



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:56 AM UTC

PDB ID : 1E5N / pdb_00001e5n
Title : E246C mutant of *P fluorescens* subsp. *cellulosa* xylanase A in complex with xylopentaose
Authors : Lo Leggio, L.; Jenkins, J.A.; Harris, G.W.; Pickersgill, R.W.
Deposited on : 2000-07-27
Resolution : 3.20 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

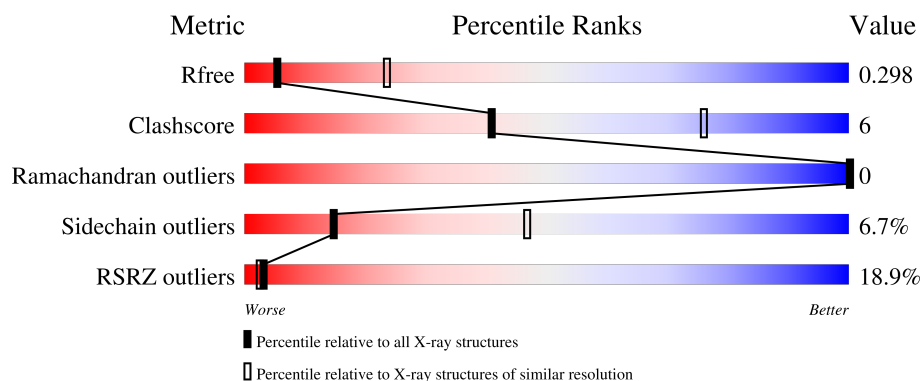
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	348	4% 78% 18% • •
1	B	348	33% 77% 18% • •
2	C	5	80% 20%
2	D	5	60% 40%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5494 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 ENDO-1,4-BETA-XYLANASE A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	346	Total	C	N	O	S	0	0	0
			2700	1693	480	518	9			
1	B	346	Total	C	N	O	S	0	0	0
			2700	1693	480	518	9			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	246	CYS	GLU	engineered mutation	UNP P14768
B	246	CYS	GLU	engineered mutation	UNP P14768

- Molecule 2 is an oligosaccharide called beta-D-xylopyranose-(1-4)-beta-D-xylopyranose-(1-4)-beta-D-xylopyranose-(1-4)-beta-D-xylopyranose-(1-4)-beta-D-xylopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	5	Total	C	O	0	0	0
			46	25	21			
2	D	5	Total	C	O	0	0	0
			46	25	21			

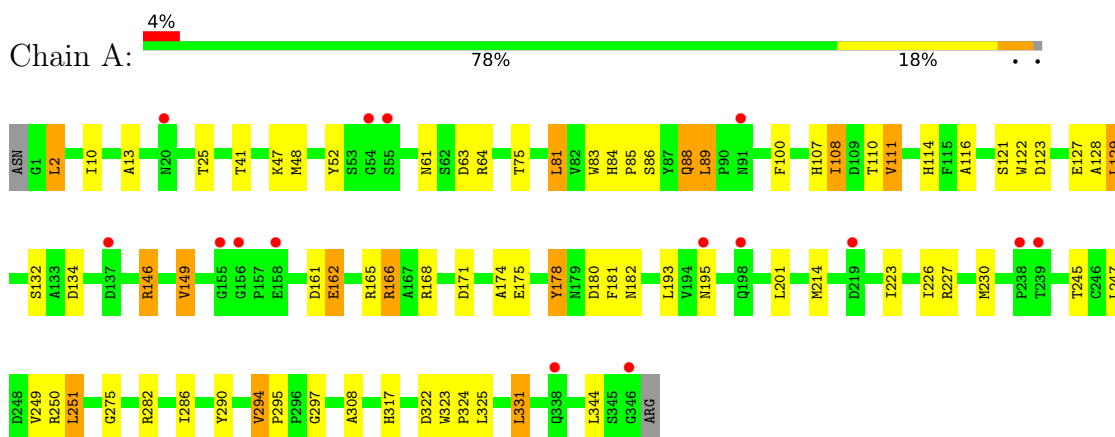
- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	0
			1	1		
3	B	1	Total	Ca	0	0
			1	1		

