



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 29, 2025 – 08:50 am BST

PDB ID : 9FHT / pdb_00009fht
Title : Bacteroides ovatus polysaccharide lyase family 38 (BoPL38) wild type in complex hexaguluronic acid at pH 3.5
Authors : Tandrup, T.; Wilkens, C.
Deposited on : 2024-05-28
Resolution : 2.05 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 Phenix2.0rc1
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.0rc1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.44

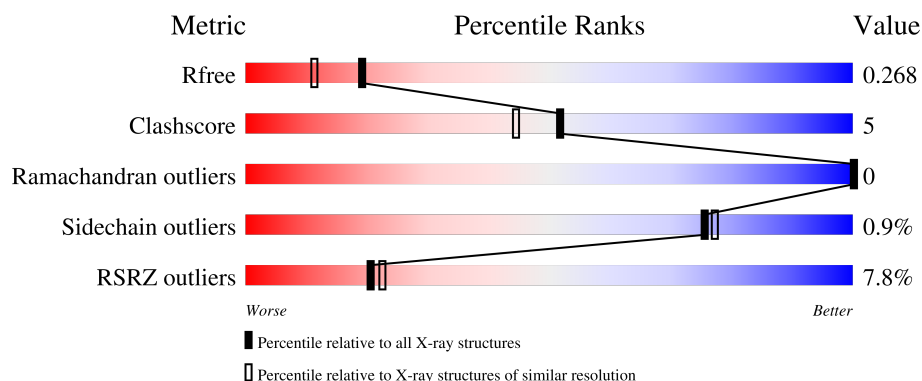
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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.



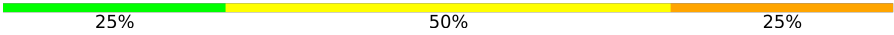


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	164625	2096 (2.04-2.04)
Clashscore	180529	2229 (2.04-2.04)
Ramachandran outliers	177936	2217 (2.04-2.04)
Sidechain outliers	177891	2217 (2.04-2.04)
RSRZ outliers	164620	2096 (2.04-2.04)

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	404	<div> <div>7%</div> <div> <div></div> <div>84%</div> <div>9%</div> <div>6%</div> </div> </div>
1	B	404	<div> <div>5%</div> <div> <div></div> <div>83%</div> <div>11%</div> <div>6%</div> </div> </div>
1	C	404	<div> <div>6%</div> <div> <div></div> <div>86%</div> <div>7%</div> <div>6%</div> </div> </div>
1	D	404	<div> <div>11%</div> <div> <div></div> <div>81%</div> <div>13%</div> <div>6%</div> </div> </div>
2	E	4	<div> <div></div> <div> <div>50%</div> <div>50%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
2	F	4	 25% 50% 25%
2	G	4	 75% 25%
2	H	4	 75% 25%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13069 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 Alginate lyase family protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	379	Total	C	N	O	S	0	0	0
			3046	1949	515	569	13			
1	B	380	Total	C	N	O	S	0	0	0
			3052	1952	516	571	13			
1	C	380	Total	C	N	O	S	0	0	0
			3052	1952	516	571	13			
1	D	380	Total	C	N	O	S	0	1	0
			3058	1957	517	571	13			

There are 92 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	initiating methionine	UNP A0A5M5BWR5
A	-10	GLY	-	expression tag	UNP A0A5M5BWR5
A	-9	SER	-	expression tag	UNP A0A5M5BWR5
A	-8	SER	-	expression tag	UNP A0A5M5BWR5
A	-7	HIS	-	expression tag	UNP A0A5M5BWR5
A	-6	HIS	-	expression tag	UNP A0A5M5BWR5
A	-5	HIS	-	expression tag	UNP A0A5M5BWR5
A	-4	HIS	-	expression tag	UNP A0A5M5BWR5
A	-3	HIS	-	expression tag	UNP A0A5M5BWR5
A	-2	HIS	-	expression tag	UNP A0A5M5BWR5
A	-1	SER	-	expression tag	UNP A0A5M5BWR5
A	0	SER	-	expression tag	UNP A0A5M5BWR5
A	1	GLY	-	expression tag	UNP A0A5M5BWR5
A	2	LEU	-	expression tag	UNP A0A5M5BWR5
A	3	VAL	-	expression tag	UNP A0A5M5BWR5
A	4	PRO	-	expression tag	UNP A0A5M5BWR5
A	5	ARG	-	expression tag	UNP A0A5M5BWR5
A	6	GLY	-	expression tag	UNP A0A5M5BWR5
A	7	SER	-	expression tag	UNP A0A5M5BWR5
A	8	HIS	-	expression tag	UNP A0A5M5BWR5
A	9	MET	-	expression tag	UNP A0A5M5BWR5

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Chain	Residue	Modelled	Actual	Comment	Reference
A	10	ALA	-	expression tag	UNP A0A5M5BWR5
A	11	SER	-	expression tag	UNP A0A5M5BWR5
B	-11	MET	-	initiating methionine	UNP A0A5M5BWR5
B	-10	GLY	-	expression tag	UNP A0A5M5BWR5
B	-9	SER	-	expression tag	UNP A0A5M5BWR5
B	-8	SER	-	expression tag	UNP A0A5M5BWR5
B	-7	HIS	-	expression tag	UNP A0A5M5BWR5
B	-6	HIS	-	expression tag	UNP A0A5M5BWR5
B	-5	HIS	-	expression tag	UNP A0A5M5BWR5
B	-4	HIS	-	expression tag	UNP A0A5M5BWR5
B	-3	HIS	-	expression tag	UNP A0A5M5BWR5
B	-2	HIS	-	expression tag	UNP A0A5M5BWR5
B	-1	SER	-	expression tag	UNP A0A5M5BWR5
B	0	SER	-	expression tag	UNP A0A5M5BWR5
B	1	GLY	-	expression tag	UNP A0A5M5BWR5
B	2	LEU	-	expression tag	UNP A0A5M5BWR5
B	3	VAL	-	expression tag	UNP A0A5M5BWR5
B	4	PRO	-	expression tag	UNP A0A5M5BWR5
B	5	ARG	-	expression tag	UNP A0A5M5BWR5
B	6	GLY	-	expression tag	UNP A0A5M5BWR5
B	7	SER	-	expression tag	UNP A0A5M5BWR5
B	8	HIS	-	expression tag	UNP A0A5M5BWR5
B	9	MET	-	expression tag	UNP A0A5M5BWR5
B	10	ALA	-	expression tag	UNP A0A5M5BWR5
B	11	SER	-	expression tag	UNP A0A5M5BWR5
C	-11	MET	-	initiating methionine	UNP A0A5M5BWR5
C	-10	GLY	-	expression tag	UNP A0A5M5BWR5
C	-9	SER	-	expression tag	UNP A0A5M5BWR5
C	-8	SER	-	expression tag	UNP A0A5M5BWR5
C	-7	HIS	-	expression tag	UNP A0A5M5BWR5
C	-6	HIS	-	expression tag	UNP A0A5M5BWR5
C	-5	HIS	-	expression tag	UNP A0A5M5BWR5
C	-4	HIS	-	expression tag	UNP A0A5M5BWR5
C	-3	HIS	-	expression tag	UNP A0A5M5BWR5
C	-2	HIS	-	expression tag	UNP A0A5M5BWR5
C	-1	SER	-	expression tag	UNP A0A5M5BWR5
C	0	SER	-	expression tag	UNP A0A5M5BWR5
C	1	GLY	-	expression tag	UNP A0A5M5BWR5
C	2	LEU	-	expression tag	UNP A0A5M5BWR5
C	3	VAL	-	expression tag	UNP A0A5M5BWR5
C	4	PRO	-	expression tag	UNP A0A5M5BWR5
C	5	ARG	-	expression tag	UNP A0A5M5BWR5

