



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 9, 2020 – 02:23 AM BST

PDB ID : 1TG7
Title : Native structure of beta-galactosidase from *Penicillium* sp.
Authors : Rojas, A.L.; Nagem, R.A.P.; Neustroev, K.N.; Arand, M.; Adamska, M.; Eneyskaya, E.V.; Kulminskaya, A.A.; Garratt, R.C.; Golubev, A.M.; Polikarpov, I.
Deposited on : 2004-05-28
Resolution : 1.90 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
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) : 2.13.1

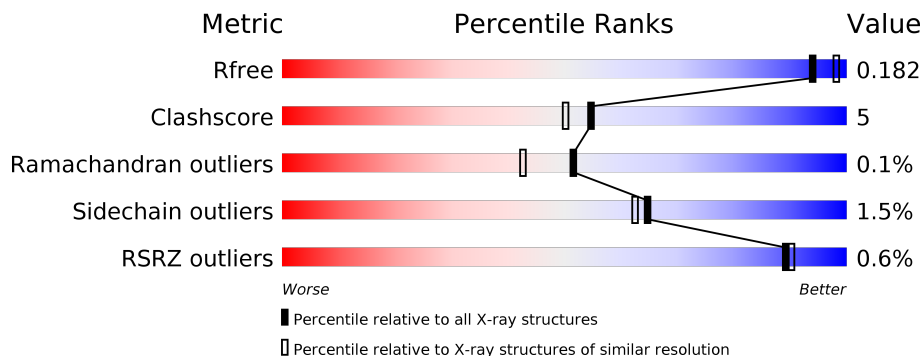
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 on 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	971	 88% 11%
2	B	3	 100%
3	C	9	 78% 22%
4	D	2	 50% 50%
5	E	5	 40% 60%
6	F	7	 100%

2 Entry composition i

There are 11 unique types of molecules in this entry. The entry contains 9154 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 beta-galactosidase.

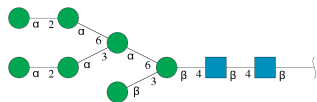
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	971	7499	4790	1241	1458	10	21	3	0

- Molecule 2 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			
2	B	3	39	22	2	15	0	0	0

- Molecule 3 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[beta-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			
3	C	9	105	58	2	45	0	0	0

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



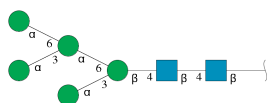
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
4	D	2	28	16	2	10	0	0	0

- Molecule 5 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			
5	E	5	61	34	2	25	0	0	0

- Molecule 6 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[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
			Total	C	N	O			
6	F	7	83	46	2	35	0	0	0

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).

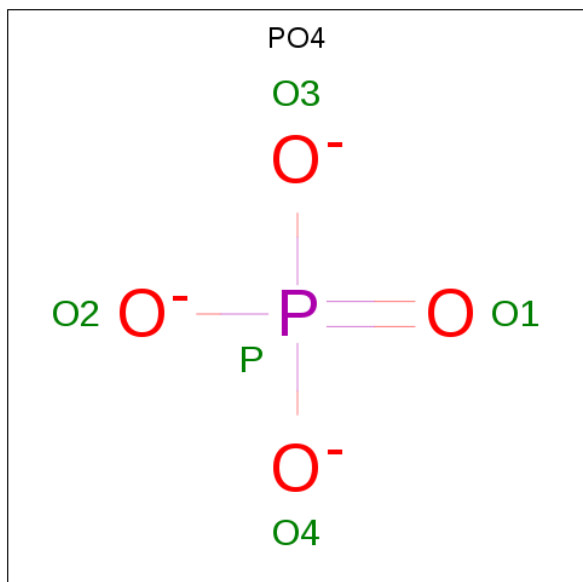


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			14	8	1	5		
7	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 8 is SODIUM ION (three-letter code: NA) (formula: Na).

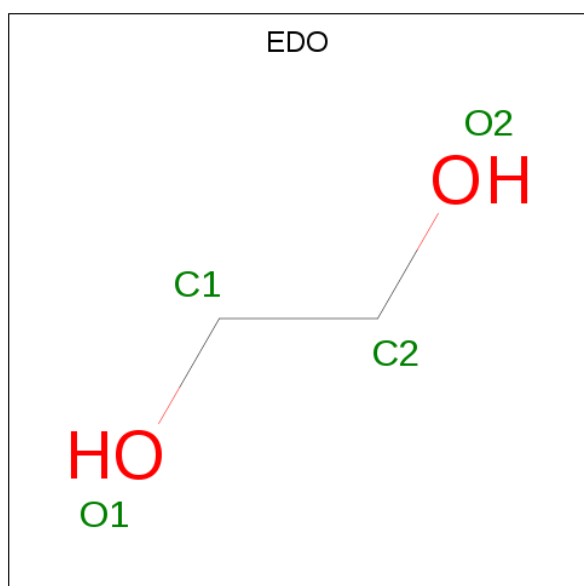
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	3	Total	Na	0	0
			3	3		

- Molecule 9 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	1	Total O P 5 4 1	0	0
9	A	1	Total O P 5 4 1	0	0
9	A	1	Total O P 5 4 1	0	0
9	A	1	Total O P 5 4 1	0	0

- Molecule 10 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

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

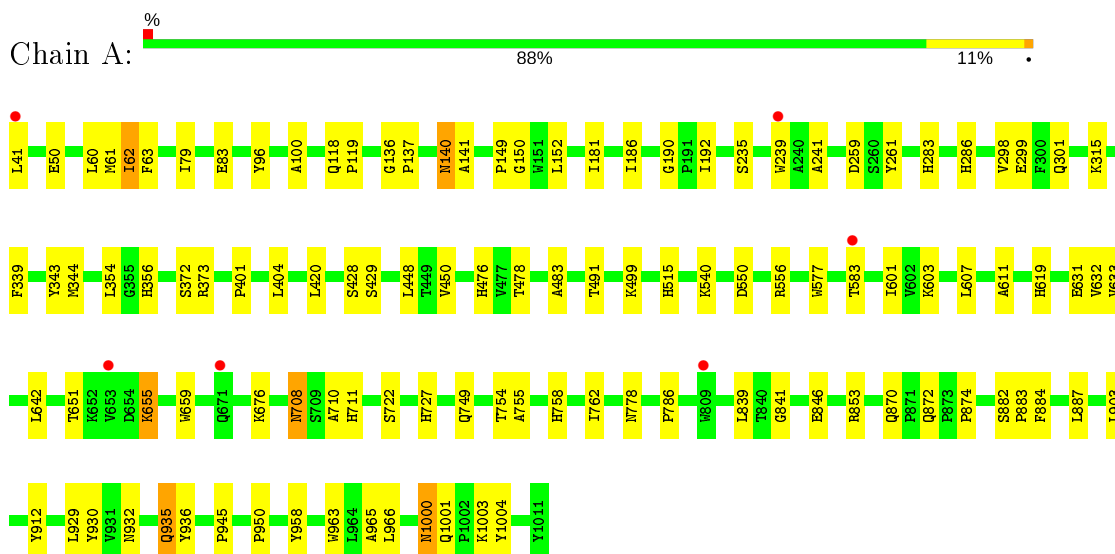
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	1252	Total	O	0	0
			1252	1252		

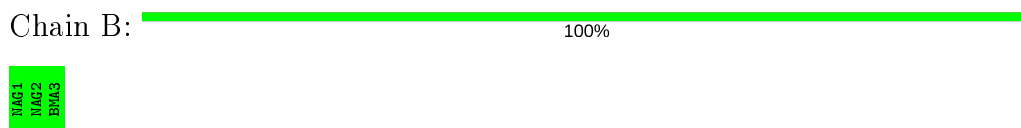
3 Residue-property plots

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.

