



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 22, 2025 – 12:59 AM EDT

PDB ID : 8W0V / pdb_00008w0v
Title : Crystal structure of broadly neutralizing antibody hcab55 in complex with Hepatitis C virus envelope glycoprotein E2 ectodomain
Authors : Flyak, A.I.; Wilcox, X.E.
Deposited on : 2024-02-14
Resolution : 2.59 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0rc1
EDS	:	FAILED
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.42

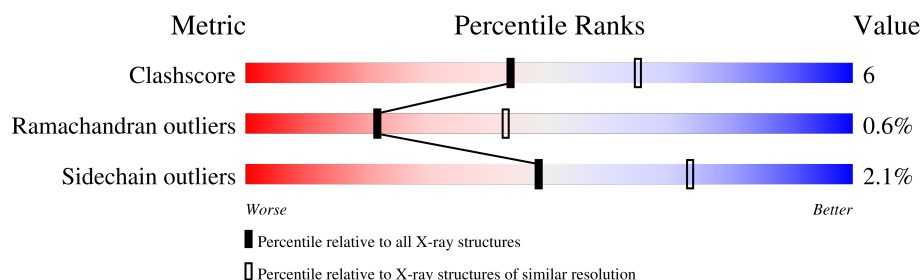
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	180529	4181 (2.60-2.60)
Ramachandran outliers	177936	4129 (2.60-2.60)
Sidechain outliers	177891	4129 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 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 $\leq 5\%$.

Note EDS failed to run properly.

Mol	Chain	Length	Quality of chain
1	C	262	
2	H	237	
3	L	215	
4	A	4	
5	B	2	
5	D	2	
6	E	4	

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 5203 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 Envelope glycoprotein E2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	219	Total	C	N	O	S	0	0	0
			1716	1090	303	304	19			

- Molecule 2 is a protein called hcab55 Fab Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	224	Total	C	N	O	S	0	0	0
			1677	1056	287	327	7			

- Molecule 3 is a protein called hcab55 Fab Light Chain.

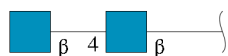
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	213	Total	C	N	O	S	0	0	0
			1640	1030	279	327	4			

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-6)-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
4	A	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 5 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



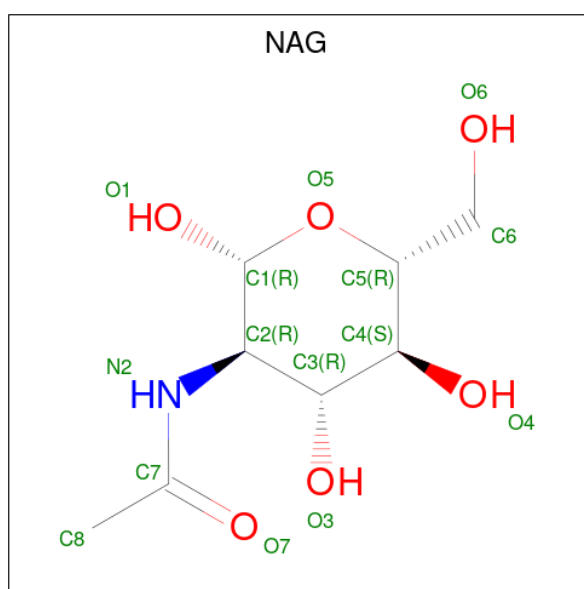
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	B	2	Total	C	N	O	0	0	0
			28	16	2	10			
5	D	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 6 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-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
6	E	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



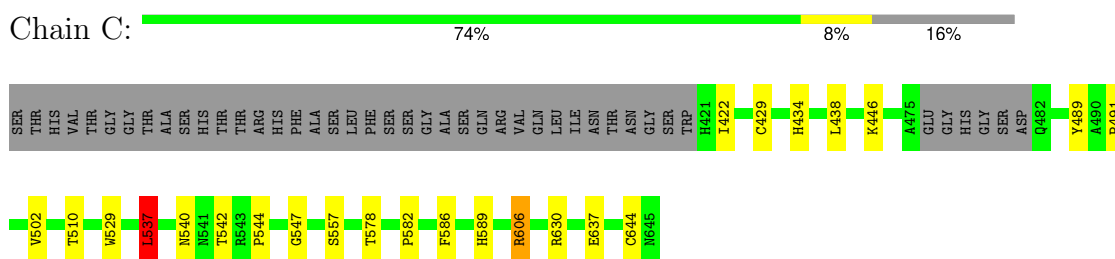
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	C	1	Total	C	N	O	0	0
			14	8	1	5		

3 Residue-property plots

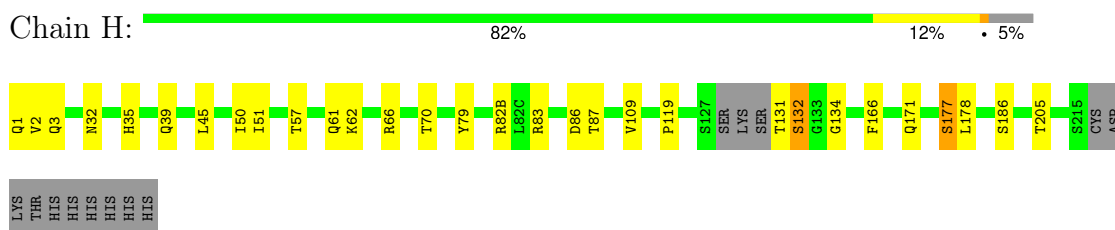
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS failed to run properly.

