



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 04:23 AM UTC

PDB ID : 6E3L / pdb_00006e3l
Title : Interferon gamma signalling complex with IFNGR1 and IFNGR2
Authors : Jude, K.M.; Mendoza, J.L.; Garcia, K.C.
Deposited on : 2018-07-14
Resolution : 3.80 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

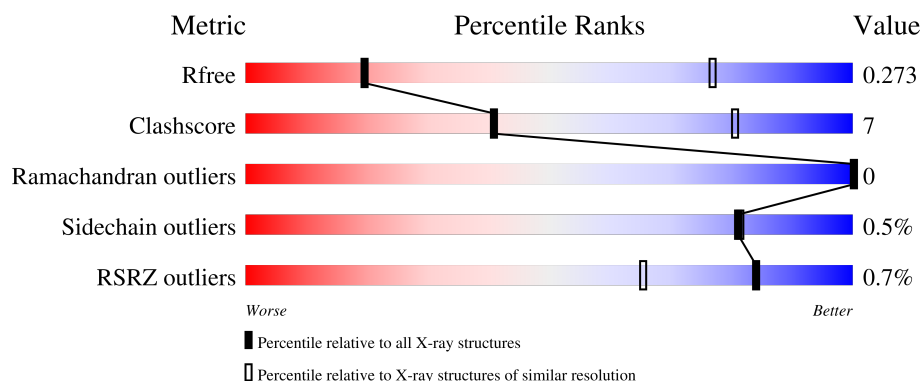
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

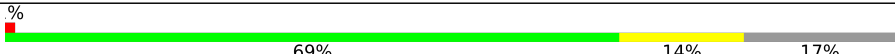
The reported resolution of this entry is 3.80 Å.

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	1065 (3.96-3.64)
Clashscore	190562	1012 (3.94-3.66)
Ramachandran outliers	187476	1048 (3.96-3.64)
Sidechain outliers	187428	1043 (3.96-3.64)
RSRZ outliers	180081	1064 (3.96-3.64)

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	148	
1	B	148	
2	C	242	
2	D	242	
3	E	233	

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Mol	Chain	Length	Quality of chain
3	I	233	 78% 13% 9%
4	F	2	 100%
4	J	2	 100%
4	M	2	 100%
5	G	3	 100%
5	H	3	 67% 33%
5	K	3	 67% 33%
6	L	6	 17% 33% 50%

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 8976 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 Interferon gamma.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	125	Total	C	N	O	S	0	0	0
			1027	653	173	198	3			
1	B	126	Total	C	N	O	S	0	0	0
			1020	647	171	199	3			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP P01579
A	-2	PRO	-	expression tag	UNP P01579
A	-1	GLY	-	expression tag	UNP P01579
A	0	SER	-	expression tag	UNP P01579
A	134	ALA	-	expression tag	UNP P01579
A	135	ALA	-	expression tag	UNP P01579
A	136	ALA	-	expression tag	UNP P01579
A	137	HIS	-	expression tag	UNP P01579
A	138	HIS	-	expression tag	UNP P01579
A	139	HIS	-	expression tag	UNP P01579
A	140	HIS	-	expression tag	UNP P01579
A	141	HIS	-	expression tag	UNP P01579
A	142	HIS	-	expression tag	UNP P01579
A	143	HIS	-	expression tag	UNP P01579
A	144	HIS	-	expression tag	UNP P01579
B	-3	GLY	-	expression tag	UNP P01579
B	-2	PRO	-	expression tag	UNP P01579
B	-1	GLY	-	expression tag	UNP P01579
B	0	SER	-	expression tag	UNP P01579
B	134	ALA	-	expression tag	UNP P01579
B	135	ALA	-	expression tag	UNP P01579
B	136	ALA	-	expression tag	UNP P01579
B	137	HIS	-	expression tag	UNP P01579
B	138	HIS	-	expression tag	UNP P01579
B	139	HIS	-	expression tag	UNP P01579

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Chain	Residue	Modelled	Actual	Comment	Reference
B	140	HIS	-	expression tag	UNP P01579
B	141	HIS	-	expression tag	UNP P01579
B	142	HIS	-	expression tag	UNP P01579
B	143	HIS	-	expression tag	UNP P01579
B	144	HIS	-	expression tag	UNP P01579

- Molecule 2 is a protein called Interferon gamma receptor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	204	Total	C	N	O	S	0	0	0
			1618	1029	271	307	11			
2	D	202	Total	C	N	O	S	0	0	0
			1572	997	260	305	10			

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-1	GLY	-	expression tag	UNP P15260
C	0	SER	-	expression tag	UNP P15260
C	149	ILE	THR	engineered mutation	UNP P15260
C	161	LYS	MET	engineered mutation	UNP P15260
C	167	LYS	GLN	engineered mutation	UNP P15260
C	174	ASN	LYS	engineered mutation	UNP P15260
C	182	ARG	GLN	engineered mutation	UNP P15260
C	205	ASN	HIS	engineered mutation	UNP P15260
C	230	ALA	-	expression tag	UNP P15260
C	231	ALA	-	expression tag	UNP P15260
C	232	ALA	-	expression tag	UNP P15260
C	233	HIS	-	expression tag	UNP P15260
C	234	HIS	-	expression tag	UNP P15260
C	235	HIS	-	expression tag	UNP P15260
C	236	HIS	-	expression tag	UNP P15260
C	237	HIS	-	expression tag	UNP P15260
C	238	HIS	-	expression tag	UNP P15260
C	239	HIS	-	expression tag	UNP P15260
C	240	HIS	-	expression tag	UNP P15260
D	-1	GLY	-	expression tag	UNP P15260
D	0	SER	-	expression tag	UNP P15260
D	149	ILE	THR	engineered mutation	UNP P15260
D	161	LYS	MET	engineered mutation	UNP P15260
D	167	LYS	GLN	engineered mutation	UNP P15260
D	174	ASN	LYS	engineered mutation	UNP P15260

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Chain	Residue	Modelled	Actual	Comment	Reference
D	182	ARG	GLN	engineered mutation	UNP P15260
D	205	ASN	HIS	engineered mutation	UNP P15260
D	230	ALA	-	expression tag	UNP P15260
D	231	ALA	-	expression tag	UNP P15260
D	232	ALA	-	expression tag	UNP P15260
D	233	HIS	-	expression tag	UNP P15260
D	234	HIS	-	expression tag	UNP P15260
D	235	HIS	-	expression tag	UNP P15260
D	236	HIS	-	expression tag	UNP P15260
D	237	HIS	-	expression tag	UNP P15260
D	238	HIS	-	expression tag	UNP P15260
D	239	HIS	-	expression tag	UNP P15260
D	240	HIS	-	expression tag	UNP P15260

- Molecule 3 is a protein called Interferon gamma receptor 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	213	Total	C	N	O	S	0	0	0
			1684	1082	281	312	9			
3	I	213	Total	C	N	O	S	0	0	0
			1698	1090	286	313	9			

There are 26 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	26	GLY	-	expression tag	UNP P38484
E	27	SER	-	expression tag	UNP P38484
E	248	ALA	-	expression tag	UNP P38484
E	249	ALA	-	expression tag	UNP P38484
E	250	ALA	-	expression tag	UNP P38484
E	251	HIS	-	expression tag	UNP P38484
E	252	HIS	-	expression tag	UNP P38484
E	253	HIS	-	expression tag	UNP P38484
E	254	HIS	-	expression tag	UNP P38484
E	255	HIS	-	expression tag	UNP P38484
E	256	HIS	-	expression tag	UNP P38484
E	257	HIS	-	expression tag	UNP P38484
E	258	HIS	-	expression tag	UNP P38484
I	26	GLY	-	expression tag	UNP P38484
I	27	SER	-	expression tag	UNP P38484
I	248	ALA	-	expression tag	UNP P38484
I	249	ALA	-	expression tag	UNP P38484

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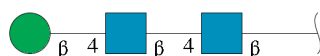
Chain	Residue	Modelled	Actual	Comment	Reference
I	250	ALA	-	expression tag	UNP P38484
I	251	HIS	-	expression tag	UNP P38484
I	252	HIS	-	expression tag	UNP P38484
I	253	HIS	-	expression tag	UNP P38484
I	254	HIS	-	expression tag	UNP P38484
I	255	HIS	-	expression tag	UNP P38484
I	256	HIS	-	expression tag	UNP P38484
I	257	HIS	-	expression tag	UNP P38484
I	258	HIS	-	expression tag	UNP P38484

- 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	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	J	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	M	2	Total	C	N	O	0	0	0
			28	16	2	10			

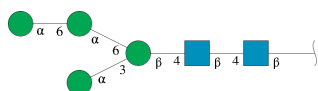
- Molecule 5 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	G	3	Total	C	N	O	0	0	0
			39	22	2	15			
5	H	3	Total	C	N	O	0	0	0
			39	22	2	15			
5	K	3	Total	C	N	O	0	0	0
			39	22	2	15			

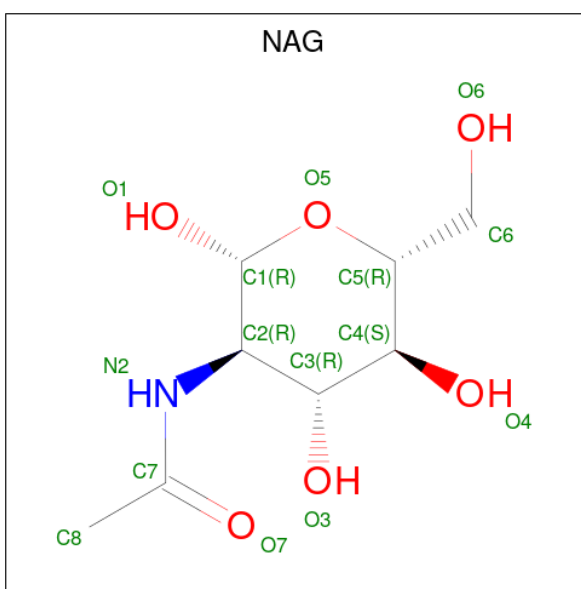
- Molecule 6 is an oligosaccharide called alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-

