



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 11:44 PM UTC

PDB ID : 6WRW / pdb_00006wrw
Title : Crystal structure of computationally designed protein 2DS25.5 in complex with the human Transferrin receptor ectodomain
Authors : Abraham, J.; Coscia, A.; Olal, D.; Sahtoe, D.D.; Baker, D.; Clark, L.
Deposited on : 2020-04-30
Resolution : 2.84 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

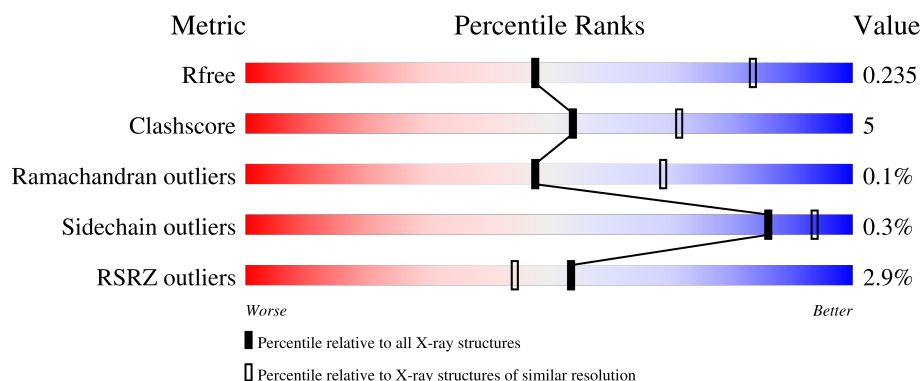
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.84 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}	180053	1520 (2.86-2.82)
Clashscore	190562	1559 (2.86-2.82)
Ramachandran outliers	187476	1517 (2.86-2.82)
Sidechain outliers	187428	1518 (2.86-2.82)
RSRZ outliers	180081	1521 (2.86-2.82)

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 $\leq 5\%$. 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	640	2% 86% 14%
1	B	640	3% 87% 12%
2	C	88	3% 84% 15% .
2	D	88	5% 88% 11% .
3	E	2	100%

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Mol	Chain	Length	Quality of chain
3	G	2	 100%
3	H	2	 100%
3	J	2	 100%
4	F	3	 100%
5	I	5	 60%  40%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11757 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 Transferrin receptor protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	639	Total	C	N	O	S	0	0	0
			5052	3239	850	949	14			
1	B	640	Total	C	N	O	S	0	0	0
			5056	3242	851	949	14			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	142	SER	GLY	conflict	UNP P02786
B	142	SER	GLY	conflict	UNP P02786

- Molecule 2 is a protein called Computationally designed protein 2DS25.5.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	88	Total	C	N	O	0	0	0
			717	457	123	137			
2	D	88	Total	C	N	O	0	0	0
			718	457	123	138			

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



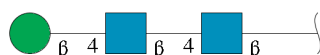
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	E	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	G	2	Total	C	N	O	0	0	0
			28	16	2	10			

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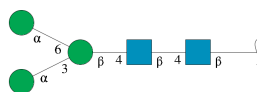
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	H	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	J	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 4 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
4	F	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 5 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[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
5	I	5	Total	C	N	O	0	0	0
			61	34	2	25			

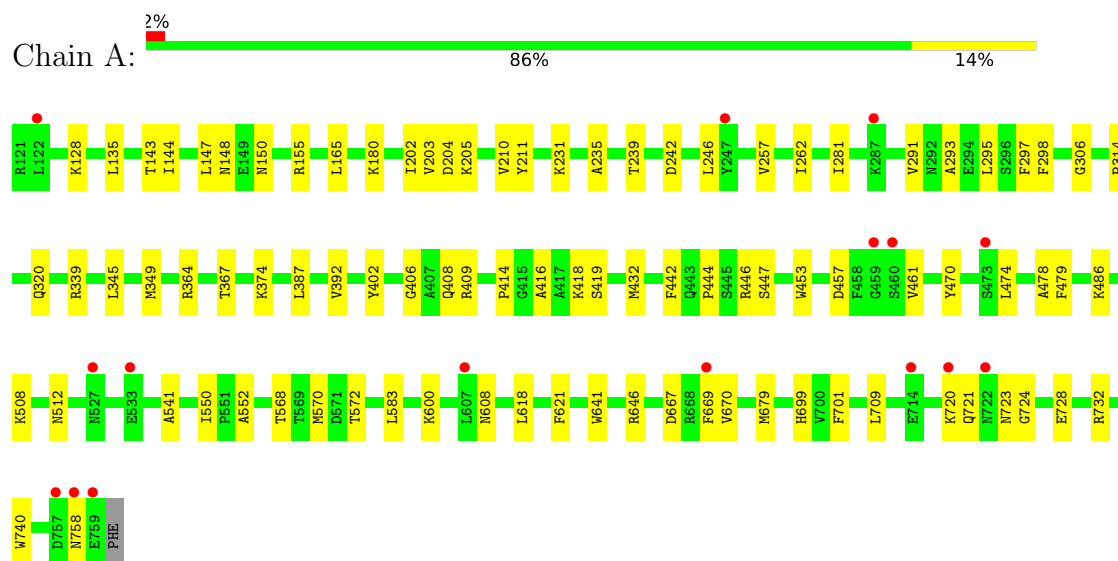
- Molecule 6 is CALCIUM ION (CCD ID: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

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

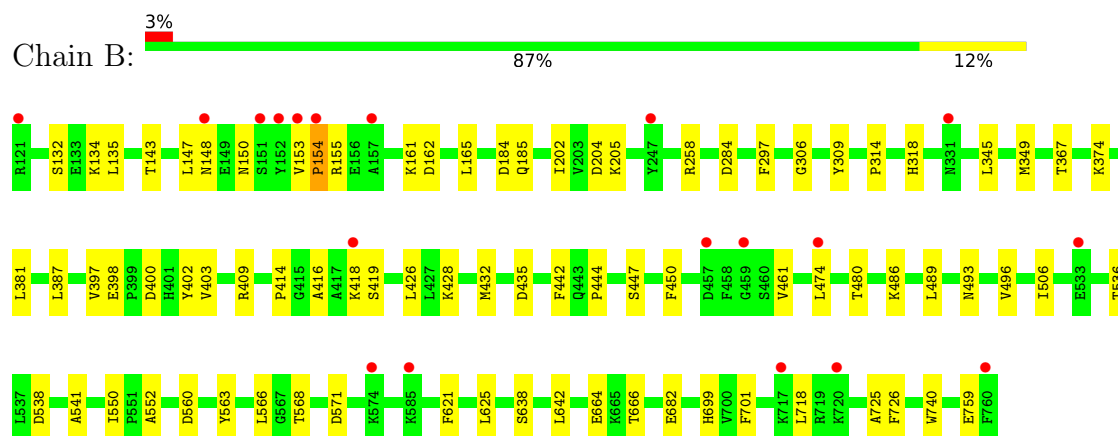
3 Residue-property plots [i](#)

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.

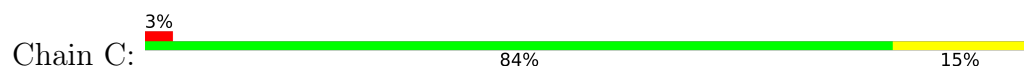
- Molecule 1: Transferrin receptor protein 1



- Molecule 1: Transferrin receptor protein 1

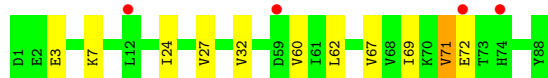
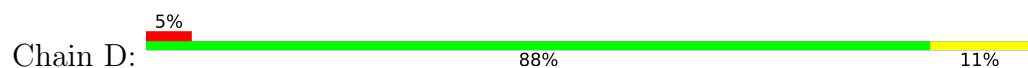


- Molecule 2: Computationally designed protein 2DS25.5





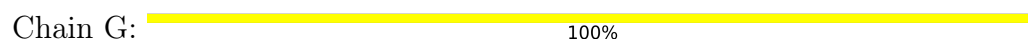
- Molecule 2: Computationally designed protein 2DS25.5



- 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



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



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



- Molecule 5: alpha-D-mannopyranose-(1-3)-[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

Chain I:



MAG1	MAG2	MAN4	MAN5
EM13			

4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	138.40Å 138.40Å 279.85Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	69.20 – 2.84 69.20 – 2.84	Depositor EDS
% Data completeness (in resolution range)	97.7 (69.20-2.84) 97.7 (69.20-2.84)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.60 (at 2.86Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.211 , 0.235 0.214 , 0.235	Depositor DCC
R_{free} test set	3632 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	88.4	Xtriage
Anisotropy	0.173	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 50.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.011 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11757	wwPDB-VP
Average B, all atoms (Å ²)	88.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.23% 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: CA, NAG, MAN, BMA

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	0.10	0/5172	0.29	0/7014
1	B	0.10	0/5176	0.28	0/7019
2	C	0.09	0/726	0.27	0/982
2	D	0.10	0/727	0.28	0/982
All	All	0.10	0/11801	0.28	0/15997

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	150	ASN	Peptide

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	5052	0	4983	54	0
1	B	5056	0	4984	54	0
2	C	717	0	731	8	0
2	D	718	0	731	6	0
3	E	28	0	25	0	0
3	G	28	0	25	0	0
3	H	28	0	25	0	0
3	J	28	0	25	1	0
4	F	39	0	34	0	0
5	I	61	0	52	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
All	All	11757	0	11615	115	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 5.

