



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:28 PM UTC

PDB ID : 8RZK / pdb_00008rzk
Title : The Michaelis complex of ZgGH129 D486N from Zobellia galactanivorans with neo-b/k-oligo-carrageenan tetrasaccharide (beta-kappa neo-oligo-carrageenan DP4).
Authors : Roret, T.; Czjzek, M.; Ficko-Blean, E.
Deposited on : 2024-02-12
Resolution : 2.29 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

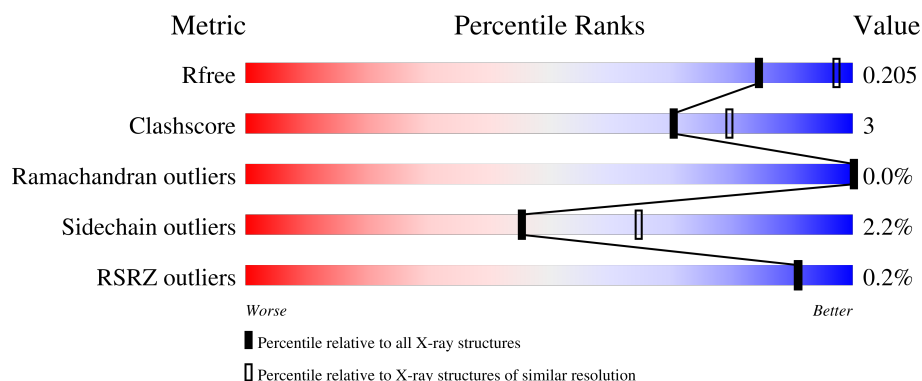
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.29 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	676	
1	B	676	
1	C	676	
1	D	676	
2	E	4	

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

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 22347 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 Conserved hypothetical periplasmic protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	659	Total	C	N	O	S	0	0	0
			5238	3347	889	979	23			
1	B	659	Total	C	N	O	S	0	0	0
			5241	3348	889	981	23			
1	C	659	Total	C	N	O	S	0	0	0
			5234	3345	886	980	23			
1	D	659	Total	C	N	O	S	0	0	0
			5238	3347	888	980	23			

There are 60 discrepancies between the modelled and reference sequences:

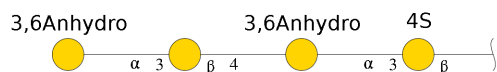
Chain	Residue	Modelled	Actual	Comment	Reference
A	18	MET	-	initiating methionine	UNP G0L004
A	19	GLY	-	expression tag	UNP G0L004
A	20	SER	-	expression tag	UNP G0L004
A	21	SER	-	expression tag	UNP G0L004
A	22	HIS	-	expression tag	UNP G0L004
A	23	HIS	-	expression tag	UNP G0L004
A	24	HIS	-	expression tag	UNP G0L004
A	25	HIS	-	expression tag	UNP G0L004
A	26	HIS	-	expression tag	UNP G0L004
A	27	HIS	-	expression tag	UNP G0L004
A	28	GLY	-	expression tag	UNP G0L004
A	29	SER	-	expression tag	UNP G0L004
A	30	LEU	-	expression tag	UNP G0L004
A	31	ASP	-	expression tag	UNP G0L004
A	486	ASN	ASP	engineered mutation	UNP G0L004
B	18	MET	-	initiating methionine	UNP G0L004
B	19	GLY	-	expression tag	UNP G0L004
B	20	SER	-	expression tag	UNP G0L004
B	21	SER	-	expression tag	UNP G0L004
B	22	HIS	-	expression tag	UNP G0L004
B	23	HIS	-	expression tag	UNP G0L004

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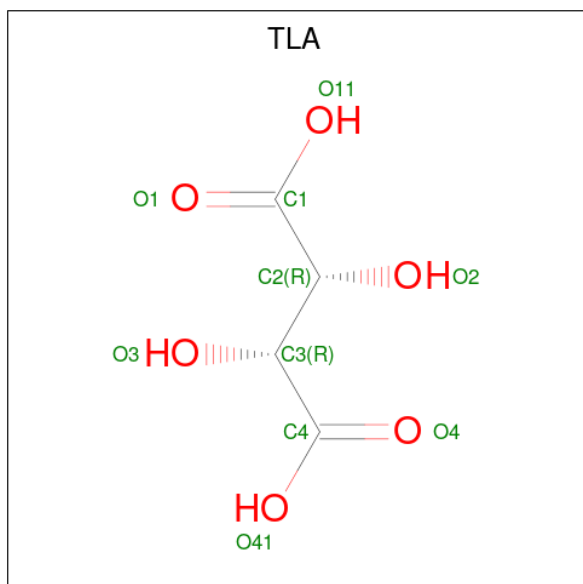
Chain	Residue	Modelled	Actual	Comment	Reference
B	24	HIS	-	expression tag	UNP G0L004
B	25	HIS	-	expression tag	UNP G0L004
B	26	HIS	-	expression tag	UNP G0L004
B	27	HIS	-	expression tag	UNP G0L004
B	28	GLY	-	expression tag	UNP G0L004
B	29	SER	-	expression tag	UNP G0L004
B	30	LEU	-	expression tag	UNP G0L004
B	31	ASP	-	expression tag	UNP G0L004
B	486	ASN	ASP	engineered mutation	UNP G0L004
C	18	MET	-	initiating methionine	UNP G0L004
C	19	GLY	-	expression tag	UNP G0L004
C	20	SER	-	expression tag	UNP G0L004
C	21	SER	-	expression tag	UNP G0L004
C	22	HIS	-	expression tag	UNP G0L004
C	23	HIS	-	expression tag	UNP G0L004
C	24	HIS	-	expression tag	UNP G0L004
C	25	HIS	-	expression tag	UNP G0L004
C	26	HIS	-	expression tag	UNP G0L004
C	27	HIS	-	expression tag	UNP G0L004
C	28	GLY	-	expression tag	UNP G0L004
C	29	SER	-	expression tag	UNP G0L004
C	30	LEU	-	expression tag	UNP G0L004
C	31	ASP	-	expression tag	UNP G0L004
C	486	ASN	ASP	engineered mutation	UNP G0L004
D	18	MET	-	initiating methionine	UNP G0L004
D	19	GLY	-	expression tag	UNP G0L004
D	20	SER	-	expression tag	UNP G0L004
D	21	SER	-	expression tag	UNP G0L004
D	22	HIS	-	expression tag	UNP G0L004
D	23	HIS	-	expression tag	UNP G0L004
D	24	HIS	-	expression tag	UNP G0L004
D	25	HIS	-	expression tag	UNP G0L004
D	26	HIS	-	expression tag	UNP G0L004
D	27	HIS	-	expression tag	UNP G0L004
D	28	GLY	-	expression tag	UNP G0L004
D	29	SER	-	expression tag	UNP G0L004
D	30	LEU	-	expression tag	UNP G0L004
D	31	ASP	-	expression tag	UNP G0L004
D	486	ASN	ASP	engineered mutation	UNP G0L004

- Molecule 2 is an oligosaccharide called 3,6-anhydro-alpha-D-galactopyranose-(1-3)-beta-D-galactopyranose-(1-4)-3,6-anhydro-alpha-D-galactopyranose-(1-3)-4-O-sulfo-beta-D-galactopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	4	Total	C	O	S	0	0	0
			47	24	22	1			
2	F	4	Total	C	O	S	0	0	0
			47	24	22	1			
2	G	4	Total	C	O	S	0	0	0
			47	24	22	1			
2	H	4	Total	C	O	S	0	0	0
			47	24	22	1			

- Molecule 3 is L(+)-TARTARIC ACID (CCD ID: TLA) (formula: C₄H₆O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			10	4	6		
3	A	1	Total	C	O	0	0
			10	4	6		
3	B	1	Total	C	O	0	0
			10	4	6		
3	C	1	Total	C	O	0	0
			10	4	6		
3	D	1	Total	C	O	0	0
			10	4	6		

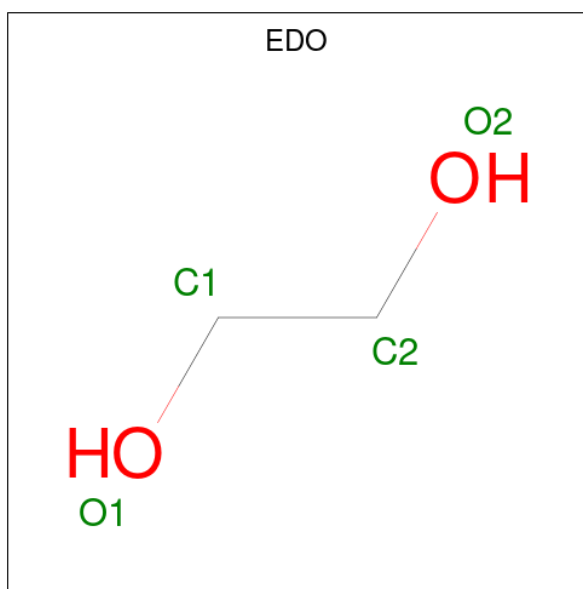
- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Na	0	0
			1	1		
4	B	1	Total	Na	0	0
			1	1		
4	C	1	Total	Na	0	0
			1	1		
4	D	1	Total	Na	0	0
			1	1		

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Cl	0	0
			1	1		
5	B	1	Total	Cl	0	0
			1	1		
5	C	1	Total	Cl	0	0
			1	1		
5	D	1	Total	Cl	0	0
			1	1		