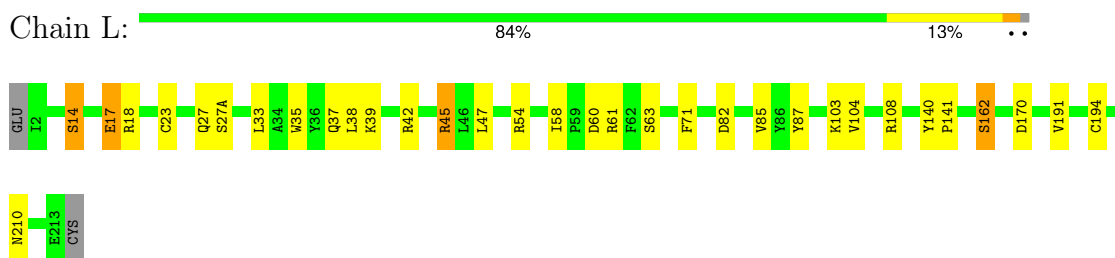
- Molecule 1: Envelope glycoprotein E2



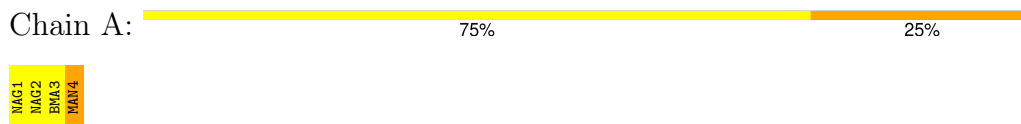
- Molecule 2: hcab55 Fab Heavy Chain



- Molecule 3: hcab55 Fab Light Chain



- Molecule 4: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose




- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain B:  50% 50%

MAG1
MAG2

- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:  100%

MAG1
MAG2

- Molecule 6: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:  25% 50% 25%

MAG1
MAG2
BMA3
MAN4

4 Data and refinement statistics

EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.02Å 88.84Å 175.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.67 – 2.59	Depositor
% Data completeness (in resolution range)	98.6 (45.67-2.59)	Depositor
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 2.58Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.202 , 0.257	Depositor
Wilson B-factor (Å ²)	51.4	Xtriage
Anisotropy	0.448	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5203	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.02% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, BMA, MAN

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	C	0.46	0/1778	0.65	1/2435 (0.0%)
2	H	0.49	0/1716	0.66	0/2337
3	L	0.50	1/1677 (0.1%)	0.68	0/2278
All	All	0.48	1/5171 (0.0%)	0.66	1/7050 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	L	194	CYS	CB-SG	-6.11	1.71	1.82

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	537	LEU	CA-CB-CG	-5.03	103.74	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1716	0	1591	17	0
2	H	1677	0	1655	25	0
3	L	1640	0	1600	26	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	50	0	43	1	0
5	B	28	0	25	0	0
5	D	28	0	25	0	0
6	E	50	0	43	1	0
7	C	14	0	13	0	0
All	All	5203	0	4995	61	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 6.