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Chain	Residue	Modelled	Actual	Comment	Reference
C	6	GLY	-	expression tag	UNP A0A5M5BWR5
C	7	SER	-	expression tag	UNP A0A5M5BWR5
C	8	HIS	-	expression tag	UNP A0A5M5BWR5
C	9	MET	-	expression tag	UNP A0A5M5BWR5
C	10	ALA	-	expression tag	UNP A0A5M5BWR5
C	11	SER	-	expression tag	UNP A0A5M5BWR5
D	-11	MET	-	initiating methionine	UNP A0A5M5BWR5
D	-10	GLY	-	expression tag	UNP A0A5M5BWR5
D	-9	SER	-	expression tag	UNP A0A5M5BWR5
D	-8	SER	-	expression tag	UNP A0A5M5BWR5
D	-7	HIS	-	expression tag	UNP A0A5M5BWR5
D	-6	HIS	-	expression tag	UNP A0A5M5BWR5
D	-5	HIS	-	expression tag	UNP A0A5M5BWR5
D	-4	HIS	-	expression tag	UNP A0A5M5BWR5
D	-3	HIS	-	expression tag	UNP A0A5M5BWR5
D	-2	HIS	-	expression tag	UNP A0A5M5BWR5
D	-1	SER	-	expression tag	UNP A0A5M5BWR5
D	0	SER	-	expression tag	UNP A0A5M5BWR5
D	1	GLY	-	expression tag	UNP A0A5M5BWR5
D	2	LEU	-	expression tag	UNP A0A5M5BWR5
D	3	VAL	-	expression tag	UNP A0A5M5BWR5
D	4	PRO	-	expression tag	UNP A0A5M5BWR5
D	5	ARG	-	expression tag	UNP A0A5M5BWR5
D	6	GLY	-	expression tag	UNP A0A5M5BWR5
D	7	SER	-	expression tag	UNP A0A5M5BWR5
D	8	HIS	-	expression tag	UNP A0A5M5BWR5
D	9	MET	-	expression tag	UNP A0A5M5BWR5
D	10	ALA	-	expression tag	UNP A0A5M5BWR5
D	11	SER	-	expression tag	UNP A0A5M5BWR5

- Molecule 2 is an oligosaccharide called alpha-L-gulopyranuronic acid-(1-4)-alpha-L-gulopyranuronic acid-(1-4)-alpha-L-gulopyranuronic acid-(1-4)-alpha-L-gulopyranuronic acid.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	E	4	Total	C	O	0	0	0
			49	24	25			
2	F	4	Total	C	O	0	0	0
			49	24	25			

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	G	4	Total	C	O	0	0	0
			49	24	25			
2	H	4	Total	C	O	0	0	0
			49	24	25			

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

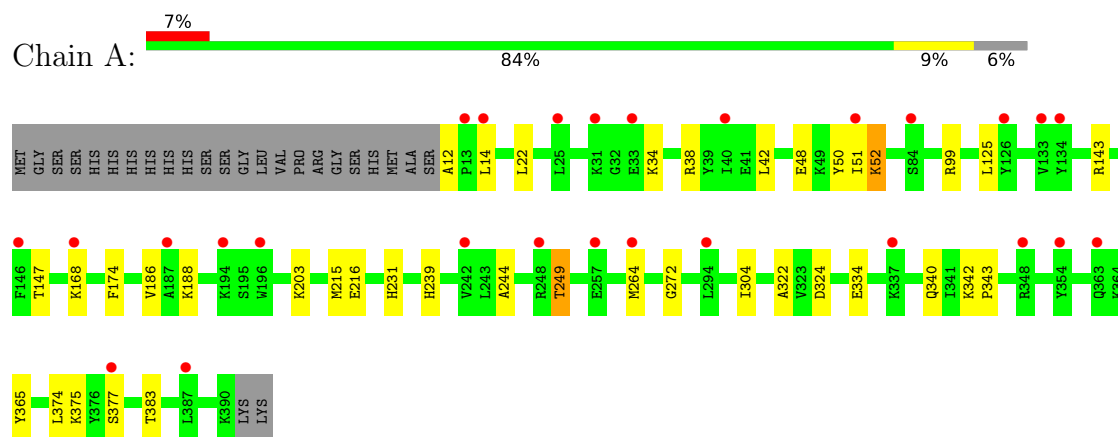
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	193	Total	O	0	0
			193	193		
4	B	201	Total	O	0	0
			201	201		
4	C	133	Total	O	0	0
			133	133		
4	D	123	Total	O	0	0
			123	123		

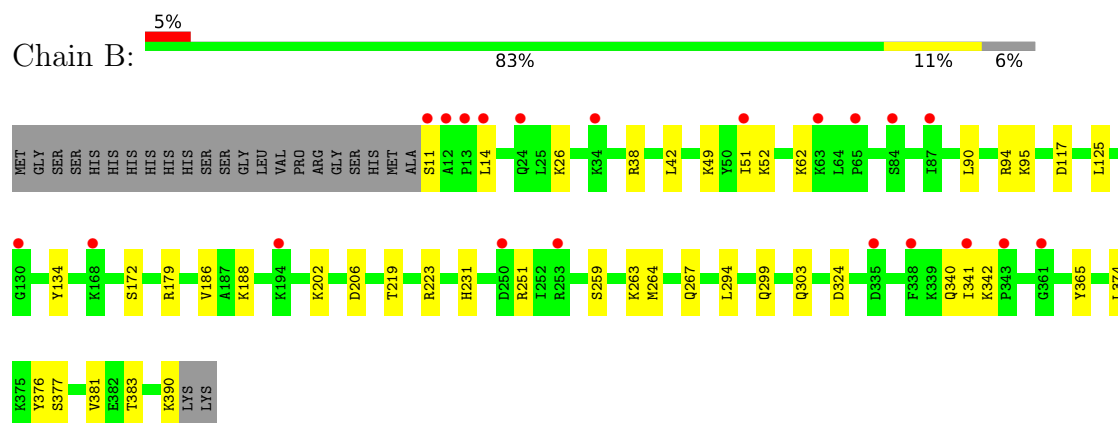
3 Residue-property plots [i](#)

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.

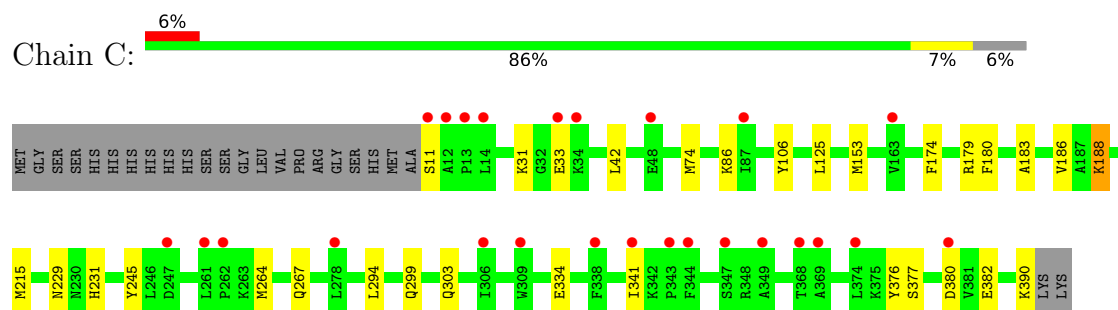
- Molecule 1: Alginate lyase family protein



