



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:27 AM UTC

PDB ID : 5E65 / pdb_00005e65
Title : The complex structure of Hemagglutinin-esterase-fusion mutant protein from the influenza D virus with receptor analog 9-O-Ac-3'SLN (Tr322)
Authors : Song, H.; Qi, J.; Shi, Y.; Gao, G.F.
Deposited on : 2015-10-09
Resolution : 2.20 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

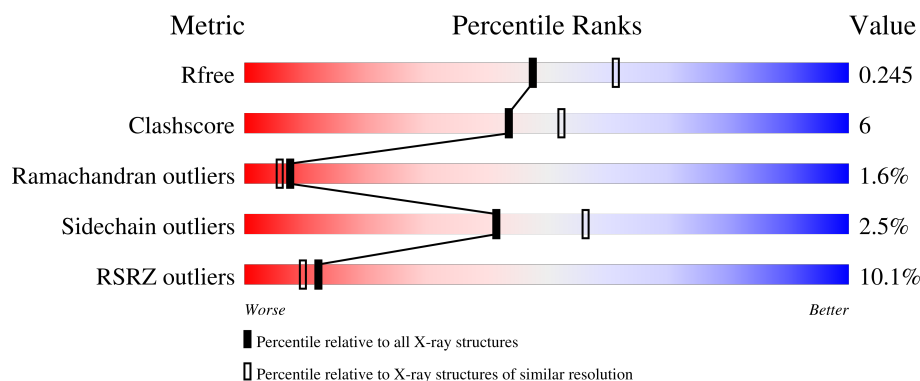
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	427	
1	C	427	
2	B	166	
2	D	166	
3	E	6	

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Mol	Chain	Length	Quality of chain
3	F	6	 17% 33% 50%
3	J	6	 50% 50%
4	G	2	 50% 50%
4	I	2	 50% 50%
4	K	2	 50% 50%
4	L	2	 100%
4	N	2	 50% 50%
5	H	2	 50% 50%
5	M	2	 100%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 9829 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Hemagglutinin-esterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	424	Total	C	N	O	S	0	0	0
			3240	2039	540	639	22			
1	C	424	Total	C	N	O	S	0	0	0
			3240	2039	540	639	22			

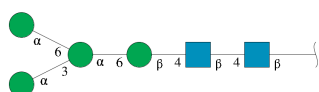
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	57	ALA	SER	engineered mutation	UNP K9LG83
A	356	ALA	ASP	engineered mutation	UNP K9LG83
A	359	ALA	HIS	engineered mutation	UNP K9LG83
C	57	ALA	SER	engineered mutation	UNP K9LG83
C	356	ALA	ASP	engineered mutation	UNP K9LG83
C	359	ALA	HIS	engineered mutation	UNP K9LG83

- Molecule 2 is a protein called Hemagglutinin-esterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	149	Total	C	N	O	S	0	0	0
			1110	695	194	217	4			
2	D	149	Total	C	N	O	S	0	0	0
			1110	695	194	217	4			

- Molecule 3 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]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.



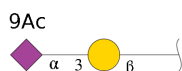
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	E	6	Total	C	N	O	0	0	0
			72	40	2	30			
3	F	6	Total	C	N	O	0	0	0
			72	40	2	30			
3	J	6	Total	C	N	O	0	0	0
			72	40	2	30			

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



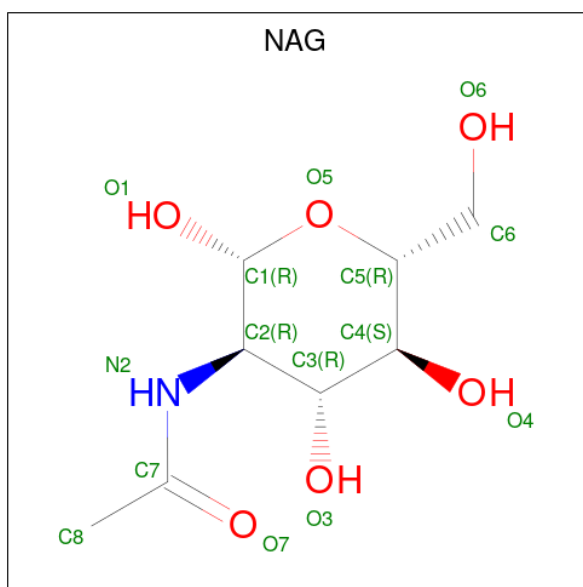
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	I	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	K	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	L	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	N	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 5 is an oligosaccharide called 9-O-acetyl-5-acetamido-3,5-dideoxy-D-glycero-alpha-D-galacto-non-2-ulopyranosonic acid-(2-3)-beta-D-galactopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	H	2	Total	C	N	O	0	0	0
			35	19	1	15			
5	M	2	Total	C	N	O	0	0	0
			35	19	1	15			

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

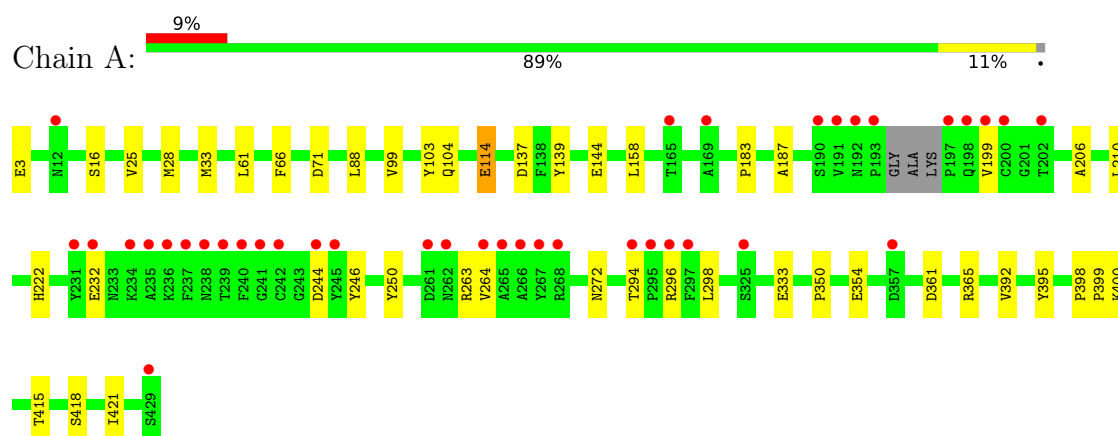
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	264	Total	O	0	0
			264	264		
7	B	68	Total	O	0	0
			68	68		
7	C	304	Total	O	0	0
			304	304		
7	D	39	Total	O	0	0
			39	39		

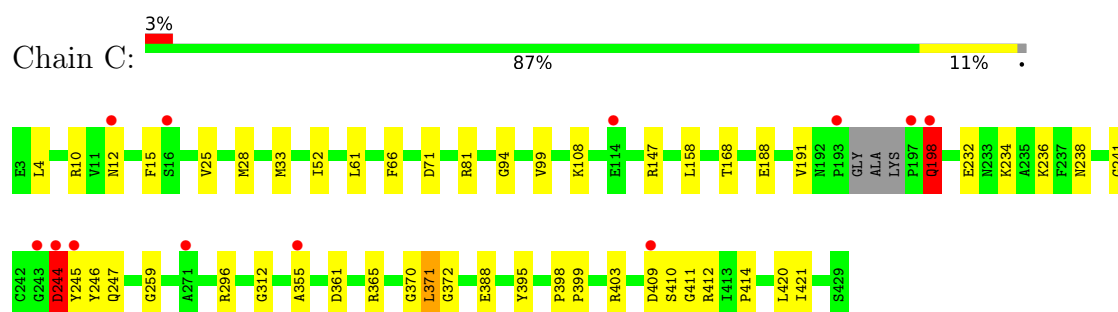
3 Residue-property plots [i](#)

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

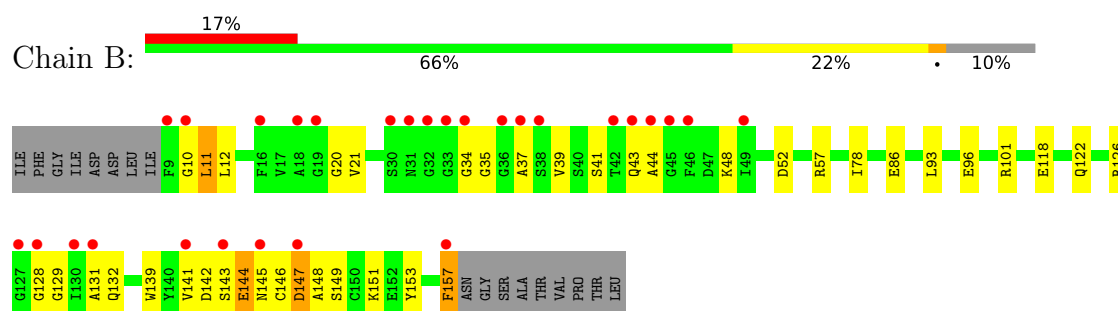
• Molecule 1: Hemagglutinin-esterase



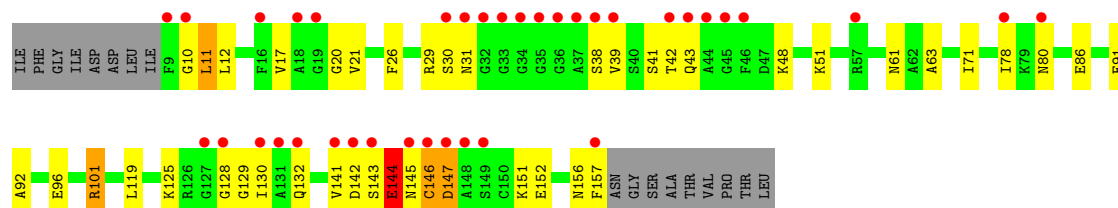
• Molecule 1: Hemagglutinin-esterase



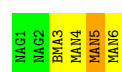
• Molecule 2: Hemagglutinin-esterase



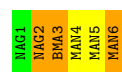
• Molecule 2: Hemagglutinin-esterase



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



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



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

Chain K:  50% 50%

 MAG1
MAG2

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

Chain L:  100%

 MAG1
MAG2

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

Chain N:  50% 50%

 MAG1
MAG2

- Molecule 5: 9-O-acetyl-5-acetamido-3,5-dideoxy-D-glycero-alpha-D-galacto-non-2-ulopyranosonic acid-(2-3)-beta-D-galactopyranose

Chain H:  50% 50%

 GAL1
SUG2

- Molecule 5: 9-O-acetyl-5-acetamido-3,5-dideoxy-D-glycero-alpha-D-galacto-non-2-ulopyranosonic acid-(2-3)-beta-D-galactopyranose

Chain M:  100%

 GAL1
SUG2

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α , β , γ	164.90Å 164.90Å 164.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.60 – 2.20 47.60 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.5 (47.60-2.20) 99.6 (47.60-2.20)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.30 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.203 , 0.243 0.208 , 0.245	Depositor DCC
R_{free} test set	3800 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	33.4	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.021 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9829	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MAN, NAG, 5N6, BMA, GAL

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	0.34	0/3315	0.70	1/4494 (0.0%)
1	C	0.36	0/3315	0.74	2/4494 (0.0%)
2	B	0.36	0/1127	0.76	0/1515
2	D	0.40	0/1127	0.77	0/1515
All	All	0.36	0/8884	0.73	3/12018 (0.0%)

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	C	0	1
2	D	0	1
All	All	0	2

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	371	LEU	N-CA-C	8.75	123.77	113.18
1	C	198	GLN	CA-CB-CG	6.01	126.12	114.10
1	A	199	VAL	N-CA-C	5.11	115.33	108.17

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	244	ASP	Peptide

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Mol	Chain	Res	Type	Group
2	D	144	GLU	Peptide

5.2 Too-close contacts ⓘ

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	3240	0	3114	31	0
1	C	3240	0	3114	34	0
2	B	1110	0	1067	23	0
2	D	1110	0	1067	30	0
3	E	72	0	61	1	0
3	F	72	0	61	2	0
3	J	72	0	61	0	0
4	G	28	0	25	1	0
4	I	28	0	25	1	0
4	K	28	0	25	0	0
4	L	28	0	25	0	0
4	N	28	0	25	0	0
5	H	35	0	11	1	0
5	M	35	0	11	1	0
6	B	14	0	13	0	0
6	D	14	0	13	1	0
7	A	264	0	0	9	1
7	B	68	0	0	2	0
7	C	304	0	0	4	0
7	D	39	0	0	2	0
All	All	9829	0	8718	109	1