All (61) 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)
2:H:70:THR:HG22	2:H:79:TYR:HB2	1.55	0.89
3:L:108:ARG:NH1	3:L:170:ASP:O	2.09	0.85
1:C:446:LYS:NZ	2:H:32:ASN:HD21	1.85	0.74
3:L:17:GLU:HG3	3:L:18:ARG:H	1.52	0.74
2:H:66:ARG:NH2	2:H:86:ASP:OD2	2.22	0.72
2:H:39:GLN:NE2	3:L:38:LEU:HD21	2.07	0.70
1:C:446:LYS:HZ2	2:H:32:ASN:HD21	1.41	0.66
3:L:61:ARG:HD2	3:L:82:ASP:OD2	1.98	0.64
1:C:429:CYS:HB3	1:C:438:LEU:HD21	1.80	0.63
3:L:54:ARG:HE	3:L:60:ASP:HA	1.63	0.62
3:L:38:LEU:HD23	3:L:87:TYR:CE1	2.35	0.61
2:H:39:GLN:HE22	3:L:38:LEU:HD21	1.65	0.60
3:L:39:LYS:HD3	3:L:42:ARG:NH1	2.15	0.60
3:L:108:ARG:HH12	3:L:170:ASP:HB2	1.67	0.59
1:C:446:LYS:NZ	2:H:32:ASN:ND2	2.52	0.57
2:H:1:GLN:OE1	2:H:1:GLN:N	2.38	0.57
2:H:131:THR:HG22	2:H:132:SER:H	1.70	0.57
2:H:61:GLN:O	2:H:62:LYS:HB3	2.07	0.55
3:L:47:LEU:HA	3:L:58:ILE:HG13	1.89	0.54
2:H:1:GLN:N	2:H:1:GLN:CD	2.61	0.54
3:L:54:ARG:HG3	3:L:54:ARG:HH11	1.72	0.54
2:H:35:HIS:CE1	2:H:50:ILE:HD12	2.42	0.54
2:H:171:GLN:NE2	2:H:177:SER:HB2	2.25	0.52
3:L:54:ARG:HG3	3:L:54:ARG:NH1	2.25	0.52
1:C:557:SER:HB2	6:E:1:NAG:H82	1.91	0.52
2:H:119:PRO:HD2	2:H:205:THR:HG21	1.93	0.51
1:C:446:LYS:HZ3	2:H:32:ASN:ND2	2.09	0.51
2:H:66:ARG:HG2	2:H:82(B):ARG:NH2	2.27	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:37:GLN:OE1	3:L:45:ARG:HD3	2.13	0.49
3:L:17:GLU:CG	3:L:18:ARG:H	2.23	0.48
3:L:27:GLN:HG2	3:L:27(A):SER:H	1.78	0.48
3:L:14:SER:O	3:L:17:GLU:HB3	2.14	0.48
1:C:542:THR:O	1:C:547:GLY:HA3	2.14	0.48
2:H:131:THR:HG21	2:H:134:GLY:HA3	1.97	0.46
3:L:85:VAL:HG22	3:L:103:LYS:HG2	1.98	0.45
2:H:1:GLN:CD	2:H:1:GLN:H3	2.19	0.45
1:C:502:VAL:HG21	1:C:537:LEU:HD13	1.99	0.45
3:L:27:GLN:HG2	3:L:27(A):SER:N	2.31	0.45
3:L:140:TYR:CD1	3:L:141:PRO:HA	2.52	0.44
3:L:103:LYS:HD3	3:L:104:VAL:N	2.31	0.44
2:H:87:THR:HA	2:H:109:VAL:O	2.17	0.44
1:C:630:ARG:HD3	1:C:637:GLU:OE1	2.16	0.44
1:C:434:HIS:O	2:H:2:VAL:HG23	2.18	0.44
3:L:37:GLN:HB2	3:L:47:LEU:HD11	1.98	0.44
2:H:166:PHE:HB3	3:L:162:SER:OG	2.18	0.43
2:H:39:GLN:HB2	2:H:45:LEU:HD23	2.00	0.43
2:H:51:ILE:HG13	2:H:57:THR:HG22	2.00	0.43
3:L:191:VAL:HG22	3:L:210:ASN:OD1	2.18	0.43
1:C:489:TYR:O	1:C:491:PRO:HD3	2.18	0.42
1:C:606:ARG:O	1:C:644:CYS:HA	2.19	0.42
1:C:542:THR:HG22	1:C:547:GLY:HA3	2.01	0.42
4:A:3:BMA:H62	4:A:4:MAN:H3	2.01	0.42
1:C:438:LEU:HA	1:C:438:LEU:HD23	1.85	0.42
3:L:33:LEU:HD13	3:L:71:PHE:CD1	2.54	0.42
2:H:35:HIS:NE2	2:H:50:ILE:HD12	2.34	0.42
3:L:23:CYS:HB2	3:L:35:TRP:CH2	2.54	0.42
1:C:586:PHE:HA	1:C:589:HIS:O	2.20	0.41
3:L:17:GLU:CG	3:L:18:ARG:N	2.82	0.41
1:C:422:ILE:HG22	1:C:529:TRP:CH2	2.56	0.41
2:H:131:THR:HB	2:H:186:SER:OG	2.21	0.41
1:C:544:PRO:HA	1:C:547:GLY:H	1.85	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	C	215/262 (82%)	200 (93%)	12 (6%)	3 (1%)	9	19
2	H	220/237 (93%)	213 (97%)	6 (3%)	1 (0%)	25	47
3	L	211/215 (98%)	204 (97%)	7 (3%)	0	100	100
All	All	646/714 (90%)	617 (96%)	25 (4%)	4 (1%)	22	43

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	132	SER
1	C	540	ASN
1	C	582	PRO
1	C	510	THR

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	C	185/219 (84%)	182 (98%)	3 (2%)	58	79
2	H	190/203 (94%)	186 (98%)	4 (2%)	48	73
3	L	185/187 (99%)	180 (97%)	5 (3%)	40	66
All	All	560/609 (92%)	548 (98%)	12 (2%)	48	73

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	537	LEU
1	C	578	THR
1	C	606	ARG
2	H	3	GLN
2	H	83	ARG
2	H	177	SER
2	H	178	LEU
3	L	14	SER
3	L	17	GLU
3	L	45	ARG
3	L	63	SER
3	L	162	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	458	HIS
1	C	467	GLN
1	C	493	GLN
2	H	32	ASN
2	H	39	GLN

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 ⓘ

12 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
4	NAG	A	1	4,1	14,14,15	0.38	0	17,19,21	1.31	2 (11%)
4	NAG	A	2	4	14,14,15	0.25	0	17,19,21	1.02	1 (5%)
4	BMA	A	3	4	11,11,12	0.91	0	15,15,17	0.64	0
4	MAN	A	4	4	11,11,12	2.42	5 (45%)	15,15,17	1.54	3 (20%)
5	NAG	B	1	5,1	14,14,15	0.39	0	17,19,21	1.04	1 (5%)
5	NAG	B	2	5	14,14,15	0.66	0	17,19,21	0.75	0
5	NAG	D	1	5,1	14,14,15	0.31	0	17,19,21	1.42	2 (11%)
5	NAG	D	2	5	14,14,15	0.54	0	17,19,21	0.82	1 (5%)
6	NAG	E	1	6,1	14,14,15	0.44	0	17,19,21	1.17	2 (11%)
6	NAG	E	2	6	14,14,15	0.39	0	17,19,21	0.80	0
6	BMA	E	3	6	11,11,12	1.01	0	15,15,17	1.02	1 (6%)
6	MAN	E	4	6	11,11,12	1.45	2 (18%)	15,15,17	1.01	1 (6%)

