



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 13, 2024 – 06:35 AM EST

PDB ID : 4MMY
Title : Integrin AlphaVBeta3 ectodomain bound to the tenth domain of Fibronectin with the IAKGDWND motif
Authors : van Agthoven, J.; Xiong, J.; Arnaout, M.A.
Deposited on : 2013-09-09
Resolution : 3.18 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)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

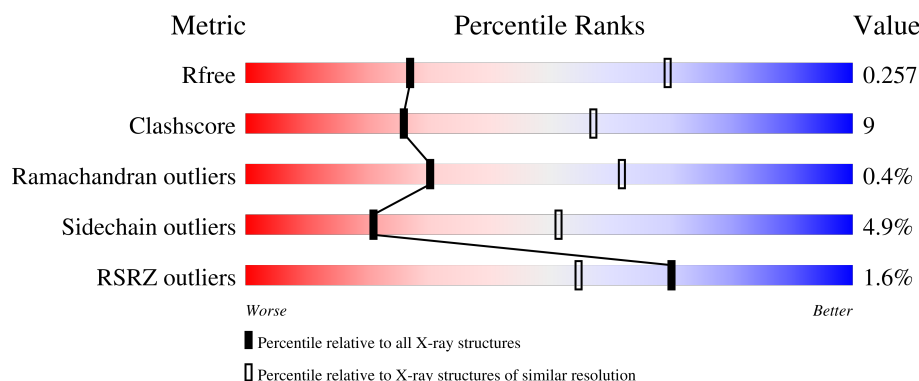
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.18 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}	164625	1851 (3.20-3.16)
Clashscore	180529	1999 (3.20-3.16)
Ramachandran outliers	177936	1961 (3.20-3.16)
Sidechain outliers	177891	1960 (3.20-3.16)
RSRZ outliers	164620	1852 (3.20-3.16)

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	959	
2	B	692	
3	C	98	
4	D	4	
4	G	4	

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Mol	Chain	Length	Quality of chain
5	E	2	 100%
5	H	2	 100%
5	I	2	 50%  50%
5	K	2	 50%  50%
6	F	6	 50%  50%
7	J	3	 33%  67%
7	L	3	 33%  67%

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 13000 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 Integrin alpha-V.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	924	Total	C	N	O	S	0	0	0
			7196	4556	1221	1384	35			

- Molecule 2 is a protein called Integrin beta-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	690	Total	C	N	O	S	0	0	0
			5294	3250	904	1070	70			

- Molecule 3 is a protein called Fibronectin.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	8	Total	C	N	O	0	0	0
			65	41	13	11			

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	1491	ILE	THR	engineered mutation	UNP P02751
C	1492	ALA	GLY	engineered mutation	UNP P02751
C	1496	TRP	SER	engineered mutation	UNP P02751
C	1497	ASN	PRO	engineered mutation	UNP P02751
C	1498	ASP	ALA	engineered mutation	UNP P02751
C	1499	GLY	SER	engineered mutation	UNP P02751
C	1510	GLY	-	expression tag	UNP P02751
C	1511	LYS	-	expression tag	UNP P02751
C	1512	LYS	-	expression tag	UNP P02751
C	1513	GLY	-	expression tag	UNP P02751
C	1514	LYS	-	expression tag	UNP P02751

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-beta-D-mannopyranos e-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucos e

pyranose.



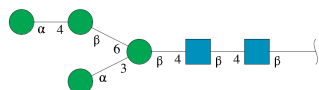
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	D	4	Total	C	N	O	0	0	0
			50	28	2	20			
4	G	4	Total	C	N	O	0	0	0
			50	28	2	20			

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



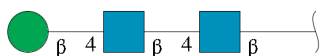
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	E	2	Total	C	N	O	0	0	0
			28	16	2	10			
5	H	2	Total	C	N	O	0	0	0
			28	16	2	10			
5	I	2	Total	C	N	O	0	0	0
			28	16	2	10			
5	K	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 6 is an oligosaccharide called alpha-D-mannopyranose-(1-4)-beta-D-mannopyranose-(1-6)-[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.



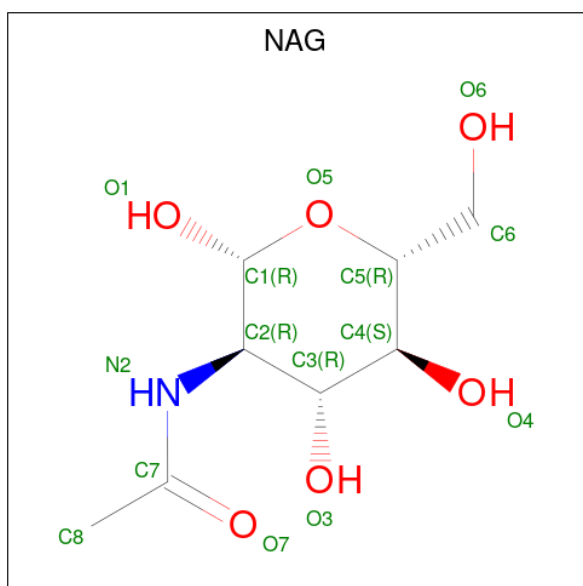
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	F	6	Total	C	N	O	0	0	0
			72	40	2	30			

- Molecule 7 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
7	J	3	Total	C	N	O	0	0	0
			39	22	2	15			
7	L	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 8 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 9 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	5	Total 5	Mn 5	0	0
9	B	3	Total 3	Mn 3	0	0

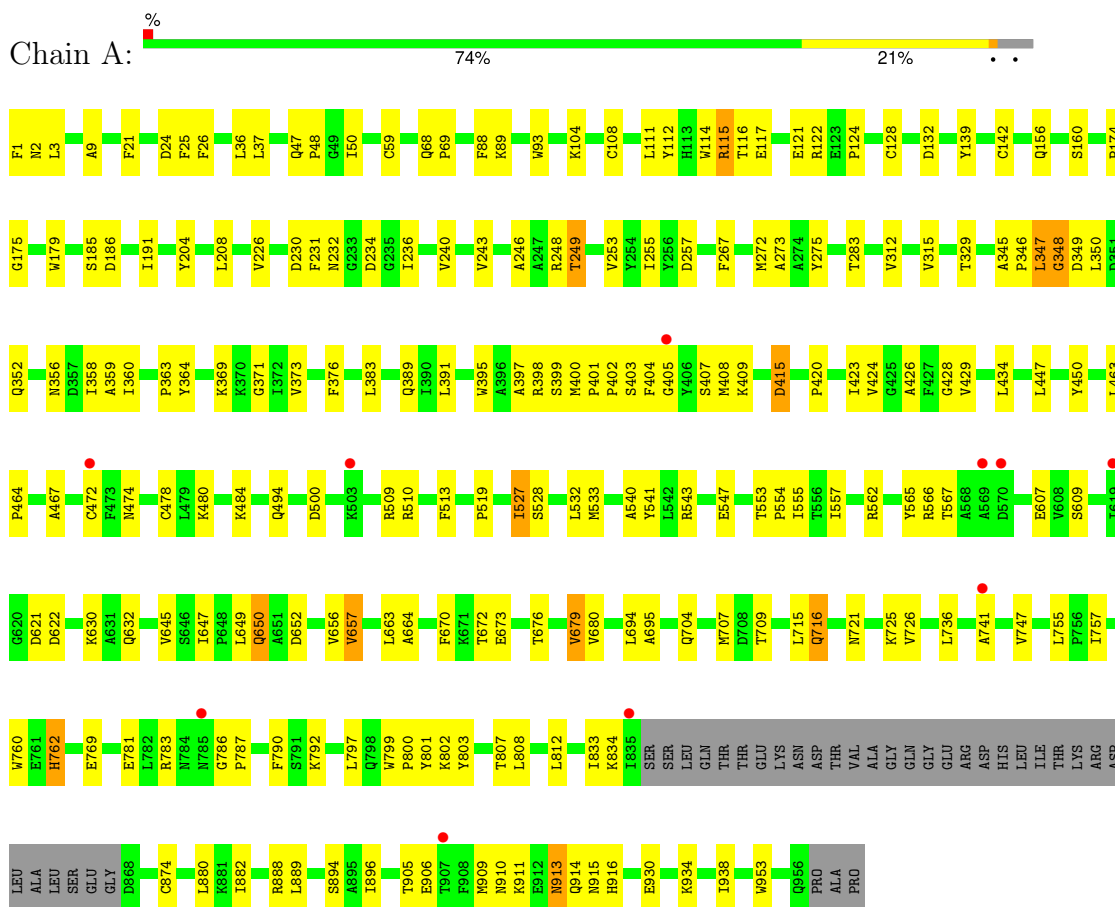
- Molecule 10 is water.

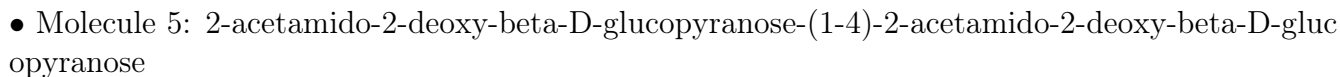
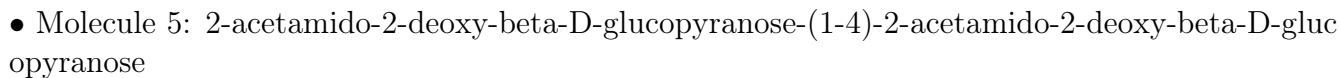
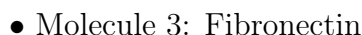
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	3	Total 3	O 3	0	0
10	B	2	Total 2	O 2	0	0

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: Integrin alpha-V





Chain H:  100%

MAG1
MAG2

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

Chain I:  50% 50%

MAG1
MAG2

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

Chain K:  50% 50%

MAG1
MAG2

- Molecule 6: alpha-D-mannopyranose-(1-4)-beta-D-mannopyranose-(1-6)-[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 F:  50% 50%

MAG1
MAG2
BMA3
BMA4
MAN5
MAN6

- Molecule 7: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  33% 67%

MAG1
MAG2
BMA3

- Molecule 7: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain L:  33% 67%

MAG1
MAG2
BMA3

4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	130.02Å 130.02Å 308.20Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	41.02 – 3.18 41.02 – 3.18	Depositor EDS
% Data completeness (in resolution range)	99.4 (41.02-3.18) 99.4 (41.02-3.18)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.27 (at 3.19Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.211 , 0.252 0.217 , 0.257	Depositor DCC
R_{free} test set	2535 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	95.8	Xtriage
Anisotropy	0.269	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 92.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.028 for -h,-k,l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	13000	wwPDB-VP
Average B, all atoms (Å ²)	117.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.60% 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: MAN, MN, BMA, NAG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.23	0/7352	0.44	0/9967
2	B	0.22	0/5390	0.47	0/7289
3	C	0.17	0/66	0.49	0/89
All	All	0.23	0/12808	0.45	0/17345

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	7196	0	7013	129	0
2	B	5294	0	5024	109	0
3	C	65	0	60	1	0
4	D	50	0	43	0	0
4	G	50	0	43	0	0
5	E	28	0	25	0	0
5	H	28	0	25	1	0
5	I	28	0	25	0	0
5	K	28	0	25	2	0
6	F	72	0	61	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	J	39	0	34	0	0
7	L	39	0	34	0	0
8	A	42	0	39	0	0
8	B	28	0	26	2	0
9	A	5	0	0	0	0
9	B	3	0	0	0	0
10	A	3	0	0	0	0
10	B	2	0	0	0	0
All	All	13000	0	12477	237	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 9.