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 (109) 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)
5:H:2:5N6:O10	5:H:2:5N6:O7	1.89	0.90
3:F:2:NAG:O3	3:F:3:BMA:H2	1.81	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:63:ALA:O	7:D:801:HOH:O	2.02	0.76
1:A:415:THR:O	2:B:101:ARG:NH1	2.19	0.75
7:B:801:HOH:O	4:I:2:NAG:O4	2.07	0.73
1:A:250:TYR:OH	7:A:801:HOH:O	1.89	0.70
2:B:48:LYS:NZ	2:B:52:ASP:OD2	2.24	0.68
1:C:388:GLU:OE1	7:C:802:HOH:O	2.12	0.68
1:A:33:MET:HE1	2:B:101:ARG:HB2	1.74	0.68
1:C:188:GLU:OE1	7:C:801:HOH:O	2.12	0.68
2:B:126:ARG:NE	2:B:153:TYR:O	2.27	0.66
2:B:96:GLU:OE2	7:B:802:HOH:O	2.14	0.66
1:A:16:SER:OG	7:A:802:HOH:O	2.14	0.65
1:A:144:GLU:OE1	7:A:803:HOH:O	2.14	0.64
1:A:333:GLU:OE1	7:A:804:HOH:O	2.15	0.63
2:D:142:ASP:O	2:D:144:GLU:N	2.33	0.61
2:D:141:VAL:HG23	2:D:145:ASN:HD21	1.65	0.61
2:B:142:ASP:O	2:B:144:GLU:N	2.25	0.61
1:C:371:LEU:N	1:C:372:GLY:HA3	2.16	0.60
2:D:30:SER:OG	2:D:31:ASN:N	2.31	0.60
2:D:41:SER:OG	2:D:42:THR:N	2.35	0.60
1:C:25:VAL:HG13	1:C:421:ILE:HG23	1.84	0.60
2:D:130:ILE:CG2	2:D:145:ASN:HD22	2.16	0.58
1:A:114:GLU:OE2	1:A:139:TYR:OH	2.11	0.58
1:C:33:MET:HE1	2:D:101:ARG:HB2	1.84	0.58
7:A:805:HOH:O	3:F:6:MAN:O2	2.17	0.58
1:A:361:ASP:O	1:A:365:ARG:HG3	2.04	0.58
1:C:147:ARG:NH1	7:C:814:HOH:O	2.37	0.57
1:A:350:PRO:HD3	1:C:410:SER:HB3	1.86	0.57
2:B:142:ASP:C	2:B:144:GLU:H	2.13	0.57
2:D:10:GLY:O	2:D:12:LEU:N	2.31	0.56
1:C:409:ASP:OD1	1:C:411:GLY:N	2.31	0.56
1:A:104:GLN:NE2	7:A:814:HOH:O	2.39	0.55
2:D:141:VAL:CG2	2:D:145:ASN:HD21	2.19	0.55
2:B:145:ASN:OD1	2:B:146:CYS:N	2.39	0.54
2:D:78:ILE:HG22	2:D:80:ASN:H	1.73	0.54
2:D:51:LYS:N	2:D:51:LYS:HD3	2.22	0.54
1:A:137:ASP:OD2	7:A:806:HOH:O	2.18	0.53
1:A:244:ASP:OD2	1:A:263:ARG:HG3	2.09	0.53
2:D:92:ALA:O	2:D:96:GLU:HG3	2.09	0.53
1:A:25:VAL:HG13	1:A:421:ILE:HG23	1.90	0.53
1:A:418:SER:HA	2:B:57:ARG:HD2	1.91	0.53
2:D:29:ARG:O	2:D:38:SER:OG	2.26	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:12:ASN:H	1:C:15:PHE:HD1	1.56	0.53
1:A:88:LEU:N	1:A:114:GLU:OE2	2.39	0.52
2:D:156:ASN:O	2:D:157:PHE:HB2	2.09	0.52
2:D:26:PHE:HD1	2:D:43:GLN:N	2.07	0.52
1:C:412:ARG:NH2	2:D:91:GLU:HA	2.24	0.52
1:C:94:GLY:HA2	7:C:906:HOH:O	2.09	0.52
1:A:28:MET:SD	4:G:1:NAG:H5	2.50	0.51
5:M:1:GAL:H4	5:M:2:5N6:C1	2.40	0.51
2:D:43:GLN:NE2	7:D:803:HOH:O	2.44	0.51
2:D:146:CYS:SG	2:D:151:LYS:N	2.83	0.51
1:C:370:GLY:C	1:C:372:GLY:HA3	2.36	0.51
1:C:28:MET:HE2	1:C:420:LEU:HB3	1.92	0.51
1:C:52:ILE:HD11	1:C:81:ARG:HG2	1.93	0.50
1:C:33:MET:HE2	1:C:414:PRO:HB2	1.94	0.49
1:A:88:LEU:HD23	1:A:103:TYR:CD2	2.48	0.48
1:C:71:ASP:HA	1:C:365:ARG:HG2	1.94	0.48
2:D:61:ASN:HD22	6:D:703:NAG:H61	1.77	0.48
1:A:395:TYR:CD2	2:B:86:GLU:HG2	2.48	0.48
2:B:10:GLY:O	2:B:12:LEU:N	2.43	0.48
1:C:234:LYS:NZ	1:C:238:ASN:HD21	2.11	0.48
1:A:246:TYR:H	1:A:294:THR:HG23	1.79	0.48
1:C:108:LYS:HD2	1:C:312:GLY:HA2	1.95	0.48
2:D:145:ASN:CG	2:D:146:CYS:HB3	2.38	0.48
1:C:4:LEU:H	2:D:30:SER:HB3	1.78	0.48
1:C:10:ARG:HE	2:D:17:VAL:HG11	1.79	0.47
2:B:147:ASP:OD1	2:B:148:ALA:N	2.48	0.47
1:C:238:ASN:HA	1:C:241:GLY:O	2.14	0.47
1:C:395:TYR:HD2	2:D:86:GLU:HG2	1.80	0.46
1:C:395:TYR:CD2	2:D:86:GLU:HG2	2.50	0.46
2:D:26:PHE:HD1	2:D:43:GLN:H	1.63	0.46
2:B:44:ALA:HA	2:B:157:PHE:CE2	2.51	0.46
1:A:61:LEU:HA	1:A:66:PHE:CD1	2.51	0.46
1:C:234:LYS:HZ3	1:C:238:ASN:HD21	1.64	0.46
1:C:198:GLN:HE21	1:C:198:GLN:HB3	1.27	0.46
2:B:146:CYS:SG	2:B:151:LYS:N	2.89	0.45
1:A:3:GLU:N	2:B:141:VAL:O	2.50	0.45
1:A:398:PRO:HA	1:A:399:PRO:C	2.42	0.45
2:B:128:GLY:HA2	2:B:129:GLY:HA2	1.61	0.45
2:B:118:GLU:O	2:B:122:GLN:HG3	2.17	0.44
2:D:48:LYS:HD2	2:D:119:LEU:HD13	1.98	0.44
7:A:807:HOH:O	3:E:5:MAN:O6	2.21	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:244:ASP:HB3	1:C:246:TYR:CD2	2.53	0.43
1:A:206:ALA:HB2	1:A:298:LEU:HD21	2.01	0.43
1:C:403:ARG:HB3	2:D:71:ILE:HB	2.00	0.43
1:C:191:VAL:HB	1:C:296:ARG:HD3	2.01	0.43
1:A:71:ASP:HA	1:A:365:ARG:HG2	2.00	0.43
2:B:131:ALA:HB1	2:B:139:TRP:CD1	2.54	0.43
1:C:61:LEU:HA	1:C:66:PHE:CD1	2.54	0.43
1:C:398:PRO:HB2	2:D:78:ILE:O	2.19	0.43
2:B:145:ASN:OD1	2:B:146:CYS:HB2	2.19	0.43
2:D:128:GLY:HA2	2:D:129:GLY:HA2	1.70	0.42
1:C:232:GLU:OE1	1:C:236:LYS:NZ	2.31	0.42
1:C:247:GLN:O	1:C:259:GLY:HA3	2.20	0.42
1:A:187:ALA:HA	1:A:296:ARG:O	2.19	0.42
1:C:361:ASP:O	1:C:365:ARG:HG3	2.20	0.41
1:A:183:PRO:HB3	7:A:834:HOH:O	2.20	0.41
1:A:210:LEU:HB3	1:A:222:HIS:CD2	2.56	0.41
1:A:244:ASP:OD2	1:A:263:ARG:N	2.54	0.41
1:A:354:GLU:OE2	1:A:365:ARG:NH2	2.47	0.41
2:B:34:GLY:HA2	2:B:35:GLY:HA2	1.68	0.41
1:C:398:PRO:HA	1:C:399:PRO:C	2.46	0.41
2:D:125:LYS:HB3	2:D:125:LYS:HE2	1.76	0.41
1:A:88:LEU:H	1:A:114:GLU:CD	2.29	0.40
1:A:400:LYS:NZ	2:B:86:GLU:OE2	2.53	0.40
2:B:147:ASP:OD1	2:B:149:SER:N	2.54	0.40
2:B:93:LEU:HD12	2:B:93:LEU:HA	1.90	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:1061:HOH:O	7:A:1063:HOH:O[9_555]	2.04	0.16

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	420/427 (98%)	403 (96%)	17 (4%)	0	100	100
1	C	420/427 (98%)	398 (95%)	20 (5%)	2 (0%)	24	27
2	B	147/166 (89%)	122 (83%)	15 (10%)	10 (7%)	1	0
2	D	147/166 (89%)	124 (84%)	17 (12%)	6 (4%)	2	1
All	All	1134/1186 (96%)	1047 (92%)	69 (6%)	18 (2%)	7	6

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	147	ASP
1	C	245	TYR
2	D	11	LEU
2	D	143	SER
2	B	11	LEU
2	B	39	VAL
2	B	143	SER
1	C	355	ALA
2	D	147	ASP
2	B	41	SER
2	B	37	ALA
2	B	43	GLN
2	D	21	VAL
2	B	21	VAL
2	B	132	GLN
2	D	132	GLN
2	D	20	GLY
2	B	20	GLY

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	357/358 (100%)	350 (98%)	7 (2%)	48 64
1	C	357/358 (100%)	352 (99%)	5 (1%)	59 75
2	B	109/123 (89%)	105 (96%)	4 (4%)	30 41
2	D	109/123 (89%)	102 (94%)	7 (6%)	16 19
All	All	932/962 (97%)	909 (98%)	23 (2%)	42 56

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	99	VAL
1	A	114	GLU
1	A	158	LEU
1	A	232	GLU
1	A	264	VAL
1	A	272	ASN
1	A	392	VAL
2	B	11	LEU
2	B	78	ILE
2	B	144	GLU
2	B	157	PHE
1	C	99	VAL
1	C	158	LEU
1	C	168	THR
1	C	198	GLN
1	C	244	ASP
2	D	11	LEU
2	D	39	VAL
2	D	101	ARG
2	D	144	GLU
2	D	146	CYS
2	D	147	ASP
2	D	152	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	104	GLN

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Mol	Chain	Res	Type
1	A	238	ASN
1	A	255	ASN
1	A	272	ASN
2	B	55	GLN
2	B	73	HIS
2	B	121	ASN
1	C	9	GLN
1	C	104	GLN
1	C	198	GLN
1	C	204	GLN
1	C	238	ASN
1	C	253	ASN
2	D	61	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