All (115) 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:153:VAL:HG23	1:B:154:PRO:HD3	1.62	0.81
1:B:154:PRO:HD2	1:B:161:LYS:HD2	1.63	0.79
1:A:135:LEU:HG	1:A:432:MET:HE1	1.71	0.72
1:B:135:LEU:HG	1:B:432:MET:HE1	1.76	0.68
1:A:306:GLY:HA2	1:A:461:VAL:HA	1.75	0.67
1:B:306:GLY:HA2	1:B:461:VAL:HA	1.77	0.67
1:B:474:LEU:HD13	1:B:550:ILE:HD11	1.77	0.66
1:A:600:LYS:O	1:A:608:ASN:ND2	2.29	0.66
1:A:479:PHE:HB3	1:A:679:MET:HE3	1.79	0.65
1:A:728:GLU:OE2	1:A:732:ARG:NH1	2.30	0.64
1:A:387:LEU:O	1:A:409:ARG:NH2	2.31	0.61
1:B:541:ALA:HB1	1:B:552:ALA:HB1	1.82	0.61
1:A:320:GLN:NE2	1:B:638:SER:O	2.32	0.60
1:A:408:GLN:NE2	1:A:457:ASP:OD1	2.31	0.60
1:B:387:LEU:O	1:B:409:ARG:NH2	2.35	0.60
1:A:203:VAL:HG12	1:A:210:VAL:HG12	1.84	0.59
2:D:3:GLU:O	2:D:7:LYS:HG2	2.01	0.59
1:B:162:ASP:OD1	1:B:409:ARG:NH1	2.35	0.59
1:A:297:PHE:HB3	1:A:568:THR:HG22	1.84	0.58
1:B:402:TYR:HB3	1:B:447:SER:HB2	1.85	0.58
1:A:402:TYR:HB3	1:A:447:SER:HB2	1.86	0.58
1:B:155:ARG:NH2	1:B:419:SER:OG	2.37	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:493:ASN:ND2	1:B:560:ASP:OD1	2.35	0.57
1:B:150:ASN:HA	1:B:153:VAL:HG22	1.88	0.56
2:C:54:TYR:O	2:C:79:ARG:NH2	2.40	0.55
1:B:418:LYS:HA	1:B:486:LYS:HG2	1.89	0.55
1:A:295:LEU:HD22	1:A:570:MET:HE2	1.89	0.55
1:B:132:SER:HA	1:B:135:LEU:HD12	1.88	0.54
1:A:508:LYS:O	1:A:512:ASN:ND2	2.41	0.54
1:B:202:ILE:HD11	1:B:345:LEU:HD23	1.90	0.54
1:B:297:PHE:HB3	1:B:568:THR:HG22	1.90	0.53
1:A:541:ALA:HB1	1:A:552:ALA:HB1	1.91	0.52
1:A:618:LEU:HB3	1:A:646:ARG:HG3	1.92	0.52
1:A:298:PHE:HE1	1:A:457:ASP:HB3	1.74	0.52
1:B:496:VAL:HG11	1:B:506:ILE:HD13	1.92	0.52
1:A:235:ALA:HA	1:A:246:LEU:HD11	1.92	0.51
1:A:202:ILE:HD11	1:A:345:LEU:HD23	1.93	0.51
1:A:239:THR:OG1	1:A:242:ASP:OD1	2.29	0.51
1:B:664:GLU:HG3	1:B:666:THR:H	1.77	0.50
1:A:349:MET:HG3	1:A:364:ARG:HB2	1.93	0.50
1:A:418:LYS:HA	1:A:486:LYS:HG2	1.94	0.50
1:B:258:ARG:NH1	1:B:284:ASP:OD2	2.44	0.50
1:A:128:LYS:NZ	1:A:512:ASN:HB2	2.27	0.50
1:B:349:MET:HG2	1:B:367:THR:HG22	1.93	0.49
1:A:144:ILE:HD12	1:A:583:LEU:HD21	1.93	0.49
1:A:758:ASN:ND2	1:B:184:ASP:OD1	2.45	0.49
1:B:621:PHE:HE2	1:B:642:LEU:HD21	1.77	0.49
1:A:314:PRO:HG3	1:B:740:TRP:CE2	2.48	0.48
2:C:62:LEU:HD13	2:C:67:VAL:HG22	1.95	0.48
1:A:721:GLN:HG3	1:A:723:ASN:H	1.79	0.48
1:A:349:MET:HB3	1:A:367:THR:HA	1.96	0.47
1:A:740:TRP:CE2	1:B:314:PRO:HG3	2.48	0.47
1:A:621:PHE:CE1	1:A:709:LEU:HB3	2.49	0.47
2:D:60:VAL:HG22	2:D:69:ILE:HG23	1.97	0.46
1:B:309:TYR:CZ	1:B:381:LEU:HD21	2.51	0.46
1:B:442:PHE:CE2	1:B:444:PRO:HG3	2.51	0.46
1:B:148:ASN:HD21	1:B:416:ALA:HB2	1.80	0.46
2:C:27:VAL:HA	2:C:32:VAL:HG22	1.96	0.46
1:B:400:ASP:OD1	1:B:400:ASP:N	2.47	0.46
2:C:60:VAL:HG22	2:C:69:ILE:HG23	1.98	0.46
1:A:204:ASP:OD1	1:A:205:LYS:N	2.47	0.46
1:A:720:LYS:HA	1:A:720:LYS:HD2	1.77	0.46
1:B:349:MET:HA	1:B:367:THR:HA	1.96	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:231:LYS:NZ	1:A:367:THR:OG1	2.49	0.45
1:A:293:ALA:HB2	1:A:339:ARG:NH2	2.32	0.45
1:A:699:HIS:CE1	1:A:701:PHE:HB2	2.51	0.45
2:D:62:LEU:HD13	2:D:67:VAL:HG22	1.98	0.45
1:A:406:GLY:HA3	1:A:453:TRP:HE1	1.82	0.45
2:D:27:VAL:HA	2:D:32:VAL:HG22	1.99	0.45
1:A:446:ARG:HD2	1:A:479:PHE:CE1	2.52	0.45
1:A:148:ASN:HD21	1:A:416:ALA:HB2	1.83	0.44
1:A:374:LYS:HE3	2:C:24:ILE:HD13	2.00	0.44
1:A:442:PHE:CE2	1:A:444:PRO:HG3	2.52	0.44
1:B:718:LEU:HD13	1:B:725:ALA:HB1	1.99	0.44
1:A:155:ARG:NH1	1:A:419:SER:OG	2.44	0.44
1:B:563:TYR:HB3	1:B:566:LEU:HG	2.00	0.44
1:B:185:GLN:HG2	1:B:387:LEU:HD23	1.99	0.43
1:B:143:THR:O	1:B:147:LEU:HG	2.19	0.43
1:B:147:LEU:HD22	1:B:165:LEU:HD11	1.99	0.43
2:C:71:VAL:HG12	2:C:72:GLU:H	1.82	0.43
3:J:1:NAG:H61	3:J:2:NAG:C7	2.48	0.43
1:A:474:LEU:HD13	1:A:550:ILE:HD11	2.00	0.43
2:D:71:VAL:HG12	2:D:72:GLU:H	1.83	0.43
1:B:143:THR:HG21	1:B:428:LYS:HE2	2.01	0.43
1:B:204:ASP:OD1	1:B:205:LYS:N	2.50	0.43
1:A:414:PRO:HG2	1:A:572:THR:HG22	2.00	0.43
1:A:470:TYR:O	1:A:474:LEU:HG	2.19	0.42
1:B:536:THR:OG1	1:B:538:ASP:OD1	2.35	0.42
1:B:403:VAL:HG12	1:B:480:THR:HB	2.00	0.42
1:A:242:ASP:O	1:A:246:LEU:HG	2.20	0.42
1:B:718:LEU:HB2	1:B:726:PHE:HB2	2.01	0.42
1:A:667:ASP:HB3	1:A:670:VAL:HG12	2.01	0.42
1:A:641:TRP:CG	1:B:318:HIS:HA	2.55	0.41
1:B:414:PRO:HB2	1:B:571:ASP:O	2.20	0.41
1:A:723:ASN:OD1	1:A:724:GLY:N	2.53	0.41
1:B:153:VAL:CG2	1:B:154:PRO:HD3	2.39	0.41
1:B:682:GLU:OE2	1:B:699:HIS:NE2	2.40	0.41
1:B:625:LEU:HD12	1:B:625:LEU:HA	1.91	0.41
1:B:759:GLU:H	1:B:759:GLU:CD	2.28	0.41
1:A:165:LEU:HD23	1:A:409:ARG:HG3	2.03	0.41
1:B:718:LEU:H	1:B:718:LEU:HG	1.68	0.41
1:A:143:THR:O	1:A:147:LEU:HG	2.20	0.41
1:A:402:TYR:CE2	1:A:478:ALA:HB2	2.56	0.41
1:B:374:LYS:HE3	2:D:24:ILE:HD13	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:397:VAL:HG12	1:B:398:GLU:HG3	2.03	0.41
1:B:426:LEU:HD11	1:B:450:PHE:HB3	2.03	0.41
1:A:211:TYR:HB2	2:C:81:TRP:CZ3	2.56	0.40
1:A:239:THR:HG22	1:A:262:ILE:HD13	2.03	0.40
1:A:257:VAL:O	1:A:281:ILE:HA	2.21	0.40
1:B:134:LYS:HE3	1:B:435:ASP:HB3	2.03	0.40
1:B:642:LEU:HD12	1:B:642:LEU:HA	1.89	0.40
2:C:3:GLU:O	2:C:7:LYS:HG2	2.21	0.40
1:B:699:HIS:CE1	1:B:701:PHE:HB2	2.56	0.40
1:A:180:LYS:O	1:A:392:VAL:HG12	2.21	0.40
1:B:486:LYS:HB3	1:B:489:LEU:HD11	2.02	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	637/640 (100%)	613 (96%)	24 (4%)	0	100	100
1	B	638/640 (100%)	621 (97%)	16 (2%)	1 (0%)	43	62
2	C	86/88 (98%)	85 (99%)	1 (1%)	0	100	100
2	D	86/88 (98%)	85 (99%)	1 (1%)	0	100	100
All	All	1447/1456 (99%)	1404 (97%)	42 (3%)	1 (0%)	48	69

