



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:50 AM UTC

PDB ID : 5X7P / pdb_00005x7p
Title : Crystal structure of Paenibacillus sp. 598K alpha-1,6-glucosyltransferase complexed with acarbose
Authors : Fujimoto, Z.; Kishine, N.; Suzuki, N.; Momma, M.; Ichinose, H.; Kimura, A.; Funane, K.
Deposited on : 2017-02-27
Resolution : 2.40 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

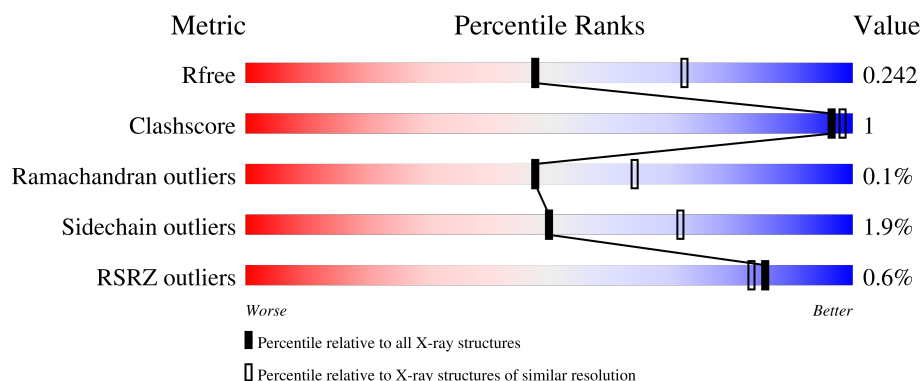
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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.



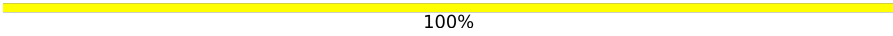
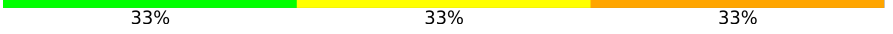
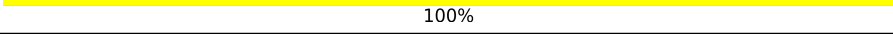
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	1263	% 94% 5% .
1	B	1263	% 94% . .
2	C	3	100%
2	E	3	33% 67%
2	G	3	33% 33% 33%

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Mol	Chain	Length	Quality of chain
2	H	3	 67%33%
2	I	3	 100%
2	J	3	 33%67%
2	L	3	 33%33%33%
3	D	2	 100%
4	F	2	 50%50%
4	K	2	 50%50%

2 Entry composition

There are 11 unique types of molecules in this entry. The entry contains 20523 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 Glycoside hydrolase family 31 alpha-glucosidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1247	Total	C	N	O	S	0	0	0
			9601	6023	1633	1923	22			
1	B	1247	Total	C	N	O	S	0	1	0
			9612	6033	1634	1923	22			

There are 34 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	19	HIS	-	expression tag	UNP A0A193PKW5
A	20	HIS	-	expression tag	UNP A0A193PKW5
A	21	HIS	-	expression tag	UNP A0A193PKW5
A	22	HIS	-	expression tag	UNP A0A193PKW5
A	23	HIS	-	expression tag	UNP A0A193PKW5
A	24	HIS	-	expression tag	UNP A0A193PKW5
A	25	SER	-	expression tag	UNP A0A193PKW5
A	26	SER	-	expression tag	UNP A0A193PKW5
A	27	GLY	-	expression tag	UNP A0A193PKW5
A	28	LEU	-	expression tag	UNP A0A193PKW5
A	29	VAL	-	expression tag	UNP A0A193PKW5
A	30	PRO	-	expression tag	UNP A0A193PKW5
A	31	ARG	-	expression tag	UNP A0A193PKW5
A	32	GLY	-	expression tag	UNP A0A193PKW5
A	33	SER	-	expression tag	UNP A0A193PKW5
A	34	HIS	-	expression tag	UNP A0A193PKW5
A	35	MET	-	expression tag	UNP A0A193PKW5
B	19	HIS	-	expression tag	UNP A0A193PKW5
B	20	HIS	-	expression tag	UNP A0A193PKW5
B	21	HIS	-	expression tag	UNP A0A193PKW5
B	22	HIS	-	expression tag	UNP A0A193PKW5
B	23	HIS	-	expression tag	UNP A0A193PKW5
B	24	HIS	-	expression tag	UNP A0A193PKW5
B	25	SER	-	expression tag	UNP A0A193PKW5
B	26	SER	-	expression tag	UNP A0A193PKW5

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Chain	Residue	Modelled	Actual	Comment	Reference
B	27	GLY	-	expression tag	UNP A0A193PKW5
B	28	LEU	-	expression tag	UNP A0A193PKW5
B	29	VAL	-	expression tag	UNP A0A193PKW5
B	30	PRO	-	expression tag	UNP A0A193PKW5
B	31	ARG	-	expression tag	UNP A0A193PKW5
B	32	GLY	-	expression tag	UNP A0A193PKW5
B	33	SER	-	expression tag	UNP A0A193PKW5
B	34	HIS	-	expression tag	UNP A0A193PKW5
B	35	MET	-	expression tag	UNP A0A193PKW5

- Molecule 2 is an oligosaccharide called 4,6-dideoxy-4-[(1S,4R,5S,6S)-4,5,6-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino}-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	3	Total	C	N	O	0	0	0
			44	25	1	18			
2	E	3	Total	C	N	O	0	0	1
			33	19	1	13			
2	G	3	Total	C	N	O	0	0	1
			33	19	1	13			
2	H	3	Total	C	N	O	0	0	0
			44	25	1	18			
2	I	3	Total	C	N	O	0	0	0
			33	18	1	14			
2	J	3	Total	C	N	O	0	0	1
			33	19	1	13			
2	L	3	Total	C	N	O	0	0	1
			33	19	1	13			

- Molecule 3 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	D	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 4 is an oligosaccharide called 4,6-dideoxy-4-[(1S,4R,5S,6S)-4,5,6-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino}-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	F	2	Total	C	N	O	0	0	1
			22	13	1	8			
4	K	2	Total	C	N	O	0	0	1
			22	13	1	8			

- Molecule 5 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

- Molecule 6 is NICKEL (II) ION (CCD ID: NI) (formula: Ni).

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

- Molecule 7 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

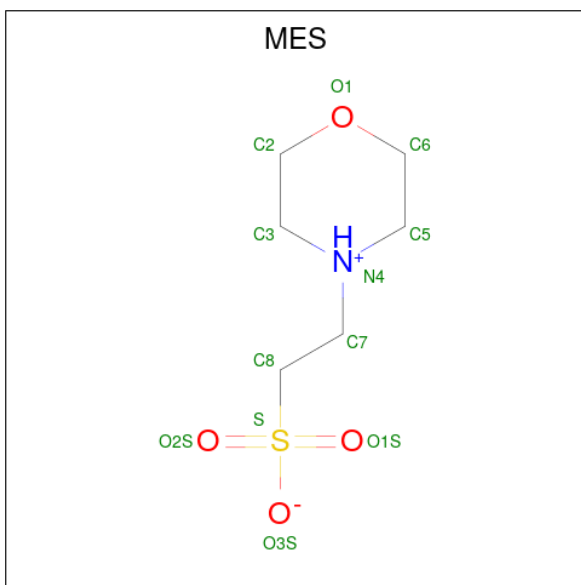
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	6	Total	Mg	0	0
			6	6		
7	B	6	Total	Mg	0	0
			6	6		

- Molecule 8 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	O	S	0	0
			5	4	1		
8	A	1	Total	O	S	0	0
			5	4	1		
8	A	1	Total	O	S	0	0
			5	4	1		
8	A	1	Total	O	S	0	0
			5	4	1		
8	B	1	Total	O	S	0	0
			5	4	1		
8	B	1	Total	O	S	0	0
			5	4	1		
8	B	1	Total	O	S	0	0
			5	4	1		
8	B	1	Total	O	S	0	0
			5	4	1		
8	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 9 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
9	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
9	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
9	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
9	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
9	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
9	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 10 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	A	1	Total	C	O	0	0
			4	2	2		
10	A	1	Total	C	O	0	0
			4	2	2		
10	A	1	Total	C	O	0	0
			4	2	2		
10	A	1	Total	C	O	0	0
			4	2	2		
10	B	1	Total	C	O	0	0
			4	2	2		
10	B	1	Total	C	O	0	0
			4	2	2		
10	B	1	Total	C	O	0	0
			4	2	2		
10	B	1	Total	C	O	0	0
			4	2	2		

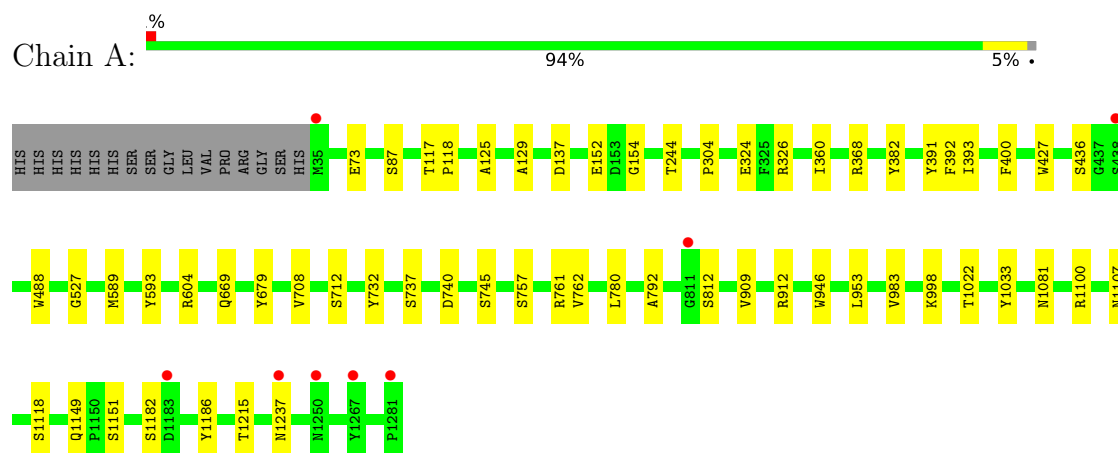
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	391	Total	O	0	0
			391	391		
11	B	403	Total	O	0	0
			403	403		

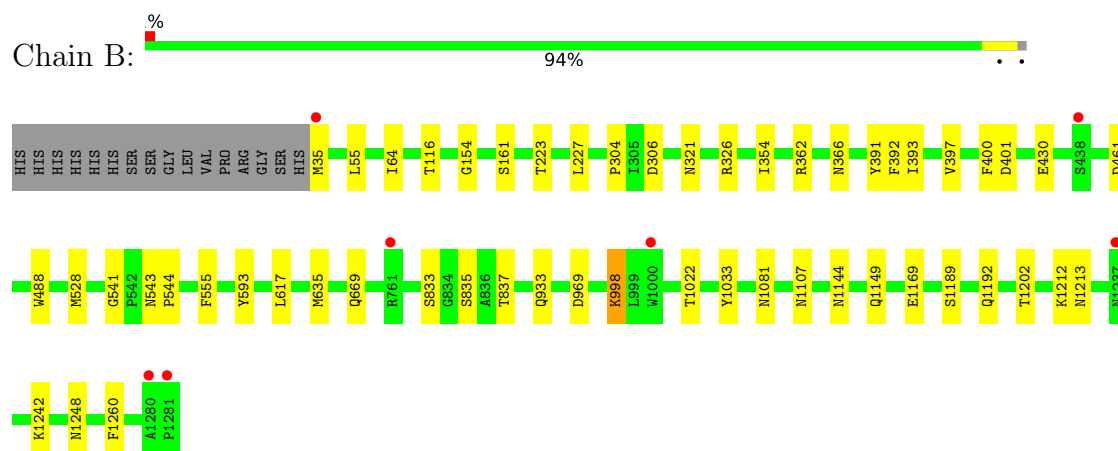
3 Residue-property plots

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

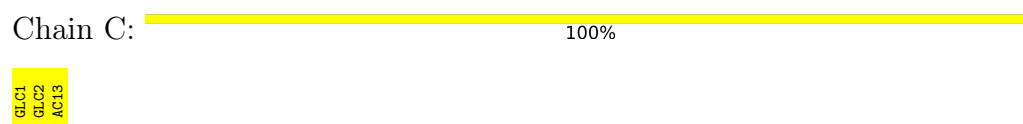
- Molecule 1: Glycoside hydrolase family 31 alpha-glucosidase



- Molecule 1: Glycoside hydrolase family 31 alpha-glucosidase



- Molecule 2: 4,6-dideoxy-4-([(1S,4R,5S,6S)-4,5,6-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