32 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	0.36	0	17,19,21	0.43	0
3	NAG	E	2	3	14,14,15	0.55	0	17,19,21	0.56	0
3	BMA	E	3	3	11,11,12	0.73	0	15,15,17	1.20	2 (13%)
3	MAN	E	4	3	11,11,12	0.88	1 (9%)	15,15,17	1.04	2 (13%)
3	MAN	E	5	3	11,11,12	0.98	0	15,15,17	1.06	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	MAN	E	6	3	11,11,12	0.85	0	15,15,17	1.08	1 (6%)
3	NAG	F	1	3,1	14,14,15	0.30	0	17,19,21	0.39	0
3	NAG	F	2	3	14,14,15	0.21	0	17,19,21	0.87	1 (5%)
3	BMA	F	3	3	11,11,12	1.33	2 (18%)	15,15,17	1.40	2 (13%)
3	MAN	F	4	3	11,11,12	0.91	1 (9%)	15,15,17	1.94	2 (13%)
3	MAN	F	5	3	11,11,12	0.86	1 (9%)	15,15,17	0.90	1 (6%)
3	MAN	F	6	3	11,11,12	1.15	1 (9%)	15,15,17	0.81	0
4	NAG	G	1	4,1	14,14,15	0.32	0	17,19,21	0.49	0
4	NAG	G	2	4	14,14,15	0.36	0	17,19,21	0.54	0
5	GAL	H	1	5	12,12,12	1.14	1 (8%)	17,17,17	1.28	3 (17%)
5	5N6	H	2	5	23,23,24	1.93	9 (39%)	25,32,35	1.53	5 (20%)
4	NAG	I	1	4,2	14,14,15	0.48	0	17,19,21	0.50	0
4	NAG	I	2	4	14,14,15	0.22	0	17,19,21	0.66	1 (5%)
3	NAG	J	1	3,1	14,14,15	0.37	0	17,19,21	0.44	0
3	NAG	J	2	3	14,14,15	0.67	0	17,19,21	0.57	0
3	BMA	J	3	3	11,11,12	0.75	0	15,15,17	1.07	1 (6%)
3	MAN	J	4	3	11,11,12	0.90	0	15,15,17	1.31	1 (6%)
3	MAN	J	5	3	11,11,12	0.81	0	15,15,17	0.84	0
3	MAN	J	6	3	11,11,12	0.92	1 (9%)	15,15,17	0.91	1 (6%)
4	NAG	K	1	4,1	14,14,15	0.34	0	17,19,21	0.58	0
4	NAG	K	2	4	14,14,15	0.21	0	17,19,21	0.64	1 (5%)
4	NAG	L	1	4,1	14,14,15	0.42	0	17,19,21	0.59	0
4	NAG	L	2	4	14,14,15	0.36	0	17,19,21	0.56	0
5	GAL	M	1	5	12,12,12	1.39	3 (25%)	17,17,17	1.16	1 (5%)
5	5N6	M	2	5	23,23,24	2.35	12 (52%)	25,32,35	1.39	3 (12%)
4	NAG	N	1	4,2	14,14,15	0.43	0	17,19,21	0.57	0
4	NAG	N	2	4	14,14,15	0.29	0	17,19,21	0.65	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	-	0/6/23/26	0/1/1/1
3	NAG	E	2	3	-	0/6/23/26	0/1/1/1
3	BMA	E	3	3	-	0/2/19/22	0/1/1/1
3	MAN	E	4	3	-	0/2/19/22	0/1/1/1
3	MAN	E	5	3	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MAN	E	6	3	-	0/2/19/22	0/1/1/1
3	NAG	F	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	F	2	3	-	0/6/23/26	0/1/1/1
3	BMA	F	3	3	-	2/2/19/22	0/1/1/1
3	MAN	F	4	3	-	0/2/19/22	0/1/1/1
3	MAN	F	5	3	-	0/2/19/22	0/1/1/1
3	MAN	F	6	3	-	0/2/19/22	0/1/1/1
4	NAG	G	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	G	2	4	-	2/6/23/26	0/1/1/1
5	GAL	H	1	5	-	2/2/22/22	0/1/1/1
5	5N6	H	2	5	-	10/21/37/41	0/1/1/1
4	NAG	I	1	4,2	-	0/6/23/26	0/1/1/1
4	NAG	I	2	4	-	0/6/23/26	0/1/1/1
3	NAG	J	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	J	2	3	-	0/6/23/26	0/1/1/1
3	BMA	J	3	3	-	0/2/19/22	0/1/1/1
3	MAN	J	4	3	-	0/2/19/22	0/1/1/1
3	MAN	J	5	3	-	2/2/19/22	0/1/1/1
3	MAN	J	6	3	-	2/2/19/22	0/1/1/1
4	NAG	K	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	K	2	4	-	0/6/23/26	0/1/1/1
4	NAG	L	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	L	2	4	-	0/6/23/26	0/1/1/1
5	GAL	M	1	5	-	2/2/22/22	0/1/1/1
5	5N6	M	2	5	-	1/21/37/41	0/1/1/1
4	NAG	N	1	4,2	-	0/6/23/26	0/1/1/1
4	NAG	N	2	4	-	0/6/23/26	0/1/1/1

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	M	2	5N6	O4-C4	-5.03	1.32	1.43
5	M	2	5N6	C4-C5	-3.99	1.49	1.53
5	M	2	5N6	C7-C6	-3.91	1.48	1.52
3	F	6	MAN	O5-C1	-3.39	1.38	1.43
5	H	2	5N6	O9-C9	-3.36	1.37	1.45
5	H	2	5N6	C6-C5	-3.32	1.47	1.53
5	M	2	5N6	C8-C7	-3.30	1.47	1.53
5	H	2	5N6	C3-C2	-3.27	1.47	1.52
5	H	2	5N6	O4-C4	-3.02	1.37	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	M	2	5N6	O6-C6	-2.89	1.39	1.44
5	M	2	5N6	O6-C2	-2.81	1.38	1.43
5	M	2	5N6	C6-C5	-2.80	1.48	1.53
3	F	3	BMA	O5-C5	2.76	1.48	1.43
5	H	2	5N6	C3-C4	-2.74	1.47	1.52
5	H	2	5N6	O8-C8	-2.69	1.37	1.43
5	M	1	GAL	C4-C5	-2.62	1.47	1.53
5	M	2	5N6	C9-C8	-2.52	1.48	1.51
5	H	2	5N6	O1B-C1	-2.49	1.22	1.30
3	F	4	MAN	C1-C2	2.46	1.58	1.52
5	M	2	5N6	O1B-C1	-2.45	1.22	1.30
3	F	3	BMA	C2-C3	-2.35	1.48	1.52
5	M	2	5N6	O8-C8	-2.35	1.38	1.43
5	M	1	GAL	O4-C4	-2.31	1.37	1.43
5	H	2	5N6	C9-C8	-2.27	1.48	1.51
3	F	5	MAN	O5-C1	-2.26	1.39	1.43
5	M	2	5N6	C3-C4	-2.18	1.48	1.52
3	E	4	MAN	O5-C1	-2.14	1.40	1.43
5	M	1	GAL	O2-C2	-2.14	1.37	1.43
5	M	2	5N6	O9-C9	-2.11	1.40	1.45
5	H	1	GAL	C3-C2	-2.07	1.47	1.52
5	H	2	5N6	C8-C7	-2.04	1.49	1.53
3	J	6	MAN	O5-C1	-2.03	1.40	1.43

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	4	MAN	C1-O5-C5	6.00	120.23	112.19
3	J	4	MAN	C1-O5-C5	3.86	117.36	112.19
5	H	2	5N6	C9-O9-CAG	-3.86	107.56	117.08
3	F	3	BMA	C2-C3-C4	-3.82	104.14	110.86
5	M	2	5N6	C6-C5-N5	-3.21	105.79	110.91
3	J	3	BMA	C1-O5-C5	3.14	116.40	112.19
3	E	3	BMA	C1-O5-C5	3.02	116.23	112.19
5	H	2	5N6	C9-C8-C7	-2.87	106.81	112.22
5	H	2	5N6	O8-C8-C9	-2.80	103.82	109.99
3	E	6	MAN	C1-O5-C5	2.78	115.91	112.19
5	H	2	5N6	C4-C5-N5	-2.73	105.06	110.44
5	M	2	5N6	C9-C8-C7	-2.56	107.39	112.22
3	F	5	MAN	O2-C2-C3	-2.46	105.05	110.15
5	M	1	GAL	O5-C5-C6	-2.46	100.36	106.44
3	E	4	MAN	C1-O5-C5	2.45	115.47	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	3	BMA	O2-C2-C3	-2.36	105.25	110.15
3	F	4	MAN	O2-C2-C3	-2.28	105.43	110.15
5	M	2	5N6	O10-C10-C11	-2.24	118.06	122.05
3	E	3	BMA	O2-C2-C3	-2.21	105.58	110.15
3	E	4	MAN	O2-C2-C3	-2.20	105.58	110.15
5	H	2	5N6	O1A-C1-C2	-2.19	118.12	122.85
5	H	1	GAL	C4-C3-C2	2.16	114.62	110.83
3	E	5	MAN	C1-O5-C5	2.15	115.06	112.19
5	H	1	GAL	O4-C4-C5	2.12	114.56	109.32
4	N	2	NAG	C1-O5-C5	2.12	115.03	112.19
3	F	2	NAG	O4-C4-C5	-2.07	104.23	109.32
3	J	6	MAN	C1-O5-C5	2.06	114.95	112.19
5	H	1	GAL	O5-C1-C2	-2.05	106.69	110.30
4	I	2	NAG	C1-O5-C5	2.03	114.90	112.19
4	K	2	NAG	C1-O5-C5	2.00	114.87	112.19

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	H	2	5N6	C7-C8-C9-O9
5	H	2	5N6	O8-C8-C9-O9
5	H	2	5N6	O6-C6-C7-C8
5	H	2	5N6	O6-C6-C7-O7
5	H	2	5N6	CAF-CAG-O9-C9
5	H	2	5N6	OBJ-CAG-O9-C9
4	G	1	NAG	O5-C5-C6-O6
5	H	1	GAL	C4-C5-C6-O6
5	H	1	GAL	O5-C5-C6-O6
4	G	1	NAG	C4-C5-C6-O6
4	G	2	NAG	O5-C5-C6-O6
3	J	5	MAN	O5-C5-C6-O6
5	M	1	GAL	O5-C5-C6-O6
4	G	2	NAG	C4-C5-C6-O6
3	J	5	MAN	C4-C5-C6-O6
5	M	1	GAL	C4-C5-C6-O6
3	F	3	BMA	O5-C5-C6-O6
5	H	2	5N6	C4-C5-N5-C10
3	J	6	MAN	C4-C5-C6-O6
5	H	2	5N6	C5-C6-C7-O7
5	H	2	5N6	C5-C6-C7-C8
5	M	2	5N6	O1A-C1-C2-O6

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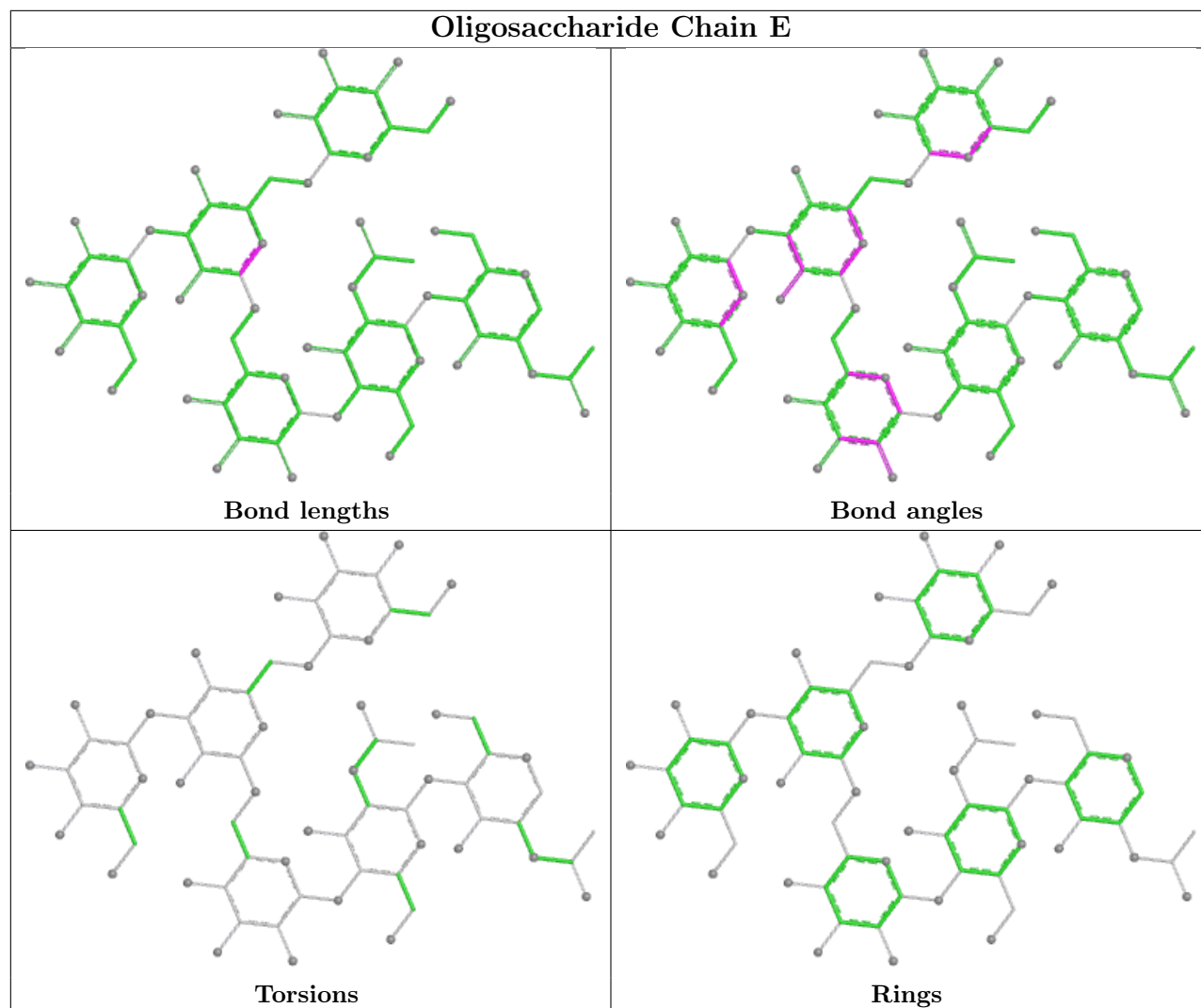
Mol	Chain	Res	Type	Atoms
5	H	2	5N6	C6-C5-N5-C10
3	J	6	MAN	O5-C5-C6-O6
3	F	3	BMA	C4-C5-C6-O6