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	154	PRO

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	547/549 (100%)	545 (100%)	2 (0%)	84	92
1	B	547/549 (100%)	547 (100%)	0	100	100
2	C	77/77 (100%)	76 (99%)	1 (1%)	61	79
2	D	77/77 (100%)	76 (99%)	1 (1%)	61	79
All	All	1248/1252 (100%)	1244 (100%)	4 (0%)	86	94

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	291	VAL
1	A	669	PHE
2	C	71	VAL
2	D	71	VAL

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

Mol	Chain	Res	Type
1	A	148	ASN
1	A	300	HIS
1	A	603	HIS
1	A	743	GLN
1	B	148	ASN
1	B	331	ASN
1	B	372	ASN
1	B	512	ASN
1	B	640	GLN
1	B	715	ASN
1	B	733	ASN
1	B	743	GLN
2	C	74	HIS

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 ⓘ

16 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	3,1	14,14,15	0.23	0	17,19,21	0.46	0
3	NAG	E	2	3	14,14,15	0.25	0	17,19,21	0.46	0
4	NAG	F	1	4,1	14,14,15	0.21	0	17,19,21	0.44	0
4	NAG	F	2	4	14,14,15	0.21	0	17,19,21	0.43	0
4	BMA	F	3	4	11,11,12	0.51	0	15,15,17	0.73	0
3	NAG	G	1	3,1	14,14,15	0.24	0	17,19,21	0.72	1 (5%)
3	NAG	G	2	3	14,14,15	0.65	1 (7%)	17,19,21	0.44	0
3	NAG	H	1	3,1	14,14,15	0.27	0	17,19,21	0.52	0
3	NAG	H	2	3	14,14,15	0.26	0	17,19,21	0.40	0
5	NAG	I	1	5,1	14,14,15	0.23	0	17,19,21	0.42	0
5	NAG	I	2	5	14,14,15	0.16	0	17,19,21	0.53	0
5	BMA	I	3	5	11,11,12	0.39	0	15,15,17	0.72	0
5	MAN	I	4	5	11,11,12	0.62	0	15,15,17	0.96	2 (13%)
5	MAN	I	5	5	11,11,12	0.61	0	15,15,17	0.97	2 (13%)
3	NAG	J	1	3,1	14,14,15	0.21	0	17,19,21	0.66	1 (5%)
3	NAG	J	2	3	14,14,15	0.34	0	17,19,21	0.75	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
3	NAG	E	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	E	2	3	-	4/6/23/26	0/1/1/1
4	NAG	F	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	F	2	4	-	2/6/23/26	0/1/1/1
4	BMA	F	3	4	-	0/2/19/22	0/1/1/1
3	NAG	G	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	G	2	3	-	2/6/23/26	0/1/1/1
3	NAG	H	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	H	2	3	-	0/6/23/26	0/1/1/1
5	NAG	I	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	I	2	5	-	0/6/23/26	0/1/1/1
5	BMA	I	3	5	-	2/2/19/22	0/1/1/1
5	MAN	I	4	5	-	0/2/19/22	0/1/1/1
5	MAN	I	5	5	-	0/2/19/22	0/1/1/1
3	NAG	J	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	J	2	3	-	1/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	2	NAG	C1-C2	2.03	1.55	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	J	2	NAG	C1-O5-C5	2.71	115.82	112.19
5	I	5	MAN	C1-O5-C5	2.35	115.33	112.19
5	I	4	MAN	O2-C2-C3	-2.33	105.33	110.15
5	I	5	MAN	O2-C2-C3	-2.31	105.36	110.15
5	I	4	MAN	C1-O5-C5	2.25	115.20	112.19
3	J	1	NAG	C1-O5-C5	2.19	115.12	112.19
3	G	1	NAG	C1-O5-C5	2.06	114.95	112.19

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	I	3	BMA	O5-C5-C6-O6
5	I	3	BMA	C4-C5-C6-O6
3	E	2	NAG	O5-C5-C6-O6
3	E	2	NAG	C4-C5-C6-O6

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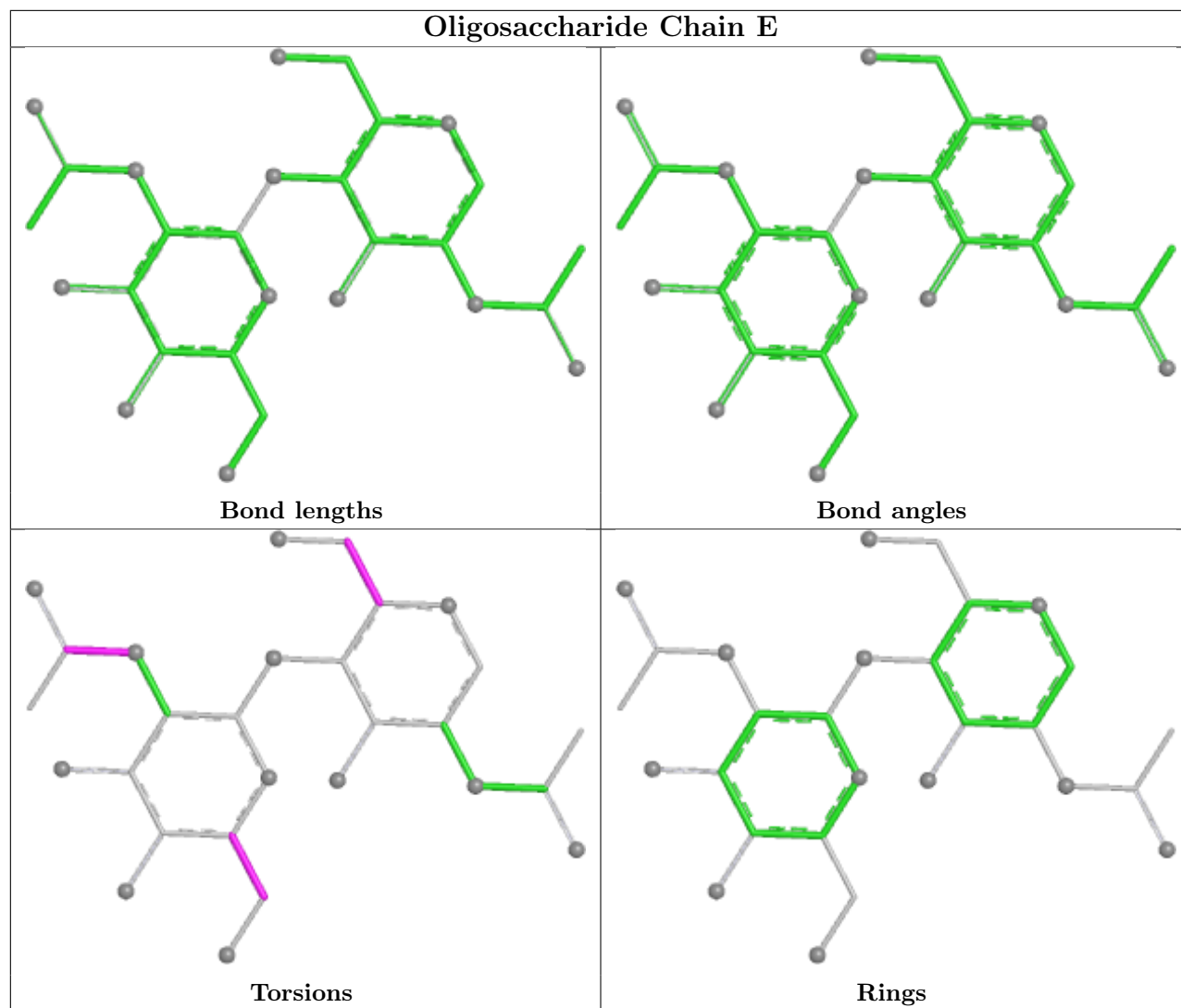
Mol	Chain	Res	Type	Atoms
3	E	2	NAG	C8-C7-N2-C2
3	E	2	NAG	O7-C7-N2-C2
3	G	1	NAG	O5-C5-C6-O6
3	G	2	NAG	C4-C5-C6-O6
3	G	2	NAG	O5-C5-C6-O6
4	F	2	NAG	C4-C5-C6-O6
4	F	2	NAG	O5-C5-C6-O6
3	E	1	NAG	C4-C5-C6-O6
3	E	1	NAG	O5-C5-C6-O6
3	J	2	NAG	C1-C2-N2-C7
3	G	1	NAG	C4-C5-C6-O6