All (237) 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:346:PRO:C	1:A:347:LEU:HD23	1.66	1.15
1:A:347:LEU:HD23	1:A:347:LEU:N	1.50	1.08
2:B:27:SER:O	2:B:28:ASP:HB2	1.30	1.07
2:B:30:ALA:O	2:B:31:LEU:HD23	1.61	1.00
1:A:347:LEU:N	1:A:347:LEU:CD2	2.30	0.94
2:B:27:SER:O	2:B:28:ASP:CB	2.16	0.91
2:B:27:SER:HB2	2:B:53:SER:HB2	1.50	0.91
2:B:30:ALA:O	2:B:31:LEU:HB3	1.79	0.83
2:B:28:ASP:O	2:B:29:GLU:HB2	1.77	0.83
2:B:30:ALA:O	2:B:31:LEU:CB	2.30	0.80
2:B:510:SER:HB3	2:B:511:SER:HA	1.65	0.78
2:B:31:LEU:HD12	2:B:31:LEU:O	1.84	0.77
2:B:30:ALA:O	2:B:31:LEU:CD2	2.32	0.76
2:B:455:PHE:HA	2:B:460:CYS:HA	1.69	0.74
1:A:415:ASP:OD1	1:A:415:ASP:N	2.16	0.73
1:A:707:MET:HG2	1:A:934:LYS:H	1.53	0.73
1:A:609:SER:HB2	1:A:630:LYS:HB3	1.70	0.73
1:A:348:GLY:O	1:A:356:ASN:CB	2.36	0.73
2:B:141:GLN:HE21	2:B:345:VAL:HG23	1.53	0.73
1:A:664:ALA:HB3	1:A:695:ALA:HB2	1.73	0.71
1:A:348:GLY:O	1:A:356:ASN:HB3	1.90	0.70
2:B:616:GLU:HG3	2:B:622:ARG:HB3	1.75	0.68
2:B:169:PRO:HD2	2:B:172:ALA:HB2	1.76	0.68
1:A:645:VAL:HB	1:A:679:VAL:HG13	1.78	0.66
1:A:402:PRO:HA	1:A:428:GLY:HA3	1.78	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:498:ARG:HD3	2:B:516:ILE:HD13	1.77	0.66
1:A:371:GLY:HA3	1:A:404:PHE:HB3	1.78	0.64
1:A:565:TYR:O	1:A:566:ARG:NH1	2.28	0.64
1:A:741:ALA:H	1:A:786:GLY:HA3	1.61	0.64
2:B:417:LYS:NZ	2:B:421:PHE:O	2.29	0.64
2:B:404:ARG:NH1	2:B:550:ASP:OD1	2.28	0.64
1:A:650:GLN:HE21	1:A:704:GLN:HB2	1.62	0.64
2:B:487:SER:HB2	2:B:492:GLN:HB3	1.80	0.63
2:B:292:LEU:HD22	2:B:325:ILE:HD11	1.81	0.63
1:A:348:GLY:O	1:A:356:ASN:HA	2.00	0.61
2:B:28:ASP:O	2:B:29:GLU:CB	2.48	0.61
1:A:348:GLY:O	1:A:356:ASN:CA	2.49	0.61
1:A:725:LYS:HG3	1:A:726:VAL:HG13	1.81	0.61
2:B:473:CYS:SG	2:B:474:SER:N	2.71	0.60
2:B:383:LEU:HG	2:B:385:SER:H	1.66	0.60
1:A:248:ARG:NH1	1:A:272:MET:SD	2.74	0.60
1:A:346:PRO:C	1:A:347:LEU:CD2	2.58	0.60
1:A:345:ALA:O	1:A:347:LEU:HD21	2.02	0.59
1:A:402:PRO:HB3	1:A:429:VAL:HG13	1.85	0.59
2:B:30:ALA:O	2:B:31:LEU:CG	2.51	0.59
2:B:137:LYS:HB3	2:B:341:LEU:HD21	1.85	0.59
2:B:318:TYR:HA	2:B:321:TYR:HB2	1.85	0.58
1:A:3:LEU:HG	1:A:350:LEU:HD21	1.85	0.58
1:A:802:LYS:HG2	1:A:807:THR:HA	1.84	0.58
2:B:129:TRP:HZ3	2:B:339:ASN:HD21	1.52	0.58
2:B:134:LEU:O	2:B:137:LYS:N	2.37	0.58
1:A:395:TRP:HB3	1:A:429:VAL:HG11	1.86	0.57
1:A:543:ARG:NH1	1:A:547:GLU:OE1	2.34	0.57
2:B:411:GLU:HG2	2:B:430:THR:HG22	1.86	0.57
2:B:99:ASN:HD22	8:B:701:NAG:H62	1.69	0.57
1:A:909:MET:HA	1:A:914:GLN:HA	1.87	0.56
1:A:645:VAL:HG22	1:A:715:LEU:HD22	1.88	0.56
2:B:31:LEU:O	2:B:31:LEU:CD1	2.52	0.55
1:A:114:TRP:CD2	1:A:116:THR:HA	2.42	0.55
1:A:114:TRP:CE2	1:A:116:THR:HA	2.41	0.55
2:B:674:SER:HB3	2:B:676:LYS:HG3	1.89	0.55
1:A:397:ALA:HB2	1:A:402:PRO:HD3	1.90	0.54
1:A:913:ASN:O	1:A:913:ASN:ND2	2.40	0.54
2:B:31:LEU:O	2:B:31:LEU:CG	2.55	0.54
2:B:230:THR:HG23	2:B:304:ILE:HD12	1.89	0.54
2:B:288:ASP:OD1	2:B:289:TYR:N	2.40	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:510:SER:H	2:B:511:SER:HB3	1.73	0.54
1:A:450:TYR:HB2	1:A:474:ASN:HB2	1.89	0.54
1:A:253:VAL:HB	1:A:267:PHE:HB2	1.90	0.54
2:B:29:GLU:O	2:B:30:ALA:C	2.47	0.53
1:A:621:ASP:OD1	1:A:622:ASP:N	2.40	0.53
2:B:608:CYS:O	2:B:612:LYS:N	2.41	0.53
2:B:31:LEU:H	2:B:32:PRO:HD3	1.73	0.53
2:B:459:VAL:HG12	2:B:461:ARG:H	1.74	0.53
1:A:315:VAL:HG21	1:A:360:ILE:HD13	1.91	0.52
2:B:620:PHE:HB3	2:B:622:ARG:HE	1.75	0.52
1:A:124:PRO:HB2	1:A:156:GLN:HG2	1.91	0.52
1:A:915:ASN:HA	1:A:953:TRP:CD1	2.44	0.52
1:A:607:GLU:OE1	1:A:632:GLN:NE2	2.42	0.52
2:B:120:LEU:HD12	2:B:188:PHE:HZ	1.75	0.52
2:B:530:ARG:HG2	2:B:535:MET:HA	1.90	0.52
2:B:144:LYS:HG3	2:B:145:LEU:HD13	1.91	0.52
1:A:349:ASP:O	1:A:420:PRO:HG3	2.09	0.52
2:B:580:LYS:HG2	2:B:589:ILE:HD11	1.92	0.52
1:A:906:GLU:O	1:A:910:ASN:ND2	2.42	0.51
2:B:529:VAL:HG11	2:B:556:TYR:HE1	1.76	0.51
1:A:231:PHE:N	1:A:231:PHE:CD2	2.77	0.51
2:B:133:ASN:O	2:B:204:ASN:ND2	2.43	0.51
1:A:801:TYR:HB2	1:A:880:LEU:HB2	1.93	0.51
2:B:494:VAL:HG22	2:B:495:CYS:H	1.75	0.51
1:A:21:PHE:HE1	2:B:266:VAL:HG11	1.75	0.51
1:A:50:ILE:HD12	1:A:89:LYS:HB2	1.93	0.51
2:B:316:ASN:HB3	8:B:702:NAG:H82	1.93	0.51
1:A:345:ALA:O	1:A:347:LEU:CD2	2.59	0.51
1:A:510:ARG:NH2	1:A:553:THR:O	2.25	0.50
2:B:215:ASN:O	3:C:1495:ASP:HB3	2.12	0.50
2:B:50:ALA:N	2:B:51:PRO:HD2	2.27	0.49
1:A:36:LEU:HB2	1:A:59:CYS:HB2	1.94	0.49
1:A:185:SER:HB3	1:A:208:LEU:HB2	1.95	0.49
1:A:104:LYS:HD3	1:A:132:ASP:HB2	1.93	0.49
1:A:649:LEU:HD13	1:A:704:GLN:HG3	1.94	0.49
1:A:24:ASP:OD1	1:A:25:PHE:N	2.45	0.49
1:A:26:PHE:HB2	1:A:37:LEU:HG	1.93	0.49
1:A:232:ASN:ND2	1:A:257:ASP:OD1	2.45	0.49
1:A:513:PHE:HA	1:A:540:ALA:HA	1.95	0.49
1:A:909:MET:HA	1:A:914:GLN:HG3	1.94	0.49
1:A:408:MET:HG2	1:A:424:VAL:HG22	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:PHE:HA	1:A:389:GLN:HB2	1.95	0.49
2:B:362:LEU:HD12	2:B:363:PRO:HD2	1.94	0.48
2:B:442:GLU:HG3	2:B:443:PRO:HD2	1.95	0.48
1:A:423:ILE:HG12	1:A:434:LEU:HD23	1.94	0.48
2:B:400:GLU:HB2	5:K:1:NAG:H83	1.94	0.48
2:B:358:GLU:N	2:B:417:LYS:O	2.42	0.48
2:B:124:MET:HG2	2:B:251:ASP:HB2	1.95	0.48
1:A:249:THR:HG23	1:A:273:ALA:H	1.79	0.48
1:A:349:ASP:O	1:A:420:PRO:CG	2.61	0.48
1:A:2:ASN:OD1	1:A:2:ASN:N	2.47	0.47
2:B:338:SER:HB3	2:B:342:GLN:HB2	1.96	0.47
2:B:479:ARG:O	2:B:481:SER:N	2.45	0.47
2:B:608:CYS:HB3	2:B:655:CYS:HB3	1.58	0.47
5:K:1:NAG:H62	5:K:2:NAG:H83	1.96	0.47
2:B:116:TYR:HA	2:B:247:VAL:HG13	1.97	0.47
1:A:480:LYS:HB3	1:A:533:MET:HG2	1.96	0.47
2:B:133:ASN:ND2	2:B:137:LYS:HE3	2.29	0.47
2:B:66:ASP:HB3	2:B:85:PRO:HB3	1.97	0.47
2:B:319:GLN:HA	2:B:330:VAL:HG21	1.96	0.47
2:B:134:LEU:O	2:B:138:LEU:N	2.39	0.46
1:A:792:LYS:HB2	1:A:930:GLU:HG2	1.97	0.46
1:A:405:GLY:C	1:A:407:SER:H	2.19	0.46
1:A:243:VAL:HG22	1:A:246:ALA:HB2	1.97	0.46
2:B:158:ASP:HB3	2:B:187:MET:HE1	1.98	0.46
1:A:447:LEU:HD21	1:A:557:ILE:HG22	1.98	0.46
2:B:157:VAL:O	2:B:220:GLU:HG3	2.16	0.46
1:A:108:CYS:HA	1:A:128:CYS:HA	1.98	0.46
2:B:135:GLY:HA2	2:B:138:LEU:HB2	1.97	0.46
1:A:24:ASP:HA	1:A:409:LYS:HG2	1.98	0.45
1:A:400:MET:HB2	1:A:401:PRO:HD2	1.96	0.45
2:B:137:LYS:O	2:B:140:THR:HG22	2.16	0.45
1:A:527:ILE:HG22	1:A:528:SER:H	1.82	0.45
1:A:803:TYR:HB2	1:A:808:LEU:HD11	1.98	0.45
1:A:139:TYR:OH	1:A:186:ASP:OD2	2.31	0.45
1:A:797:LEU:HD23	1:A:882:ILE:HD12	1.98	0.45
1:A:240:VAL:HG22	1:A:255:ILE:HG12	1.98	0.45
2:B:41:LYS:HD2	2:B:41:LYS:HA	1.69	0.45
2:B:590:GLN:CD	2:B:591:PRO:HD2	2.36	0.45
1:A:9:ALA:HB3	1:A:434:LEU:HB3	1.98	0.45
1:A:652:ASP:OD1	1:A:652:ASP:N	2.50	0.45
1:A:472:CYS:HA	1:A:541:TYR:HA	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:160:SER:HB2	1:A:226:VAL:HG22	1.98	0.45
2:B:305:ASN:HB3	2:B:351:ILE:HD13	1.98	0.45
2:B:371:ASN:HB2	2:B:398:SER:OG	2.17	0.45
1:A:93:TRP:CD1	1:A:111:LEU:HD22	2.52	0.45
2:B:249:THR:HA	2:B:309:ALA:O	2.17	0.45
1:A:494:GLN:HB2	1:A:562:ARG:HB3	1.99	0.45
1:A:769:GLU:HG2	1:A:812:LEU:HD21	1.99	0.44
2:B:133:ASN:HD22	2:B:137:LYS:HE3	1.82	0.44
1:A:352:GLN:O	1:A:484:LYS:NZ	2.49	0.44
2:B:244:HIS:HB2	2:B:304:ILE:HA	1.99	0.44
1:A:350:LEU:HD12	1:A:350:LEU:HA	1.74	0.44
1:A:500:ASP:HB2	1:A:555:ILE:HG23	1.98	0.44
1:A:888:ARG:HA	1:A:888:ARG:HD2	1.73	0.44
1:A:364:TYR:HB3	1:A:369:LYS:HE3	2.00	0.44
2:B:105:ARG:HB2	2:B:394:THR:HG23	1.98	0.44
1:A:398:ARG:HG3	1:A:399:SER:H	1.82	0.44
2:B:125:LYS:HG3	2:B:212:VAL:HG21	1.99	0.44
1:A:175:GLY:HA2	1:A:179:TRP:CD1	2.53	0.44
2:B:21:PRO:HB3	2:B:96:ASP:HB2	2.00	0.44
1:A:799:TRP:HA	1:A:800:PRO:HD3	1.79	0.43
2:B:497:GLN:HB3	2:B:498:ARG:H	1.54	0.43
1:A:762:HIS:ND1	1:A:906:GLU:OE1	2.46	0.43
1:A:783:ARG:HG3	1:A:894:SER:HB2	1.99	0.43
2:B:466:TRP:CD2	2:B:471:CYS:HB3	2.53	0.43
2:B:494:VAL:HG21	2:B:500:GLU:HB2	1.99	0.43
1:A:88:PHE:CE2	1:A:122:ARG:HG2	2.54	0.43
1:A:347:LEU:HD11	1:A:359:ALA:HB2	1.99	0.43
2:B:173:LEU:O	2:B:174:GLU:HB3	2.18	0.43
5:H:1:NAG:H4	5:H:2:NAG:H2	1.83	0.43
1:A:47:GLN:HA	1:A:48:PRO:HD3	1.88	0.43
2:B:462:CYS:HB3	2:B:467:LEU:H	1.83	0.43
2:B:141:GLN:O	2:B:144:LYS:HG2	2.17	0.43
2:B:407:PRO:HB2	2:B:409:GLU:HG2	2.00	0.43
1:A:509:ARG:HH11	1:A:519:PRO:HB2	1.83	0.43
2:B:86:GLN:O	2:B:425:LEU:HD12	2.19	0.43
2:B:131:ILE:HD12	2:B:131:ILE:HA	1.87	0.43
1:A:234:ASP:HB2	1:A:236:ILE:HD12	2.00	0.43
1:A:747:VAL:HG13	2:B:603:THR:HB	2.01	0.43
2:B:39:ASP:HB3	2:B:40:LEU:H	1.59	0.43
1:A:347:LEU:O	1:A:349:ASP:N	2.51	0.43
2:B:70:SER:HB2	2:B:107:VAL:HG11	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:647:ASP:OD1	2:B:647:ASP:N	2.52	0.43
2:B:188:PHE:CZ	2:B:191:LYS:HE3	2.53	0.42
1:A:364:TYR:O	1:A:369:LYS:HG3	2.19	0.42
1:A:115:ARG:O	1:A:115:ARG:HD3	2.19	0.42
2:B:150:ARG:HG3	2:B:239:ARG:CZ	2.49	0.42
2:B:4:ILE:HD12	2:B:4:ILE:H	1.84	0.42
2:B:108:GLU:HG2	2:B:391:ILE:HG22	2.01	0.42
1:A:346:PRO:HA	1:A:358:ILE:HG13	2.02	0.42
2:B:137:LYS:O	2:B:141:GLN:HG2	2.18	0.42
1:A:68:GLN:HA	1:A:69:PRO:HD3	1.92	0.42
1:A:716:GLN:HE21	1:A:716:GLN:HB2	1.62	0.42
2:B:134:LEU:HD12	2:B:134:LEU:HA	1.92	0.42
1:A:174:PRO:HB2	2:B:262:LEU:HD21	2.02	0.42
1:A:709:THR:O	1:A:736:LEU:HB2	2.20	0.42
2:B:71:ASP:OD1	2:B:71:ASP:N	2.45	0.42
1:A:403:SER:O	1:A:426:ALA:HA	2.20	0.41
1:A:621:ASP:HB2	1:A:787:PRO:HB3	2.02	0.41
2:B:80:VAL:HG11	2:B:110:TYR:HE1	1.85	0.41
1:A:230:ASP:OD1	1:A:234:ASP:O	2.38	0.41
2:B:334:SER:OG	2:B:337:SER:OG	2.37	0.41
1:A:911:LYS:C	1:A:913:ASN:H	2.22	0.41
2:B:406:CYS:HA	2:B:407:PRO:HD3	1.81	0.41
1:A:191:ILE:HA	1:A:204:TYR:CE2	2.54	0.41
1:A:464:PRO:HG2	1:A:467:ALA:HB3	2.01	0.41
2:B:141:GLN:HG3	2:B:142:MET:H	1.85	0.41
1:A:760:TRP:CZ2	1:A:905:THR:HB	2.56	0.41
1:A:112:TYR:O	1:A:124:PRO:HA	2.21	0.41
1:A:373:VAL:HB	1:A:391:LEU:HB2	2.02	0.41
1:A:553:THR:HA	1:A:554:PRO:HD3	1.93	0.41
2:B:283:ALA:HB1	2:B:287:MET:HG2	2.02	0.41
1:A:376:PHE:HB3	1:A:383:LEU:HD11	2.02	0.41
1:A:647:ILE:HD11	1:A:670:PHE:HE1	1.86	0.41
1:A:790:PHE:CZ	1:A:889:LEU:HB2	2.55	0.41
1:A:721:ASN:O	1:A:725:LYS:HE2	2.20	0.41
1:A:755:LEU:HD23	1:A:757:ILE:HD13	2.03	0.41
1:A:781:GLU:HG3	1:A:896:ILE:HG13	2.03	0.41
2:B:126:ASP:OD1	2:B:126:ASP:N	2.52	0.41
2:B:495:CYS:HB3	2:B:520:TYR:O	2.21	0.41
1:A:115:ARG:HA	1:A:121:GLU:HB2	2.03	0.41
1:A:657:VAL:HG21	1:A:663:LEU:HD13	2.03	0.41
1:A:232:ASN:OD1	1:A:232:ASN:N	2.53	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:478:CYS:HB3	1:A:533:MET:HB2	2.03	0.40
1:A:363:PRO:HA	1:A:404:PHE:O	2.21	0.40
2:B:117:LEU:HD23	2:B:117:LEU:HA	1.95	0.40
2:B:150:ARG:CZ	2:B:239:ARG:HD3	2.51	0.40
2:B:370:PHE:CG	2:B:384:LYS:HG2	2.56	0.40
1:A:673:GLU:O	1:A:676:THR:OG1	2.31	0.40
1:A:797:LEU:HB3	1:A:882:ILE:HB	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	920/959 (96%)	853 (93%)	65 (7%)	2 (0%)	44	73
2	B	688/692 (99%)	595 (86%)	89 (13%)	4 (1%)	22	55
3	C	6/98 (6%)	5 (83%)	1 (17%)	0	100	100
All	All	1614/1749 (92%)	1453 (90%)	155 (10%)	6 (0%)	30	62