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	A	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	A	2	4	-	2/6/23/26	0/1/1/1
4	BMA	A	3	4	-	1/2/19/22	0/1/1/1
4	MAN	A	4	4	-	0/2/19/22	0/1/1/1
5	NAG	B	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	B	2	5	-	0/6/23/26	0/1/1/1
5	NAG	D	1	5,1	-	4/6/23/26	0/1/1/1
5	NAG	D	2	5	-	0/6/23/26	0/1/1/1
6	NAG	E	1	6,1	-	0/6/23/26	0/1/1/1
6	NAG	E	2	6	-	2/6/23/26	0/1/1/1
6	BMA	E	3	6	-	0/2/19/22	0/1/1/1
6	MAN	E	4	6	-	0/2/19/22	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	4	MAN	C1-C2	4.87	1.63	1.52
4	A	4	MAN	C2-C3	3.46	1.57	1.52
4	A	4	MAN	O5-C5	3.09	1.49	1.43
6	E	4	MAN	C4-C5	2.47	1.58	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	4	MAN	O4-C4	2.27	1.48	1.43
4	A	4	MAN	C4-C5	2.20	1.57	1.53
6	E	4	MAN	O5-C5	2.01	1.47	1.43

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	1	NAG	C2-N2-C7	-4.33	117.10	122.90
4	A	4	MAN	C1-C2-C3	4.17	115.71	109.64
6	E	3	BMA	C1-O5-C5	3.04	116.27	112.19
4	A	1	NAG	C1-O5-C5	2.99	116.19	112.19
6	E	1	NAG	O5-C1-C2	-2.67	107.17	111.29
4	A	2	NAG	C4-C3-C2	-2.62	107.18	111.02
5	D	1	NAG	O4-C4-C3	-2.61	104.23	110.38
6	E	4	MAN	O3-C3-C2	2.42	114.99	110.05
5	B	1	NAG	C2-N2-C7	-2.22	119.92	122.90
5	D	2	NAG	C2-N2-C7	2.14	125.77	122.90
4	A	1	NAG	C2-N2-C7	-2.14	120.03	122.90
4	A	4	MAN	C2-C3-C4	2.04	114.44	110.86
6	E	1	NAG	O4-C4-C3	-2.03	105.60	110.38
4	A	4	MAN	O2-C2-C3	-2.01	105.98	110.15

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	D	1	NAG	C8-C7-N2-C2
5	D	1	NAG	O7-C7-N2-C2
4	A	2	NAG	C8-C7-N2-C2
4	A	2	NAG	O7-C7-N2-C2
5	D	1	NAG	C4-C5-C6-O6
4	A	3	BMA	O5-C5-C6-O6
6	E	2	NAG	C4-C5-C6-O6
5	D	1	NAG	O5-C5-C6-O6
6	E	2	NAG	O5-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 2 short contacts:

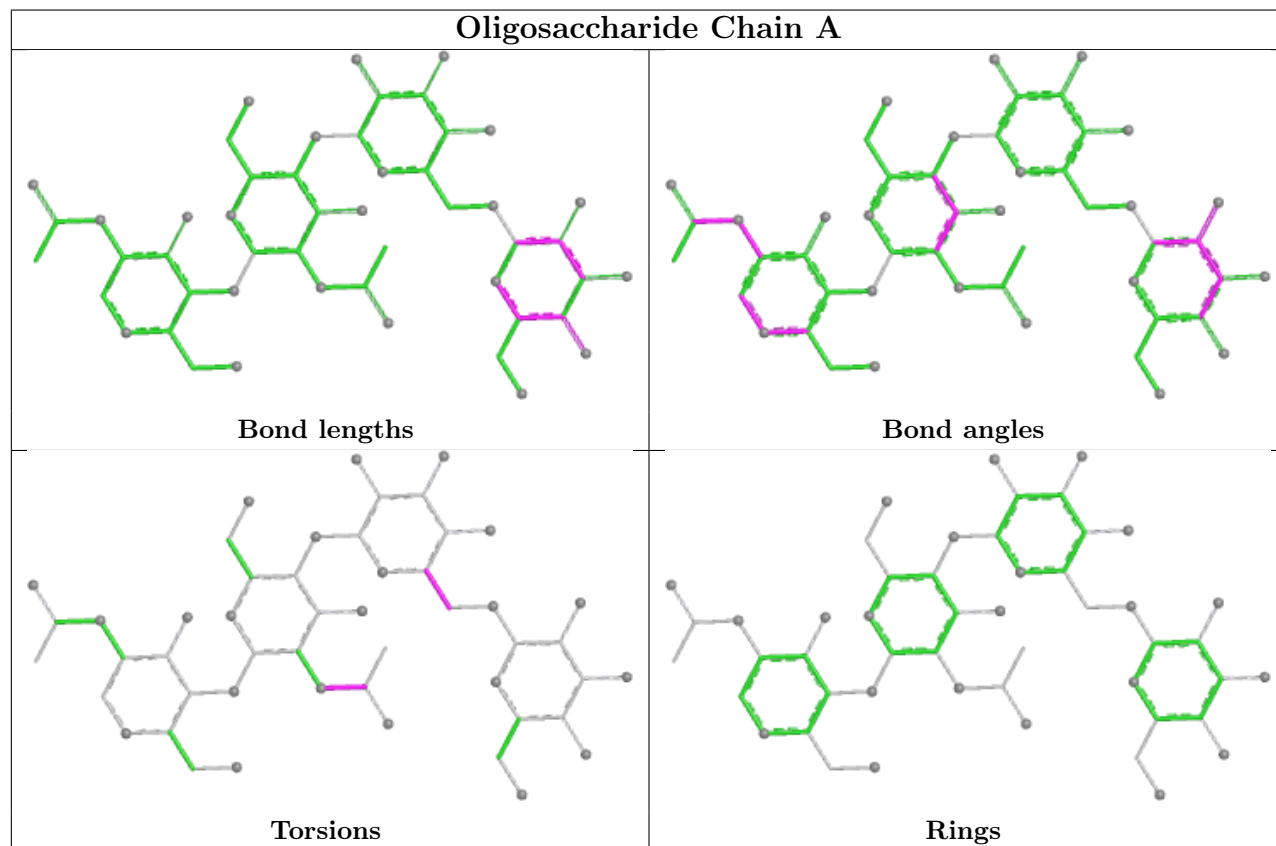
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	E	1	NAG	1	0