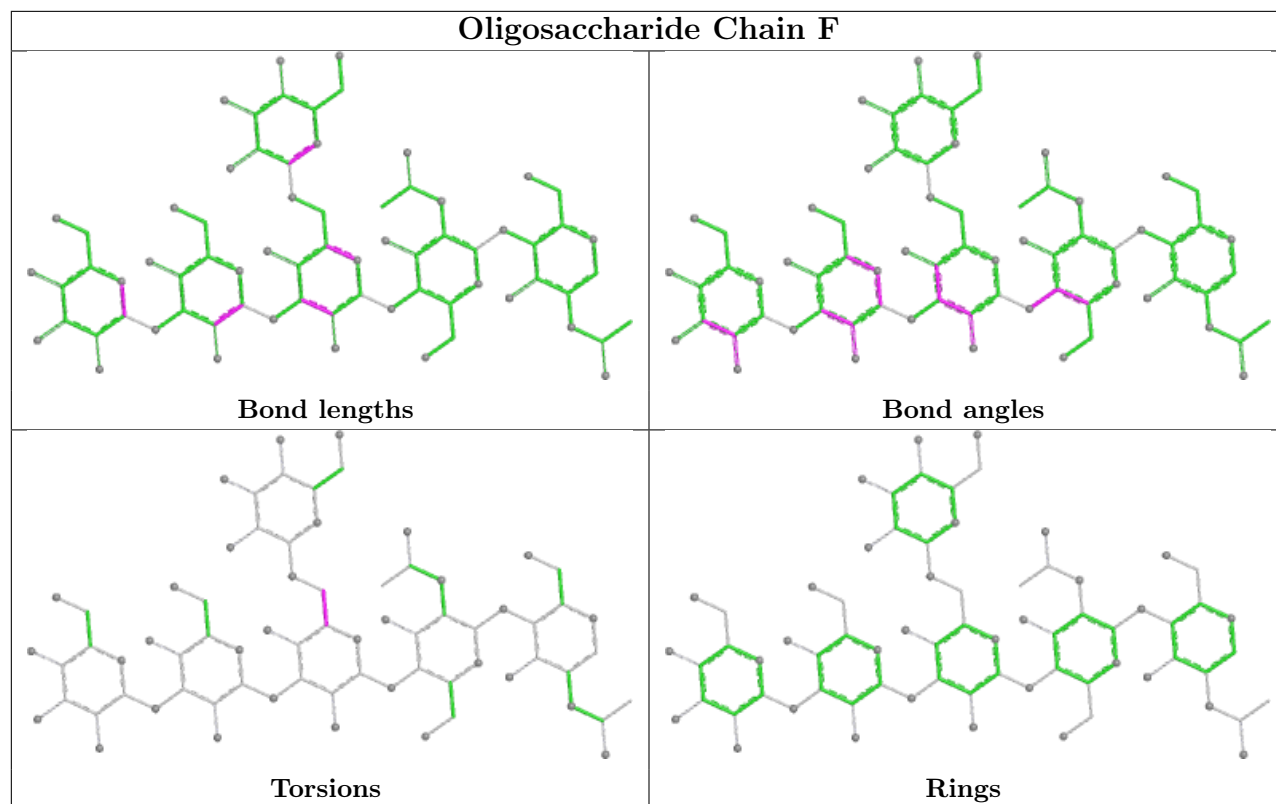
There are no ring outliers.

9 monomers are involved in 7 short contacts:

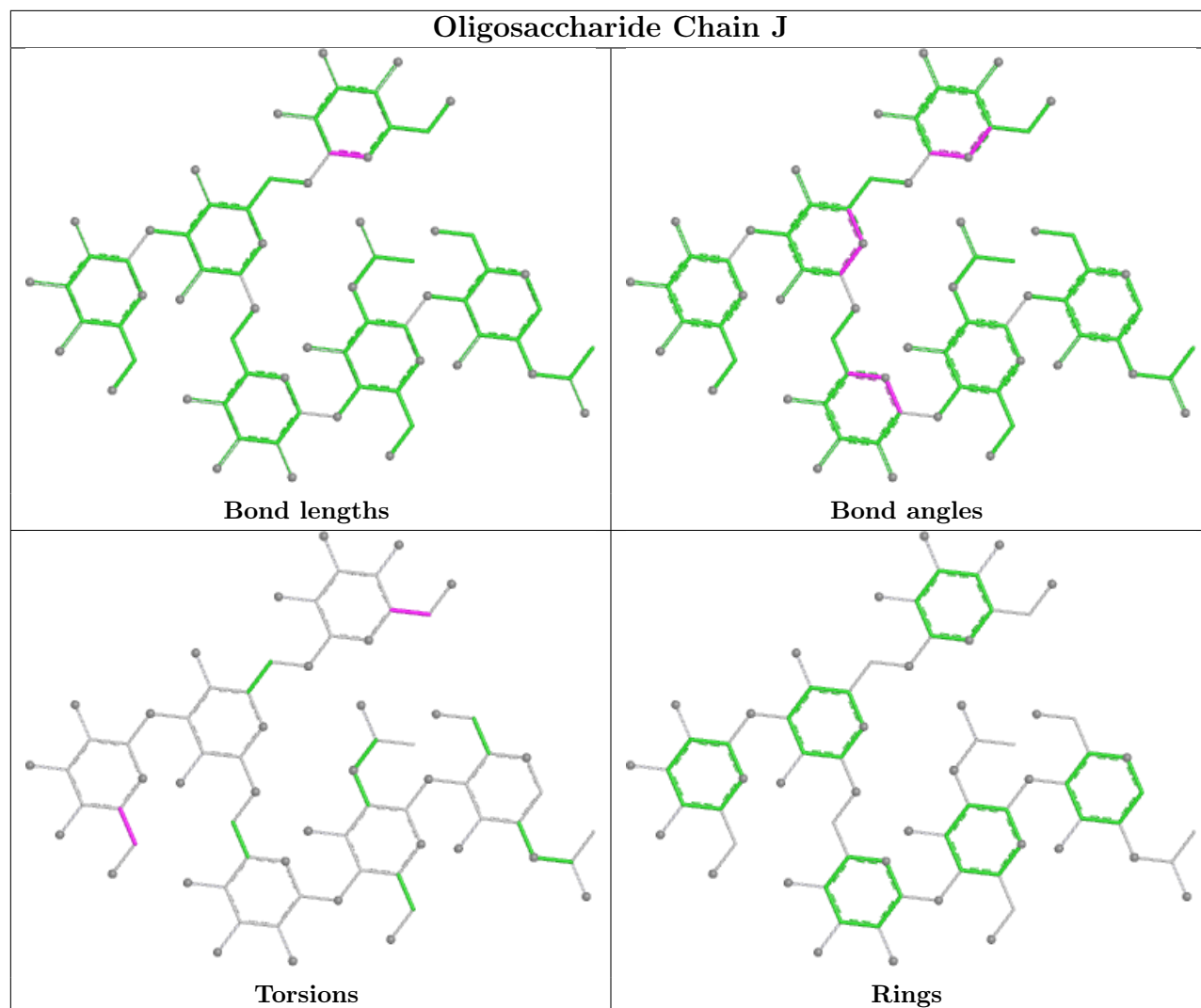
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	6	MAN	1	0
5	M	1	GAL	1	0
3	F	2	NAG	1	0
3	F	3	BMA	1	0
3	E	5	MAN	1	0
5	H	2	5N6	1	0
5	M	2	5N6	1	0
4	I	2	NAG	1	0
4	G	1	NAG	1	0

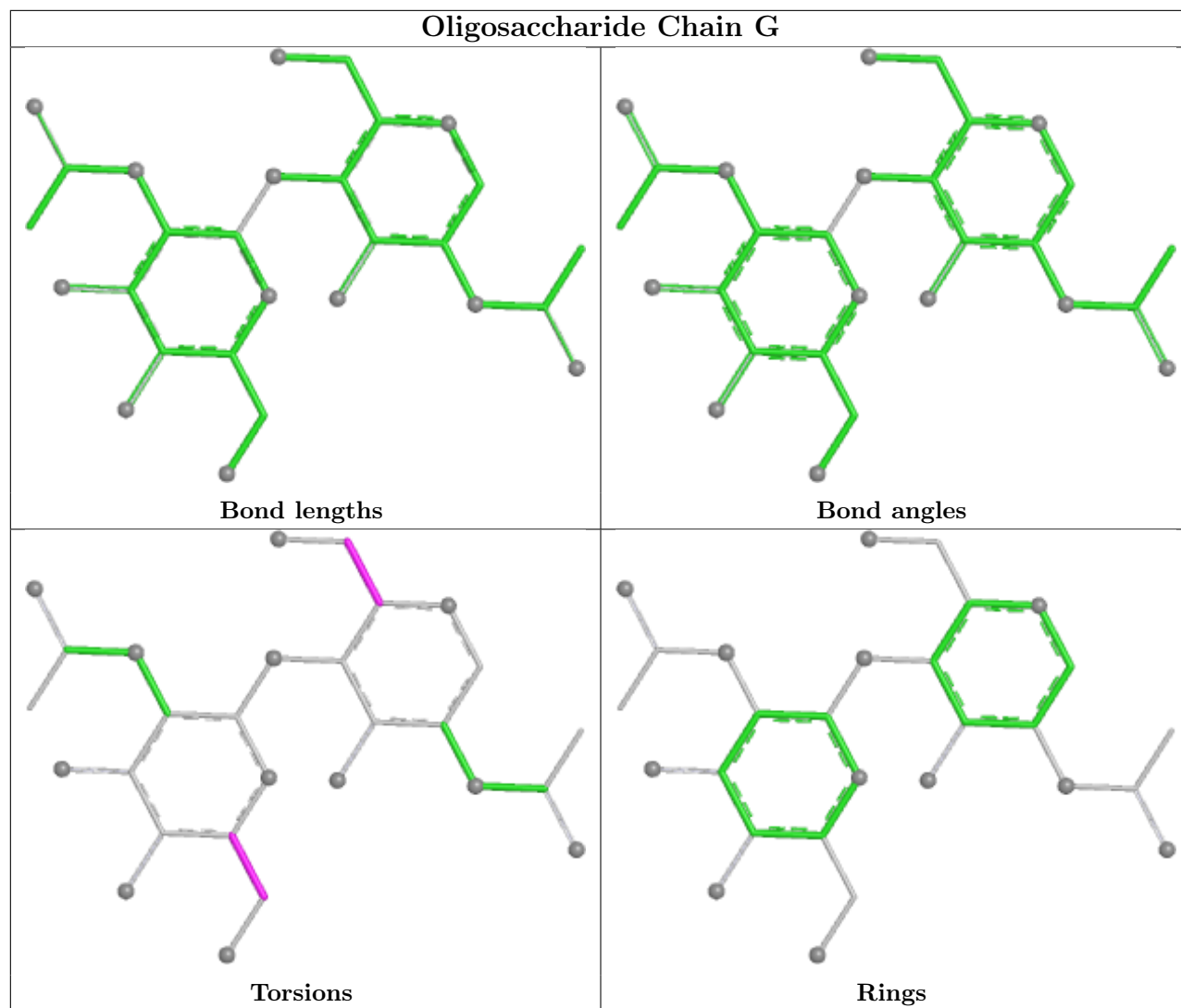
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