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

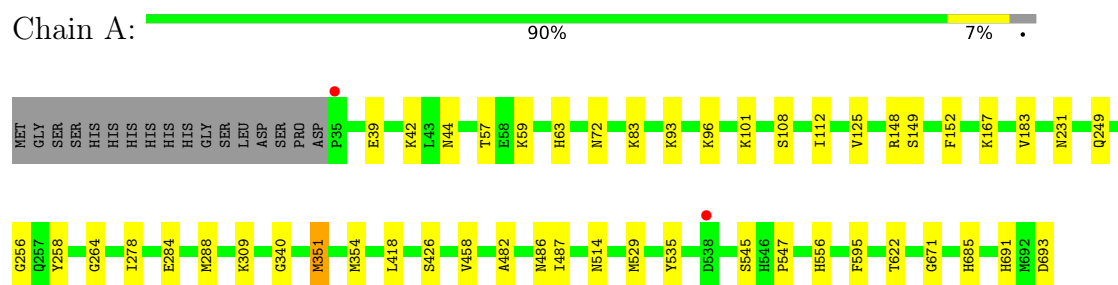
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	310	Total 310	O 310	0	0
7	B	290	Total 290	O 290	0	0
7	C	281	Total 281	O 281	0	0
7	D	265	Total 265	O 265	0	0

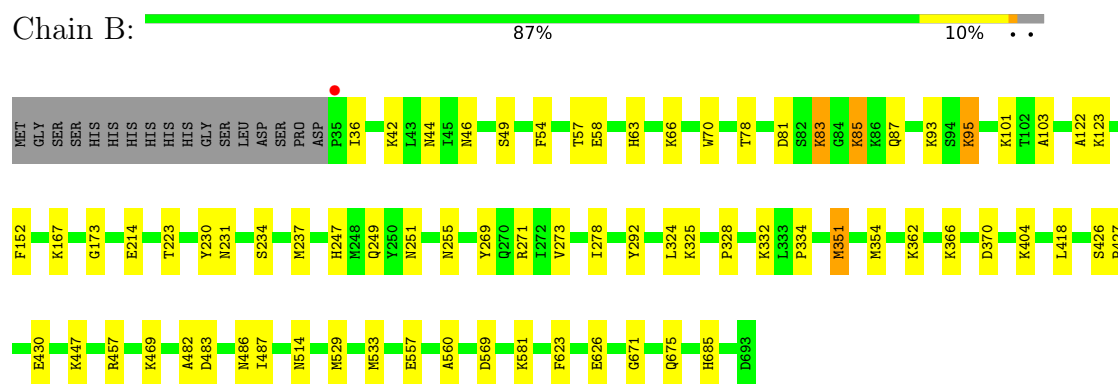
3 Residue-property plots

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.

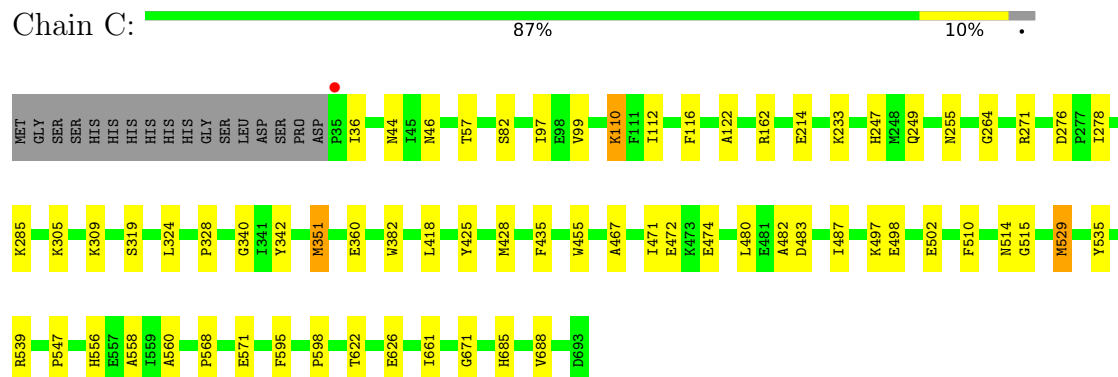
• Molecule 1: Conserved hypothetical periplasmic protein



• Molecule 1: Conserved hypothetical periplasmic protein

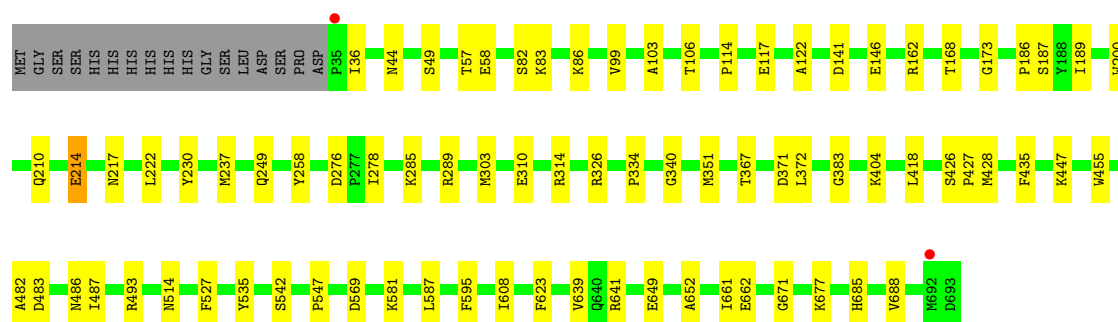


• Molecule 1: Conserved hypothetical periplasmic protein



- Molecule 1: Conserved hypothetical periplasmic protein

Chain D:  86% 12%



- Molecule 2: 3,6-anhydro-alpha-D-galactopyranose-(1-3)-beta-D-galactopyranose-(1-4)-3,6-anhydro-alpha-D-galactopyranose-(1-3)-4-O-sulfo-beta-D-galactopyranose

Chain E:  50% 50%



- Molecule 2: 3,6-anhydro-alpha-D-galactopyranose-(1-3)-beta-D-galactopyranose-(1-4)-3,6-anhydro-alpha-D-galactopyranose-(1-3)-4-O-sulfo-beta-D-galactopyranose

Chain F:  50% 50%



- Molecule 2: 3,6-anhydro-alpha-D-galactopyranose-(1-3)-beta-D-galactopyranose-(1-4)-3,6-anhydro-alpha-D-galactopyranose-(1-3)-4-O-sulfo-beta-D-galactopyranose

Chain G:  75% 25%



- Molecule 2: 3,6-anhydro-alpha-D-galactopyranose-(1-3)-beta-D-galactopyranose-(1-4)-3,6-anhydro-alpha-D-galactopyranose-(1-3)-4-O-sulfo-beta-D-galactopyranose

Chain H:  75% 25%



4 Data and refinement statistics

Property	Value	Source
Space group	I 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	162.73Å 105.94Å 213.66Å 90.00° 111.23° 90.00°	Depositor
Resolution (Å)	47.82 – 2.29 47.82 – 2.29	Depositor EDS
% Data completeness (in resolution range)	100.0 (47.82-2.29) 100.0 (47.82-2.29)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.64 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.16_3549	Depositor
R, R_{free}	0.169 , 0.202 0.171 , 0.205	Depositor DCC
R_{free} test set	7610 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	42.1	Xtriage
Anisotropy	0.169	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 39.1	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	0.96	EDS
Total number of atoms	22347	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.63% 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: G4S, GAL, NA, 9RN, TLA, CL, EDO

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	0.24	0/5372	0.47	0/7262
1	B	0.25	0/5375	0.45	0/7266
1	C	0.23	0/5367	0.46	0/7255
1	D	0.24	0/5372	0.46	0/7262
All	All	0.24	0/21486	0.46	0/29045

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	5238	0	5129	23	0
1	B	5241	0	5131	38	0
1	C	5234	0	5124	38	0
1	D	5238	0	5127	39	0
2	E	47	0	20	3	0
2	F	47	0	20	3	0
2	G	47	0	20	2	0
2	H	47	0	20	2	0
3	A	20	0	8	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	10	0	4	0	0
3	C	10	0	4	0	0
3	D	10	0	4	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	1	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
6	B	4	0	6	0	0
7	A	310	0	0	3	0
7	B	290	0	0	7	0
7	C	281	0	0	4	0
7	D	265	0	0	3	0
All	All	22347	0	20617	143	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 3.

