



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 15, 2024 – 02:49 pm GMT

PDB ID : 6QP8
Title : Drosophila Semaphorin 2b
Authors : Robinson, R.A.; Rozbesky, D.; Harlos, K.; Siebold, C.; Jones, E.Y.
Deposited on : 2019-02-13
Resolution : 2.33 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 : 1.8.4, CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

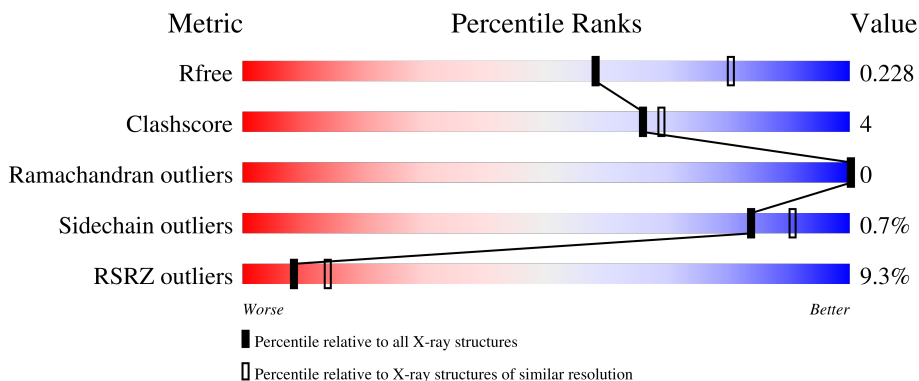
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.33 Å.

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



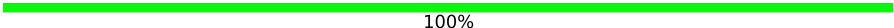

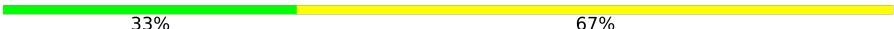



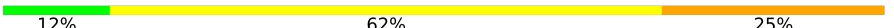

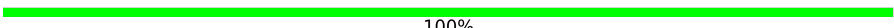
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2096 (2.36-2.32)
Clashscore	141614	2193 (2.36-2.32)
Ramachandran outliers	138981	2159 (2.36-2.32)
Sidechain outliers	138945	2160 (2.36-2.32)
RSRZ outliers	127900	2067 (2.36-2.32)

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	657	 8% 86% 9% 5%
1	B	657	 10% 86% 8% 6%
2	C	2	 100%
2	D	2	 100%
2	I	2	 100%

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Mol	Chain	Length	Quality of chain
2	J	2	 100%
3	E	5	 60% 40%
4	F	9	 33% 67%
5	G	6	 50% 50%
6	H	4	 75% 25%
6	K	4	 75% 25%
7	L	8	 12% 62% 25%
8	M	4	 25% 75%
9	N	3	 100%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NAG	C	2	-	-	-	X

2 Entry composition i

There are 13 unique types of molecules in this entry. The entry contains 11335 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 IP13724p.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	622	4974	3151	862	929	32	0	0	0
1	B	620	4959	3142	858	927	32	0	0	0

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	32	GLU	-	expression tag	UNP Q29QV5
A	33	THR	-	expression tag	UNP Q29QV5
A	34	GLY	-	expression tag	UNP Q29QV5
A	680	GLY	-	expression tag	UNP Q29QV5
A	681	THR	-	expression tag	UNP Q29QV5
A	682	LYS	-	expression tag	UNP Q29QV5
A	683	HIS	-	expression tag	UNP Q29QV5
A	684	HIS	-	expression tag	UNP Q29QV5
A	685	HIS	-	expression tag	UNP Q29QV5
A	686	HIS	-	expression tag	UNP Q29QV5
A	687	HIS	-	expression tag	UNP Q29QV5
A	688	HIS	-	expression tag	UNP Q29QV5
B	32	GLU	-	expression tag	UNP Q29QV5
B	33	THR	-	expression tag	UNP Q29QV5
B	34	GLY	-	expression tag	UNP Q29QV5
B	680	GLY	-	expression tag	UNP Q29QV5
B	681	THR	-	expression tag	UNP Q29QV5
B	682	LYS	-	expression tag	UNP Q29QV5
B	683	HIS	-	expression tag	UNP Q29QV5
B	684	HIS	-	expression tag	UNP Q29QV5
B	685	HIS	-	expression tag	UNP Q29QV5
B	686	HIS	-	expression tag	UNP Q29QV5
B	687	HIS	-	expression tag	UNP Q29QV5
B	688	HIS	-	expression tag	UNP Q29QV5

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



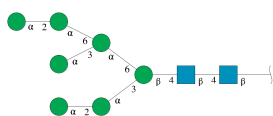
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	C	2	28	16	2	10	0	0	0
2	D	2	28	16	2	10	0	0	0
2	I	2	28	16	2	10	0	0	0
2	J	2	28	16	2	10	0	0	0

- Molecule 3 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	E	5	61	34	2	25	0	0	0

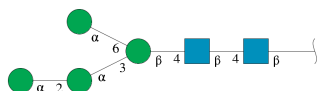
- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
4	F	9	105	58	2	45	0	0	0

- Molecule 5 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-

beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



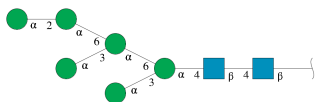
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
5	G	6	72	40	2	30	0	0	0

- Molecule 6 is an oligosaccharide called alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
6	H	4	50	28	2	20	0	0	0
6	K	4	50	28	2	20	0	0	0

- Molecule 7 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



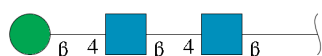
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
7	L	8	94	52	2	40	0	0	0

- Molecule 8 is an oligosaccharide called alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



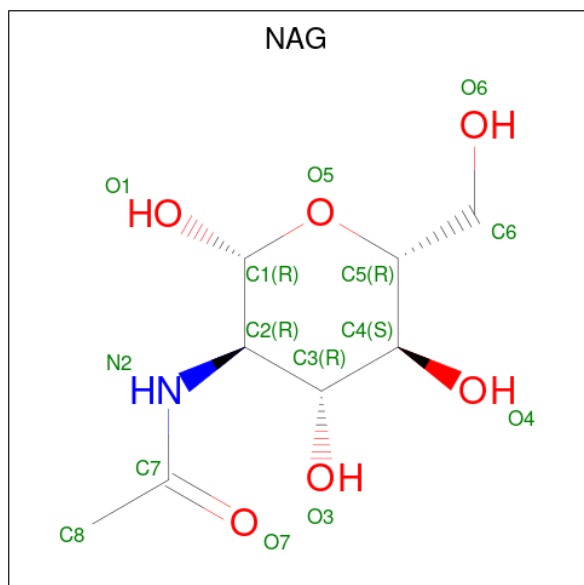
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
8	M	4	50	28	2	20	0	0	0

- Molecule 9 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
			Total	C	N	O			
9	N	3	39	22	2	15	0	0	0

- Molecule 10 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



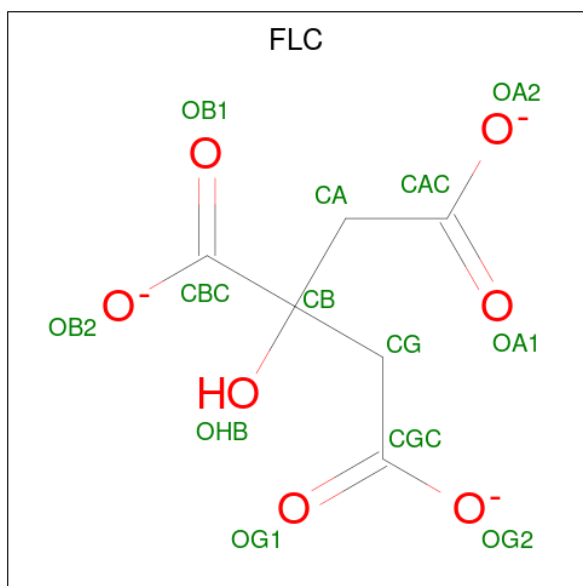
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
10	A	1	14	8	1	5	0	0
10	B	1	14	8	1	5	0	0

- Molecule 11 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
11	A	1	Total C O 6 3 3	0	0
11	B	1	Total C O 6 3 3	0	0
11	B	1	Total C O 6 3 3	0	0

- Molecule 12 is CITRATE ANION (three-letter code: FLC) (formula: $C_6H_5O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
12	B	1	Total	C	O	0	0
			13	6	7		

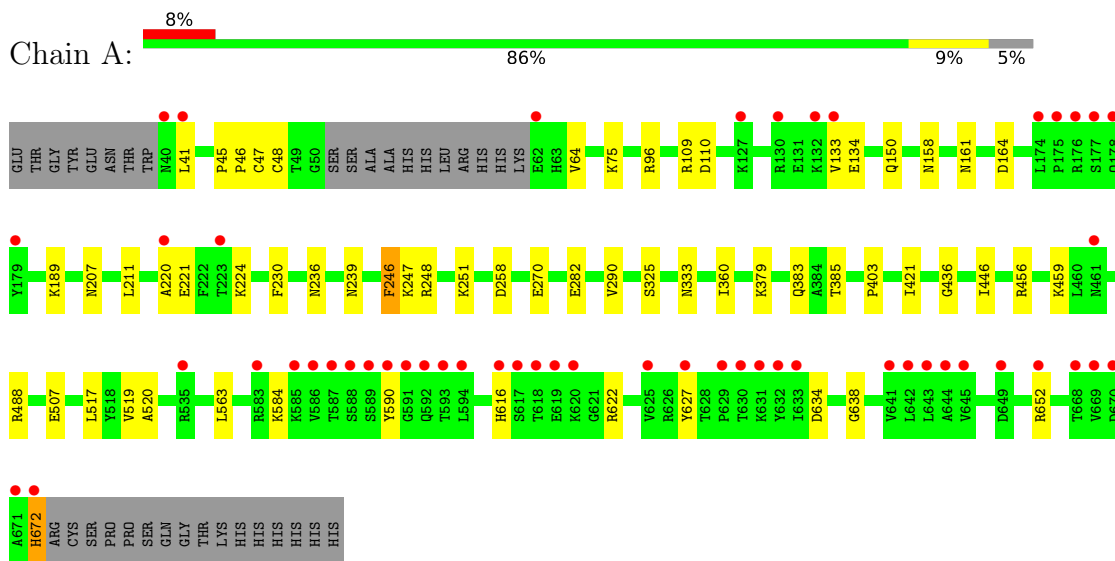
- Molecule 13 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
13	A	357	Total	O	0	0
			357	357		
13	B	353	Total	O	0	0
			353	353		

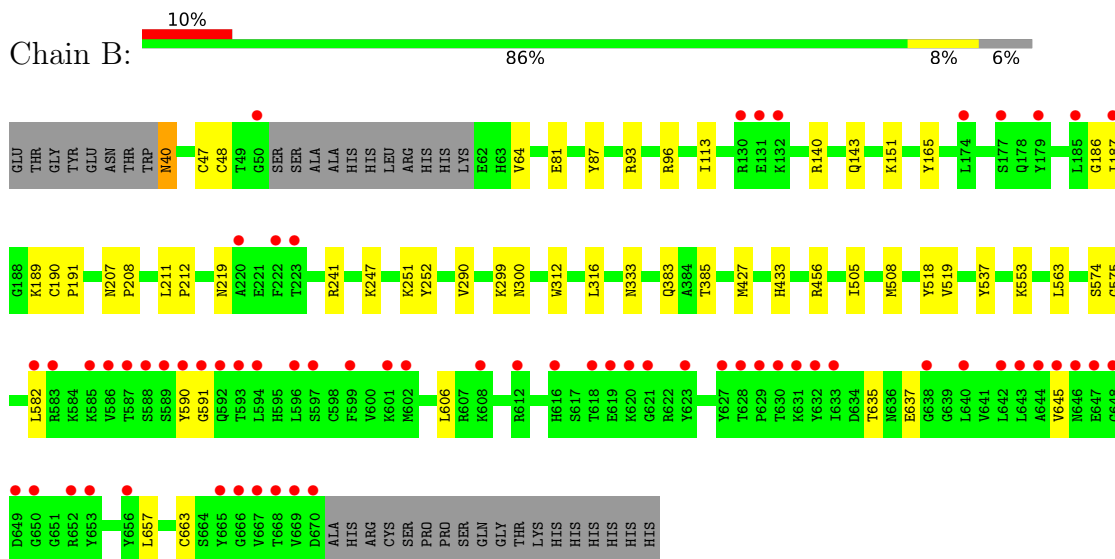
3 Residue-property plots [i](#)

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: IP13724p



- Molecule 1: IP13724p



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

Chain C:  100%

WAG1
WAG2

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

Chain D:  100%

WAG1
WAG2

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

Chain I:  100%

WAG1
WAG2

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

Chain J:  100%

WAG1
WAG2

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

Chain E:  60% 40%

WAG1
WAG2
BMA3
MAN4
MAN5

- Molecule 4: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-2)-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:  33% 67%

WAG1
WAG2
BMA3
MAN4
MAN5
MAN6
MAN7
MAN8
MAN9

- Molecule 5: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:  50% 50%


MAG1
MAG2
BMA3
MAN4
MAN5
MAN6

- Molecule 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

Chain H:  75% 25%

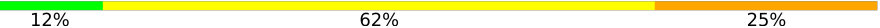
MAG1
MAG2
BMA3
MAN4

- Molecule 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

Chain K:  75% 25%

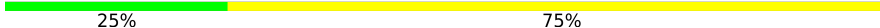
MAG1
MAG2
BMA3
MAN4

- Molecule 7: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain L:  12% 62% 25%

MAG1
MAG2
MAN3
MAN4
MAN5
MAN6
MAN7
MAN8

- Molecule 8: alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M:  25% 75%

MAG1
MAG2
MAN3
MAN4

- Molecule 9: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N:  100%

MAG1
MAG2
BMA3

4 Data and refinement statistics i

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	107.99Å 145.58Å 152.20Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	65.67 – 2.33 86.73 – 2.33	Depositor EDS
% Data completeness (in resolution range)	98.4 (65.67-2.33) 98.4 (86.73-2.33)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 2.32Å)	Xtrriage
Refinement program	PHENIX (dev_3386: ???)	Depositor
R, R_{free}	0.195 , 0.228 0.196 , 0.228	Depositor DCC
R_{free} test set	5076 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	40.7	Xtrriage
Anisotropy	0.304	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 48.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.001 for -h,l,k	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11335	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.16% 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: BMA, FLC, MAN, GOL, 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	0.26	0/5095	0.46	0/6902
1	B	0.26	0/5079	0.46	0/6880
All	All	0.26	0/10174	0.46	0/13782

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	4974	0	4815	45	0
1	B	4959	0	4803	37	0
2	C	28	0	25	0	0
2	D	28	0	25	0	0
2	I	28	0	25	0	0
2	J	28	0	25	0	0
3	E	61	0	52	0	0
4	F	105	0	88	0	0
5	G	72	0	61	0	0
6	H	50	0	43	0	0
6	K	50	0	43	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	L	94	0	79	2	0
8	M	50	0	43	1	0
9	N	39	0	34	0	0
10	A	14	0	13	0	0
10	B	14	0	13	0	0
11	A	6	0	8	1	0
11	B	12	0	16	0	0
12	B	13	0	5	0	0
13	A	357	0	0	14	1
13	B	353	0	0	18	1
All	All	11335	0	10216	84	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 4.