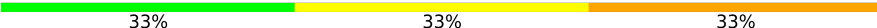


- Molecule 2: 4,6-dideoxy-4- $\{[(1S,4R,5S,6S)-4,5,6\text{-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino}\}$ - α -D-glucopyranose-(1-4)- α -D-glucopyranose-(1-4)- α -D-glucopyranose

Chain E:  33% 67%

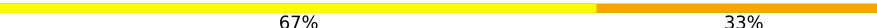


- Molecule 2: 4,6-dideoxy-4- $\{[(1S,4R,5S,6S)-4,5,6\text{-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino}\}$ - α -D-glucopyranose-(1-4)- α -D-glucopyranose-(1-4)- α -D-glucopyranose

Chain G:  33% 33% 33%

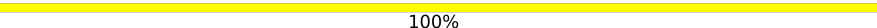


- Molecule 2: 4,6-dideoxy-4- $\{[(1S,4R,5S,6S)-4,5,6\text{-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino}\}$ - α -D-glucopyranose-(1-4)- α -D-glucopyranose-(1-4)- α -D-glucopyranose

Chain H:  67% 33%



- Molecule 2: 4,6-dideoxy-4- $\{[(1S,4R,5S,6S)-4,5,6\text{-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino}\}$ - α -D-glucopyranose-(1-4)- α -D-glucopyranose-(1-4)- α -D-glucopyranose

Chain I:  100%

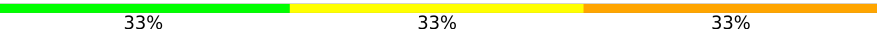


- Molecule 2: 4,6-dideoxy-4- $\{[(1S,4R,5S,6S)-4,5,6\text{-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino}\}$ - α -D-glucopyranose-(1-4)- α -D-glucopyranose-(1-4)- α -D-glucopyranose

Chain J:  33% 67%

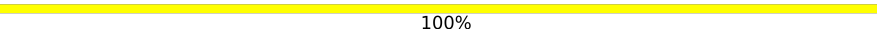


- Molecule 2: 4,6-dideoxy-4- $\{[(1S,4R,5S,6S)-4,5,6\text{-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino}\}$ - α -D-glucopyranose-(1-4)- α -D-glucopyranose-(1-4)- α -D-glucopyranose

Chain L:  33% 33% 33%



- Molecule 3: α -D-glucopyranose-(1-4)- α -D-glucopyranose

Chain D:  100%

GLC1
GLC2

- Molecule 4: 4,6-dideoxy-4-{[(1S,4R,5S,6S)-4,5,6-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino}-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain F:  50% 50%

GLC1
AC12

- Molecule 4: 4,6-dideoxy-4-{[(1S,4R,5S,6S)-4,5,6-trihydroxy-3-(hydroxymethyl)cyclohex-2-en-1-yl]amino}-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain K:  50% 50%

GLC1
AC12

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	184.29Å 271.52Å 133.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	152.48 – 2.40 152.48 – 2.40	Depositor EDS
% Data completeness (in resolution range)	97.1 (152.48-2.40) 97.1 (152.48-2.40)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.64 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.180 , 0.237 0.190 , 0.242	Depositor DCC
R_{free} test set	6536 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	39.0	Xtriage
Anisotropy	0.727	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 32.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.37$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	20523	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.45% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: GLC, EDO, CA, AC1, MES, NI, SO4, MG

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	0.94	1/9863 (0.0%)	0.94	5/13473 (0.0%)
1	B	0.94	1/9879 (0.0%)	0.96	2/13496 (0.0%)
All	All	0.94	2/19742 (0.0%)	0.95	7/26969 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	679	TYR	C-O	-6.12	1.20	1.25
1	B	998	LYS	N-CA	5.15	1.52	1.46

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	154	GLY	N-CA-C	6.74	119.21	110.45
1	A	154	GLY	N-CA-C	5.98	118.22	110.45
1	A	708	VAL	N-CA-C	5.51	117.94	111.05
1	A	983	VAL	CA-C-N	-5.50	114.30	119.85
1	A	983	VAL	C-N-CA	-5.50	114.30	119.85
1	B	541	GLY	N-CA-C	5.39	118.47	112.00
1	A	812	SER	N-CA-C	5.08	117.24	110.53

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within

the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	9601	0	8954	23	1
1	B	9612	0	8964	22	0
2	C	44	0	30	0	0
2	E	33	0	19	0	0
2	G	33	0	19	1	0
2	H	44	0	30	1	0
2	I	33	0	21	0	0
2	J	33	0	19	0	0
2	L	33	0	19	1	0
3	D	23	0	21	0	0
4	F	22	0	10	0	0
4	K	22	0	10	0	1
5	A	3	0	0	0	0
5	B	3	0	0	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
7	A	6	0	0	0	0
7	B	6	0	0	0	0
8	A	25	0	0	0	0
8	B	35	0	0	0	0
9	A	24	0	26	0	0
9	B	60	0	65	0	0
10	A	16	0	24	0	0
10	B	16	0	24	0	0
11	A	391	0	0	0	0
11	B	403	0	0	0	0
All	All	20523	0	18255	38	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 1.

All (38) 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:362:ARG:HD3	1:B:366:ASN:OD1	2.06	0.54
1:B:161:SER:HB2	1:B:223:THR:O	2.10	0.52
1:A:360:ILE:HD13	1:B:397:VAL:HG12	1.93	0.51
1:A:909:VAL:HG21	1:A:953:LEU:HD12	1.91	0.51
1:B:430:GLU:OE2	2:H:3:AC1:O3	2.30	0.50
1:A:593:TYR:CE1	1:B:304:PRO:HB2	2.46	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:391:TYR:HB3	1:B:391:TYR:HB3	1.93	0.49
1:B:1169:GLU:OE1	2:L:3:AC1:O3	2.30	0.49
1:B:543:ASN:HB3	1:B:544:PRO:HD2	1.95	0.48
1:A:1182:SER:HB3	1:A:1186:TYR:CD2	2.49	0.47
1:B:528:MET:HE1	1:B:555:PHE:CE2	2.50	0.47
1:A:669:GLN:HG2	1:A:1033:TYR:CE1	2.49	0.47
1:B:1189:SER:HA	1:B:1260:PHE:O	2.15	0.47
1:A:604:ARG:HG3	1:A:732:TYR:CD2	2.50	0.46
1:B:1144:ASN:ND2	1:B:1192:GLN:HG3	2.30	0.46
1:A:137:ASP:OD1	1:A:137:ASP:C	2.59	0.46
1:A:304:PRO:HB2	1:B:593:TYR:CZ	2.51	0.45
1:B:669:GLN:HG2	1:B:1033:TYR:CE1	2.52	0.45
1:B:321:ASN:O	1:B:326:ARG:NH1	2.50	0.44
1:A:117:THR:HB	1:A:118:PRO:HD2	1.99	0.44
1:A:368:ARG:HE	2:G:2:GLC:H61	1.82	0.44
1:A:393:ILE:HD12	1:B:393:ILE:HD12	2.00	0.44
1:A:912:ARG:NH2	1:A:946:TRP:HB2	2.32	0.44
1:A:762:VAL:HG11	1:A:780:LEU:HD13	2.00	0.43
1:A:1215:THR:HA	1:A:1237:ASN:HD22	1.83	0.43
1:B:1144:ASN:HD22	1:B:1192:GLN:HG3	1.84	0.42
1:B:528:MET:HE1	1:B:555:PHE:CD2	2.54	0.42
1:A:125:ALA:O	1:A:129:ALA:HA	2.18	0.42
1:B:969:ASP:OD2	1:B:1242:LYS:NZ	2.51	0.42
1:A:87:SER:HG	1:A:244:THR:H	1.68	0.41
1:A:1022:THR:O	1:A:1100:ARG:NH1	2.50	0.41
1:B:1202:THR:CG2	1:B:1248:ASN:HA	2.50	0.41
1:A:740:ASP:HB3	1:A:745:SER:HB3	2.02	0.41
1:A:427:TRP:HB3	1:B:354:ILE:HA	2.02	0.41
1:A:382:TYR:HA	1:B:401:ASP:HB2	2.01	0.41
1:A:488:TRP:HA	1:A:527:GLY:O	2.20	0.41
1:B:617:LEU:HD23	1:B:635:MET:HG2	2.04	0.40
1:A:604:ARG:HG3	1:A:732:TYR:CE2	2.57	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)
1:A:1215:THR:OG1	4:K:2:AC1:O6B[6_555]	2.02	0.18

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	1245/1263 (99%)	1207 (97%)	37 (3%)	1 (0%)	48	64
1	B	1246/1263 (99%)	1206 (97%)	39 (3%)	1 (0%)	48	64
All	All	2491/2526 (99%)	2413 (97%)	76 (3%)	2 (0%)	48	64

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	1213	ASN
1	A	792	ALA

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	997/1011 (99%)	979 (98%)	18 (2%)	51	73
1	B	998/1011 (99%)	978 (98%)	20 (2%)	48	70
All	All	1995/2022 (99%)	1957 (98%)	38 (2%)	50	71

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	73	GLU
1	A	152	GLU
1	A	324	GLU
1	A	326	ARG

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Mol	Chain	Res	Type
1	A	392	PHE
1	A	400	PHE
1	A	436	SER
1	A	589	MET
1	A	712	SER
1	A	737	SER
1	A	757	SER
1	A	761	ARG
1	A	998	LYS
1	A	1081	ASN
1	A	1107	ASN
1	A	1118	SER
1	A	1149	GLN
1	A	1151	SER
1	B	35	MET
1	B	55	LEU
1	B	64	ILE
1	B	116	THR
1	B	227	LEU
1	B	306	ASP
1	B	392	PHE
1	B	400	PHE
1	B	461	ASP
1	B	488	TRP
1	B	833	SER
1	B	835	SER
1	B	837	THR
1	B	933	GLN
1	B	998	LYS
1	B	1022	THR
1	B	1081	ASN
1	B	1107	ASN
1	B	1149	GLN
1	B	1212	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	190	HIS
1	A	302	ASN
1	A	503	ASN
1	A	949	GLN

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Mol	Chain	Res	Type
1	A	977	ASN
1	A	1107	ASN
1	A	1149	GLN
1	A	1237	ASN
1	B	339	GLN
1	B	440	GLN
1	B	776	GLN
1	B	777	HIS
1	B	977	ASN
1	B	1009	HIS
1	B	1107	ASN
1	B	1144	ASN
1	B	1200	ASN
1	B	1213	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

Of 27 monosaccharides modelled in this entry, 21 were used for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	GLC	C	1	2	12,12,12	0.52	0	17,17,17	0.85	1 (5%)
2	GLC	C	2	2	11,11,12	0.72	0	15,15,17	1.07	1 (6%)
2	AC1	C	3	2	21,22,23	0.93	0	22,32,34	1.45	4 (18%)
3	GLC	D	1	3	12,12,12	0.76	0	17,17,17	1.43	3 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GLC	D	2	3	11,11,12	1.29	1 (9%)	15,15,17	1.83	3 (20%)
2	GLC	E	2	2	11,11,12	0.71	0	15,15,17	1.48	5 (33%)
2	AC1	E	3	2	21,22,23	1.55	3 (14%)	22,32,34	1.00	1 (4%)
4	AC1	F	2	4	21,22,23	0.97	1 (4%)	22,32,34	1.32	2 (9%)
2	GLC	G	2	2	11,11,12	0.62	0	15,15,17	1.61	3 (20%)
2	AC1	G	3	2	21,22,23	1.29	3 (14%)	22,32,34	1.58	4 (18%)
2	GLC	H	1	2	12,12,12	0.70	0	17,17,17	0.96	1 (5%)
2	GLC	H	2	2	11,11,12	0.61	0	15,15,17	1.05	1 (6%)
2	AC1	H	3	2	21,22,23	1.03	1 (4%)	22,32,34	1.73	6 (27%)
2	GLC	I	1	2	12,12,12	0.98	0	17,17,17	1.40	2 (11%)
2	GLC	I	2	2	11,11,12	0.80	0	15,15,17	1.35	2 (13%)
2	AC1	I	3	2	10,10,23	1.18	0	12,14,34	1.80	4 (33%)
2	GLC	J	2	2	11,11,12	0.75	0	15,15,17	1.42	2 (13%)
2	AC1	J	3	2	21,22,23	1.04	1 (4%)	22,32,34	1.36	3 (13%)
4	AC1	K	2	4	21,22,23	1.00	1 (4%)	22,32,34	1.59	4 (18%)
2	GLC	L	2	2	11,11,12	0.81	0	15,15,17	2.02	3 (20%)
2	AC1	L	3	2	21,22,23	0.77	1 (4%)	22,32,34	1.04	1 (4%)