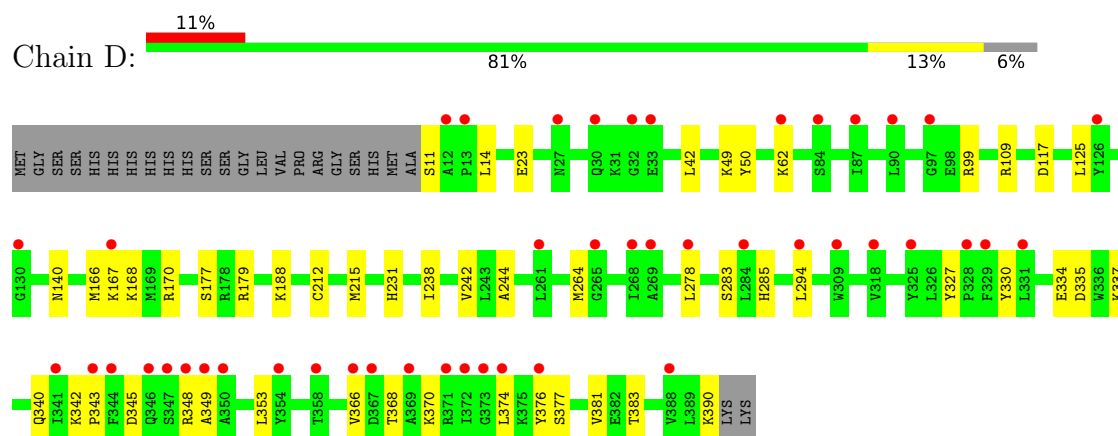
- Molecule 1: Alginate lyase family protein



- Molecule 1: Alginate lyase family protein



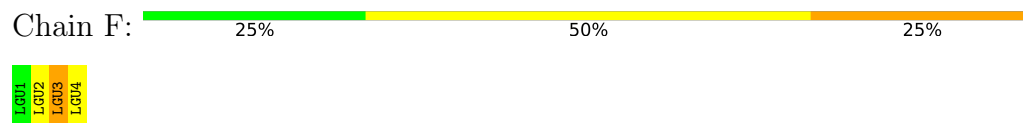
- Molecule 1: Alginate lyase family protein



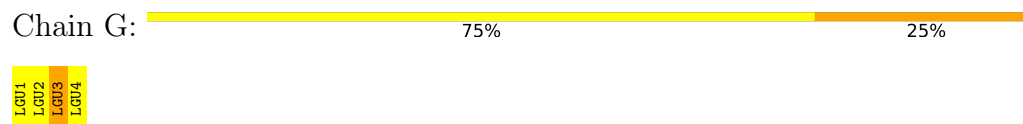
- Molecule 2: alpha-L-gulopyranuronic acid-(1-4)-alpha-L-gulopyranuronic acid-(1-4)-alpha-L-gulopyranuronic acid-(1-4)-alpha-L-gulopyranuronic acid



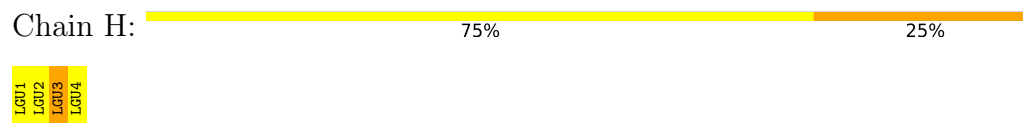
- Molecule 2: alpha-L-gulopyranuronic acid-(1-4)-alpha-L-gulopyranuronic acid-(1-4)-alpha-L-gulopyranuronic acid-(1-4)-alpha-L-gulopyranuronic acid



- Molecule 2: alpha-L-gulopyranuronic acid-(1-4)-alpha-L-gulopyranuronic acid-(1-4)-alpha-L-gulopyranuronic acid-(1-4)-alpha-L-gulopyranuronic acid



- Molecule 2: alpha-L-gulopyranuronic acid-(1-4)-alpha-L-gulopyranuronic acid-(1-4)-alpha-L-gulopyranuronic acid-(1-4)-alpha-L-gulopyranuronic acid



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	197.21Å 88.85Å 146.58Å 90.00° 120.36° 90.00°	Depositor
Resolution (Å)	42.54 – 2.05 42.54 – 2.05	Depositor EDS
% Data completeness (in resolution range)	98.1 (42.54-2.05) 98.1 (42.54-2.05)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.44 (at 2.05Å)	Xtriage
Refinement program	PHENIX 1.21_5207	Depositor
R, R_{free}	0.234 , 0.267 0.234 , 0.268	Depositor DCC
R_{free} test set	6794 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	47.3	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 38.4	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	0.94	EDS
Total number of atoms	13069	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.92% 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: SO4, LGU

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	0.38	0/3124	0.57	0/4234
1	B	0.38	0/3130	0.57	0/4242
1	C	0.35	0/3130	0.54	0/4242
1	D	0.33	0/3139	0.51	0/4253
All	All	0.36	0/12523	0.55	0/16971

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3046	0	3017	36	0
1	B	3052	0	3022	31	0
1	C	3052	0	3022	21	0
1	D	3058	0	3035	37	0
2	E	49	0	27	2	0
2	F	49	0	27	1	0
2	G	49	0	27	1	0
2	H	49	0	27	1	0
3	A	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	5	0	0	0	0
3	C	5	0	0	0	0
4	A	193	0	0	8	1
4	B	201	0	0	7	1
4	C	133	0	0	3	0
4	D	123	0	0	2	0
All	All	13069	0	12204	123	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 5.