All (84) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:312:TRP:O	13:B:802:HOH:O	1.89	0.89
1:B:456:ARG:NH1	13:B:805:HOH:O	2.06	0.87
1:B:151:LYS:O	13:B:803:HOH:O	2.00	0.79
1:A:134:GLU:N	13:A:804:HOH:O	2.15	0.79
1:B:316:LEU:HD22	1:B:427:MET:HE1	1.64	0.79
1:A:220:ALA:O	13:A:801:HOH:O	2.01	0.78
1:B:212:PRO:O	13:B:804:HOH:O	2.03	0.77
1:B:143:GLN:OE1	13:B:806:HOH:O	2.08	0.69
13:B:807:HOH:O	7:L:8:MAN:O6	2.09	0.69
1:A:459:LYS:HG2	1:A:563:LEU:HD11	1.74	0.68
1:A:248:ARG:HH12	1:A:251:LYS:HB2	1.59	0.68
1:A:325:SER:OG	1:A:436:GLY:O	2.12	0.68
1:A:282:GLU:O	13:A:803:HOH:O	2.12	0.66
1:A:164:ASP:OD2	13:A:802:HOH:O	2.12	0.66
1:B:563:LEU:O	13:B:808:HOH:O	2.13	0.66
1:B:635:THR:HG22	1:B:637:GLU:H	1.60	0.66
1:A:488:ARG:NH1	13:A:806:HOH:O	2.19	0.64
1:A:246:PHE:CE1	1:A:248:ARG:HD2	2.33	0.63
1:B:219:ASN:OD1	13:B:809:HOH:O	2.15	0.62
1:A:590:TYR:HB3	1:A:672:HIS:HB3	1.82	0.61
1:B:48:CYS:HB3	1:B:64:VAL:HG21	1.82	0.61
1:A:627:TYR:CZ	1:A:634:ASP:HB3	2.36	0.61
1:A:270:GLU:OE1	13:A:805:HOH:O	2.17	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:B:810:HOH:O	8:M:1:NAG:O6	2.16	0.59
1:B:81:GLU:OE2	13:B:811:HOH:O	2.17	0.58
1:A:133:VAL:N	13:A:804:HOH:O	2.34	0.57
1:A:246:PHE:HE1	1:A:248:ARG:HD2	1.68	0.56
1:A:150:GLN:HB3	11:A:730:GOL:H31	1.87	0.56
1:A:96:ARG:HD3	13:A:825:HOH:O	2.06	0.56
1:A:221:GLU:HB2	1:A:224:LYS:HG2	1.89	0.55
1:A:48:CYS:HB3	1:A:64:VAL:HG21	1.92	0.52
1:A:41:LEU:HD21	1:A:584:LYS:HB2	1.91	0.52
1:B:574:SER:O	13:B:812:HOH:O	2.19	0.52
1:A:379:LYS:NZ	13:A:820:HOH:O	2.43	0.52
1:A:189:LYS:HA	1:A:230:PHE:HZ	1.76	0.50
1:A:507:GLU:HB2	1:A:520:ALA:HB3	1.94	0.50
1:B:241:ARG:NH1	13:B:831:HOH:O	2.45	0.49
1:B:251:LYS:HE2	1:B:252:TYR:CZ	2.47	0.49
1:B:575:GLY:N	13:B:825:HOH:O	2.40	0.48
1:A:75:LYS:HD2	1:A:507:GLU:HG2	1.95	0.47
1:A:517:LEU:HD21	1:A:519:VAL:HG23	1.97	0.47
1:B:300:ASN:ND2	13:B:832:HOH:O	2.46	0.47
1:A:456:ARG:CZ	13:A:828:HOH:O	2.63	0.47
1:B:93:ARG:HG2	1:B:113:ILE:HD13	1.96	0.47
1:B:208:PRO:HD3	1:B:299:LYS:O	2.14	0.47
1:A:360:ILE:HD13	1:A:446:ILE:HD11	1.96	0.47
1:A:517:LEU:C	1:A:517:LEU:HD23	2.35	0.46
1:B:186:GLY:HA2	1:B:189:LYS:HD2	1.97	0.46
1:A:96:ARG:NH1	13:A:825:HOH:O	2.48	0.46
1:A:248:ARG:HH12	1:A:251:LYS:CB	2.26	0.46
1:A:109:ARG:HD2	1:A:110:ASP:OD1	2.15	0.46
1:A:248:ARG:NH1	1:A:251:LYS:HB2	2.28	0.46
1:B:207:ASN:HB2	1:B:211:LEU:O	2.16	0.45
1:B:383:GLN:HG2	1:B:385:THR:O	2.16	0.45
1:B:582:LEU:HB3	1:B:663:CYS:HB2	1.97	0.45
1:B:553:LYS:HE2	1:B:553:LYS:HB3	1.85	0.45
1:A:456:ARG:NH2	13:A:828:HOH:O	2.49	0.45
1:B:140:ARG:NE	13:B:816:HOH:O	2.26	0.45
1:B:87:TYR:CE1	1:B:96:ARG:HD3	2.51	0.45
1:B:190:CYS:HB2	1:B:191:PRO:HD2	1.99	0.44
1:A:207:ASN:HB2	1:A:211:LEU:O	2.19	0.43
1:B:47:CYS:HA	1:B:48:CYS:HA	1.65	0.43
1:A:652:ARG:NH2	13:A:832:HOH:O	2.50	0.43
1:A:638:GLY:C	13:A:814:HOH:O	2.57	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:616:HIS:ND1	1:A:622:ARG:HB2	2.34	0.43
1:B:187:ILE:H	1:B:187:ILE:HD12	1.83	0.43
1:B:537:TYR:O	13:B:813:HOH:O	2.21	0.43
1:B:40:ASN:N	1:B:40:ASN:HD22	2.17	0.43
1:A:290:VAL:O	1:A:333:ASN:HB3	2.19	0.42
1:B:574:SER:N	13:B:825:HOH:O	2.52	0.42
1:A:158:ASN:HB3	1:A:161:ASN:O	2.19	0.42
1:B:606:LEU:HB3	1:B:657:LEU:HD21	2.02	0.42
1:B:290:VAL:O	1:B:333:ASN:HB3	2.20	0.42
1:A:47:CYS:HA	1:A:48:CYS:HA	1.65	0.41
1:A:563:LEU:HD23	1:A:563:LEU:HA	1.87	0.41
1:A:383:GLN:HG2	1:A:385:THR:O	2.21	0.41
1:B:591:GLY:HA2	1:B:645:VAL:HB	2.02	0.41
1:B:505:ILE:HG23	1:B:519:VAL:HG13	2.02	0.41
1:A:236:ASN:HB3	1:A:239:ASN:OD1	2.21	0.41
1:B:508:MET:HA	1:B:518:TYR:O	2.19	0.41
1:A:258:ASP:OD2	7:L:7:MAN:H62	2.21	0.41
1:A:403:PRO:HB3	1:A:421:ILE:HB	2.03	0.41
1:B:187:ILE:O	13:B:814:HOH:O	2.22	0.40
1:A:45:PRO:HA	1:A:46:PRO:HD3	1.99	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)
13:A:1045:HOH:O	13:B:1061:HOH:O[2_675]	2.05	0.15

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	618/657 (94%)	594 (96%)	24 (4%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	616/657 (94%)	594 (96%)	22 (4%)	0	100	100
All	All	1234/1314 (94%)	1188 (96%)	46 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	549/580 (95%)	546 (100%)	3 (0%)	88	93
1	B	548/580 (94%)	543 (99%)	5 (1%)	78	87
All	All	1097/1160 (95%)	1089 (99%)	8 (1%)	84	90

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	246	PHE
1	A	247	LYS
1	A	672	HIS
1	B	40	ASN
1	B	165	TYR
1	B	247	LYS
1	B	433	HIS
1	B	590	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	300	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

51 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	C	1	1,2	14,14,15	0.28	0	17,19,21	0.46	0
2	NAG	C	2	2	14,14,15	0.24	0	17,19,21	0.40	0
2	NAG	D	1	1,2	14,14,15	0.39	0	17,19,21	0.48	0
2	NAG	D	2	2	14,14,15	0.27	0	17,19,21	0.40	0
3	NAG	E	1	1,3	14,14,15	0.19	0	17,19,21	0.38	0
3	NAG	E	2	3	14,14,15	0.22	0	17,19,21	0.38	0
3	BMA	E	3	3	11,11,12	0.70	0	15,15,17	0.74	0
3	MAN	E	4	3	11,11,12	0.59	0	15,15,17	1.13	2 (13%)
3	MAN	E	5	3	11,11,12	0.60	0	15,15,17	1.03	2 (13%)
4	NAG	F	1	1,4	14,14,15	0.28	0	17,19,21	0.46	0
4	NAG	F	2	4	14,14,15	0.14	0	17,19,21	0.41	0
4	BMA	F	3	4	11,11,12	0.53	0	15,15,17	0.70	0
4	MAN	F	4	4	11,11,12	0.65	0	15,15,17	1.17	2 (13%)
4	MAN	F	5	4	11,11,12	0.77	0	15,15,17	0.99	2 (13%)
4	MAN	F	6	4	11,11,12	0.73	0	15,15,17	0.97	2 (13%)
4	MAN	F	7	4	11,11,12	0.81	0	15,15,17	1.02	2 (13%)
4	MAN	F	8	4	11,11,12	0.66	0	15,15,17	0.99	2 (13%)
4	MAN	F	9	4	11,11,12	0.66	0	15,15,17	1.05	2 (13%)
5	NAG	G	1	1,5	14,14,15	0.30	0	17,19,21	0.41	0
5	NAG	G	2	5	14,14,15	0.30	0	17,19,21	0.39	0
5	BMA	G	3	5	11,11,12	0.68	0	15,15,17	0.73	0
5	MAN	G	4	5	11,11,12	0.69	0	15,15,17	1.01	2 (13%)
5	MAN	G	5	5	11,11,12	0.73	0	15,15,17	1.00	2 (13%)
5	MAN	G	6	5	11,11,12	0.58	0	15,15,17	1.08	2 (13%)
6	NAG	H	1	1,6	14,14,15	0.23	0	17,19,21	0.57	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NAG	H	2	6	14,14,15	0.21	0	17,19,21	0.56	0
6	BMA	H	3	6	11,11,12	0.61	0	15,15,17	0.76	0
6	MAN	H	4	6	11,11,12	0.73	0	15,15,17	1.09	2 (13%)
2	NAG	I	1	1,2	14,14,15	0.28	0	17,19,21	0.45	0
2	NAG	I	2	2	14,14,15	0.28	0	17,19,21	0.36	0
2	NAG	J	1	1,2	14,14,15	0.31	0	17,19,21	0.46	0
2	NAG	J	2	2	14,14,15	0.28	0	17,19,21	0.37	0
6	NAG	K	1	1,6	14,14,15	0.22	0	17,19,21	0.33	0
6	NAG	K	2	6	14,14,15	0.25	0	17,19,21	0.35	0
6	BMA	K	3	6	11,11,12	0.68	0	15,15,17	0.94	0
6	MAN	K	4	6	11,11,12	0.70	0	15,15,17	0.99	2 (13%)
7	NAG	L	1	1,7	14,14,15	0.32	0	17,19,21	0.42	0
7	NAG	L	2	7	14,14,15	0.60	1 (7%)	17,19,21	0.47	0
7	MAN	L	3	7	11,11,12	0.47	0	15,15,17	1.24	2 (13%)
7	MAN	L	4	7	11,11,12	0.85	1 (9%)	15,15,17	1.02	2 (13%)
7	MAN	L	5	7	11,11,12	0.73	0	15,15,17	0.96	1 (6%)
7	MAN	L	6	7	11,11,12	0.75	0	15,15,17	0.99	2 (13%)
7	MAN	L	7	7	11,11,12	0.75	1 (9%)	15,15,17	1.04	2 (13%)
7	MAN	L	8	7	11,11,12	0.82	0	15,15,17	1.10	2 (13%)
8	NAG	M	1	1,8	14,14,15	0.38	0	17,19,21	0.38	0
8	NAG	M	2	8	14,14,15	0.29	0	17,19,21	0.34	0
8	MAN	M	3	8	11,11,12	0.76	0	15,15,17	0.99	2 (13%)
8	MAN	M	4	8	11,11,12	0.92	0	15,15,17	1.23	2 (13%)
9	NAG	N	1	1,9	14,14,15	0.32	0	17,19,21	0.46	0
9	NAG	N	2	9	14,14,15	0.23	0	17,19,21	0.38	0
9	BMA	N	3	9	11,11,12	0.71	0	15,15,17	0.76	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	C	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
2	NAG	D	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	D	2	2	-	2/6/23/26	0/1/1/1
3	NAG	E	1	1,3	-	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