All (143) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:4:9RN:O5	2:G:4:9RN:C1	1.68	1.42
2:F:4:9RN:C1	2:F:4:9RN:O5	1.68	1.41
2:E:4:9RN:O5	2:E:4:9RN:C1	1.67	1.40
2:H:4:9RN:O5	2:H:4:9RN:C1	1.67	1.39
2:G:4:9RN:C1	2:G:4:9RN:C5	2.57	0.76
1:C:44:ASN:HB3	1:C:57:THR:HB	1.68	0.75
2:F:4:9RN:C1	2:F:4:9RN:C5	2.58	0.74
1:B:351:MET:HG2	1:B:354:MET:HB2	1.73	0.69
5:B:704:CL:CL	7:B:977:HOH:O	2.48	0.69
1:D:114:PRO:HD2	1:D:122:ALA:HB3	1.80	0.64
2:E:4:9RN:C1	2:E:4:9RN:C5	2.57	0.64
2:H:4:9RN:C1	2:H:4:9RN:C5	2.58	0.64
1:C:264:GLY:HA3	1:D:447:LYS:HB3	1.81	0.63
1:B:255:ASN:HB3	1:B:271:ARG:HG3	1.81	0.61
1:B:469:LYS:NZ	7:B:802:HOH:O	2.34	0.60
1:A:44:ASN:HB2	1:A:57:THR:HB	1.84	0.60
1:B:486:ASN:ND2	7:B:803:HOH:O	2.34	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:418:LEU:HD22	1:A:482:ALA:HB2	1.85	0.58
1:C:44:ASN:OD1	1:C:46:ASN:ND2	2.38	0.57
1:C:309:LYS:NZ	7:C:804:HOH:O	2.37	0.56
1:A:249:GLN:HB2	1:A:278:ILE:HD11	1.88	0.56
1:B:487:ILE:HG23	2:F:3:GAL:H62	1.89	0.56
1:A:96:LYS:HB2	1:A:112:ILE:HB	1.89	0.55
1:C:498:GLU:H	1:C:498:GLU:CD	2.15	0.55
1:A:39:GLU:OE1	1:A:59:LYS:NZ	2.29	0.55
1:B:36:ILE:HD13	1:B:49:SER:HB3	1.88	0.54
1:D:367:THR:HG23	1:D:371:ASP:OD2	2.07	0.54
1:B:332:LYS:NZ	7:B:805:HOH:O	2.40	0.54
1:A:351:MET:HG2	1:A:354:MET:HB2	1.90	0.53
1:C:539:ARG:NH2	7:C:807:HOH:O	2.41	0.53
1:B:418:LEU:HD22	1:B:482:ALA:HB2	1.91	0.53
1:B:81:ASP:OD2	1:B:85:LYS:HB3	2.09	0.52
1:C:515:GLY:HA3	1:C:529:MET:HE2	1.92	0.52
1:C:474:GLU:HG2	1:C:480:LEU:HD11	1.91	0.52
1:C:535:TYR:CZ	1:C:547:PRO:HB3	2.44	0.52
1:B:101:LYS:HG2	1:B:103:ALA:O	2.10	0.51
1:D:486:ASN:ND2	7:D:807:HOH:O	2.43	0.50
1:C:455:TRP:HB3	1:C:487:ILE:HD13	1.92	0.50
1:A:183:VAL:HB	1:A:545:SER:HB2	1.94	0.49
1:D:514:ASN:ND2	7:D:802:HOH:O	2.29	0.49
1:C:233:LYS:NZ	7:C:809:HOH:O	2.44	0.49
1:D:237:MET:HG2	1:D:303:MET:HG2	1.93	0.49
1:A:535:TYR:CZ	1:A:547:PRO:HB3	2.48	0.49
1:C:688:VAL:HG22	1:D:688:VAL:HG22	1.95	0.49
1:A:101:LYS:O	3:A:702:TLA:H2	2.13	0.49
1:D:455:TRP:HB3	1:D:487:ILE:HD13	1.95	0.49
1:C:324:LEU:HD12	1:C:328:PRO:HA	1.94	0.49
1:A:691:HIS:ND1	7:A:803:HOH:O	2.35	0.48
1:B:249:GLN:HB2	1:B:278:ILE:HD11	1.94	0.48
1:B:483:ASP:O	1:B:514:ASN:HA	2.13	0.48
1:C:418:LEU:HD22	1:C:482:ALA:HB2	1.95	0.48
1:D:514:ASN:OD1	1:D:527:PHE:HA	2.14	0.48
1:B:58:GLU:OE2	1:B:234:SER:HB3	2.14	0.47
1:B:83:LYS:H	1:B:83:LYS:HG3	1.48	0.47
1:B:223:THR:HG21	1:B:533:MET:SD	2.55	0.47
1:D:36:ILE:HD13	1:D:99:VAL:HG12	1.95	0.47
1:B:78:THR:HA	1:B:87:GLN:O	2.15	0.47
1:D:141:ASP:OD2	1:D:289:ARG:HD2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:671:GLY:HA2	1:D:685:HIS:HA	1.97	0.47
1:D:418:LEU:HD22	1:D:482:ALA:HB2	1.96	0.47
1:A:256:GLY:HA2	1:A:258:TYR:CE1	2.51	0.46
1:B:70:TRP:CD1	1:B:269:TYR:HB3	2.50	0.46
1:C:671:GLY:HA2	1:C:685:HIS:HA	1.97	0.46
1:D:44:ASN:HB3	1:D:57:THR:HB	1.98	0.46
1:D:103:ALA:HB3	1:D:106:THR:HB	1.97	0.46
1:D:535:TYR:CZ	1:D:547:PRO:HB3	2.51	0.46
1:B:95:LYS:HD2	1:B:95:LYS:HA	1.52	0.45
1:D:168:THR:HA	1:D:187:SER:O	2.17	0.45
1:B:324:LEU:HD12	1:B:328:PRO:HA	1.99	0.45
1:D:162:ARG:H	1:D:276:ASP:CG	2.25	0.45
1:C:342:TYR:HA	1:C:382:TRP:HB2	1.98	0.45
1:A:264:GLY:HA3	1:B:447:LYS:HB3	1.98	0.45
1:A:284:GLU:HG2	1:A:288:MET:HG2	1.99	0.45
1:B:66:LYS:HD3	7:B:912:HOH:O	2.16	0.45
1:D:587:LEU:HD11	1:D:652:ALA:HB3	1.98	0.45
1:C:110:LYS:HG2	1:C:112:ILE:HD11	1.99	0.45
1:D:285:LYS:HD2	1:D:285:LYS:HA	1.78	0.45
1:A:486:ASN:ND2	7:A:818:HOH:O	2.50	0.44
1:C:249:GLN:HB2	1:C:278:ILE:HD11	1.97	0.44
1:D:569:ASP:HA	1:D:581:LYS:HD3	2.00	0.44
1:C:116:PHE:CE2	1:C:122:ALA:HB2	2.53	0.44
1:C:351:MET:HE3	1:D:189:ILE:HG23	2.00	0.44
1:B:366:LYS:NZ	1:B:370:ASP:OD2	2.26	0.44
1:B:122:ALA:HA	1:B:152:PHE:CZ	2.53	0.44
1:D:186:PRO:HD2	1:D:200:TRP:CZ2	2.53	0.44
1:C:342:TYR:O	1:C:598:PRO:HD3	2.18	0.44
1:B:430:GLU:CD	1:B:457:ARG:HH12	2.26	0.44
1:C:483:ASP:O	1:C:514:ASN:HA	2.17	0.43
1:C:556:HIS:HE1	1:C:622:THR:O	2.01	0.43
1:C:529:MET:HA	1:C:560:ALA:O	2.18	0.43
1:D:639:VAL:HG11	1:D:661:ILE:HG21	1.99	0.43
1:A:125:VAL:HA	1:A:148:ARG:O	2.18	0.43
1:D:249:GLN:HB2	1:D:278:ILE:HD11	2.01	0.43
1:D:428:MET:HG3	1:D:435:PHE:CE2	2.53	0.43
1:D:334:PRO:HG3	1:D:623:PHE:CD2	2.53	0.43
1:B:557:GLU:CD	1:B:626:GLU:HB2	2.43	0.43
1:A:671:GLY:HA2	1:A:685:HIS:HA	2.01	0.42
1:D:641:ARG:HD2	1:D:649:GLU:OE2	2.19	0.42
1:B:426:SER:H	1:B:427:PRO:HD3	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:319:SER:HA	1:C:558:ALA:HA	2.00	0.42
1:B:334:PRO:HG3	1:B:623:PHE:CD1	2.54	0.42
1:D:173:GLY:HA3	1:D:230:TYR:O	2.19	0.42
1:D:186:PRO:HD2	1:D:200:TRP:HZ2	1.83	0.42
1:D:483:ASP:O	1:D:514:ASN:HA	2.19	0.42
1:A:63:HIS:CD2	1:A:231:ASN:HD21	2.38	0.42
1:C:162:ARG:H	1:C:276:ASP:CG	2.27	0.42
1:A:149:SER:HB3	1:A:152:PHE:O	2.19	0.42
1:C:305:LYS:NZ	1:C:626:GLU:OE2	2.45	0.42
1:B:237:MET:O	1:B:292:TYR:HA	2.20	0.42
1:B:251:ASN:O	1:B:273:VAL:HA	2.19	0.42
1:C:247:HIS:HB2	1:C:278:ILE:HB	2.02	0.42
1:B:514:ASN:HB2	7:B:853:HOH:O	2.20	0.42
1:D:310:GLU:O	1:D:314:ARG:HG3	2.20	0.42
1:A:340:GLY:HA3	1:A:595:PHE:CD1	2.55	0.42
1:B:247:HIS:HB2	1:B:278:ILE:HB	2.02	0.42
1:C:36:ILE:HD12	1:C:99:VAL:HG12	2.02	0.42
1:B:569:ASP:HA	1:B:581:LYS:HD3	2.02	0.41
1:D:426:SER:N	1:D:427:PRO:HD3	2.35	0.41
1:D:214:GLU:HG2	1:D:217:ASN:HD22	1.85	0.41
1:D:340:GLY:HA3	1:D:595:PHE:HD1	1.85	0.41
1:D:493:ARG:HD2	7:D:962:HOH:O	2.20	0.41
1:D:372:LEU:HD13	1:D:608:ILE:HG21	2.03	0.41
1:B:44:ASN:HB3	1:B:57:THR:HB	2.03	0.41
1:B:671:GLY:HA2	1:B:685:HIS:HA	2.02	0.41
1:C:428:MET:HG3	1:C:435:PHE:CD2	2.56	0.41
1:B:560:ALA:HB1	7:B:866:HOH:O	2.20	0.41
1:C:340:GLY:HA3	1:C:595:PHE:HD1	1.86	0.41
1:C:471:ILE:HD12	1:C:471:ILE:HA	1.88	0.41
1:D:426:SER:H	1:D:427:PRO:HD3	1.85	0.41
1:C:255:ASN:HB3	1:C:271:ARG:HG3	2.02	0.41
1:B:63:HIS:CD2	1:B:231:ASN:HD21	2.39	0.40
1:C:425:TYR:CE1	1:C:467:ALA:HA	2.56	0.40
1:C:472:GLU:HG2	1:C:510:PHE:CE1	2.56	0.40
1:C:514:ASN:HB2	7:C:854:HOH:O	2.21	0.40
1:B:46:ASN:O	1:B:54:PHE:HA	2.22	0.40
1:D:210:GLN:HG2	1:D:258:TYR:CD1	2.57	0.40
1:A:426:SER:HA	1:A:458:VAL:HB	2.01	0.40
1:C:568:PRO:HD2	1:C:571:GLU:HB2	2.03	0.40
1:A:556:HIS:HE1	1:A:622:THR:O	2.04	0.40
1:C:214:GLU:HG2	1:C:247:HIS:CD2	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:487:ILE:HG23	2:E:3:GAL:H62	2.04	0.40
1:A:514:ASN:HB2	7:A:926:HOH:O	2.21	0.40
1:B:173:GLY:HA3	1:B:230:TYR:O	2.21	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	657/676 (97%)	627 (95%)	30 (5%)	0	100	100
1	B	657/676 (97%)	629 (96%)	28 (4%)	0	100	100
1	C	657/676 (97%)	629 (96%)	28 (4%)	0	100	100
1	D	657/676 (97%)	629 (96%)	27 (4%)	1 (0%)	43	55
All	All	2628/2704 (97%)	2514 (96%)	113 (4%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	383	GLY