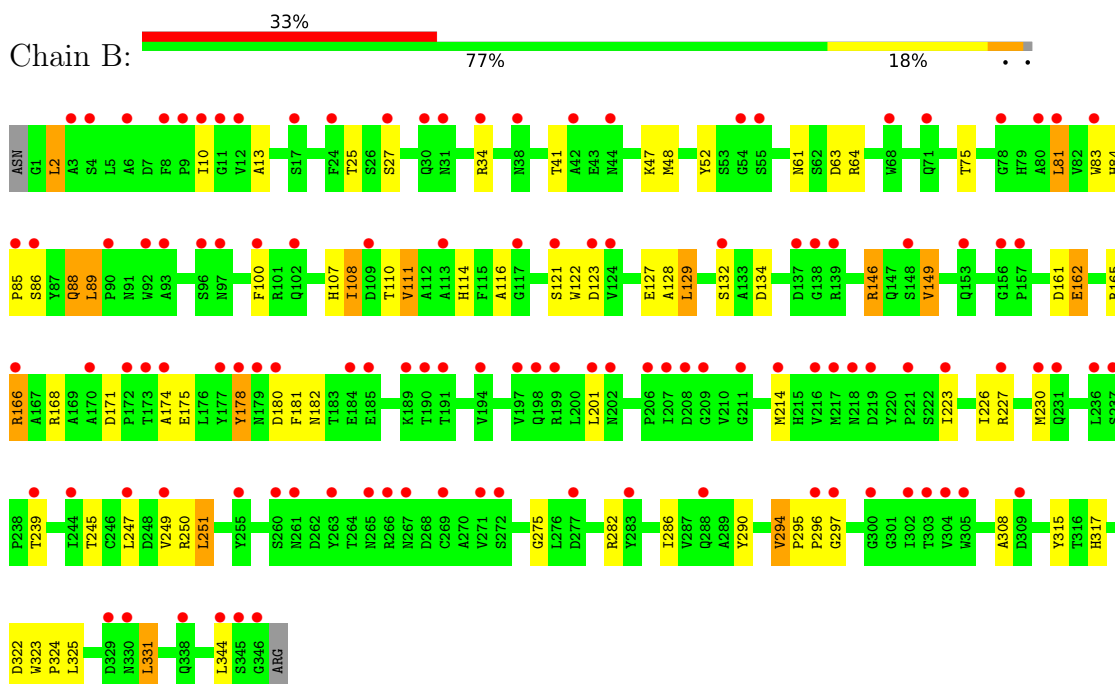
3 Residue-property plots [i](#)

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

• Molecule 1: ENDO-1,4-BETA-XYLANASE A

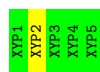


• Molecule 1: ENDO-1,4-BETA-XYLANASE A



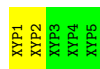
• Molecule 2: beta-D-xylopyranose-(1-4)-beta-D-xylopyranose-(1-4)-beta-D-xylopyranose-(1-4)-beta-D-xylopyranose-(1-4)-beta-D-xylopyranose

Chain C:  80% 20%



- Molecule 2: beta-D-xylopyranose-(1-4)-beta-D-xylopyranose-(1-4)-beta-D-xylopyranose-(1-4)-beta-D-xylopyranose-(1-4)-beta-D-xylopyranose

Chain D:  60% 40%



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	96.70Å 96.70Å 152.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.70 – 3.20 29.70 – 3.20	Depositor EDS
% Data completeness (in resolution range)	90.7 (29.70-3.20) 90.8 (29.70-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.76 (at 3.18Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.190 , 0.245 0.272 , 0.298	Depositor DCC
R_{free} test set	773 reflections (6.78%)	wwPDB-VP
Wilson B-factor (Å ²)	39.4	Xtriage
Anisotropy	0.145	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 56.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.82	EDS
Total number of atoms	5494	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.21% 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: CA, XYP

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	0.80	0/2767	1.22	21/3770 (0.6%)
1	B	0.80	0/2767	1.22	22/3770 (0.6%)
All	All	0.80	0/5534	1.22	43/7540 (0.6%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	324	PRO	N-CA-C	9.97	126.30	114.03
1	A	324	PRO	N-CA-C	9.93	126.25	114.03
1	A	149	VAL	CB-CA-C	-9.37	99.77	112.04
1	B	149	VAL	CB-CA-C	-9.35	99.79	112.04
1	A	84	HIS	N-CA-C	7.94	127.36	109.81
1	B	84	HIS	N-CA-C	7.93	127.33	109.81
1	B	123	ASP	N-CA-C	-7.38	95.54	108.23
1	B	294	VAL	CA-C-N	7.38	124.97	119.66
1	B	294	VAL	C-N-CA	7.38	124.97	119.66
1	A	123	ASP	N-CA-C	-7.37	95.56	108.23
1	A	294	VAL	CA-C-N	7.36	124.96	119.66
1	A	294	VAL	C-N-CA	7.36	124.96	119.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	178	TYR	N-CA-C	-6.82	98.13	109.24
1	B	178	TYR	N-CA-C	-6.82	98.13	109.24
1	A	110	THR	N-CA-C	6.75	118.42	111.14
1	B	110	THR	N-CA-C	6.72	118.40	111.14
1	B	323	TRP	CA-C-N	6.71	126.85	119.87
1	B	323	TRP	C-N-CA	6.71	126.85	119.87
1	A	323	TRP	CA-C-N	6.70	126.84	119.87
1	A	323	TRP	C-N-CA	6.70	126.84	119.87
1	A	48	MET	N-CA-C	6.53	122.07	111.37
1	B	48	MET	N-CA-C	6.51	122.06	111.37
1	B	245	THR	N-CA-C	5.84	120.57	112.45
1	A	245	THR	N-CA-C	5.83	120.56	112.45
1	B	100	PHE	N-CA-C	5.64	117.87	111.11
1	A	100	PHE	N-CA-C	5.63	117.86	111.11
1	B	75	THR	N-CA-C	-5.59	101.69	110.36
1	A	75	THR	N-CA-C	-5.58	101.72	110.36
1	B	325	LEU	N-CA-C	5.55	118.73	109.96
1	A	325	LEU	N-CA-C	5.54	118.72	109.96
1	B	116	ALA	N-CA-C	5.44	122.39	110.80
1	A	116	ALA	N-CA-C	5.43	122.38	110.80
1	B	247	LEU	N-CA-C	5.42	118.17	110.10
1	A	247	LEU	N-CA-C	5.39	118.13	110.10
1	B	297	GLY	N-CA-C	-5.17	108.50	115.21
1	A	297	GLY	N-CA-C	-5.15	108.52	115.21
1	B	249	VAL	N-CA-C	5.09	115.79	107.24
1	B	127	GLU	N-CA-C	5.09	118.64	111.52
1	A	127	GLU	N-CA-C	5.08	118.62	111.52
1	A	249	VAL	N-CA-C	5.07	115.76	107.24
1	B	52	TYR	N-CA-C	5.06	117.50	109.50
1	A	52	TYR	N-CA-C	5.06	117.50	109.50
1	B	315	TYR	N-CA-C	-5.01	105.95	111.71