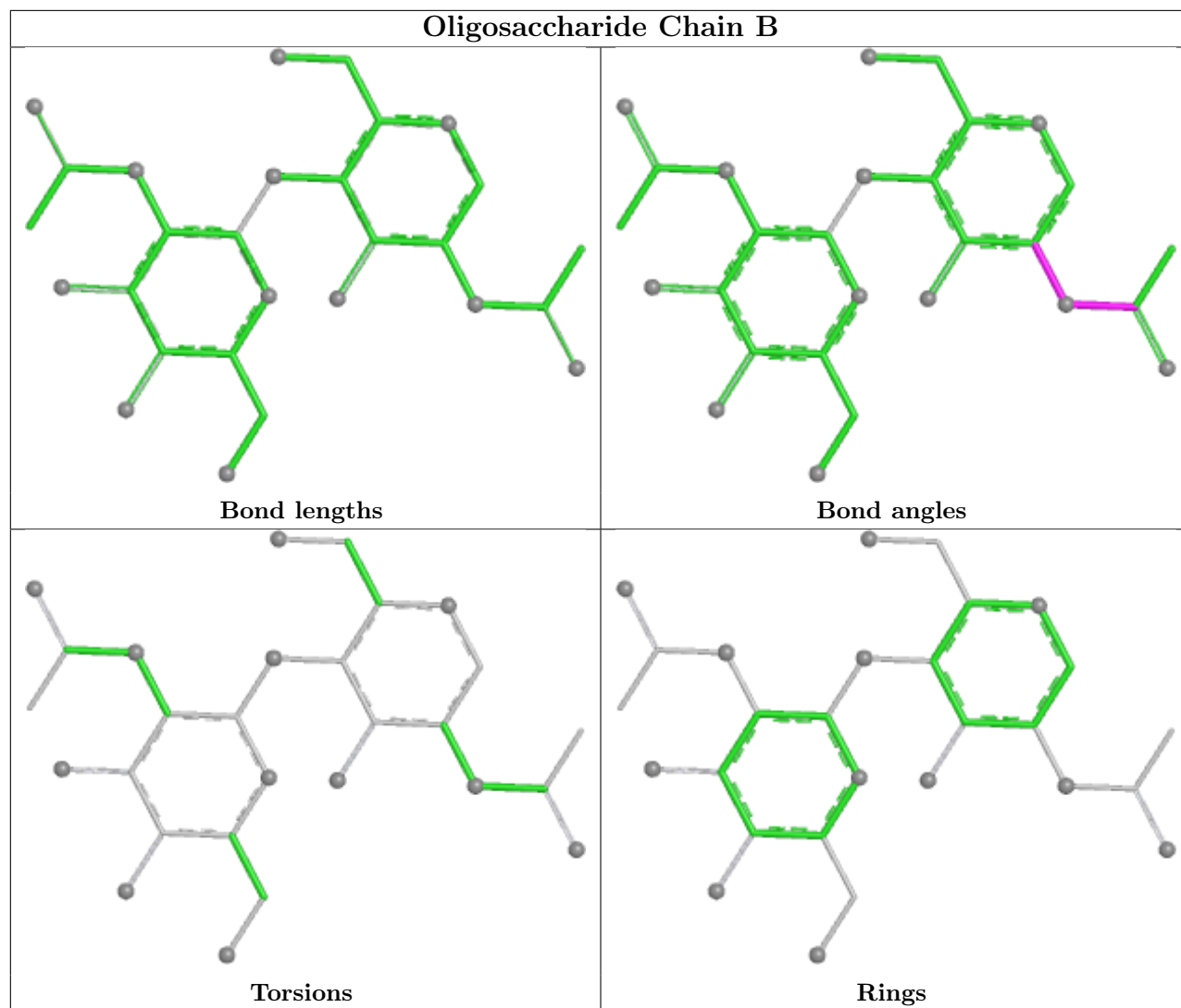
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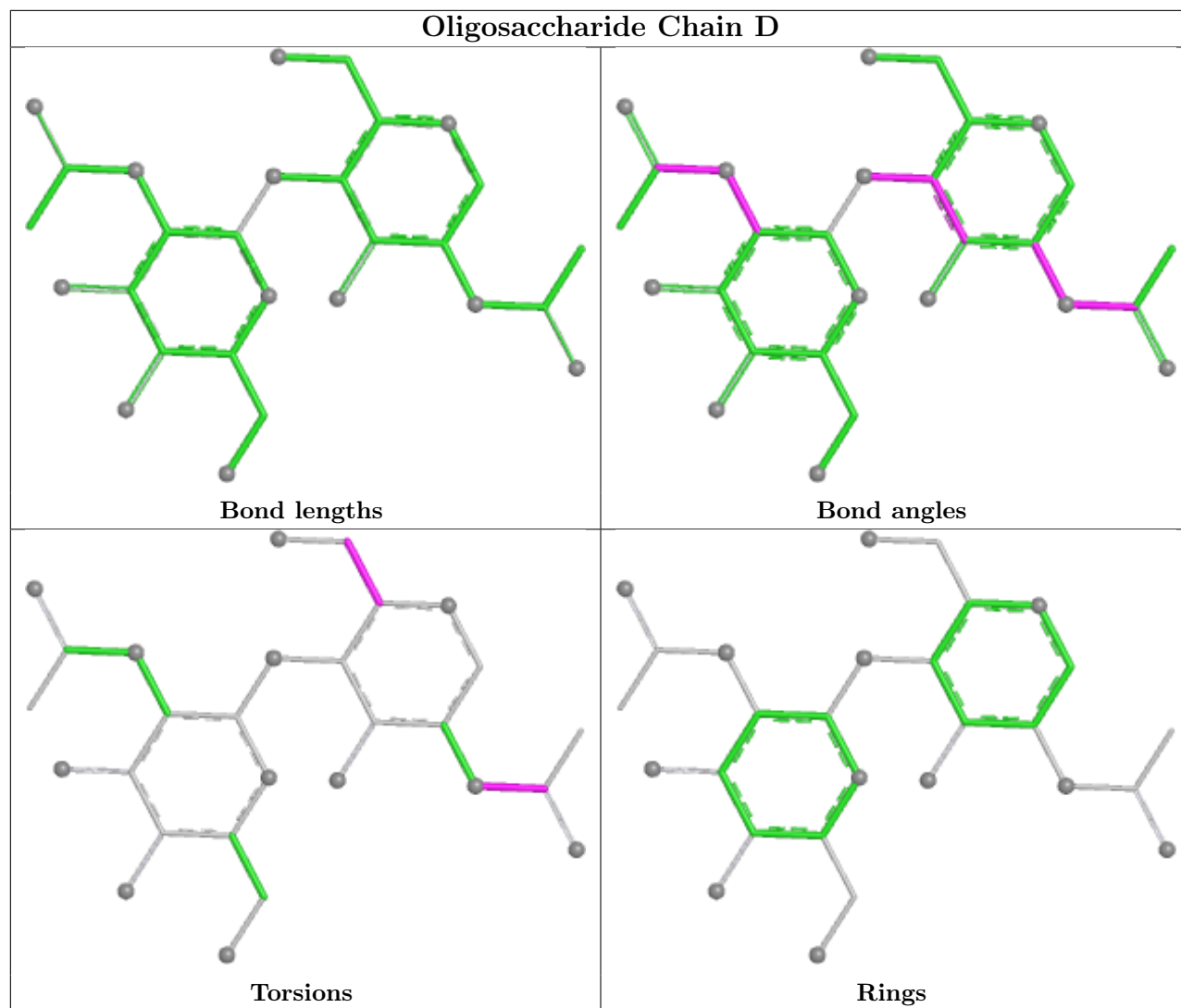