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	561/577 (97%)	551 (98%)	10 (2%)	51	70
1	B	562/577 (97%)	548 (98%)	14 (2%)	42	60
1	C	560/577 (97%)	550 (98%)	10 (2%)	51	70
1	D	561/577 (97%)	546 (97%)	15 (3%)	39	58
All	All	2244/2308 (97%)	2195 (98%)	49 (2%)	45	65

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	LYS
1	A	72	ASN
1	A	83	LYS
1	A	93	LYS
1	A	108	SER
1	A	167	LYS
1	A	309	LYS
1	A	351	MET
1	A	529	MET
1	A	693	ASP
1	B	42	LYS
1	B	83	LYS
1	B	85	LYS
1	B	93	LYS
1	B	95	LYS
1	B	123	LYS
1	B	167	LYS
1	B	214	GLU
1	B	325	LYS
1	B	351	MET
1	B	362	LYS
1	B	404	LYS
1	B	529	MET
1	B	675	GLN
1	C	82	SER
1	C	97	ILE
1	C	110	LYS
1	C	285	LYS
1	C	351	MET
1	C	360	GLU
1	C	497	LYS
1	C	502	GLU
1	C	529	MET

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Mol	Chain	Res	Type
1	C	661	ILE
1	D	49	SER
1	D	58	GLU
1	D	82	SER
1	D	83	LYS
1	D	86	LYS
1	D	117	GLU
1	D	146	GLU
1	D	214	GLU
1	D	222	LEU
1	D	326	ARG
1	D	351	MET
1	D	404	LYS
1	D	542	SER
1	D	662	GLU
1	D	677	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	44	ASN
1	B	247	HIS
1	B	486	ASN
1	C	63	HIS
1	C	147	HIS
1	C	308	GLN
1	C	331	ASN
1	C	450	ASN
1	D	46	ASN
1	D	147	HIS
1	D	363	ASN
1	D	387	ASN
1	D	486	ASN
1	D	640	GLN
1	D	675	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	G4S	E	1	2	16,16,16	1.64	4 (25%)	21,24,24	1.17	2 (9%)
2	9RN	E	2	2	11,11,12	5.64	7 (63%)	13,16,18	2.53	4 (30%)
2	GAL	E	3	2	11,11,12	1.39	2 (18%)	15,15,17	1.15	1 (6%)
2	9RN	E	4	2	11,11,12	6.82	7 (63%)	13,16,18	2.35	4 (30%)
2	G4S	F	1	2	16,16,16	1.62	4 (25%)	21,24,24	1.05	1 (4%)
2	9RN	F	2	2	11,11,12	5.66	7 (63%)	13,16,18	2.37	5 (38%)
2	GAL	F	3	2	11,11,12	1.43	2 (18%)	15,15,17	1.18	2 (13%)
2	9RN	F	4	2	11,11,12	6.98	7 (63%)	13,16,18	2.49	5 (38%)
2	G4S	G	1	2	16,16,16	1.61	4 (25%)	21,24,24	0.88	0
2	9RN	G	2	2	11,11,12	5.68	7 (63%)	13,16,18	2.60	6 (46%)
2	GAL	G	3	2	11,11,12	1.35	2 (18%)	15,15,17	1.06	1 (6%)
2	9RN	G	4	2	11,11,12	7.03	7 (63%)	13,16,18	2.36	4 (30%)
2	G4S	H	1	2	16,16,16	1.59	4 (25%)	21,24,24	1.28	3 (14%)
2	9RN	H	2	2	11,11,12	5.62	7 (63%)	13,16,18	2.48	5 (38%)
2	GAL	H	3	2	11,11,12	1.39	2 (18%)	15,15,17	0.99	1 (6%)
2	9RN	H	4	2	11,11,12	6.86	8 (72%)	13,16,18	2.54	5 (38%)

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	G4S	E	1	2	-	5/7/27/27	0/1/1/1
2	9RN	E	2	2	-	-	0/3/2/2
2	GAL	E	3	2	-	0/2/19/22	0/1/1/1
2	9RN	E	4	2	-	-	0/3/2/2
2	G4S	F	1	2	-	5/7/27/27	0/1/1/1
2	9RN	F	2	2	-	-	0/3/2/2
2	GAL	F	3	2	-	0/2/19/22	0/1/1/1
2	9RN	F	4	2	-	-	0/3/2/2
2	G4S	G	1	2	-	4/7/27/27	0/1/1/1
2	9RN	G	2	2	-	-	0/3/2/2
2	GAL	G	3	2	-	0/2/19/22	0/1/1/1
2	9RN	G	4	2	-	-	0/3/2/2
2	G4S	H	1	2	-	4/7/27/27	0/1/1/1
2	9RN	H	2	2	-	-	0/3/2/2
2	GAL	H	3	2	-	0/2/19/22	0/1/1/1
2	9RN	H	4	2	-	-	0/3/2/2

All (81) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	4	9RN	O5-C1	14.94	1.68	1.43
2	G	4	9RN	O5-C1	14.80	1.68	1.43
2	H	4	9RN	O5-C1	14.38	1.67	1.43
2	E	4	9RN	O5-C1	14.34	1.67	1.43
2	G	4	9RN	O3-C6	-9.95	1.22	1.43
2	H	4	9RN	O3-C6	-9.78	1.22	1.43
2	F	2	9RN	O5-C1	9.73	1.60	1.43
2	E	4	9RN	O3-C6	-9.72	1.23	1.43
2	G	2	9RN	O5-C1	9.66	1.59	1.43
2	F	4	9RN	O3-C6	-9.61	1.23	1.43
2	H	2	9RN	O5-C1	9.59	1.59	1.43
2	E	2	9RN	C4-C3	-9.50	1.32	1.53
2	H	2	9RN	C4-C3	-9.48	1.32	1.53
2	E	2	9RN	O5-C1	9.47	1.59	1.43
2	F	2	9RN	C4-C3	-9.42	1.32	1.53
2	G	2	9RN	C4-C3	-9.30	1.32	1.53
2	G	4	9RN	C1-C2	8.73	1.73	1.52
2	F	4	9RN	C1-C2	8.61	1.72	1.52
2	E	4	9RN	C1-C2	8.58	1.72	1.52
2	G	2	9RN	C4-C5	-8.42	1.31	1.53
2	H	4	9RN	C1-C2	8.38	1.72	1.52
2	E	2	9RN	C4-C5	-8.30	1.31	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	2	9RN	C4-C5	-8.25	1.32	1.53
2	H	2	9RN	C4-C5	-8.10	1.32	1.53
2	G	2	9RN	O3-C6	6.96	1.58	1.43
2	H	2	9RN	O3-C6	6.89	1.58	1.43
2	F	2	9RN	O3-C6	6.72	1.58	1.43
2	E	2	9RN	O3-C6	6.70	1.58	1.43
2	G	4	9RN	C4-C3	-6.60	1.38	1.53
2	H	4	9RN	C4-C3	-6.53	1.38	1.53
2	F	4	9RN	C4-C3	-6.47	1.38	1.53
2	E	4	9RN	C4-C3	-6.36	1.39	1.53
2	H	4	9RN	C4-C5	-6.35	1.36	1.53
2	G	4	9RN	C4-C5	-6.30	1.37	1.53
2	F	4	9RN	C4-C5	-6.28	1.37	1.53
2	E	4	9RN	C4-C5	-6.17	1.37	1.53
2	G	4	9RN	C6-C5	-5.91	1.41	1.52
2	F	4	9RN	C6-C5	-5.90	1.41	1.52
2	E	2	9RN	C1-C2	5.84	1.66	1.52
2	H	4	9RN	C6-C5	-5.72	1.41	1.52
2	E	4	9RN	C6-C5	-5.72	1.41	1.52
2	G	2	9RN	C1-C2	5.65	1.65	1.52
2	F	2	9RN	C1-C2	5.62	1.65	1.52
2	H	2	9RN	C1-C2	5.34	1.65	1.52
2	G	4	9RN	C2-C3	-5.00	1.41	1.52
2	F	4	9RN	C2-C3	-4.92	1.41	1.52
2	H	4	9RN	C2-C3	-4.80	1.41	1.52
2	E	4	9RN	C2-C3	-4.76	1.41	1.52
2	G	2	9RN	O3-C3	3.43	1.51	1.43
2	H	1	G4S	O4-S	3.39	1.67	1.57
2	E	2	9RN	O3-C3	3.36	1.51	1.43
2	H	2	9RN	O3-C3	3.35	1.51	1.43
2	F	2	9RN	O3-C3	3.31	1.51	1.43
2	F	2	9RN	C2-C3	-3.28	1.45	1.52
2	H	2	9RN	C2-C3	-3.25	1.45	1.52
2	F	1	G4S	O4-S	3.23	1.67	1.57
2	G	2	9RN	C2-C3	-3.11	1.45	1.52
2	G	1	G4S	O4-C4	-3.05	1.39	1.46
2	G	1	G4S	O4-S	3.00	1.66	1.57
2	E	3	GAL	O5-C5	2.95	1.49	1.43
2	E	1	G4S	O4-C4	-2.93	1.39	1.46
2	E	2	9RN	C2-C3	-2.91	1.45	1.52
2	E	1	G4S	O4-S	2.91	1.66	1.57
2	F	1	G4S	O4-C4	-2.77	1.40	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	3	GAL	O5-C5	2.77	1.48	1.43
2	H	3	GAL	O5-C5	2.62	1.48	1.43
2	H	1	G4S	O4-C4	-2.53	1.40	1.46
2	H	1	G4S	O5-C1	2.49	1.49	1.42
2	G	3	GAL	O5-C5	2.46	1.48	1.43
2	F	1	G4S	O5-C1	2.46	1.48	1.42
2	E	1	G4S	O5-C1	2.44	1.48	1.42
2	H	3	GAL	C2-C3	-2.33	1.48	1.52
2	F	1	G4S	C3-C2	-2.27	1.46	1.52
2	G	3	GAL	C2-C3	-2.23	1.49	1.52
2	E	1	G4S	C3-C2	-2.23	1.46	1.52
2	F	3	GAL	C2-C3	-2.22	1.49	1.52
2	G	1	G4S	C3-C2	-2.18	1.46	1.52
2	G	1	G4S	O5-C1	2.14	1.48	1.42
2	E	3	GAL	C2-C3	-2.10	1.49	1.52
2	H	4	9RN	O3-C3	-2.06	1.39	1.43
2	H	1	G4S	C3-C2	-2.06	1.47	1.52