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	178	TYR	Sidechain
1	B	178	TYR	Sidechain

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	2700	0	2562	34	1
1	B	2700	0	2562	35	5
2	C	46	0	0	0	0
2	D	46	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
All	All	5494	0	5124	69	5

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (69) 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:161:ASP:O	1:B:165:ARG:HG3	1.94	0.67
1:A:161:ASP:O	1:A:165:ARG:HG3	1.94	0.66
1:B:230:MET:HE1	1:B:290:TYR:HD1	1.64	0.62
1:A:230:MET:HE1	1:A:290:TYR:HD1	1.64	0.61
1:A:108:ILE:HG21	1:A:166:ARG:HB3	1.84	0.60
1:B:61:ASN:HD22	1:B:64:ARG:HH12	1.50	0.60
1:A:61:ASN:HD22	1:A:64:ARG:HH12	1.50	0.58
1:B:107:HIS:O	1:B:111:VAL:HG13	2.03	0.58
1:B:108:ILE:HG21	1:B:166:ARG:HB3	1.84	0.58
1:A:107:HIS:O	1:A:111:VAL:HG13	2.03	0.58
1:A:47:LYS:HB3	1:A:88:GLN:HE21	1.70	0.57
1:B:47:LYS:HB3	1:B:88:GLN:HE21	1.70	0.56
1:B:47:LYS:HD3	1:B:83:TRP:CZ3	2.41	0.56
1:A:83:TRP:CZ2	1:A:85:PRO:HG2	2.41	0.56
1:A:47:LYS:HD3	1:A:83:TRP:CZ3	2.40	0.55
1:B:83:TRP:CZ2	1:B:85:PRO:HG2	2.41	0.55
1:A:47:LYS:HD3	1:A:83:TRP:HZ3	1.72	0.55
1:B:47:LYS:HD3	1:B:83:TRP:HZ3	1.72	0.53
1:B:134:ASP:O	1:B:146:ARG:NH2	2.42	0.53
1:A:134:ASP:O	1:A:146:ARG:NH2	2.42	0.53
1:A:251:LEU:HD13	1:A:275:GLY:HA3	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:251:LEU:HD13	1:B:275:GLY:HA3	1.92	0.51
1:A:308:ALA:HB2	1:A:331:LEU:CD1	2.42	0.50
1:B:308:ALA:HB2	1:B:331:LEU:CD1	2.42	0.49
1:A:317:HIS:HE1	1:A:322:ASP:OD2	1.95	0.49
1:B:317:HIS:HE1	1:B:322:ASP:OD2	1.95	0.48
1:B:308:ALA:HB2	1:B:331:LEU:HD12	1.96	0.47
1:A:308:ALA:HB2	1:A:331:LEU:HD12	1.96	0.47
1:B:181:PHE:O	1:B:182:ASN:HB2	2.16	0.46
1:B:129:LEU:HD12	1:B:129:LEU:HA	1.68	0.46
1:A:63:ASP:OD1	1:A:114:HIS:HE1	1.99	0.46
1:B:121:SER:HB3	1:B:175:GLU:HB2	1.98	0.46
1:A:121:SER:HB3	1:A:175:GLU:HB2	1.98	0.45
1:A:171:ASP:OD2	1:A:174:ALA:HB2	2.17	0.45
1:A:181:PHE:O	1:A:182:ASN:HB2	2.15	0.45
1:A:214:MET:HE1	1:A:226:ILE:HG23	1.99	0.45
1:B:214:MET:HE1	1:B:226:ILE:HG23	1.99	0.44
1:A:294:VAL:HA	1:A:295:PRO:HD3	1.75	0.44
1:B:63:ASP:OD1	1:B:114:HIS:HE1	1.99	0.44
1:B:171:ASP:OD2	1:B:174:ALA:HB2	2.17	0.44
1:B:295:PRO:HA	1:B:296:PRO:HD3	1.90	0.44
1:B:162:GLU:O	1:B:166:ARG:HB2	2.18	0.43
1:A:201:LEU:HD23	1:A:201:LEU:HA	1.82	0.43
1:A:2:LEU:HB3	1:A:10:ILE:HG21	2.00	0.43
1:A:129:LEU:HD12	1:A:129:LEU:HA	1.68	0.43
1:A:86:SER:HA	1:A:89:LEU:HD22	2.01	0.42
1:A:162:GLU:O	1:A:166:ARG:HB2	2.18	0.42
1:B:61:ASN:ND2	1:B:64:ARG:HH12	2.16	0.42
1:A:81:LEU:HD13	1:A:122:TRP:CZ3	2.55	0.42
1:B:2:LEU:HB3	1:B:10:ILE:HG21	2.00	0.42
1:B:250:ARG:HG2	1:B:317:HIS:CE1	2.55	0.42
1:A:61:ASN:ND2	1:A:64:ARG:HH12	2.16	0.42
1:B:81:LEU:HD13	1:B:122:TRP:CZ3	2.55	0.42
1:A:331:LEU:HD12	1:A:331:LEU:HA	1.81	0.42
1:A:128:ALA:O	1:A:146:ARG:HB3	2.20	0.41
1:A:129:LEU:HB2	1:A:180:ASP:OD2	2.21	0.41
1:A:250:ARG:HG2	1:A:317:HIS:CE1	2.55	0.41
1:B:86:SER:HA	1:B:89:LEU:HD22	2.01	0.41
1:B:129:LEU:HB2	1:B:180:ASP:OD2	2.21	0.41
1:B:294:VAL:HA	1:B:295:PRO:HD3	1.75	0.41
1:B:128:ALA:O	1:B:146:ARG:HB3	2.20	0.41
1:B:201:LEU:HA	1:B:201:LEU:HD23	1.82	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:282:ARG:O	1:B:286:ILE:HG13	2.21	0.40
1:A:13:ALA:HA	1:A:41:THR:O	2.21	0.40
1:A:129:LEU:HD21	1:A:193:LEU:HA	2.04	0.40
1:B:331:LEU:HD12	1:B:331:LEU:HA	1.81	0.40
1:A:282:ARG:O	1:A:286:ILE:HG13	2.21	0.40
1:B:13:ALA:HA	1:B:41:THR:O	2.22	0.40
1:B:61:ASN:HD22	1:B:64:ARG:NH1	2.17	0.40