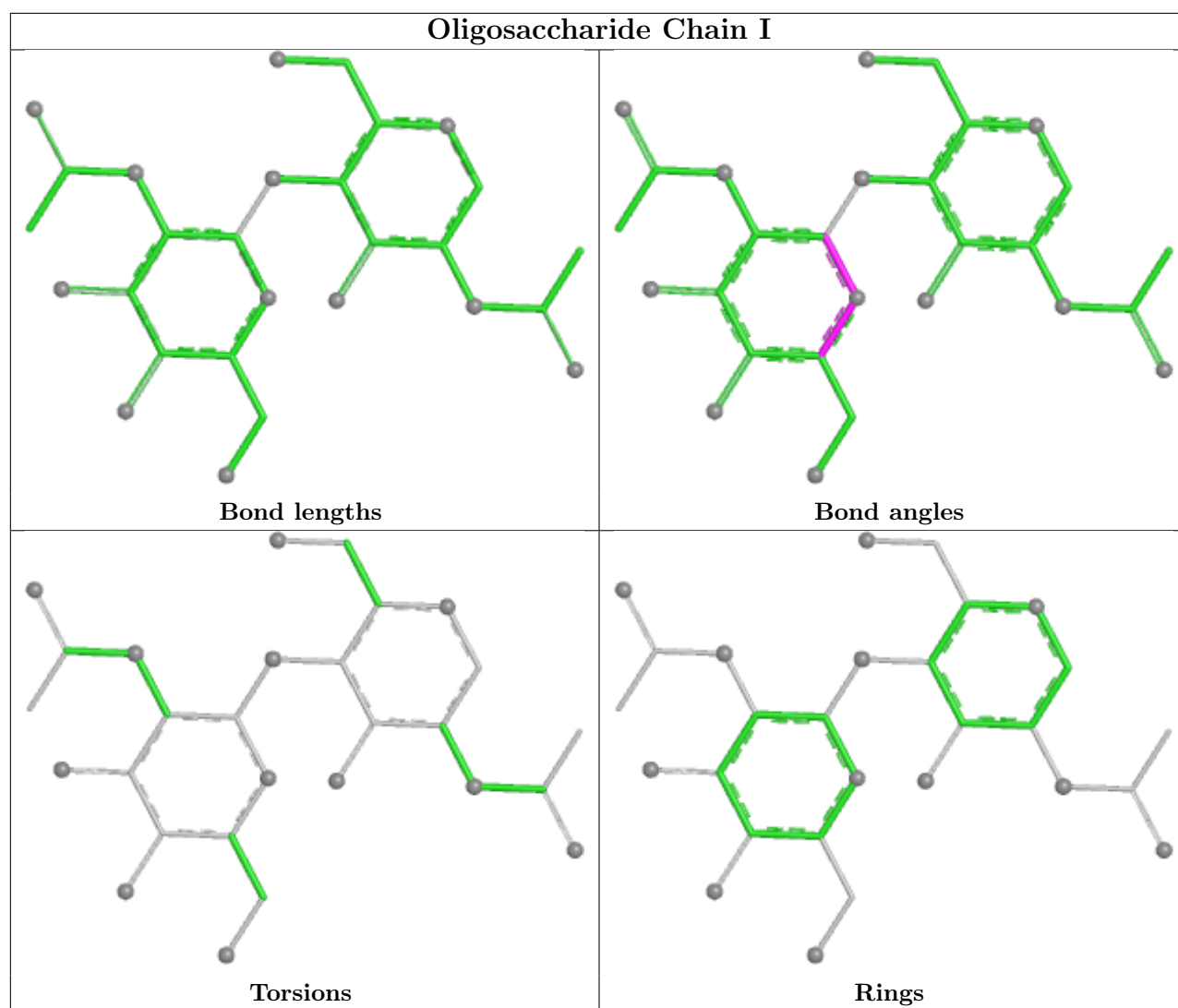


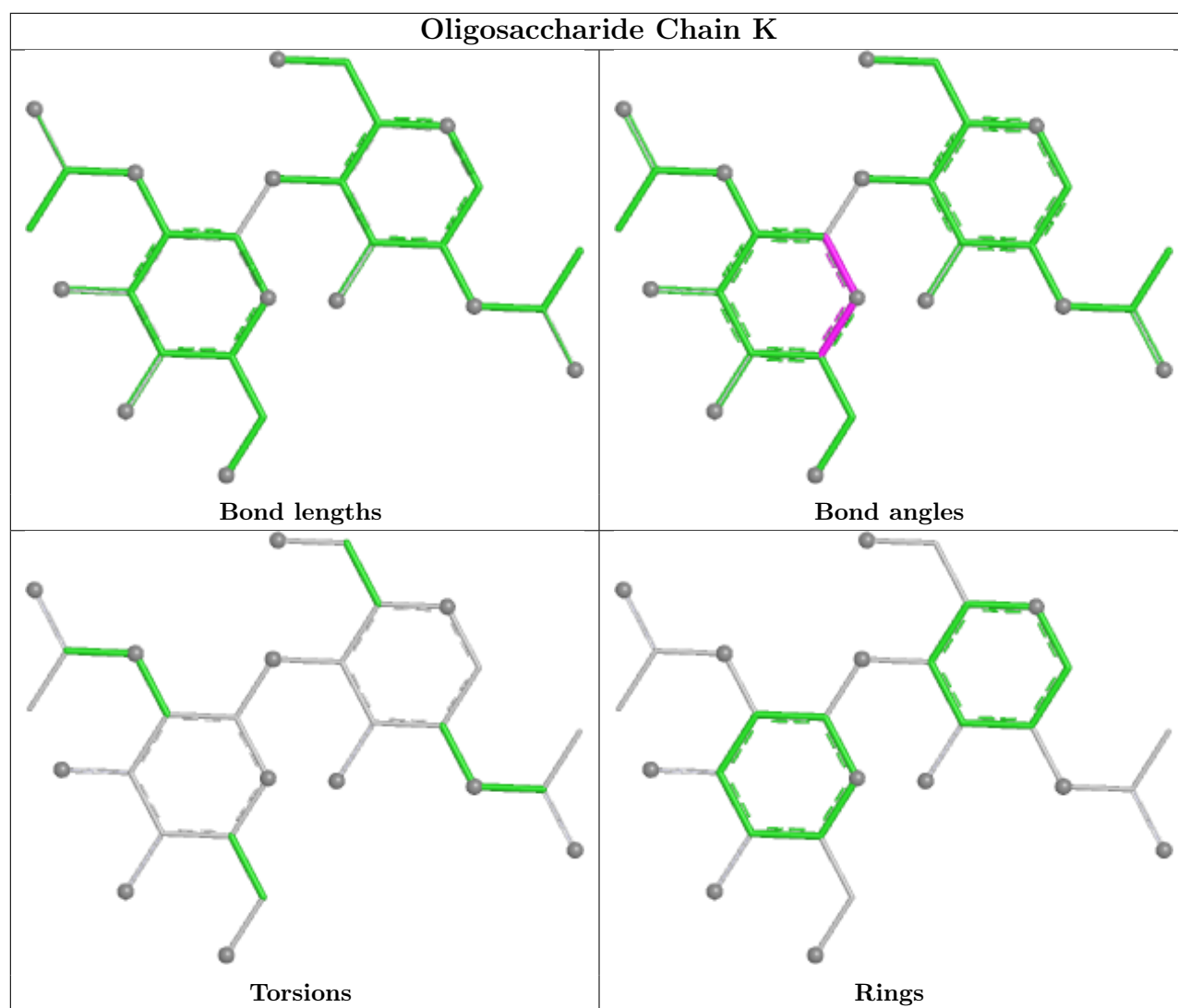


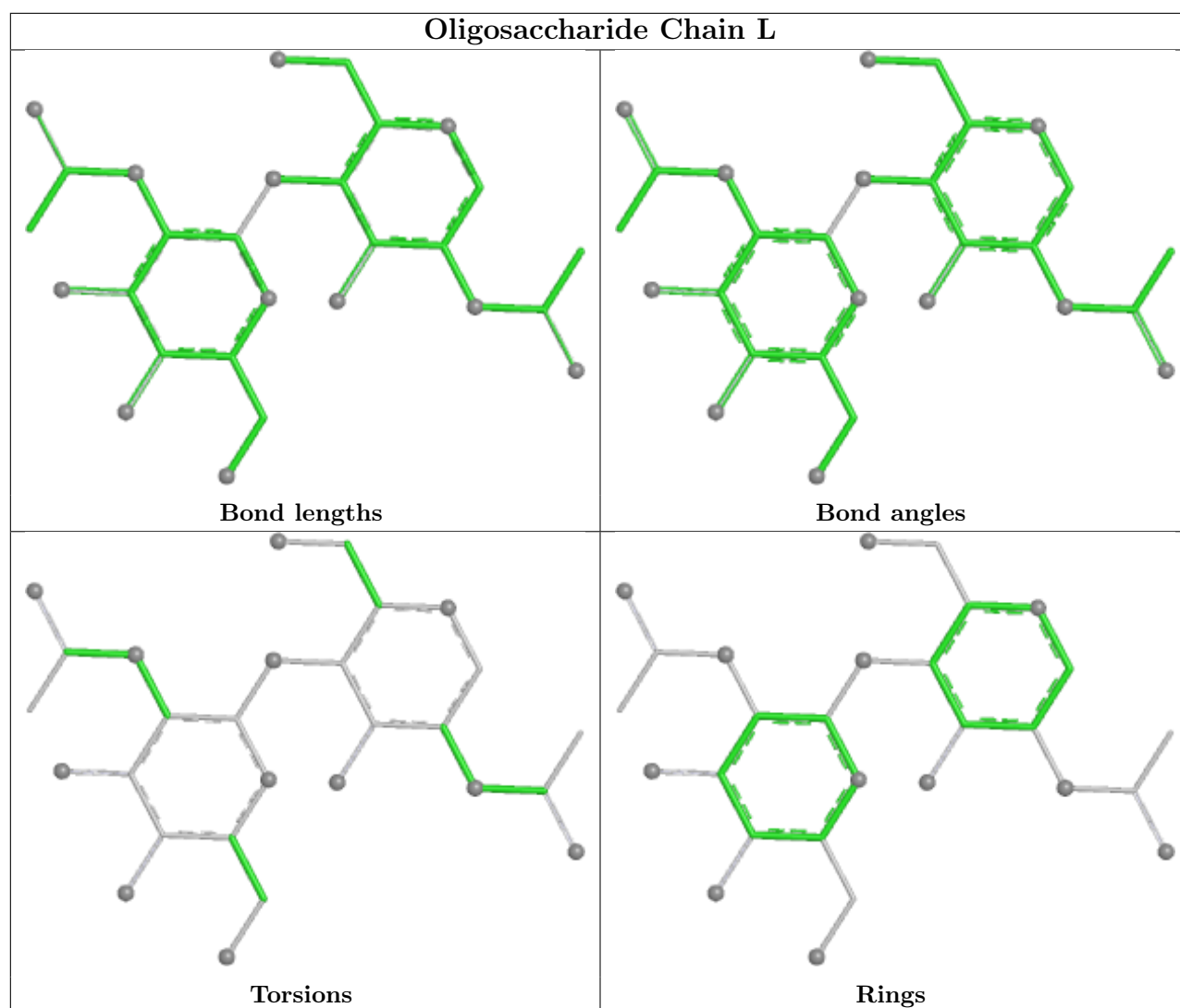
Oligosaccharide Chain J

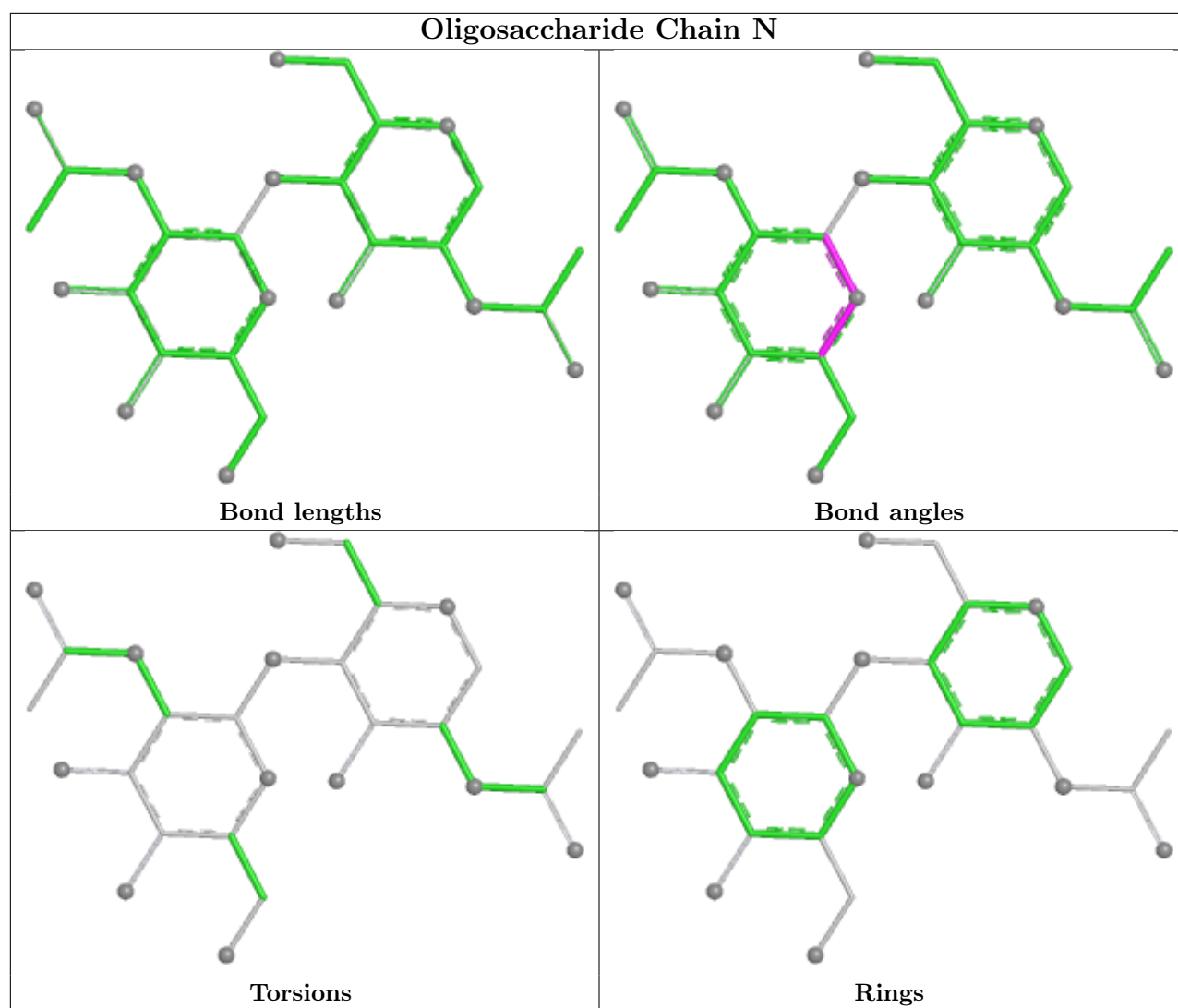




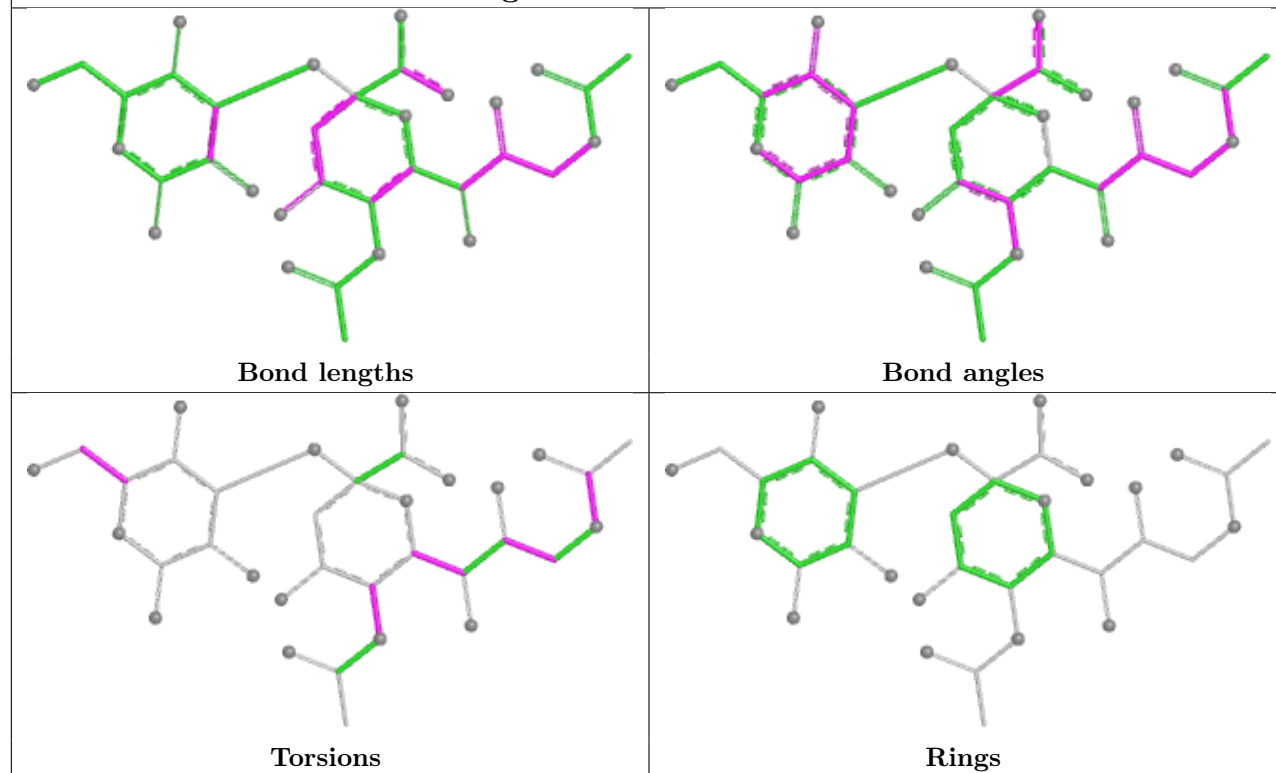




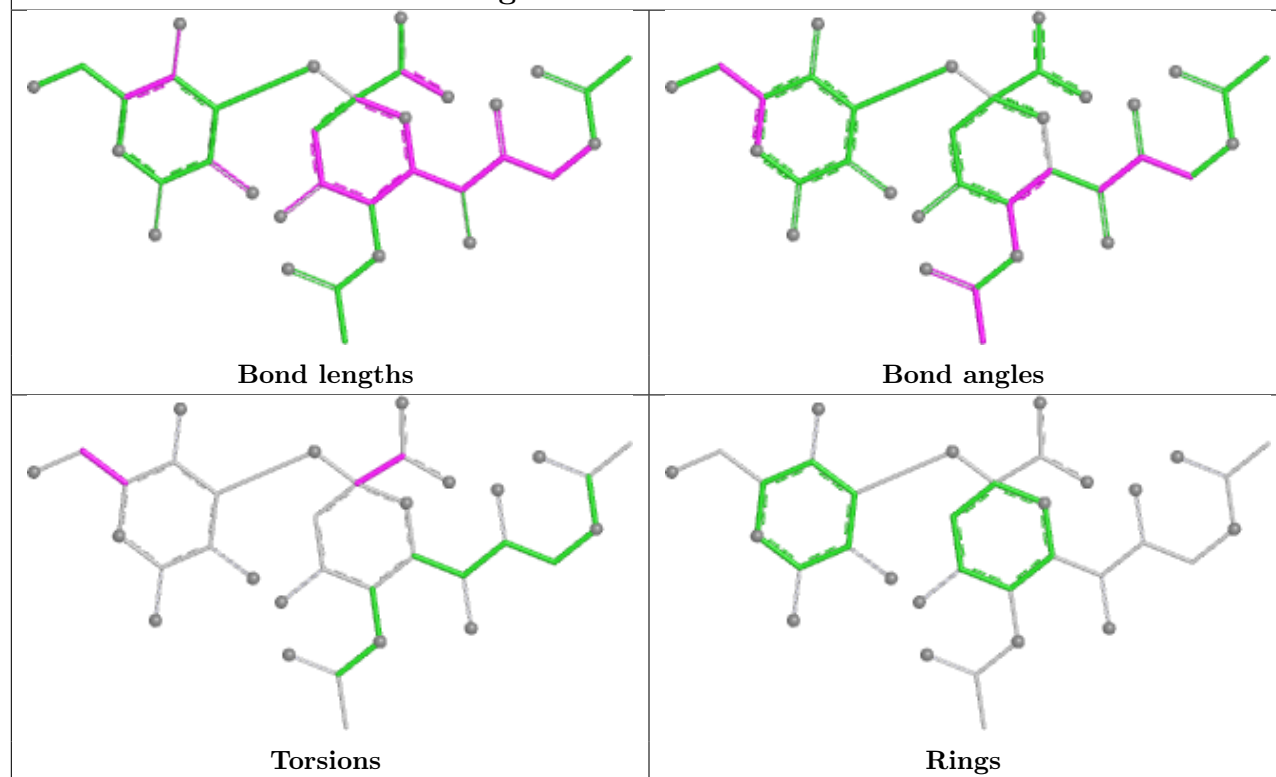




Oligosaccharide Chain H



Oligosaccharide Chain M



5.6 Ligand geometry

2 ligands 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	$\# Z > 2$
6	NAG	B	703	2	14,14,15	0.28	0	17,19,21	0.59	0
6	NAG	D	703	2	14,14,15	0.52	0	17,19,21	1.22	2 (11%)

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
6	NAG	B	703	2	-	2/6/23/26	0/1/1/1
6	NAG	D	703	2	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	D	703	NAG	C1-O5-C5	3.68	117.12	112.19
6	D	703	NAG	O5-C5-C6	-2.14	103.49	107.66

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	703	NAG	C4-C5-C6-O6
6	B	703	NAG	O5-C5-C6-O6
6	D	703	NAG	O5-C5-C6-O6
6	D	703	NAG	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	703	NAG	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	424/427 (99%)	0.28	39 (9%)	14 12	24, 35, 85, 131	0
1	C	424/427 (99%)	-0.01	12 (2%)	55 52	20, 33, 58, 89	0
2	B	149/166 (89%)	1.06	28 (18%)	3 2	22, 48, 128, 201	0
2	D	149/166 (89%)	1.11	37 (24%)	2 1	26, 49, 141, 173	0
All	All	1146/1186 (96%)	0.38	116 (10%)	12 10	20, 37, 91, 201	0

All (116) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	131	ALA	8.4
2	D	44	ALA	6.7
2	D	145	ASN	6.4
2	B	157	PHE	6.4
2	B	44	ALA	6.3
1	A	191	VAL	5.9
2	D	157	PHE	5.3
2	B	34	GLY	5.2
2	B	145	ASN	5.1
2	B	141	VAL	5.1
1	A	265	ALA	5.1
2	D	131	ALA	5.1
1	A	193	PRO	5.0
1	A	199	VAL	4.9
2	D	9	PHE	4.9
1	C	244	ASP	4.9
1	C	243	GLY	4.9
2	B	9	PHE	4.9
1	A	244	ASP	4.8
1	A	197	PRO	4.6
1	A	264	VAL	4.3