All (123) 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:C:267:GLN:NE2	4:C:501:HOH:O	2.01	0.92
1:D:345:ASP:H	1:D:348:ARG:HH21	1.14	0.92
1:D:212:CYS:HA	1:D:215:MET:HE3	1.52	0.91
1:D:345:ASP:HB3	1:D:348:ARG:HE	1.45	0.80
1:C:31:LYS:NZ	1:C:33:GLU:OE2	2.15	0.78
1:B:324:ASP:OD1	1:B:365:TYR:OH	2.07	0.72
1:C:229:ASN:HA	1:C:267:GLN:NE2	2.07	0.69
1:D:349:ALA:HB1	1:D:353:LEU:HD23	1.75	0.67
1:A:324:ASP:OD1	1:A:365:TYR:OH	2.12	0.67
1:B:179:ARG:O	4:B:501:HOH:O	2.12	0.66
1:B:49:LYS:NZ	4:B:505:HOH:O	2.28	0.66
1:A:38:ARG:NH1	4:A:504:HOH:O	2.31	0.63
1:D:345:ASP:HB3	1:D:348:ARG:NE	2.13	0.63
1:A:375:LYS:H	1:A:375:LYS:CD	2.12	0.62
1:B:179:ARG:NH1	4:B:502:HOH:O	2.21	0.62
1:D:215:MET:HE1	1:D:242:VAL:HG21	1.82	0.62
1:B:38:ARG:O	1:B:38:ARG:HD3	1.99	0.61
1:A:12:ALA:N	4:A:505:HOH:O	2.33	0.60
1:A:34:LYS:O	1:A:38:ARG:HG3	2.01	0.60
1:C:86:LYS:NZ	4:C:503:HOH:O	2.29	0.60
1:B:263:LYS:NZ	4:B:508:HOH:O	2.32	0.59
1:A:375:LYS:H	1:A:375:LYS:CE	2.16	0.58
1:B:231:HIS:NE2	2:F:3:LGU:H5	2.19	0.58
1:A:168:LYS:HE3	4:A:639:HOH:O	2.03	0.58
1:A:375:LYS:HD3	1:A:375:LYS:N	2.20	0.57
1:C:183:ALA:O	1:C:186:VAL:HG12	2.05	0.57
1:D:188:LYS:NZ	1:D:244:ALA:O	2.35	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:375:LYS:H	1:A:375:LYS:HE2	1.69	0.57
1:A:42:LEU:HD21	1:A:377:SER:HA	1.87	0.56
1:D:14:LEU:HG	1:D:244:ALA:HB1	1.85	0.56
1:D:264:MET:HE2	1:D:294:LEU:HD13	1.88	0.56
1:D:370:LYS:O	4:D:401:HOH:O	2.18	0.56
1:A:147:THR:HG21	4:A:550:HOH:O	2.06	0.55
1:D:166:MET:SD	1:D:170:ARG:HD3	2.47	0.55
1:A:147:THR:HG23	4:A:642:HOH:O	2.06	0.55
1:D:231:HIS:NE2	2:H:3:LGU:H5	2.22	0.55
1:C:42:LEU:HD21	1:C:377:SER:HA	1.88	0.54
1:D:327:TYR:OH	1:D:368:THR:OG1	2.23	0.54
1:D:374:LEU:HD13	1:D:383:THR:HB	1.88	0.54
1:A:99:ARG:HH11	1:A:99:ARG:HG3	1.73	0.54
1:C:174:PHE:CZ	1:C:215:MET:HG2	2.42	0.54
1:B:51:ILE:HG22	1:B:52:LYS:HD3	1.89	0.54
1:A:48:GLU:HG3	4:A:650:HOH:O	2.07	0.54
1:C:153:MET:HE1	1:C:180:PHE:CZ	2.43	0.54
1:C:188:LYS:HE3	1:C:245:TYR:O	2.09	0.52
1:A:375:LYS:CD	1:A:375:LYS:N	2.72	0.52
1:A:249:THR:HG22	1:A:304:ILE:HG21	1.91	0.52
1:B:202:LYS:HE3	1:B:206:ASP:OD2	2.09	0.52
1:B:264:MET:SD	1:B:294:LEU:HD13	2.50	0.52
1:A:14:LEU:HG	1:A:244:ALA:HB1	1.92	0.51
1:A:375:LYS:H	1:A:375:LYS:HD3	1.75	0.51
1:D:334:GLU:OE2	1:D:334:GLU:N	2.25	0.51
1:A:174:PHE:CZ	1:A:215:MET:HG2	2.45	0.51
1:D:349:ALA:HB1	1:D:353:LEU:CD2	2.40	0.51
1:A:374:LEU:HD13	1:A:383:THR:HB	1.92	0.50
1:B:374:LEU:HD13	1:B:383:THR:HB	1.93	0.49
1:C:380:ASP:HB3	1:C:382:GLU:OE2	2.11	0.49
1:C:376:TYR:HA	1:C:390:LYS:HB3	1.94	0.49
1:D:330:TYR:CD2	1:D:353:LEU:HD21	2.47	0.49
1:A:188:LYS:HG2	4:A:595:HOH:O	2.11	0.49
1:B:299:GLN:O	1:B:303:GLN:NE2	2.43	0.49
1:D:376:TYR:HA	1:D:390:LYS:HB3	1.93	0.49
1:B:14:LEU:O	1:B:188:LYS:HE3	2.15	0.47
1:A:99:ARG:NH1	2:E:4:LGU:O6B	2.47	0.47
1:A:168:LYS:HE2	1:A:168:LYS:HB3	1.85	0.47
1:C:264:MET:SD	1:C:294:LEU:HD13	2.55	0.47
1:A:264:MET:HE2	1:A:264:MET:HB2	1.62	0.46
1:B:376:TYR:HA	1:B:390:LYS:HB3	1.96	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:99:ARG:HH11	1:D:109:ARG:NH2	2.13	0.46
1:D:335:ASP:O	1:D:337:LYS:HE3	2.15	0.46
1:B:259:SER:O	1:B:263:LYS:HB2	2.15	0.46
1:D:177:SER:HB2	1:D:238:ILE:HD13	1.97	0.46
1:D:340:GLN:HG3	1:D:342:LYS:O	2.16	0.46
1:B:94:ARG:O	1:B:95:LYS:HE3	2.16	0.46
1:D:334:GLU:H	1:D:334:GLU:CD	2.15	0.45
1:A:216:GLU:OE2	1:A:239:HIS:NE2	2.38	0.45
1:B:263:LYS:HG3	4:B:604:HOH:O	2.16	0.45
1:B:340:GLN:HG3	1:B:342:LYS:O	2.17	0.45
1:D:285:HIS:ND1	1:D:348:ARG:NH1	2.64	0.45
1:B:125:LEU:HD23	1:B:125:LEU:HA	1.78	0.45
1:C:179:ARG:O	4:C:502:HOH:O	2.21	0.45
1:A:51:ILE:HG22	1:A:52:LYS:HE3	1.97	0.45
1:D:167:LYS:HD3	1:D:168:LYS:H	1.82	0.44
1:D:117:ASP:HA	1:D:381:VAL:HG21	1.98	0.44
1:B:51:ILE:HD11	1:B:134:TYR:CE1	2.52	0.44
1:D:168:LYS:NZ	1:D:170:ARG:HG2	2.33	0.44
1:D:179:ARG:HD2	4:D:458:HOH:O	2.16	0.44
1:A:334:GLU:OE1	1:A:334:GLU:N	2.32	0.44
1:D:42:LEU:HD21	1:D:377:SER:HA	1.99	0.43
1:D:278:LEU:HA	1:D:283:SER:HB2	1.99	0.43
1:A:147:THR:OG1	1:A:203:LYS:HB3	2.19	0.43
1:A:340:GLN:HG3	1:A:342:LYS:O	2.18	0.43
1:B:42:LEU:HD21	1:B:377:SER:HA	2.00	0.43
1:D:125:LEU:HD23	1:D:125:LEU:HA	1.85	0.43
1:B:219:THR:O	1:B:223:ARG:HG3	2.18	0.43
1:B:342:LYS:HB3	1:B:342:LYS:HE2	1.77	0.43
1:A:50:TYR:OH	4:A:501:HOH:O	2.18	0.43
1:A:272:GLY:O	1:A:322:ALA:HA	2.19	0.43
1:B:117:ASP:HA	1:B:381:VAL:HG21	2.01	0.43
1:B:264:MET:HE3	1:B:264:MET:HA	2.01	0.43
1:C:334:GLU:OE1	1:C:334:GLU:N	2.34	0.43
1:D:49:LYS:HE2	1:D:50:TYR:CZ	2.54	0.43
1:C:299:GLN:O	1:C:303:GLN:NE2	2.51	0.42
1:A:143:ARG:HA	1:A:147:THR:HG22	2.02	0.42
1:A:343:PRO:HD3	1:B:341:ILE:HB	2.01	0.42
1:D:11:SER:HA	1:D:23:GLU:OE2	2.19	0.42
1:B:263:LYS:O	1:B:267:GLN:HG3	2.20	0.42
1:C:74:MET:HE3	1:C:106:TYR:HE2	1.84	0.42
1:A:22:LEU:HD23	1:A:22:LEU:HA	1.86	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:366:VAL:HG12	1:D:370:LYS:NZ	2.35	0.42
1:B:179:ARG:HD2	4:B:502:HOH:O	2.18	0.41
1:A:231:HIS:NE2	2:E:3:LGU:H5	2.34	0.41
1:A:125:LEU:HD23	1:A:125:LEU:HA	1.78	0.41
1:C:125:LEU:HD23	1:C:125:LEU:HA	1.74	0.41
1:B:26:LYS:NZ	4:B:516:HOH:O	2.47	0.41
1:D:167:LYS:HD3	1:D:168:LYS:N	2.34	0.41
1:C:341:ILE:HB	1:D:343:PRO:HD3	2.02	0.41
1:B:38:ARG:HD3	1:B:38:ARG:C	2.46	0.41
1:C:74:MET:HE3	1:C:106:TYR:CE2	2.56	0.41
1:C:231:HIS:NE2	2:G:3:LGU:H5	2.36	0.41
1:B:90:LEU:HD23	1:B:90:LEU:HA	1.90	0.41
1:D:212:CYS:HA	1:D:215:MET:CE	2.38	0.40
1:C:174:PHE:CZ	1:C:215:MET:HE3	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)
4:A:559:HOH:O	4:B:612:HOH:O[4_455]	2.08	0.12