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



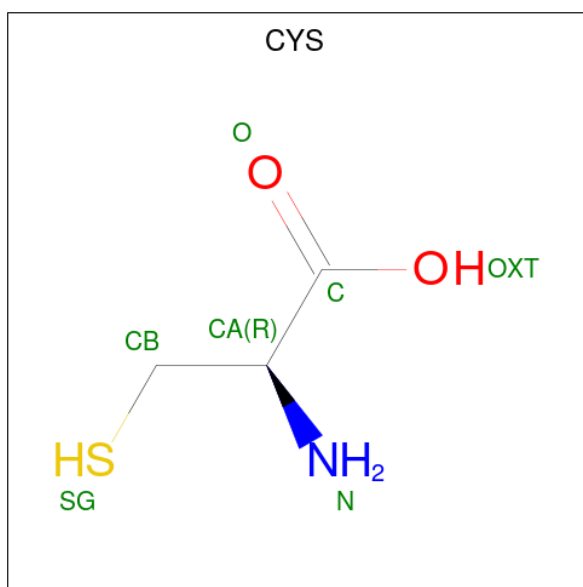
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	L	6	Total	C	N	O	0	0	0
			72	40	2	30			

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



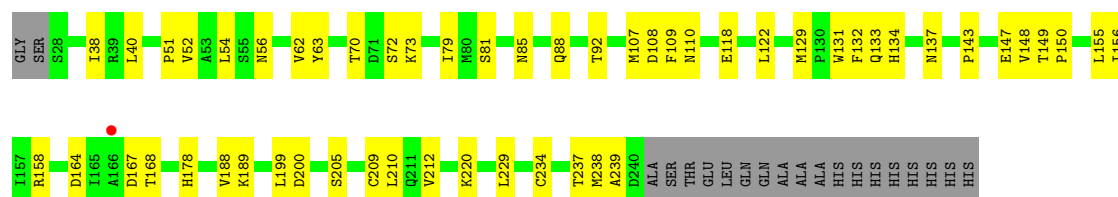
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			14	8	1	5		
7	B	1	Total	C	N	O	0	0
			14	8	1	5		
7	D	1	Total	C	N	O	0	0
			14	8	1	5		
7	E	1	Total	C	N	O	0	0
			14	8	1	5		
7	E	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 8 is CYSTEINE (CCD ID: CYS) (formula: $C_3H_7NO_2S$).



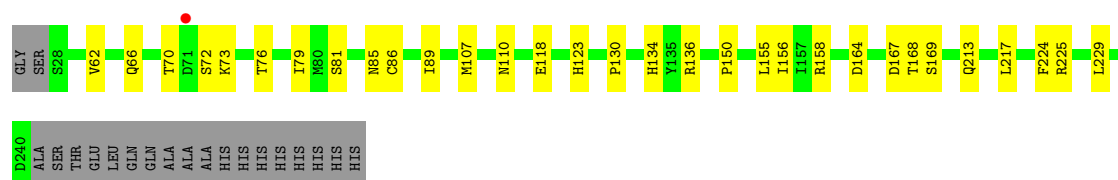
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	E	1	Total	C	N	O	S	0	0
			7	3	1	2	1		
8	I	1	Total	C	N	O	S	0	0
			7	3	1	2	1		

Chain E:  68% 23% 9%



- Molecule 3: Interferon gamma receptor 2

Chain I:  78% 13% 9%



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

Chain F:  100%

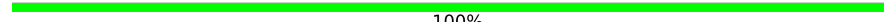


- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  100%

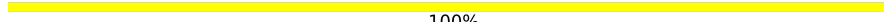


- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M:  100%



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

Chain G:  100%



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

Chain H:  67% 33%

MAG1
MAG2
BMA3

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

Chain K:  67% 33%

MAG1
MAG2
BMA3

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

Chain L:  17% 33% 50%

MAG1
MAG2
BMA3
MAN4
MAN5
MAN6

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.69Å 150.21Å 212.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.74 – 3.80 48.74 – 3.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.74-3.80) 88.5 (48.74-3.80)	Depositor EDS
R_{merge}	0.27	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.71 (at 3.77Å)	Xtriage
Refinement program	PHENIX 1.14_3211	Depositor
R, R_{free}	0.248 , 0.272 0.252 , 0.273	Depositor DCC
R_{free} test set	1693 reflections (6.61%)	wwPDB-VP
Wilson B-factor (Å ²)	133.4	Xtriage
Anisotropy	0.540	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 168.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8976	wwPDB-VP
Average B, all atoms (Å ²)	216.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.48% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.14	0/1045	0.32	0/1399
1	B	0.18	0/1038	0.34	0/1394
2	C	0.13	0/1654	0.35	0/2250
2	D	0.18	0/1607	0.41	0/2194
3	E	0.24	0/1735	0.57	0/2372
3	I	0.17	0/1749	0.42	0/2388
All	All	0.18	0/8828	0.42	0/11997