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	28	ASP
2	B	31	LEU
1	A	833	ILE
2	B	457	CYS
1	A	348	GLY
2	B	379	VAL

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	784/813 (96%)	756 (96%)	28 (4%)	30	59
2	B	612/614 (100%)	572 (94%)	40 (6%)	14	42
3	C	6/81 (7%)	5 (83%)	1 (17%)	2	8
All	All	1402/1508 (93%)	1333 (95%)	69 (5%)	21	51

All (69) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	115	ARG
1	A	117	GLU
1	A	142	CYS
1	A	249	THR
1	A	275	TYR
1	A	283	THR
1	A	312	VAL
1	A	329	THR
1	A	347	LEU
1	A	415	ASP
1	A	463	LEU
1	A	527	ILE
1	A	532	LEU
1	A	567	THR
1	A	650	GLN
1	A	656	VAL
1	A	657	VAL
1	A	672	THR
1	A	679	VAL
1	A	680	VAL
1	A	694	LEU
1	A	716	GLN
1	A	762	HIS
1	A	834	LYS
1	A	874	CYS
1	A	913	ASN

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Mol	Chain	Res	Type
1	A	916	HIS
1	A	938	ILE
2	B	31	LEU
2	B	40	LEU
2	B	42	GLU
2	B	44	LEU
2	B	116	TYR
2	B	128	LEU
2	B	140	THR
2	B	168	SER
2	B	179	ASP
2	B	182	THR
2	B	212	VAL
2	B	215	ASN
2	B	220	GLU
2	B	316	ASN
2	B	321	TYR
2	B	340	VAL
2	B	342	GLN
2	B	376	ASN
2	B	448	CYS
2	B	460	CYS
2	B	462	CYS
2	B	466	TRP
2	B	467	LEU
2	B	470	GLN
2	B	472	GLU
2	B	485	GLU
2	B	501	CYS
2	B	509	HIS
2	B	512	ASP
2	B	561	THR
2	B	562	THR
2	B	589	ILE
2	B	590	GLN
2	B	608	CYS
2	B	622	ARG
2	B	645	LEU
2	B	647	ASP
2	B	648	THR
2	B	653	VAL
2	B	662	ASP

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Mol	Chain	Res	Type
3	C	1493	ARG

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

Mol	Chain	Res	Type
1	A	650	GLN
2	B	141	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 ⓘ

28 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	D	1	1,4	14,14,15	0.53	0	17,19,21	0.48	0
4	NAG	D	2	4	14,14,15	0.18	0	17,19,21	0.41	0
4	BMA	D	3	4	11,11,12	0.59	0	15,15,17	0.70	0
4	MAN	D	4	4	11,11,12	0.69	0	15,15,17	0.98	2 (13%)
5	NAG	E	1	1,5	14,14,15	0.26	0	17,19,21	0.43	0
5	NAG	E	2	5	14,14,15	0.24	0	17,19,21	0.43	0
6	NAG	F	1	6,1	14,14,15	0.30	0	17,19,21	0.41	0
6	NAG	F	2	6	14,14,15	0.21	0	17,19,21	0.46	0
6	BMA	F	3	6	11,11,12	0.67	0	15,15,17	0.70	0
6	BMA	F	4	6	11,11,12	1.31	3 (27%)	15,15,17	1.10	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	MAN	F	5	6	11,11,12	0.70	0	15,15,17	0.94	2 (13%)
6	MAN	F	6	6	11,11,12	0.66	0	15,15,17	0.99	2 (13%)
4	NAG	G	1	1,4	14,14,15	0.36	0	17,19,21	0.42	0
4	NAG	G	2	4	14,14,15	0.32	0	17,19,21	0.42	0
4	BMA	G	3	4	11,11,12	0.78	0	15,15,17	0.89	1 (6%)
4	MAN	G	4	4	11,11,12	0.79	0	15,15,17	0.93	1 (6%)
5	NAG	H	1	1,5	14,14,15	0.20	0	17,19,21	0.46	0
5	NAG	H	2	5	14,14,15	0.26	0	17,19,21	0.40	0
5	NAG	I	1	1,5	14,14,15	0.29	0	17,19,21	0.51	0
5	NAG	I	2	5	14,14,15	0.52	0	17,19,21	0.69	1 (5%)
7	NAG	J	1	1,7	14,14,15	0.25	0	17,19,21	0.46	0
7	NAG	J	2	7	14,14,15	0.75	1 (7%)	17,19,21	0.94	1 (5%)
7	BMA	J	3	7	11,11,12	1.18	2 (18%)	15,15,17	1.57	4 (26%)
5	NAG	K	1	2,5	14,14,15	0.24	0	17,19,21	0.41	0
5	NAG	K	2	5	14,14,15	0.72	1 (7%)	17,19,21	0.92	1 (5%)
7	NAG	L	1	2,7	14,14,15	0.28	0	17,19,21	0.48	0
7	NAG	L	2	7	14,14,15	0.70	1 (7%)	17,19,21	0.90	1 (5%)
7	BMA	L	3	7	11,11,12	0.76	0	15,15,17	1.11	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	D	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	D	2	4	-	4/6/23/26	0/1/1/1
4	BMA	D	3	4	-	0/2/19/22	0/1/1/1
4	MAN	D	4	4	-	0/2/19/22	0/1/1/1
5	NAG	E	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	E	2	5	-	2/6/23/26	0/1/1/1
6	NAG	F	1	6,1	-	1/6/23/26	0/1/1/1
6	NAG	F	2	6	-	0/6/23/26	0/1/1/1
6	BMA	F	3	6	-	2/2/19/22	0/1/1/1
6	BMA	F	4	6	-	0/2/19/22	0/1/1/1
6	MAN	F	5	6	-	0/2/19/22	0/1/1/1
6	MAN	F	6	6	-	1/2/19/22	0/1/1/1
4	NAG	G	1	1,4	-	1/6/23/26	0/1/1/1
4	NAG	G	2	4	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	BMA	G	3	4	-	1/2/19/22	0/1/1/1
4	MAN	G	4	4	-	2/2/19/22	0/1/1/1
5	NAG	H	1	1,5	-	2/6/23/26	0/1/1/1
5	NAG	H	2	5	-	1/6/23/26	0/1/1/1
5	NAG	I	1	1,5	-	1/6/23/26	0/1/1/1
5	NAG	I	2	5	-	2/6/23/26	0/1/1/1
7	NAG	J	1	1,7	-	0/6/23/26	0/1/1/1
7	NAG	J	2	7	-	2/6/23/26	0/1/1/1
7	BMA	J	3	7	-	1/2/19/22	0/1/1/1
5	NAG	K	1	2,5	-	1/6/23/26	0/1/1/1
5	NAG	K	2	5	-	4/6/23/26	0/1/1/1
7	NAG	L	1	2,7	-	3/6/23/26	0/1/1/1
7	NAG	L	2	7	-	2/6/23/26	0/1/1/1
7	BMA	L	3	7	-	1/2/19/22	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	J	3	BMA	C2-C3	2.78	1.56	1.52
6	F	4	BMA	C4-C5	2.66	1.58	1.53
5	K	2	NAG	O5-C1	2.43	1.47	1.43
7	J	3	BMA	C1-C2	2.29	1.57	1.52
6	F	4	BMA	O5-C1	-2.28	1.39	1.43
7	L	2	NAG	O5-C1	2.27	1.47	1.43
7	J	2	NAG	O5-C1	2.21	1.47	1.43
6	F	4	BMA	C4-C3	2.02	1.57	1.52