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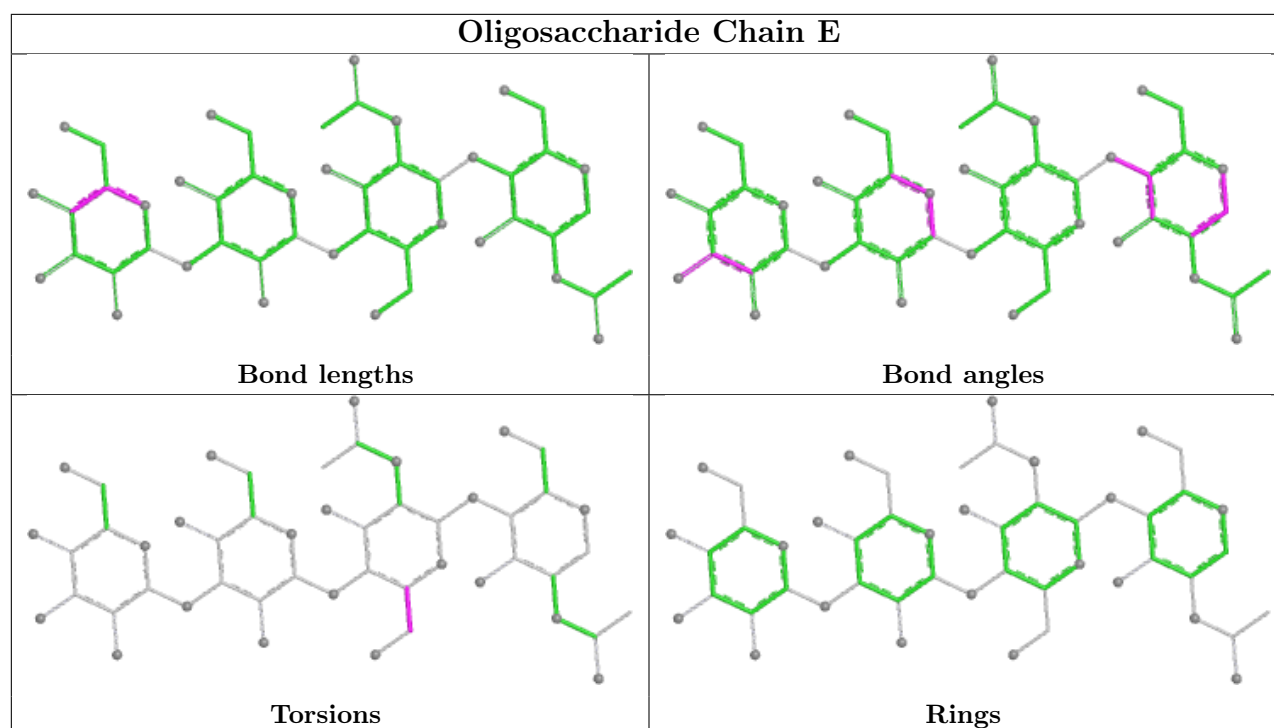
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	3	BMA	1	0
4	A	4	MAN	1	0