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	1027	0	1021	11	0
1	B	1020	0	993	17	0
2	C	1618	0	1573	18	0
2	D	1572	0	1490	21	0
3	E	1684	0	1605	45	0
3	I	1698	0	1634	25	0
4	F	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	J	28	0	25	1	0
4	M	28	0	25	0	0
5	G	39	0	34	1	0
5	H	39	0	34	0	0
5	K	39	0	34	2	0
6	L	72	0	61	2	0
7	A	14	0	13	0	0
7	B	14	0	13	0	0
7	D	14	0	13	0	0
7	E	28	0	26	2	0
8	E	7	0	3	0	0
8	I	7	0	3	0	0
All	All	8976	0	8625	115	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (115) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:27:PRO:HG2	2:D:70:ILE:HG13	1.74	0.69
2:D:28:ILE:HD11	5:G:1:NAG:H82	1.77	0.66
2:C:170:ILE:HD11	2:C:185:LEU:HD13	1.81	0.64
2:D:171:LEU:HD21	3:E:168:THR:HB	1.80	0.63
3:I:79:ILE:HG13	3:I:85:ASN:HB2	1.81	0.63
3:E:131:TRP:HZ3	3:E:133:GLN:HG3	1.63	0.62
3:E:149:THR:HB	3:E:156:ILE:HD11	1.82	0.62
1:B:84:SER:HB3	3:E:133:GLN:HE22	1.66	0.60
3:E:54:LEU:HD23	7:E:307:NAG:HN2	1.67	0.60
1:B:79:VAL:HG11	3:E:108:ASP:HB3	1.83	0.58
3:E:38:ILE:HG13	3:E:129:MET:HE2	1.83	0.58
2:D:110:ILE:HD12	2:D:204:LEU:HD11	1.84	0.58
3:E:62:VAL:HG21	3:E:88:GLN:OE1	2.03	0.58
3:E:189:LYS:NZ	3:E:200:ASP:OD1	2.36	0.58
3:E:62:VAL:HG23	3:E:118:GLU:HG3	1.86	0.58
2:C:110:ILE:HB	2:C:211:THR:HG22	1.86	0.57
1:A:22:VAL:HG12	2:C:49:TYR:HE2	1.68	0.57
1:B:78:ASN:HA	1:B:82:PHE:HD2	1.70	0.57
3:E:205:SER:OG	3:E:239:ALA:HA	2.04	0.57
2:D:22:SER:HB2	2:D:27:PRO:HB3	1.88	0.56
2:C:27:PRO:HG2	2:C:70:ILE:HG13	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:51:VAL:O	2:C:54:SER:OG	2.25	0.55
2:C:188:PRO:HB2	3:I:156:ILE:HD12	1.89	0.54
3:E:54:LEU:CD2	7:E:307:NAG:HN2	2.20	0.54
1:A:119:GLU:HB3	1:B:42:ARG:HH12	1.72	0.54
2:C:204:LEU:HD22	2:C:209:VAL:HB	1.89	0.53
2:D:110:ILE:HB	2:D:211:THR:HG22	1.90	0.53
2:C:169:LYS:NZ	3:I:164:ASP:OD1	2.28	0.53
1:B:9:GLU:HA	1:B:12:LYS:HB2	1.90	0.53
2:D:169:LYS:NZ	3:E:164:ASP:OD1	2.39	0.53
3:I:150:PRO:HA	3:I:155:LEU:HD23	1.90	0.53
1:B:72:THR:HG23	3:E:109:PHE:HE2	1.75	0.52
1:A:42:ARG:HH12	1:B:119:GLU:HB3	1.73	0.52
1:A:113:LEU:HD21	1:B:73:ILE:HD13	1.93	0.51
2:C:55:GLU:N	2:C:55:GLU:OE1	2.44	0.51
2:C:171:LEU:HD21	3:I:168:THR:HB	1.92	0.51
1:B:68:LYS:NZ	3:E:81:SER:O	2.43	0.50
1:A:18:GLY:O	2:C:82:TRP:NE1	2.38	0.50
3:I:70:THR:OG1	3:I:110:ASN:HB3	2.12	0.50
2:D:170:ILE:HD11	2:D:185:LEU:HD13	1.94	0.50
3:I:79:ILE:CG1	3:I:85:ASN:HB2	2.42	0.50
3:I:86:CYS:HA	3:I:89:ILE:HD13	1.93	0.50
2:D:153:ARG:NH2	3:E:167:ASP:HA	2.27	0.49
3:E:209:CYS:HA	3:E:234:CYS:HA	1.94	0.49
3:E:40:LEU:HD12	3:E:137:ASN:HB2	1.94	0.49
1:B:27:THR:CG2	2:D:50:GLY:HA2	2.43	0.49
2:C:153:ARG:NH2	3:I:167:ASP:HA	2.28	0.49
3:I:70:THR:HG1	3:I:110:ASN:HB3	1.78	0.49
3:E:143:PRO:HG2	3:E:212:VAL:HG12	1.94	0.48
3:E:148:VAL:HG21	3:E:210:LEU:HD21	1.95	0.47
1:B:83:ASN:ND2	3:E:220:LYS:HG2	2.29	0.47
3:E:150:PRO:HA	3:E:155:LEU:HD23	1.95	0.47
1:B:16:ASN:HB3	1:B:19:HIS:HD2	1.80	0.47
3:I:118:GLU:HA	3:I:123:HIS:HA	1.97	0.47
3:E:149:THR:O	3:E:156:ILE:HG12	2.14	0.47
3:E:178:HIS:HD2	3:E:229:LEU:HD21	1.80	0.47
3:E:107:MET:HG2	3:E:134:HIS:CD2	2.50	0.47
3:E:79:ILE:HG13	3:E:85:ASN:HB2	1.97	0.46
2:D:159:VAL:HG22	2:D:198:VAL:HG22	1.97	0.46
1:B:83:ASN:HA	3:E:220:LYS:HD3	1.98	0.46
3:I:213:GLN:HB2	3:I:229:LEU:HG	1.97	0.46
3:E:56:ASN:OD1	3:E:56:ASN:N	2.36	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:217:LEU:HD23	3:I:224:PHE:HB3	1.96	0.46
3:E:178:HIS:CD2	3:E:229:LEU:HD21	2.51	0.46
3:I:62:VAL:HG11	5:K:2:NAG:C8	2.46	0.46
3:E:72:SER:OG	3:E:73:LYS:N	2.49	0.46
3:E:178:HIS:ND1	3:E:188:VAL:HG22	2.31	0.45
3:I:66:GLN:HG2	3:I:76:THR:HA	1.97	0.45
1:A:68:LYS:HE3	3:I:81:SER:O	2.17	0.45
3:E:79:ILE:HD12	3:E:79:ILE:H	1.81	0.45
3:I:62:VAL:HG11	5:K:2:NAG:H82	1.98	0.45
2:D:189:VAL:HG12	2:D:196:TYR:CZ	2.52	0.45
1:A:27:THR:CG2	2:C:50:GLY:HA2	2.47	0.45
2:D:76:ASP:OD1	2:D:148:THR:OG1	2.34	0.45
3:I:118:GLU:HB3	3:I:123:HIS:HB3	1.98	0.44
1:B:72:THR:HG23	3:E:109:PHE:CE2	2.52	0.44
3:E:147:GLU:HB2	3:E:158:ARG:HG3	1.99	0.44
2:D:40:PRO:HB3	2:D:89:VAL:HG12	2.00	0.44
2:D:114:LYS:HB2	2:D:129:PHE:HB2	2.00	0.43
3:E:150:PRO:O	3:E:238:MET:HE3	2.18	0.43
3:E:79:ILE:HD11	3:E:85:ASN:HB2	2.00	0.43
3:I:167:ASP:OD1	3:I:169:SER:OG	2.28	0.43
1:A:114:ILE:H	1:A:114:ILE:HD12	1.84	0.43
3:E:51:PRO:HB3	3:E:92:THR:HG21	1.99	0.43
3:I:72:SER:OG	3:I:73:LYS:N	2.51	0.43
1:A:86:LYS:NZ	1:A:90:ASP:OD2	2.50	0.43
1:A:112:GLU:OE1	1:B:30:LEU:HG	2.19	0.43
2:C:166:ILE:HG12	3:I:158:ARG:NH2	2.34	0.43
2:D:160:ARG:HA	2:D:165:GLU:HA	2.00	0.42
2:C:181:ILE:HG13	2:C:182:ARG:HG2	2.01	0.42
2:D:12:VAL:HG21	2:D:93:GLU:C	2.44	0.42
3:E:205:SER:N	3:E:237:THR:OG1	2.52	0.42
3:I:107:MET:HG2	3:I:134:HIS:CD2	2.54	0.42
1:A:117:MET:HE3	1:A:117:MET:HB3	1.87	0.42
1:B:117:MET:HE3	1:B:117:MET:HB3	1.91	0.42
4:J:1:NAG:O3	4:J:2:NAG:O5	2.36	0.42
2:D:47:LYS:HB2	2:D:56:TRP:CE3	2.54	0.42
1:B:114:ILE:H	1:B:114:ILE:HD12	1.84	0.42
3:E:88:GLN:NE2	3:E:118:GLU:OE1	2.50	0.42
3:I:217:LEU:HD23	3:I:217:LEU:HA	1.85	0.41
2:C:22:SER:HA	2:C:27:PRO:HA	2.01	0.41
2:C:189:VAL:HG12	2:C:196:TYR:CZ	2.55	0.41
3:E:40:LEU:HD11	3:E:132:PHE:CD1	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:103:ALA:HB3	2:C:106:ARG:HB3	2.02	0.41
3:E:70:THR:OG1	3:E:110:ASN:HB3	2.20	0.41
6:L:3:BMA:H61	6:L:4:MAN:H2	1.75	0.41
3:E:62:VAL:HG22	3:E:118:GLU:O	2.21	0.41
3:E:199:LEU:HD23	3:E:199:LEU:HA	1.98	0.41
3:I:136:ARG:O	3:I:225:ARG:HD3	2.21	0.41
3:E:79:ILE:CD1	3:E:85:ASN:HB2	2.51	0.41
2:D:219:ILE:HD12	2:D:219:ILE:HA	1.99	0.40
3:E:52:VAL:HG23	3:E:63:TYR:HE2	1.86	0.40
2:D:134:PHE:CE2	2:D:152:ILE:HD11	2.56	0.40
2:D:179:ASP:OD1	2:D:181:ILE:HG22	2.21	0.40
3:I:130:PRO:HB2	6:L:1:NAG:H61	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	123/148 (83%)	121 (98%)	2 (2%)	0	100	100
1	B	124/148 (84%)	123 (99%)	1 (1%)	0	100	100
2	C	200/242 (83%)	195 (98%)	5 (2%)	0	100	100
2	D	198/242 (82%)	193 (98%)	5 (2%)	0	100	100
3	E	211/233 (91%)	203 (96%)	8 (4%)	0	100	100
3	I	211/233 (91%)	204 (97%)	7 (3%)	0	100	100
All	All	1067/1246 (86%)	1039 (97%)	28 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	115/132 (87%)	115 (100%)	0	100	100
1	B	112/132 (85%)	112 (100%)	0	100	100
2	C	87/218 (40%)	87 (100%)	0	100	100
2	D	176/218 (81%)	173 (98%)	3 (2%)	53	67
3	E	185/203 (91%)	184 (100%)	1 (0%)	81	81
3	I	188/203 (93%)	188 (100%)	0	100	100
All	All	863/1106 (78%)	859 (100%)	4 (0%)	81	81

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

Mol	Chain	Res	Type
2	D	135	VAL
2	D	136	ASN
2	D	152	ILE
3	E	122	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	ASN
1	A	19	HIS
1	A	83	ASN
1	A	104	ASN
1	B	1	GLN
1	B	19	HIS
1	B	46	GLN
2	C	195	GLN
2	D	123	GLN
2	D	195	GLN
3	E	186	GLN
3	I	29	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

21 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	NAG	F	1	2,4	14,14,15	0.46	0	17,19,21	0.70	0
4	NAG	F	2	4	14,14,15	0.47	0	17,19,21	0.60	0
5	NAG	G	1	2,5	14,14,15	0.59	0	17,19,21	0.66	0
5	NAG	G	2	5	14,14,15	0.66	1 (7%)	17,19,21	0.95	1 (5%)
5	BMA	G	3	5	11,11,12	1.08	1 (9%)	15,15,17	0.91	1 (6%)
5	NAG	H	1	3,5	14,14,15	0.64	1 (7%)	17,19,21	0.69	0
5	NAG	H	2	5	14,14,15	0.27	0	17,19,21	0.47	0
5	BMA	H	3	5	11,11,12	0.53	0	15,15,17	0.62	0
4	NAG	J	1	4,3	14,14,15	0.45	0	17,19,21	0.83	1 (5%)
4	NAG	J	2	4	14,14,15	0.44	0	17,19,21	0.73	1 (5%)
5	NAG	K	1	3,5	14,14,15	0.35	0	17,19,21	0.61	0
5	NAG	K	2	5	14,14,15	0.29	0	17,19,21	0.53	0
5	BMA	K	3	5	11,11,12	0.56	0	15,15,17	0.62	0
6	NAG	L	1	6,3	14,14,15	1.08	1 (7%)	17,19,21	1.68	3 (17%)
6	NAG	L	2	6	14,14,15	0.29	0	17,19,21	0.60	0
6	BMA	L	3	6	11,11,12	0.68	0	15,15,17	1.06	1 (6%)
6	MAN	L	4	6	11,11,12	0.64	0	15,15,17	0.96	2 (13%)
6	MAN	L	5	6	11,11,12	1.20	1 (9%)	15,15,17	0.92	0
6	MAN	L	6	6	11,11,12	0.75	1 (9%)	15,15,17	1.27	2 (13%)
4	NAG	M	1	4,3	14,14,15	0.37	0	17,19,21	0.64	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	M	2	4	14,14,15	0.25	0	17,19,21	0.51	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
4	NAG	F	1	2,4	-	0/6/23/26	0/1/1/1
4	NAG	F	2	4	-	0/6/23/26	0/1/1/1
5	NAG	G	1	2,5	-	0/6/23/26	0/1/1/1
5	NAG	G	2	5	-	1/6/23/26	0/1/1/1
5	BMA	G	3	5	-	2/2/19/22	0/1/1/1
5	NAG	H	1	3,5	-	2/6/23/26	0/1/1/1
5	NAG	H	2	5	-	2/6/23/26	0/1/1/1
5	BMA	H	3	5	-	0/2/19/22	0/1/1/1
4	NAG	J	1	4,3	-	2/6/23/26	0/1/1/1
4	NAG	J	2	4	-	0/6/23/26	0/1/1/1
5	NAG	K	1	3,5	-	2/6/23/26	0/1/1/1
5	NAG	K	2	5	-	2/6/23/26	0/1/1/1
5	BMA	K	3	5	-	0/2/19/22	0/1/1/1
6	NAG	L	1	6,3	-	1/6/23/26	0/1/1/1
6	NAG	L	2	6	-	0/6/23/26	0/1/1/1
6	BMA	L	3	6	-	0/2/19/22	0/1/1/1
6	MAN	L	4	6	-	2/2/19/22	0/1/1/1
6	MAN	L	5	6	-	0/2/19/22	0/1/1/1
6	MAN	L	6	6	-	2/2/19/22	0/1/1/1
4	NAG	M	1	4,3	-	0/6/23/26	0/1/1/1
4	NAG	M	2	4	-	2/6/23/26	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	L	1	NAG	O5-C1	-3.45	1.37	1.43
5	G	3	BMA	C1-C2	2.85	1.59	1.52
6	L	5	MAN	O5-C5	2.63	1.48	1.43
5	G	2	NAG	C1-C2	2.17	1.55	1.52
6	L	6	MAN	C1-C2	2.16	1.57	1.52
5	H	1	NAG	C1-C2	2.06	1.55	1.52