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	J	3	BMA	C1-C2-C3	3.68	115.00	109.64
5	K	2	NAG	C1-O5-C5	3.52	116.90	112.19
7	J	2	NAG	C1-O5-C5	3.47	116.83	112.19
7	L	2	NAG	C1-O5-C5	3.42	116.77	112.19
7	L	3	BMA	C1-O5-C5	2.76	115.89	112.19
5	I	2	NAG	C1-O5-C5	2.54	115.59	112.19
7	J	3	BMA	C2-C3-C4	2.34	114.97	110.86
6	F	4	BMA	C3-C4-C5	2.26	114.34	110.23
4	G	4	MAN	O2-C2-C3	-2.26	105.47	110.15
4	D	4	MAN	O2-C2-C3	-2.24	105.51	110.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	J	3	BMA	O5-C5-C4	-2.23	105.41	110.83
6	F	6	MAN	C1-O5-C5	2.21	115.14	112.19
6	F	5	MAN	O2-C2-C3	-2.18	105.63	110.15
6	F	6	MAN	O2-C2-C3	-2.14	105.71	110.15
7	J	3	BMA	O5-C1-C2	2.10	115.81	110.79
4	G	3	BMA	O3-C3-C2	2.09	114.32	110.05
4	D	4	MAN	C1-O5-C5	2.05	114.94	112.19
6	F	5	MAN	C1-O5-C5	2.05	114.94	112.19
6	F	4	BMA	C2-C3-C4	2.01	114.39	110.86

There are no chirality outliers.

All (38) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	K	2	NAG	C4-C5-C6-O6
4	G	2	NAG	O5-C5-C6-O6
5	K	2	NAG	O5-C5-C6-O6
7	L	1	NAG	O5-C5-C6-O6
7	L	2	NAG	C4-C5-C6-O6
4	D	2	NAG	O5-C5-C6-O6
4	G	2	NAG	C4-C5-C6-O6
6	F	3	BMA	O5-C5-C6-O6
4	D	2	NAG	C4-C5-C6-O6
4	G	4	MAN	O5-C5-C6-O6
7	J	2	NAG	O5-C5-C6-O6
7	L	2	NAG	O5-C5-C6-O6
7	L	1	NAG	C4-C5-C6-O6
7	J	2	NAG	C4-C5-C6-O6
4	D	2	NAG	C8-C7-N2-C2
4	D	2	NAG	O7-C7-N2-C2
5	E	2	NAG	C8-C7-N2-C2
5	E	2	NAG	O7-C7-N2-C2
5	I	2	NAG	C8-C7-N2-C2
5	I	2	NAG	O7-C7-N2-C2
5	K	2	NAG	C8-C7-N2-C2
5	K	2	NAG	O7-C7-N2-C2
4	D	1	NAG	O5-C5-C6-O6
5	H	1	NAG	O5-C5-C6-O6
4	G	4	MAN	C4-C5-C6-O6
4	G	1	NAG	O5-C5-C6-O6
7	L	3	BMA	O5-C5-C6-O6
5	K	1	NAG	O5-C5-C6-O6

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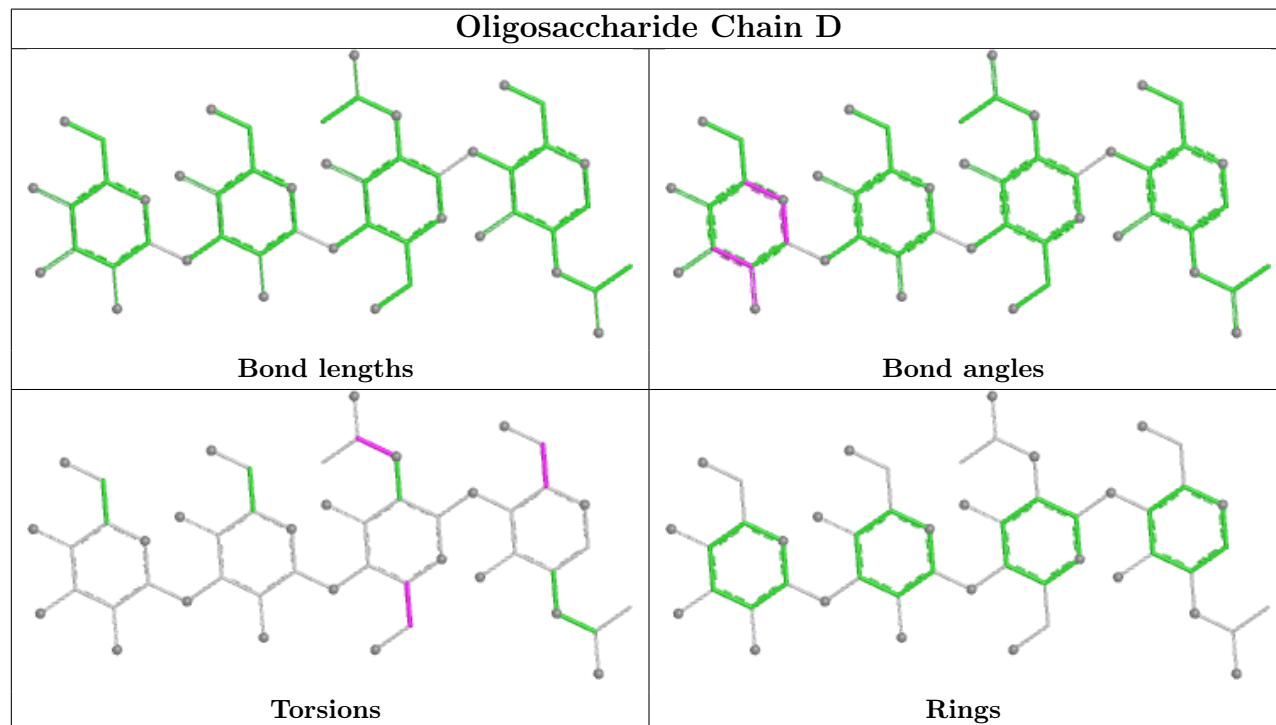
Mol	Chain	Res	Type	Atoms
4	G	3	BMA	O5-C5-C6-O6
6	F	6	MAN	O5-C5-C6-O6
7	J	3	BMA	O5-C5-C6-O6
4	D	1	NAG	C4-C5-C6-O6
5	H	1	NAG	C4-C5-C6-O6
6	F	3	BMA	C4-C5-C6-O6
5	I	1	NAG	O5-C5-C6-O6
5	H	2	NAG	O5-C5-C6-O6
6	F	1	NAG	C1-C2-N2-C7
7	L	1	NAG	C1-C2-N2-C7

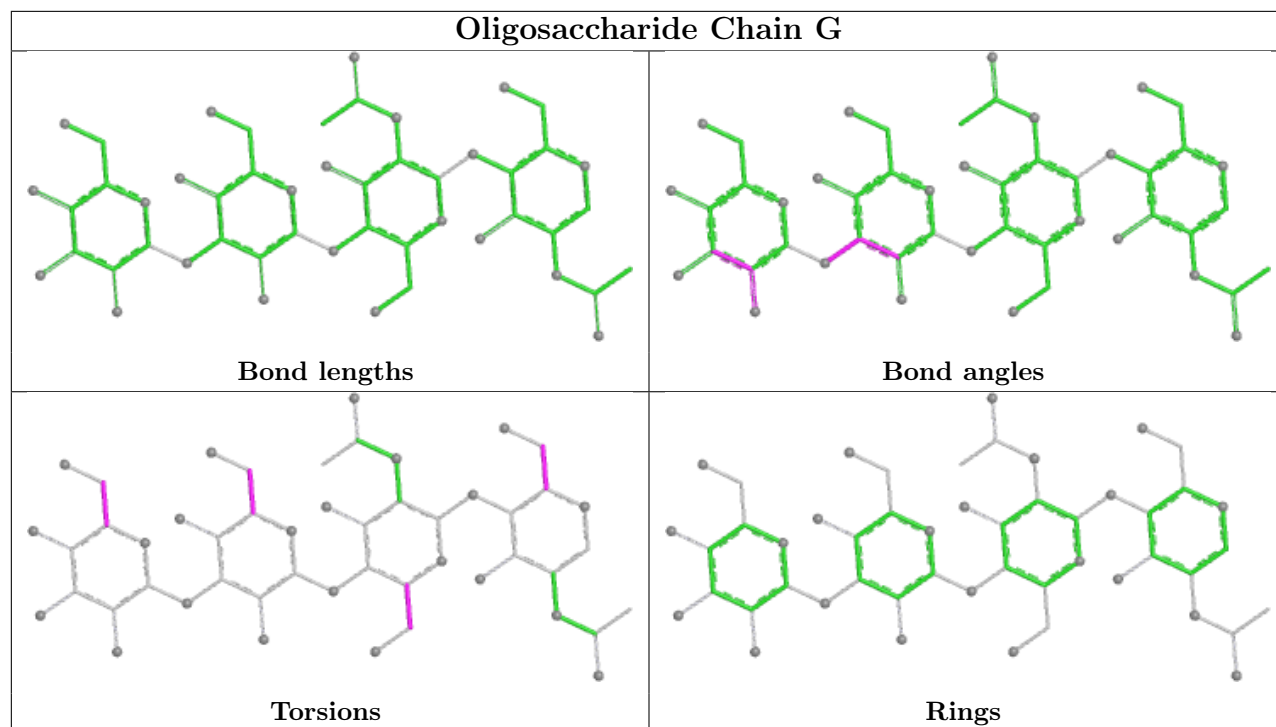
There are no ring outliers.

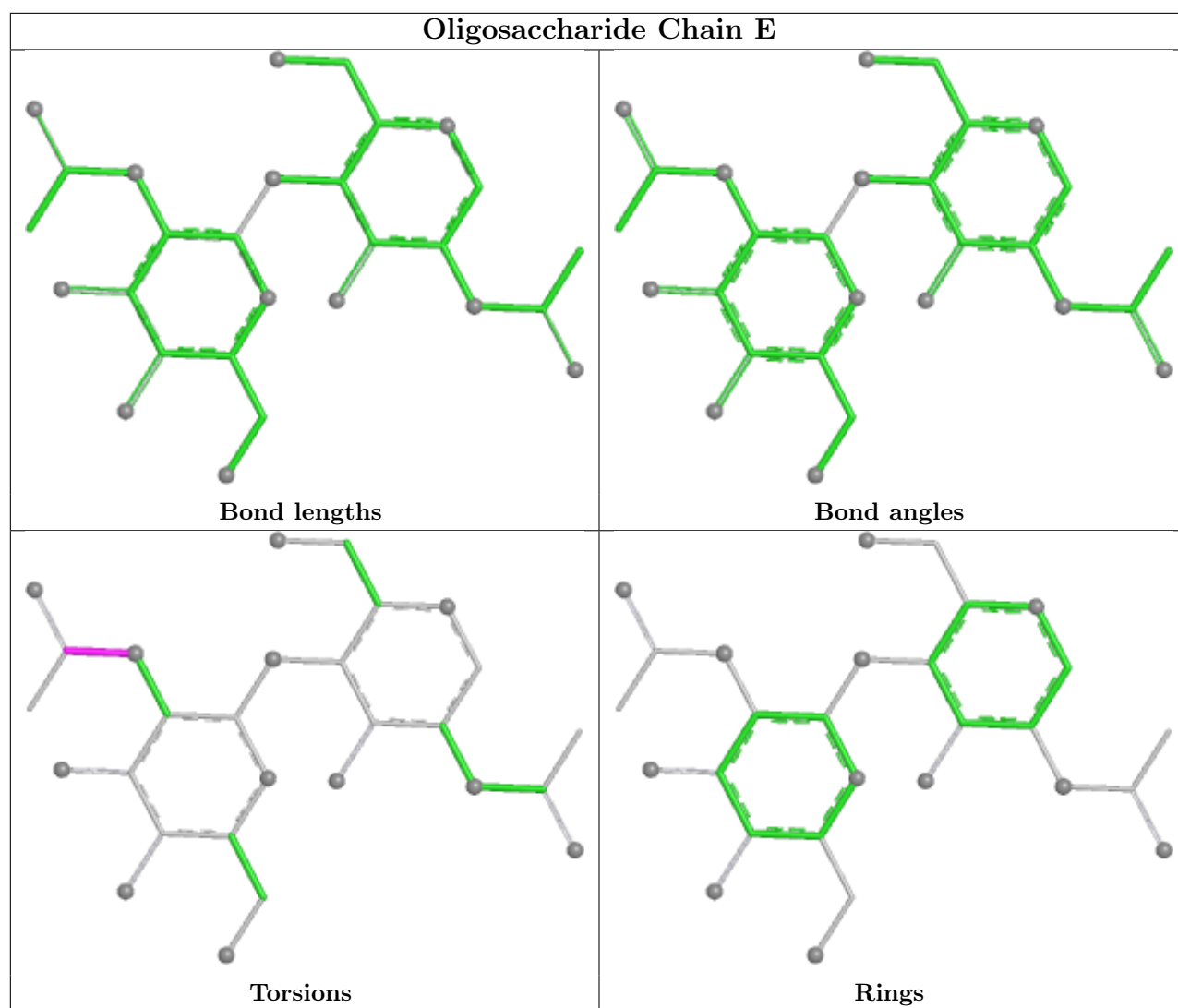
4 monomers are involved in 3 short contacts:

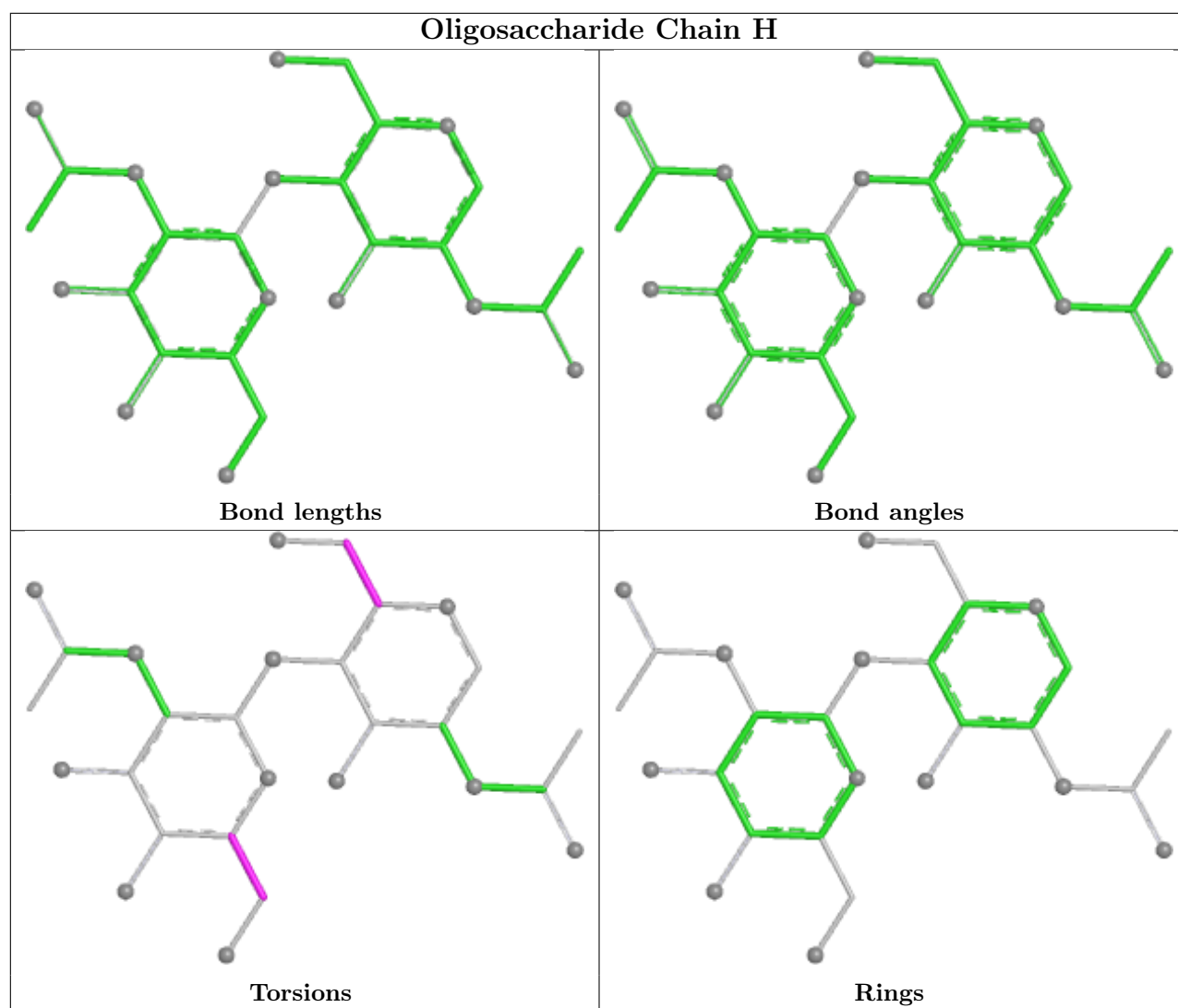
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	K	1	NAG	2	0
5	H	2	NAG	1	0
5	K	2	NAG	1	0
5	H	1	NAG	1	0

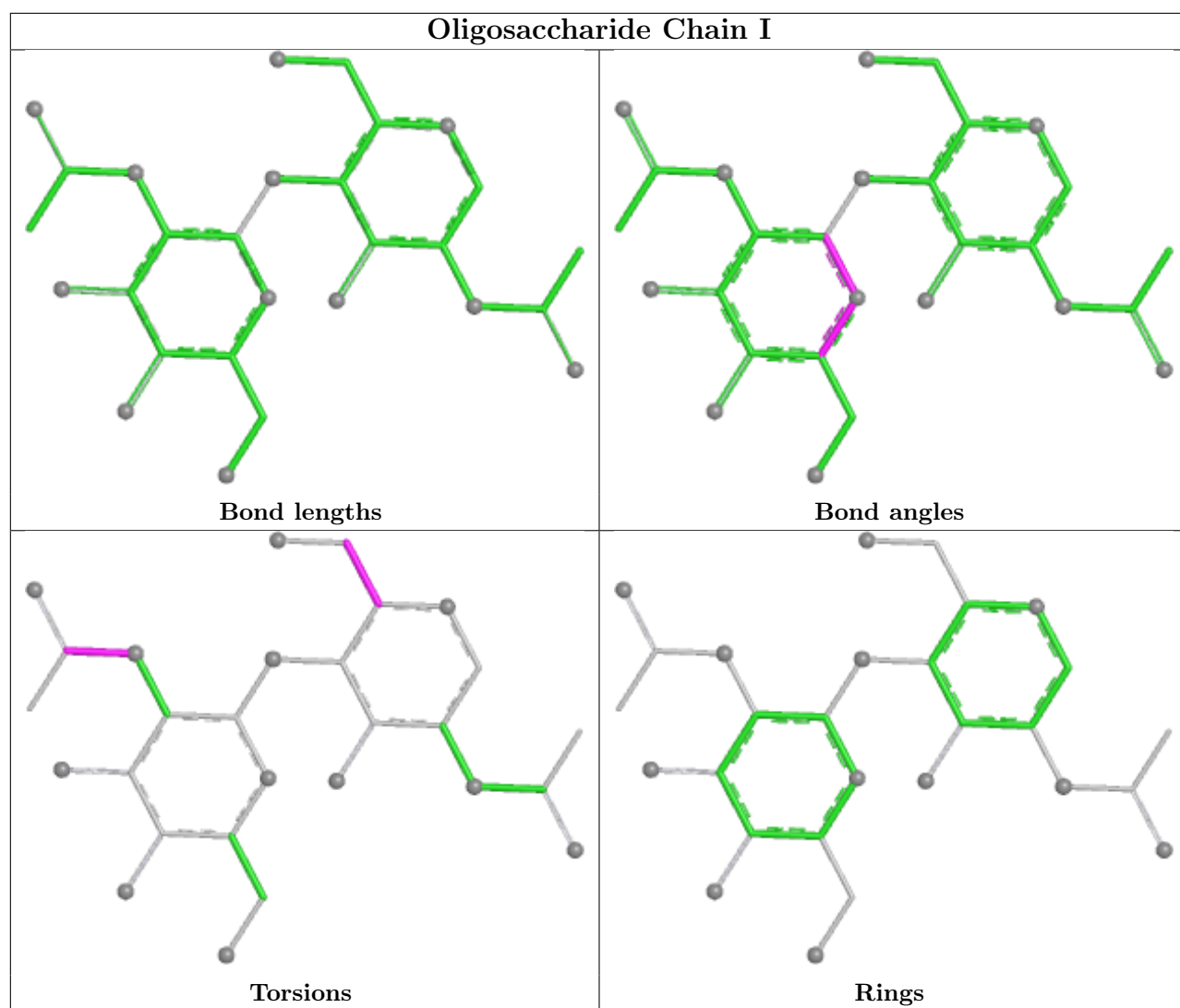
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