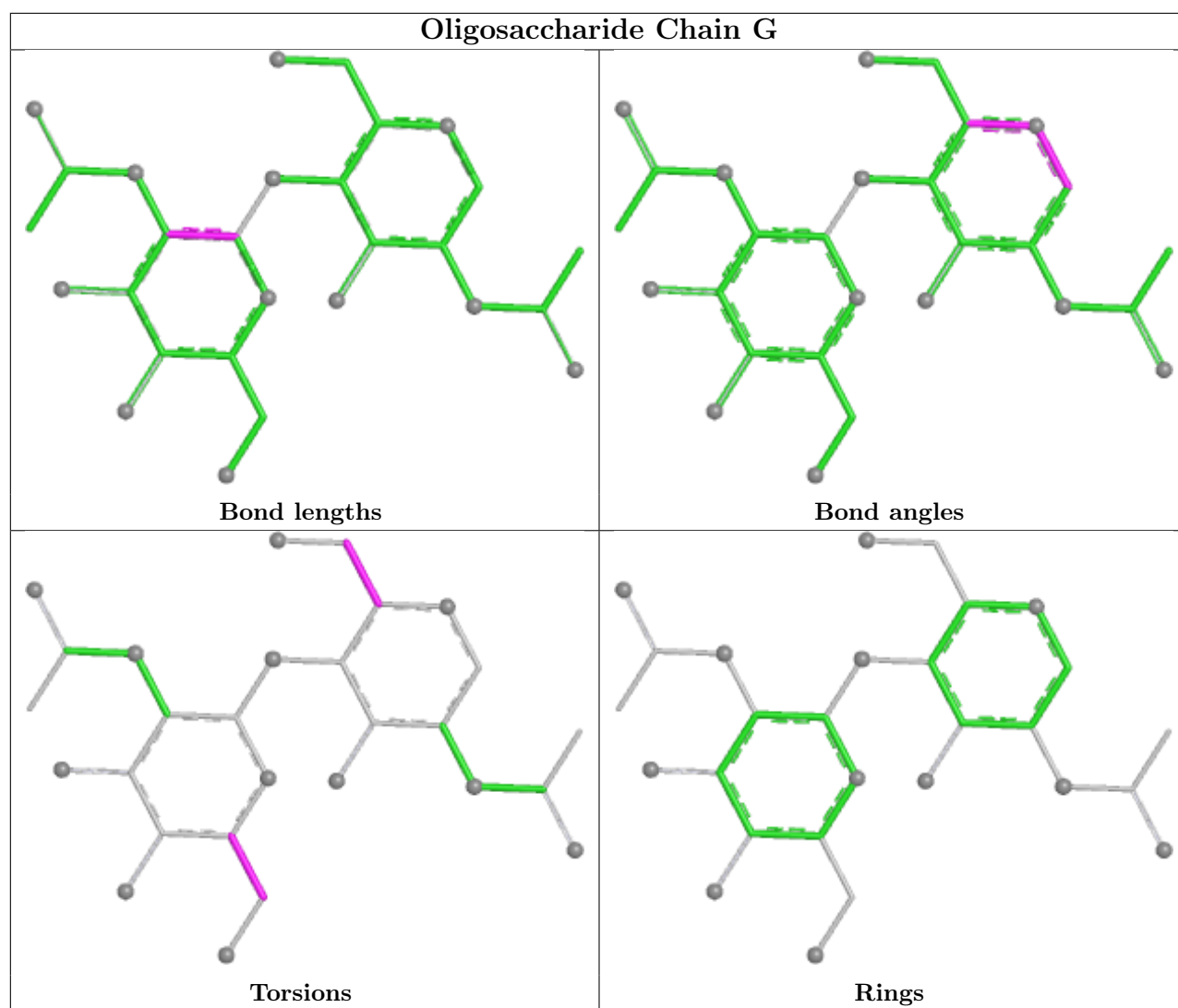
There are no ring outliers.

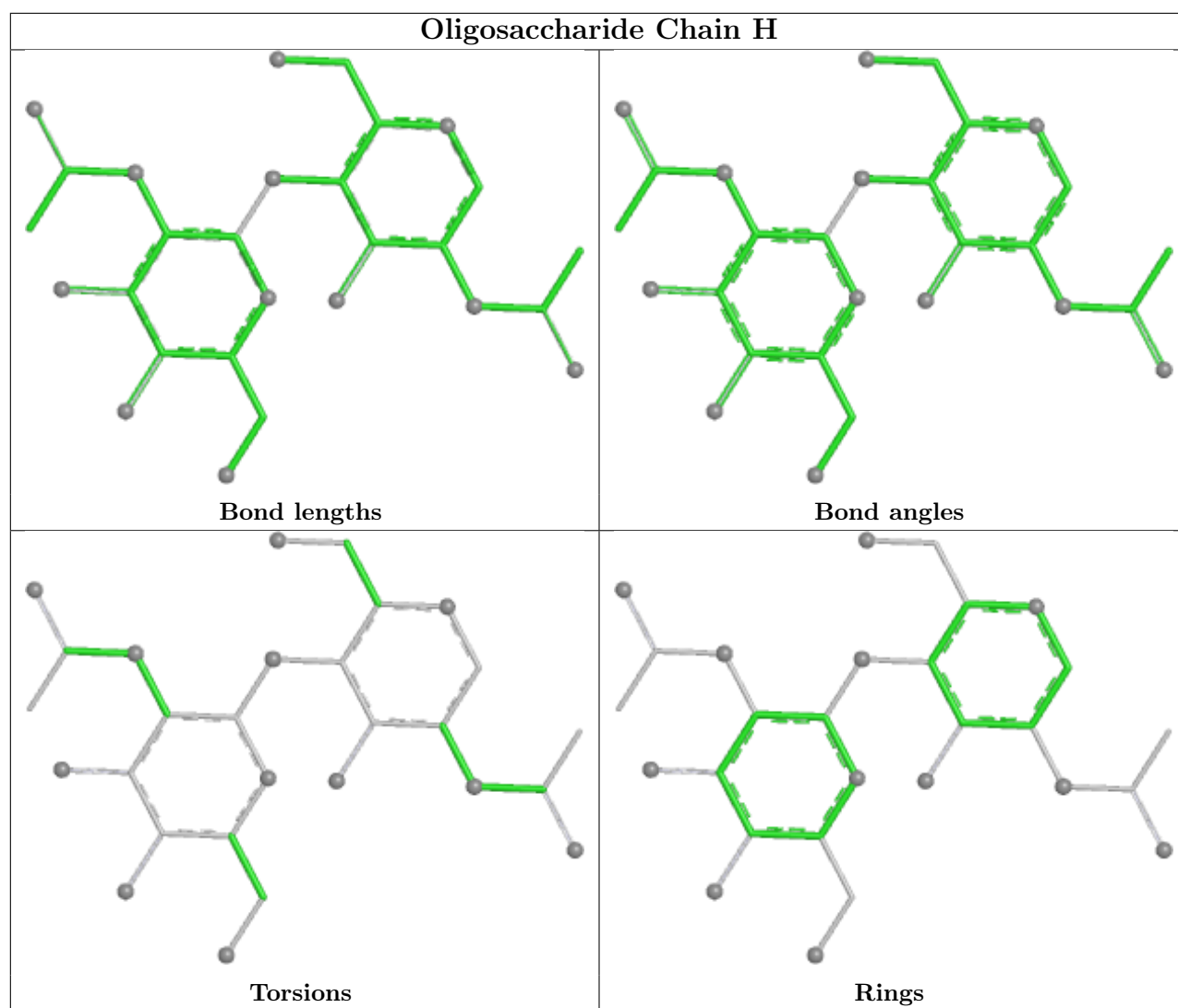
2 monomers are involved in 1 short contact:

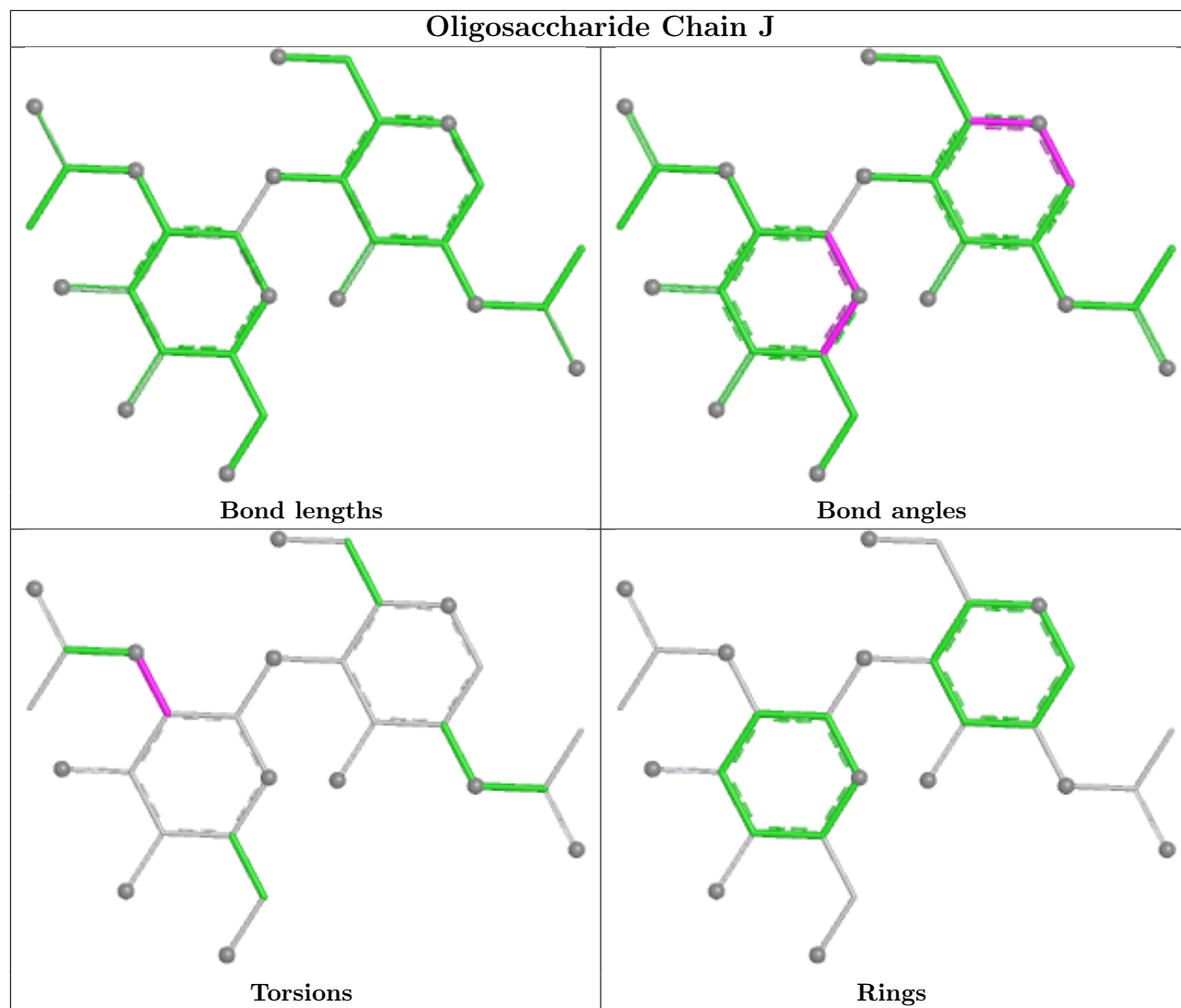
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	J	2	NAG	1	0
3	J	1	NAG	1	0

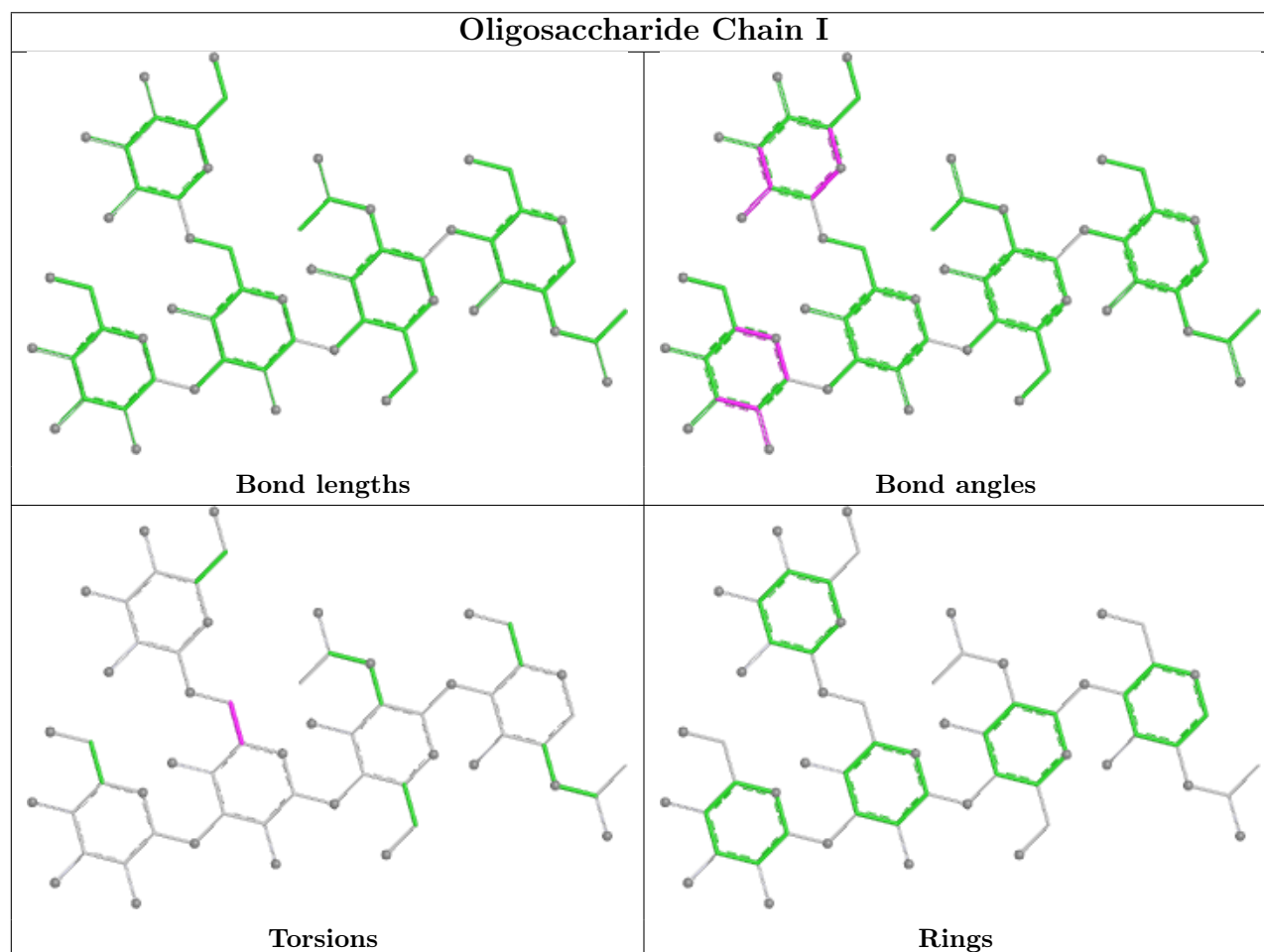
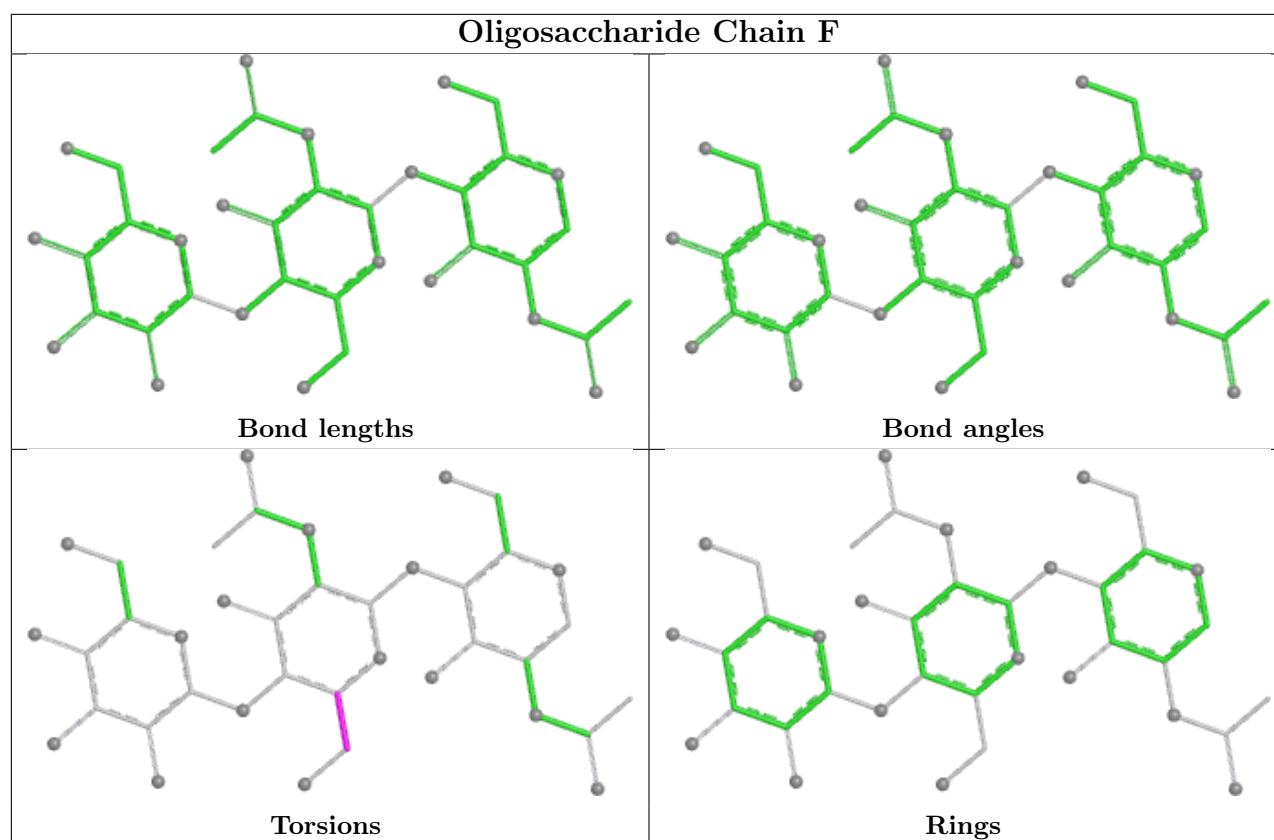
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