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	377/404 (93%)	370 (98%)	7 (2%)	0	100	100
1	B	378/404 (94%)	370 (98%)	8 (2%)	0	100	100
1	C	378/404 (94%)	370 (98%)	8 (2%)	0	100	100
1	D	379/404 (94%)	371 (98%)	8 (2%)	0	100	100
All	All	1512/1616 (94%)	1481 (98%)	31 (2%)	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	323/344 (94%)	320 (99%)	3 (1%)	75	77
1	B	324/344 (94%)	319 (98%)	5 (2%)	60	60
1	C	324/344 (94%)	322 (99%)	2 (1%)	84	86
1	D	325/344 (94%)	322 (99%)	3 (1%)	75	77
All	All	1296/1376 (94%)	1283 (99%)	13 (1%)	75	74

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	LYS
1	A	186	VAL
1	A	249	THR
1	B	11	SER
1	B	62	LYS
1	B	172	SER
1	B	186	VAL
1	B	251	ARG
1	C	11	SER
1	C	188	LYS
1	D	62[A]	LYS
1	D	62[B]	LYS
1	D	140	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	276	GLN
1	C	24	GLN
1	C	267	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

16 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	LGU	E	1	2	13,13,13	0.96	1 (7%)	18,19,19	1.03	1 (5%)
2	LGU	E	2	2	12,12,13	1.09	1 (8%)	14,17,19	1.66	4 (28%)
2	LGU	E	3	2	12,12,13	1.19	2 (16%)	14,17,19	1.55	2 (14%)
2	LGU	E	4	2	12,12,13	1.10	1 (8%)	14,17,19	1.08	2 (14%)
2	LGU	F	1	2	13,13,13	0.93	0	18,19,19	0.95	0
2	LGU	F	2	2	12,12,13	1.01	1 (8%)	14,17,19	1.11	0
2	LGU	F	3	2	12,12,13	1.13	0	14,17,19	1.56	2 (14%)
2	LGU	F	4	2	12,12,13	1.08	2 (16%)	14,17,19	1.16	1 (7%)
2	LGU	G	1	2	13,13,13	0.90	1 (7%)	18,19,19	1.08	0
2	LGU	G	2	2	12,12,13	1.02	1 (8%)	14,17,19	1.35	2 (14%)
2	LGU	G	3	2	12,12,13	1.20	0	14,17,19	1.63	2 (14%)
2	LGU	G	4	2	12,12,13	1.10	1 (8%)	14,17,19	1.39	2 (14%)
2	LGU	H	1	2	13,13,13	0.94	1 (7%)	18,19,19	1.04	1 (5%)
2	LGU	H	2	2	12,12,13	1.06	0	14,17,19	1.45	3 (21%)
2	LGU	H	3	2	12,12,13	1.19	1 (8%)	14,17,19	1.07	1 (7%)
2	LGU	H	4	2	12,12,13	1.04	1 (8%)	14,17,19	0.88	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	LGU	E	1	2	-	0/4/24/24	0/1/1/1
2	LGU	E	2	2	-	0/4/21/24	0/1/1/1
2	LGU	E	3	2	-	2/4/21/24	0/1/1/1
2	LGU	E	4	2	-	0/4/21/24	0/1/1/1
2	LGU	F	1	2	-	0/4/24/24	0/1/1/1
2	LGU	F	2	2	-	0/4/21/24	0/1/1/1
2	LGU	F	3	2	-	0/4/21/24	0/1/1/1
2	LGU	F	4	2	-	0/4/21/24	0/1/1/1
2	LGU	G	1	2	-	0/4/24/24	0/1/1/1
2	LGU	G	2	2	-	2/4/21/24	0/1/1/1
2	LGU	G	3	2	-	1/4/21/24	0/1/1/1
2	LGU	G	4	2	-	0/4/21/24	0/1/1/1
2	LGU	H	1	2	-	0/4/24/24	0/1/1/1
2	LGU	H	2	2	-	0/4/21/24	0/1/1/1
2	LGU	H	3	2	-	0/4/21/24	0/1/1/1
2	LGU	H	4	2	-	0/4/21/24	0/1/1/1

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	4	LGU	O6B-C6	-2.43	1.22	1.30
2	H	4	LGU	O6B-C6	-2.16	1.23	1.30
2	G	2	LGU	O6B-C6	-2.15	1.23	1.30
2	E	3	LGU	C1-C2	2.13	1.57	1.52
2	E	3	LGU	O6B-C6	-2.07	1.23	1.30
2	E	1	LGU	O6B-C6	-2.05	1.23	1.30
2	F	2	LGU	O6B-C6	-2.03	1.23	1.30
2	F	4	LGU	O6B-C6	-2.03	1.23	1.30
2	H	1	LGU	O6B-C6	-2.03	1.23	1.30
2	G	4	LGU	C1-C2	2.02	1.56	1.52
2	E	2	LGU	O6B-C6	-2.02	1.23	1.30
2	H	3	LGU	C5-C6	2.01	1.57	1.53
2	G	1	LGU	O6B-C6	-2.01	1.24	1.30
2	F	4	LGU	C1-C2	2.00	1.56	1.52

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	3	LGU	O4-C4-C5	3.63	117.89	109.74
2	G	4	LGU	O2-C2-C1	3.25	115.81	109.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	3	LGU	O4-C4-C5	3.19	116.89	109.74
2	G	3	LGU	C1-C2-C3	3.12	113.51	109.67
2	E	2	LGU	O2-C2-C1	2.88	115.04	109.15
2	F	4	LGU	O2-C2-C1	2.83	114.94	109.15
2	F	3	LGU	O3-C3-C2	2.71	115.19	109.99
2	F	3	LGU	O4-C4-C5	2.63	115.64	109.74
2	G	4	LGU	O2-C2-C3	-2.45	105.23	110.14
2	E	2	LGU	C2-C3-C4	-2.43	106.69	110.89
2	H	2	LGU	O2-C2-C1	2.38	114.03	109.15
2	E	4	LGU	C1-C2-C3	2.29	112.48	109.67
2	E	3	LGU	O2-C2-C1	2.25	113.76	109.15
2	E	2	LGU	O6A-C6-C5	-2.23	112.65	120.81
2	H	3	LGU	O2-C2-C1	2.21	113.68	109.15
2	E	2	LGU	O2-C2-C3	-2.16	105.82	110.14
2	E	4	LGU	O2-C2-C1	2.16	113.56	109.15
2	H	2	LGU	O6B-C6-C5	2.11	121.39	113.65
2	H	2	LGU	O6A-C6-C5	-2.03	113.37	120.81
2	G	2	LGU	O6A-C6-C5	-2.03	113.37	120.81
2	E	1	LGU	C4-C3-C2	2.02	114.35	110.82
2	H	1	LGU	C1-C2-C3	2.01	114.49	110.31
2	G	2	LGU	C1-C2-C3	2.01	112.14	109.67

There are no chirality outliers.