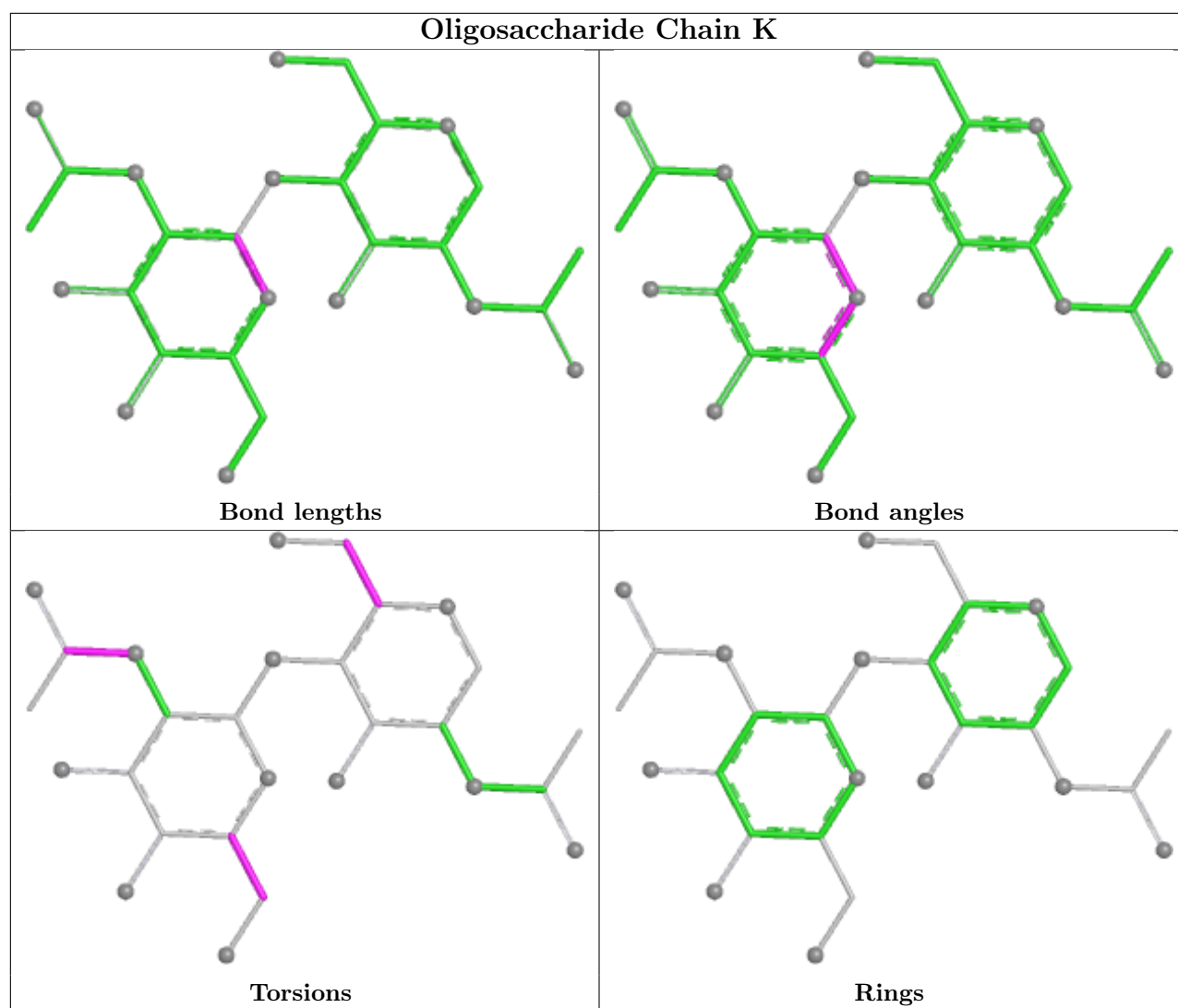




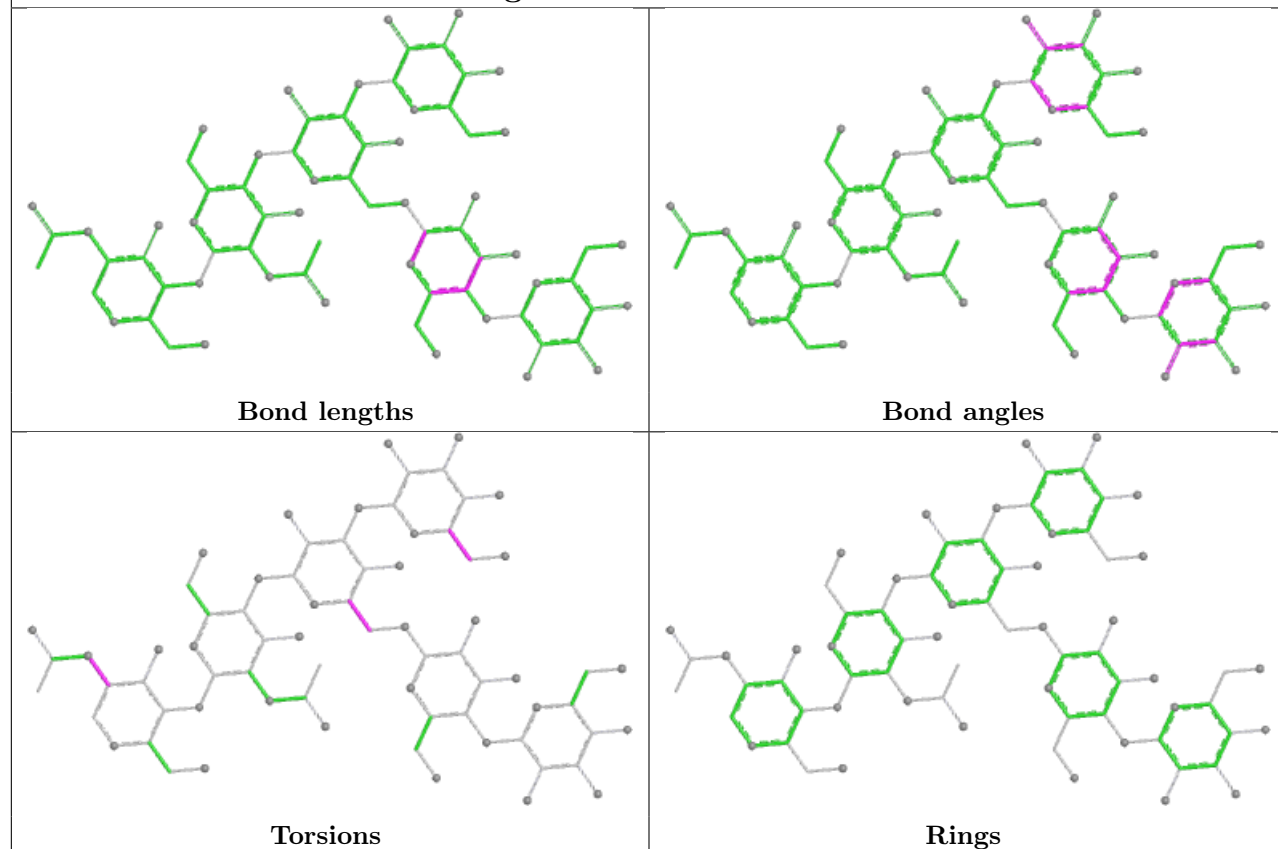




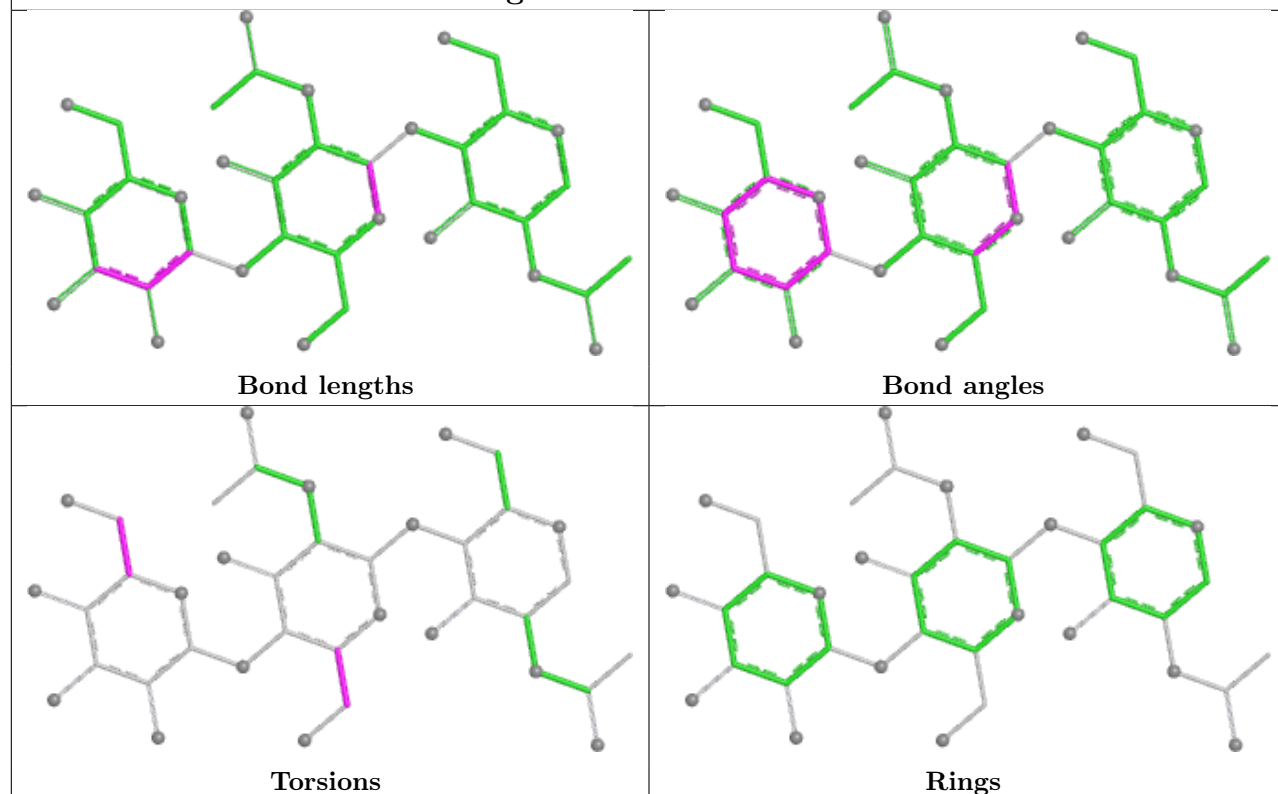


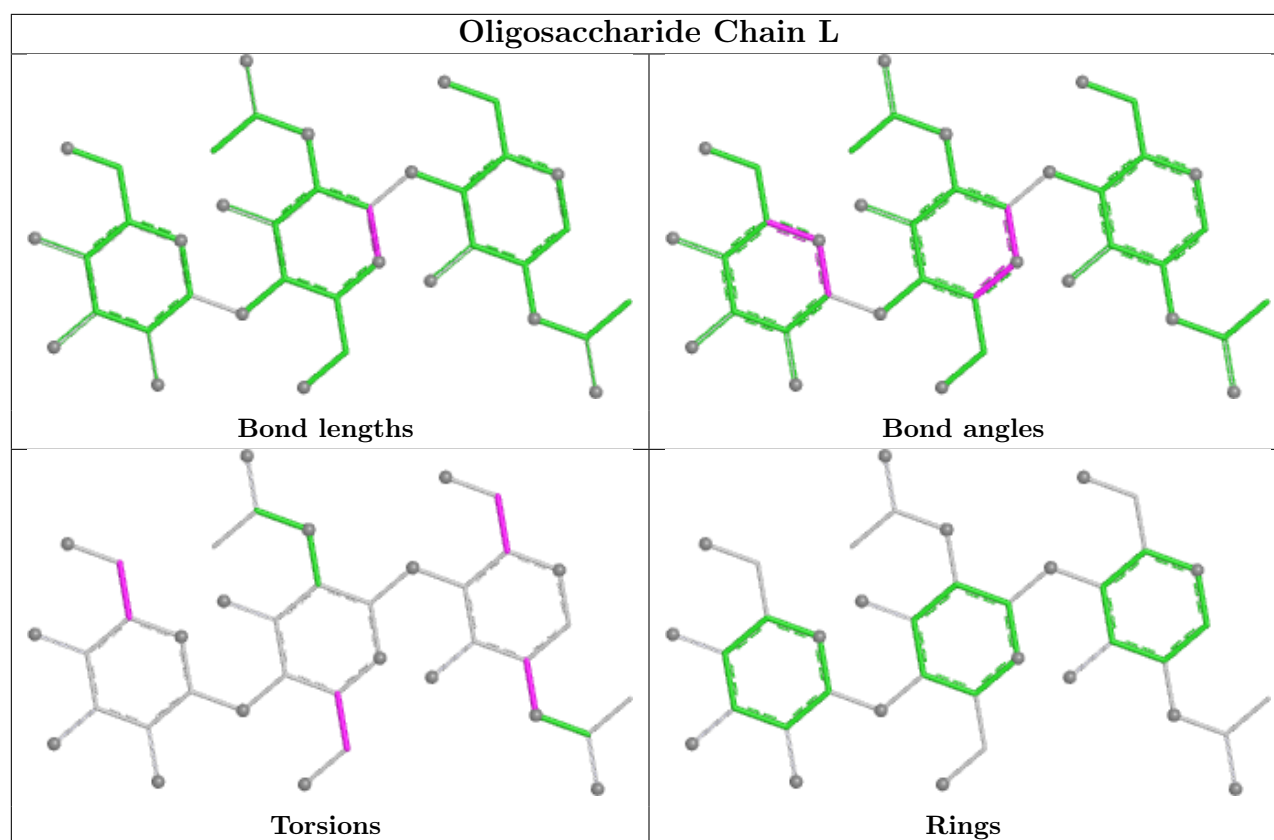


Oligosaccharide Chain F



Oligosaccharide Chain J





5.6 Ligand geometry [i](#)

Of 13 ligands modelled in this entry, 8 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	$\# Z > 2$
8	NAG	A	1020	1	14,14,15	0.27	0	17,19,21	0.47	0
8	NAG	A	1021	1	14,14,15	0.24	0	17,19,21	0.44	0
8	NAG	B	701	2	14,14,15	0.69	1 (7%)	17,19,21	0.83	1 (5%)
8	NAG	B	702	2	14,14,15	0.36	0	17,19,21	0.49	0
8	NAG	A	1017	1	14,14,15	0.28	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
8	NAG	A	1020	1	-	1/6/23/26	0/1/1/1
8	NAG	A	1021	1	-	1/6/23/26	0/1/1/1
8	NAG	B	701	2	-	1/6/23/26	0/1/1/1
8	NAG	B	702	2	-	2/6/23/26	0/1/1/1
8	NAG	A	1017	1	-	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)
8	B	701	NAG	O5-C1	2.30	1.47	1.43

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	B	701	NAG	C1-O5-C5	3.18	116.45	112.19

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	B	702	NAG	O5-C5-C6-O6
8	A	1017	NAG	O5-C5-C6-O6
8	B	701	NAG	O5-C5-C6-O6
8	B	702	NAG	C4-C5-C6-O6
8	A	1020	NAG	O5-C5-C6-O6
8	A	1021	NAG	C1-C2-N2-C7

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	B	701	NAG	1	0
8	B	702	NAG	1	0

5.7 Other polymers [i](#)

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	924/959 (96%)	-0.13	10 (1%)	77 62	56, 105, 149, 190	0
2	B	690/692 (99%)	0.10	15 (2%)	62 46	58, 118, 211, 259	1 (0%)
3	C	8/98 (8%)	0.74	1 (12%)	9 6	99, 135, 152, 157	0
All	All	1622/1749 (92%)	-0.03	26 (1%)	70 55	56, 110, 186, 259	1 (0%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	785	ASN	5.0
2	B	557	TYR	4.3
2	B	456	GLU	3.7
2	B	499	GLY	3.6
2	B	29	GLU	3.3
2	B	556	TYR	3.2
1	A	835	ILE	2.8
2	B	661	ASP	2.8
1	A	405	GLY	2.7
1	A	472	CYS	2.7
2	B	340	VAL	2.7
2	B	591	PRO	2.6
1	A	741	ALA	2.5
2	B	37	ARG	2.5
2	B	36	PRO	2.5
2	B	670	TYR	2.4
2	B	535	MET	2.4
1	A	619	ILE	2.4
2	B	443	PRO	2.3
2	B	411	GLU	2.3
1	A	570	ASP	2.2
1	A	907	THR	2.2
3	C	1491	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	569	ALA	2.1
1	A	503	LYS	2.1
2	B	458	GLY	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.