5.6 Ligand geometry

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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

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, 95th 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 ²)	Q<0.9
1	A	639/640 (99%)	0.20	16 (2%) 58 49	55, 83, 123, 185	0
1	B	640/640 (100%)	0.21	19 (2%) 52 43	55, 79, 118, 186	0
2	C	88/88 (100%)	0.59	3 (3%) 48 38	74, 108, 147, 160	0
2	D	88/88 (100%)	0.32	4 (4%) 38 30	66, 92, 139, 175	0
All	All	1455/1456 (99%)	0.24	42 (2%) 53 44	55, 83, 126, 186	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	760	PHE	8.5
1	B	153	VAL	5.3
1	B	457	ASP	4.7
1	A	247	TYR	4.5
1	B	459	GLY	4.3
1	A	714	GLU	4.1
1	A	607	LEU	3.9
1	B	474	LEU	3.5
1	B	720	LYS	3.2
1	A	669	PHE	3.2
1	A	757	ASP	3.1
1	B	121	ARG	3.0
1	A	758	ASN	3.0
1	B	585	LYS	2.9
1	A	759	GLU	2.8
1	B	533	GLU	2.8
1	B	418	LYS	2.7
1	B	717	LYS	2.7
1	B	157	ALA	2.6
2	C	39	GLN	2.5
1	B	574	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
2	D	59	ASP	2.4
1	A	460	SER	2.4
1	A	720	LYS	2.4
2	C	1	ASP	2.4
2	C	40	ASP	2.3
1	A	459	GLY	2.3
2	D	12	LEU	2.3
2	D	74	HIS	2.3
2	D	72	GLU	2.3
1	B	151	SER	2.3
1	A	122	LEU	2.2
1	B	148	ASN	2.2
1	B	154	PRO	2.2
1	A	287	LYS	2.2
1	A	722	ASN	2.2
1	A	533	GLU	2.1
1	A	527	ASN	2.1
1	B	247	TYR	2.1
1	B	331	ASN	2.1
1	A	473	SER	2.0
1	B	152	TYR	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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

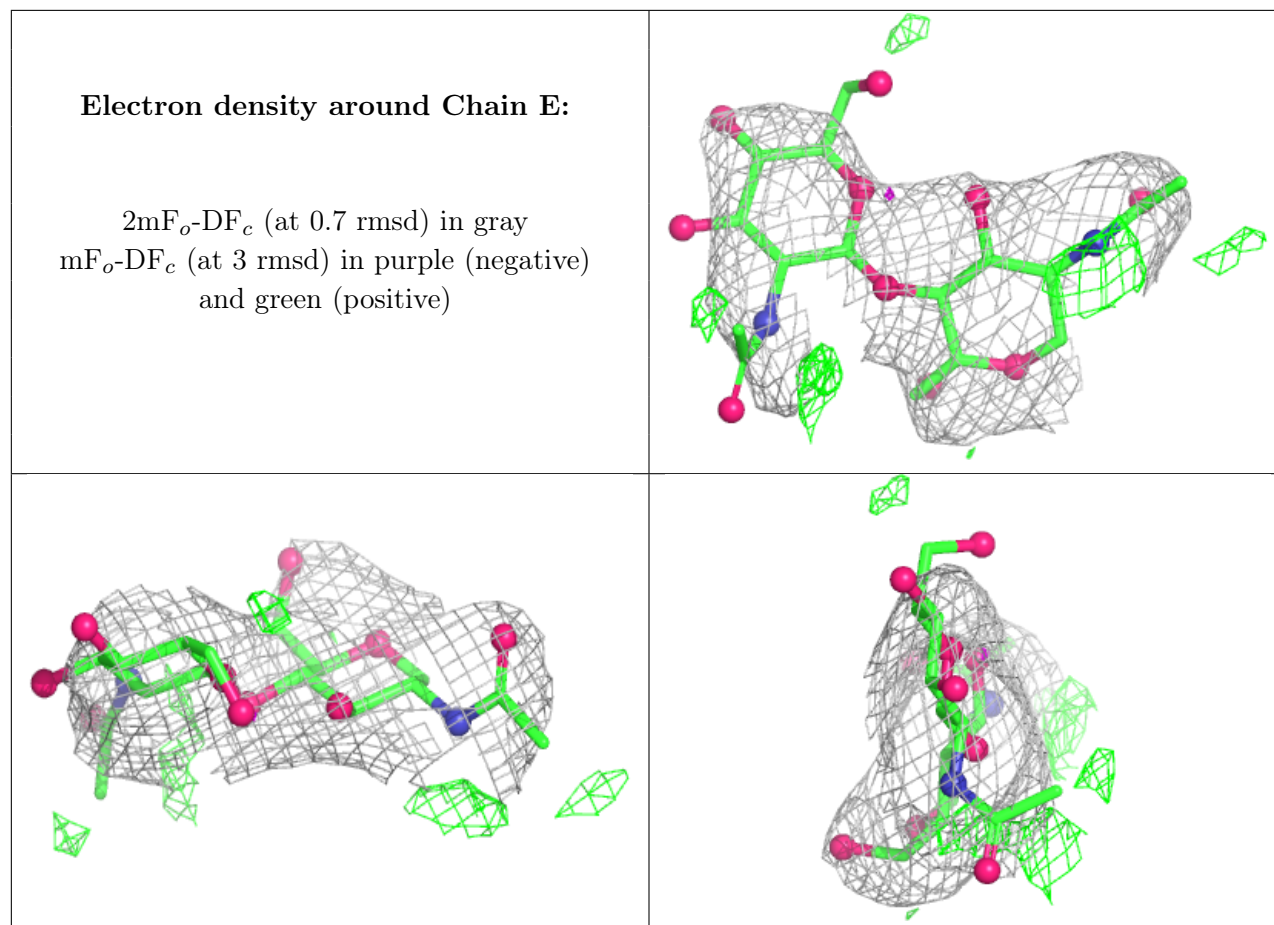
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	G	2	14/15	0.45	0.15	165,176,178,178	0
4	BMA	F	3	11/12	0.53	0.13	141,160,165,171	0
5	MAN	I	5	11/12	0.53	0.16	137,156,173,174	0
3	NAG	E	2	14/15	0.56	0.20	150,157,162,163	0
3	NAG	H	2	14/15	0.56	0.14	136,162,169,169	0
5	MAN	I	4	11/12	0.66	0.22	130,140,157,158	0
3	NAG	J	2	14/15	0.71	0.16	133,169,179,183	0

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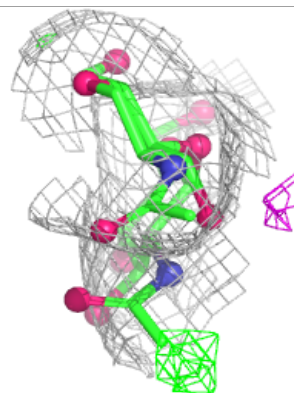
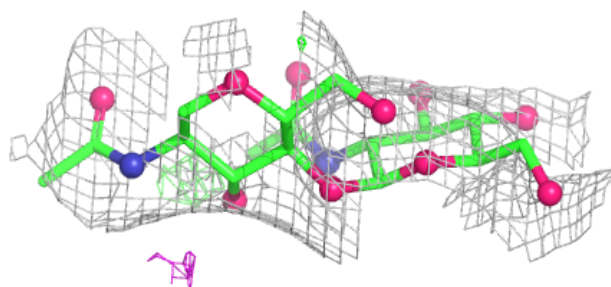
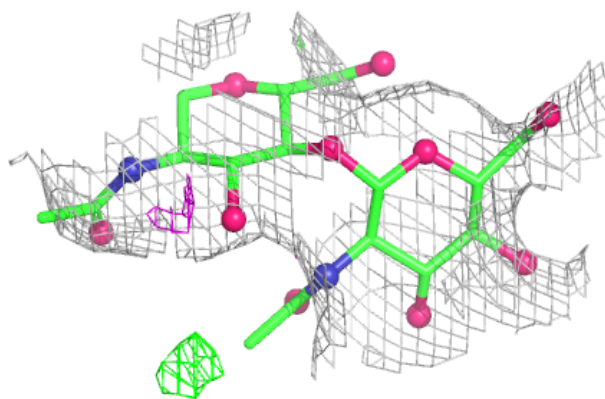
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	BMA	I	3	11/12	0.72	0.17	125,150,166,174	0
3	NAG	G	1	14/15	0.75	0.13	127,140,163,172	0
3	NAG	E	1	14/15	0.84	0.14	88,109,125,136	0
3	NAG	H	1	14/15	0.86	0.12	102,122,145,158	0
4	NAG	F	2	14/15	0.88	0.10	93,137,149,157	0
3	NAG	J	1	14/15	0.89	0.12	89,107,121,145	0
5	NAG	I	2	14/15	0.91	0.10	95,102,123,143	0
4	NAG	F	1	14/15	0.92	0.10	76,97,120,128	0
5	NAG	I	1	14/15	0.95	0.08	63,76,94,99	0