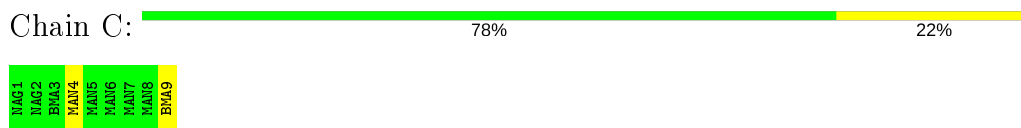
- Molecule 1: beta-galactosidase



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



- Molecule 3: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[beta-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



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

Chain D:  50% 50%

MAG1
MAG2

- Molecule 5: 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:  40% 60%

MAG1
MAG2
BGL3
MAN4
MAN5

- Molecule 6: alpha-D-mannopyranose-(1-3)-[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 F:  100%

MAG1
MAG2
BGL3
MAN4
MAN5
MAN6
MAN7

4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	110.96Å 110.96Å 161.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	22.30 – 1.90 110.96 – 1.90	Depositor EDS
% Data completeness (in resolution range)	(Not available) (22.30-1.90) 99.9 (110.96-1.90)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.03 (at 1.90Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.166 , 0.185 0.160 , 0.182	Depositor DCC
R_{free} test set	1543 reflections (1.01%)	wwPDB-VP
Wilson B-factor (Å ²)	20.7	Xtrriage
Anisotropy	0.333	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 58.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.024 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9154	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.58% 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, NAG, NA, PO4, EDO, MAN

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	0.32	0/7726	0.64	3/10540 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	62[A]	ILE	CA-C-N	-5.49	105.13	117.20
1	A	62[B]	ILE	CA-C-N	-5.49	105.13	117.20
1	A	841	GLY	N-CA-C	-5.26	99.94	113.10

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	7499	0	7143	76	0
2	B	39	0	34	0	0
3	C	105	0	88	0	0
4	D	28	0	25	1	0
5	E	61	0	52	3	0
6	F	83	0	70	0	0
7	A	28	0	26	0	0
8	A	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	A	20	0	0	0	0
10	A	36	0	54	3	0
11	A	1252	0	0	13	0
All	All	9154	0	7492	79	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 5.

All (79) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:778:ASN:HD21	5:E:1:NAG:HN2	1.30	0.78
1:A:853:ARG:HH12	1:A:870:GLN:HE21	1.38	0.71
1:A:259:ASP:CG	1:A:299:GLU:HB2	2.14	0.68
1:A:708:ASN:HD22	1:A:710:ALA:H	1.46	0.62
1:A:932:ASN:ND2	1:A:963:TRP:H	1.96	0.61
1:A:778:ASN:ND2	5:E:1:NAG:HN2	2.00	0.60
1:A:239:TRP:NE1	1:A:241:ALA:HB2	2.15	0.60
1:A:41:LEU:HD11	11:A:8281:HOH:O	2.02	0.58
1:A:428:SER:HB2	11:A:9942:HOH:O	2.05	0.55
1:A:882:SER:HB2	1:A:883:PRO:HD2	1.89	0.55
1:A:722:SER:OG	1:A:727:HIS:HD2	1.89	0.55
1:A:140:ASN:HB3	10:A:9229:EDO:H22	1.89	0.55
1:A:96:TYR:CE2	1:A:141:ALA:HB2	2.42	0.54
1:A:100:ALA:HA	1:A:149:PRO:HG3	1.90	0.54
1:A:286:HIS:HE1	11:A:9371:HOH:O	1.91	0.54
1:A:711:HIS:HE1	11:A:9444:HOH:O	1.90	0.53
1:A:286:HIS:HD2	11:A:9425:HOH:O	1.92	0.53
1:A:476:HIS:HD2	1:A:491:THR:O	1.93	0.52
1:A:261:TYR:HA	1:A:301:GLN:HB2	1.91	0.52
1:A:930[B]:TYR:HB2	1:A:965:ALA:HB3	1.91	0.52
1:A:540:LYS:HE3	11:A:8300:HOH:O	2.09	0.51
1:A:283:HIS:HE1	11:A:8255:HOH:O	1.93	0.51
1:A:749:GLN:HE21	1:A:778:ASN:HD22	1.58	0.50
1:A:283:HIS:HD2	11:A:9466:HOH:O	1.93	0.50
1:A:499:LYS:HE3	11:A:9983:HOH:O	2.12	0.49
11:A:8357:HOH:O	4:D:1:NAG:H83	2.13	0.49
5:E:2:NAG:H83	5:E:4:MAN:O4	2.13	0.48
1:A:603:LYS:HB3	1:A:631:GLU:HB2	1.95	0.48
1:A:401:PRO:HA	1:A:420:LEU:HD23	1.95	0.48
1:A:708:ASN:ND2	1:A:710:ALA:H	2.11	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:762:ILE:O	1:A:762:ILE:HG23	2.13	0.48
1:A:749:GLN:HE21	1:A:778:ASN:ND2	2.12	0.48
1:A:874:PRO:HD3	1:A:1004:TYR:CD2	2.48	0.48
1:A:932:ASN:HD21	1:A:963:TRP:H	1.62	0.47
1:A:515:HIS:HE1	11:A:9463:HOH:O	1.96	0.47
10:A:9228:EDO:O1	10:A:9229:EDO:H21	2.14	0.47
1:A:515:HIS:HD2	1:A:550:ASP:OD1	1.97	0.47
1:A:1000:ASN:HD22	1:A:1001:GLN:N	2.12	0.46
1:A:118:GLN:HB3	1:A:119:PRO:HD3	1.97	0.46
1:A:483:ALA:HB1	1:A:499:LYS:HD2	1.97	0.46
1:A:181:ILE:HG23	1:A:192:ILE:HD13	1.97	0.46
1:A:611:ALA:HA	1:A:619:HIS:O	2.15	0.46
1:A:676:LYS:HA	1:A:884:PHE:CE2	2.50	0.46
1:A:136:GLY:HA3	1:A:137:PRO:C	2.36	0.46
1:A:930[A]:TYR:CE1	1:A:935:GLN:HB2	2.50	0.46
1:A:60:LEU:HD23	1:A:62[B]:ILE:HD11	1.98	0.46
1:A:577:TRP:CZ2	1:A:607:LEU:HA	2.52	0.45
1:A:601:ILE:HB	1:A:633:VAL:HB	1.99	0.45
1:A:727:HIS:HE1	1:A:846:GLU:OE2	2.00	0.44
1:A:149:PRO:HD2	1:A:152:LEU:HD23	2.00	0.43
1:A:708:ASN:C	1:A:708:ASN:HD22	2.22	0.43
1:A:887:LEU:N	1:A:887:LEU:HD12	2.33	0.43
1:A:655:LYS:C	1:A:655:LYS:HD3	2.38	0.43
1:A:903:LEU:HD12	1:A:958:TYR:CD2	2.54	0.43
1:A:872:GLN:HE22	1:A:1003:LYS:NZ	2.17	0.43
1:A:140:ASN:HB3	1:A:141:ALA:H	1.73	0.43
1:A:754:THR:O	1:A:755:ALA:HB3	2.19	0.42
1:A:186:ILE:HA	1:A:190:GLY:O	2.19	0.42
1:A:912:TYR:CE2	1:A:950:PRO:HB3	2.54	0.42
1:A:632:VAL:HB	1:A:659:TRP:HB2	2.02	0.42
1:A:315:LYS:HE2	11:A:8018:HOH:O	2.20	0.41
1:A:429:SER:HB2	1:A:478:THR:OG1	2.19	0.41
1:A:642:LEU:HB3	1:A:651:THR:HG21	2.03	0.41
1:A:404:LEU:HD12	1:A:404:LEU:C	2.41	0.41
1:A:50:GLU:HG2	11:A:8348:HOH:O	2.19	0.41
1:A:79:ILE:O	1:A:83:GLU:HG3	2.21	0.41
1:A:929:LEU:HA	1:A:966:LEU:HD23	2.03	0.41
1:A:448:LEU:HG	1:A:450:VAL:HG13	2.03	0.41
1:A:727:HIS:CE1	1:A:846:GLU:OE2	2.74	0.41
1:A:150:GLY:HA3	1:A:727:HIS:CG	2.56	0.41
1:A:758:HIS:C	1:A:758:HIS:CD2	2.94	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:354:LEU:C	1:A:354:LEU:HD12	2.42	0.41
1:A:372:SER:O	1:A:373:ARG:HB2	2.21	0.41
1:A:60:LEU:HD23	1:A:62[B]:ILE:CD1	2.51	0.41
1:A:235:SER:O	1:A:298:VAL:HG21	2.20	0.40
1:A:343:TYR:HA	1:A:344:MET:HA	1.88	0.40
1:A:786:PRO:HG2	10:A:9226:EDO:H12	2.02	0.40
1:A:96:TYR:CD2	1:A:141:ALA:HB2	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	972/971 (100%)	939 (97%)	32 (3%)	1 (0%)	51 43