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MAN	E	4	3	-	0/2/19/22	0/1/1/1
3	MAN	E	5	3	-	1/2/19/22	0/1/1/1
4	NAG	F	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	F	2	4	-	0/6/23/26	0/1/1/1
4	BMA	F	3	4	-	0/2/19/22	0/1/1/1
4	MAN	F	4	4	-	0/2/19/22	0/1/1/1
4	MAN	F	5	4	-	0/2/19/22	0/1/1/1
4	MAN	F	6	4	-	2/2/19/22	0/1/1/1
4	MAN	F	7	4	-	2/2/19/22	0/1/1/1
4	MAN	F	8	4	-	0/2/19/22	0/1/1/1
4	MAN	F	9	4	-	2/2/19/22	0/1/1/1
5	NAG	G	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	G	2	5	-	0/6/23/26	0/1/1/1
5	BMA	G	3	5	-	0/2/19/22	0/1/1/1
5	MAN	G	4	5	-	0/2/19/22	0/1/1/1
5	MAN	G	5	5	-	0/2/19/22	0/1/1/1
5	MAN	G	6	5	-	2/2/19/22	0/1/1/1
6	NAG	H	1	1,6	-	3/6/23/26	0/1/1/1
6	NAG	H	2	6	-	1/6/23/26	0/1/1/1
6	BMA	H	3	6	-	2/2/19/22	0/1/1/1
6	MAN	H	4	6	-	0/2/19/22	0/1/1/1
2	NAG	I	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	I	2	2	-	0/6/23/26	0/1/1/1
2	NAG	J	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	J	2	2	-	0/6/23/26	0/1/1/1
6	NAG	K	1	1,6	-	0/6/23/26	0/1/1/1
6	NAG	K	2	6	-	2/6/23/26	0/1/1/1
6	BMA	K	3	6	-	2/2/19/22	0/1/1/1
6	MAN	K	4	6	-	0/2/19/22	0/1/1/1
7	NAG	L	1	1,7	-	0/6/23/26	0/1/1/1
7	NAG	L	2	7	-	0/6/23/26	0/1/1/1
7	MAN	L	3	7	-	0/2/19/22	0/1/1/1
7	MAN	L	4	7	-	0/2/19/22	0/1/1/1
7	MAN	L	5	7	-	0/2/19/22	0/1/1/1
7	MAN	L	6	7	-	2/2/19/22	0/1/1/1
7	MAN	L	7	7	-	2/2/19/22	0/1/1/1
7	MAN	L	8	7	-	1/2/19/22	0/1/1/1
8	NAG	M	1	1,8	-	0/6/23/26	0/1/1/1
8	NAG	M	2	8	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	MAN	M	3	8	-	2/2/19/22	0/1/1/1
8	MAN	M	4	8	-	2/2/19/22	0/1/1/1
9	NAG	N	1	1,9	-	0/6/23/26	0/1/1/1
9	NAG	N	2	9	-	0/6/23/26	0/1/1/1
9	BMA	N	3	9	-	0/2/19/22	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	L	4	MAN	O5-C1	-2.32	1.40	1.43
7	L	2	NAG	O5-C1	-2.14	1.40	1.43
7	L	7	MAN	O5-C1	-2.09	1.40	1.43

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	L	3	MAN	C1-O5-C5	3.18	116.50	112.19
3	E	4	MAN	C1-O5-C5	2.95	116.19	112.19
7	L	7	MAN	C1-O5-C5	2.72	115.88	112.19
4	F	9	MAN	C1-O5-C5	2.68	115.82	112.19
5	G	6	MAN	C1-O5-C5	2.64	115.77	112.19
7	L	4	MAN	O2-C2-C3	-2.62	104.89	110.14
4	F	4	MAN	C1-O5-C5	2.61	115.73	112.19
5	G	4	MAN	C1-O5-C5	2.58	115.68	112.19
3	E	5	MAN	C1-O5-C5	2.55	115.65	112.19
5	G	5	MAN	C1-O5-C5	2.51	115.59	112.19
3	E	5	MAN	O2-C2-C3	-2.48	105.17	110.14
4	F	8	MAN	O2-C2-C3	-2.45	105.24	110.14
4	F	9	MAN	O2-C2-C3	-2.44	105.25	110.14
3	E	4	MAN	O2-C2-C3	-2.43	105.28	110.14
4	F	5	MAN	C1-O5-C5	2.41	115.46	112.19
5	G	6	MAN	O2-C2-C3	-2.41	105.31	110.14
7	L	6	MAN	O2-C2-C3	-2.41	105.32	110.14
4	F	6	MAN	C1-O5-C5	2.41	115.45	112.19
4	F	7	MAN	C1-O5-C5	2.41	115.45	112.19
6	K	4	MAN	C1-O5-C5	2.40	115.45	112.19
4	F	4	MAN	O2-C2-C3	-2.35	105.42	110.14
7	L	8	MAN	C1-O5-C5	2.33	115.35	112.19
6	K	4	MAN	O2-C2-C3	-2.33	105.46	110.14
8	M	3	MAN	C1-O5-C5	2.33	115.34	112.19
4	F	8	MAN	C1-O5-C5	2.31	115.32	112.19
7	L	3	MAN	O2-C2-C3	-2.29	105.54	110.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	L	5	MAN	C1-O5-C5	2.29	115.30	112.19
4	F	6	MAN	O2-C2-C3	-2.29	105.55	110.14
7	L	6	MAN	C1-O5-C5	2.28	115.28	112.19
6	H	4	MAN	C1-O5-C5	2.27	115.27	112.19
8	M	4	MAN	O5-C1-C2	2.26	114.26	110.77
5	G	4	MAN	O2-C2-C3	-2.26	105.62	110.14
7	L	8	MAN	O2-C2-C3	-2.23	105.67	110.14
6	H	4	MAN	O2-C2-C3	-2.16	105.81	110.14
5	G	5	MAN	O2-C2-C3	-2.11	105.91	110.14
7	L	7	MAN	O2-C2-C3	-2.11	105.91	110.14
8	M	3	MAN	O2-C2-C3	-2.08	105.97	110.14
4	F	5	MAN	O2-C2-C3	-2.08	105.98	110.14
4	F	7	MAN	O2-C2-C3	-2.07	105.99	110.14
8	M	4	MAN	O2-C2-C3	-2.07	106.00	110.14
7	L	4	MAN	C1-O5-C5	2.01	114.92	112.19

There are no chirality outliers.

All (34) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	M	4	MAN	O5-C5-C6-O6
8	M	2	NAG	O5-C5-C6-O6
2	D	2	NAG	O5-C5-C6-O6
2	C	2	NAG	O5-C5-C6-O6
4	F	9	MAN	O5-C5-C6-O6
8	M	2	NAG	C4-C5-C6-O6
8	M	4	MAN	C4-C5-C6-O6
2	C	2	NAG	C4-C5-C6-O6
6	H	1	NAG	C8-C7-N2-C2
6	H	1	NAG	O7-C7-N2-C2
6	K	2	NAG	O5-C5-C6-O6
2	D	2	NAG	C4-C5-C6-O6
4	F	9	MAN	C4-C5-C6-O6
5	G	6	MAN	C4-C5-C6-O6
6	K	2	NAG	C4-C5-C6-O6
8	M	3	MAN	O5-C5-C6-O6
8	M	3	MAN	C4-C5-C6-O6
5	G	6	MAN	O5-C5-C6-O6
3	E	5	MAN	O5-C5-C6-O6
4	F	7	MAN	C4-C5-C6-O6
7	L	8	MAN	O5-C5-C6-O6
7	L	7	MAN	C4-C5-C6-O6

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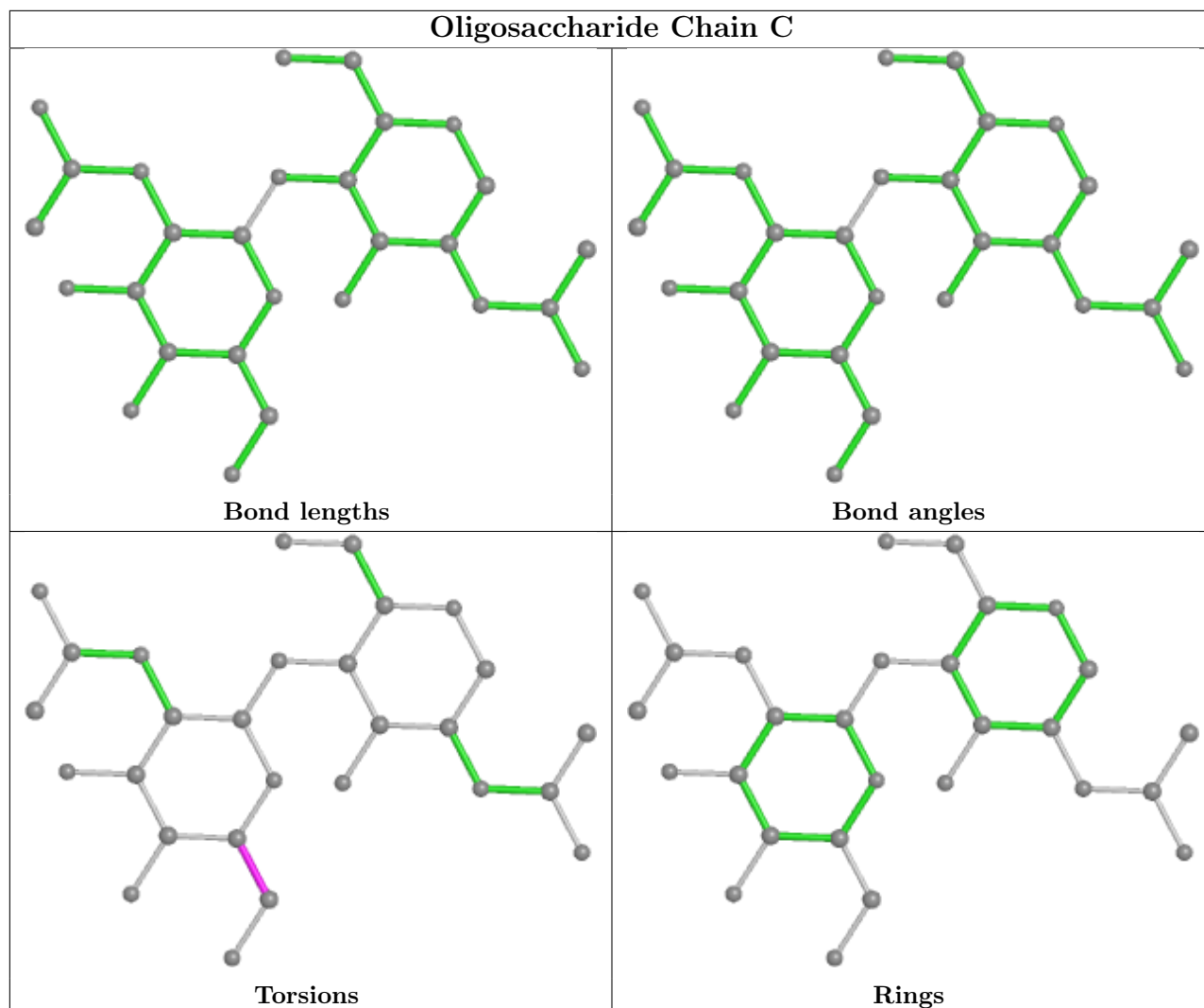
Mol	Chain	Res	Type	Atoms
7	L	6	MAN	C4-C5-C6-O6
4	F	7	MAN	O5-C5-C6-O6
6	K	3	BMA	C4-C5-C6-O6
6	K	3	BMA	O5-C5-C6-O6
6	H	1	NAG	O5-C5-C6-O6
6	H	2	NAG	C3-C2-N2-C7
7	L	6	MAN	O5-C5-C6-O6
4	F	6	MAN	C4-C5-C6-O6
6	H	3	BMA	O5-C5-C6-O6
6	H	3	BMA	C4-C5-C6-O6
7	L	7	MAN	O5-C5-C6-O6
4	F	6	MAN	O5-C5-C6-O6

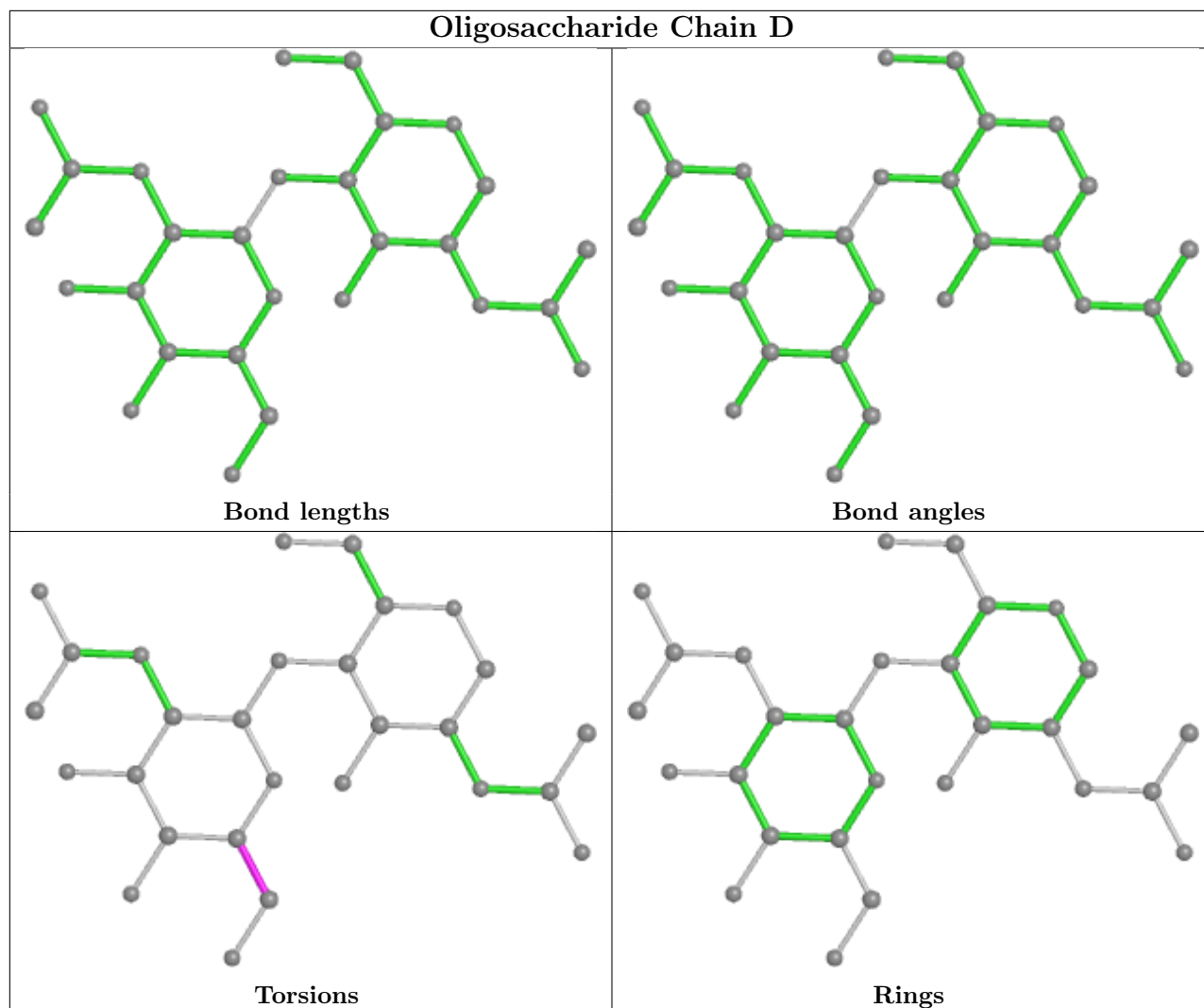
There are no ring outliers.