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	GLC	C	1	2	-	2/2/22/22	0/1/1/1
2	GLC	C	2	2	-	0/2/19/22	0/1/1/1
2	AC1	C	3	2	-	3/6/43/46	0/2/2/2
3	GLC	D	1	3	-	0/2/22/22	0/1/1/1
3	GLC	D	2	3	-	2/2/19/22	0/1/1/1
2	GLC	E	2	2	-	2/2/19/22	0/1/1/1
2	AC1	E	3	2	-	3/6/43/46	0/2/2/2
4	AC1	F	2	4	-	1/6/43/46	0/2/2/2
2	GLC	G	2	2	-	1/2/19/22	0/1/1/1
2	AC1	G	3	2	-	2/6/43/46	0/2/2/2
2	GLC	H	1	2	-	0/2/22/22	0/1/1/1
2	GLC	H	2	2	-	0/2/19/22	0/1/1/1
2	AC1	H	3	2	-	3/6/43/46	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GLC	I	1	2	-	0/2/22/22	0/1/1/1
2	GLC	I	2	2	-	2/2/19/22	0/1/1/1
2	AC1	I	3	2	-	-	0/1/1/2
2	GLC	J	2	2	-	0/2/19/22	0/1/1/1
2	AC1	J	3	2	-	3/6/43/46	0/2/2/2
4	AC1	K	2	4	-	2/6/43/46	0/2/2/2
2	GLC	L	2	2	-	1/2/19/22	0/1/1/1
2	AC1	L	3	2	-	1/6/43/46	0/2/2/2

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	3	AC1	C7B-C5B	4.08	1.38	1.32
2	E	3	AC1	C1B-C7B	2.98	1.54	1.50
2	G	3	AC1	C4A-C5B	2.91	1.53	1.51
2	H	3	AC1	C2B-C1B	2.81	1.56	1.53
2	J	3	AC1	C1B-C7B	2.71	1.54	1.50
2	G	3	AC1	C2B-C1B	2.59	1.56	1.53
3	D	2	GLC	C1-C2	2.52	1.58	1.52
2	G	3	AC1	C7B-C5B	2.51	1.36	1.32
2	E	3	AC1	C4A-C5B	2.34	1.53	1.51
4	K	2	AC1	C7B-C5B	2.26	1.36	1.32
4	F	2	AC1	C7B-C5B	2.17	1.36	1.32
2	L	3	AC1	O4-C4A	2.02	1.46	1.42

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	2	GLC	C1-O5-C5	5.88	120.06	112.19
3	D	2	GLC	C1-O5-C5	5.06	118.97	112.19
2	I	3	AC1	C1-O5-C5	4.19	122.85	112.97
4	K	2	AC1	C1-C2-C3	4.18	115.73	109.64
4	K	2	AC1	C7B-C1B-N4A	-3.90	104.94	110.68
4	F	2	AC1	C1-C2-C3	3.84	115.23	109.64
2	G	3	AC1	O2B-C2B-C1B	3.69	116.35	109.08
2	I	2	GLC	C1-O5-C5	3.68	117.11	112.19
4	F	2	AC1	C1-O5-C5	3.50	121.22	112.97
2	H	3	AC1	C7B-C1B-N4A	-3.44	105.62	110.68
2	G	2	GLC	C1-O5-C5	3.40	116.75	112.19
2	L	3	AC1	C7B-C1B-N4A	-3.29	105.84	110.68
3	D	2	GLC	O2-C2-C1	3.22	116.59	109.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	3	AC1	C1-O5-C5	3.18	120.46	112.97
2	J	2	GLC	C3-C4-C5	-3.15	104.52	110.23
2	C	3	AC1	O6B-C6B-C5B	-3.09	103.04	112.74
2	I	1	GLC	C1-O5-C5	3.01	119.47	113.65
2	G	2	GLC	C3-C4-C5	2.91	115.51	110.23
2	H	3	AC1	O2B-C2B-C1B	2.87	114.73	109.08
2	J	3	AC1	O4-C4A-C5B	-2.85	105.32	110.75
3	D	1	GLC	C1-O5-C5	2.83	119.12	113.65
2	H	3	AC1	O4-C4A-C5B	2.81	116.10	110.75
3	D	2	GLC	C1-C2-C3	2.71	113.58	109.64
2	H	3	AC1	O6B-C6B-C5B	-2.66	104.39	112.74
4	K	2	AC1	C3B-C4A-C5B	-2.64	107.53	111.67
2	J	3	AC1	O2-C2-C3	-2.54	104.88	110.15
2	L	2	GLC	O4-C4-C3	-2.49	104.51	110.38
3	D	1	GLC	O3-C3-C2	-2.45	104.60	110.38
2	H	1	GLC	C1-O5-C5	2.44	118.37	113.65
2	I	3	AC1	O2-C2-C3	2.39	115.11	110.15
2	G	3	AC1	O3B-C3B-C2B	-2.32	104.91	110.38
2	C	3	AC1	O2B-C2B-C3B	2.32	115.84	110.38
2	G	3	AC1	O2-C2-C1	2.30	114.49	109.22
2	G	2	GLC	O4-C4-C3	-2.29	104.98	110.38
2	I	3	AC1	C1-C2-C3	-2.28	106.33	109.64
2	G	3	AC1	C7B-C1B-N4A	-2.26	107.35	110.68
2	E	2	GLC	C1-O5-C5	2.25	115.20	112.19
2	L	2	GLC	O5-C5-C4	2.25	116.29	110.83
2	E	2	GLC	O4-C4-C3	-2.22	105.13	110.38
3	D	1	GLC	O5-C5-C4	2.21	113.69	109.70
2	E	2	GLC	O5-C1-C2	-2.20	105.55	110.79
2	C	2	GLC	C1-O5-C5	2.18	115.11	112.19
2	E	3	AC1	O2-C2-C1	2.16	114.17	109.22
2	E	2	GLC	O2-C2-C1	2.15	114.15	109.22
2	C	1	GLC	C1-O5-C5	2.15	117.81	113.65
2	H	3	AC1	O2-C2-C3	-2.15	105.70	110.15
2	I	3	AC1	O3-C3-C2	2.14	114.42	110.05
2	J	3	AC1	O3-C3-C4	2.13	113.81	109.58
2	H	3	AC1	C5-C4-N4A	2.10	117.52	111.74
2	C	3	AC1	C7B-C1B-N4A	-2.09	107.60	110.68
4	K	2	AC1	O6B-C6B-C5B	-2.09	106.16	112.74
2	I	2	GLC	O2-C2-C1	2.08	113.98	109.22
2	H	2	GLC	C1-O5-C5	2.08	114.97	112.19
2	J	2	GLC	O5-C5-C6	2.06	111.67	107.66
2	E	2	GLC	C2-C3-C4	-2.04	107.27	110.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	1	GLC	C1-C2-C3	2.04	114.51	110.36

There are no chirality outliers.

All (28) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	3	AC1	C4A-C5B-C6B-O6B
2	C	3	AC1	C7B-C5B-C6B-O6B
2	E	3	AC1	C4A-C5B-C6B-O6B
2	E	3	AC1	C7B-C5B-C6B-O6B
2	G	3	AC1	C4A-C5B-C6B-O6B
2	G	3	AC1	C7B-C5B-C6B-O6B
2	H	3	AC1	C4A-C5B-C6B-O6B
2	H	3	AC1	C7B-C5B-C6B-O6B
2	J	3	AC1	C4A-C5B-C6B-O6B
2	J	3	AC1	C7B-C5B-C6B-O6B
4	K	2	AC1	C7B-C5B-C6B-O6B
3	D	2	GLC	C4-C5-C6-O6
2	E	2	GLC	O5-C5-C6-O6
2	E	2	GLC	C4-C5-C6-O6
2	C	1	GLC	O5-C5-C6-O6
3	D	2	GLC	O5-C5-C6-O6
2	C	1	GLC	C4-C5-C6-O6
2	I	2	GLC	C4-C5-C6-O6
2	J	3	AC1	C5-C4-N4A-C1B
2	L	3	AC1	C3-C4-N4A-C1B
4	F	2	AC1	C5-C4-N4A-C1B
2	C	3	AC1	C7B-C1B-N4A-C4
2	H	3	AC1	C7B-C1B-N4A-C4
2	G	2	GLC	C4-C5-C6-O6
2	I	2	GLC	O5-C5-C6-O6
4	K	2	AC1	C4A-C5B-C6B-O6B
2	L	2	GLC	C4-C5-C6-O6
2	E	3	AC1	C5-C4-N4A-C1B

There are no ring outliers.

4 monomers are involved in 4 short contacts:

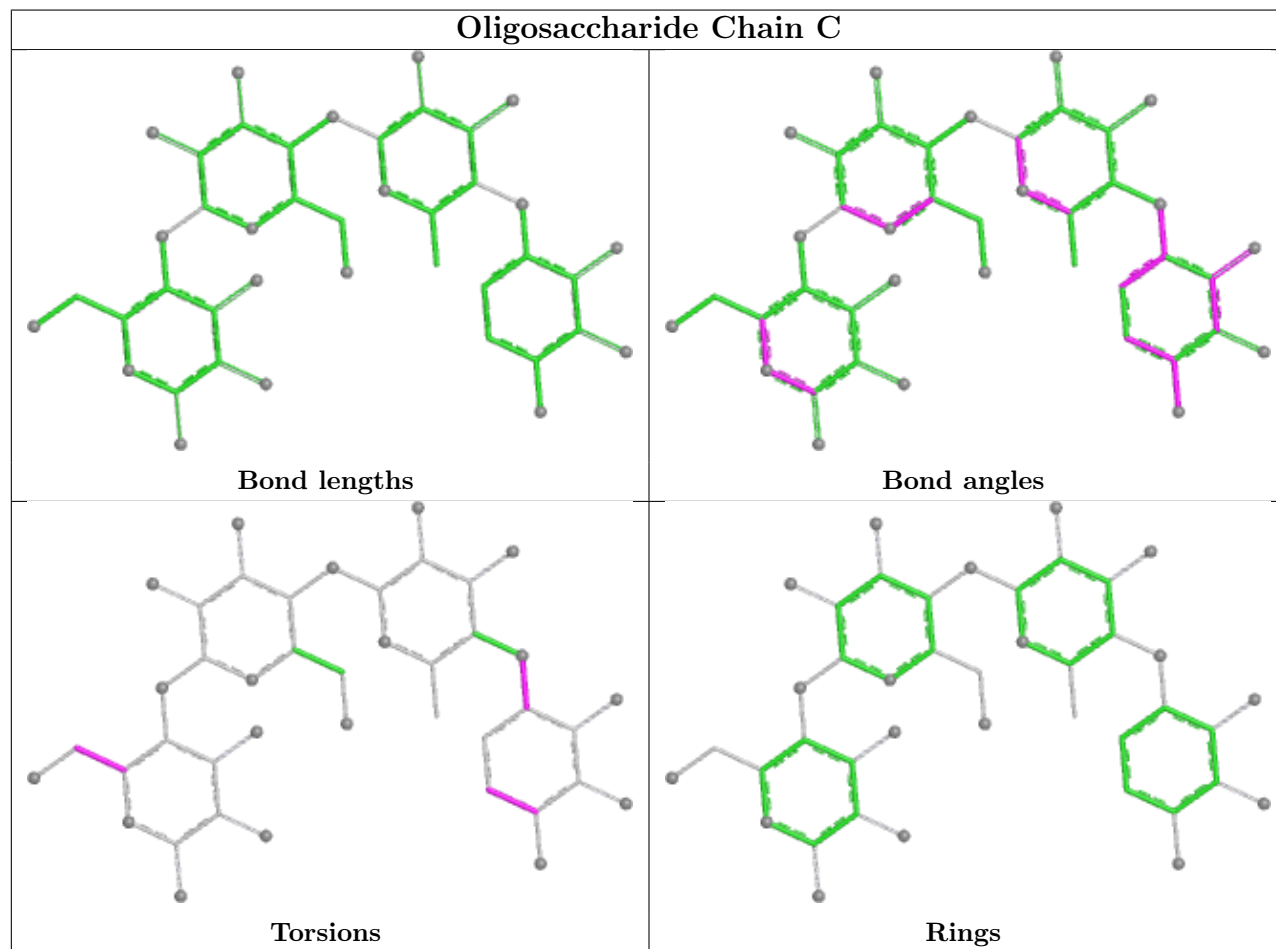
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	K	2	AC1	0	1
2	L	3	AC1	1	0