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	140	ASN

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	798/795 (100%)	786 (98%)	12 (2%)	65 62

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	61	MET
1	A	63	PHE
1	A	356	HIS
1	A	556	ARG
1	A	583	THR
1	A	655	LYS
1	A	708	ASN
1	A	839	LEU
1	A	935	GLN
1	A	936	TYR
1	A	945	PRO
1	A	1000	ASN

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

Mol	Chain	Res	Type
1	A	140	ASN
1	A	223	HIS
1	A	283	HIS
1	A	286	HIS
1	A	322	HIS
1	A	476	HIS
1	A	515	HIS
1	A	650	GLN
1	A	708	ASN
1	A	711	HIS
1	A	727	HIS
1	A	778	ASN
1	A	824	GLN
1	A	870	GLN
1	A	872	GLN
1	A	932	ASN
1	A	985	ASN
1	A	1000	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](#)

26 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	B	1	1,2	14,14,15	0.56	0	17,19,21	0.77	0
2	NAG	B	2	2	14,14,15	0.57	0	17,19,21	0.56	0
2	BMA	B	3	2	11,11,12	0.42	0	15,15,17	0.33	0
3	NAG	C	1	1,3	14,14,15	0.49	0	17,19,21	0.63	0
3	NAG	C	2	3	14,14,15	0.54	0	17,19,21	0.61	0
3	BMA	C	3	3	11,11,12	0.51	0	15,15,17	0.38	0
3	MAN	C	4	3	11,11,12	0.43	0	15,15,17	0.69	1 (6%)
3	MAN	C	5	3	11,11,12	0.45	0	15,15,17	0.46	0
3	MAN	C	6	3	11,11,12	0.44	0	15,15,17	0.60	0
3	MAN	C	7	3	11,11,12	0.52	0	15,15,17	0.65	0
3	MAN	C	8	3	11,11,12	0.49	0	15,15,17	0.60	0
3	BMA	C	9	3	11,11,12	0.58	0	15,15,17	0.83	1 (6%)
4	NAG	D	1	1,4	14,14,15	0.56	0	17,19,21	0.66	0
4	NAG	D	2	4	14,14,15	0.49	0	17,19,21	0.66	0
5	NAG	E	1	1,5	14,14,15	0.48	0	17,19,21	0.65	0
5	NAG	E	2	5	14,14,15	0.50	0	17,19,21	0.64	0
5	BMA	E	3	5	11,11,12	0.47	0	15,15,17	0.31	0
5	MAN	E	4	5	11,11,12	0.47	0	15,15,17	0.59	0
5	MAN	E	5	5	11,11,12	0.47	0	15,15,17	0.56	0
6	NAG	F	1	1,8,6	14,14,15	0.49	0	17,19,21	0.65	0
6	NAG	F	2	6	14,14,15	0.51	0	17,19,21	0.67	0
6	BMA	F	3	6	11,11,12	0.40	0	15,15,17	0.35	0
6	MAN	F	4	6	11,11,12	0.51	0	15,15,17	0.65	0
6	MAN	F	5	6	11,11,12	0.50	0	15,15,17	0.47	0
6	MAN	F	6	6	11,11,12	0.52	0	15,15,17	0.50	0
6	MAN	F	7	6	11,11,12	0.47	0	15,15,17	0.48	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	B	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	2	2	-	0/6/23/26	0/1/1/1
2	BMA	B	3	2	-	0/2/19/22	0/1/1/1
3	NAG	C	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	C	2	3	-	0/6/23/26	0/1/1/1
3	BMA	C	3	3	-	0/2/19/22	0/1/1/1
3	MAN	C	4	3	-	0/2/19/22	0/1/1/1
3	MAN	C	5	3	-	0/2/19/22	0/1/1/1
3	MAN	C	6	3	-	0/2/19/22	0/1/1/1
3	MAN	C	7	3	-	0/2/19/22	0/1/1/1
3	MAN	C	8	3	-	0/2/19/22	0/1/1/1
3	BMA	C	9	3	-	2/2/19/22	1/1/1/1
4	NAG	D	1	1,4	-	4/6/23/26	0/1/1/1
4	NAG	D	2	4	-	0/6/23/26	0/1/1/1
5	NAG	E	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	E	2	5	-	0/6/23/26	0/1/1/1
5	BMA	E	3	5	-	0/2/19/22	0/1/1/1
5	MAN	E	4	5	-	2/2/19/22	0/1/1/1
5	MAN	E	5	5	-	0/2/19/22	0/1/1/1
6	NAG	F	1	1,8,6	-	0/6/23/26	0/1/1/1
6	NAG	F	2	6	-	0/6/23/26	0/1/1/1
6	BMA	F	3	6	-	0/2/19/22	0/1/1/1
6	MAN	F	4	6	-	0/2/19/22	0/1/1/1
6	MAN	F	5	6	-	0/2/19/22	0/1/1/1
6	MAN	F	6	6	-	0/2/19/22	0/1/1/1
6	MAN	F	7	6	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	9	BMA	C1-O5-C5	2.69	115.83	112.19
3	C	4	MAN	C1-O5-C5	2.19	115.16	112.19

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	9	BMA	O5-C5-C6-O6
3	C	9	BMA	C4-C5-C6-O6
4	D	1	NAG	C4-C5-C6-O6
4	D	1	NAG	O5-C5-C6-O6
4	D	1	NAG	C8-C7-N2-C2
5	E	4	MAN	C4-C5-C6-O6
4	D	1	NAG	O7-C7-N2-C2
5	E	4	MAN	O5-C5-C6-O6