All (5) torsion outliers are listed below:

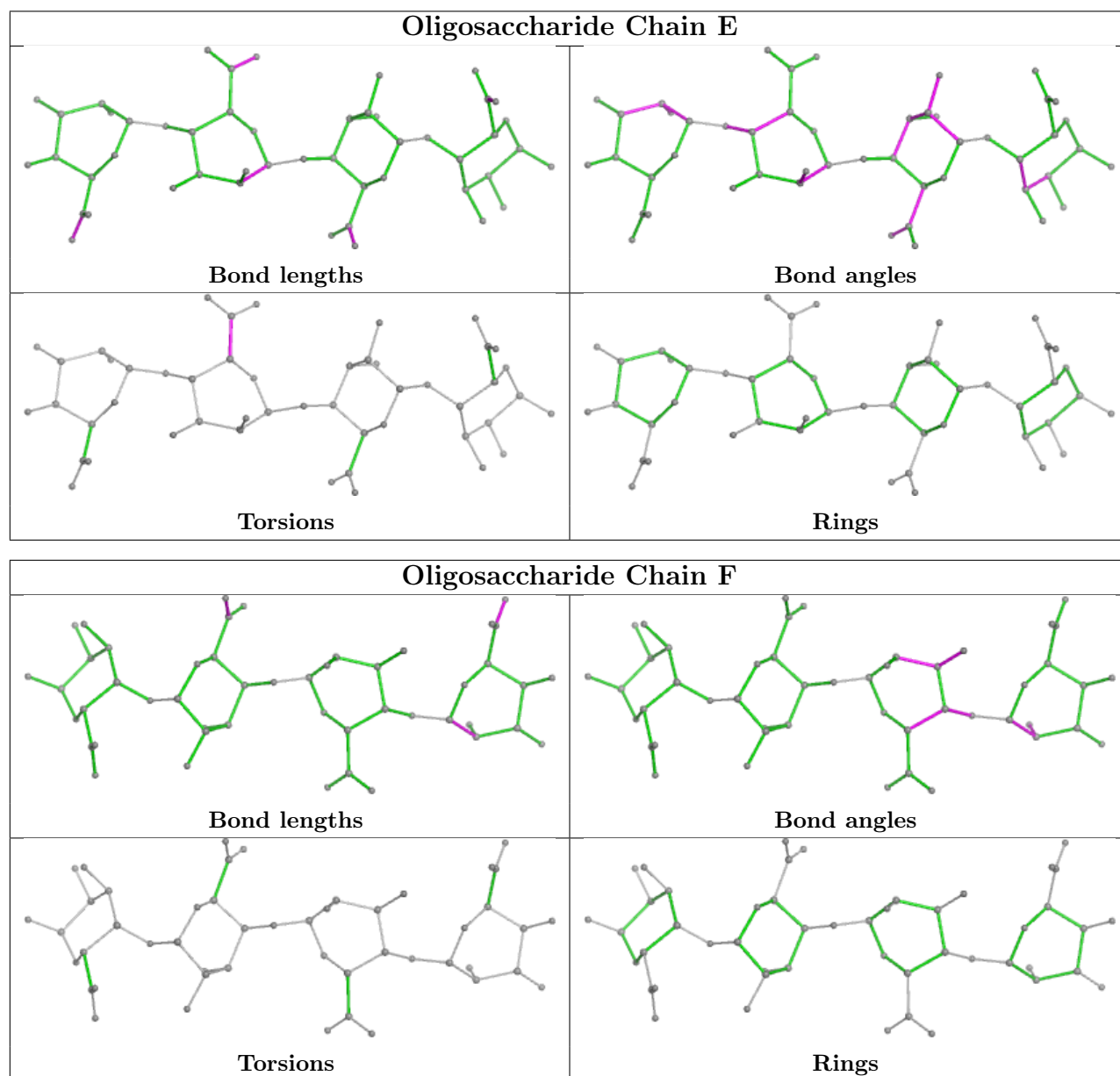
Mol	Chain	Res	Type	Atoms
2	E	3	LGU	C4-C5-C6-O6B
2	E	3	LGU	O5-C5-C6-O6B
2	G	2	LGU	O5-C5-C6-O6B
2	G	2	LGU	O5-C5-C6-O6A
2	G	3	LGU	C4-C5-C6-O6B

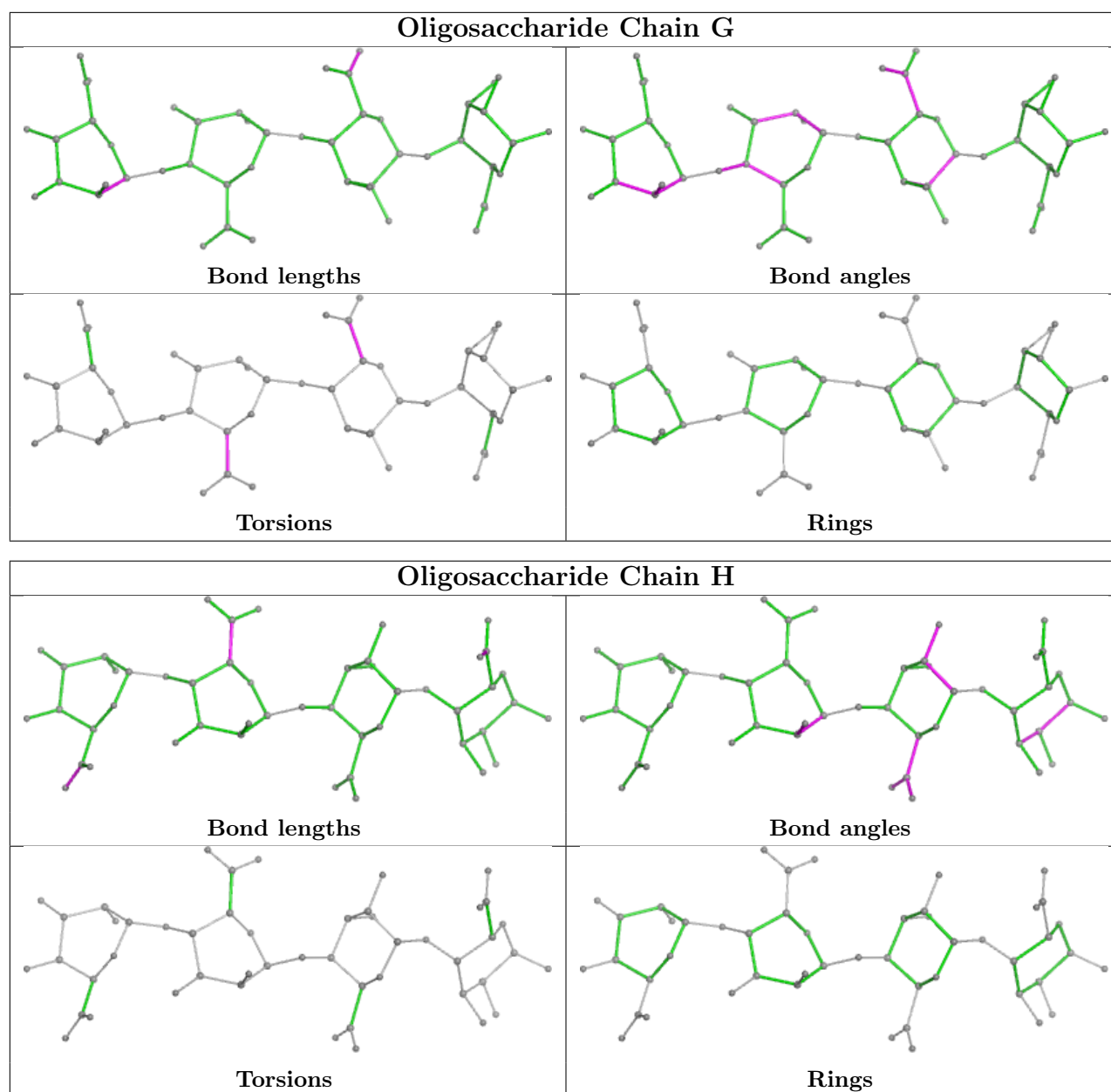
There are no ring outliers.