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	L	1	NAG	C1-O5-C5	-3.96	106.88	112.19
6	L	6	MAN	C1-O5-C5	3.90	117.41	112.19
6	L	1	NAG	C2-N2-C7	3.86	128.07	122.90
5	G	2	NAG	C1-O5-C5	2.83	115.98	112.19
6	L	4	MAN	C1-O5-C5	2.54	115.59	112.19
5	G	3	BMA	C1-O5-C5	2.30	115.27	112.19
4	J	1	NAG	C1-O5-C5	2.27	115.22	112.19
6	L	6	MAN	O2-C2-C3	-2.21	105.58	110.15
4	J	2	NAG	C1-O5-C5	2.17	115.09	112.19
6	L	1	NAG	O4-C4-C5	-2.15	104.03	109.32
6	L	4	MAN	O2-C2-C3	-2.07	105.87	110.15
6	L	3	BMA	O2-C2-C3	-2.06	105.89	110.15

There are no chirality outliers.

All (20) torsion outliers are listed below:

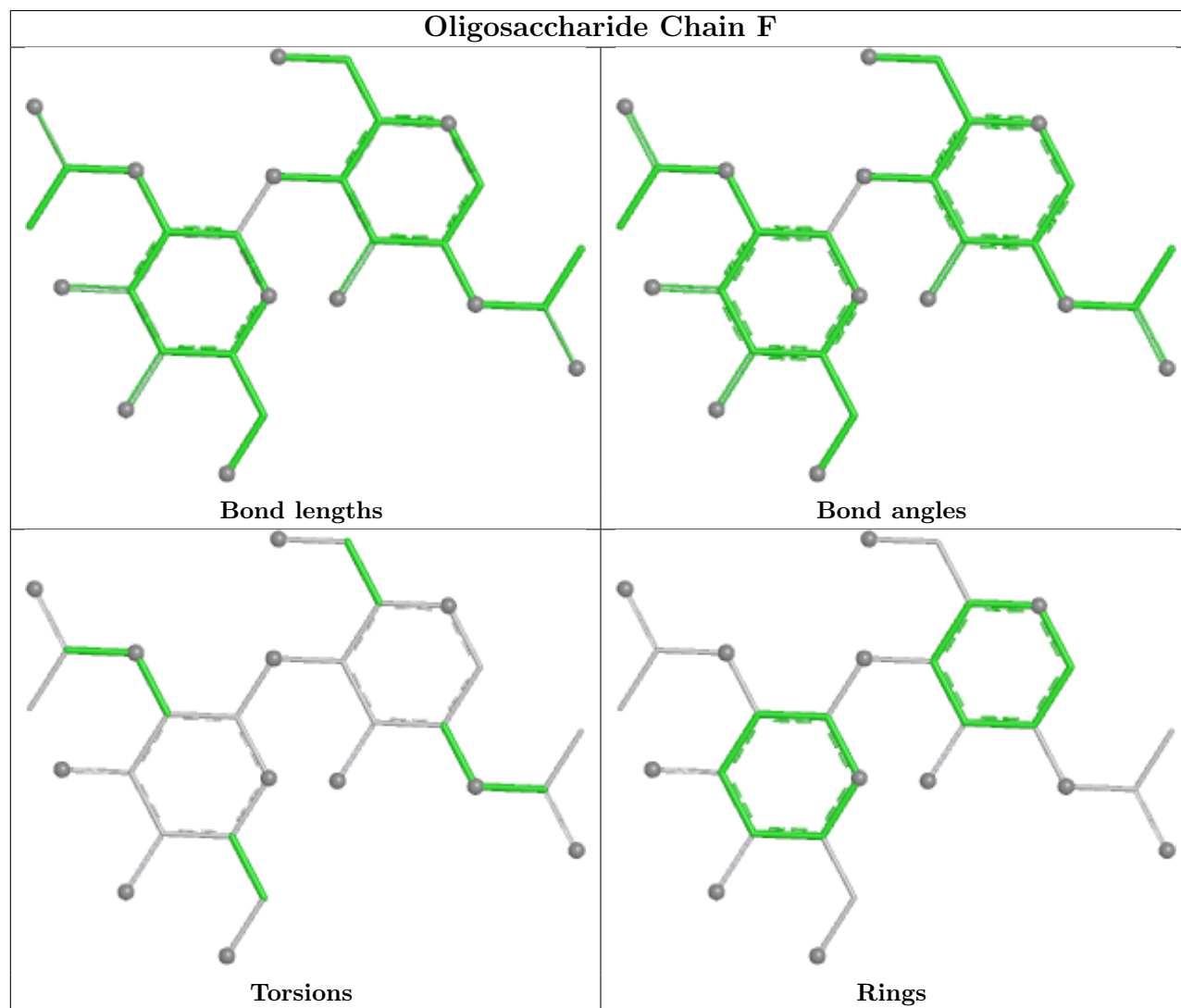
Mol	Chain	Res	Type	Atoms
5	G	2	NAG	C1-C2-N2-C7
6	L	1	NAG	C3-C2-N2-C7
5	K	1	NAG	O5-C5-C6-O6
5	H	1	NAG	O5-C5-C6-O6
5	K	1	NAG	C4-C5-C6-O6
5	H	1	NAG	C4-C5-C6-O6
5	G	3	BMA	O5-C5-C6-O6
6	L	4	MAN	O5-C5-C6-O6
4	J	1	NAG	C4-C5-C6-O6
4	J	1	NAG	O5-C5-C6-O6
6	L	6	MAN	O5-C5-C6-O6
5	G	3	BMA	C4-C5-C6-O6
6	L	4	MAN	C4-C5-C6-O6
4	M	2	NAG	C4-C5-C6-O6
4	M	2	NAG	O5-C5-C6-O6
5	H	2	NAG	C4-C5-C6-O6
5	K	2	NAG	C4-C5-C6-O6
5	H	2	NAG	O5-C5-C6-O6
5	K	2	NAG	O5-C5-C6-O6
6	L	6	MAN	C4-C5-C6-O6

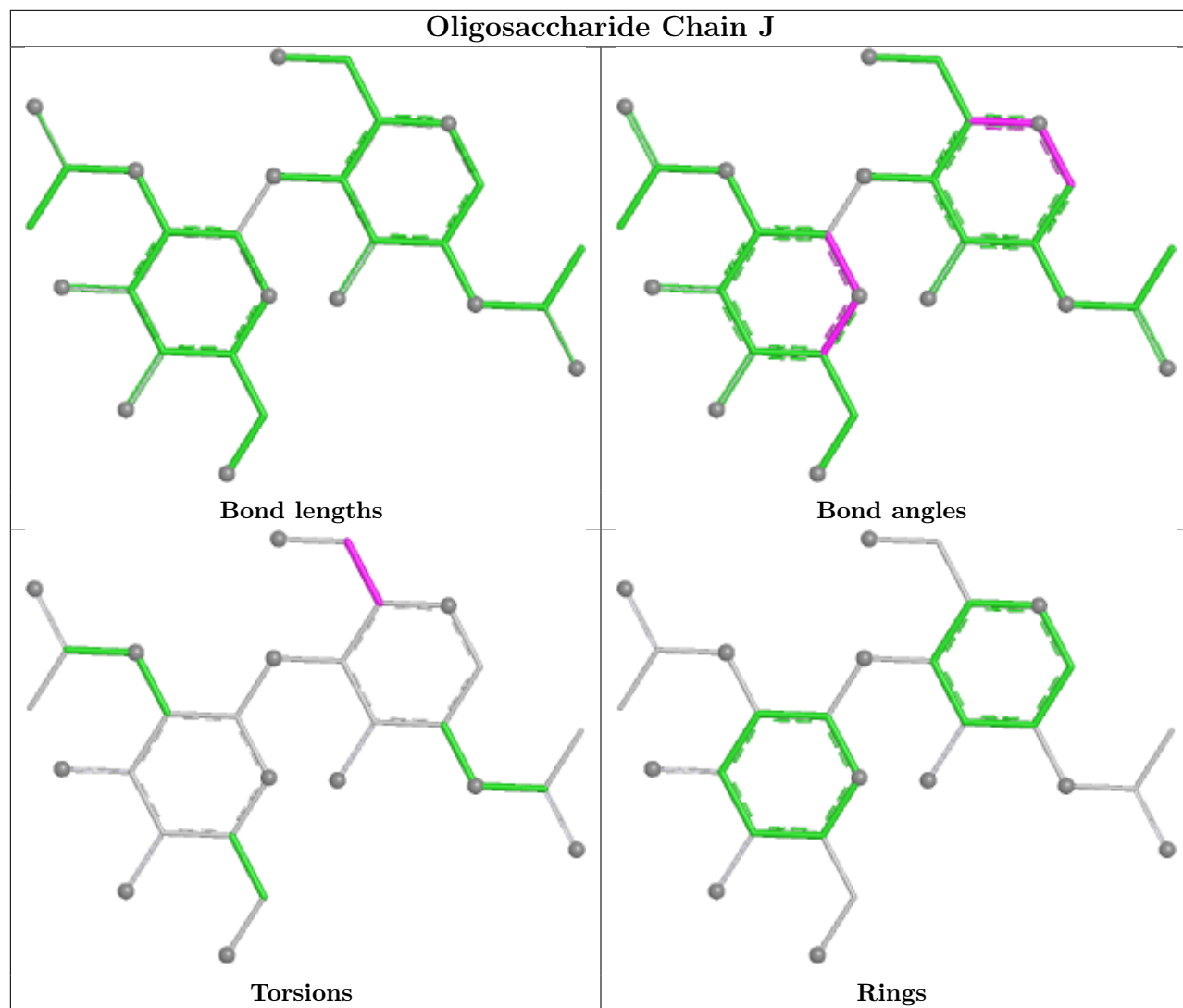
There are no ring outliers.