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



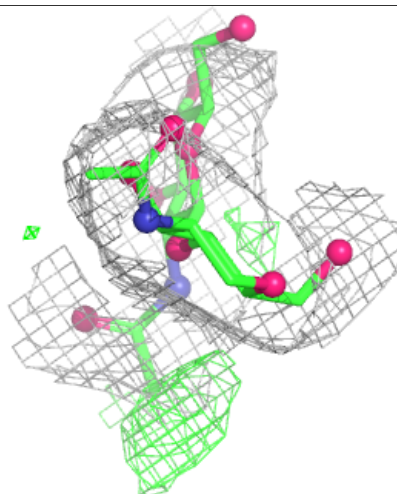
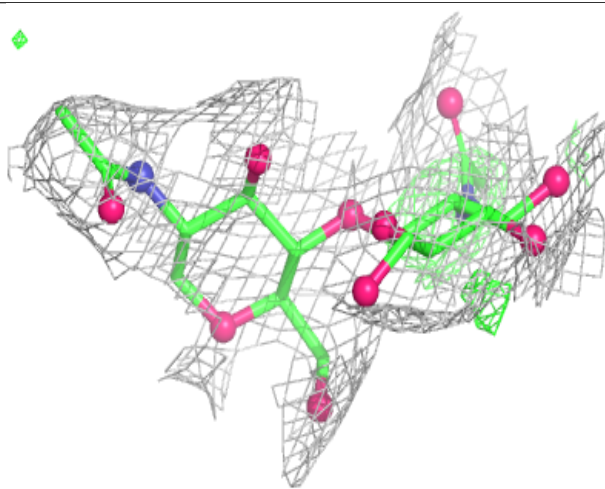
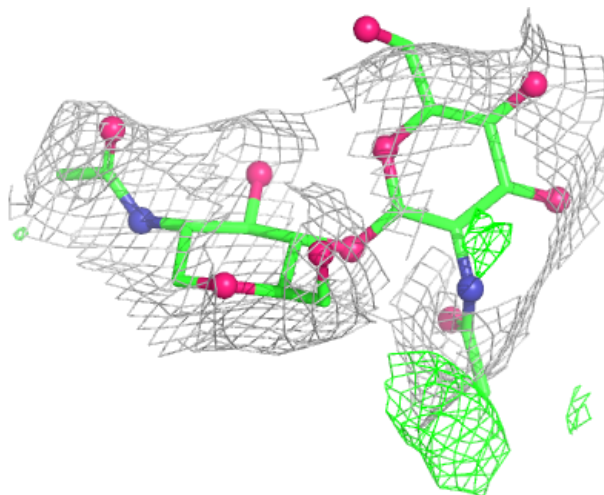
Electron density around Chain G:

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



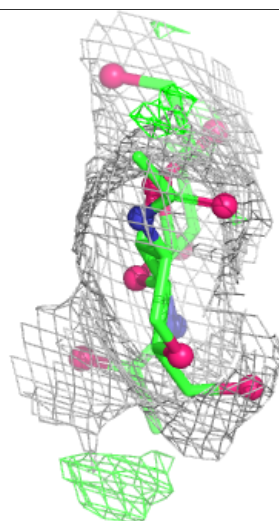
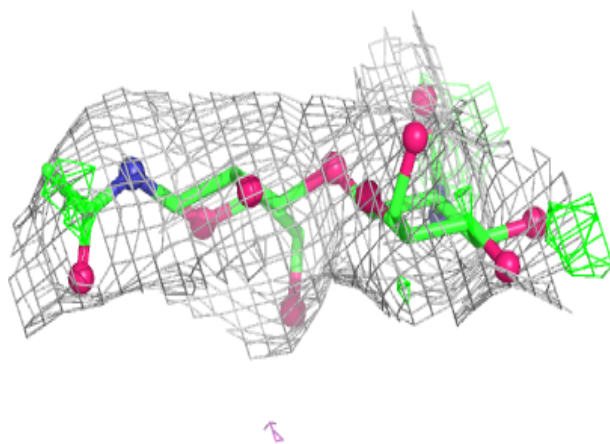
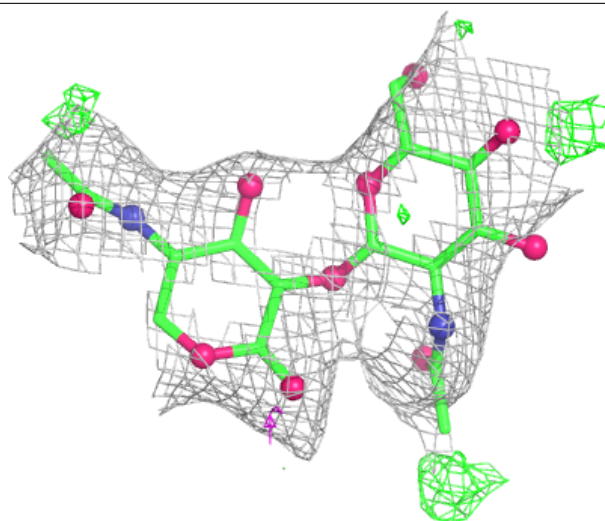
Electron density around Chain H:

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



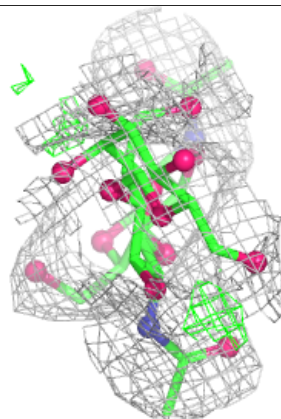
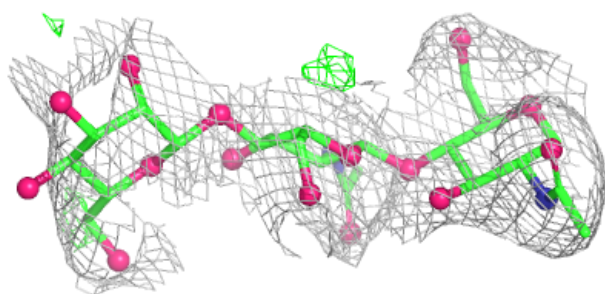
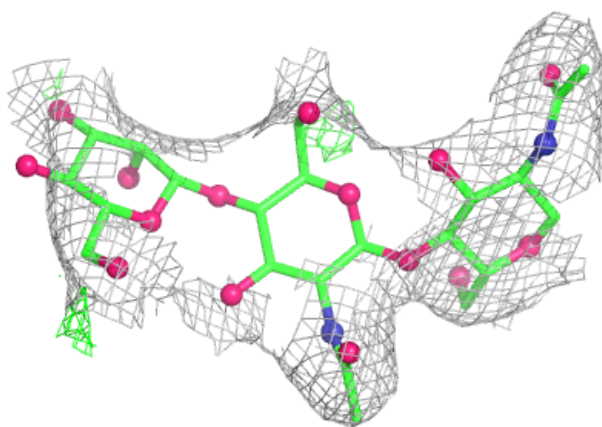
Electron density around Chain J:

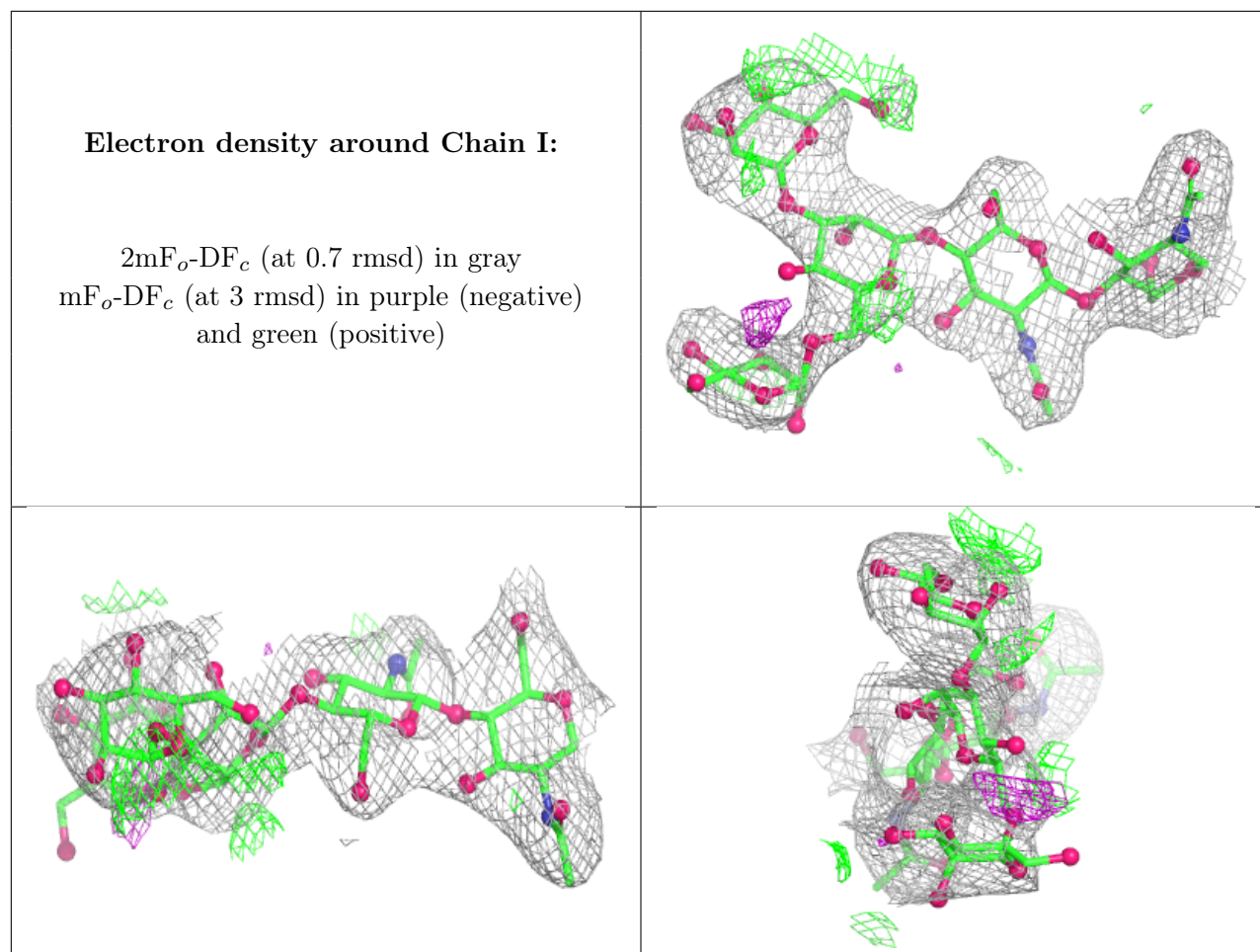
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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)