All (5) 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:195:ASN:OD1	1:B:239:THR:CG2[5_645]	1.43	0.77
1:B:27:SER:CB	1:B:34:ARG:NH1[7_556]	1.63	0.57
1:B:27:SER:OG	1:B:34:ARG:NH1[7_556]	1.78	0.42
1:B:27:SER:CB	1:B:34:ARG:CZ[7_556]	2.03	0.17
1:B:27:SER:OG	1:B:34:ARG:CZ[7_556]	2.09	0.11

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	344/348 (99%)	329 (96%)	15 (4%)	0	100	100
1	B	344/348 (99%)	329 (96%)	15 (4%)	0	100	100
All	All	688/696 (99%)	658 (96%)	30 (4%)	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	283/285 (99%)	264 (93%)	19 (7%)	15	47
1	B	283/285 (99%)	264 (93%)	19 (7%)	15	47
All	All	566/570 (99%)	528 (93%)	38 (7%)	15	47

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

Mol	Chain	Res	Type
1	A	2	LEU
1	A	25	THR
1	A	81	LEU
1	A	88	GLN
1	A	89	LEU
1	A	108	ILE
1	A	111	VAL
1	A	129	LEU
1	A	132	SER
1	A	146	ARG
1	A	149	VAL
1	A	162	GLU
1	A	166	ARG
1	A	168	ARG
1	A	223	ILE
1	A	227	ARG
1	A	251	LEU
1	A	331	LEU
1	A	344	LEU
1	B	2	LEU
1	B	25	THR
1	B	81	LEU
1	B	88	GLN
1	B	89	LEU
1	B	108	ILE
1	B	111	VAL
1	B	129	LEU

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Mol	Chain	Res	Type
1	B	132	SER
1	B	146	ARG
1	B	149	VAL
1	B	162	GLU
1	B	166	ARG
1	B	168	ARG
1	B	223	ILE
1	B	227	ARG
1	B	251	LEU
1	B	331	LEU
1	B	344	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	61	ASN
1	A	88	GLN
1	A	107	HIS
1	A	114	HIS
1	A	147	GLN
1	A	195	ASN
1	A	203	ASN
1	A	231	GLN
1	A	317	HIS
1	B	44	ASN
1	B	61	ASN
1	B	88	GLN
1	B	107	HIS
1	B	114	HIS
1	B	147	GLN
1	B	195	ASN
1	B	203	ASN
1	B	231	GLN
1	B	317	HIS