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Mol	Chain	Res	Type	RSRZ
2	B	38	SER	4.2
2	B	42	THR	4.1
1	C	193	PRO	4.1
2	D	132	GLN	4.0
1	C	197	PRO	4.0
2	D	141	VAL	4.0
2	D	146	CYS	4.0
2	D	46	PHE	3.9
1	A	237	PHE	3.9
2	B	45	GLY	3.9
2	B	128	GLY	3.8
1	A	235	ALA	3.8
2	B	43	GLN	3.7
1	A	198	GLN	3.6
2	D	33	GLY	3.6
2	B	127	GLY	3.6
2	B	46	PHE	3.6
2	B	37	ALA	3.6
2	D	37	ALA	3.5
1	A	267	TYR	3.5
2	B	33	GLY	3.4
2	D	128	GLY	3.4
1	C	12	ASN	3.4
1	A	241	GLY	3.4
1	C	355	ALA	3.3
1	A	231	TYR	3.3
2	D	143	SER	3.3
2	B	32	GLY	3.3
2	D	43	GLN	3.3
2	D	39	VAL	3.3
2	D	32	GLY	3.2
1	A	190	SER	3.2
2	D	142	ASP	3.2
2	B	18	ALA	3.1
2	B	30	SER	3.1
2	D	148	ALA	3.1
1	A	261	ASP	3.0
2	D	10	GLY	3.0
1	A	266	ALA	2.9
2	D	38	SER	2.9
2	B	49	ILE	2.9
1	A	296	ARG	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	262	ASN	2.8
2	D	16	PHE	2.8
1	A	12	ASN	2.8
2	D	30	SER	2.8
1	A	297	PHE	2.8
2	B	147	ASP	2.8
2	D	149	SER	2.8
1	C	16	SER	2.7
2	B	16	PHE	2.7
2	D	42	THR	2.7
1	A	294	THR	2.7
1	A	429	SER	2.7
1	C	271	ALA	2.6
2	D	57	ARG	2.6
2	D	147	ASP	2.6
1	A	239	THR	2.6
1	A	200	CYS	2.6
1	A	232	GLU	2.5
2	B	19	GLY	2.5
1	A	165	THR	2.5
2	D	31	ASN	2.5
2	D	34	GLY	2.4
1	A	240	PHE	2.4
2	D	80	ASN	2.4
2	B	10	GLY	2.4
1	A	236	LYS	2.4
2	B	130	ILE	2.4
1	C	245	TYR	2.3
2	D	127	GLY	2.3
2	D	19	GLY	2.3
1	A	245	TYR	2.3
2	D	78	ILE	2.3
1	C	114	GLU	2.2
1	C	409	ASP	2.2
1	A	238	ASN	2.2
1	A	202	THR	2.2
2	B	31	ASN	2.2
1	C	198	GLN	2.2
2	D	35	GLY	2.2
2	D	36	GLY	2.2
2	B	143	SER	2.2
2	D	130	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	295	PRO	2.1
2	B	36	GLY	2.1
1	A	169	ALA	2.1
2	D	18	ALA	2.1
1	A	192	ASN	2.1
1	A	357	ASP	2.1
1	A	268	ARG	2.1
1	A	325	SER	2.1
2	D	45	GLY	2.1
1	A	242	CYS	2.0
1	A	234	LYS	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.