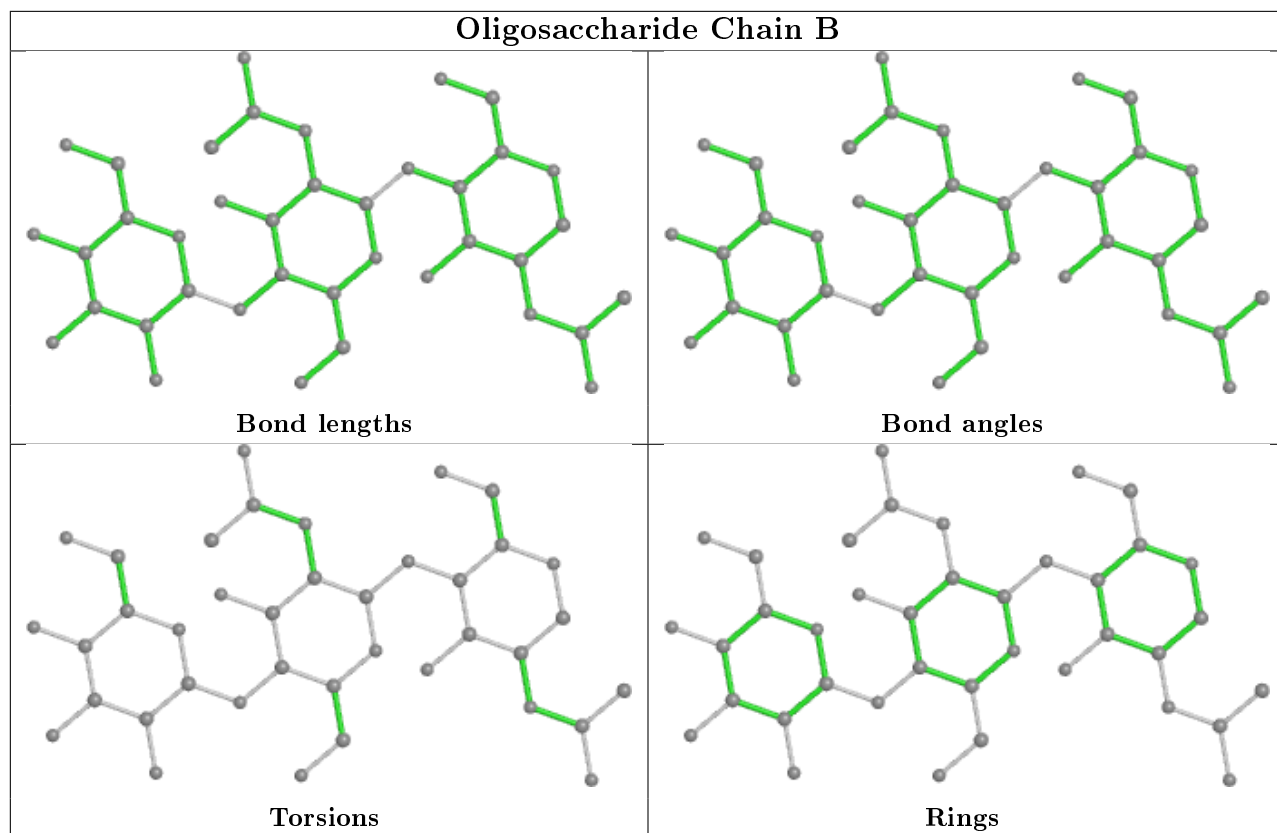
All (1) ring outliers are listed below:

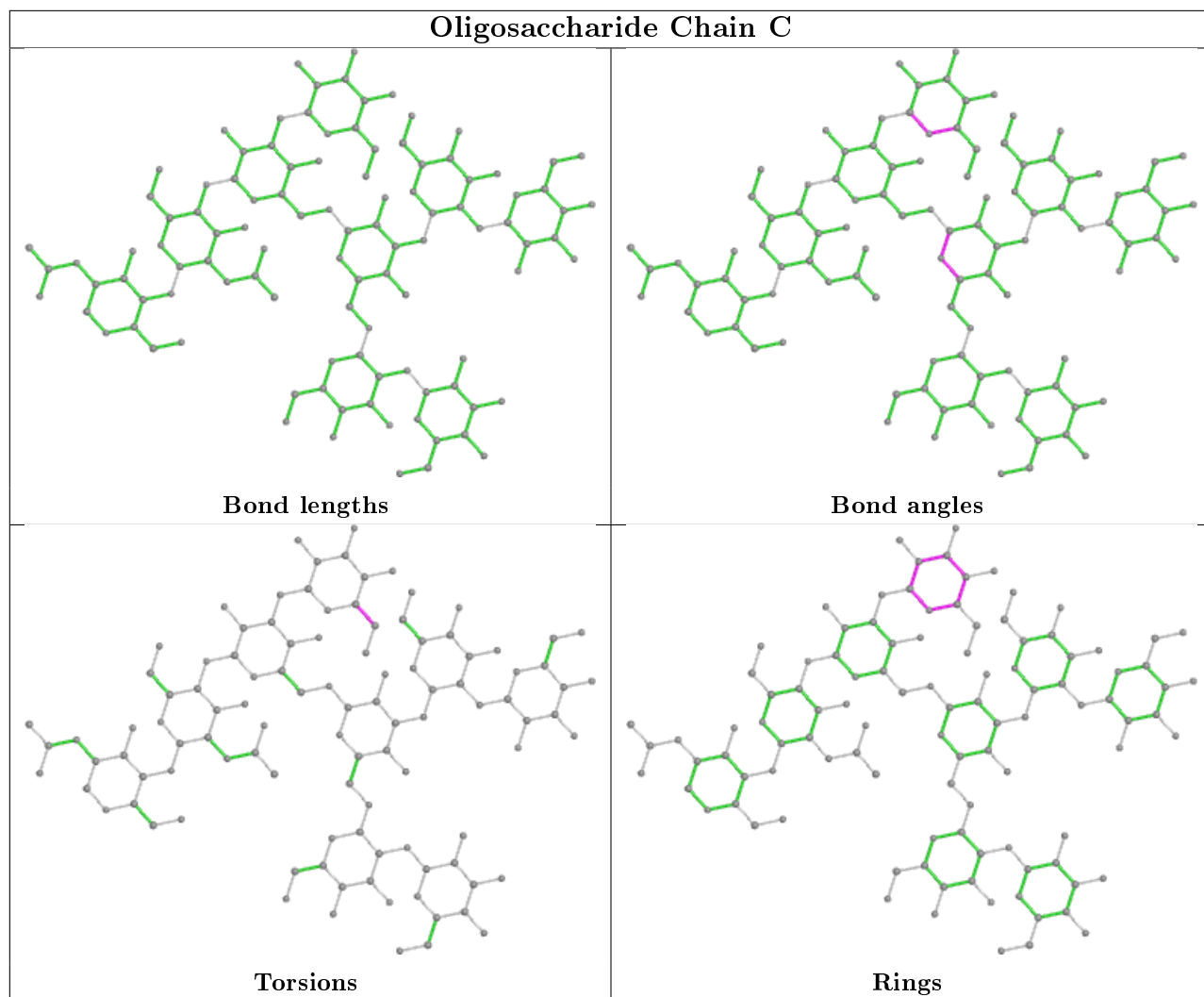
Mol	Chain	Res	Type	Atoms
3	C	9	BMA	C1-C2-C3-C4-C5-O5