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 ⓘ

10 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	XYP	C	1	2	10,10,10	0.69	0	14,14,14	0.95	0
2	XYP	C	2	2	9,9,10	1.15	1 (11%)	10,12,14	1.03	1 (10%)
2	XYP	C	3	2	9,9,10	0.69	0	10,12,14	0.48	0
2	XYP	C	4	2	9,9,10	0.86	0	10,12,14	0.83	0
2	XYP	C	5	2	9,9,10	0.46	0	10,12,14	0.64	0
2	XYP	D	1	2	10,10,10	0.69	0	14,14,14	0.95	1 (7%)
2	XYP	D	2	2	9,9,10	1.15	1 (11%)	10,12,14	1.03	1 (10%)
2	XYP	D	3	2	9,9,10	0.68	0	10,12,14	0.48	0
2	XYP	D	4	2	9,9,10	0.86	0	10,12,14	0.84	0
2	XYP	D	5	2	9,9,10	0.45	0	10,12,14	0.64	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	XYP	C	1	2	-	-	0/1/1/1
2	XYP	C	2	2	-	-	0/1/1/1
2	XYP	C	3	2	-	-	0/1/1/1
2	XYP	C	4	2	-	-	0/1/1/1
2	XYP	C	5	2	-	-	0/1/1/1
2	XYP	D	1	2	-	-	0/1/1/1
2	XYP	D	2	2	-	-	0/1/1/1
2	XYP	D	3	2	-	-	0/1/1/1
2	XYP	D	4	2	-	-	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	XYP	D	5	2	-	-	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	2	XYP	C2-C3	-2.27	1.49	1.52
2	D	2	XYP	C2-C3	-2.26	1.49	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	XYP	C1-C2-C3	-2.04	106.67	109.64
2	C	2	XYP	C1-C2-C3	-2.02	106.71	109.64
2	D	1	XYP	C5-C4-C3	-2.01	106.72	109.64

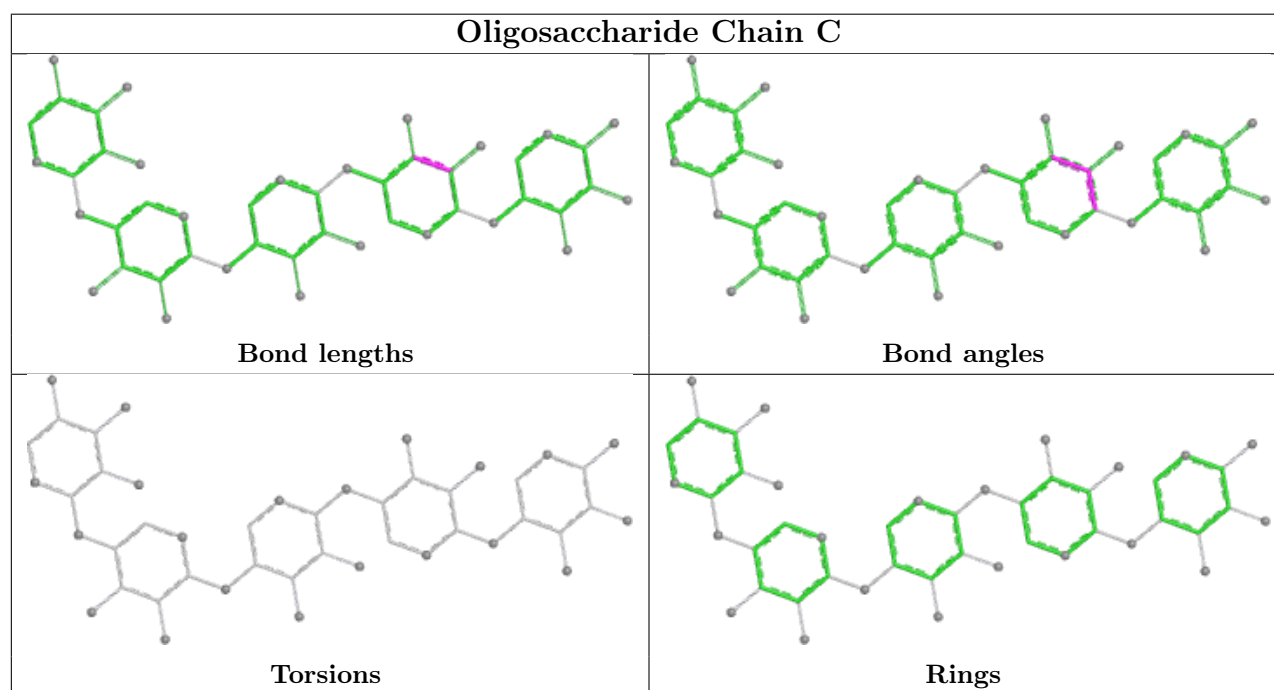
There are no chirality outliers.