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	2	9RN	O3-C3-C2	5.23	119.86	108.12
2	G	2	9RN	C3-C4-C5	4.94	112.62	101.99
2	E	2	9RN	C3-C4-C5	4.86	112.44	101.99
2	H	2	9RN	C3-C4-C5	4.85	112.42	101.99
2	F	2	9RN	C3-C4-C5	4.73	112.15	101.99
2	F	4	9RN	C3-C4-C5	4.53	111.73	101.99
2	G	4	9RN	O3-C6-C5	4.48	112.19	104.42
2	H	4	9RN	O3-C6-C5	4.42	112.09	104.42
2	E	4	9RN	O3-C6-C5	4.42	112.08	104.42
2	G	4	9RN	C3-C4-C5	4.36	111.36	101.99
2	H	4	9RN	O5-C5-C6	4.28	119.59	113.33
2	F	4	9RN	O3-C6-C5	4.26	111.81	104.42
2	H	4	9RN	C3-C4-C5	4.22	111.06	101.99
2	G	2	9RN	O3-C3-C2	4.22	117.58	108.12
2	H	2	9RN	O3-C3-C2	4.21	117.56	108.12
2	E	4	9RN	C3-C4-C5	4.19	111.00	101.99
2	H	4	9RN	O3-C3-C2	4.10	117.32	108.12
2	G	4	9RN	O3-C3-C2	4.00	117.10	108.12
2	G	2	9RN	C6-O3-C3	-3.96	99.92	107.84
2	F	4	9RN	O3-C3-C2	3.95	116.97	108.12
2	E	4	9RN	O3-C3-C2	3.93	116.94	108.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	4	9RN	O5-C5-C6	3.88	119.01	113.33
2	F	2	9RN	O3-C3-C2	3.83	116.71	108.12
2	F	2	9RN	C6-O3-C3	-3.70	100.46	107.84
2	E	2	9RN	C6-O3-C3	-3.63	100.60	107.84
2	G	2	9RN	C1-O5-C5	-3.50	107.50	112.19
2	H	2	9RN	C1-O5-C5	-3.49	107.51	112.19
2	H	2	9RN	C6-O3-C3	-3.35	101.16	107.84
2	G	2	9RN	O5-C5-C6	3.29	118.14	113.33
2	F	2	9RN	C1-O5-C5	-3.20	107.90	112.19
2	E	4	9RN	O5-C5-C6	3.14	117.92	113.33
2	E	2	9RN	O5-C5-C6	2.87	117.52	113.33
2	G	4	9RN	O5-C5-C6	2.78	117.39	113.33
2	E	3	GAL	C1-C2-C3	2.71	113.59	109.64
2	E	1	G4S	O5-C1-C2	2.68	115.00	110.30
2	F	3	GAL	C1-C2-C3	2.67	113.54	109.64
2	G	3	GAL	C1-C2-C3	2.58	113.39	109.64
2	H	1	G4S	O5-C1-C2	2.52	114.74	110.30
2	H	2	9RN	O5-C5-C6	2.50	116.98	113.33
2	F	2	9RN	O5-C5-C6	2.48	116.96	113.33
2	H	3	GAL	C1-C2-C3	2.44	113.19	109.64
2	E	1	G4S	O8-S-O7	-2.41	102.95	112.24
2	H	1	G4S	O9-S-O4	2.19	111.41	106.37
2	H	1	G4S	O9-S-O8	-2.13	101.09	108.56
2	F	3	GAL	C6-C5-C4	-2.11	107.85	113.02
2	H	4	9RN	C1-C2-C3	2.10	112.17	109.29
2	G	2	9RN	O4-C4-C5	-2.09	105.09	111.08
2	F	4	9RN	C6-O3-C3	2.07	111.98	107.84
2	F	1	G4S	C6-C5-C4	-2.00	107.75	113.38

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	F	1	G4S	C4-O4-S-O7
2	F	1	G4S	C4-O4-S-O8
2	F	1	G4S	O5-C5-C6-O6
2	E	1	G4S	O5-C5-C6-O6
2	F	1	G4S	C4-C5-C6-O6
2	G	1	G4S	C4-O4-S-O7
2	G	1	G4S	C4-O4-S-O8
2	H	1	G4S	O5-C5-C6-O6
2	E	1	G4S	C4-O4-S-O7

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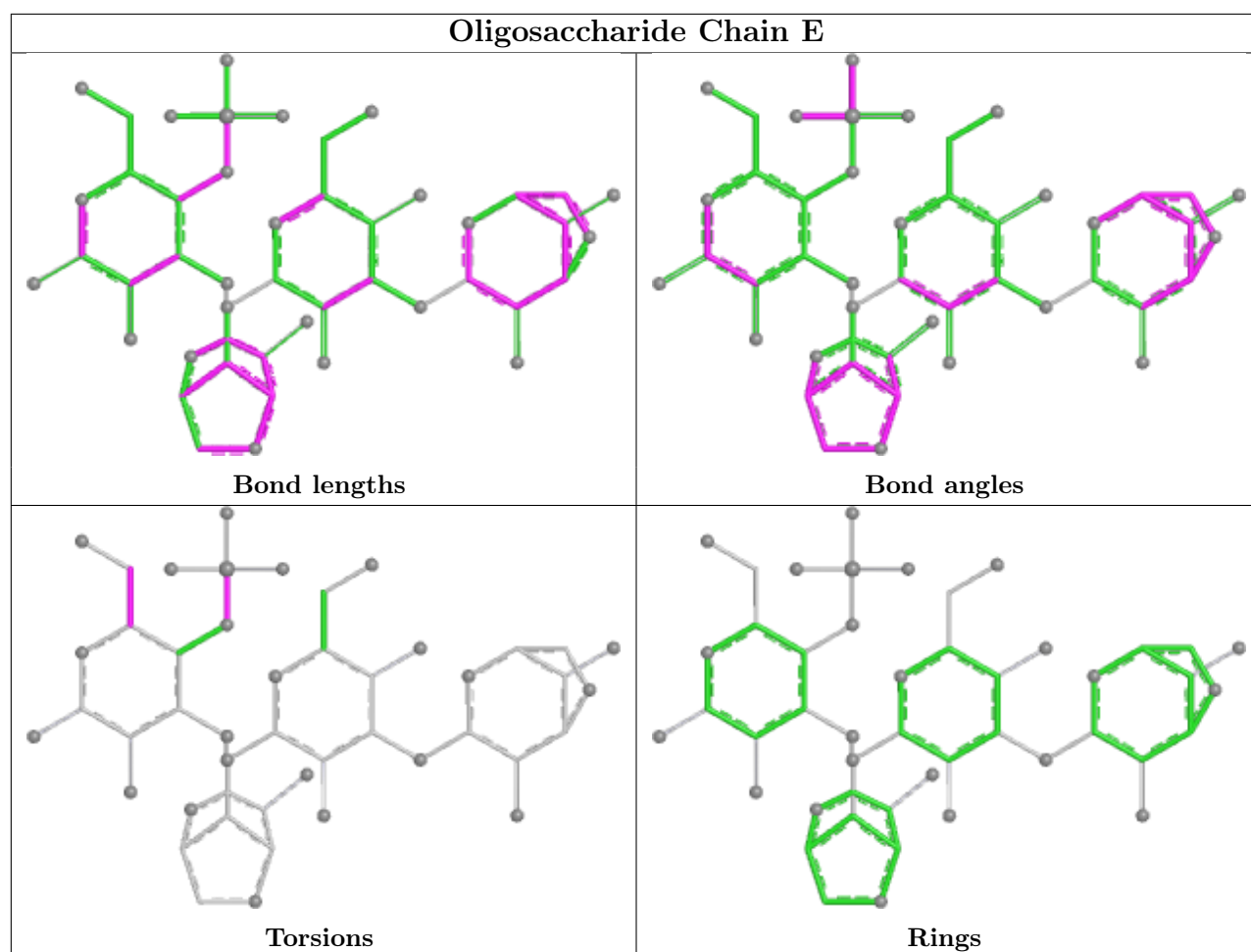
Mol	Chain	Res	Type	Atoms
2	E	1	G4S	C4-O4-S-O9
2	F	1	G4S	C4-O4-S-O9
2	G	1	G4S	C4-O4-S-O9
2	H	1	G4S	C4-O4-S-O8
2	G	1	G4S	O5-C5-C6-O6
2	E	1	G4S	C4-C5-C6-O6
2	E	1	G4S	C4-O4-S-O8
2	H	1	G4S	C3-C4-O4-S
2	H	1	G4S	C4-O4-S-O7