5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	3	LGU	1	0
2	E	4	LGU	1	0
2	H	3	LGU	1	0
2	F	3	LGU	1	0
2	E	3	LGU	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](#)

3 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	SO4	B	401	-	4,4,4	0.71	0	6,6,6	0.19	0
3	SO4	A	401	-	4,4,4	0.64	0	6,6,6	0.11	0
3	SO4	C	401	-	4,4,4	0.66	0	6,6,6	0.16	0

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 [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	379/404 (93%)	0.77	27 (7%)	23 25	31, 44, 59, 70	0
1	B	380/404 (94%)	0.77	21 (5%)	32 33	31, 44, 58, 69	0
1	C	380/404 (94%)	0.80	25 (6%)	26 27	33, 48, 62, 75	0
1	D	380/404 (94%)	0.99	46 (12%)	10 10	32, 52, 70, 76	1 (0%)
All	All	1519/1616 (93%)	0.83	119 (7%)	20 22	31, 46, 64, 76	1 (0%)

All (119) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	12	ALA	4.2
1	D	32	GLY	3.9
1	D	269	ALA	3.7
1	D	278	LEU	3.5
1	C	11	SER	3.4
1	D	348	ARG	3.4
1	A	337	LYS	3.4
1	D	268	ILE	3.2
1	D	347	SER	3.2
1	B	63	LYS	3.1
1	B	13	PRO	3.1
1	D	325	TYR	3.1
1	B	12	ALA	3.0
1	C	247	ASP	3.0
1	D	30	GLN	3.0
1	B	11	SER	3.0
1	B	84	SER	2.9
1	C	344	PHE	2.9
1	C	341	ILE	2.9
1	C	338	PHE	2.9
1	D	27	ASN	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	250	ASP	2.9
1	D	344	PHE	2.9
1	D	341	ILE	2.9
1	D	366	VAL	2.8
1	B	168	LYS	2.8
1	B	65	PRO	2.8
1	A	126	TYR	2.8
1	D	167	LYS	2.7
1	D	265	GLY	2.7
1	D	374	LEU	2.7
1	A	31	LYS	2.7
1	B	34	LYS	2.7
1	C	374	LEU	2.7
1	A	264	MET	2.7
1	B	130	GLY	2.7
1	B	341	ILE	2.7
1	B	87	ILE	2.6
1	C	13	PRO	2.6
1	D	126	TYR	2.6
1	C	14	LEU	2.6
1	C	261	LEU	2.6
1	A	13	PRO	2.6
1	D	309	TRP	2.6
1	B	343	PRO	2.5
1	C	349	ALA	2.5
1	C	34	LYS	2.5
1	A	387	LEU	2.5
1	B	253	ARG	2.4
1	D	331	LEU	2.4
1	D	371	ARG	2.4
1	C	309	TRP	2.4
1	A	242	VAL	2.4
1	D	318	VAL	2.4
1	D	130	GLY	2.4
1	C	380	ASP	2.4
1	A	33	GLU	2.4
1	A	187	ALA	2.4
1	C	48	GLU	2.4
1	A	348	ARG	2.4
1	B	14	LEU	2.4
1	D	294	LEU	2.4
1	C	368	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	62[A]	LYS	2.4
1	A	51	ILE	2.4
1	C	33	GLU	2.3
1	C	369	ALA	2.3
1	C	343	PRO	2.3
1	B	335	ASP	2.3
1	C	306	ILE	2.3
1	A	133	VAL	2.3
1	C	163	VAL	2.3
1	A	40	ILE	2.3
1	D	90	LEU	2.3
1	D	284	LEU	2.3
1	A	363	GLN	2.3
1	B	361	GLY	2.2
1	D	372	ILE	2.2
1	D	358	THR	2.2
1	A	257	GLU	2.2
1	D	97	GLY	2.2
1	D	373	GLY	2.2
1	B	338	PHE	2.2
1	D	329	PHE	2.2
1	D	87	ILE	2.2
1	A	14	LEU	2.2
1	A	354	TYR	2.2
1	D	354	TYR	2.2
1	D	376	TYR	2.2
1	D	13	PRO	2.2
1	B	24	GLN	2.2
1	A	194	LYS	2.2
1	A	146	PHE	2.1
1	D	367	ASP	2.1
1	A	25	LEU	2.1
1	C	278	LEU	2.1
1	D	261	LEU	2.1
1	A	84	SER	2.1
1	D	84	SER	2.1
1	C	12	ALA	2.1
1	A	134	TYR	2.1
1	A	196	TRP	2.1
1	B	194	LYS	2.1
1	A	377	SER	2.1
1	D	369	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	262	PRO	2.1
1	D	388	VAL	2.1
1	A	248	ARG	2.1
1	A	294	LEU	2.1
1	C	347	SER	2.1
1	B	51	ILE	2.1
1	D	343	PRO	2.0
1	A	168	LYS	2.0
1	D	346	GLN	2.0
1	C	87	ILE	2.0
1	D	33	GLU	2.0
1	D	328	PRO	2.0
1	D	349	ALA	2.0
1	D	350	ALA	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.

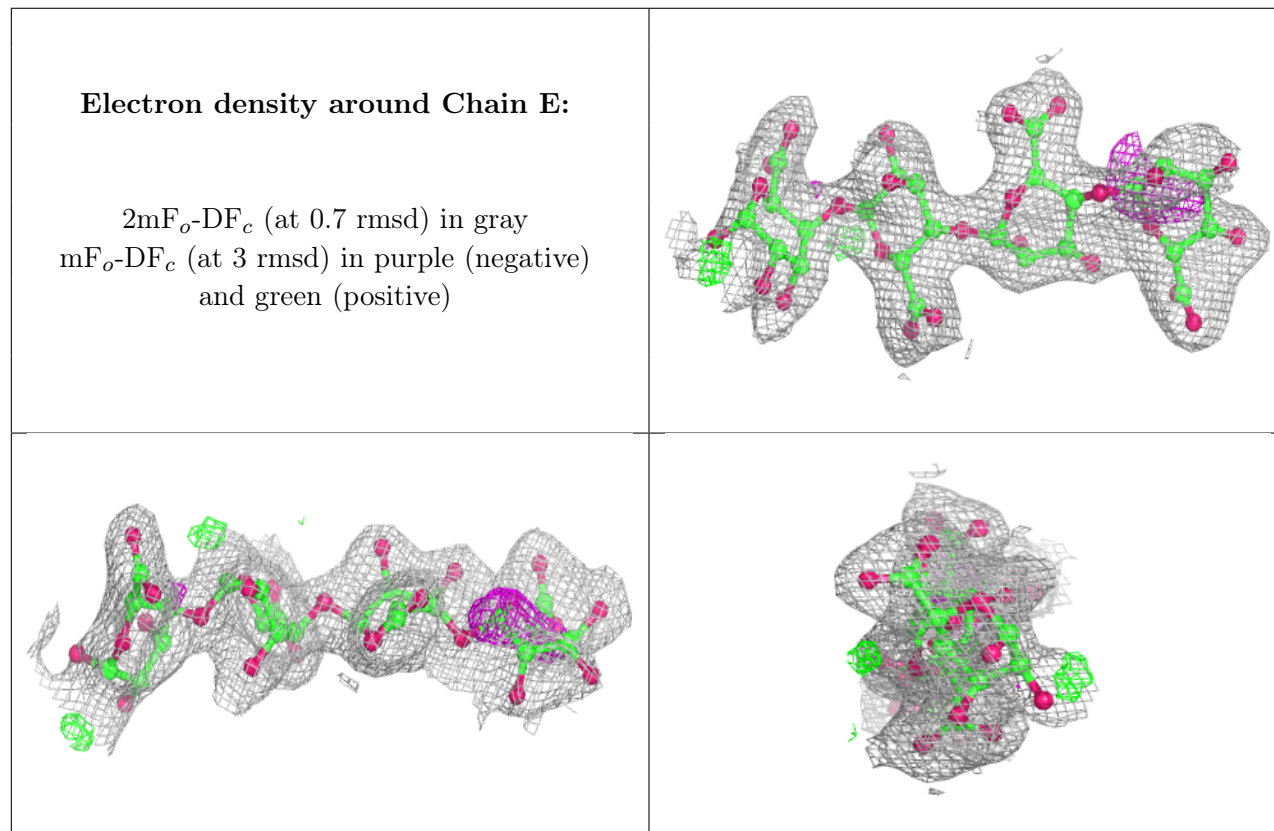
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	LGU	G	1	13/13	0.74	0.13	64,68,72,80	0
2	LGU	H	1	13/13	0.74	0.14	64,72,76,79	0
2	LGU	E	1	13/13	0.80	0.11	53,59,68,69	0
2	LGU	H	4	12/13	0.80	0.15	54,62,64,65	0
2	LGU	E	4	12/13	0.81	0.17	46,56,65,67	0
2	LGU	F	1	13/13	0.82	0.12	58,65,69,69	0
2	LGU	H	2	12/13	0.85	0.11	49,59,64,67	0
2	LGU	G	4	12/13	0.86	0.12	53,56,60,61	0
2	LGU	F	4	12/13	0.88	0.12	47,52,56,59	0
2	LGU	H	3	12/13	0.88	0.12	49,56,59,64	0
2	LGU	G	3	12/13	0.88	0.10	45,51,53,55	0
2	LGU	E	2	12/13	0.91	0.09	38,47,49,53	0
2	LGU	F	3	12/13	0.91	0.10	43,46,47,54	0
2	LGU	G	2	12/13	0.91	0.09	49,56,60,62	0