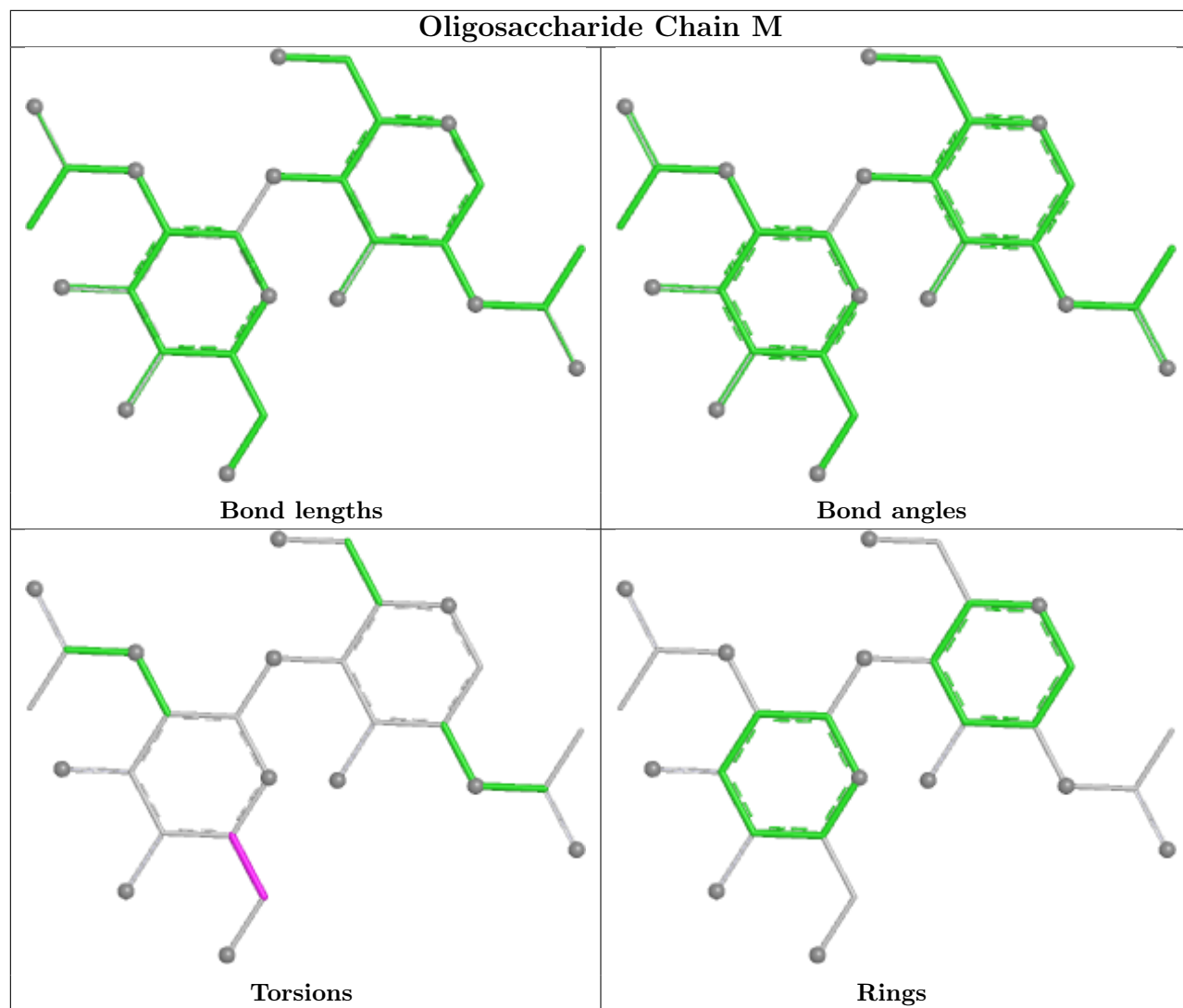
7 monomers are involved in 6 short contacts:

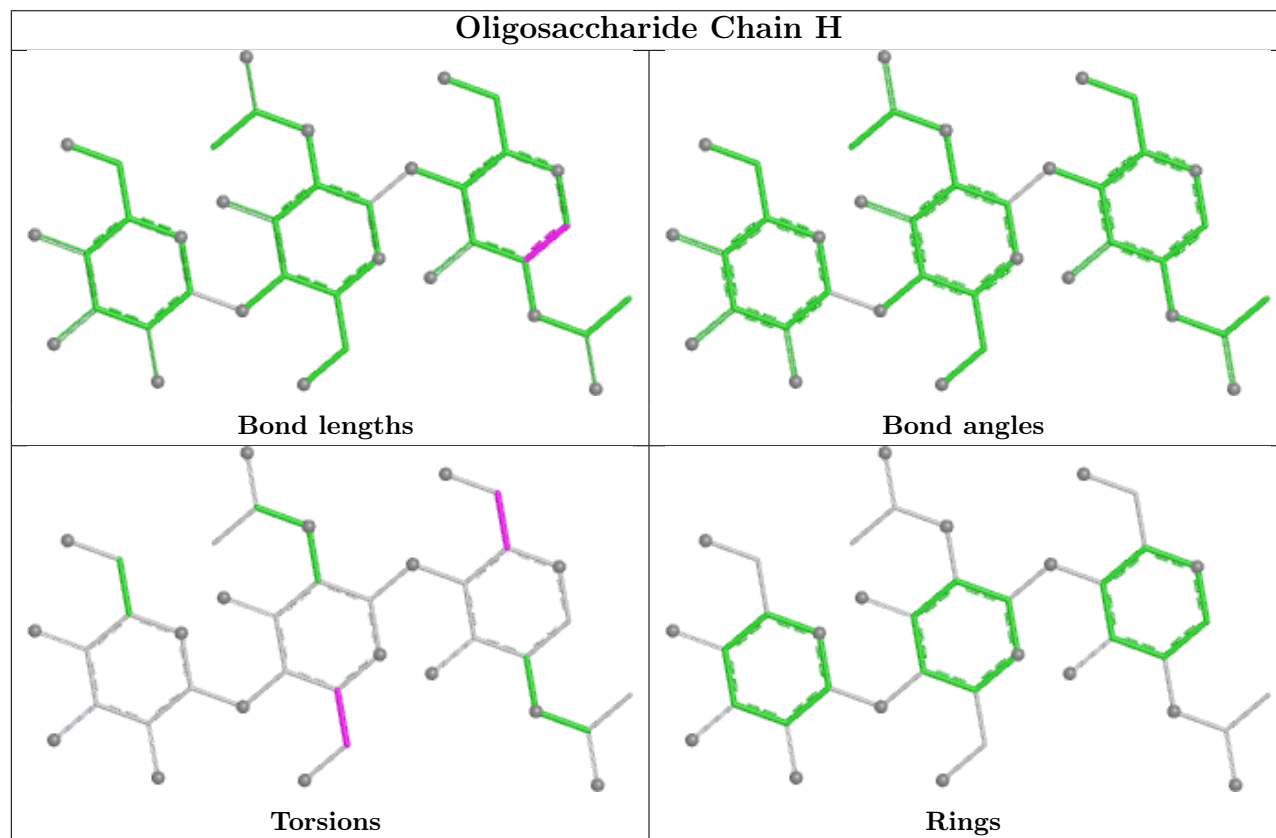
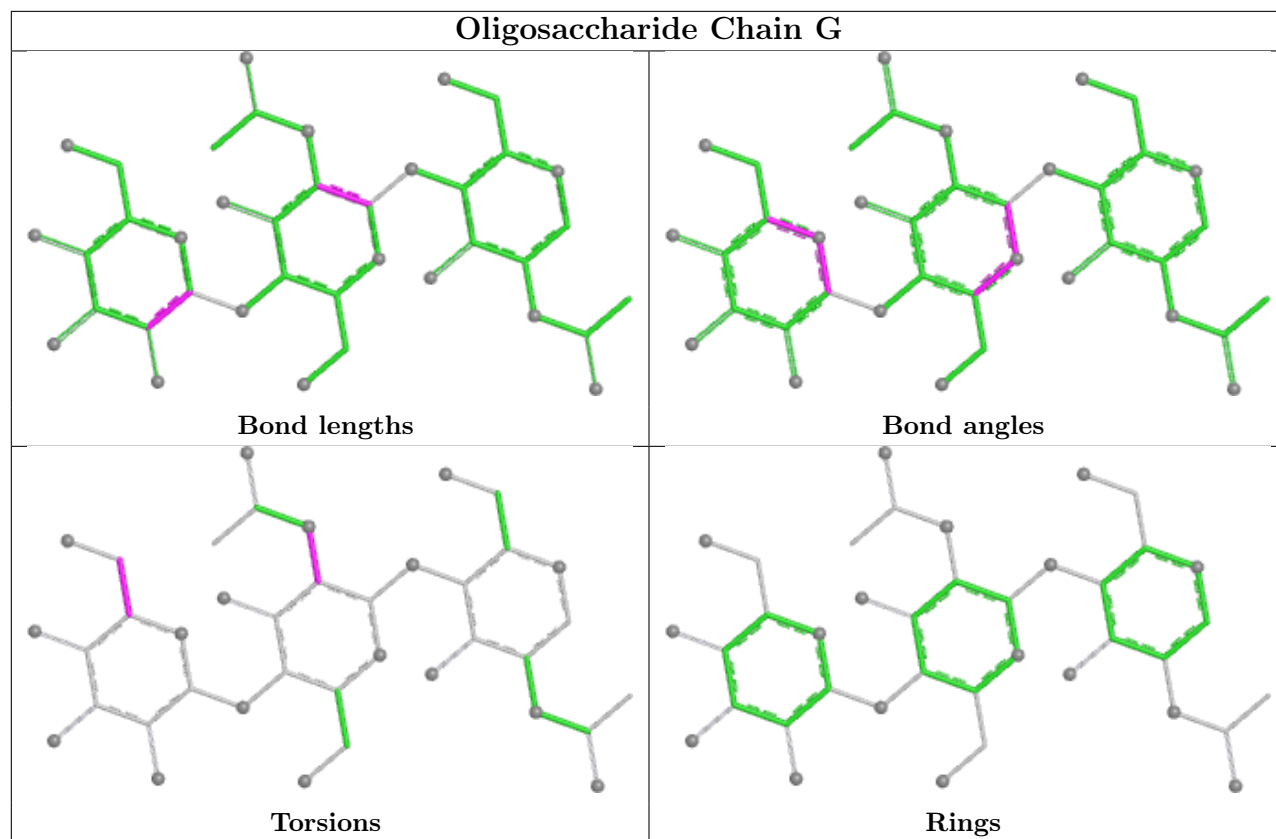
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	J	2	NAG	1	0
4	J	1	NAG	1	0
5	K	2	NAG	2	0
5	G	1	NAG	1	0
6	L	1	NAG	1	0
6	L	3	BMA	1	0
6	L	4	MAN	1	0