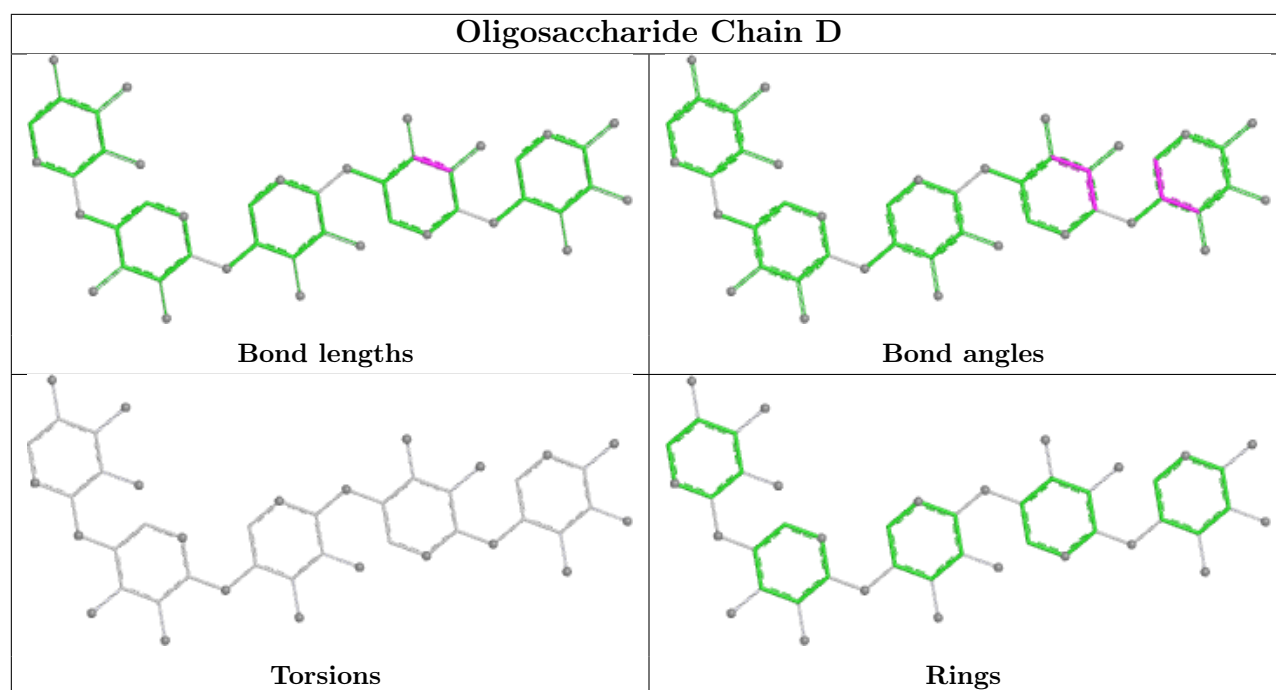
There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

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

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	346/348 (99%)	0.57	15 (4%) 40 24	7, 28, 62, 81	0
1	B	346/348 (99%)	1.76	116 (33%) 1 1	7, 28, 62, 81	0
All	All	692/696 (99%)	1.17	131 (18%) 3 2	7, 28, 63, 81	0

All (131) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	209	GLY	4.4
1	B	174	ALA	4.2
1	B	236	LEU	4.1
1	B	27	SER	4.1
1	B	346	GLY	4.0
1	B	300	GLY	3.9
1	B	30	GLN	3.9
1	A	346	GLY	3.8
1	B	10	ILE	3.6
1	B	288	GLN	3.6
1	B	93	ALA	3.4
1	B	3	ALA	3.4
1	B	345	SER	3.3
1	B	148	SER	3.3
1	B	249	VAL	3.3
1	B	271	VAL	3.2
1	B	156	GLY	3.2
1	B	283	TYR	3.2
1	B	223	ILE	3.1
1	B	263	TYR	3.1
1	B	172	PRO	3.1
1	B	338	GLN	3.0
1	B	197	VAL	3.0
1	A	137	ASP	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	124	VAL	3.0
1	B	8	PHE	3.0
1	A	54	GLY	2.9
1	B	9	PRO	2.9
1	B	4	SER	2.9
1	B	117	GLY	2.9
1	B	71	GLN	2.8
1	B	303	THR	2.8
1	B	297	GLY	2.8
1	B	230	MET	2.8
1	A	219	ASP	2.8
1	B	121	SER	2.8
1	B	296	PRO	2.8
1	B	304	VAL	2.8
1	B	309	ASP	2.8
1	B	34	ARG	2.8
1	B	214	MET	2.8
1	B	206	PRO	2.7
1	B	260	SER	2.7
1	A	198	GLN	2.7
1	B	231	GLN	2.7
1	B	11	GLY	2.7
1	B	54	GLY	2.7
1	B	199	ARG	2.7
1	B	277	ASP	2.7
1	B	227	ARG	2.7
1	B	132	SER	2.7
1	B	6	ALA	2.7
1	B	81	LEU	2.7
1	B	177	TYR	2.7
1	A	91	ASN	2.6
1	B	38	ASN	2.6
1	B	269	CYS	2.6
1	B	86	SER	2.6
1	B	12	VAL	2.6
1	B	255	TYR	2.6
1	B	178	TYR	2.5
1	B	221	PRO	2.5
1	B	267	ASN	2.5
1	B	239	THR	2.5
1	B	211	GLY	2.5
1	B	109	ASP	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	157	PRO	2.5
1	B	244	ILE	2.4
1	B	100	PHE	2.4
1	B	261	ASN	2.4
1	B	83	TRP	2.4
1	B	266	ARG	2.4
1	B	184	GLU	2.4
1	B	185	GLU	2.4
1	B	216	VAL	2.4
1	B	202	ASN	2.4
1	B	123	ASP	2.4
1	B	170	ALA	2.4
1	B	305	TRP	2.3
1	B	42	ALA	2.3
1	B	44	ASN	2.3
1	B	218	ASN	2.3
1	B	208	ASP	2.3
1	B	198	GLN	2.3
1	A	239	THR	2.3
1	B	17	SER	2.3
1	B	137	ASP	2.3
1	B	90	PRO	2.3
1	B	68	TRP	2.3
1	B	78	GLY	2.3
1	B	237	SER	2.3
1	B	80	ALA	2.3
1	B	217	MET	2.3
1	B	96	SER	2.3
1	B	265	ASN	2.3
1	B	102	GLN	2.3
1	B	55	SER	2.2
1	B	85	PRO	2.2
1	B	138	GLY	2.2
1	B	201	LEU	2.2
1	B	166	ARG	2.2
1	A	20	ASN	2.2
1	B	219	ASP	2.2
1	B	247	LEU	2.2
1	A	155	GLY	2.2
1	B	329	ASP	2.2
1	B	194	VAL	2.2
1	A	55	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	153	GLN	2.2
1	B	97	ASN	2.2
1	B	113	ALA	2.2
1	B	180	ASP	2.2
1	A	338	GLN	2.1
1	B	190	THR	2.1
1	B	272	SER	2.1
1	A	195	ASN	2.1
1	B	330	ASN	2.1
1	B	302	ILE	2.1
1	B	173	THR	2.1
1	A	156	GLY	2.1
1	B	344	LEU	2.1
1	B	179	ASN	2.0
1	B	92	TRP	2.0
1	A	158	GLU	2.0
1	B	207	ILE	2.0
1	B	31	ASN	2.0
1	B	191	THR	2.0
1	B	24	PHE	2.0
1	A	238	PRO	2.0
1	B	189	LYS	2.0
1	B	139	ARG	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.