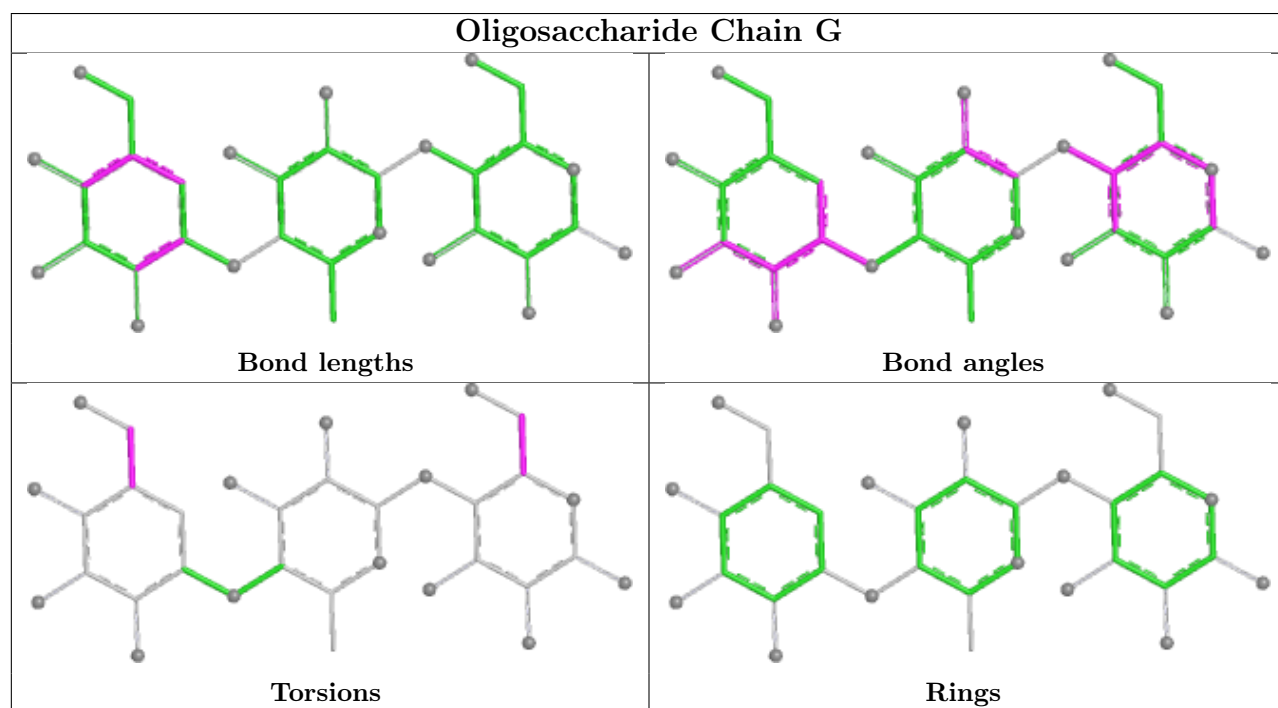
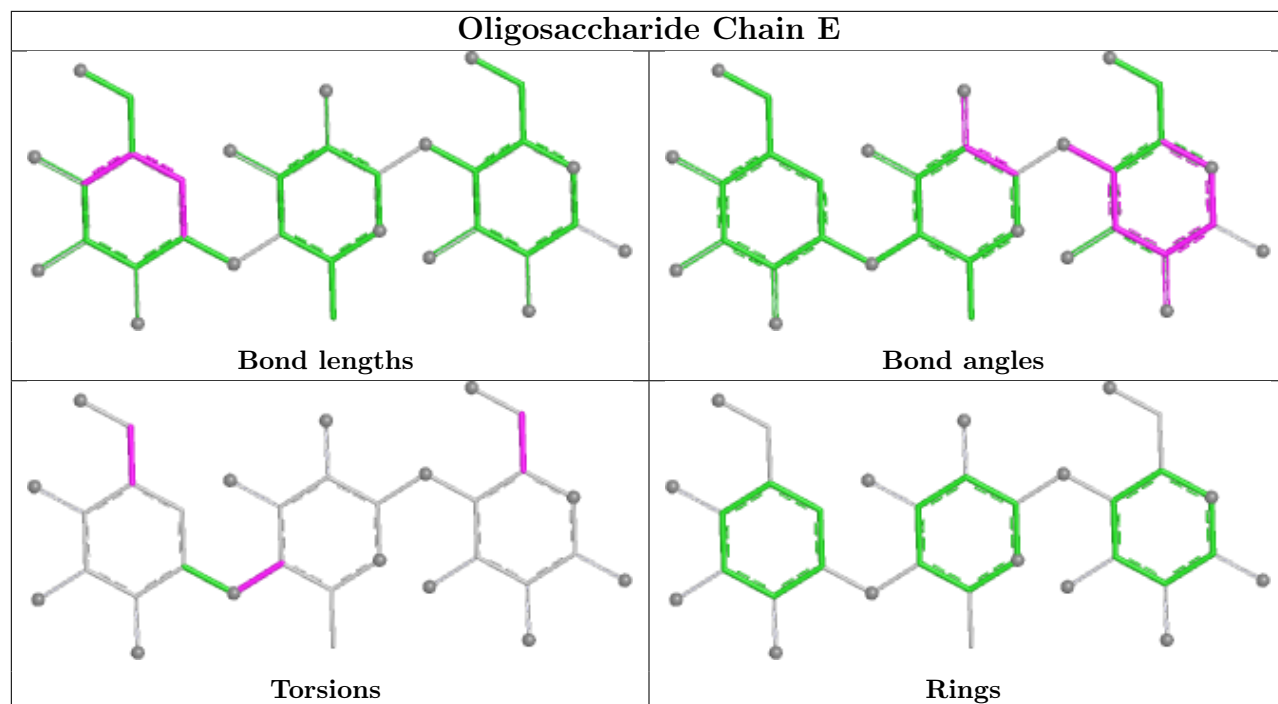
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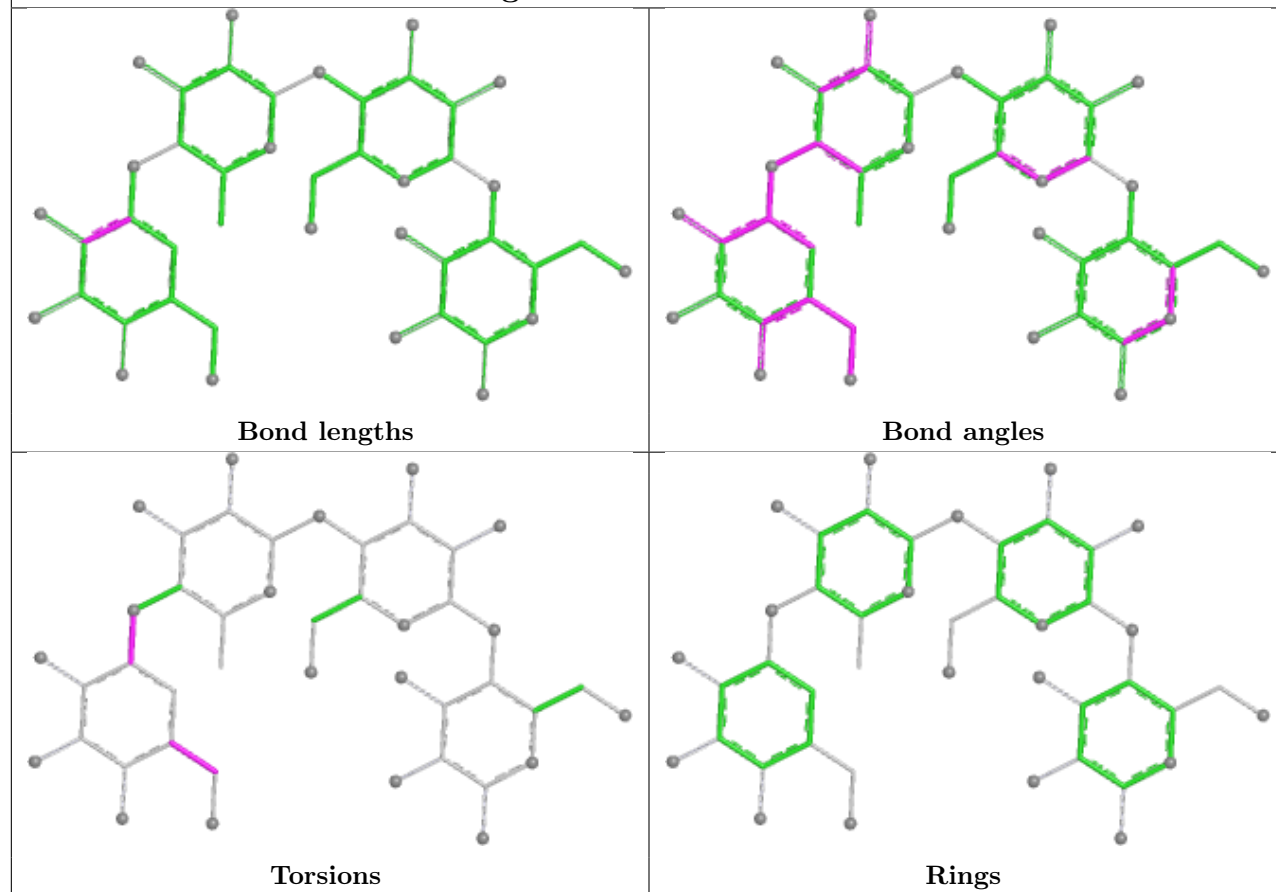
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	2	GLC	1	0
2	H	3	AC1	1	0