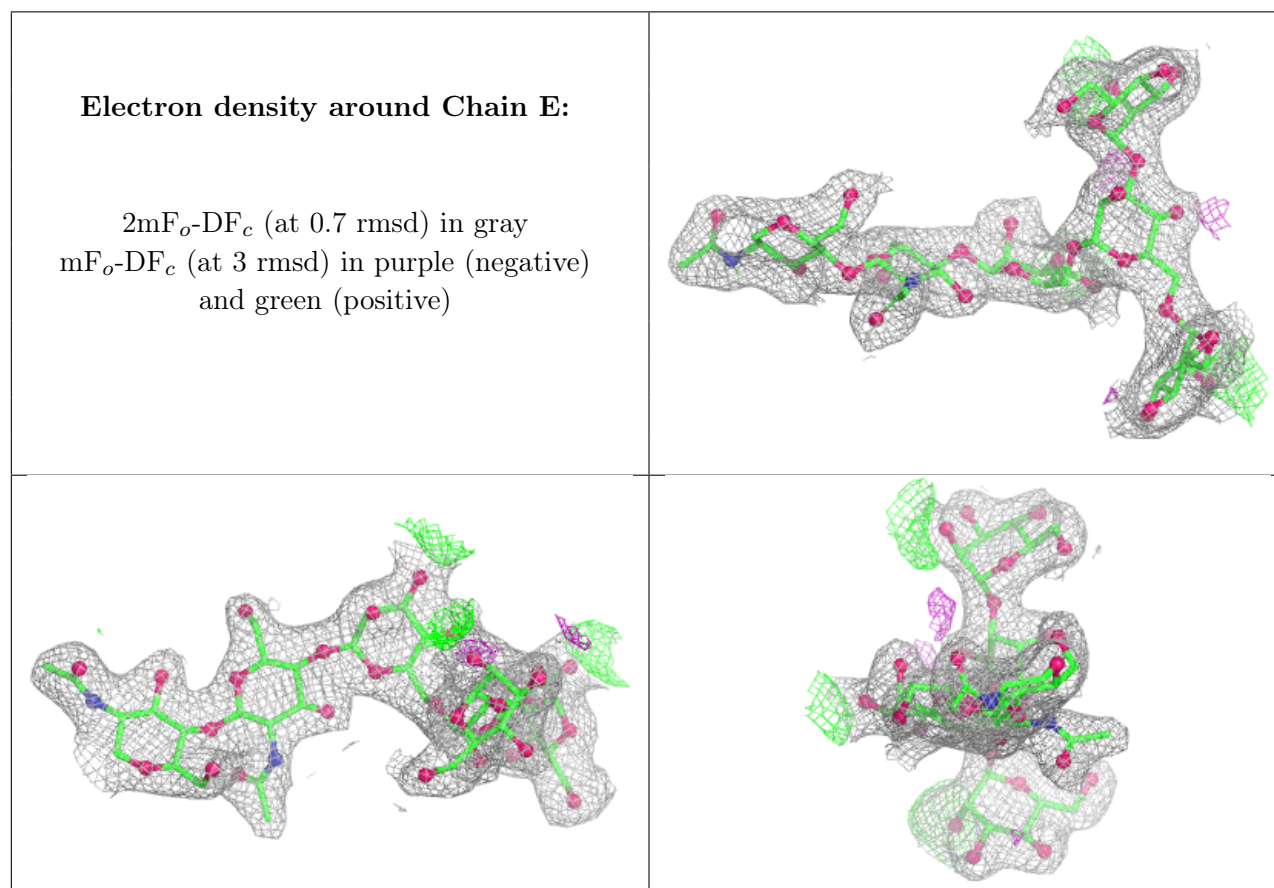
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	GAL	H	1	12/12	0.56	0.15	96,109,121,131	0
3	BMA	F	3	11/12	0.73	0.17	39,43,52,56	0
5	GAL	M	1	12/12	0.77	0.13	68,88,102,113	0
4	NAG	K	2	14/15	0.79	0.14	39,47,57,57	0
4	NAG	G	2	14/15	0.83	0.12	56,63,76,85	0
3	MAN	E	6	11/12	0.84	0.10	31,36,39,47	0
4	NAG	L	2	14/15	0.84	0.12	45,50,60,72	0
5	5N6	H	2	23/24	0.86	0.16	71,81,85,87	0
3	MAN	F	5	11/12	0.87	0.10	43,47,54,57	0
4	NAG	N	2	14/15	0.89	0.09	34,40,45,47	0
3	MAN	F	4	11/12	0.89	0.08	38,41,47,49	0
3	MAN	J	5	11/12	0.89	0.09	31,34,38,41	0
4	NAG	G	1	14/15	0.89	0.11	53,56,62,63	0
4	NAG	I	2	14/15	0.90	0.09	36,40,47,53	0
3	MAN	E	5	11/12	0.90	0.08	30,34,37,40	0
4	NAG	L	1	14/15	0.90	0.10	43,48,63,65	0
3	NAG	F	2	14/15	0.90	0.09	29,35,40,44	0

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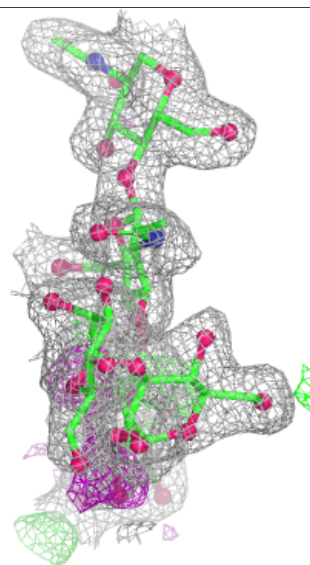
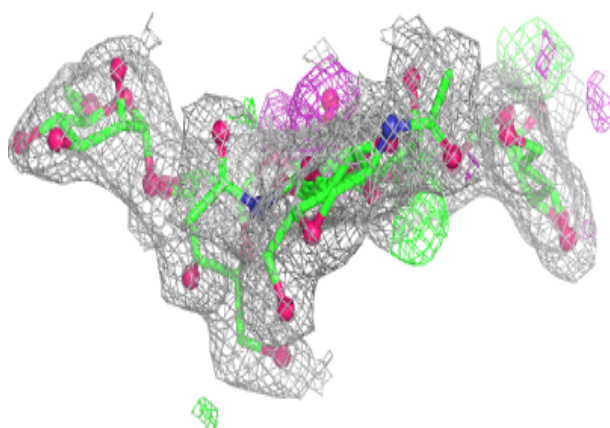
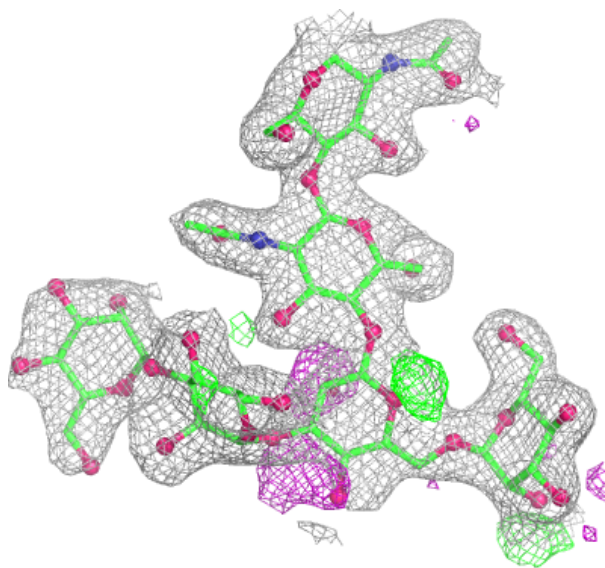
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	MAN	F	6	11/12	0.91	0.08	36,39,44,47	0
3	MAN	J	6	11/12	0.91	0.09	36,40,45,52	0
3	NAG	F	1	14/15	0.92	0.08	27,33,36,36	0
3	NAG	E	2	14/15	0.93	0.09	28,35,38,40	0
3	NAG	E	1	14/15	0.94	0.07	33,36,44,49	0
3	BMA	E	3	11/12	0.94	0.06	29,32,35,38	0
3	NAG	J	2	14/15	0.94	0.07	24,29,35,37	0
3	BMA	J	3	11/12	0.94	0.07	29,33,36,41	0
5	5N6	M	2	23/24	0.94	0.09	43,52,60,62	0
4	NAG	I	1	14/15	0.95	0.07	27,33,35,37	0
4	NAG	K	1	14/15	0.95	0.06	23,30,34,40	0
4	NAG	N	1	14/15	0.96	0.06	27,30,34,35	0
3	MAN	E	4	11/12	0.96	0.06	26,30,33,38	0
3	NAG	J	1	14/15	0.96	0.05	30,34,36,39	0
3	MAN	J	4	11/12	0.97	0.05	23,30,35,40	0

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



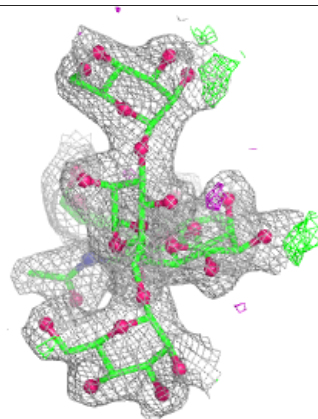
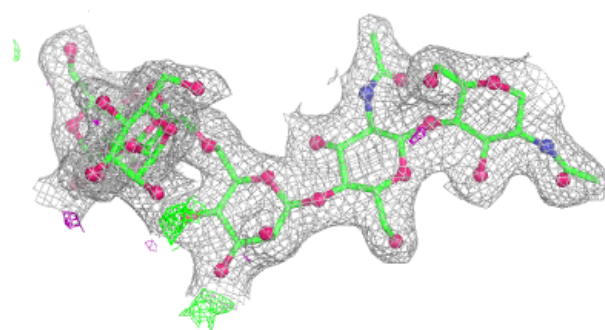
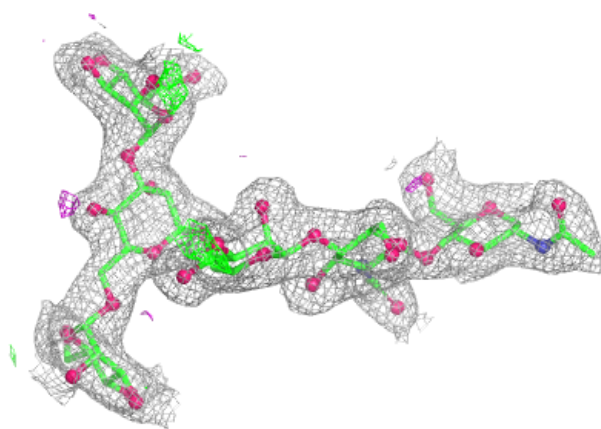
Electron density around Chain F:

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

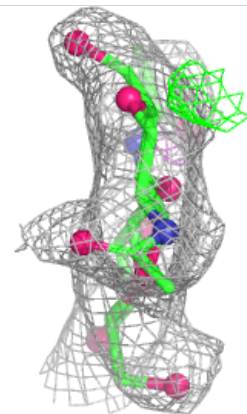
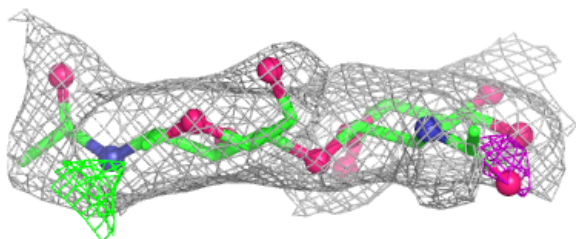
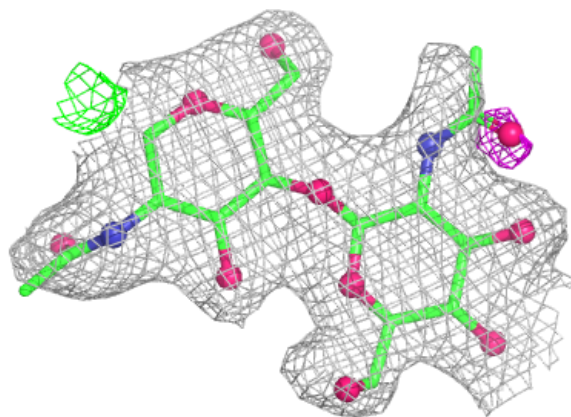


Electron density around Chain J:

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

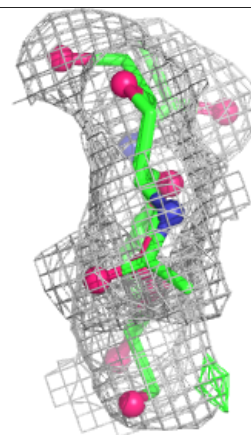
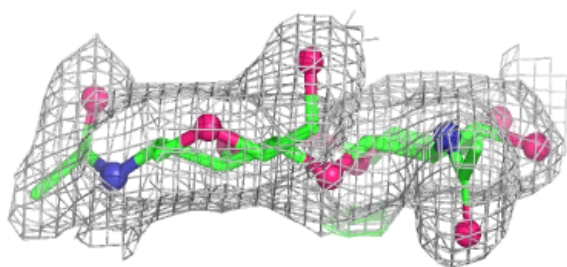
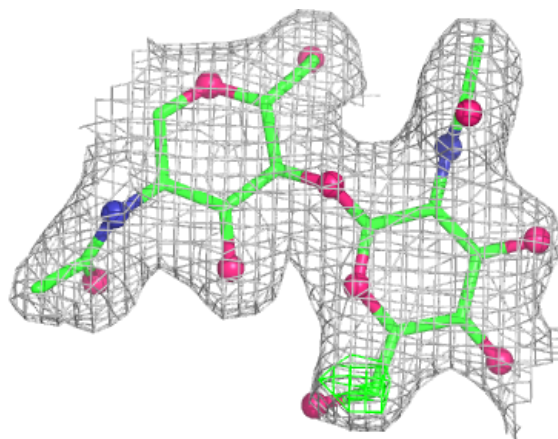
**Electron density around Chain G:**