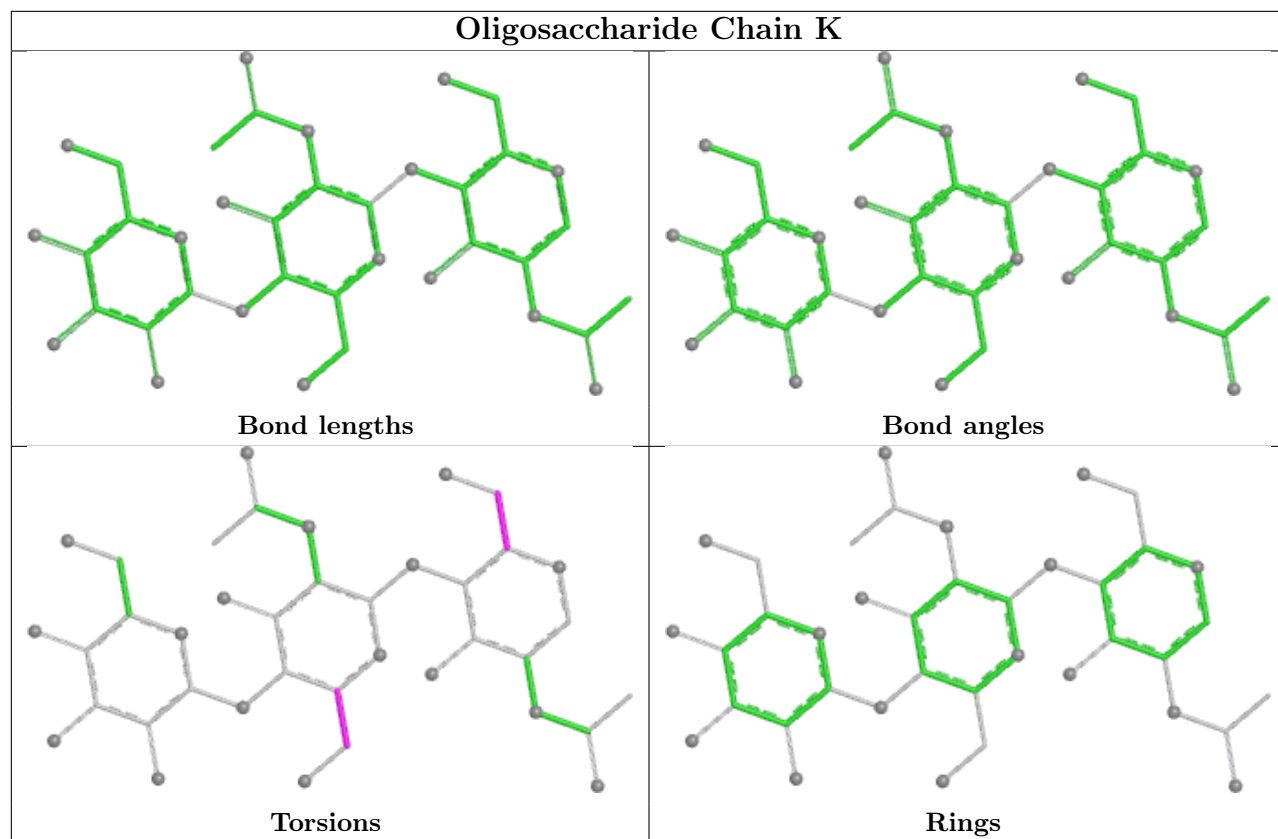
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

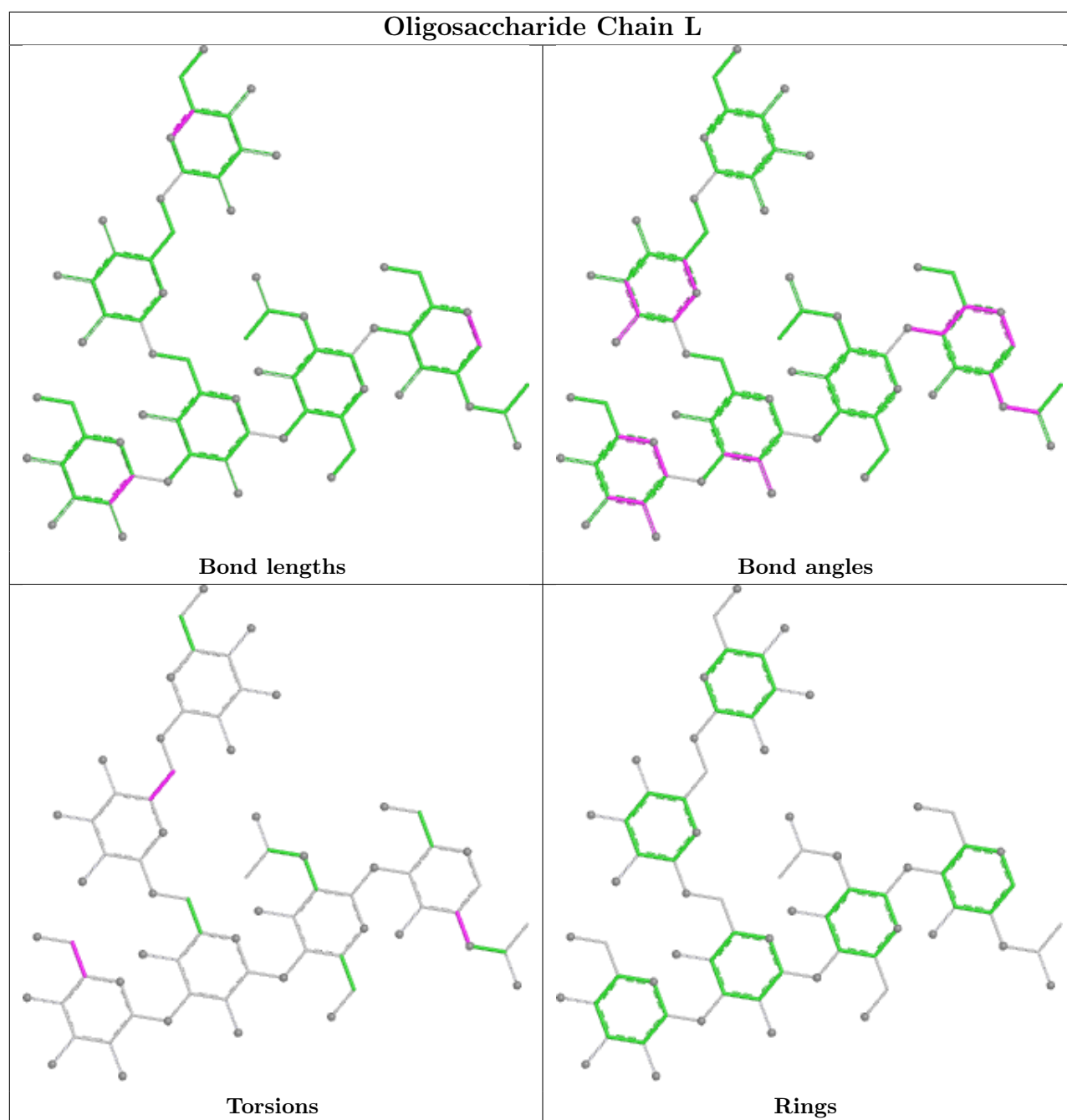












5.6 Ligand geometry [i](#)

7 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
7	NAG	B	201	1	14,14,15	0.36	0	17,19,21	0.46	0
7	NAG	E	304	3	14,14,15	0.88	1 (7%)	17,19,21	0.73	0
8	CYS	I	312	3	5,6,6	1.59	2 (40%)	3,7,7	2.63	2 (66%)
7	NAG	A	201	1	14,14,15	0.40	0	17,19,21	0.77	1 (5%)
7	NAG	D	304	2	14,14,15	0.48	0	17,19,21	0.56	0
7	NAG	E	307	3	14,14,15	0.53	0	17,19,21	0.79	0
8	CYS	E	308	3	5,6,6	1.31	0	3,7,7	1.70	1 (33%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	NAG	B	201	1	-	1/6/23/26	0/1/1/1
7	NAG	E	304	3	-	0/6/23/26	0/1/1/1
8	CYS	I	312	3	-	2/6/6/6	-
7	NAG	A	201	1	-	1/6/23/26	0/1/1/1
7	NAG	D	304	2	-	2/6/23/26	0/1/1/1
7	NAG	E	307	3	-	0/6/23/26	0/1/1/1
8	CYS	E	308	3	-	3/6/6/6	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	E	304	NAG	O5-C1	-2.77	1.39	1.43
8	I	312	CYS	OXT-C	-2.21	1.23	1.30
8	I	312	CYS	CB-CA	2.06	1.55	1.53