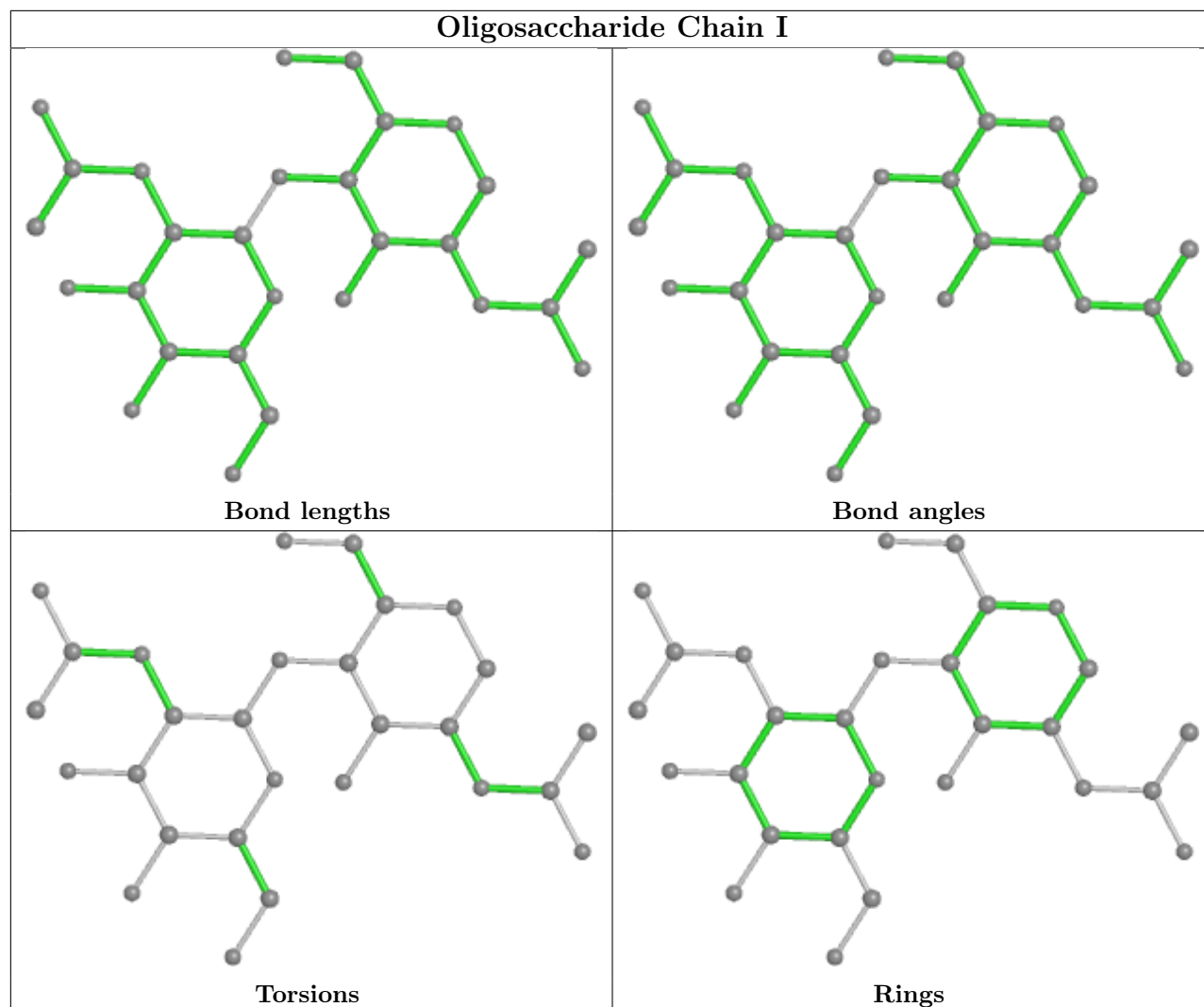
3 monomers are involved in 3 short contacts:

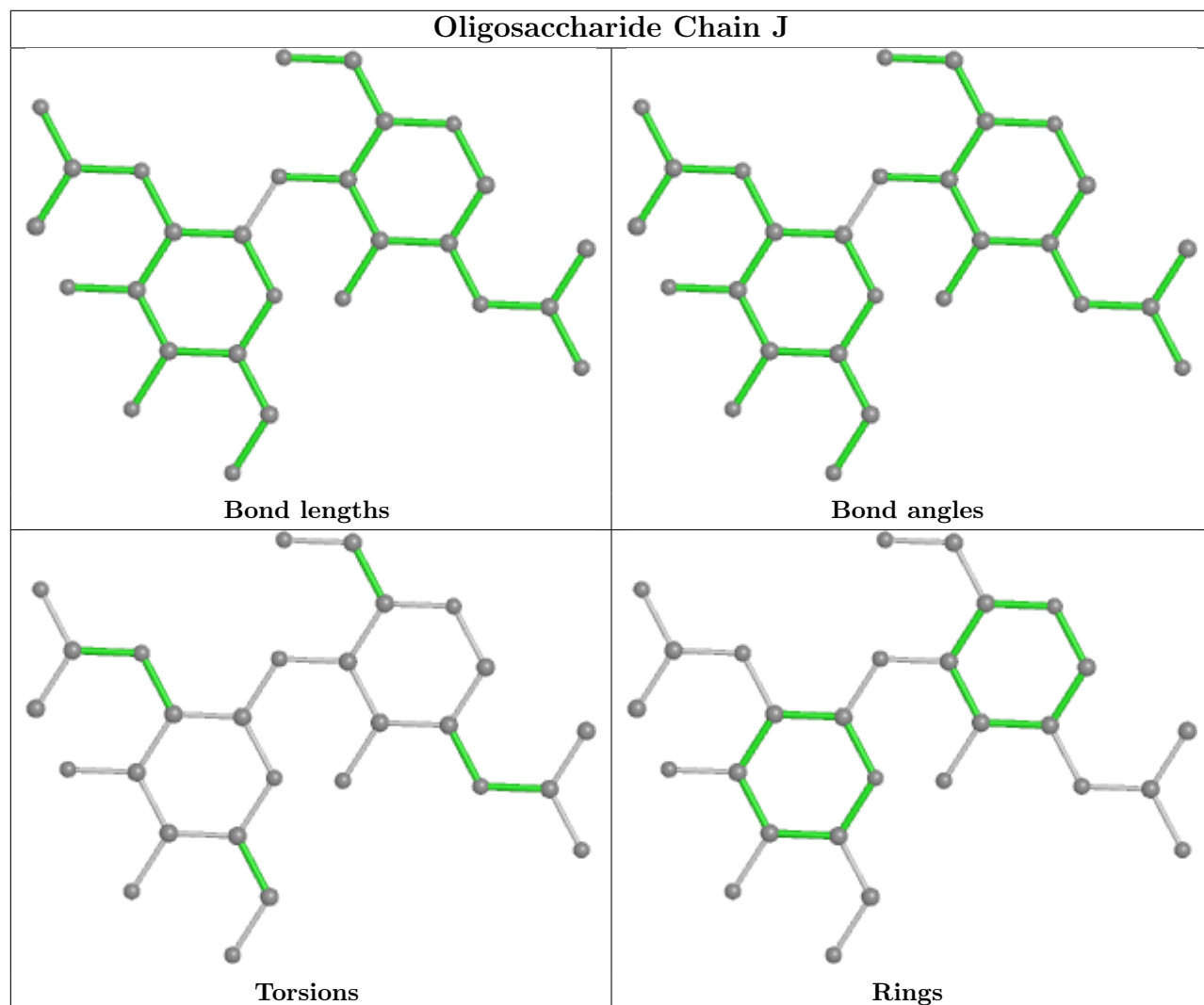
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	L	8	MAN	1	0
8	M	1	NAG	1	0
7	L	7	MAN	1	0

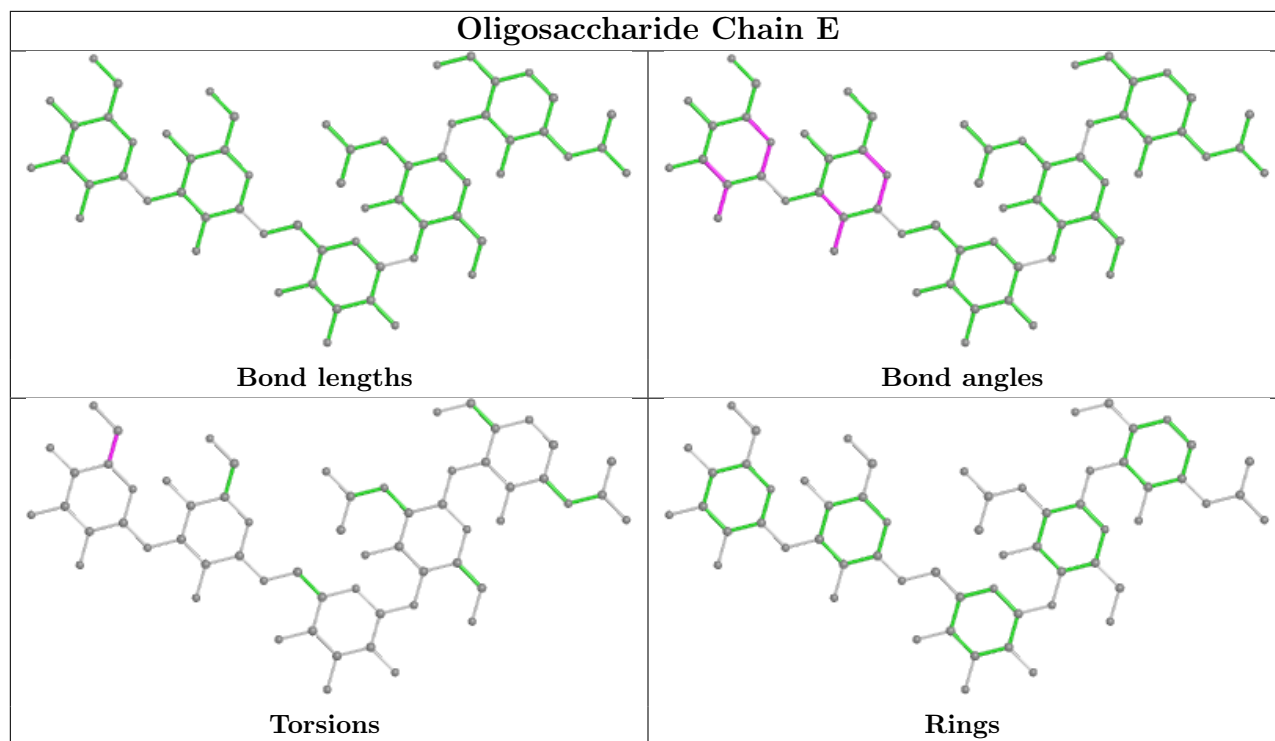
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