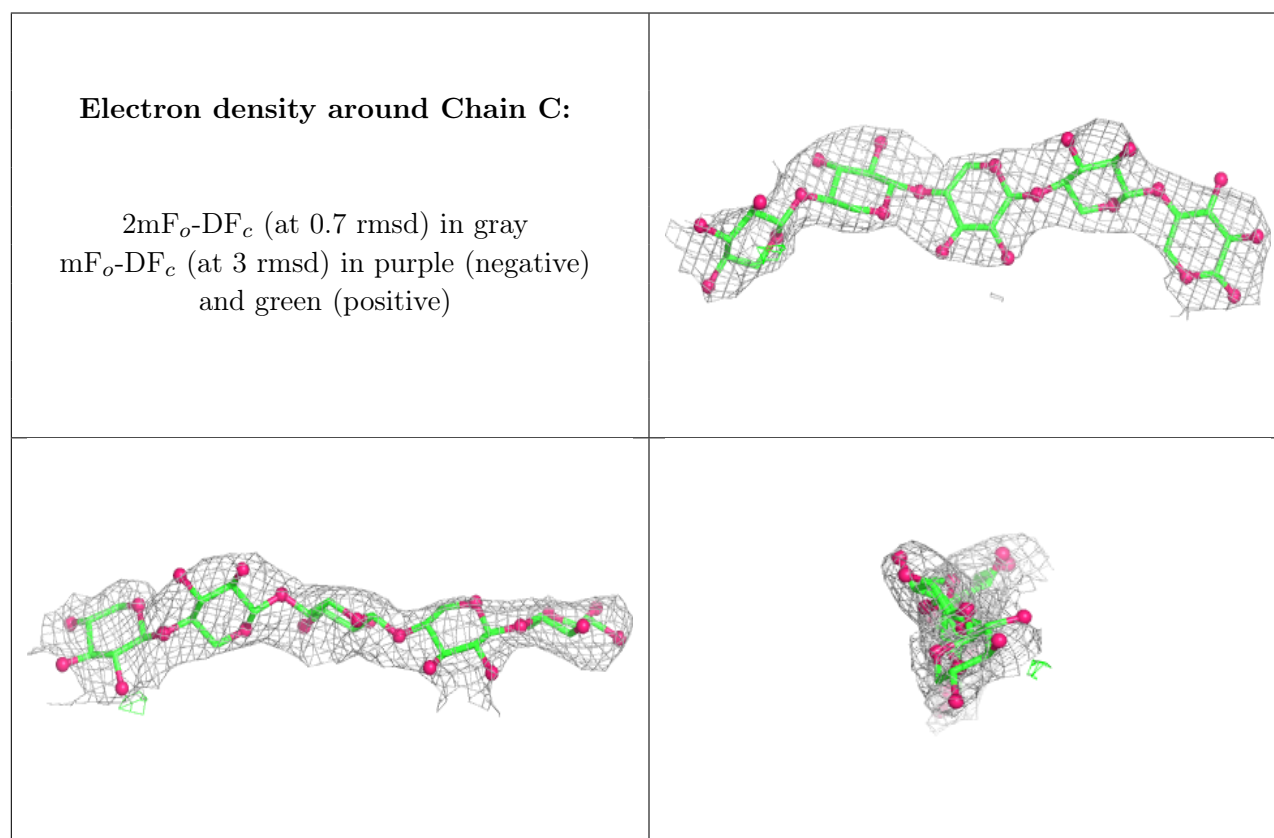
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	XYP	D	4	9/10	0.63	0.24	30,30,30,51	0
2	XYP	C	1	10/10	0.78	0.15	48,54,54,54	0
2	XYP	D	5	9/10	0.78	0.20	51,51,51,51	0
2	XYP	C	2	9/10	0.81	0.15	16,48,48,48	0
2	XYP	C	4	9/10	0.84	0.13	30,30,30,51	0

