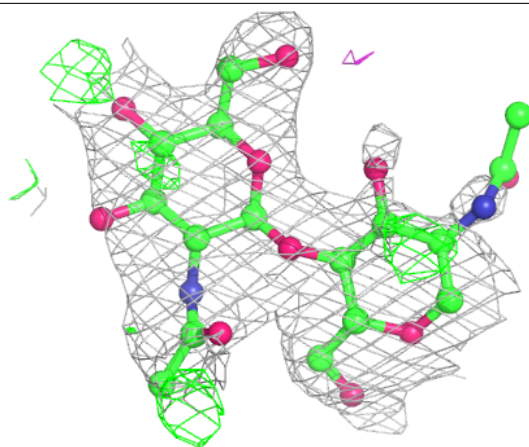
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

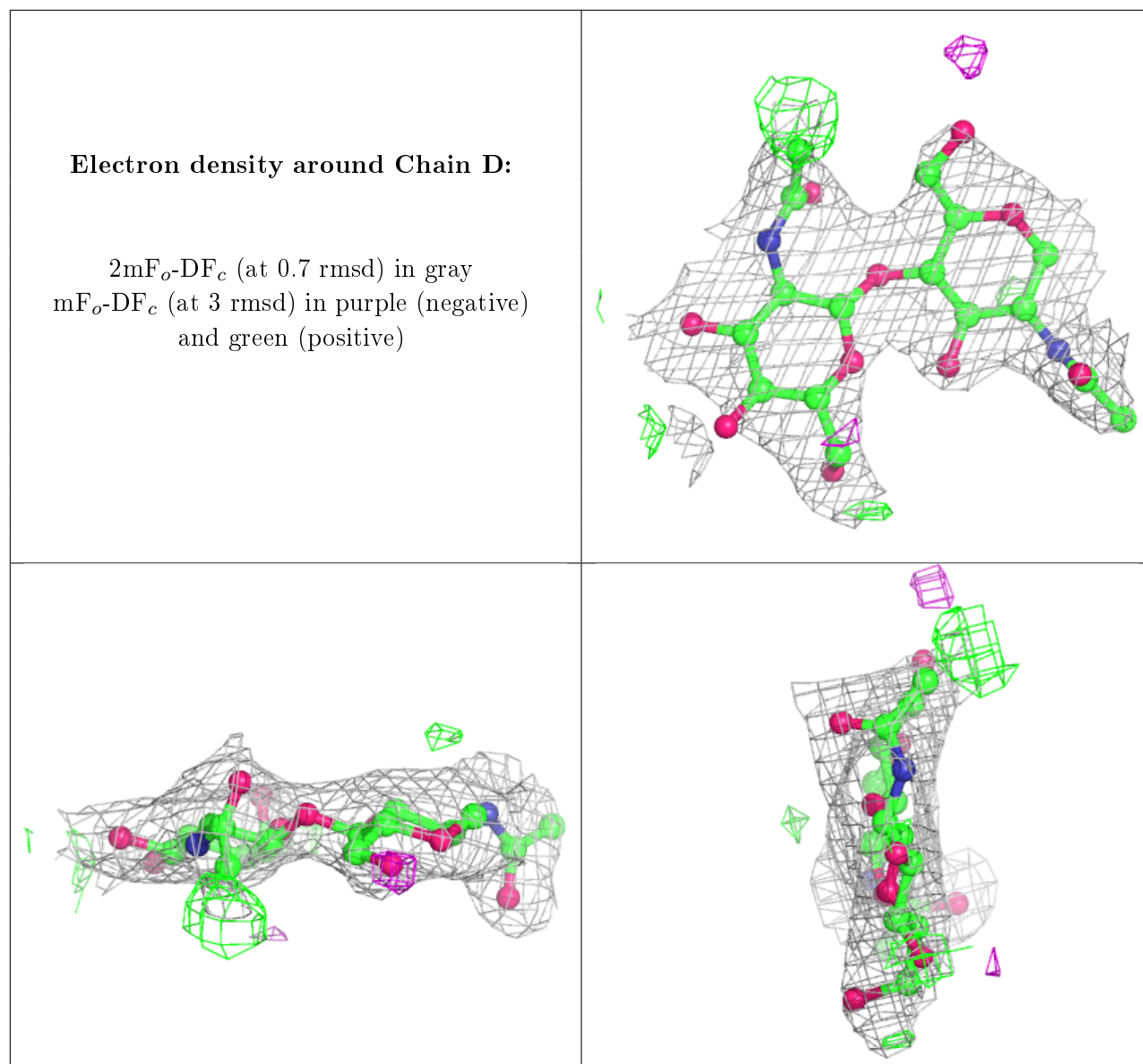
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	NAG	C	1	14/15	0.78	0.23	72,82,96,96	0
2	NAG	D	2	14/15	0.85	0.21	68,78,82,85	0
2	NAG	C	2	14/15	0.86	0.18	71,78,81,87	0
2	NAG	D	1	14/15	0.91	0.15	61,69,78,79	0

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 C:**

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





## 6.4 Ligands ⓘ

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	A	3081	14/15	0.80	0.38	71,77,81,85	0
3	NAG	B	5231	14/15	0.86	0.25	52,66,72,75	0
5	SO4	B	4002	5/5	0.88	0.21	50,59,77,79	0
3	NAG	B	3081	14/15	0.92	0.22	48,54,58,58	0
5	SO4	B	4001	5/5	0.96	0.17	77,78,85,86	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	CA	B	2001	1/1	0.98	0.04	30,30,30,30	0
4	CA	A	1001	1/1	0.99	0.06	35,35,35,35	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.