



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 10, 2020 – 07:47 AM BST

PDB ID : 6MHM
Title : Crystal structure of human acid ceramidase in covalent complex with carmofur
Authors : Dementiev, A.; Joachimiak, A.; Doan, N.
Deposited on : 2018-09-18
Resolution : 2.74 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

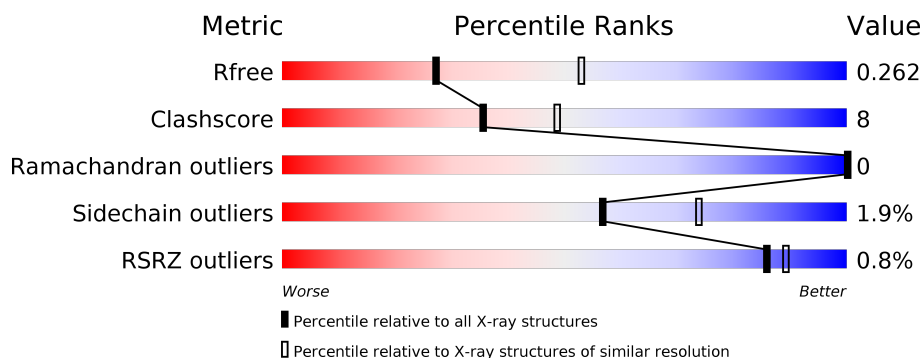
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.74 Å.

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



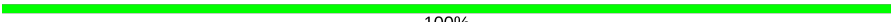
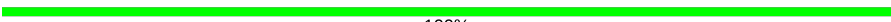
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1271 (2.76-2.72)
Clashscore	141614	1322 (2.76-2.72)
Ramachandran outliers	138981	1297 (2.76-2.72)
Sidechain outliers	138945	1298 (2.76-2.72)
RSRZ outliers	127900	1243 (2.76-2.72)

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

Mol	Chain	Length	Quality of chain
1	A	131	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 10%, orange 10%, yellow 10%, green 70%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 10% 10% 10% 74% </div> </div>
1	C	131	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 10%, orange 10%, yellow 10%, green 70%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 10% 10% 10% 74% </div> </div>
2	B	253	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 10%, orange 10%, yellow 10%, green 70%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 10% 10% 10% 79% </div> </div>
2	D	253	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 10%, orange 10%, yellow 10%, green 70%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 10% 10% 10% 79% </div> </div>
3	E	3	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 10%, orange 10%, yellow 10%, green 70%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 10% 10% 10% 33% </div> </div>
3	G	3	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red 10%, orange 10%, yellow 10%, green 70%);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> 10% 10% 10% 67% </div> </div>

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Mol	Chain	Length	Quality of chain
4	F	2	 100%
4	H	2	 100%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	SO4	B	409	-	-	X	-
7	JRY	B	407	-	X	-	-
7	JRY	D	408	-	X	-	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 6208 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 Acid ceramidase subunit alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	114	Total	C	N	O	S	0	0	0
			910	599	142	163	6			
1	C	114	Total	C	N	O	S	0	0	0
			906	596	141	163	6			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	12	ASP	-	expression tag	UNP Q13510
A	13	ARG	-	expression tag	UNP Q13510
A	14	HIS	-	expression tag	UNP Q13510
A	15	HIS	-	expression tag	UNP Q13510
A	16	HIS	-	expression tag	UNP Q13510
A	17	HIS	-	expression tag	UNP Q13510
A	18	HIS	-	expression tag	UNP Q13510
A	19	HIS	-	expression tag	UNP Q13510
A	20	LYS	-	expression tag	UNP Q13510
A	21	LEU	-	expression tag	UNP Q13510
A	72	MET	VAL	variant	UNP Q13510
A	93	VAL	ILE	variant	UNP Q13510
C	12	ASP	-	expression tag	UNP Q13510
C	13	ARG	-	expression tag	UNP Q13510
C	14	HIS	-	expression tag	UNP Q13510
C	15	HIS	-	expression tag	UNP Q13510
C	16	HIS	-	expression tag	UNP Q13510
C	17	HIS	-	expression tag	UNP Q13510
C	18	HIS	-	expression tag	UNP Q13510
C	19	HIS	-	expression tag	UNP Q13510
C	20	LYS	-	expression tag	UNP Q13510
C	21	LEU	-	expression tag	UNP Q13510
C	72	MET	VAL	variant	UNP Q13510
C	93	VAL	ILE	variant	UNP Q13510