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



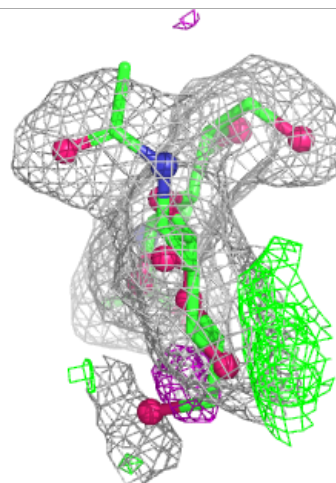
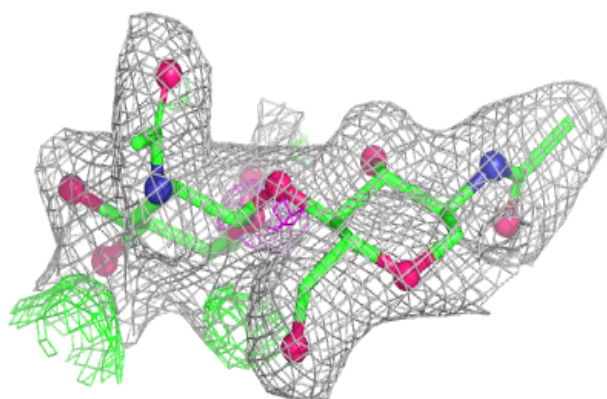
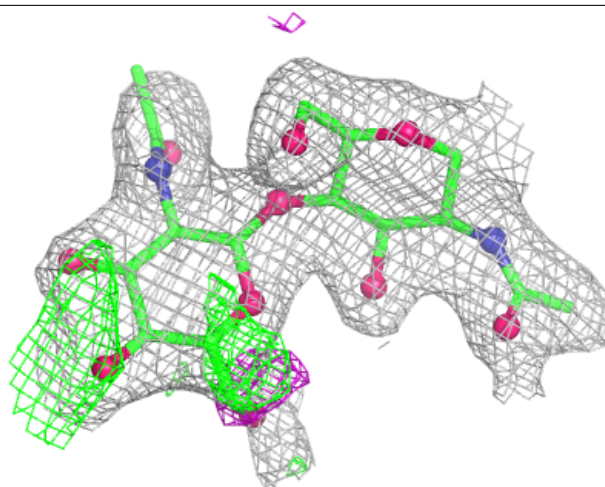
Electron density around Chain I:

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



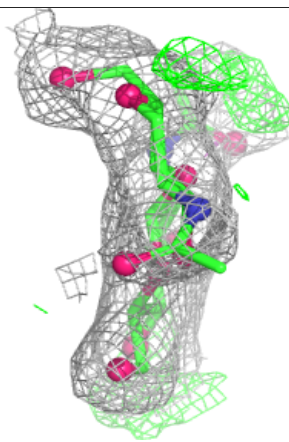
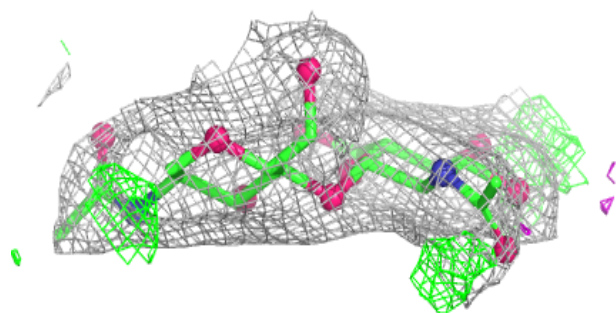
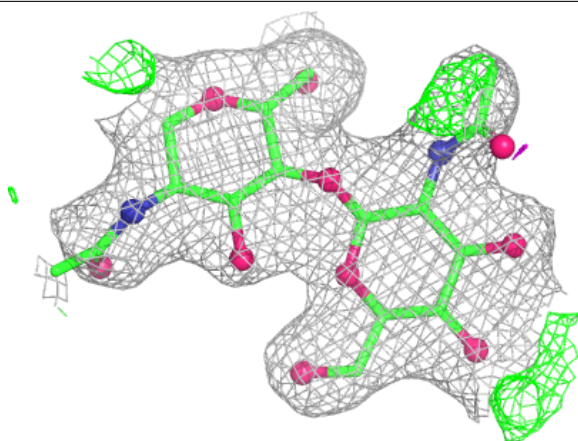
Electron density around Chain K:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



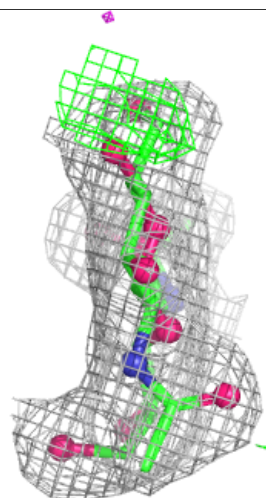
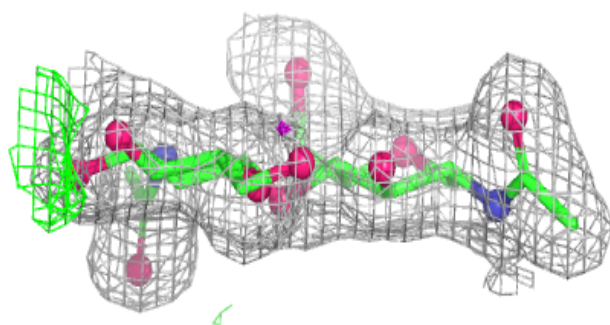
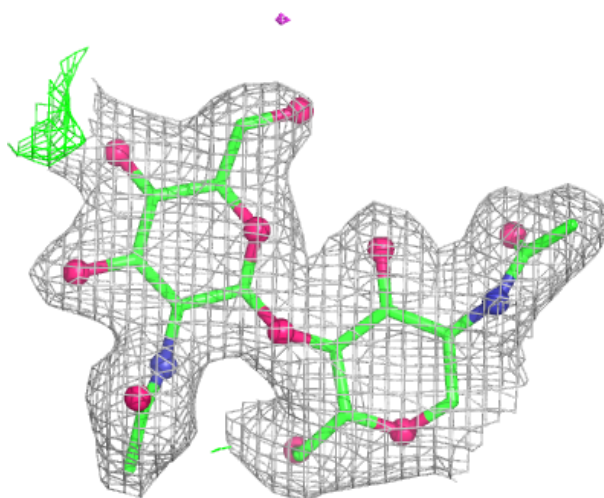
Electron density around Chain L:

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



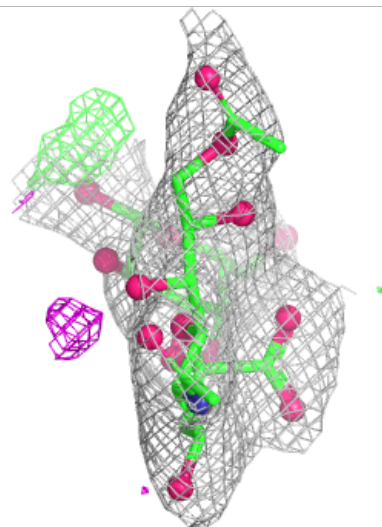
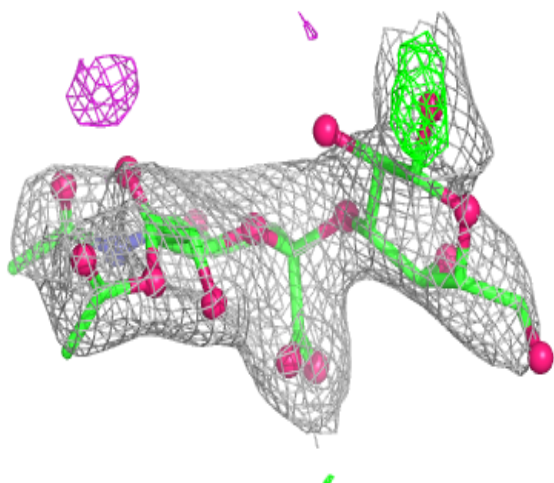
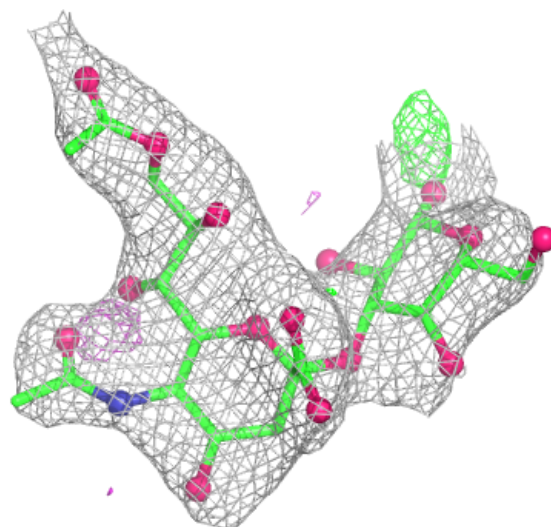
Electron density around Chain N:

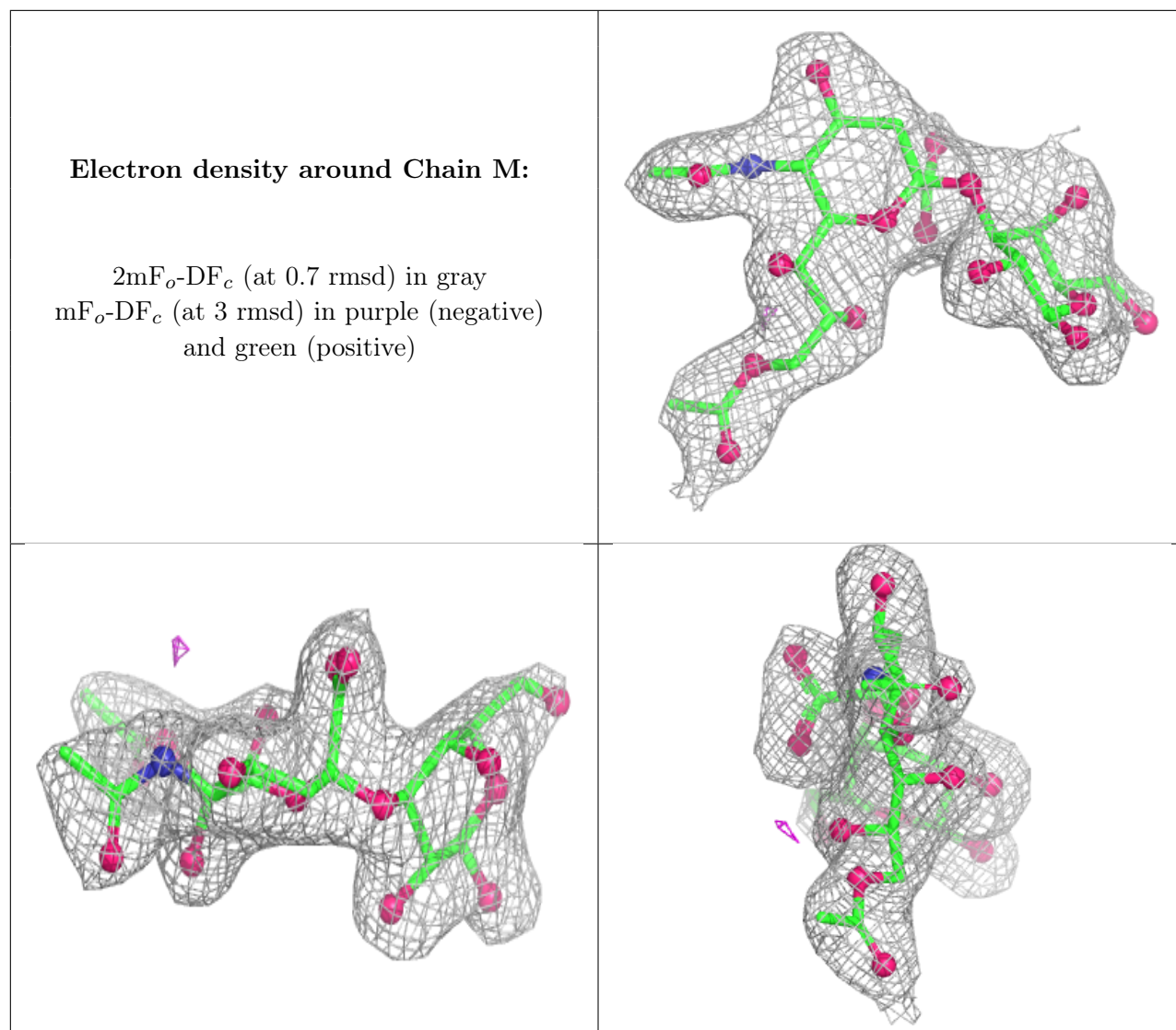
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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)





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(Å ²)	Q<0.9
6	NAG	D	703	14/15	0.46	0.16	71,79,82,83	0
6	NAG	B	703	14/15	0.69	0.14	51,59,65,66	0

6.5 Other polymers [i](#)

There are no such residues in this entry.