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	$\# Z > 2$
7	NAG	C	701	1	14,14,15	0.64	0	17,19,21	0.97	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	NAG	C	701	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.

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](#)

EDS failed to run properly - this section is therefore empty.

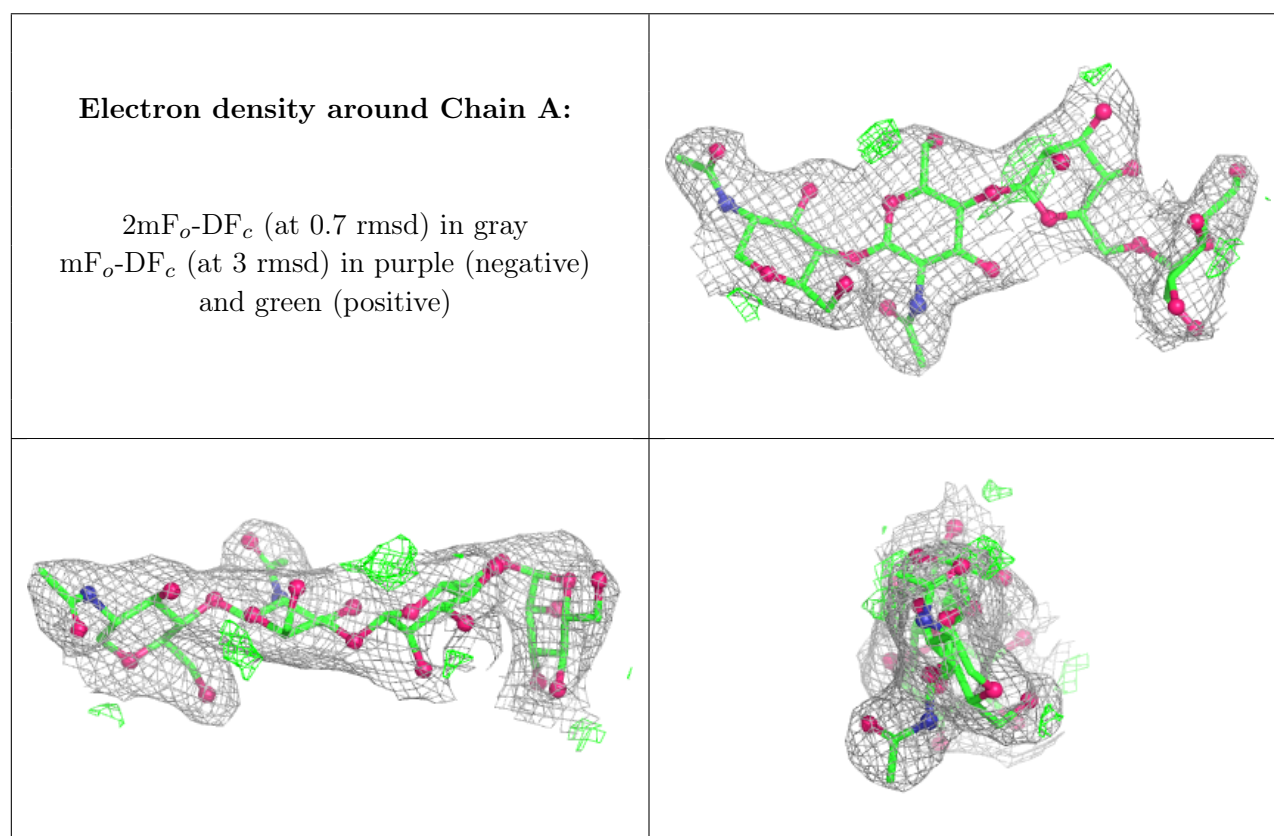
6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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6.3 Carbohydrates [i](#)

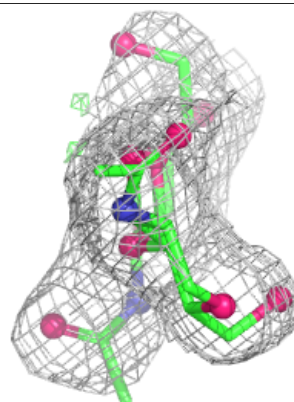
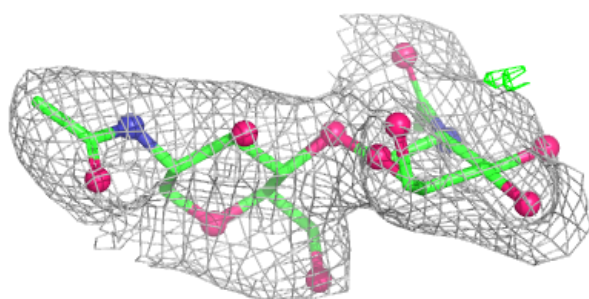
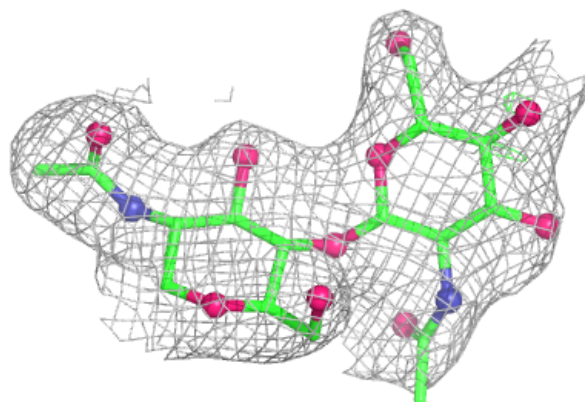
EDS failed to run properly - this section is therefore empty.

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.