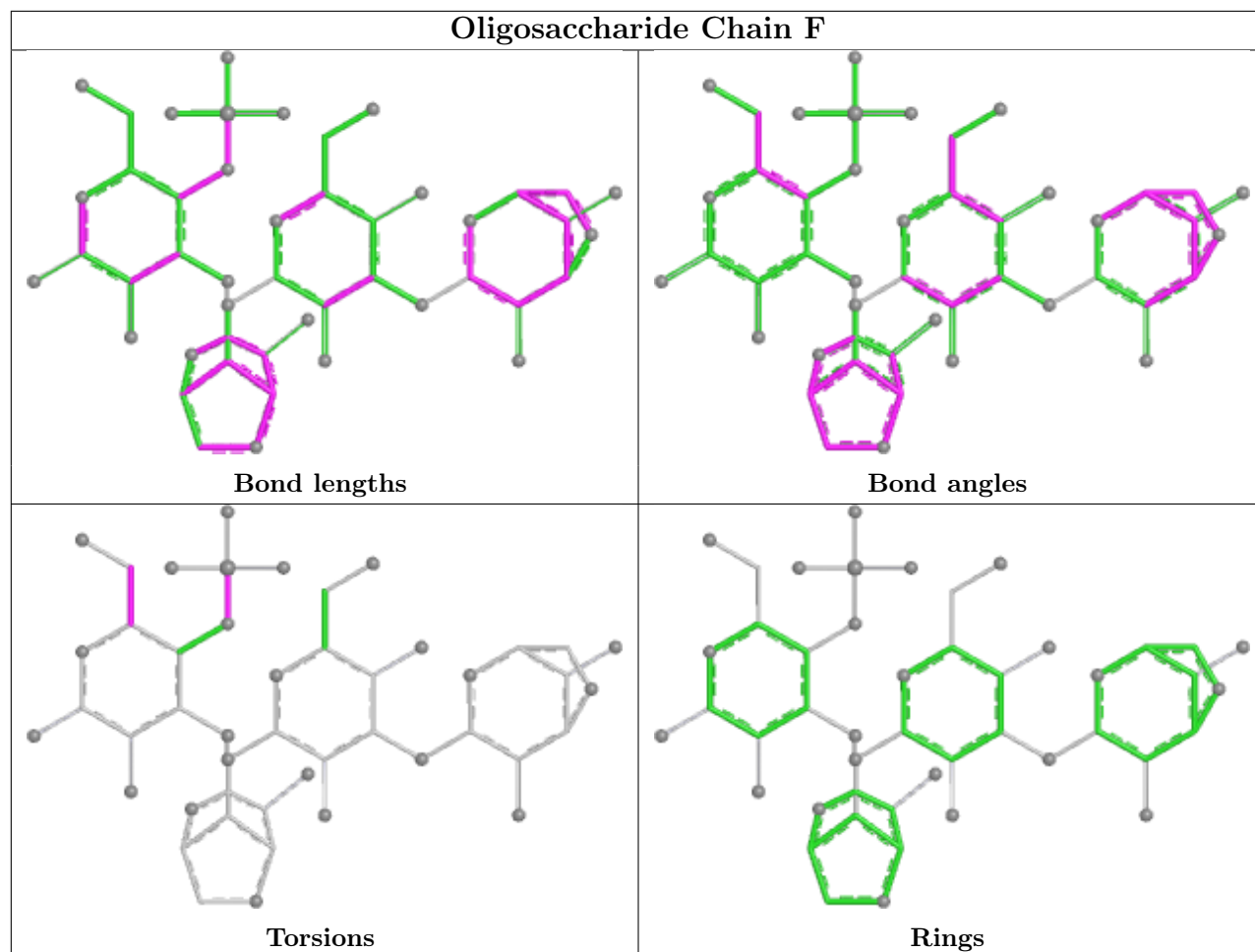
There are no ring outliers.

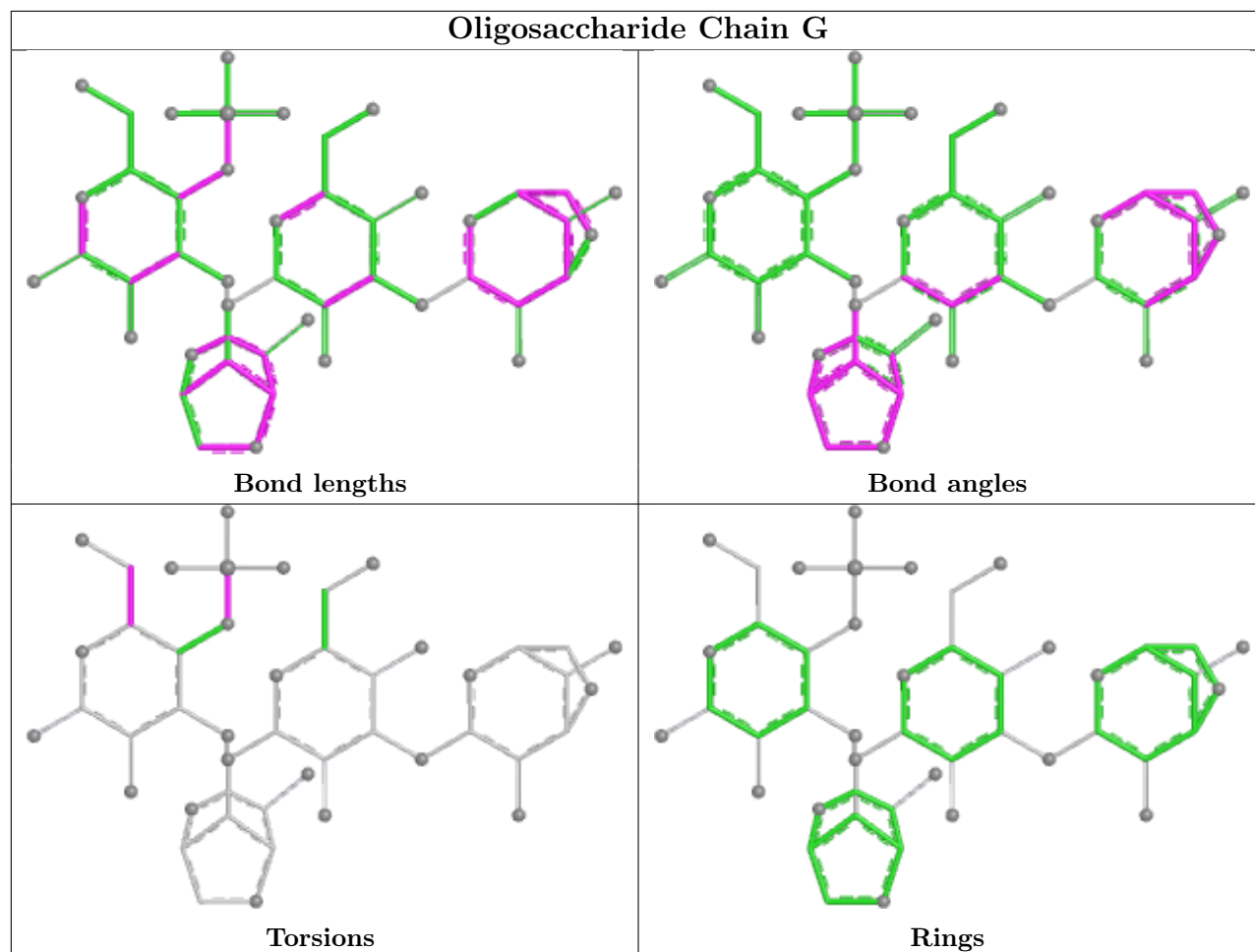
6 monomers are involved in 10 short contacts:

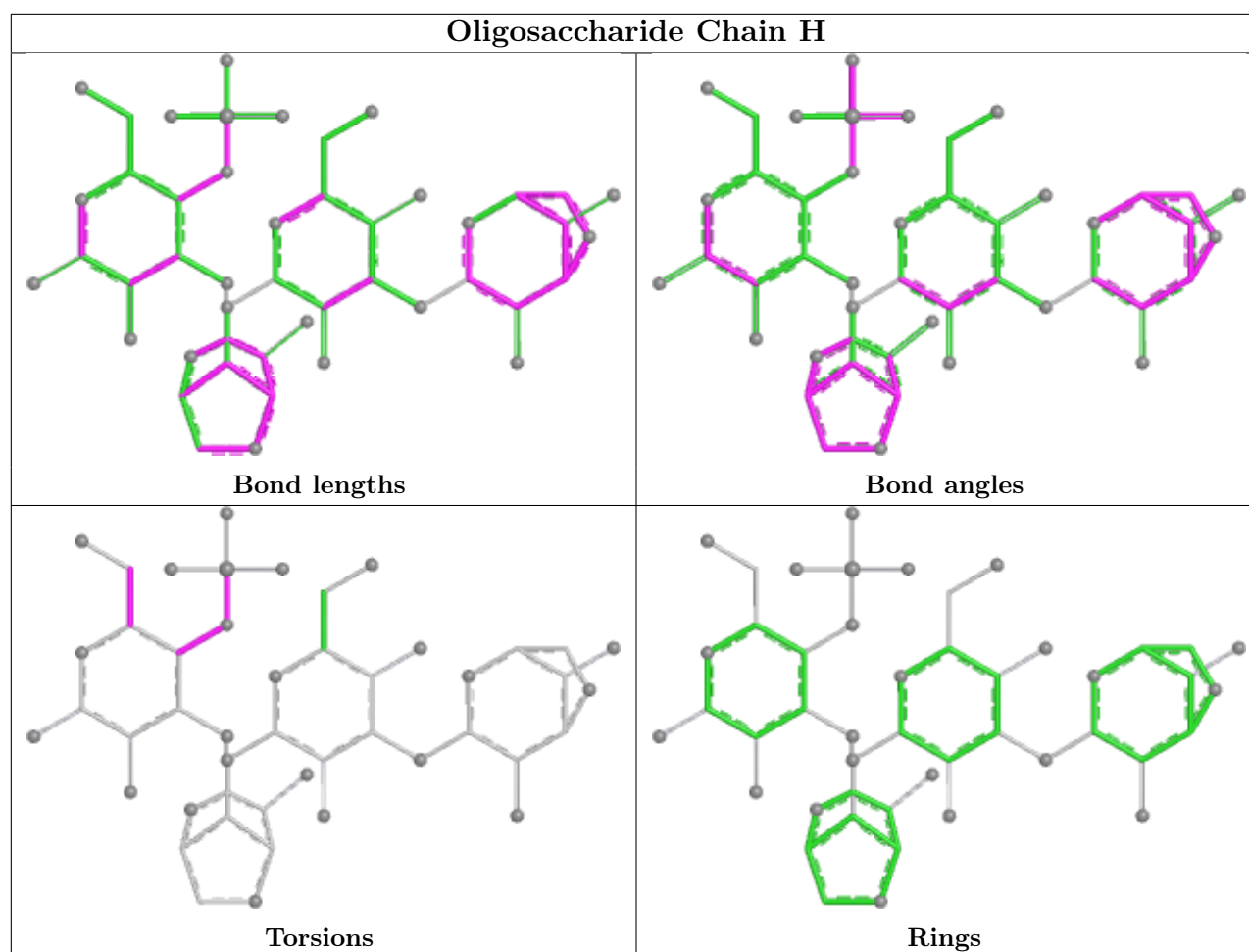
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	3	GAL	1	0
2	E	3	GAL	1	0
2	G	4	9RN	2	0
2	F	4	9RN	2	0
2	E	4	9RN	2	0
2	H	4	9RN	2	0

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 14 ligands modelled in this entry, 8 are monoatomic - leaving 6 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
3	TLA	D	701	-	9,9,9	1.03	0	12,12,12	1.42	2 (16%)
6	EDO	B	702	-	3,3,3	0.41	0	2,2,2	0.39	0
3	TLA	B	701	-	9,9,9	1.04	0	12,12,12	1.21	1 (8%)
3	TLA	A	702	-	9,9,9	1.10	0	12,12,12	2.18	5 (41%)
3	TLA	C	701	-	9,9,9	1.04	0	12,12,12	1.30	2 (16%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	TLA	A	701	-	9,9,9	1.11	0	12,12,12	1.28	2 (16%)

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
3	TLA	D	701	-	-	0/12/12/12	-
6	EDO	B	702	-	-	1/1/1/1	-
3	TLA	B	701	-	-	1/12/12/12	-
3	TLA	A	702	-	-	5/12/12/12	-
3	TLA	C	701	-	-	4/12/12/12	-
3	TLA	A	701	-	-	2/12/12/12	-

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	702	TLA	C3-C2-C1	3.65	117.90	109.82
3	A	702	TLA	C2-C3-C4	3.26	117.06	109.82
3	A	702	TLA	O11-C1-C2	3.14	122.03	113.31
3	A	702	TLA	O41-C4-C3	3.05	121.78	113.31
3	D	701	TLA	O11-C1-C2	2.68	120.75	113.31
3	C	701	TLA	O11-C1-C2	2.66	120.70	113.31
3	A	701	TLA	O41-C4-C3	2.65	120.67	113.31
3	B	701	TLA	O11-C1-C2	2.65	120.66	113.31
3	D	701	TLA	O41-C4-C3	2.60	120.54	113.31
3	A	701	TLA	O11-C1-C2	2.08	119.08	113.31
3	A	702	TLA	O11-C1-O1	-2.07	119.38	124.08
3	C	701	TLA	O41-C4-C3	2.05	119.01	113.31

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	702	TLA	C1-C2-C3-C4
3	A	702	TLA	O3-C3-C4-O4
3	A	702	TLA	O3-C3-C4-O41
3	A	702	TLA	C1-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
3	A	702	TLA	O2-C2-C3-C4
6	B	702	EDO	O1-C1-C2-O2
3	C	701	TLA	O11-C1-C2-C3
3	C	701	TLA	O1-C1-C2-C3
3	A	701	TLA	C2-C3-C4-O4
3	B	701	TLA	O1-C1-C2-C3
3	C	701	TLA	O1-C1-C2-O2
3	A	701	TLA	O3-C3-C4-O4
3	C	701	TLA	O11-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	702	TLA	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	659/676 (97%)	-0.34	2 (0%) 90 90	30, 40, 58, 83	0
1	B	659/676 (97%)	-0.31	1 (0%) 91 91	32, 42, 61, 83	0
1	C	659/676 (97%)	-0.26	1 (0%) 91 91	33, 44, 62, 87	0
1	D	659/676 (97%)	-0.24	2 (0%) 90 90	34, 44, 63, 83	0
All	All	2636/2704 (97%)	-0.29	6 (0%) 91 91	30, 43, 61, 87	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	35	PRO	3.2
1	D	35	PRO	2.7
1	D	692	MET	2.6
1	B	35	PRO	2.5
1	A	35	PRO	2.3
1	A	538	ASP	2.2

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.

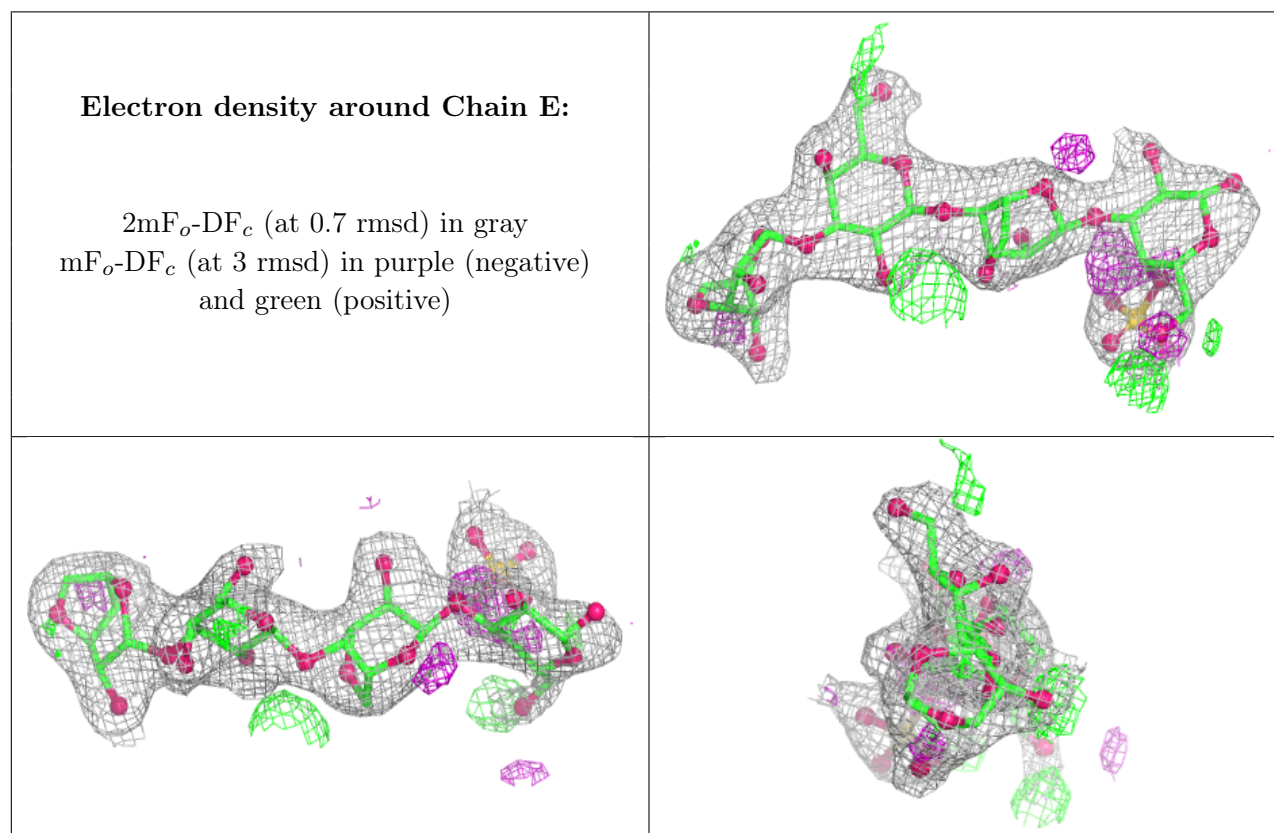
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	G4S	F	1	16/16	0.78	0.16	76,86,92,93	0