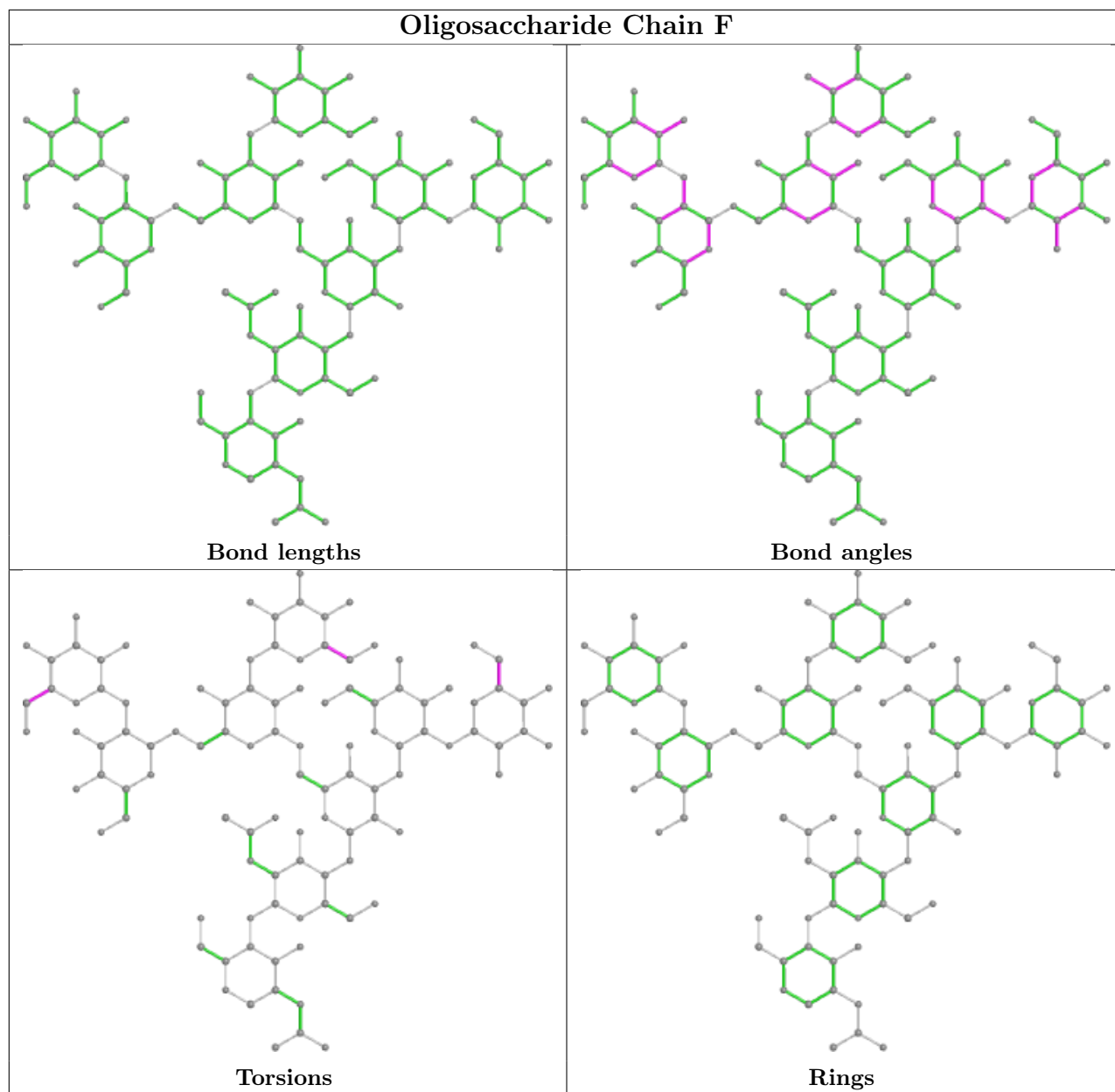


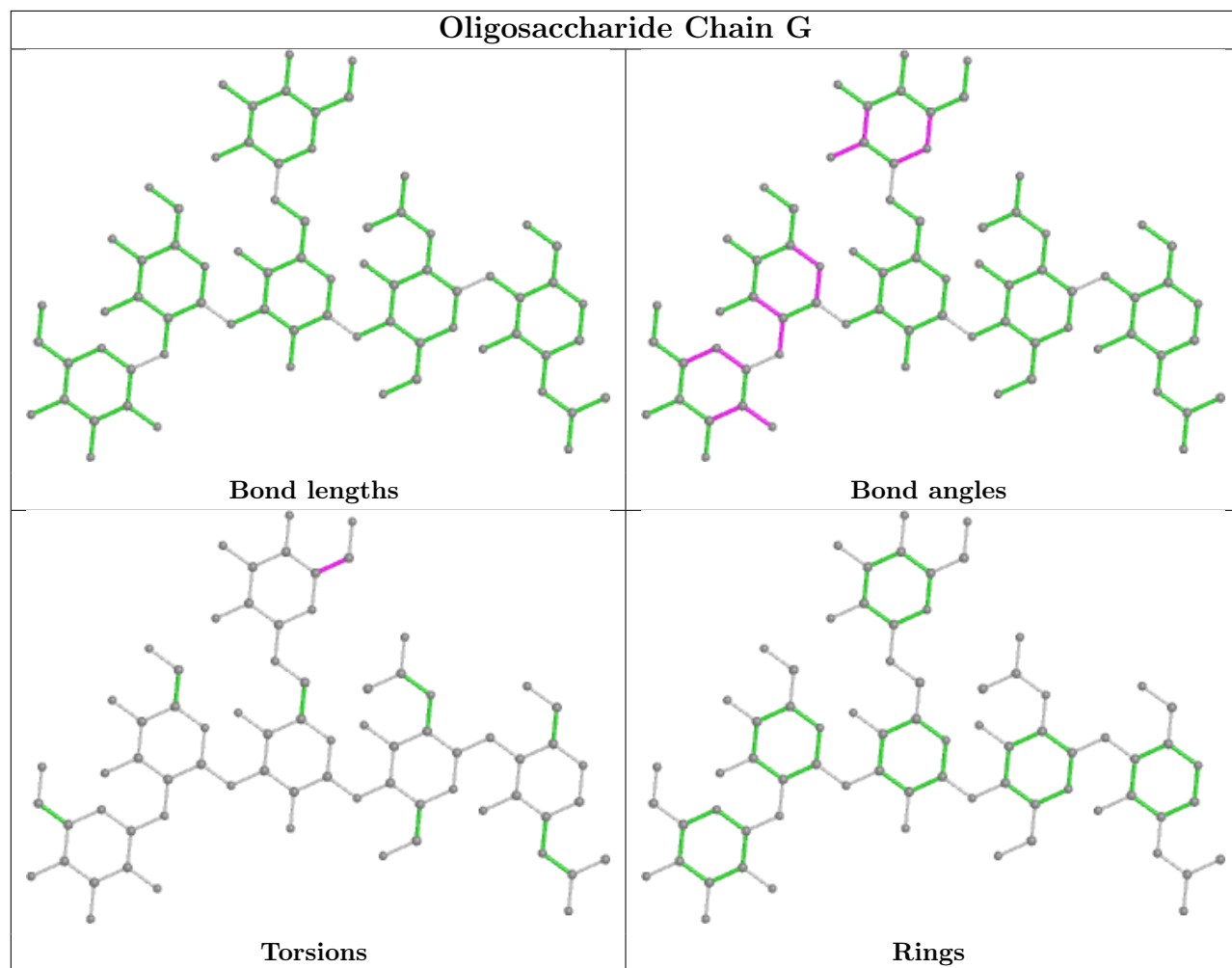


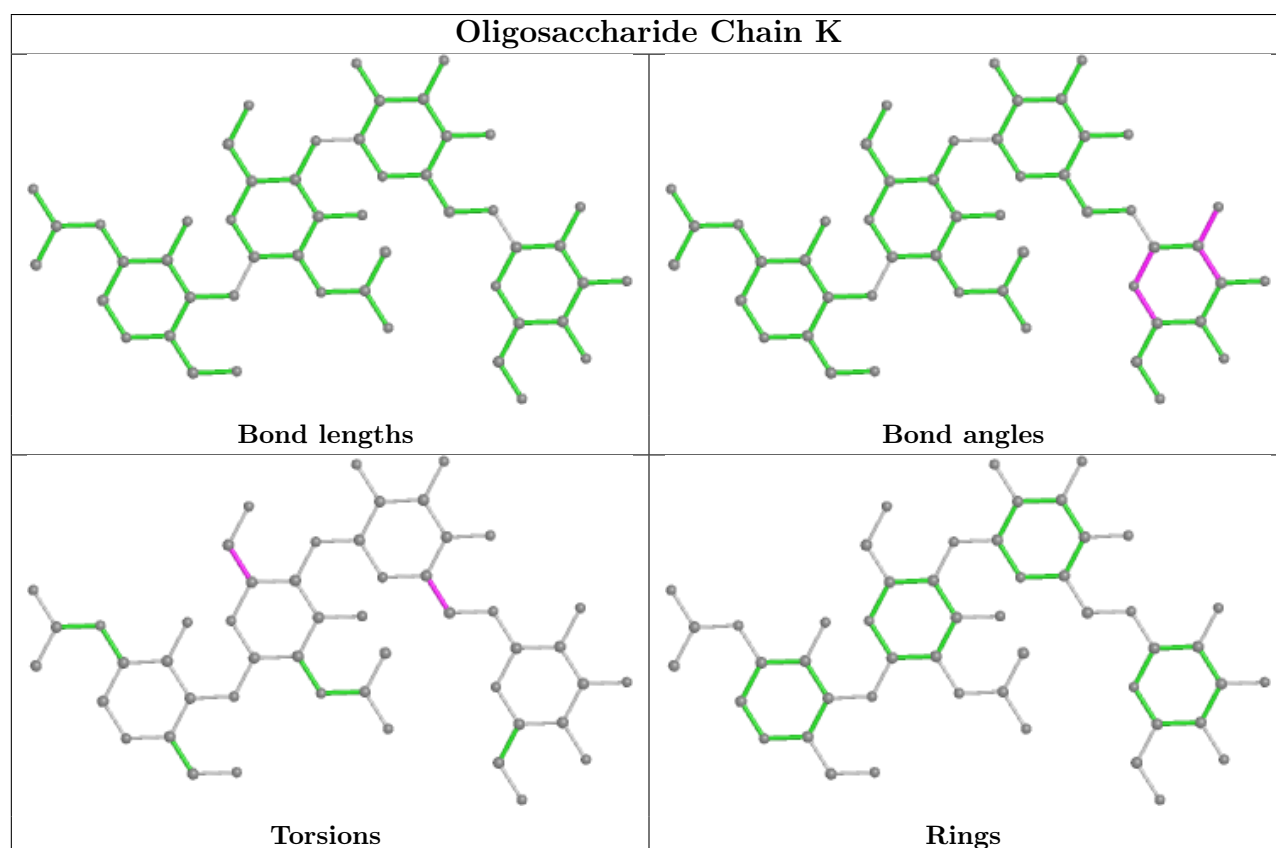
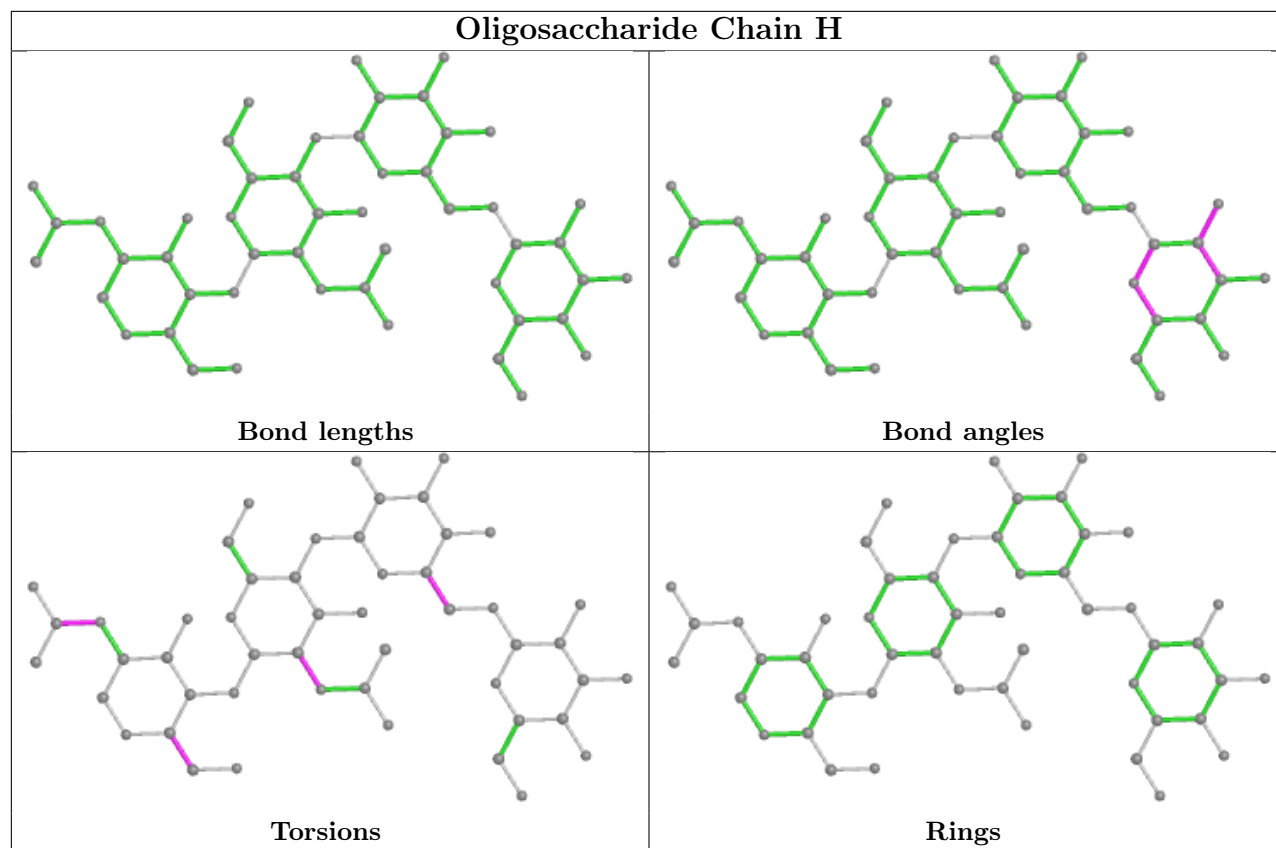