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

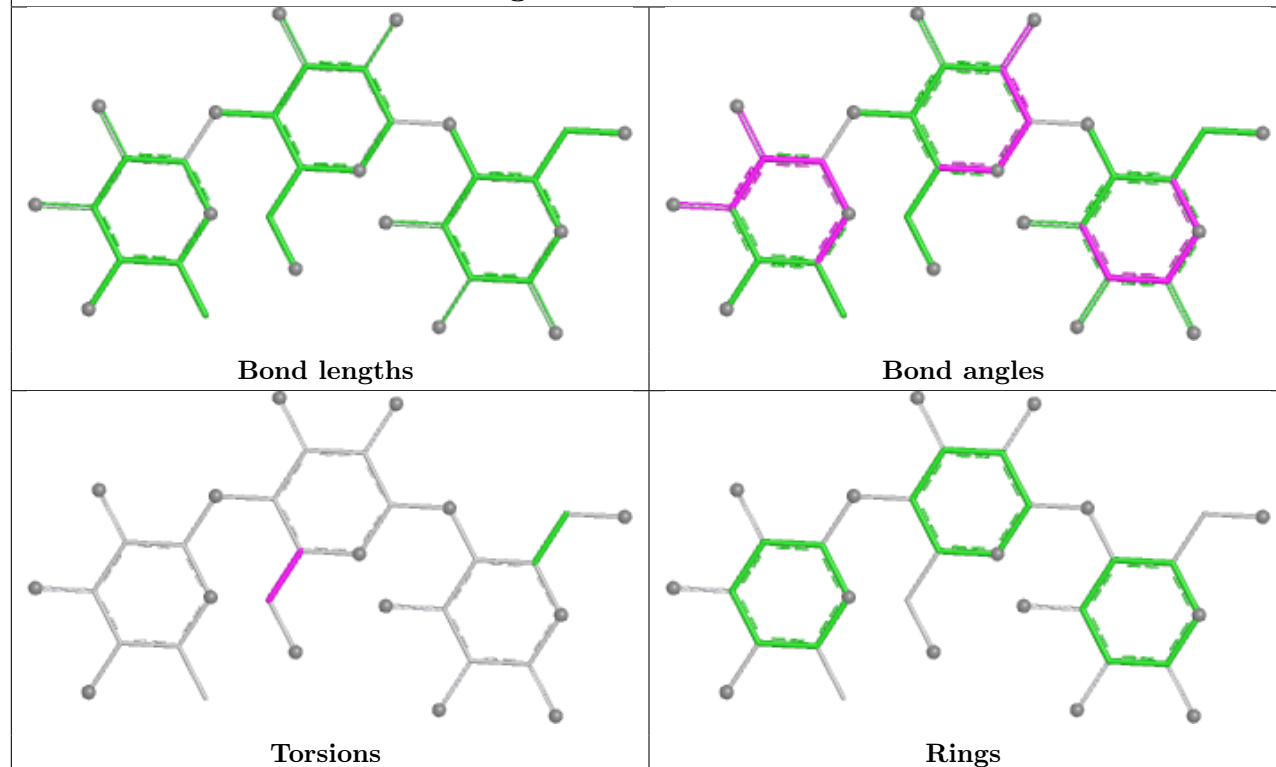




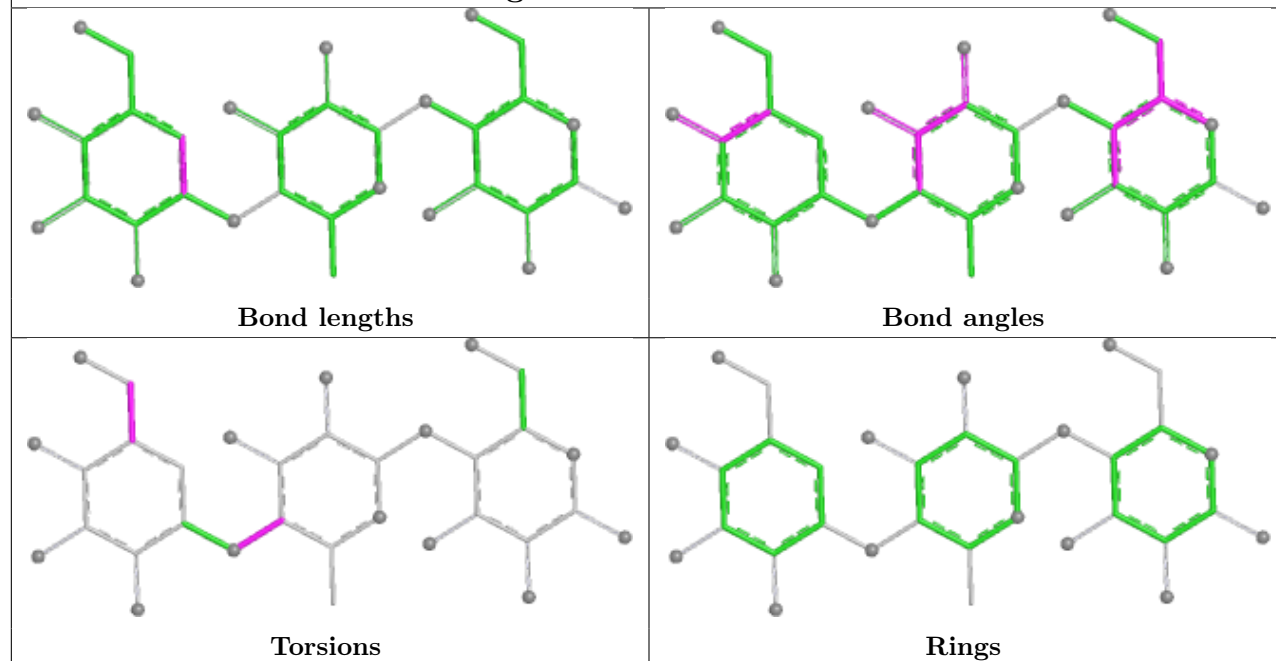
Oligosaccharide Chain H



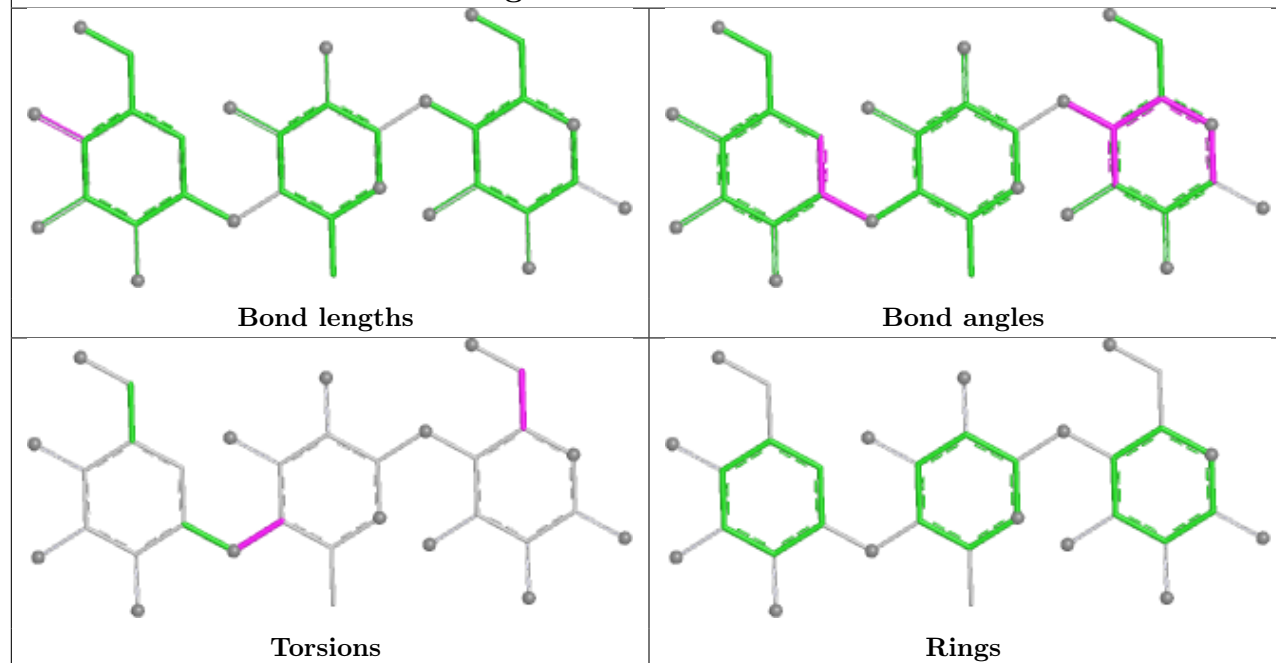
Oligosaccharide Chain I

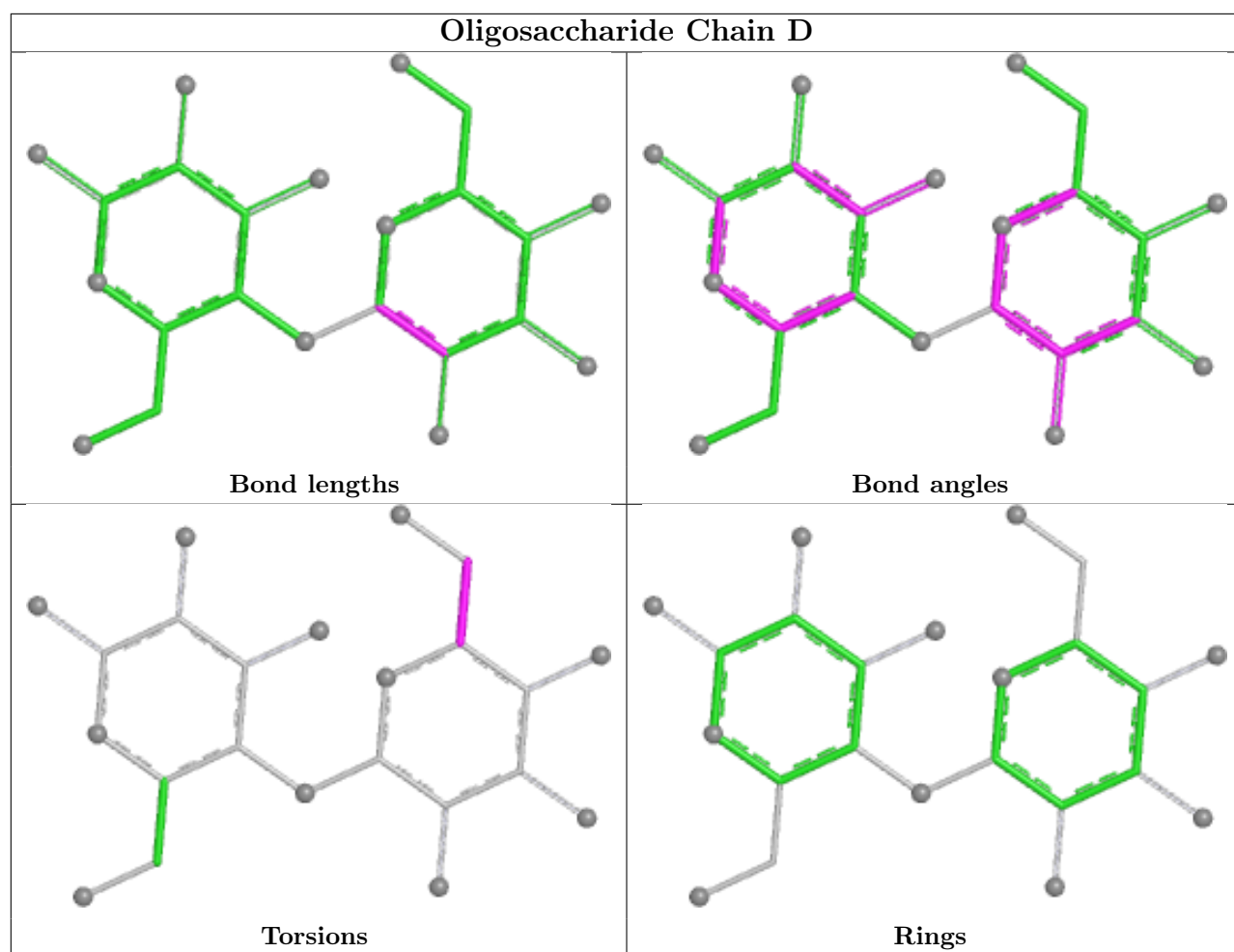


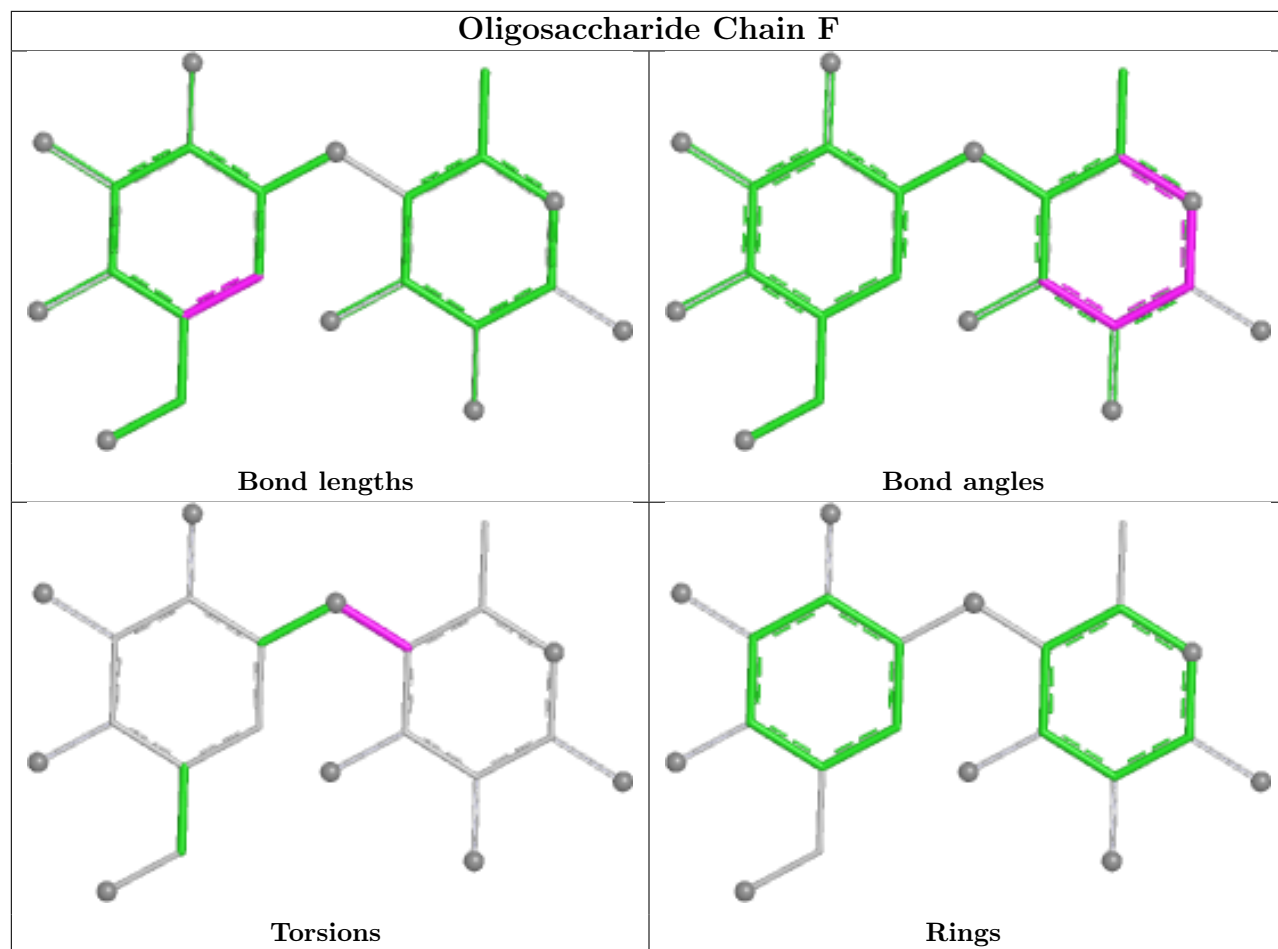
Oligosaccharide Chain J

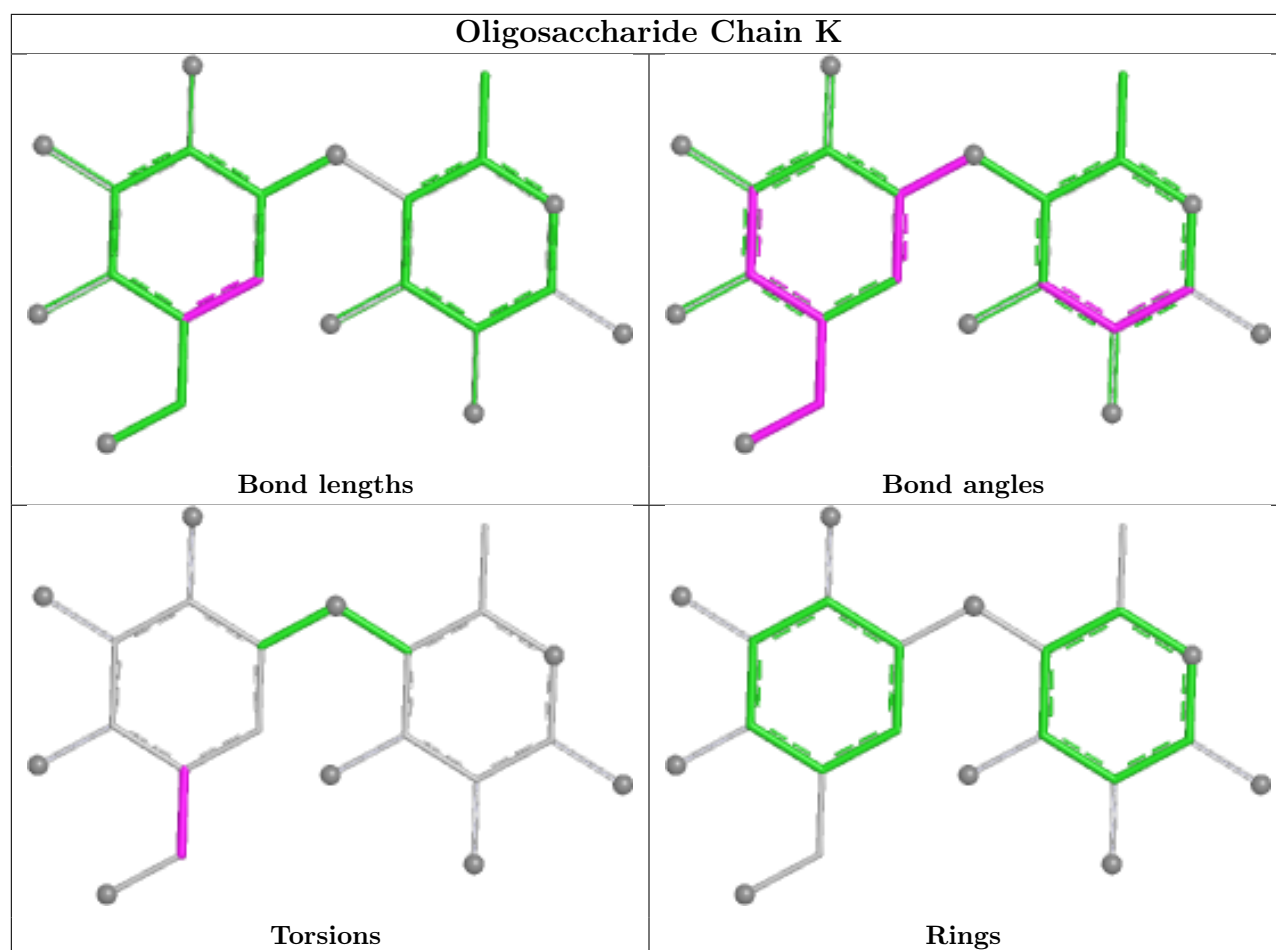


Oligosaccharide Chain L









5.6 Ligand geometry [i](#)

Of 47 ligands modelled in this entry, 20 are monoatomic - leaving 27 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	MES	B	1708	-	12,12,12	2.21	2 (16%)	15,16,16	1.06	0
10	EDO	A	1801	-	3,3,3	0.59	0	2,2,2	0.44	0
9	MES	B	1709	-	12,12,12	2.26	1 (8%)	15,16,16	1.48	2 (13%)
10	EDO	B	1802	-	3,3,3	0.40	0	2,2,2	0.73	0
8	SO4	B	1702	-	4,4,4	0.57	0	6,6,6	0.52	0
10	EDO	A	1803	-	3,3,3	0.37	0	2,2,2	0.81	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	SO4	A	1706	-	4,4,4	0.52	0	6,6,6	0.28	0
8	SO4	B	1701	-	4,4,4	0.57	0	6,6,6	0.61	0
8	SO4	B	1710	-	4,4,4	0.59	0	6,6,6	0.72	0
10	EDO	B	1801	-	3,3,3	0.61	0	2,2,2	0.37	0
8	SO4	A	1704	-	4,4,4	0.51	0	6,6,6	0.47	0
9	MES	B	1711	-	12,12,12	2.29	1 (8%)	15,16,16	1.16	2 (13%)
8	SO4	B	1707	-	4,4,4	0.47	0	6,6,6	0.82	0
10	EDO	A	1804	-	3,3,3	0.60	0	2,2,2	0.71	0
9	MES	A	1707	-	12,12,12	2.31	1 (8%)	15,16,16	1.64	4 (26%)
10	EDO	B	1803	-	3,3,3	0.20	0	2,2,2	0.73	0
9	MES	A	1705	-	12,12,12	1.80	1 (8%)	15,16,16	2.26	5 (33%)
10	EDO	B	1804	-	3,3,3	0.43	0	2,2,2	0.38	0
9	MES	B	1705	-	12,12,12	1.94	1 (8%)	15,16,16	2.23	4 (26%)
8	SO4	A	1701	-	4,4,4	0.59	0	6,6,6	0.50	0
8	SO4	A	1703	-	4,4,4	0.61	0	6,6,6	0.64	0
8	SO4	B	1704	-	4,4,4	0.52	0	6,6,6	0.42	0
8	SO4	B	1712	-	4,4,4	0.47	0	6,6,6	0.26	0
10	EDO	A	1802	-	3,3,3	0.34	0	2,2,2	0.27	0
9	MES	B	1703	-	12,12,12	2.16	1 (8%)	15,16,16	2.09	5 (33%)
8	SO4	B	1706	-	4,4,4	0.49	0	6,6,6	0.41	0
8	SO4	A	1702	-	4,4,4	0.42	0	6,6,6	0.39	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
9	MES	B	1705	-	-	0/6/14/14	0/1/1/1
9	MES	B	1708	-	-	4/6/14/14	0/1/1/1
10	EDO	A	1801	-	-	0/1/1/1	-
9	MES	B	1709	-	-	2/6/14/14	0/1/1/1
10	EDO	B	1801	-	-	0/1/1/1	-
10	EDO	B	1803	-	-	1/1/1/1	-
9	MES	A	1705	-	-	3/6/14/14	0/1/1/1