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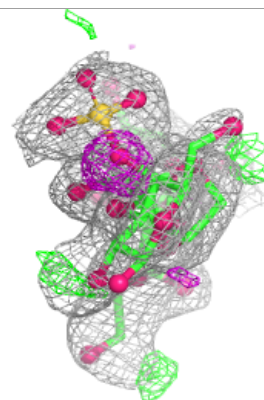
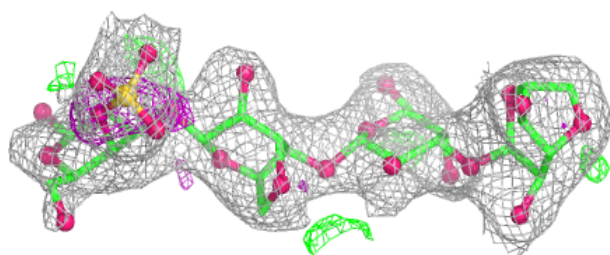
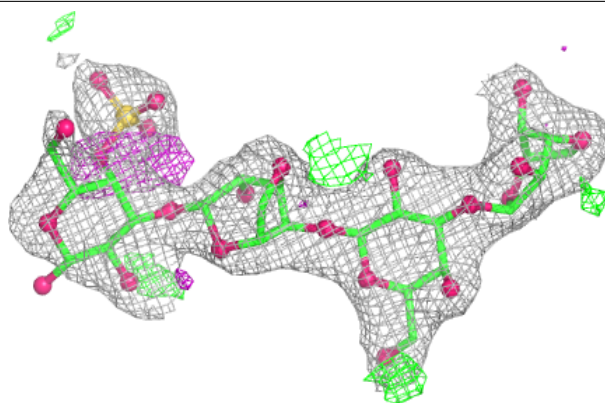
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	G4S	E	1	16/16	0.82	0.13	69,79,86,89	0
2	G4S	G	1	16/16	0.83	0.13	74,82,85,85	0
2	G4S	H	1	16/16	0.84	0.13	74,79,86,90	0
2	9RN	H	2	10/11	0.92	0.10	54,61,65,66	0
2	9RN	H	4	10/11	0.92	0.11	45,50,51,53	0
2	9RN	E	4	10/11	0.93	0.09	37,46,47,47	0
2	9RN	G	2	10/11	0.93	0.08	57,60,62,66	0
2	GAL	G	3	11/12	0.93	0.08	43,47,51,53	0
2	GAL	E	3	11/12	0.93	0.09	39,47,51,55	0
2	9RN	F	2	10/11	0.93	0.09	53,59,64,68	0
2	GAL	F	3	11/12	0.93	0.09	46,50,55,60	0
2	9RN	G	4	10/11	0.94	0.08	49,52,54,55	0
2	GAL	H	3	11/12	0.94	0.09	48,50,53,57	0
2	9RN	F	4	10/11	0.94	0.08	46,51,54,54	0
2	9RN	E	2	10/11	0.95	0.08	49,52,59,59	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.

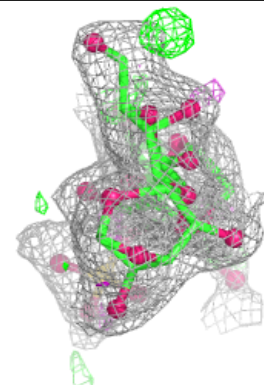
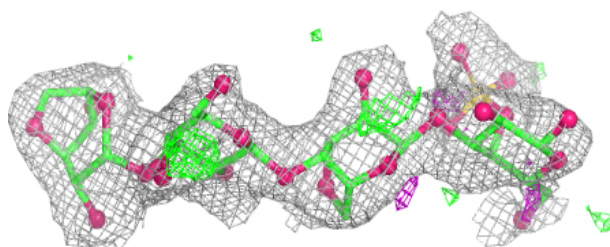
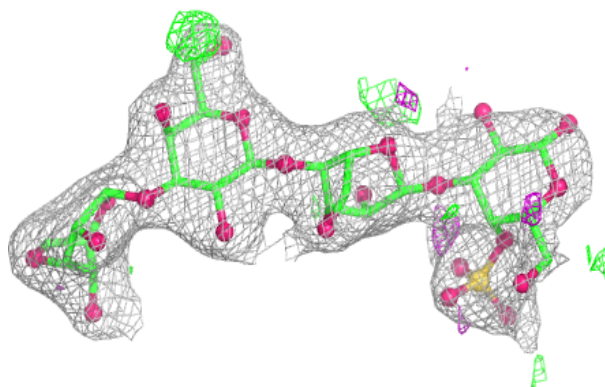


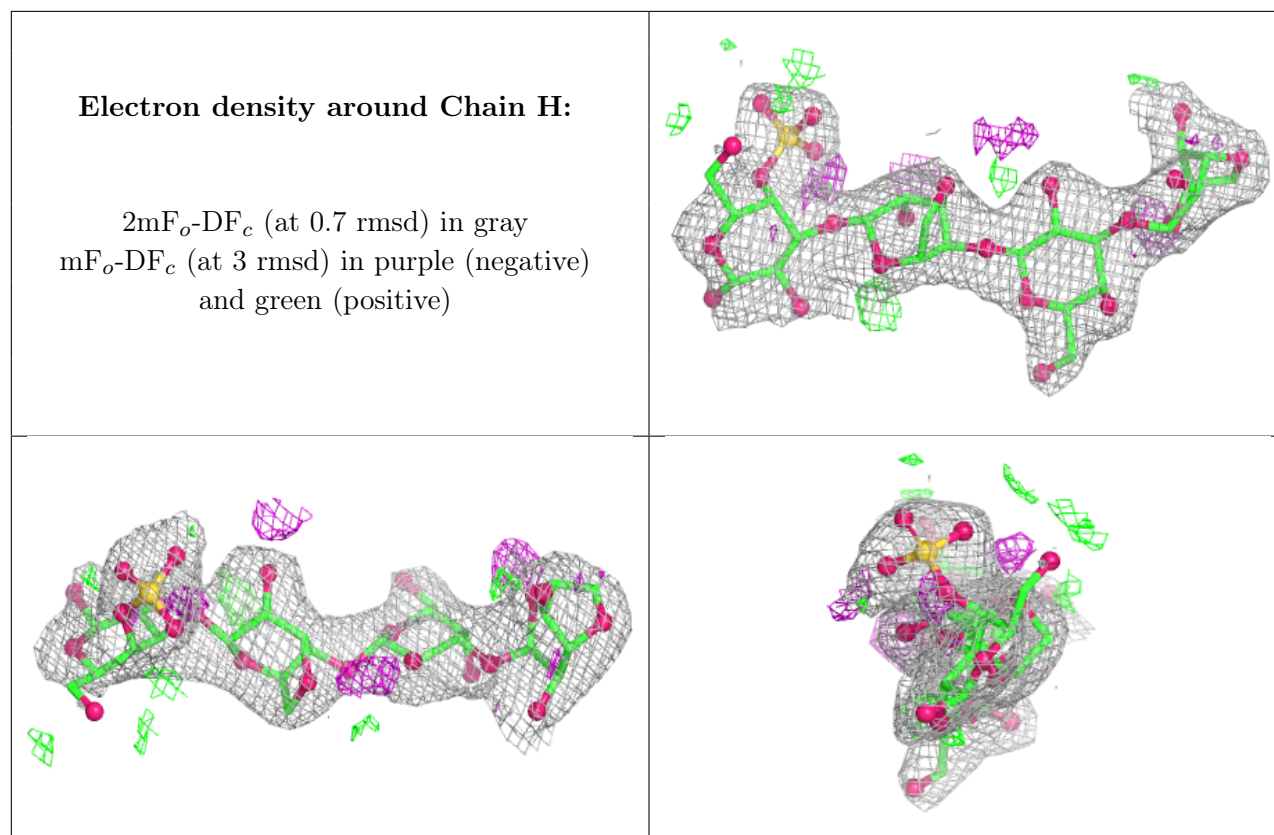
Electron density around Chain F:

$2mF_o-DF_c$ (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(Å ²)	Q<0.9
4	NA	A	703	1/1	0.64	0.30	80,80,80,80	0
4	NA	C	702	1/1	0.67	0.23	79,79,79,79	0
4	NA	B	703	1/1	0.71	0.20	78,78,78,78	0
3	TLA	A	702	10/10	0.76	0.18	66,69,72,72	0
4	NA	D	702	1/1	0.76	0.18	68,68,68,68	0
3	TLA	A	701	10/10	0.91	0.08	40,45,50,52	0
3	TLA	D	701	10/10	0.92	0.08	46,52,60,62	0
3	TLA	C	701	10/10	0.94	0.08	47,51,58,58	0
3	TLA	B	701	10/10	0.95	0.07	46,52,59,60	0
5	CL	A	704	1/1	0.95	0.09	55,55,55,55	0
6	EDO	B	702	4/4	0.95	0.12	53,55,55,56	0
5	CL	D	703	1/1	0.97	0.08	61,61,61,61	0
5	CL	B	704	1/1	0.97	0.07	57,57,57,57	0
5	CL	C	703	1/1	0.98	0.06	58,58,58,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.