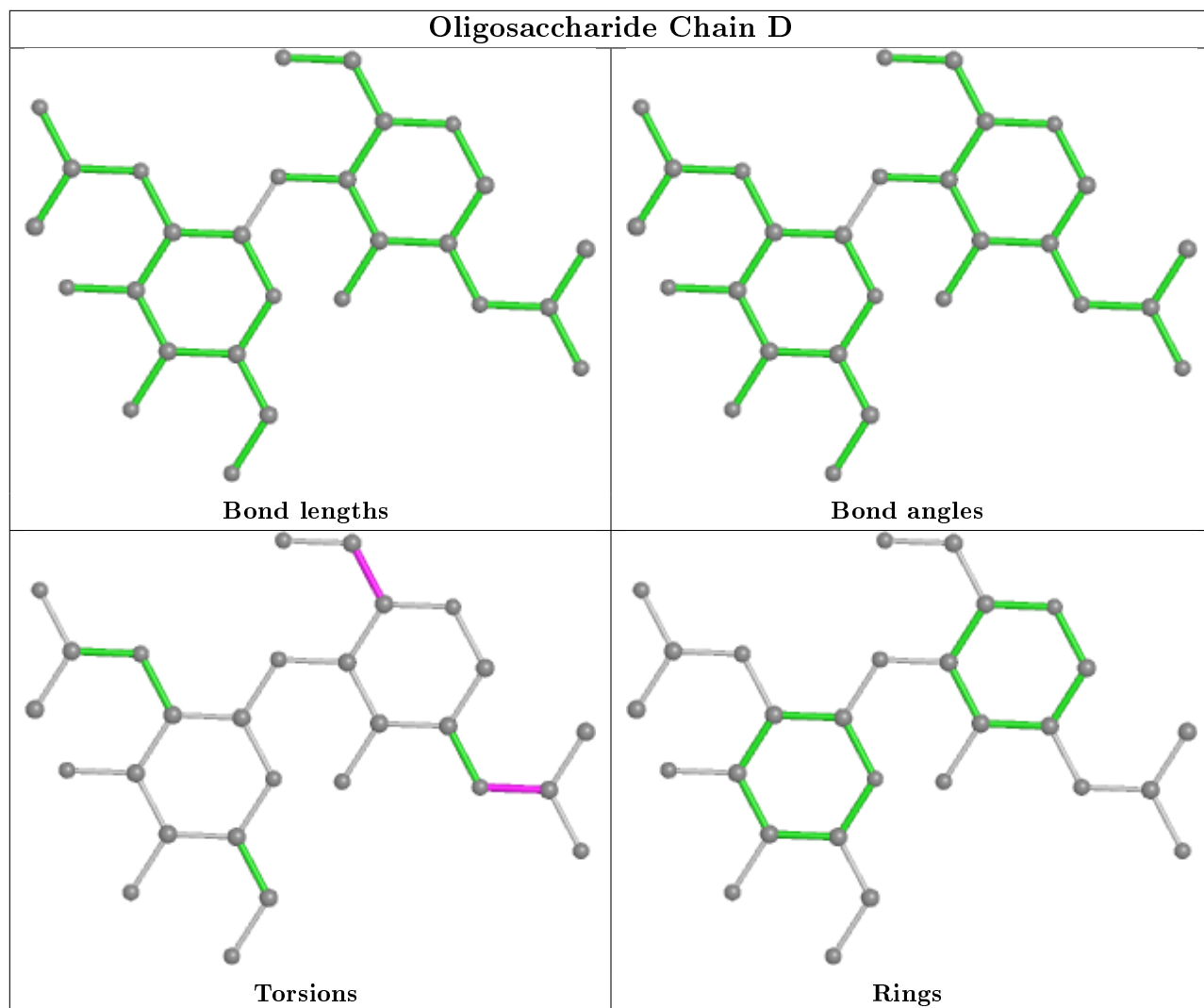
4 monomers are involved in 4 short contacts:

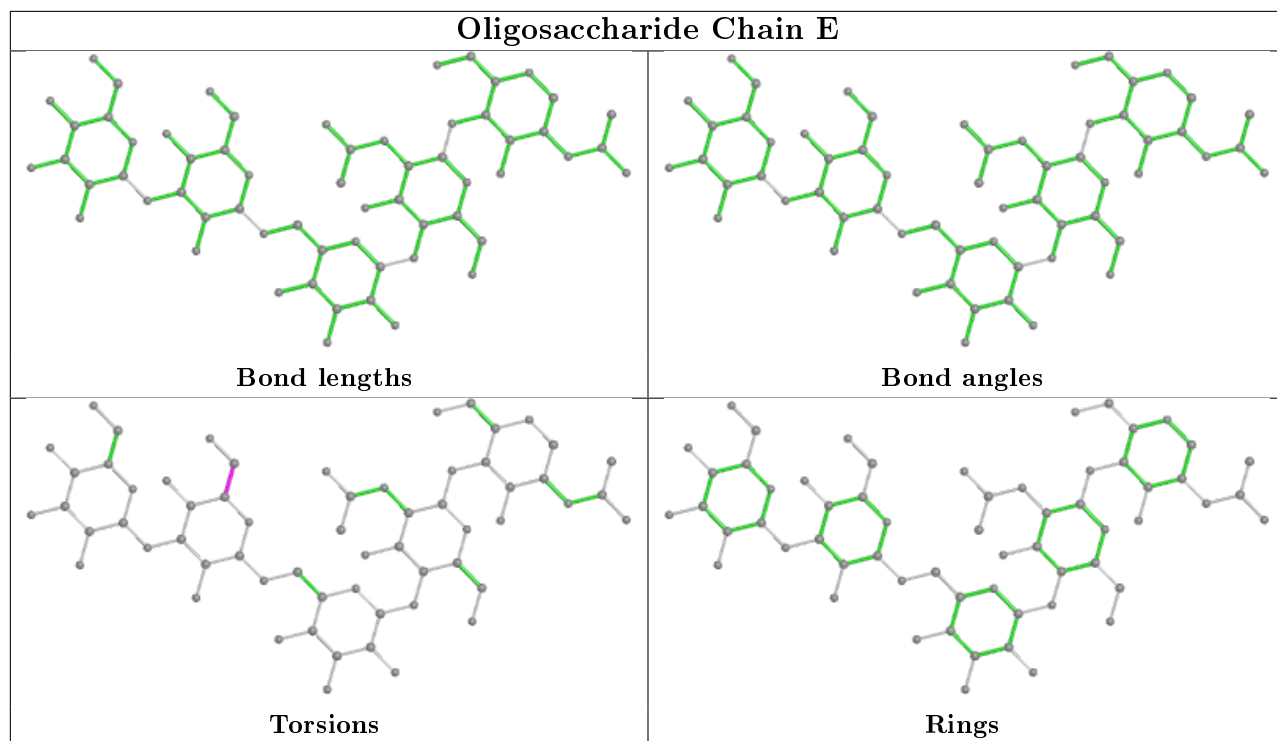
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	E	2	NAG	1	0
5	E	1	NAG	2	0
4	D	1	NAG	1	0
5	E	4	MAN	1	0