10	EDO	A	1802	-	-	0/1/1/1	-
9	MES	B	1703	-	-	5/6/14/14	0/1/1/1
9	MES	B	1711	-	-	0/6/14/14	0/1/1/1
10	EDO	A	1803	-	-	0/1/1/1	-
10	EDO	B	1802	-	-	0/1/1/1	-
10	EDO	A	1804	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	MES	A	1707	-	-	0/6/14/14	0/1/1/1
10	EDO	B	1804	-	-	1/1/1/1	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	1707	MES	C8-S	-7.65	1.66	1.77
9	B	1709	MES	C8-S	-7.47	1.67	1.77
9	B	1711	MES	C8-S	-7.36	1.67	1.77
9	B	1703	MES	C8-S	-7.10	1.67	1.77
9	B	1708	MES	C8-S	-7.02	1.67	1.77
9	A	1705	MES	C8-S	-5.49	1.69	1.77
9	B	1705	MES	C8-S	-5.45	1.70	1.77
9	B	1708	MES	O1S-S	2.01	1.50	1.45

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	1705	MES	O2S-S-C8	6.24	116.16	106.73
9	B	1705	MES	O1S-S-C8	5.39	114.88	106.73
9	B	1703	MES	C6-O1-C2	4.67	124.99	109.88
9	B	1705	MES	C6-C5-N4	3.69	115.73	110.12
9	B	1705	MES	O3S-S-O1S	-3.64	102.28	111.40
9	A	1705	MES	C6-C5-N4	3.47	115.39	110.12
9	B	1703	MES	O2S-S-C8	3.21	111.58	106.73
9	A	1707	MES	O3S-S-C8	3.20	112.27	106.00
9	B	1703	MES	O2S-S-O1S	-3.16	103.53	113.82
9	B	1703	MES	C6-C5-N4	-3.14	105.34	110.12
9	B	1709	MES	O1S-S-C8	2.97	111.22	106.73
9	A	1707	MES	O1S-S-C8	-2.92	102.31	106.73
9	A	1707	MES	C6-C5-N4	-2.88	105.74	110.12
9	B	1711	MES	O2S-S-C8	2.80	110.96	106.73
9	A	1705	MES	O2S-S-O1S	-2.71	105.00	113.82
9	B	1705	MES	O3S-S-C8	2.59	111.07	106.00
9	A	1705	MES	O1S-S-C8	2.48	110.48	106.73
9	B	1703	MES	O3S-S-C8	2.45	110.80	106.00
9	A	1705	MES	C5-N4-C3	2.19	113.55	108.84
9	A	1707	MES	O2S-S-C8	2.13	109.94	106.73
9	B	1709	MES	O3S-S-C8	2.11	110.13	106.00
9	B	1711	MES	O3S-S-O2S	-2.07	106.22	111.40

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	A	1705	MES	C7-C8-S-O1S
9	A	1705	MES	C7-C8-S-O3S
9	B	1703	MES	C7-C8-S-O1S
9	B	1703	MES	C7-C8-S-O2S
9	B	1703	MES	C7-C8-S-O3S
9	B	1703	MES	C8-C7-N4-C5
9	B	1708	MES	C8-C7-N4-C3
9	B	1708	MES	C8-C7-N4-C5
9	A	1705	MES	C7-C8-S-O2S
9	B	1708	MES	C7-C8-S-O2S
9	B	1708	MES	C7-C8-S-O3S
9	B	1703	MES	C8-C7-N4-C3
9	B	1709	MES	C8-C7-N4-C3
9	B	1709	MES	C8-C7-N4-C5
10	B	1804	EDO	O1-C1-C2-O2
10	B	1803	EDO	O1-C1-C2-O2

There are no ring outliers.

No monomer is involved in short contacts.