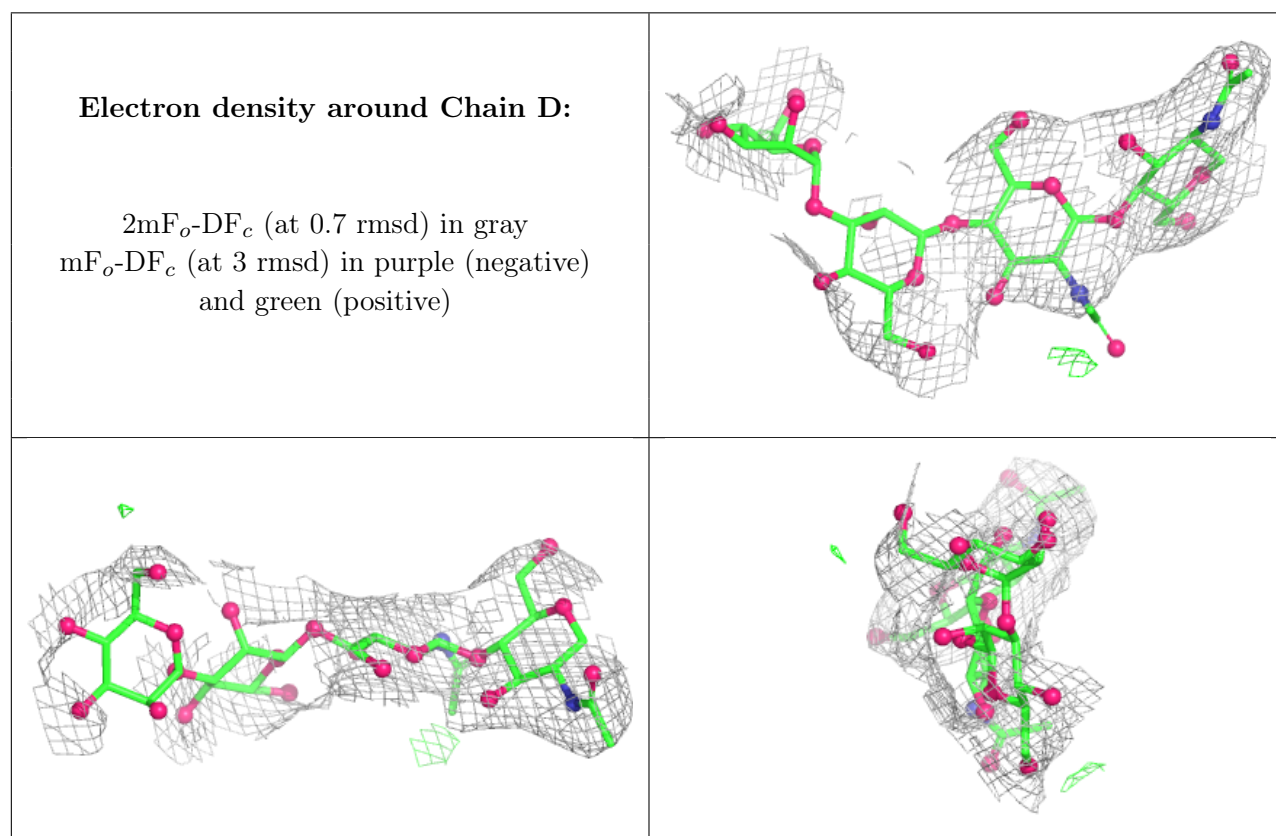
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	BMA	J	3	11/12	0.30	0.11	173,182,211,217	0
5	NAG	H	2	14/15	0.42	0.13	119,145,151,154	0
4	MAN	D	4	11/12	0.44	0.13	197,201,203,205	0
5	NAG	E	2	14/15	0.44	0.11	153,173,179,180	0
4	BMA	D	3	11/12	0.47	0.09	175,189,196,198	0
6	MAN	F	6	11/12	0.53	0.10	126,154,167,169	0
4	BMA	G	3	11/12	0.53	0.10	169,184,221,228	0
7	NAG	J	2	14/15	0.54	0.10	170,192,209,217	0
4	MAN	G	4	11/12	0.55	0.09	172,186,195,200	0
5	NAG	I	2	14/15	0.59	0.09	144,178,183,183	0
5	NAG	H	1	14/15	0.66	0.14	128,152,163,163	0
6	MAN	F	5	11/12	0.71	0.11	147,171,182,182	0
5	NAG	I	1	14/15	0.71	0.11	107,133,152,163	0
5	NAG	K	2	14/15	0.71	0.10	126,167,175,177	0
6	BMA	F	4	11/12	0.71	0.14	136,153,183,190	0
7	BMA	L	3	11/12	0.73	0.11	119,129,139,140	0
4	NAG	G	2	14/15	0.75	0.12	127,164,185,190	0
4	NAG	D	2	14/15	0.78	0.14	139,158,180,182	0
6	BMA	F	3	11/12	0.79	0.10	135,145,154,155	0
4	NAG	G	1	14/15	0.80	0.10	86,118,158,166	0
5	NAG	E	1	14/15	0.85	0.10	111,141,164,166	0
7	NAG	J	1	14/15	0.88	0.08	86,127,138,145	0
5	NAG	K	1	14/15	0.89	0.09	94,123,132,149	0
4	NAG	D	1	14/15	0.93	0.11	69,87,110,126	0
6	NAG	F	2	14/15	0.93	0.12	66,89,109,133	0

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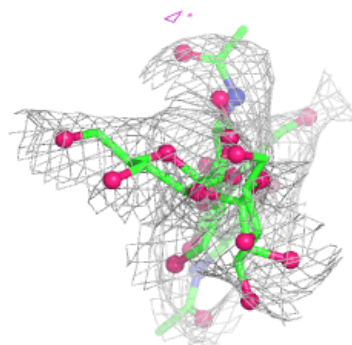
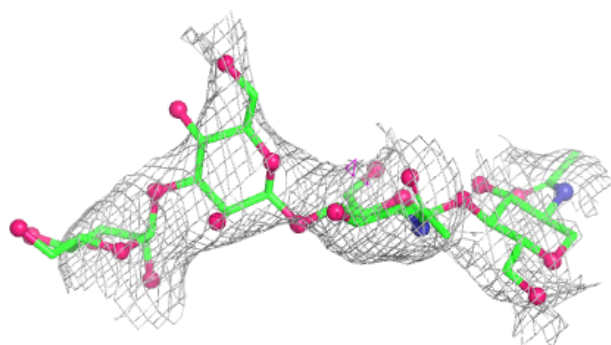
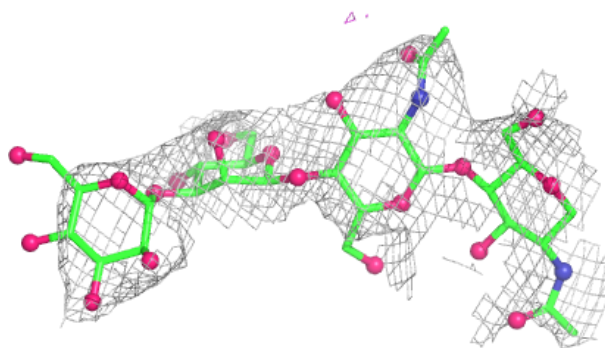
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
7	NAG	L	1	14/15	0.94	0.08	85,104,125,128	0
7	NAG	L	2	14/15	0.96	0.06	109,115,128,129	0
6	NAG	F	1	14/15	0.96	0.07	52,69,101,106	0

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.



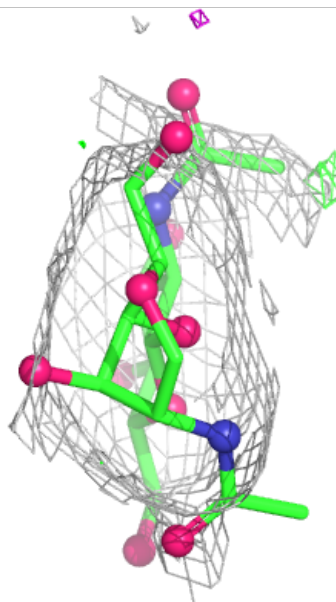
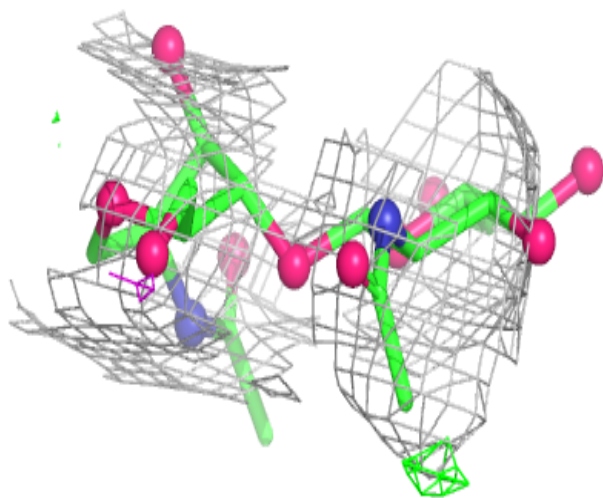
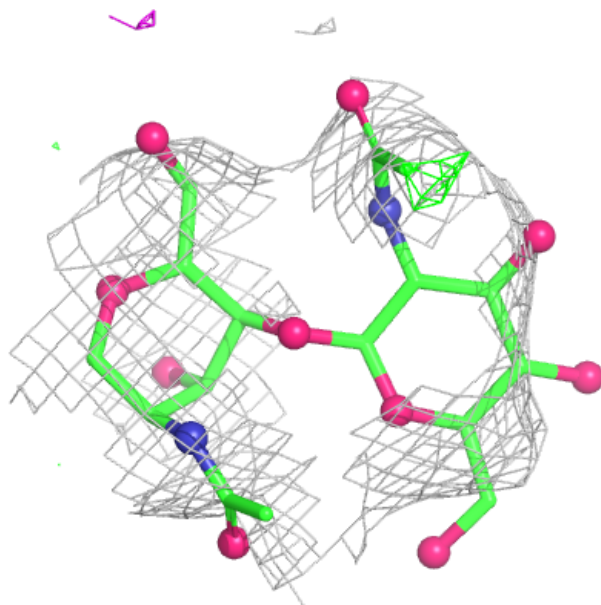
Electron density around Chain G:

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



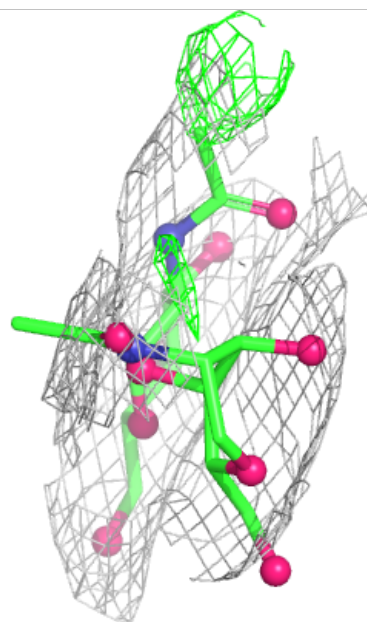
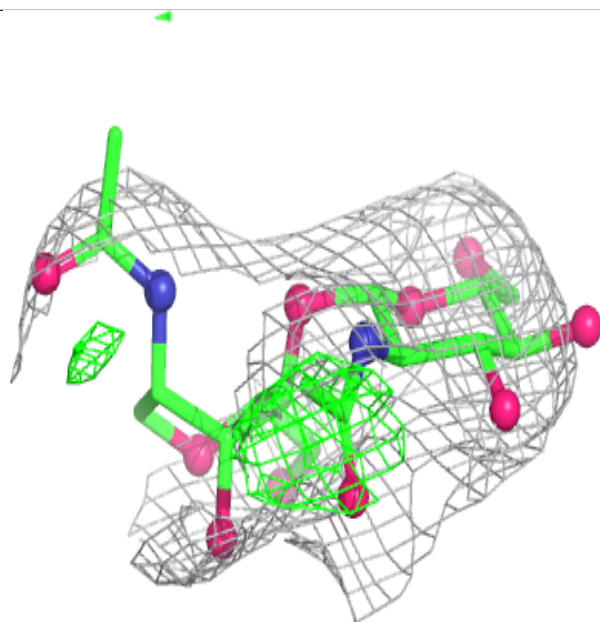
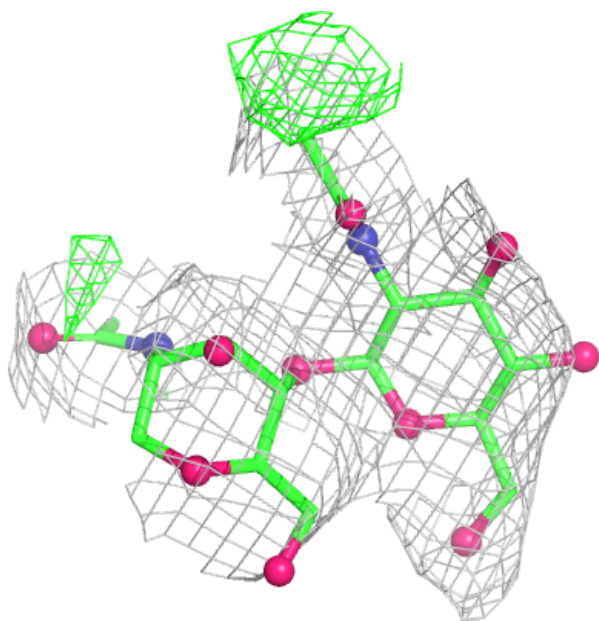
Electron density around Chain E:

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



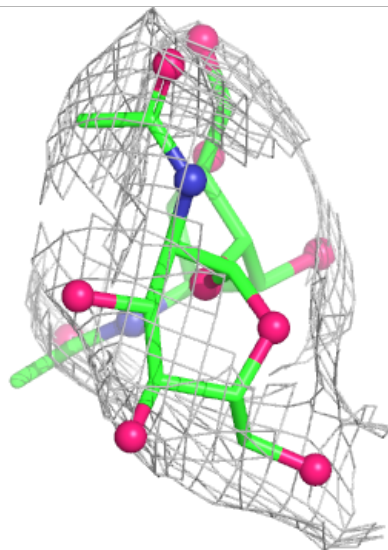
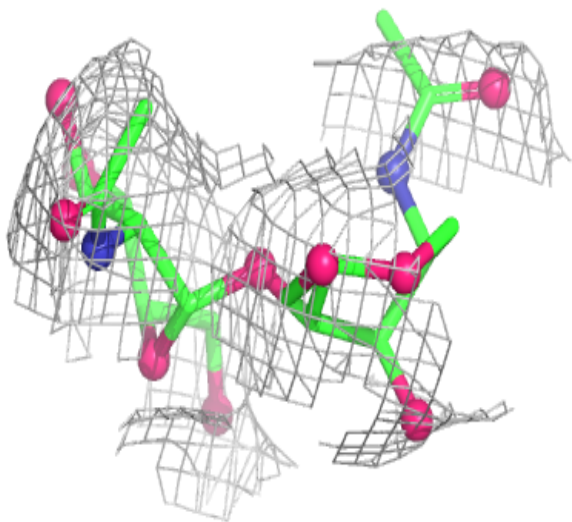
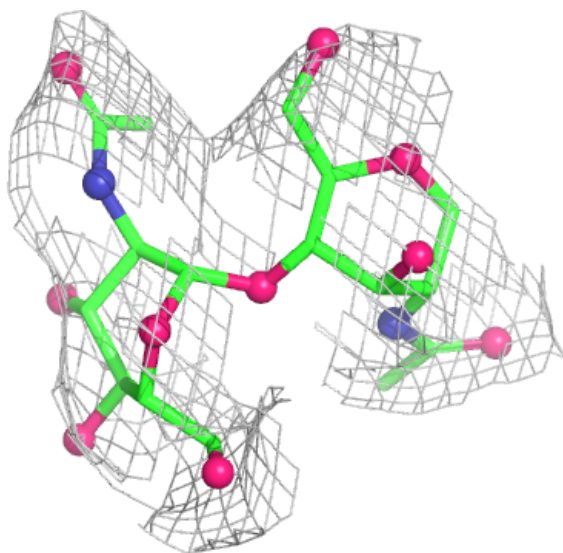
Electron density around Chain H:

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



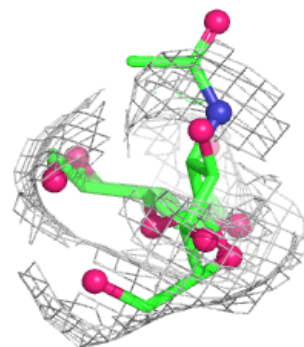
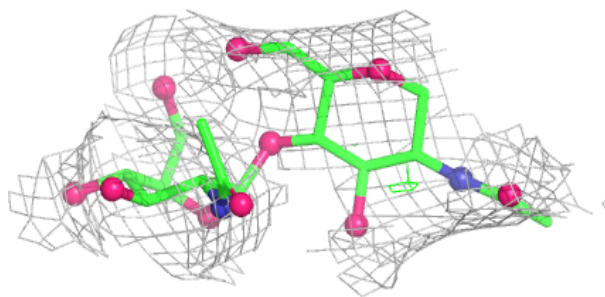
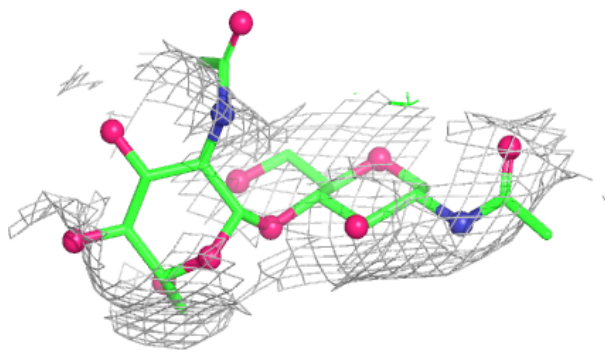
Electron density around Chain I:

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

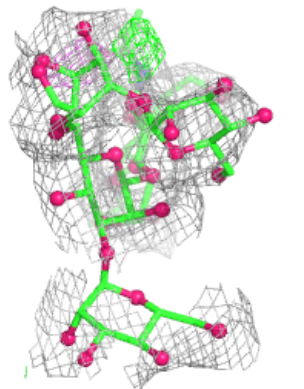
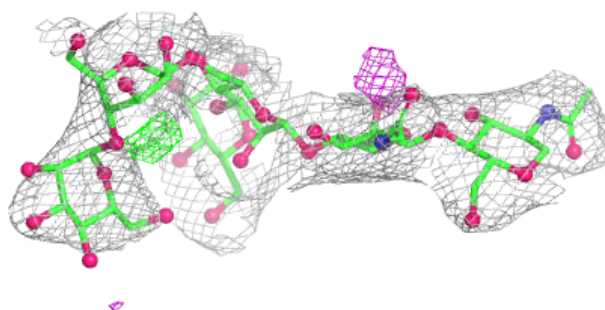
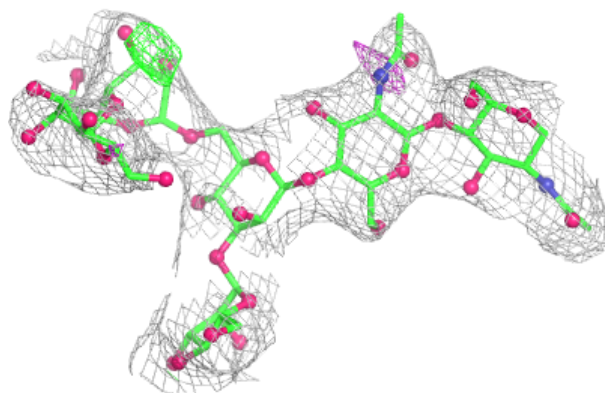


Electron density around Chain K:

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

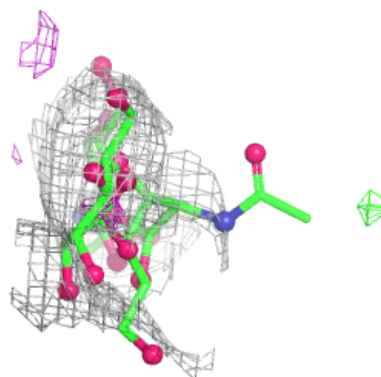
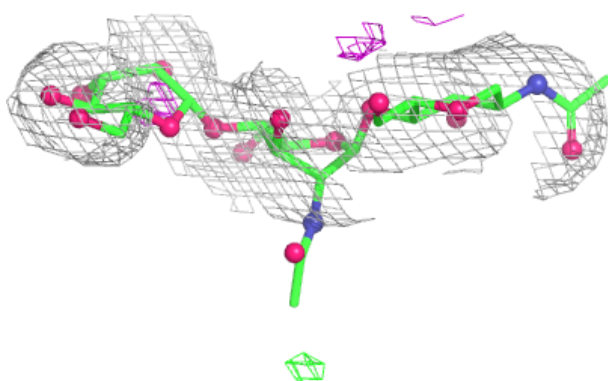
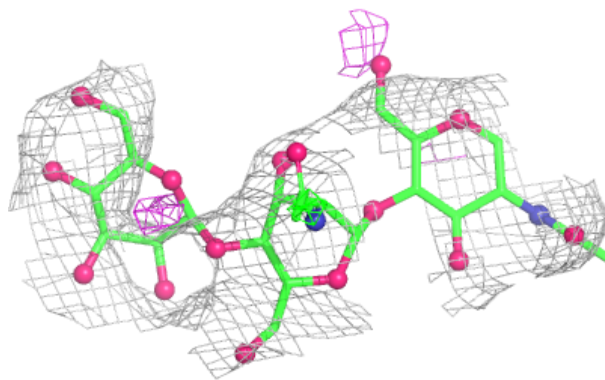
**Electron density around Chain F:**

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

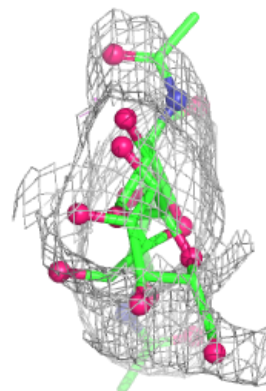
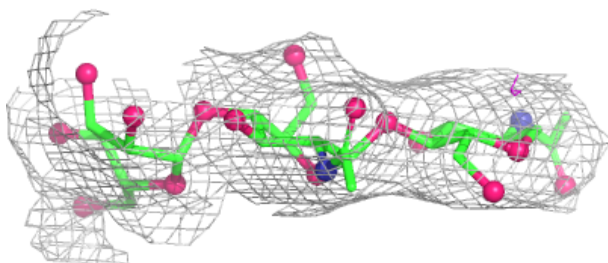
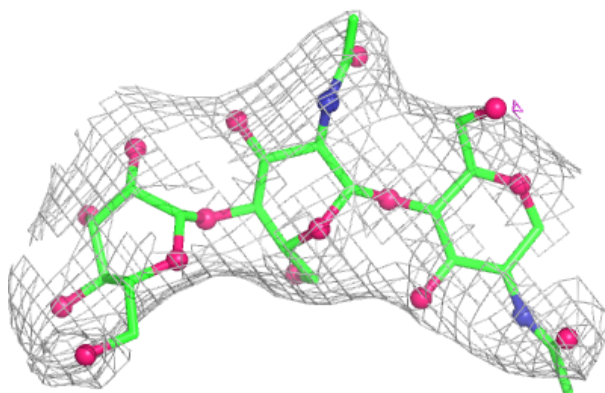


Electron density around Chain J:

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

**Electron density around Chain L:**

$2mF_o-DF_c$ (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(\AA^2)	Q<0.9
8	NAG	A	1020	14/15	0.20	0.13	156,182,185,188	0
8	NAG	A	1017	14/15	0.60	0.12	143,167,187,187	0
8	NAG	B	701	14/15	0.73	0.13	123,162,179,180	0
8	NAG	A	1021	14/15	0.77	0.11	92,119,132,144	0
8	NAG	B	702	14/15	0.78	0.10	95,128,144,150	0
9	MN	A	1027	1/1	0.97	0.04	141,141,141,141	0
9	MN	B	709	1/1	0.98	0.05	115,115,115,115	0
9	MN	A	1029	1/1	0.99	0.05	131,131,131,131	0
9	MN	A	1030	1/1	0.99	0.06	123,123,123,123	0
9	MN	A	1031	1/1	0.99	0.03	111,111,111,111	0
9	MN	A	1028	1/1	0.99	0.08	123,123,123,123	0
9	MN	B	710	1/1	0.99	0.03	99,99,99,99	0
9	MN	B	708	1/1	1.00	0.03	65,65,65,65	0

6.5 Other polymers [i](#)

There are no such residues in this entry.