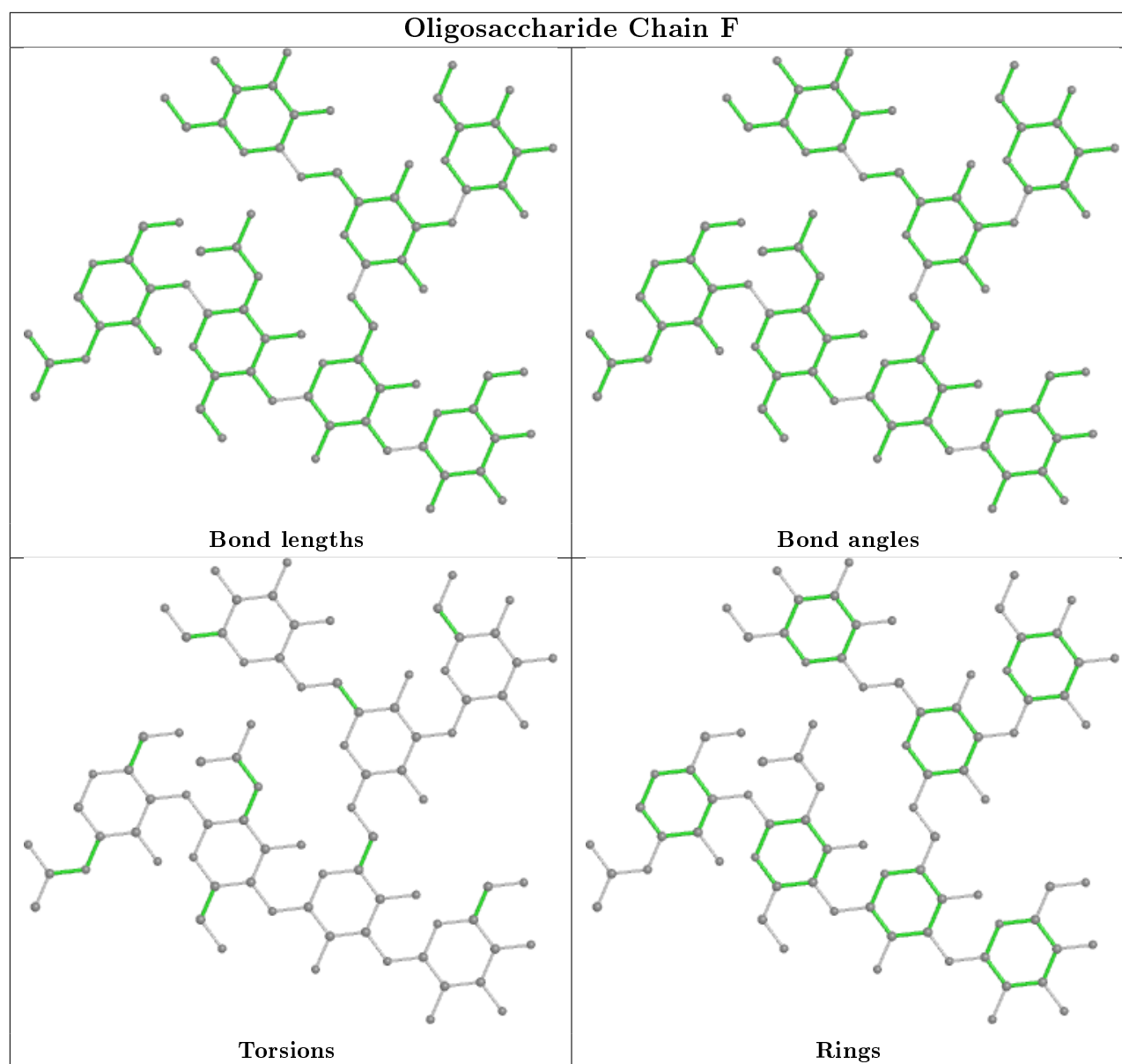
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











5.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 3 are monoatomic - leaving 15 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
10	EDO	A	9229	-	3,3,3	0.46	0	2,2,2	0.50	0
9	PO4	A	9111	-	4,4,4	1.62	0	6,6,6	0.43	0
10	EDO	A	9228	-	3,3,3	0.57	0	2,2,2	0.44	0
10	EDO	A	9226	-	3,3,3	0.57	0	2,2,2	0.42	0
10	EDO	A	9222	-	3,3,3	0.56	0	2,2,2	0.42	0
10	EDO	A	9221	-	3,3,3	0.55	0	2,2,2	0.41	0
10	EDO	A	9225	-	3,3,3	0.69	0	2,2,2	0.44	0
9	PO4	A	9112	-	4,4,4	1.64	0	6,6,6	0.43	0
7	NAG	A	4001	1	14,14,15	0.48	0	17,19,21	0.70	1 (5%)
7	NAG	A	2001	1	14,14,15	0.50	0	17,19,21	0.67	0
10	EDO	A	9223	-	3,3,3	0.57	0	2,2,2	0.42	0
10	EDO	A	9227	-	3,3,3	0.61	0	2,2,2	0.44	0
10	EDO	A	9224	-	3,3,3	0.57	0	2,2,2	0.44	0
9	PO4	A	9113	-	4,4,4	1.72	0	6,6,6	0.43	0
9	PO4	A	9114	-	4,4,4	1.70	0	6,6,6	0.44	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
10	EDO	A	9229	-	-	0/1/1/1	-
10	EDO	A	9222	-	-	0/1/1/1	-
10	EDO	A	9228	-	-	0/1/1/1	-
10	EDO	A	9226	-	-	0/1/1/1	-
10	EDO	A	9221	-	-	0/1/1/1	-
10	EDO	A	9225	-	-	0/1/1/1	-
7	NAG	A	4001	1	-	0/6/23/26	0/1/1/1
7	NAG	A	2001	1	-	5/6/23/26	0/1/1/1
10	EDO	A	9223	-	-	0/1/1/1	-
10	EDO	A	9227	-	-	0/1/1/1	-
10	EDO	A	9224	-	-	0/1/1/1	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	4001	NAG	C2-N2-C7	-2.08	119.94	122.90

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	2001	NAG	C8-C7-N2-C2
7	A	2001	NAG	O7-C7-N2-C2
7	A	2001	NAG	O5-C5-C6-O6
7	A	2001	NAG	C4-C5-C6-O6
7	A	2001	NAG	C3-C2-N2-C7

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
10	A	9229	EDO	2	0
10	A	9228	EDO	1	0
10	A	9226	EDO	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 [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	971/971 (100%)	0.01	6 (0%) 89 90	12, 19, 30, 42	30 (3%)

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	583	THR	5.0
1	A	809	TRP	3.5
1	A	41	LEU	3.2
1	A	671	GLN	2.5
1	A	239	TRP	2.1
1	A	653[A]	VAL	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
3	MAN	C	8	11/12	0.60	0.22	34,36,37,37	11
3	BMA	C	9	11/12	0.68	0.30	45,50,54,56	0
5	BMA	E	3	11/12	0.70	0.27	44,46,47,48	0
5	MAN	E	5	11/12	0.70	0.20	41,43,45,47	11
3	MAN	C	7	11/12	0.73	0.25	26,33,34,35	11
4	NAG	D	2	14/15	0.74	0.23	41,47,48,51	14