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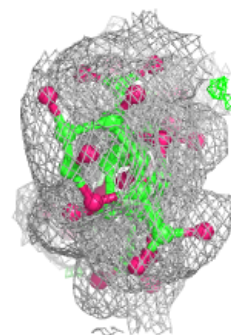
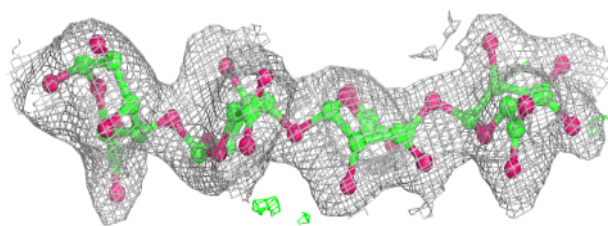
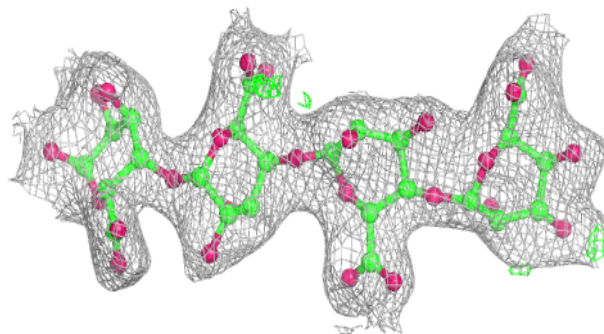
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	LGU	E	3	12/13	0.93	0.08	38,42,46,56	0
2	LGU	F	2	12/13	0.93	0.09	43,48,57,60	0

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

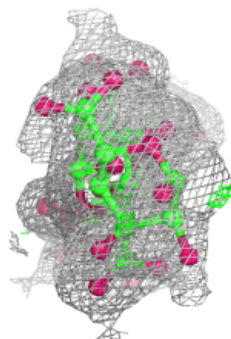
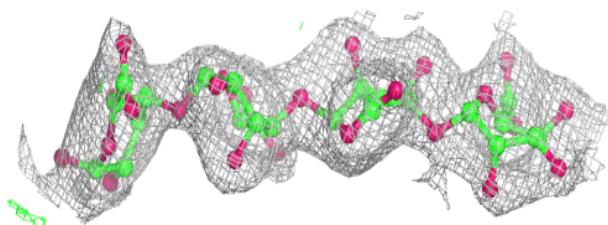
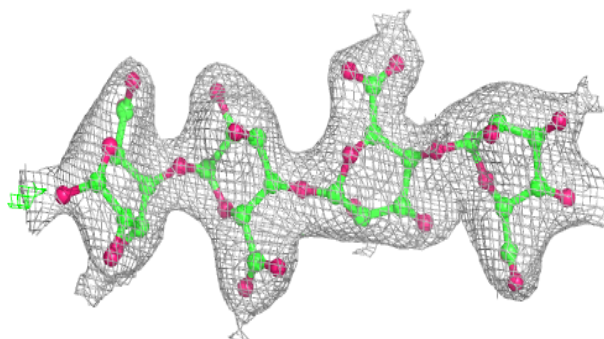


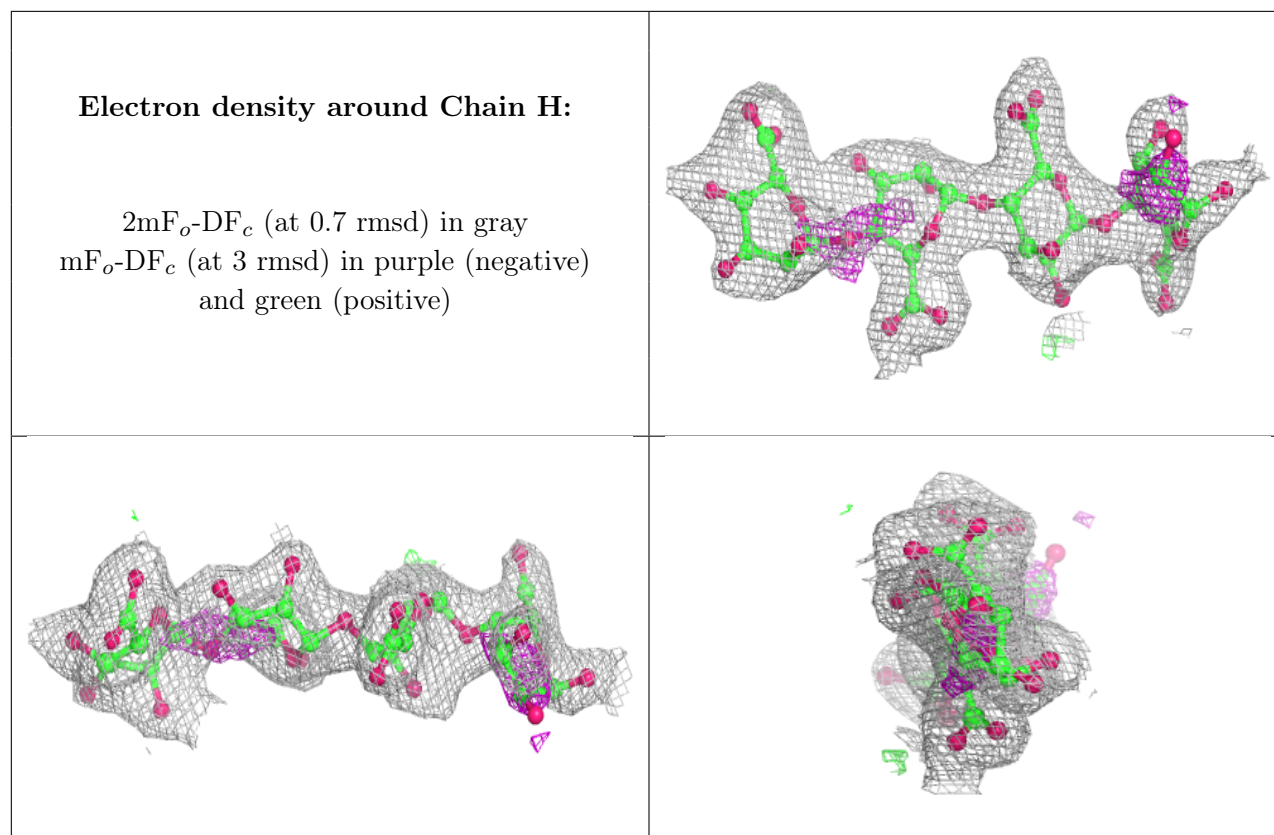
Electron density around Chain F:

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

**Electron density around Chain G:**

$2mF_o - DF_c$ (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(\AA^2)	Q<0.9
3	SO4	B	401	5/5	0.78	0.14	44,50,61,63	5
3	SO4	A	401	5/5	0.85	0.14	46,52,56,61	5
3	SO4	C	401	5/5	0.90	0.11	50,52,57,58	5

6.5 Other polymers [i](#)

There are no such residues in this entry.