6.4 Ligands [i](#)

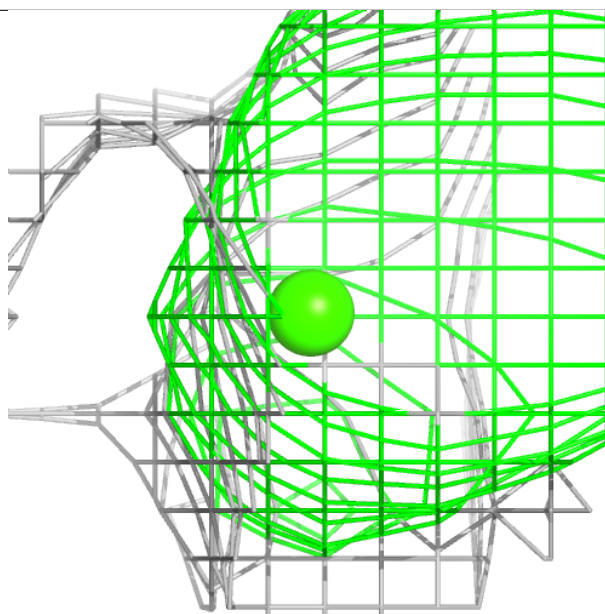
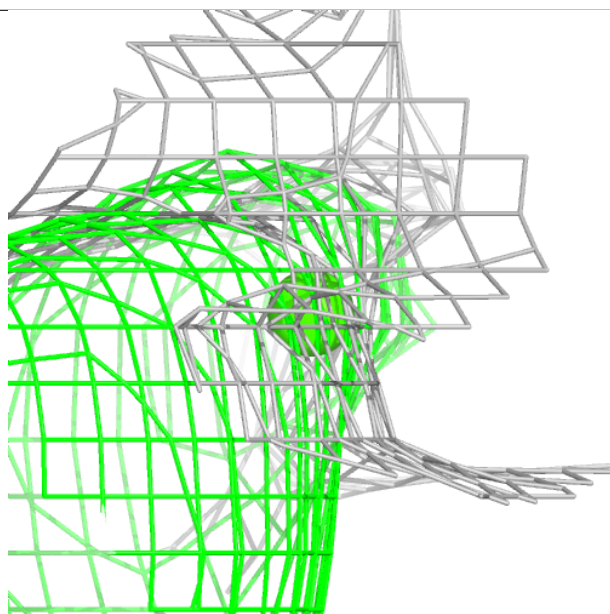
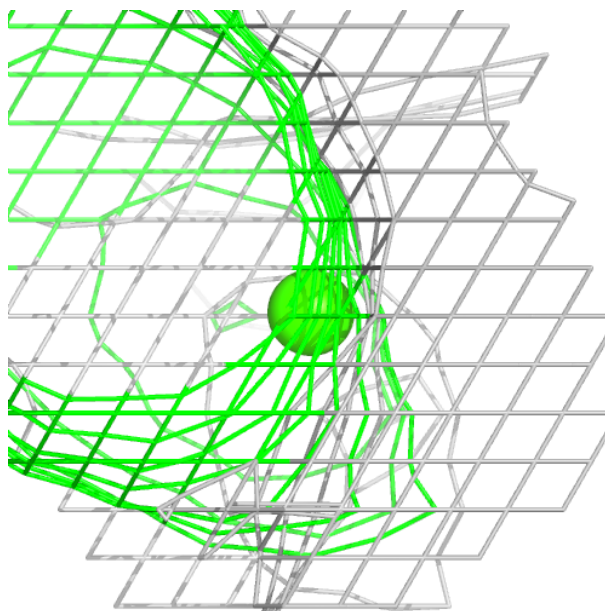
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	CA	A	801	1/1	-	-	62,62,62,62	1
6	CA	B	810	1/1	-	-	60,60,60,60	1

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.

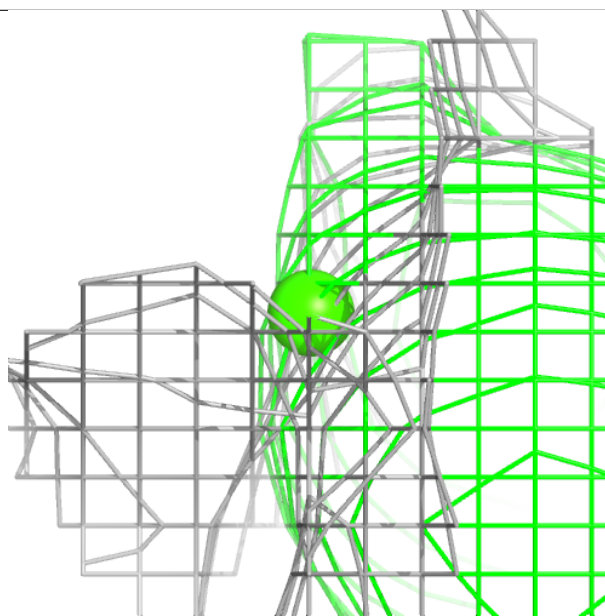
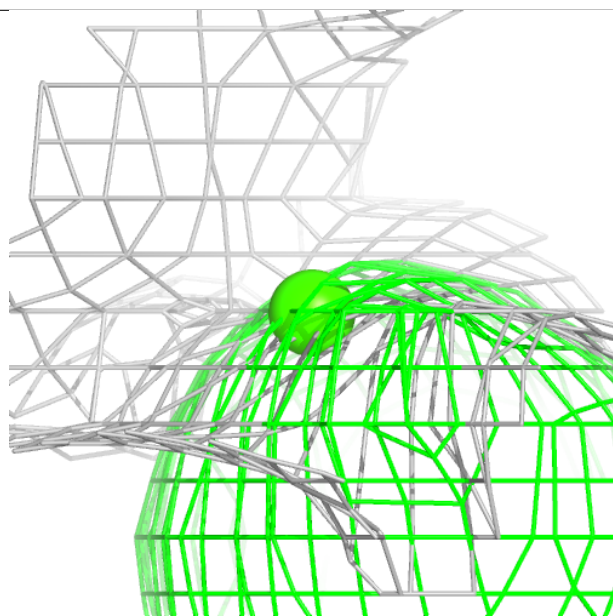
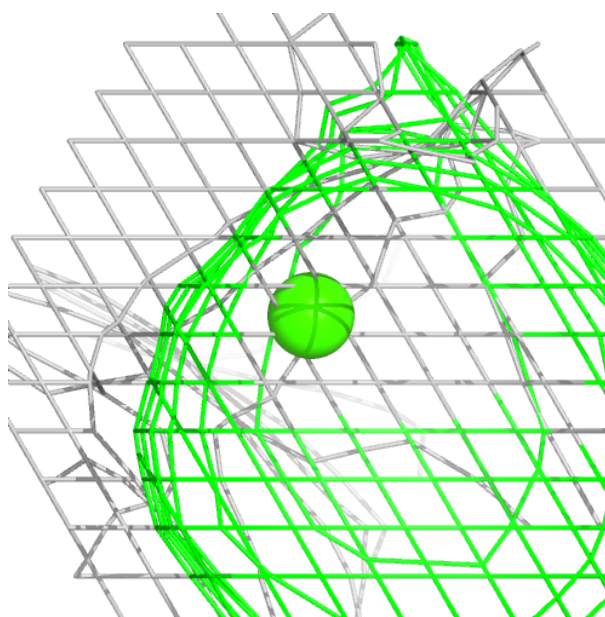
Electron density around CA A 801:

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



Electron density around CA B 810:

$2mF_o-DF_c$ (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.