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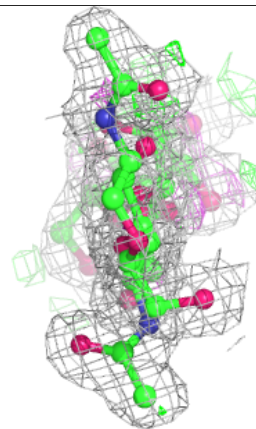
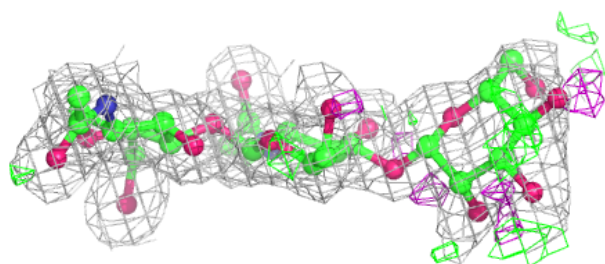
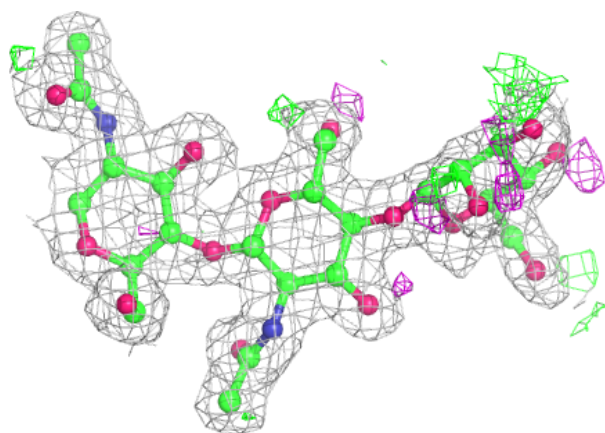
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	BMA	B	3	11/12	0.83	0.20	38,43,46,48	0
5	MAN	E	4	11/12	0.84	0.17	43,44,45,46	11
5	NAG	E	2	14/15	0.86	0.24	37,42,45,46	0
4	NAG	D	1	14/15	0.89	0.16	37,42,45,45	0
6	MAN	F	6	11/12	0.91	0.13	27,32,35,37	0
5	NAG	E	1	14/15	0.93	0.11	20,27,33,38	0
3	MAN	C	6	11/12	0.94	0.10	28,31,32,33	0
6	MAN	F	5	11/12	0.94	0.11	27,31,32,33	0
2	NAG	B	2	14/15	0.95	0.10	18,21,29,30	0
3	MAN	C	4	11/12	0.96	0.11	20,21,24,29	0
2	NAG	B	1	14/15	0.96	0.10	16,17,18,18	0
6	MAN	F	7	11/12	0.96	0.09	25,28,31,33	0
6	NAG	F	1	14/15	0.97	0.10	15,17,21,26	0
3	BMA	C	3	11/12	0.97	0.10	20,21,26,35	0
3	MAN	C	5	11/12	0.97	0.11	19,21,26,26	0
6	NAG	F	2	14/15	0.97	0.11	14,19,27,28	0
3	NAG	C	2	14/15	0.97	0.11	17,19,24,27	0
6	MAN	F	4	11/12	0.97	0.10	19,22,26,26	0
3	NAG	C	1	14/15	0.97	0.12	16,20,28,29	0
6	BMA	F	3	11/12	0.98	0.10	18,19,20,21	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.

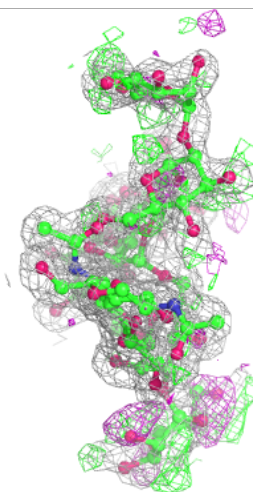
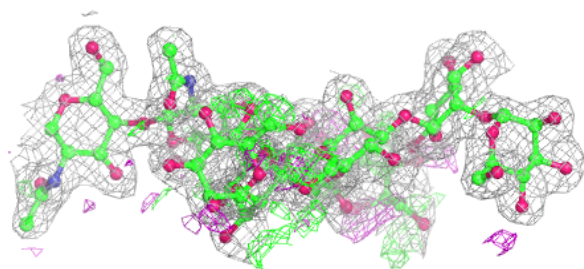
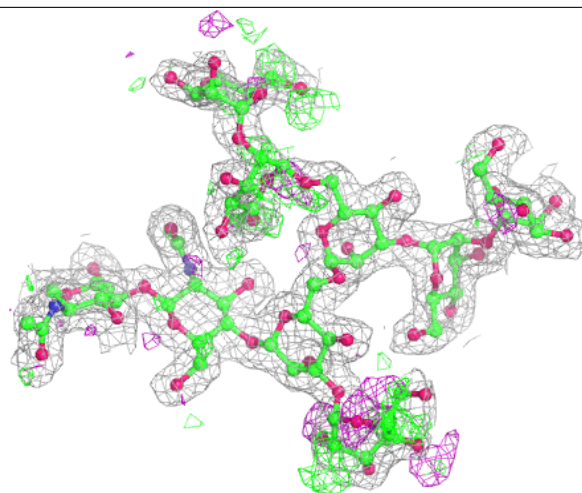
Electron density around Chain B:

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



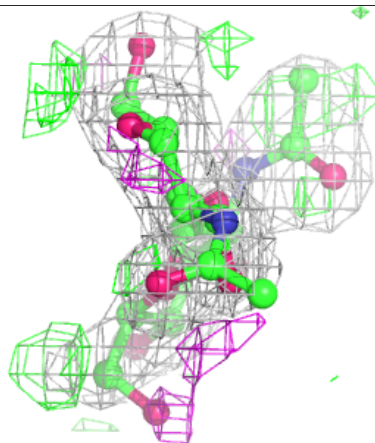
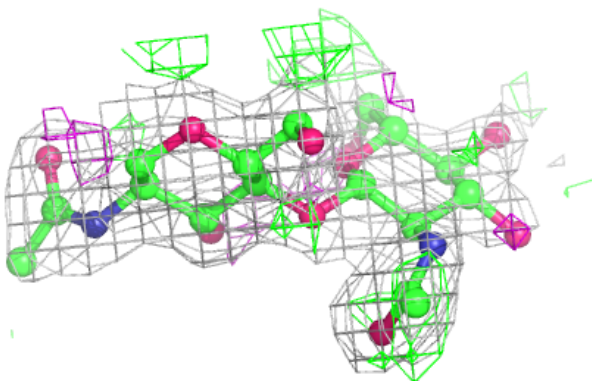
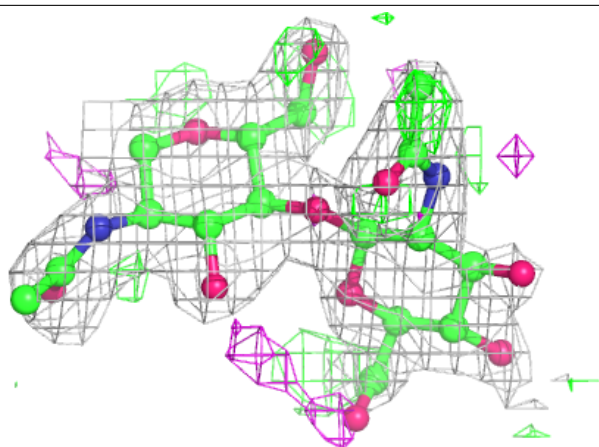
Electron density around Chain C:

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



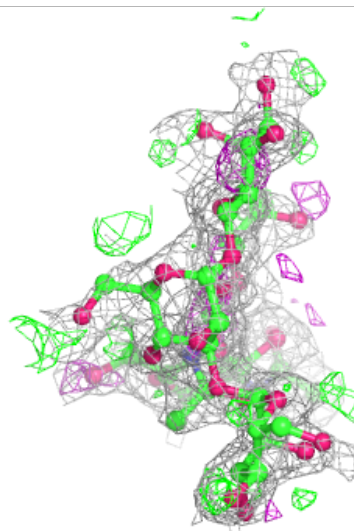
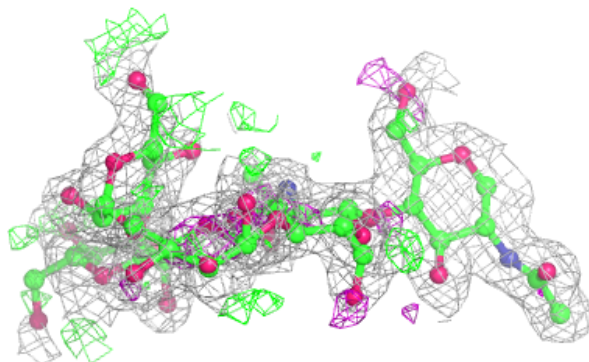
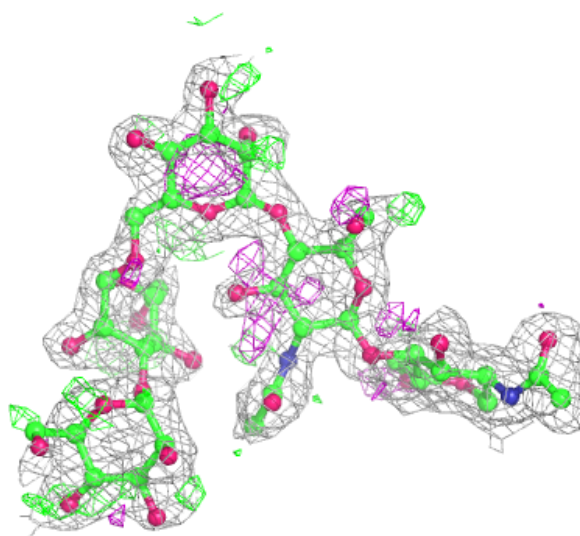
Electron density around Chain D:

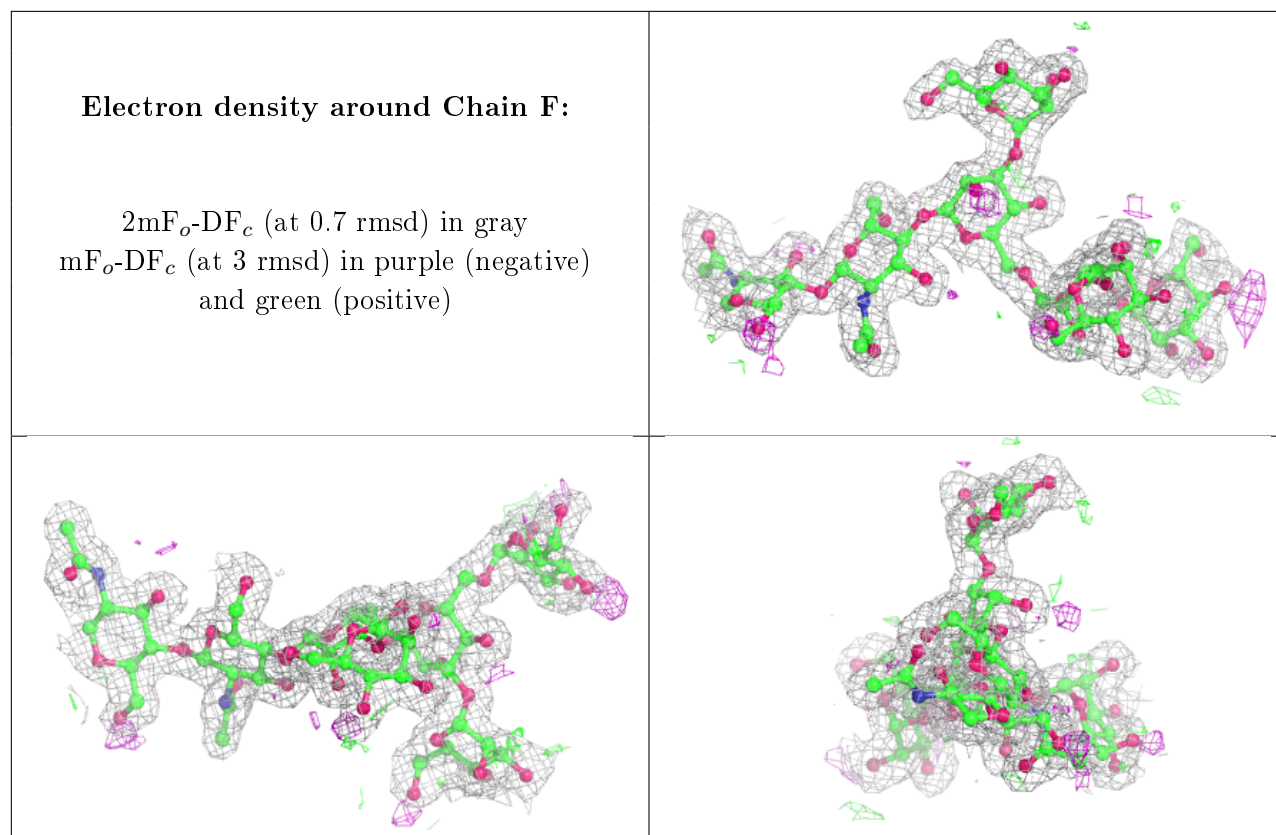
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain E:

$2mF_o-DF_c$ (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
7	NAG	A	2001	14/15	0.69	0.28	48,50,54,55	0
10	EDO	A	9225	4/4	0.83	0.16	26,32,34,34	0
10	EDO	A	9227	4/4	0.83	0.21	32,38,40,43	0
7	NAG	A	4001	14/15	0.88	0.13	26,31,36,38	0
9	PO4	A	9112	5/5	0.90	0.19	37,38,39,39	5
10	EDO	A	9229	4/4	0.90	0.20	23,25,26,30	0
10	EDO	A	9228	4/4	0.92	0.13	18,23,25,27	0
9	PO4	A	9114	5/5	0.92	0.18	53,54,55,55	5
9	PO4	A	9111	5/5	0.93	0.15	12,15,22,22	5
10	EDO	A	9226	4/4	0.94	0.15	33,33,34,37	0
10	EDO	A	9222	4/4	0.95	0.14	29,31,32,35	0
10	EDO	A	9223	4/4	0.95	0.22	34,36,37,40	0
10	EDO	A	9221	4/4	0.95	0.14	23,23,24,27	0
10	EDO	A	9224	4/4	0.96	0.15	26,27,28,30	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
8	NA	A	9003	1/1	0.97	0.12	23,23,23,23	0
8	NA	A	9001	1/1	0.98	0.15	16,16,16,16	0
9	PO4	A	9113	5/5	0.98	0.14	31,32,33,33	5
8	NA	A	9002	1/1	0.99	0.14	18,18,18,18	0

6.5 Other polymers [i](#)

There are no such residues in this entry.