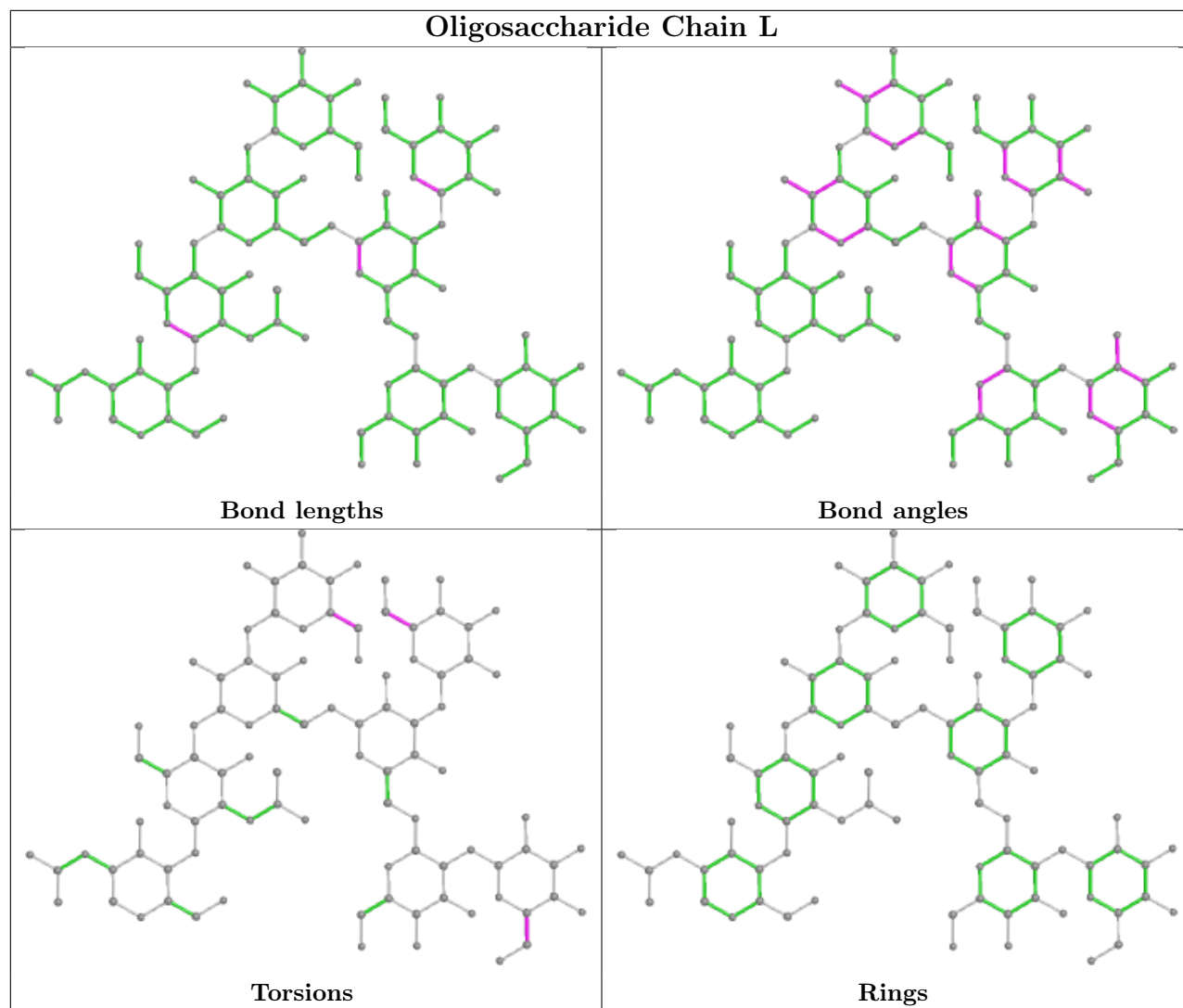


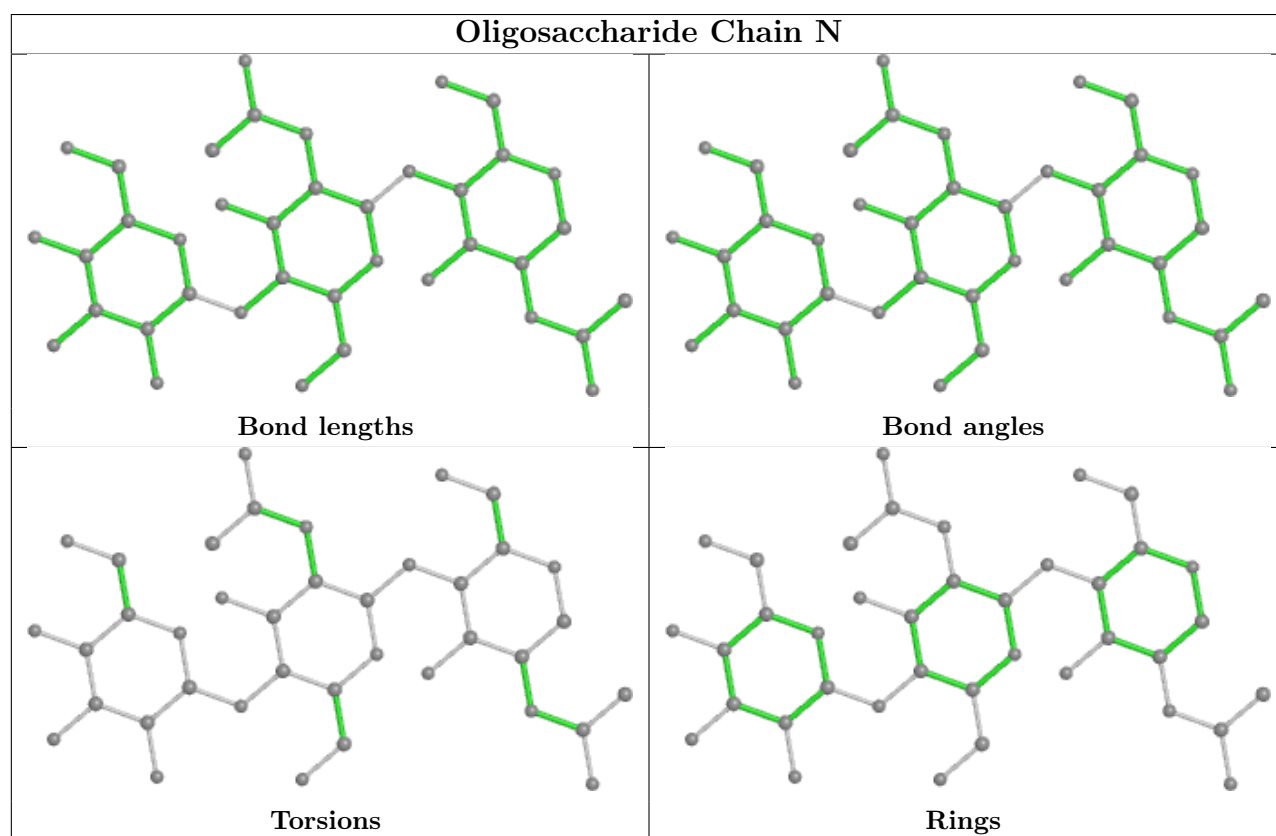
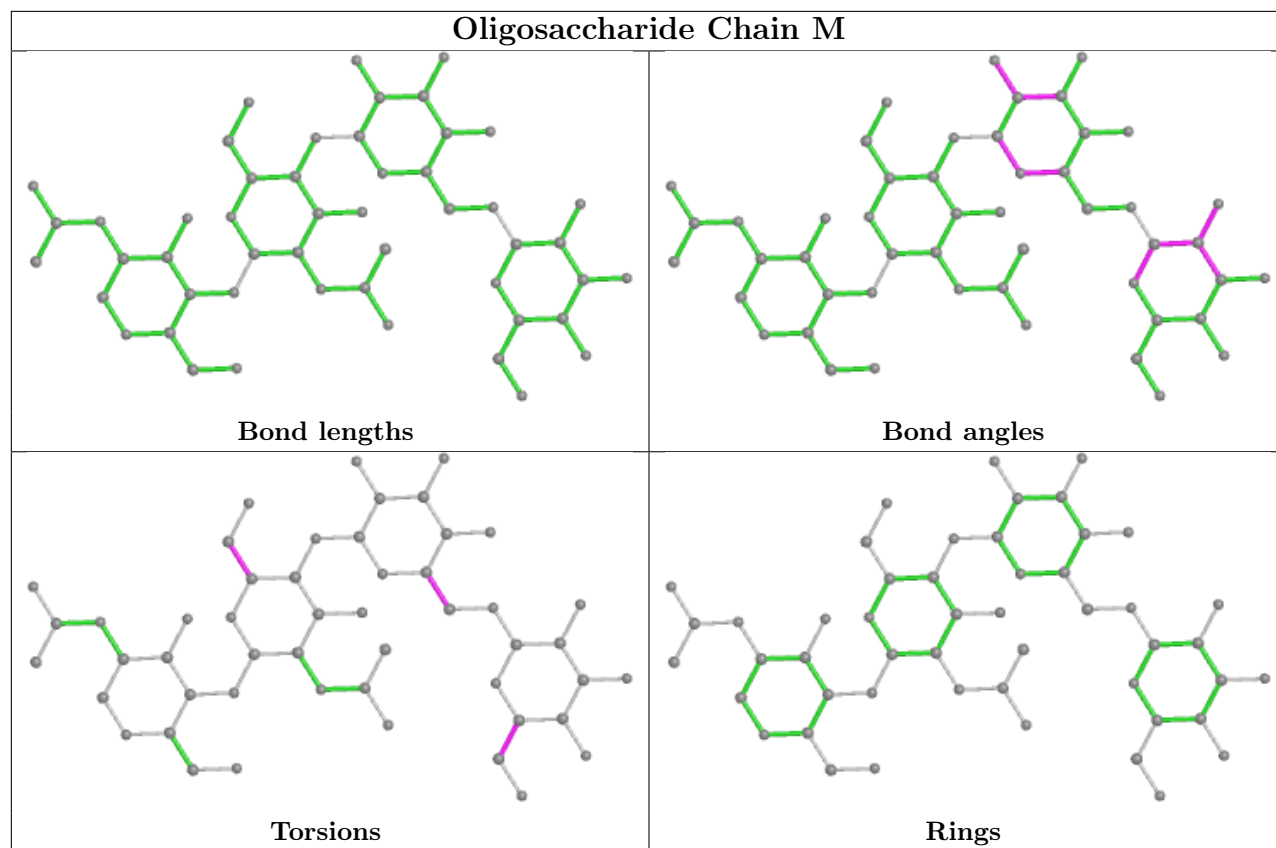












5.6 Ligand geometry

6 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
10	NAG	A	725	1	14,14,15	0.29	0	17,19,21	0.48	0
11	GOL	B	725	-	5,5,5	1.06	0	5,5,5	0.92	0
10	NAG	B	721	1	14,14,15	0.19	0	17,19,21	0.42	0
11	GOL	A	730	-	5,5,5	0.82	0	5,5,5	1.06	0
11	GOL	B	726	-	5,5,5	0.90	0	5,5,5	1.03	0
12	FLC	B	727	-	12,12,12	1.04	0	17,17,17	1.48	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
10	NAG	A	725	1	-	1/6/23/26	0/1/1/1
11	GOL	B	725	-	-	0/4/4/4	-
10	NAG	B	721	1	-	2/6/23/26	0/1/1/1
11	GOL	A	730	-	-	2/4/4/4	-
11	GOL	B	726	-	-	0/4/4/4	-
12	FLC	B	727	-	-	2/16/16/16	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	B	727	FLC	OB2-CBC-CB	3.99	119.98	113.05
12	B	727	FLC	OG2-CGC-CG	2.13	121.17	114.35

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
11	A	730	GOL	C1-C2-C3-O3
10	B	721	NAG	O5-C5-C6-O6
10	B	721	NAG	C4-C5-C6-O6
11	A	730	GOL	O2-C2-C3-O3
12	B	727	FLC	CB-CG-CGC-OG1
10	A	725	NAG	C4-C5-C6-O6
12	B	727	FLC	CB-CG-CGC-OG2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
11	A	730	GOL	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	622/657 (94%)	0.49	52 (8%) 11 16	27, 45, 112, 164	0
1	B	620/657 (94%)	0.67	64 (10%) 6 11	27, 48, 116, 158	0
All	All	1242/1314 (94%)	0.58	116 (9%) 8 14	27, 46, 112, 164	0

All (116) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	590	TYR	18.6
1	B	627	TYR	9.2
1	B	644	ALA	8.9
1	B	587	THR	8.6
1	A	618	THR	8.1
1	B	586	VAL	7.9
1	B	647	GLU	7.9
1	A	633	ILE	7.6
1	B	640	LEU	7.3
1	A	669	VAL	6.9
1	B	633	ILE	6.9
1	B	643	LEU	6.8
1	B	588	SER	6.8
1	A	643	LEU	6.7
1	A	590	TYR	6.7
1	A	671	ALA	6.5
1	A	593	THR	6.2
1	B	646	ASN	6.1
1	A	629	PRO	5.9
1	B	220	ALA	5.6
1	B	632	TYR	5.4
1	A	617	SER	5.3
1	A	132	LYS	5.2
1	B	642	LEU	5.2

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Mol	Chain	Res	Type	RSRZ
1	A	592	GLN	5.2
1	B	589	SER	5.2
1	A	589	SER	5.0
1	B	592	GLN	4.9
1	B	222	PHE	4.9
1	B	630	THR	4.8
1	B	591	GLY	4.7
1	B	223	THR	4.5
1	B	665	TYR	4.4
1	A	41	LEU	4.4
1	B	648	GLY	4.4
1	B	132	LYS	4.4
1	B	620	LYS	4.2
1	B	653	TYR	4.2
1	A	630	THR	4.1
1	B	185	LEU	4.1
1	A	620	LYS	4.1
1	A	672	HIS	4.1
1	A	670	ASP	4.1
1	B	645	VAL	4.0
1	A	625	VAL	4.0
1	A	632	TYR	3.9
1	B	631	LYS	3.9
1	B	616	HIS	3.8
1	B	667	VAL	3.8
1	B	619	GLU	3.7
1	A	619	GLU	3.6
1	B	593	THR	3.6
1	A	652	ARG	3.6
1	B	130	ARG	3.5
1	A	641	VAL	3.4
1	A	649	ASP	3.4
1	B	597	SER	3.4
1	B	668	THR	3.4
1	A	535	ARG	3.4
1	B	179	TYR	3.3
1	A	40	ASN	3.3
1	A	588	SER	3.2
1	A	644	ALA	3.2
1	A	127	LYS	3.2
1	A	178	GLN	3.2
1	B	596	LEU	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	174	LEU	3.1
1	B	650	GLY	3.1
1	B	585	LYS	3.1
1	B	50	GLY	3.1
1	A	587	THR	3.1
1	A	177	SER	3.1
1	B	582	LEU	3.0
1	A	220	ALA	3.0
1	B	629	PRO	2.9
1	A	586	VAL	2.9
1	A	594	LEU	2.9
1	A	179	TYR	2.9
1	B	131	GLU	2.9
1	B	594	LEU	2.9
1	A	176	ARG	2.9
1	B	669	VAL	2.9
1	B	618	THR	2.8
1	B	621	GLY	2.8
1	B	670	ASP	2.8
1	B	601	LYS	2.8
1	B	174	LEU	2.7
1	A	627	TYR	2.7
1	B	649	ASP	2.7
1	B	666	GLY	2.7
1	A	130	ARG	2.6
1	B	177	SER	2.6
1	A	461	ASN	2.6
1	A	642	LEU	2.5
1	A	133	VAL	2.5
1	B	608	LYS	2.4
1	B	623	TYR	2.4
1	A	583	ARG	2.4
1	B	583	ARG	2.4
1	B	612	ARG	2.4
1	A	645	VAL	2.3
1	A	223	THR	2.3
1	A	668	THR	2.3
1	B	628	THR	2.2
1	B	652	ARG	2.2
1	B	187	ILE	2.2
1	A	591	GLY	2.2
1	B	656	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	62	GLU	2.1
1	A	631	LYS	2.1
1	B	599	PHE	2.1
1	A	175	PRO	2.1
1	B	638	GLY	2.1
1	A	585	LYS	2.1
1	A	616	HIS	2.0
1	B	602	MET	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
6	MAN	H	4	11/12	0.54	0.36	140,146,147,148	0
2	NAG	C	2	14/15	0.60	0.55	102,111,125,130	0
8	MAN	M	3	11/12	0.63	0.31	101,107,118,122	0
8	MAN	M	4	11/12	0.66	0.32	95,101,108,111	0
3	MAN	E	5	11/12	0.67	0.26	110,124,125,127	0
4	MAN	F	9	11/12	0.68	0.23	109,114,117,118	0
3	BMA	E	3	11/12	0.69	0.21	119,127,148,157	0
9	BMA	N	3	11/12	0.69	0.18	114,116,118,118	0
7	MAN	L	8	11/12	0.71	0.21	93,105,110,112	0
6	BMA	H	3	11/12	0.71	0.20	125,132,139,142	0
5	MAN	G	6	11/12	0.74	0.31	114,122,131,134	0
6	BMA	K	3	11/12	0.74	0.18	104,109,117,119	0
6	MAN	K	4	11/12	0.76	0.20	111,117,121,123	0
2	NAG	I	2	14/15	0.77	0.32	94,106,112,113	0
5	BMA	G	3	11/12	0.79	0.22	121,136,152,154	0
2	NAG	D	2	14/15	0.81	0.54	115,125,127,128	0
2	NAG	J	2	14/15	0.83	0.35	110,118,121,124	0
3	MAN	E	4	11/12	0.83	0.15	113,119,126,131	0
5	MAN	G	5	11/12	0.84	0.27	81,90,96,97	0
5	MAN	G	4	11/12	0.84	0.30	78,89,103,106	0