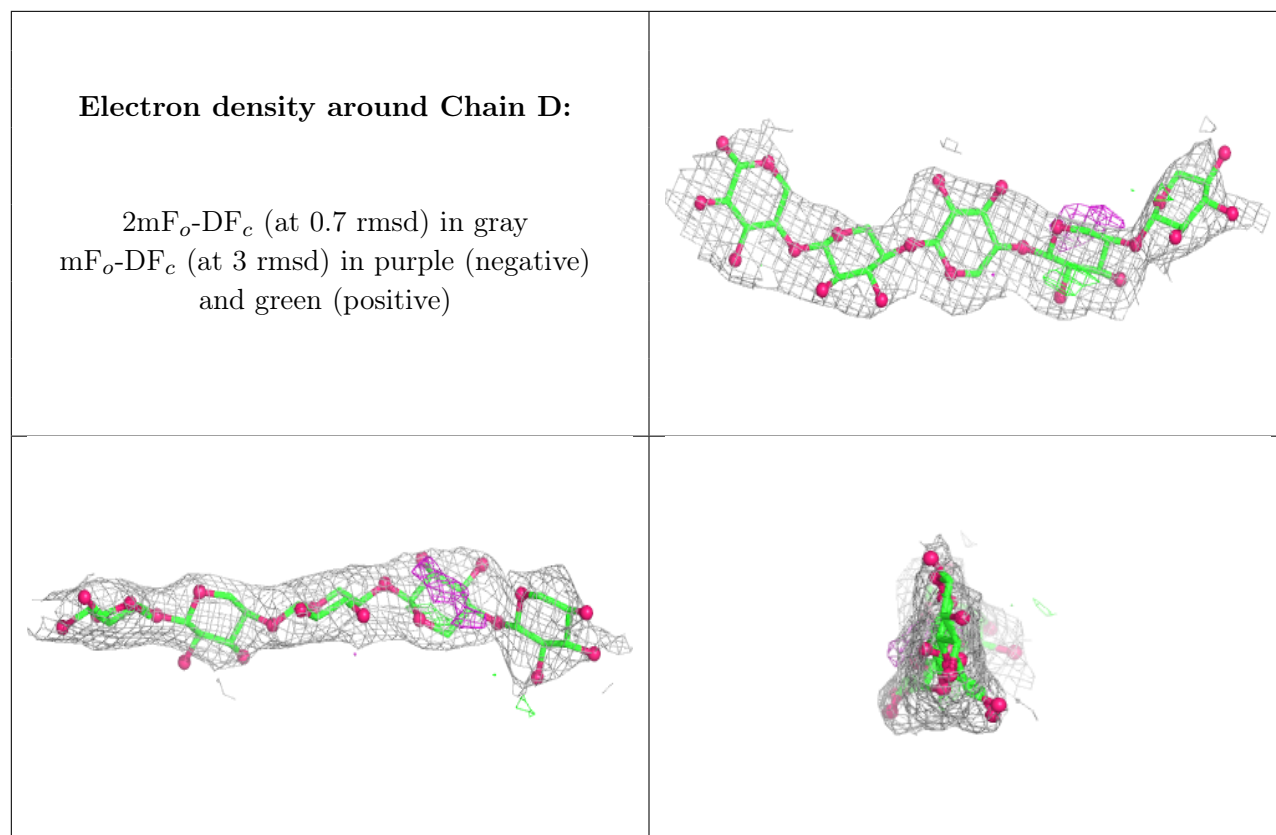
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	XYP	C	5	9/10	0.85	0.16	51,51,51,51	0
2	XYP	D	1	10/10	0.85	0.16	48,54,54,54	0
2	XYP	C	3	9/10	0.89	0.12	16,16,16,30	0
2	XYP	D	2	9/10	0.89	0.17	16,48,48,48	0
2	XYP	D	3	9/10	0.90	0.13	16,16,16,30	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.





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
3	CA	A	348	1/1	0.94	0.06	31,31,31,31	0
3	CA	B	348	1/1	0.95	0.07	31,31,31,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.