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	1247/1263 (98%)	-0.13	8 (0%) 85 83	28, 42, 61, 101	0
1	B	1247/1263 (98%)	-0.17	7 (0%) 85 83	26, 41, 59, 103	1 (0%)
All	All	2494/2526 (98%)	-0.15	15 (0%) 85 83	26, 42, 60, 103	1 (0%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1281	PRO	5.8
1	A	1237	ASN	3.9
1	A	35	MET	3.7
1	A	1281	PRO	3.4
1	B	35	MET	3.4
1	A	438	SER	3.2
1	B	1237	ASN	3.0
1	B	1280	ALA	3.0
1	B	761	ARG	2.9
1	A	1250	ASN	2.6
1	B	438	SER	2.4
1	B	1000[A]	TRP	2.3
1	A	1267	TYR	2.3
1	A	811	GLY	2.2
1	A	1183	ASP	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 ⓘ

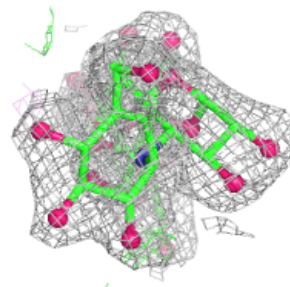
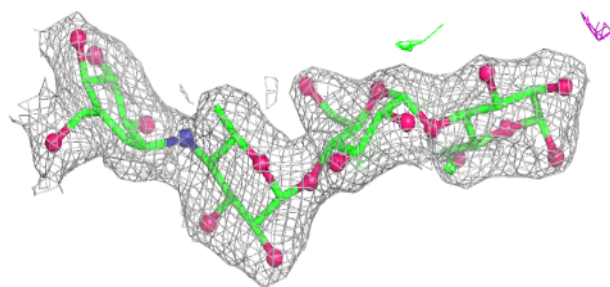
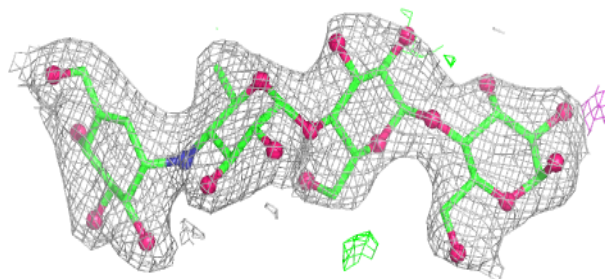
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists 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	GLC	F	1	1/12	0.39	0.19	68,68,68,68	0
4	GLC	K	1	1/12	0.47	0.17	66,66,66,66	0
2	AC1	I	3	10/22	0.53	0.25	88,101,112,116	0
2	GLC	J	2	11/12	0.69	0.17	72,89,98,101	0
2	GLC	L	1	1/12	0.71	0.20	104,104,104,104	0
2	GLC	L	2	11/12	0.73	0.18	71,86,95,98	0
2	GLC	E	2	11/12	0.73	0.15	65,81,88,89	0
2	GLC	J	1	1/12	0.73	0.13	77,77,77,77	0
2	GLC	G	2	11/12	0.77	0.15	69,82,89,93	0
2	GLC	E	1	1/12	0.80	0.12	71,71,71,71	0
3	GLC	D	2	11/12	0.81	0.15	65,79,85,86	0
3	GLC	D	1	12/12	0.82	0.15	72,86,92,92	0
2	GLC	G	1	1/12	0.83	0.20	106,106,106,106	0
2	GLC	I	1	12/12	0.86	0.13	75,83,86,87	0
2	GLC	H	1	12/12	0.86	0.11	61,68,71,72	0
2	GLC	I	2	11/12	0.87	0.14	70,80,87,87	0
2	AC1	G	3	21/22	0.87	0.13	46,62,74,78	0
2	GLC	C	1	12/12	0.89	0.10	52,60,65,66	0
2	GLC	C	2	11/12	0.90	0.11	51,61,70,70	0
2	AC1	E	3	21/22	0.91	0.11	46,57,68,73	0
4	AC1	F	2	21/22	0.91	0.12	49,61,77,82	0
2	AC1	L	3	21/22	0.91	0.13	48,61,73,74	0
4	AC1	K	2	21/22	0.91	0.10	44,58,70,71	0
2	AC1	J	3	21/22	0.92	0.10	40,50,61,63	0
2	GLC	H	2	11/12	0.92	0.09	49,60,67,72	0
2	AC1	H	3	21/22	0.93	0.08	38,43,51,52	0
2	AC1	C	3	21/22	0.94	0.08	40,42,45,46	0

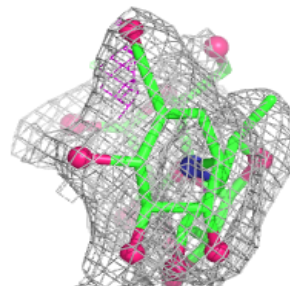
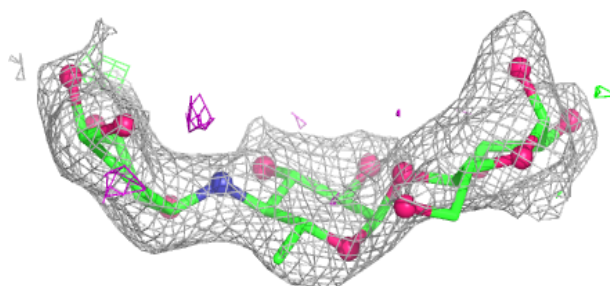
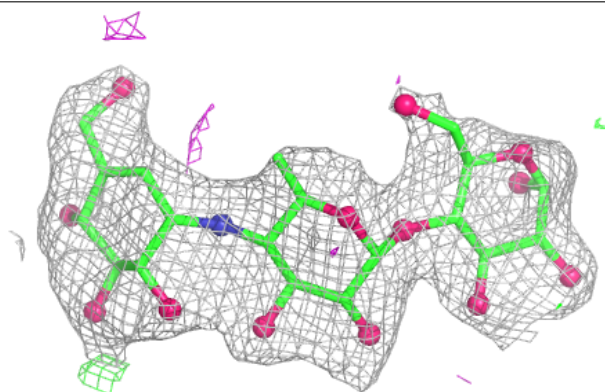
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

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)

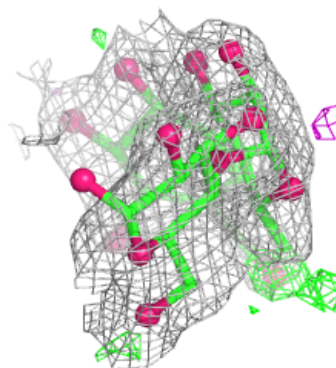
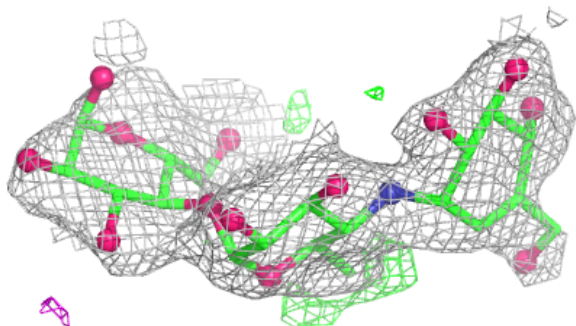
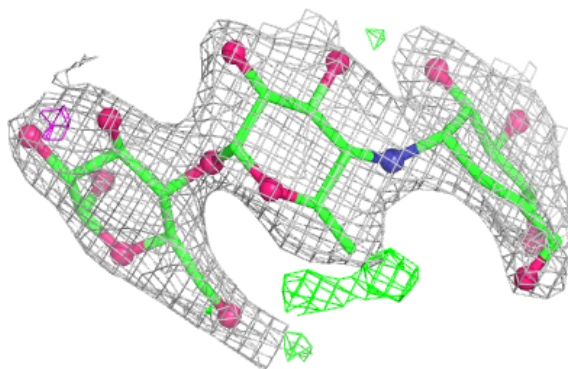
**Electron density around Chain E:**

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and green (positive)

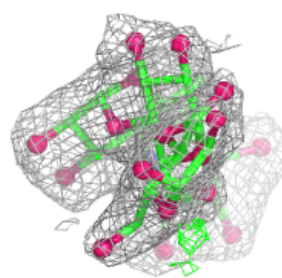
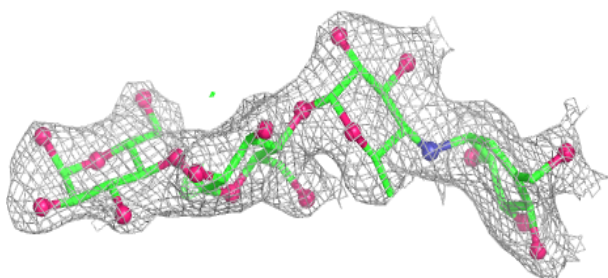
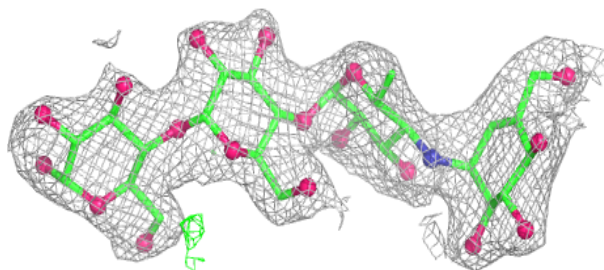


Electron density around Chain G:

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

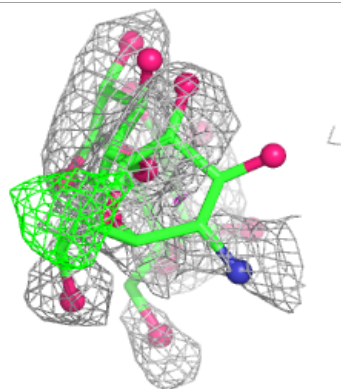
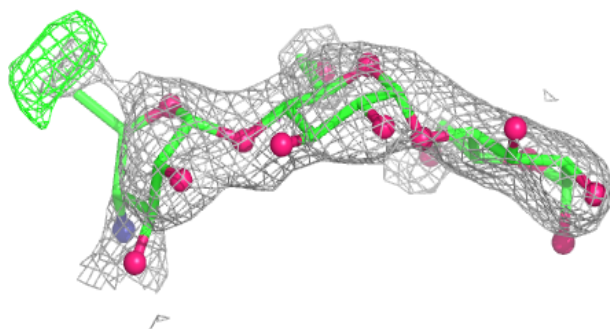
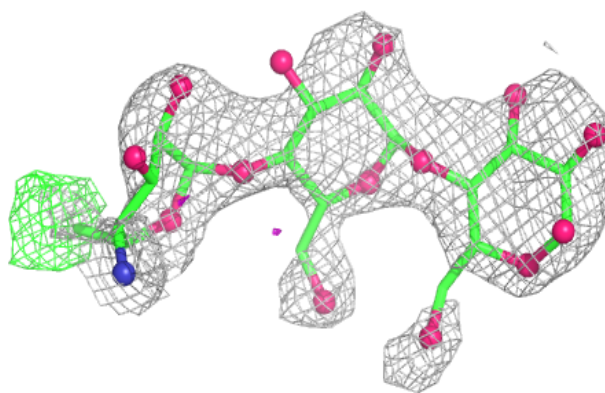
**Electron density around Chain H:**

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and green (positive)

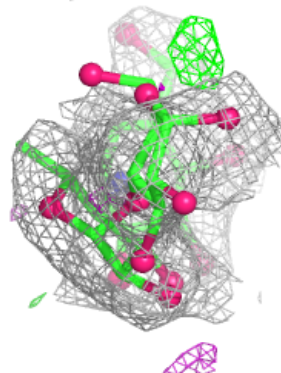
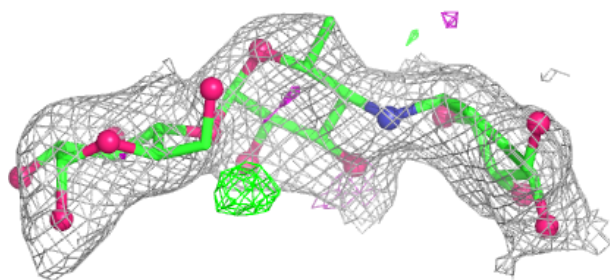
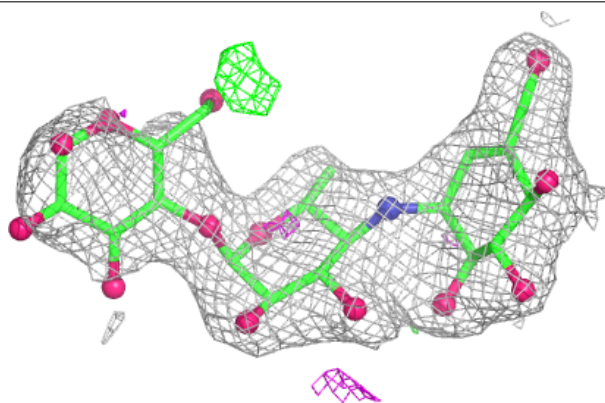


Electron density around Chain I:

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

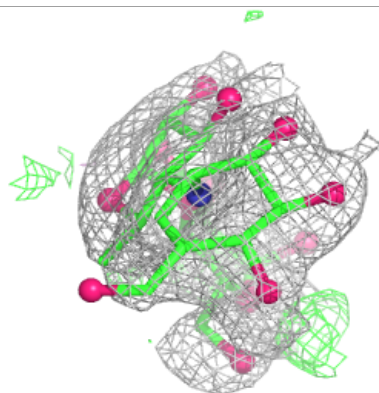
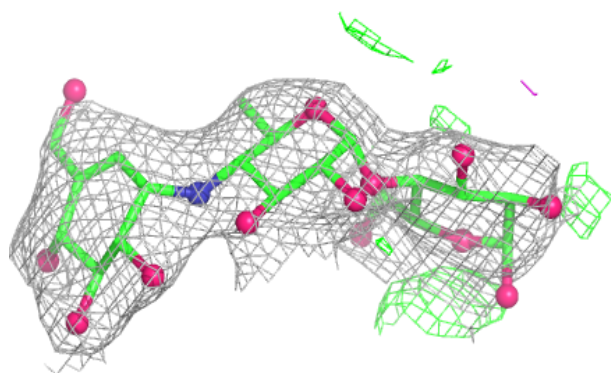
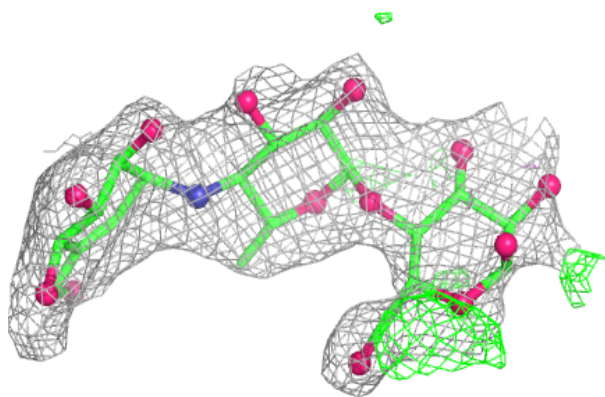
**Electron density around Chain J:**