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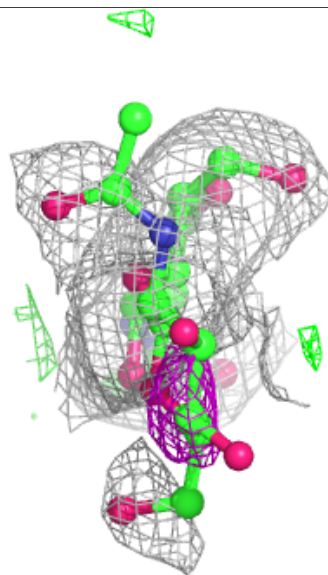
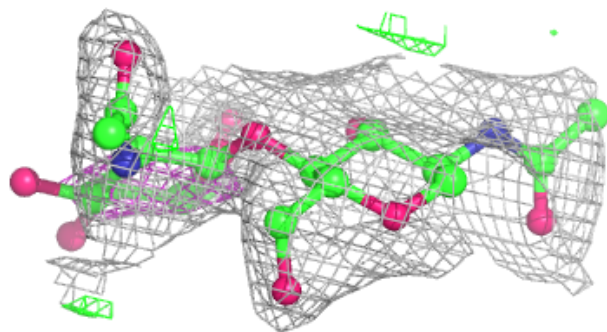
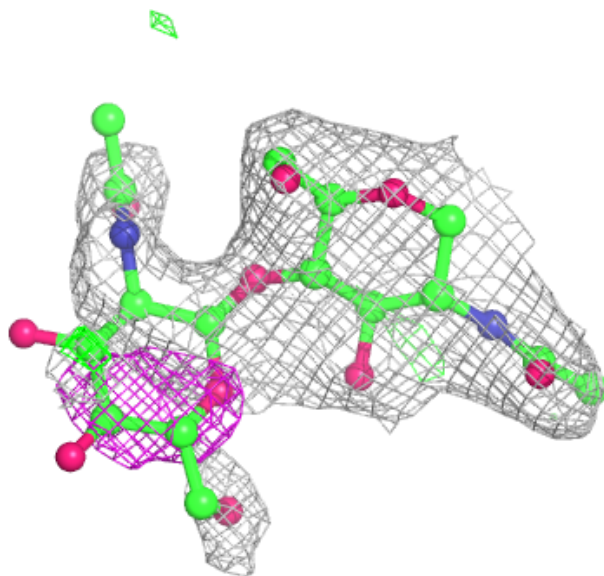
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	MAN	L	7	11/12	0.85	0.16	71,77,86,92	0
6	NAG	K	2	14/15	0.87	0.17	47,83,100,107	0
9	NAG	N	2	14/15	0.88	0.12	94,100,107,113	0
6	NAG	H	2	14/15	0.89	0.16	87,93,100,111	0
2	NAG	I	1	14/15	0.89	0.14	80,87,105,127	0
7	MAN	L	3	11/12	0.89	0.14	65,69,80,92	0
4	MAN	F	7	11/12	0.90	0.14	66,73,74,74	0
4	MAN	F	8	11/12	0.90	0.15	87,89,95,101	0
9	NAG	N	1	14/15	0.91	0.15	69,80,92,96	0
3	NAG	E	2	14/15	0.91	0.13	43,75,93,111	0
8	NAG	M	2	14/15	0.91	0.15	60,69,87,91	0
2	NAG	D	1	14/15	0.92	0.21	53,65,82,99	0
7	MAN	L	4	11/12	0.92	0.16	64,73,80,84	0
6	NAG	H	1	14/15	0.92	0.17	58,70,82,88	0
2	NAG	J	1	14/15	0.92	0.22	64,76,93,104	0
5	NAG	G	2	14/15	0.92	0.17	58,74,90,108	0
2	NAG	C	1	14/15	0.93	0.14	64,80,99,123	0
8	NAG	M	1	14/15	0.93	0.10	45,53,58,62	0
4	BMA	F	3	11/12	0.94	0.12	71,86,93,94	0
4	MAN	F	5	11/12	0.94	0.13	43,52,56,60	0
7	MAN	L	6	11/12	0.95	0.11	41,52,56,56	0
5	NAG	G	1	14/15	0.95	0.14	46,53,57,69	0
4	MAN	F	6	11/12	0.95	0.12	48,54,58,61	0
4	MAN	F	4	11/12	0.95	0.12	58,64,66,68	0
7	MAN	L	5	11/12	0.95	0.10	33,47,51,54	0
3	NAG	E	1	14/15	0.97	0.14	36,44,54,65	0
6	NAG	K	1	14/15	0.97	0.14	36,42,47,59	0
7	NAG	L	2	14/15	0.97	0.18	37,42,52,61	0
4	NAG	F	2	14/15	0.97	0.16	39,44,50,62	0
4	NAG	F	1	14/15	0.98	0.17	24,33,40,46	0
7	NAG	L	1	14/15	0.98	0.18	23,33,36,37	0

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

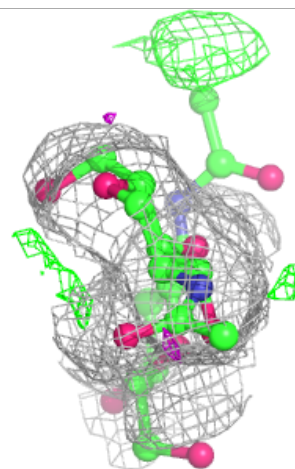
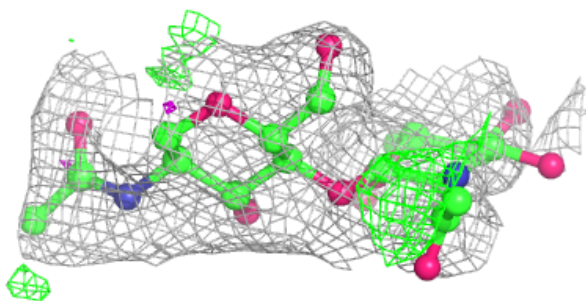
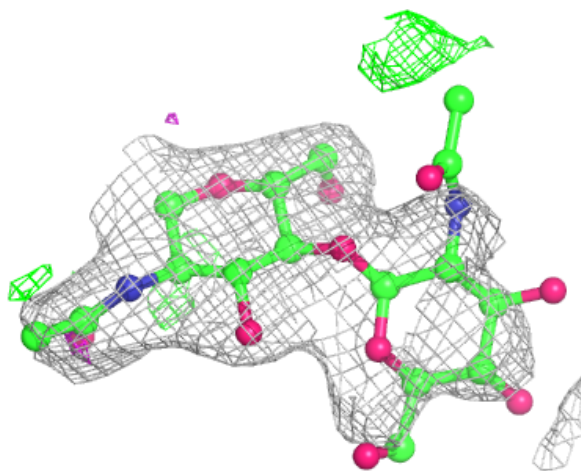
Electron density around Chain C:

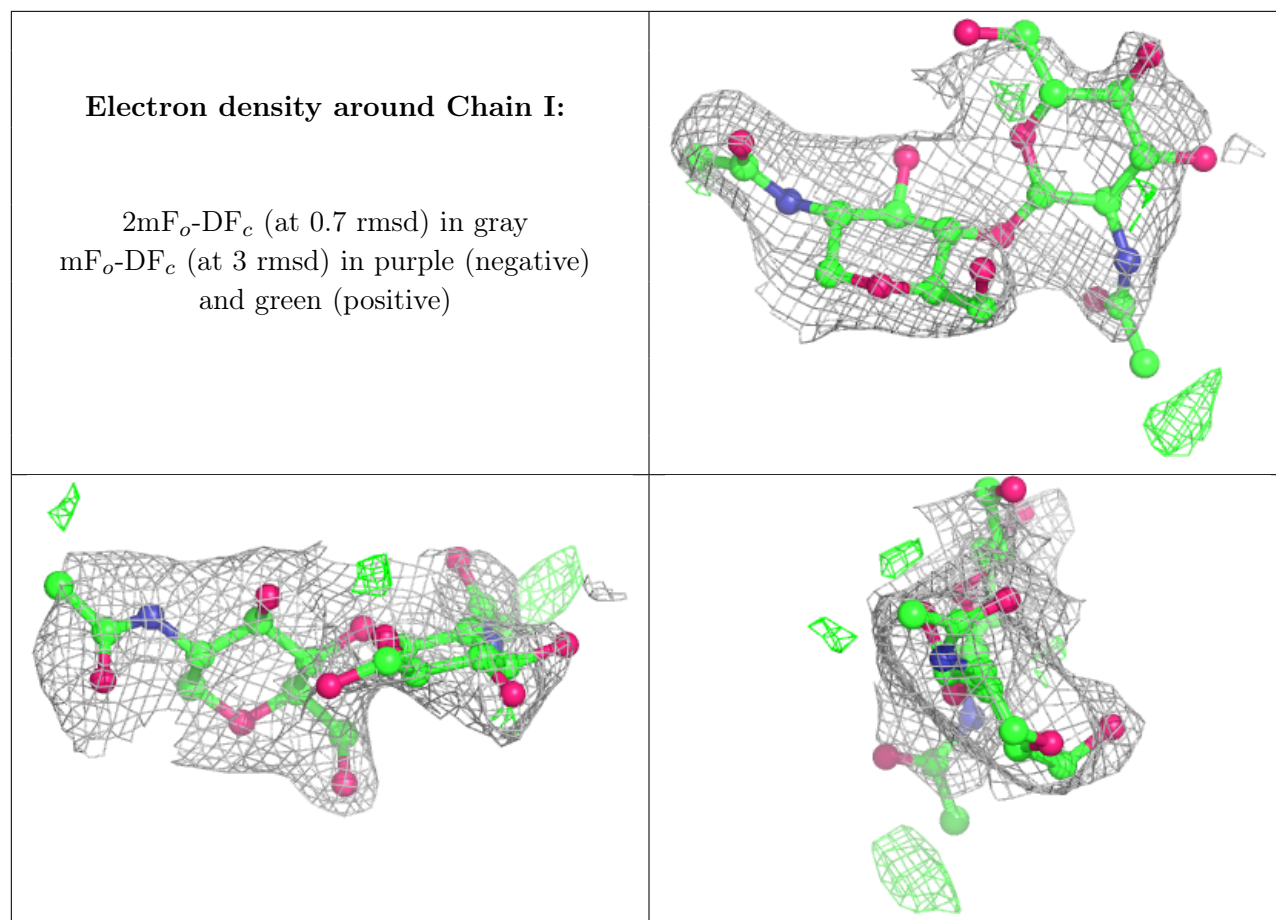
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain D:

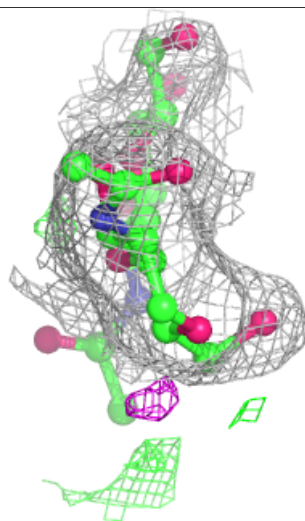
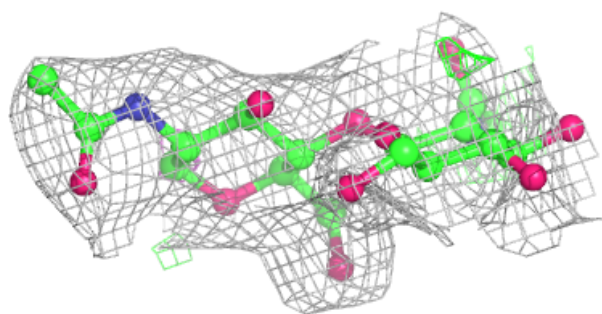
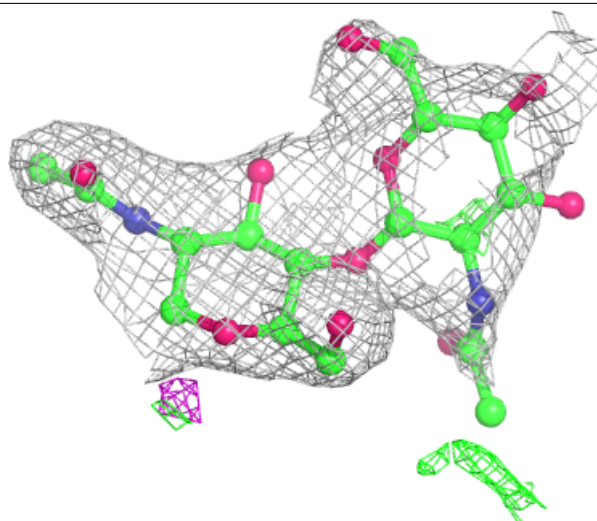
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





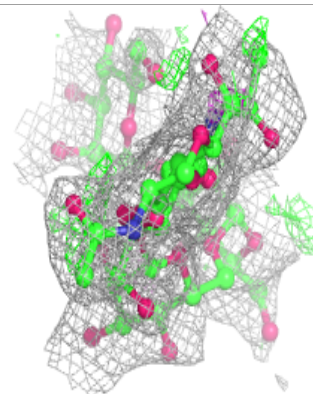
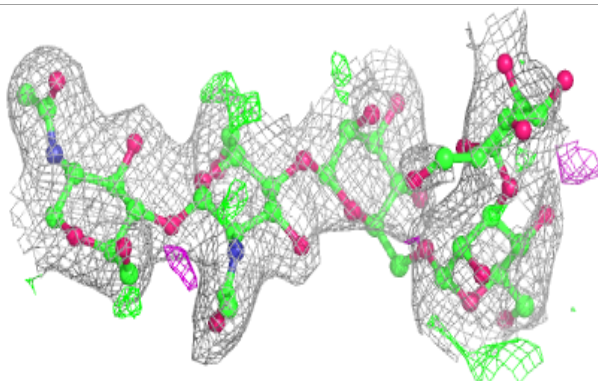
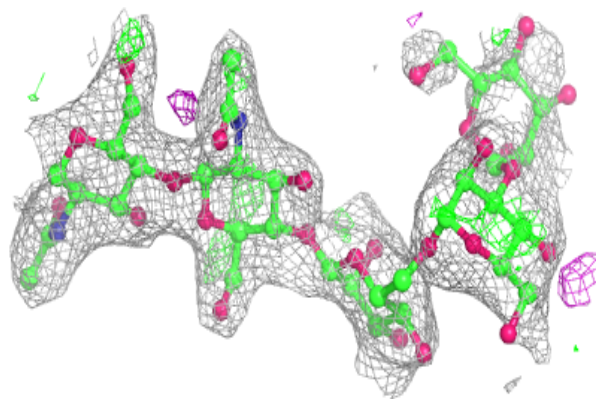
Electron density around Chain J:

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



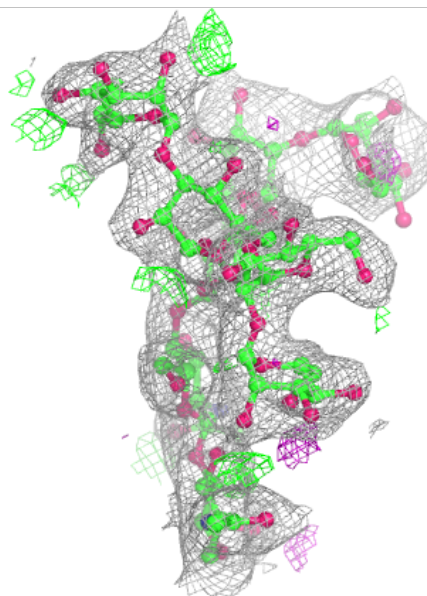
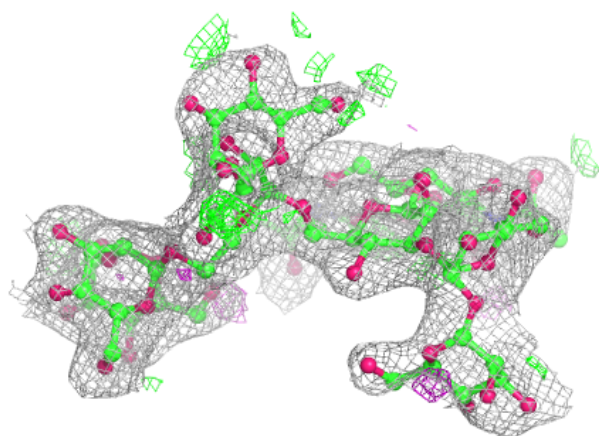
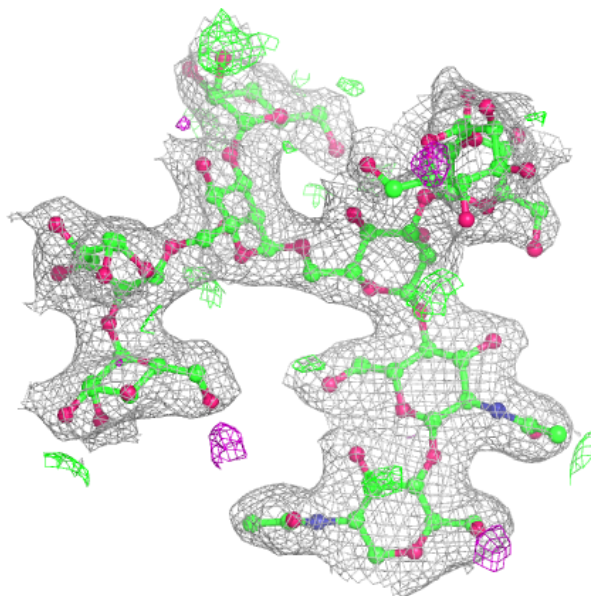
Electron density around Chain E:

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



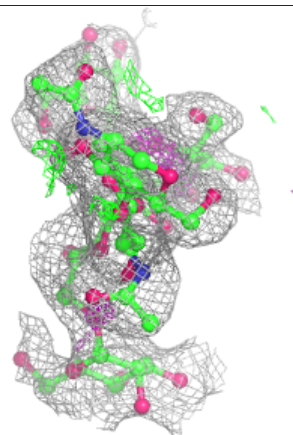
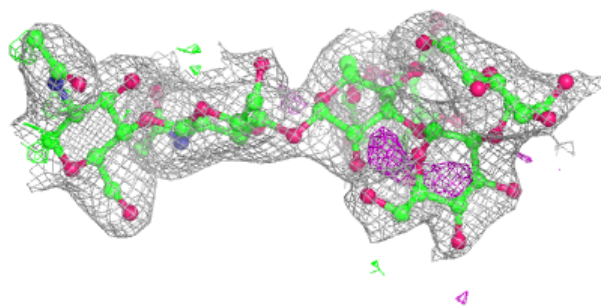
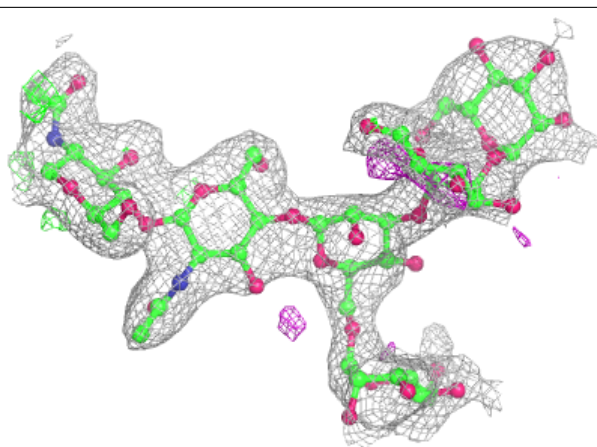
Electron density around Chain F:

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

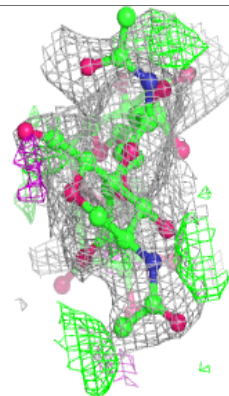
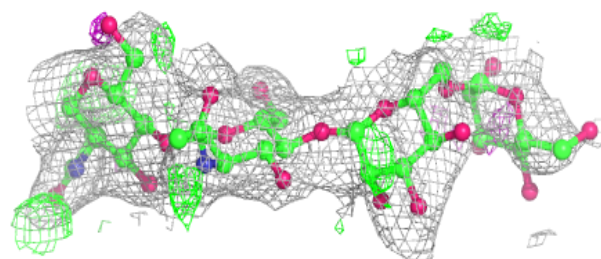
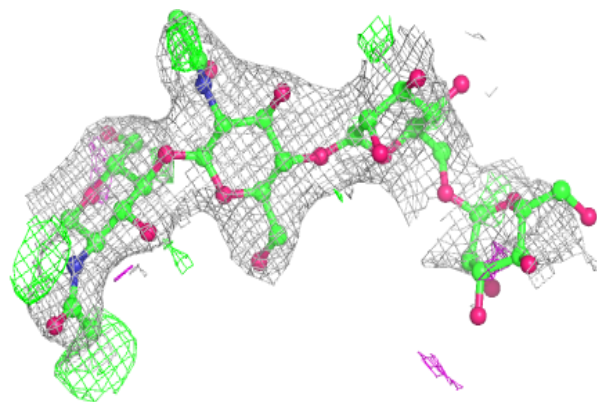


Electron density around Chain G:

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

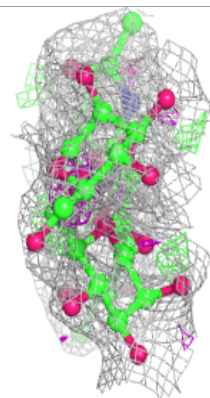
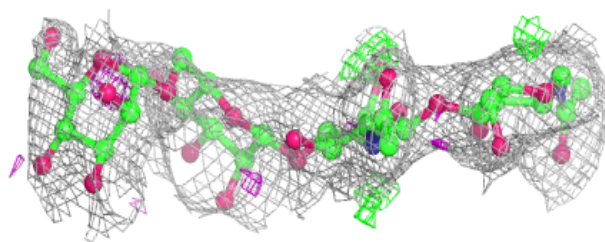
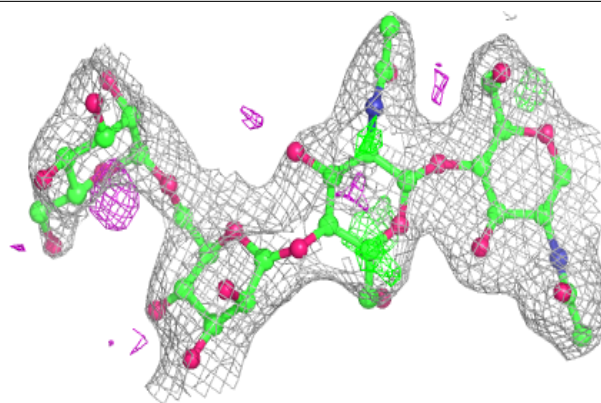
**Electron density around Chain H:**

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



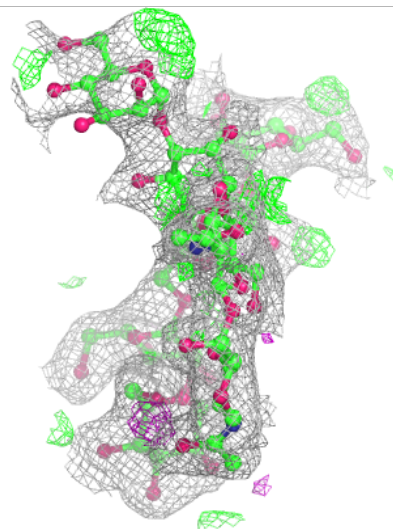
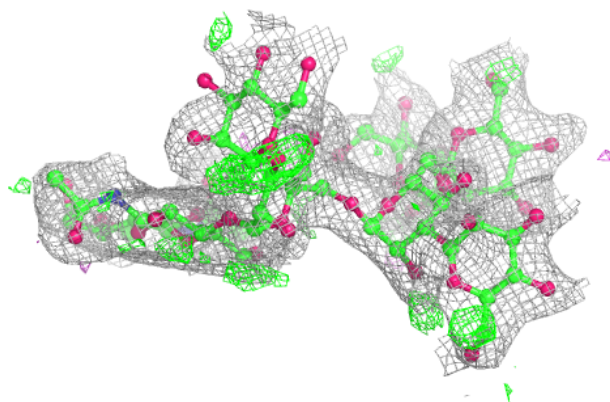
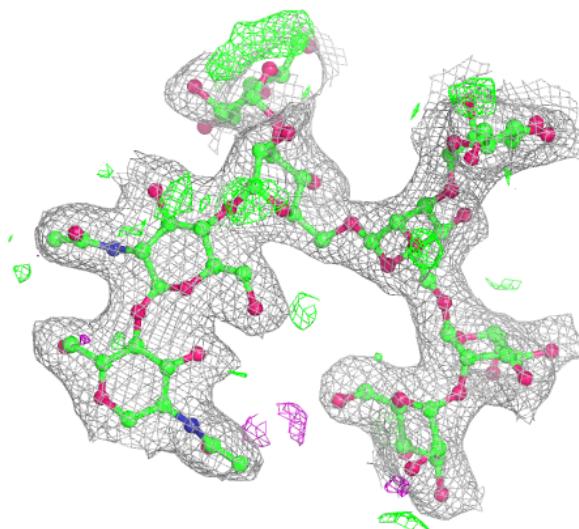
Electron density around Chain K:

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



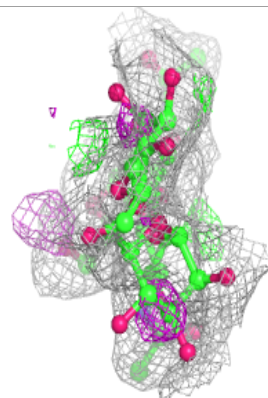
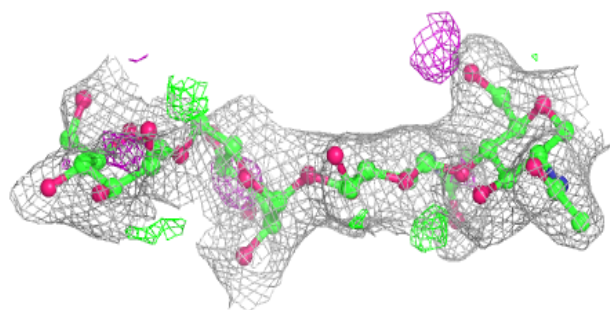
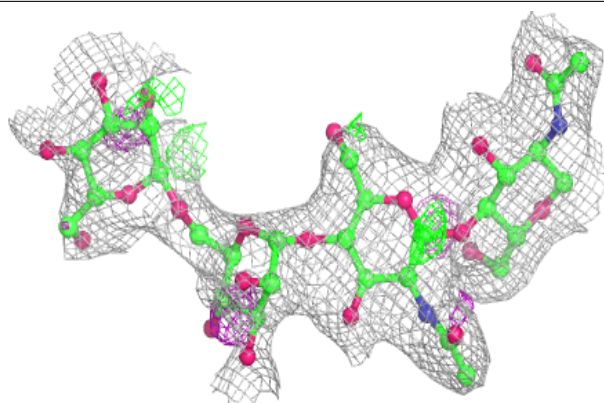
Electron density around Chain L:

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

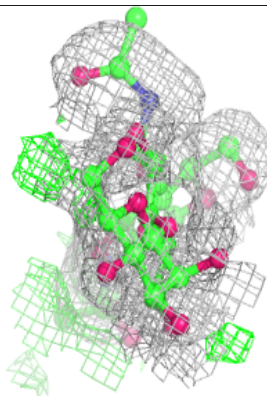
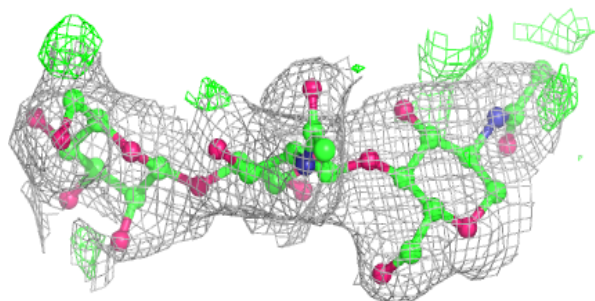
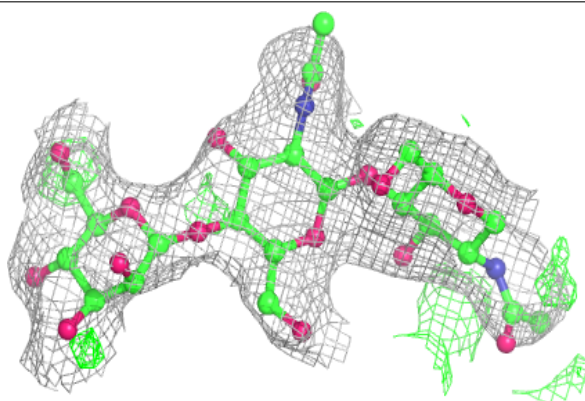


Electron density around Chain M:

$2mF_o-DF_c$ (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)



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
10	NAG	B	721	14/15	0.68	0.22	82,97,110,111	0
11	GOL	A	730	6/6	0.79	0.34	95,96,97,98	0
11	GOL	B	725	6/6	0.81	0.51	54,64,71,75	0
10	NAG	A	725	14/15	0.85	0.15	66,86,95,100	0
11	GOL	B	726	6/6	0.91	0.25	62,68,72,79	0
12	FLC	B	727	13/13	0.93	0.17	63,70,80,82	0

6.5 Other polymers [i](#)

There are no such residues in this entry.