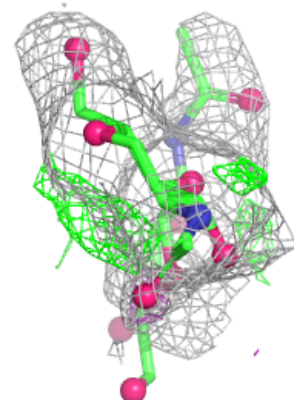
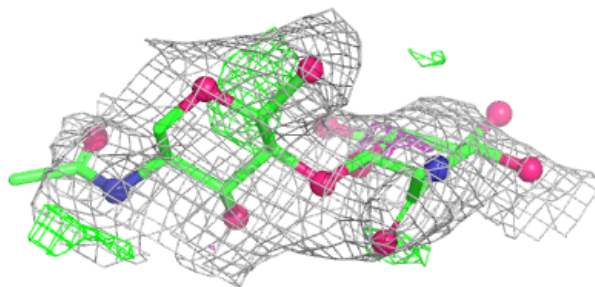
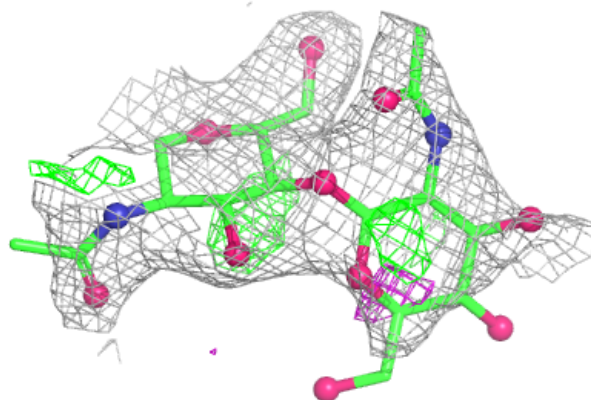


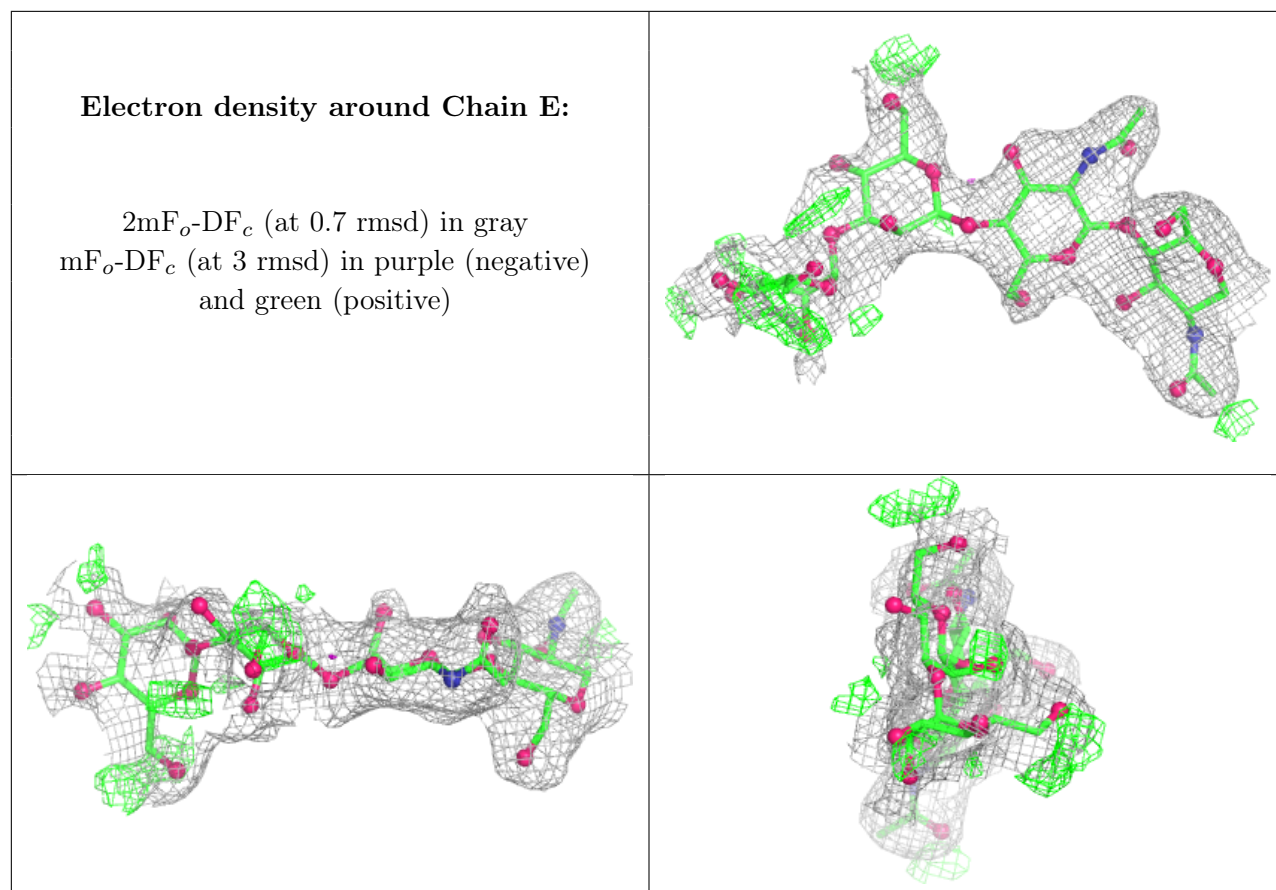
Electron density around Chain B:

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

**Electron density around Chain D:**

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





6.4 Ligands [i](#)

EDS failed to run properly - this section is therefore empty.

6.5 Other polymers [i](#)

EDS failed to run properly - this section is therefore empty.