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and green (positive)



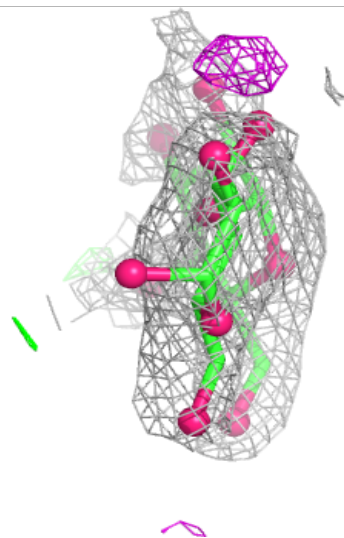
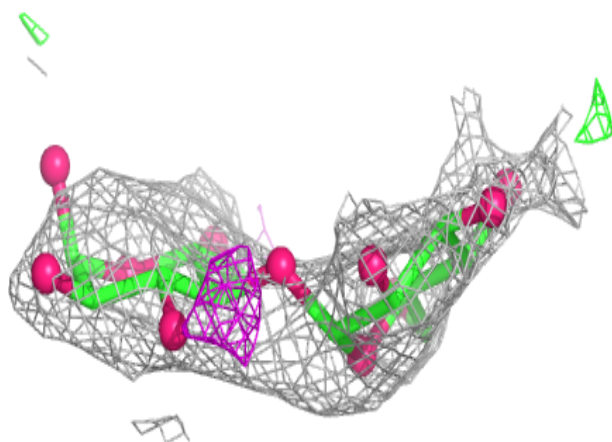
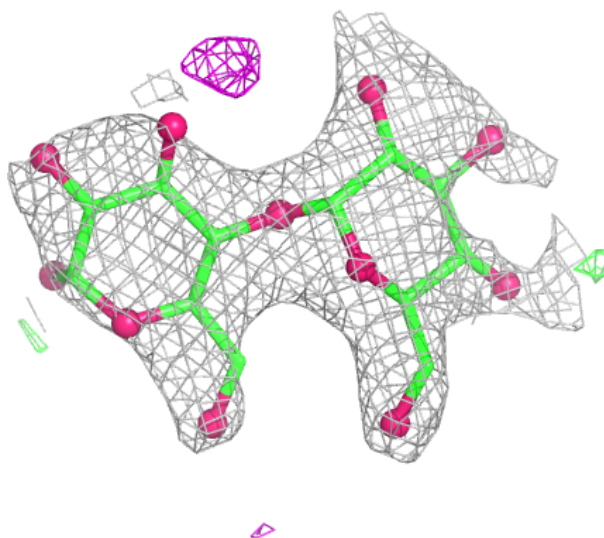
Electron density around Chain L:

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and green (positive)



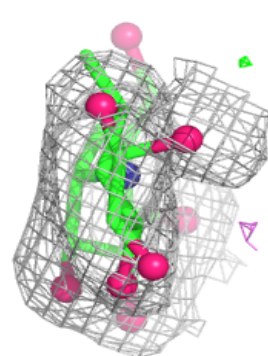
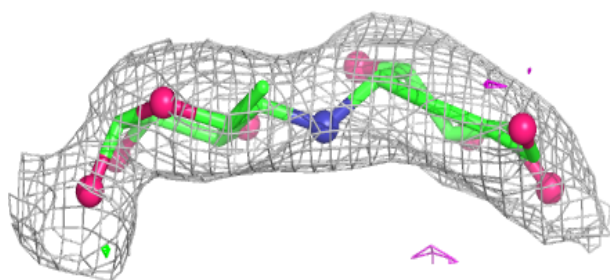
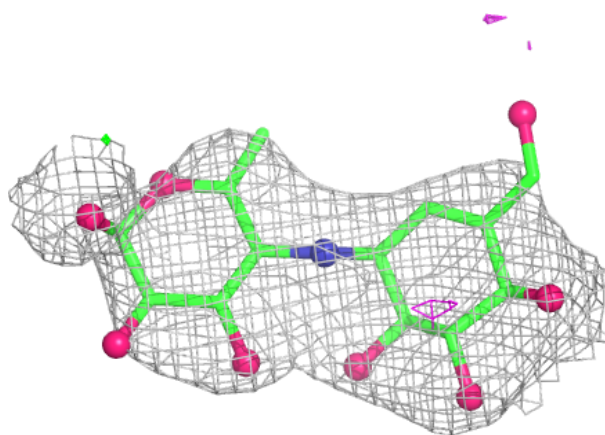
Electron density around Chain D:

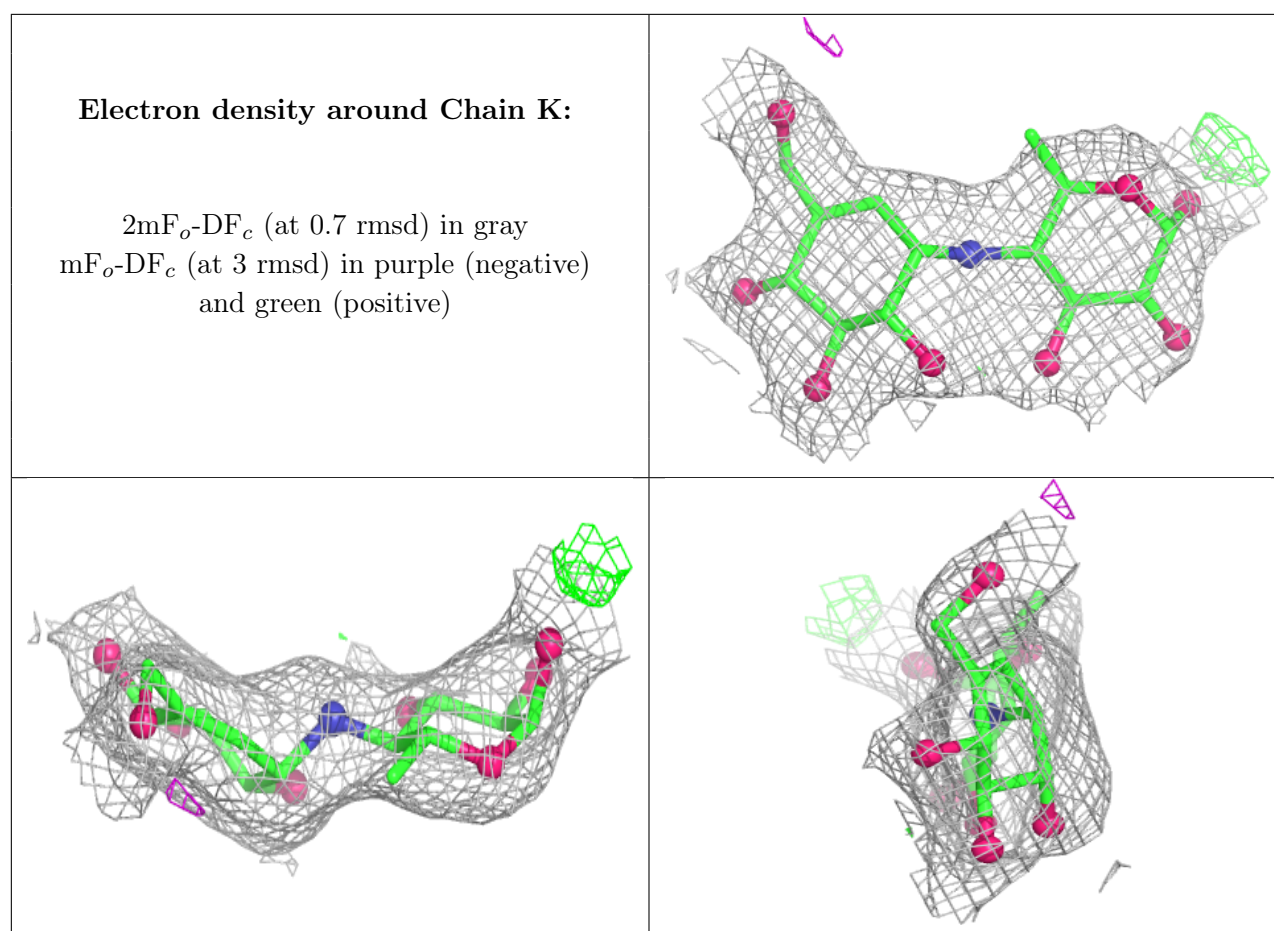
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain F:

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





6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	MG	B	1605	1/1	0.71	0.12	71,71,71,71	0
9	MES	B	1703	12/12	0.77	0.22	71,79,96,96	0
8	SO4	B	1706	5/5	0.83	0.11	75,77,84,87	0
9	MES	B	1708	12/12	0.84	0.23	80,92,97,99	0
9	MES	A	1705	12/12	0.85	0.19	61,83,88,90	0
9	MES	B	1705	12/12	0.87	0.16	64,77,81,82	0
8	SO4	B	1712	5/5	0.87	0.19	73,74,80,83	0
8	SO4	B	1710	5/5	0.88	0.16	72,73,78,87	0
10	EDO	A	1803	4/4	0.88	0.17	51,52,57,59	0
9	MES	B	1709	12/12	0.89	0.19	78,85,89,94	0
8	SO4	B	1704	5/5	0.89	0.11	70,72,80,81	0
10	EDO	B	1803	4/4	0.89	0.16	52,59,61,67	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
10	EDO	B	1802	4/4	0.90	0.14	44,48,49,52	0
8	SO4	A	1703	5/5	0.90	0.12	59,66,83,85	0
8	SO4	A	1702	5/5	0.92	0.14	58,64,68,73	0
8	SO4	A	1701	5/5	0.92	0.09	54,58,69,70	0
10	EDO	A	1804	4/4	0.92	0.12	40,47,47,52	0
8	SO4	A	1704	5/5	0.92	0.17	62,70,79,85	0
8	SO4	B	1702	5/5	0.92	0.13	60,61,69,69	0
8	SO4	B	1701	5/5	0.93	0.10	53,58,67,71	0
10	EDO	A	1802	4/4	0.93	0.12	48,48,48,51	0
10	EDO	B	1804	4/4	0.93	0.11	38,39,41,42	0
8	SO4	B	1707	5/5	0.95	0.11	51,54,58,63	0
10	EDO	A	1801	4/4	0.95	0.10	43,44,45,46	0
7	MG	A	1602	1/1	0.96	0.04	61,61,61,61	0
5	CA	A	1303	1/1	0.96	0.04	50,50,50,50	0
10	EDO	B	1801	4/4	0.96	0.09	38,38,40,41	0
9	MES	B	1711	12/12	0.97	0.07	42,44,48,50	0
8	SO4	A	1706	5/5	0.97	0.09	50,51,55,56	0
9	MES	A	1707	12/12	0.97	0.07	37,46,48,48	0
7	MG	A	1604	1/1	0.98	0.04	62,62,62,62	0
7	MG	B	1603	1/1	0.98	0.04	57,57,57,57	0
5	CA	B	1302	1/1	0.98	0.04	59,59,59,59	0
5	CA	B	1303	1/1	0.98	0.05	44,44,44,44	0
7	MG	A	1601	1/1	0.98	0.03	35,35,35,35	0
5	CA	A	1302	1/1	0.98	0.04	52,52,52,52	0
7	MG	A	1603	1/1	0.99	0.05	61,61,61,61	0
5	CA	A	1301	1/1	0.99	0.02	47,47,47,47	0
7	MG	A	1605	1/1	0.99	0.06	71,71,71,71	0
7	MG	A	1606	1/1	0.99	0.02	49,49,49,49	0
7	MG	B	1602	1/1	0.99	0.05	61,61,61,61	0
6	NI	B	1501	1/1	0.99	0.05	63,63,63,63	0
7	MG	B	1604	1/1	0.99	0.03	51,51,51,51	0
7	MG	B	1606	1/1	1.00	0.05	44,44,44,44	0
5	CA	B	1301	1/1	1.00	0.01	39,39,39,39	0
7	MG	B	1601	1/1	1.00	0.01	33,33,33,33	0
6	NI	A	1501	1/1	1.00	0.05	63,63,63,63	0

6.5 Other polymers

There are no such residues in this entry.