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	I	312	CYS	CB-CA-C	3.43	113.31	109.89
8	E	308	CYS	OXT-C-O	-2.94	117.41	124.08
8	I	312	CYS	OXT-C-O	-2.80	117.72	124.08
7	A	201	NAG	C2-N2-C7	2.12	125.75	122.90

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	201	NAG	C1-C2-N2-C7
7	D	304	NAG	C1-C2-N2-C7
8	E	308	CYS	C-CA-CB-SG
8	I	312	CYS	C-CA-CB-SG
8	I	312	CYS	OXT-C-CA-N
7	B	201	NAG	C4-C5-C6-O6
8	E	308	CYS	OXT-C-CA-N
7	D	304	NAG	C3-C2-N2-C7
8	E	308	CYS	O-C-CA-N

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	E	307	NAG	2	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	125/148 (84%)	-0.18	0	100 100	115, 159, 201, 230	0
1	B	126/148 (85%)	-0.05	2 (1%)	70 48	122, 178, 235, 267	0
2	C	204/242 (84%)	-0.24	1 (0%)	87 71	128, 170, 224, 256	0
2	D	202/242 (83%)	-0.02	3 (1%)	72 50	136, 218, 344, 355	0
3	E	213/233 (91%)	-0.02	1 (0%)	87 71	266, 360, 421, 438	0
3	I	213/233 (91%)	-0.18	1 (0%)	87 71	97, 140, 201, 271	0
All	All	1083/1246 (86%)	-0.12	8 (0%)	84 65	97, 181, 383, 438	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	124	ALA	3.9
2	D	50	GLY	3.2
1	B	123	ALA	3.1
3	E	166	ALA	2.6
3	I	71	ASP	2.4
2	D	206	VAL	2.4
2	D	25	MET	2.2
2	C	74	VAL	2.1

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,

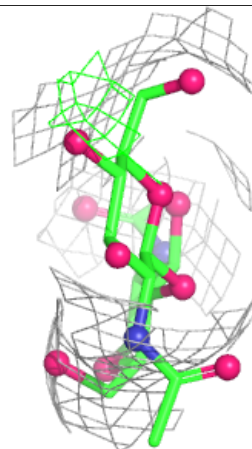
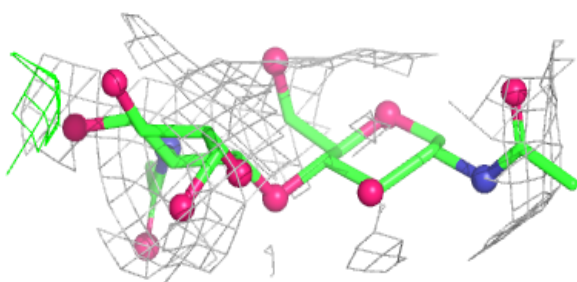
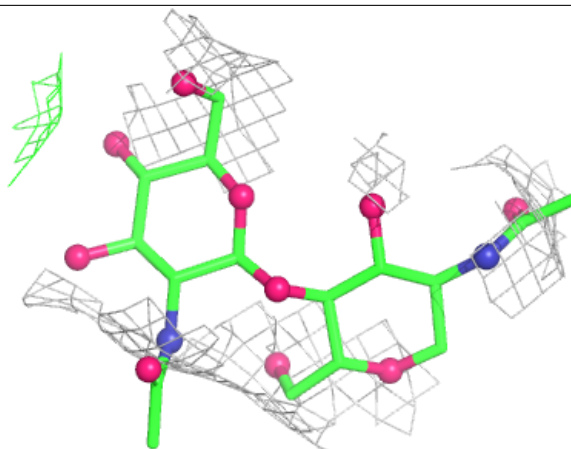
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	NAG	F	1	14/15	-	-	183,183,183,183	0
4	NAG	F	2	14/15	-	-	223,223,223,223	0
5	BMA	H	3	11/12	0.25	0.12	382,382,382,382	0
6	MAN	L	6	11/12	0.33	0.09	264,264,264,264	0
6	MAN	L	5	11/12	0.39	0.08	260,260,260,260	0
5	NAG	K	1	14/15	0.55	0.12	210,210,210,210	0
6	MAN	L	4	11/12	0.60	0.07	261,261,261,261	0
5	BMA	K	3	11/12	0.66	0.06	252,252,252,252	0
5	BMA	G	3	11/12	-	-	244,244,244,244	0
5	NAG	G	2	14/15	0.68	0.11	239,239,239,239	0
4	NAG	J	2	14/15	0.69	0.08	333,333,333,333	0
6	BMA	L	3	11/12	0.70	0.07	240,240,240,240	0
6	NAG	L	2	14/15	0.75	0.11	187,187,187,187	0
6	NAG	L	1	14/15	0.78	0.16	168,168,168,168	0
5	NAG	H	2	14/15	0.78	0.07	394,394,394,394	0
4	NAG	J	1	14/15	0.79	0.13	338,338,338,338	0
4	NAG	M	2	14/15	0.85	0.12	178,178,178,178	0
5	NAG	K	2	14/15	0.86	0.07	228,228,228,228	0
4	NAG	M	1	14/15	0.89	0.09	145,145,145,145	0
5	NAG	H	1	14/15	0.90	0.10	398,398,398,398	0
5	NAG	G	1	14/15	0.93	0.10	218,218,218,218	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.

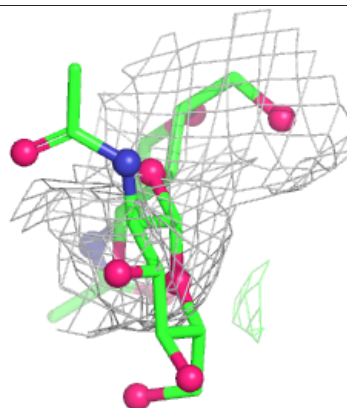
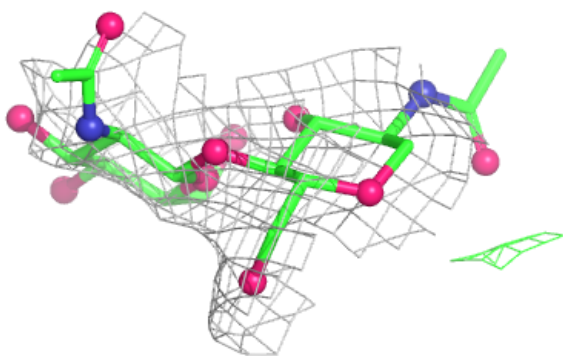
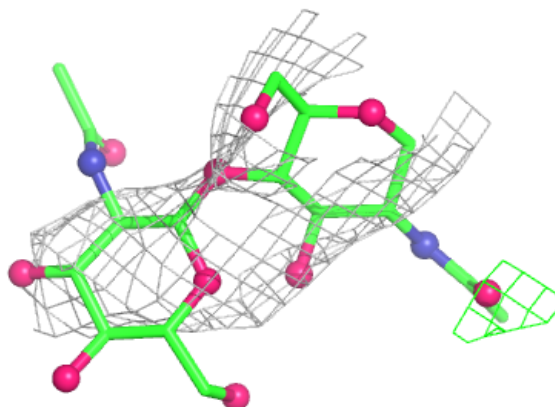
Electron density around Chain F:

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

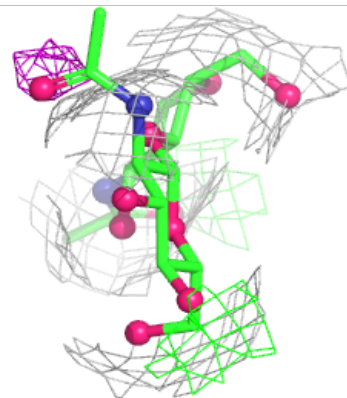
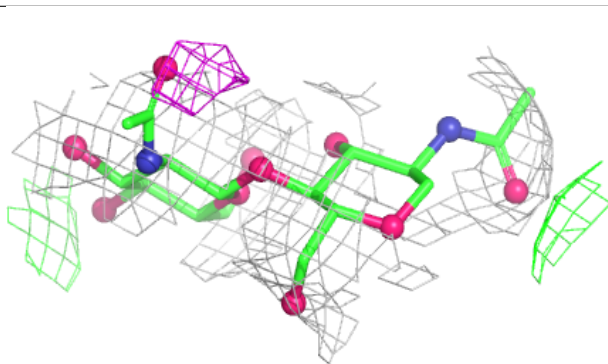
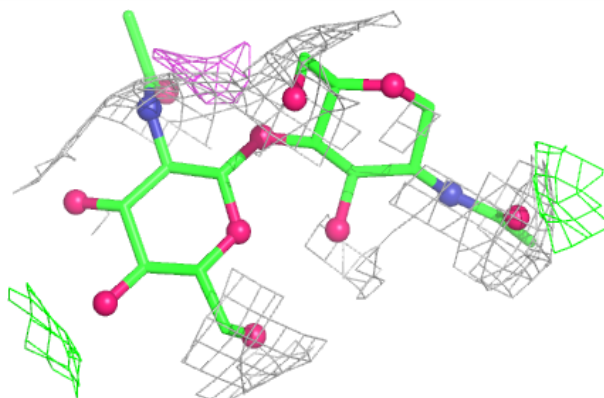


Electron density around Chain J:

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

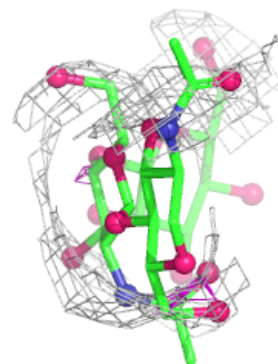
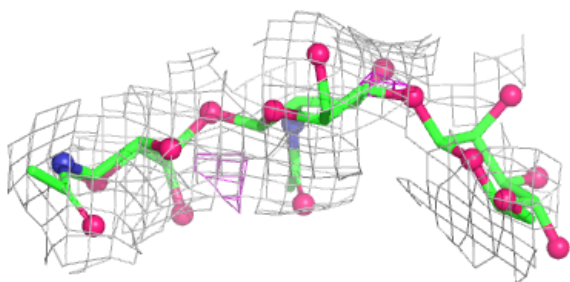
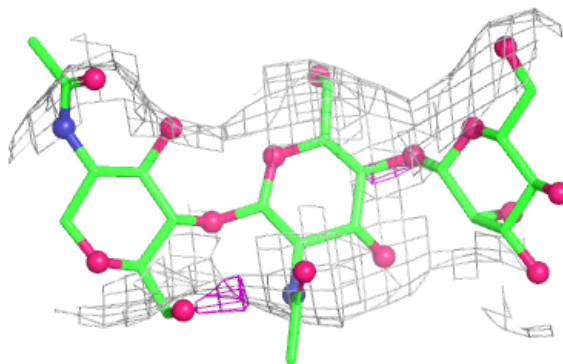
**Electron density around Chain M:**

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



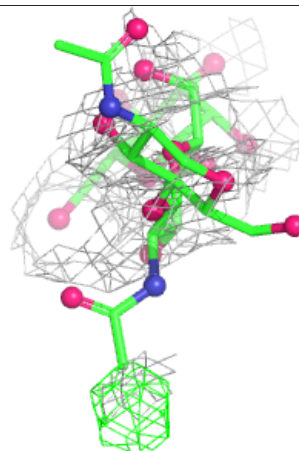
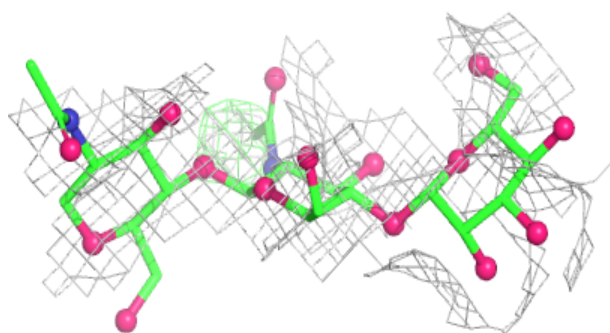
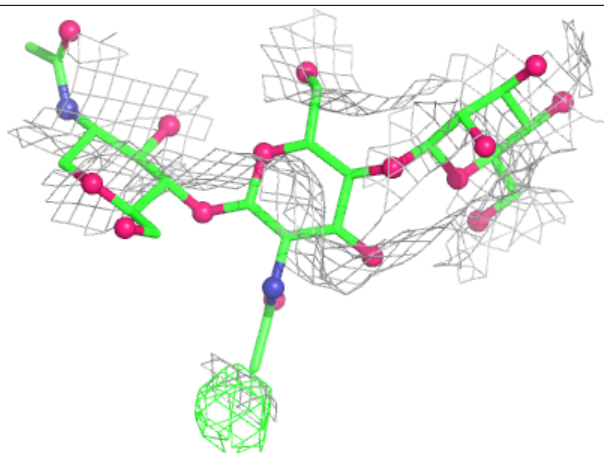
Electron density around Chain G:

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

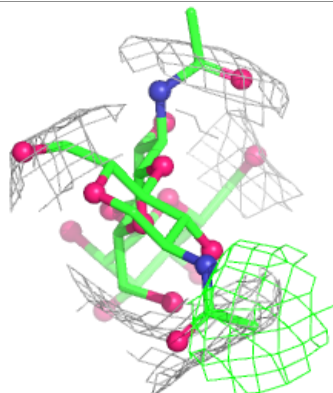
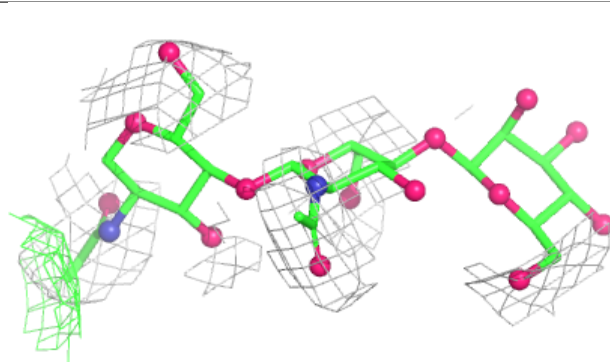
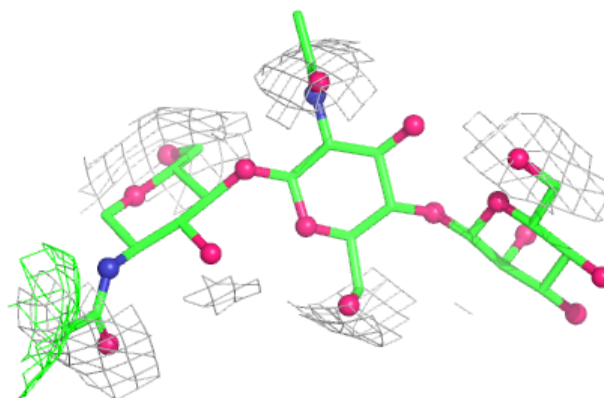


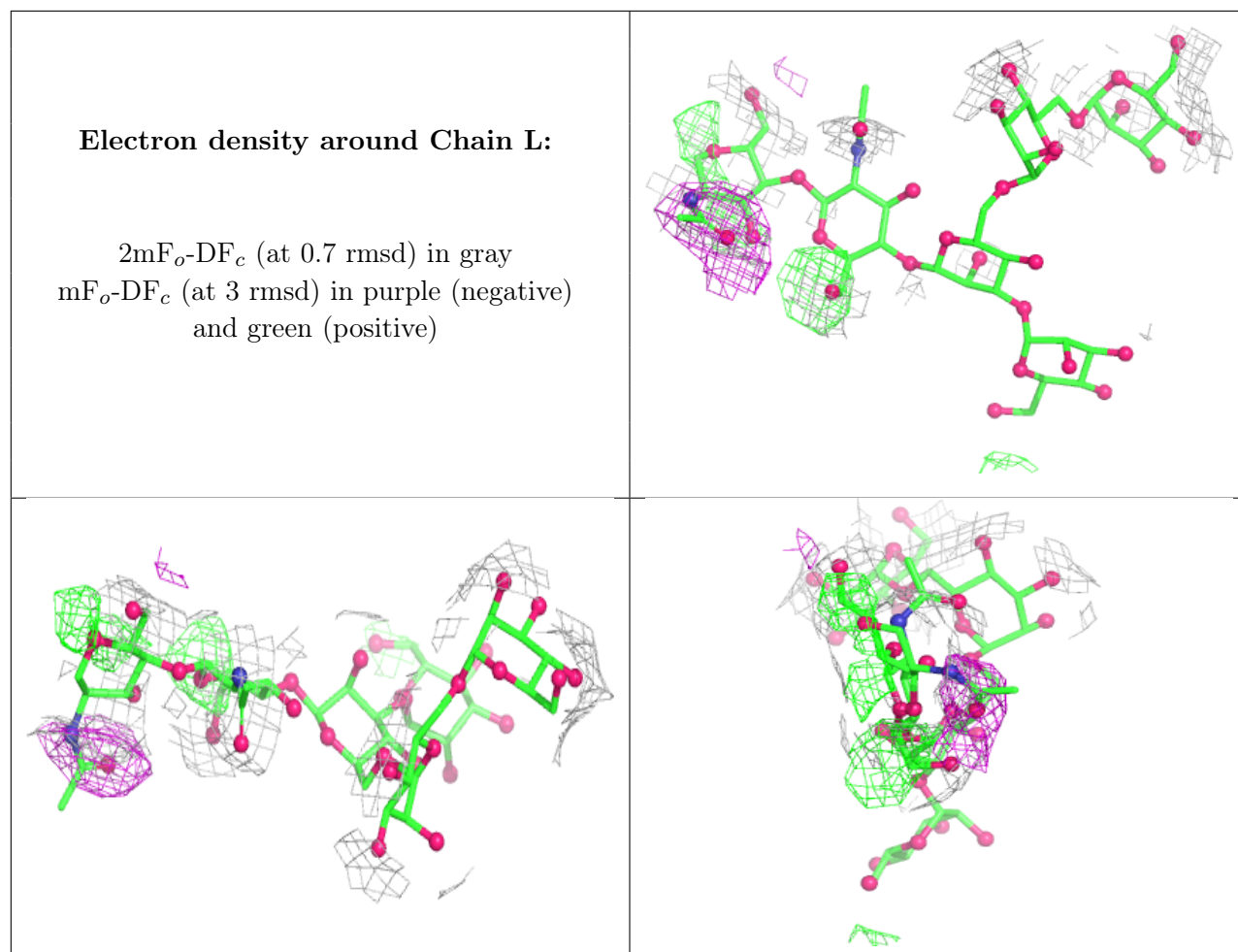
Electron density around Chain H:

$2mF_o-DF_c$ (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)





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	NAG	E	304	14/15	0.05	0.15	338,338,338,338	0
8	CYS	E	308	7/7	0.26	0.09	301,301,301,311	0
7	NAG	E	307	14/15	0.40	0.10	415,415,415,415	0
7	NAG	D	304	14/15	0.71	0.07	247,247,247,247	0
8	CYS	I	312	7/7	0.72	0.11	187,187,187,197	0
7	NAG	B	201	14/15	0.76	0.09	221,221,221,221	0
7	NAG	A	201	14/15	0.78	0.10	233,233,233,233	0

6.5 Other polymers [i](#)

There are no such residues in this entry.