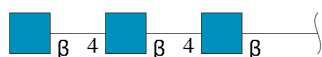
- Molecule 2 is a protein called Acid ceramidase subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	253	Total	C	N	O	S	0	0	0
			2033	1306	341	376	10			
2	D	253	Total	C	N	O	S	0	0	0
			2015	1294	335	376	10			

There are 2 discrepancies between the modelled and reference sequences:

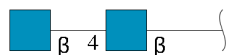
Chain	Residue	Modelled	Actual	Comment	Reference
B	246	ALA	VAL	variant	UNP Q13510
D	246	ALA	VAL	variant	UNP Q13510

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	E	3	Total	C	N	O	0	0	0
			42	24	3	15			
3	G	3	Total	C	N	O	0	0	0
			42	24	3	15			

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



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

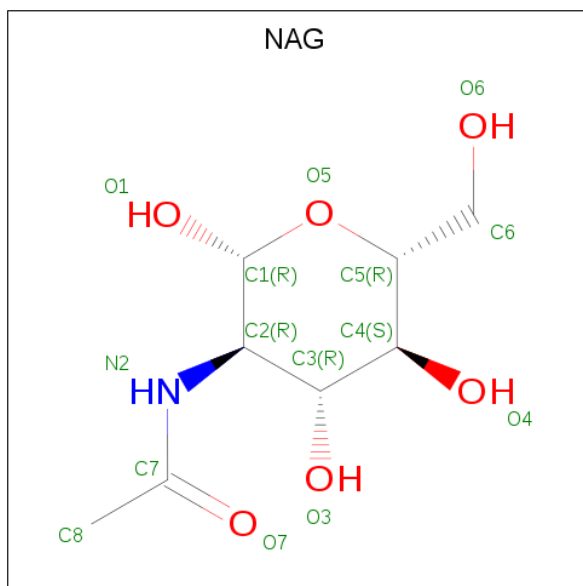
- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

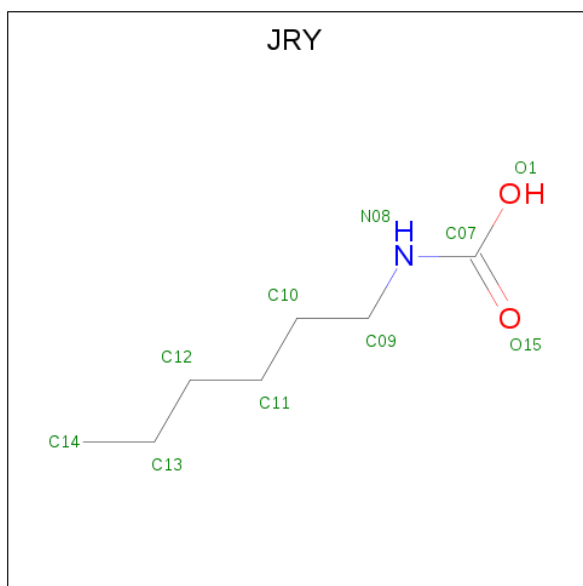
- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula:

C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	N	O	0	0
			14	8	1	5		
6	D	1	Total	C	N	O	0	0
			14	8	1	5		
6	D	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 7 is hexylcarbamic acid (three-letter code: JRY) (formula: C₇H₁₅NO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	B	1	Total	C	N	O	0	0
			9	7	1	1		
7	D	1	Total	C	N	O	0	0
			9	7	1	1		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	13	Total	O	0	0
			13	13		
8	B	29	Total	O	0	0
			29	29		
8	C	12	Total	O	0	0
			12	12		
8	D	25	Total	O	0	0
			25	25		

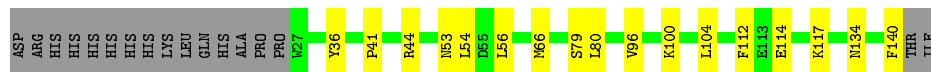
3 Residue-property plots

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.

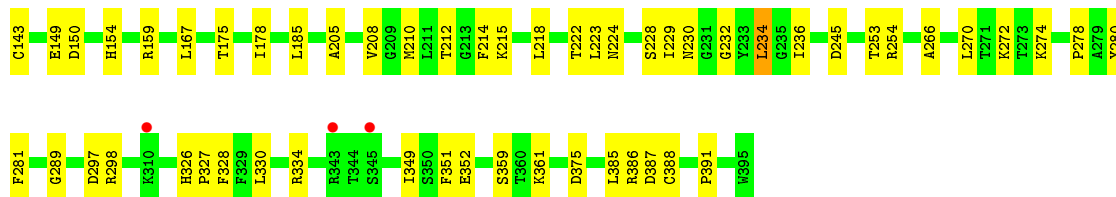
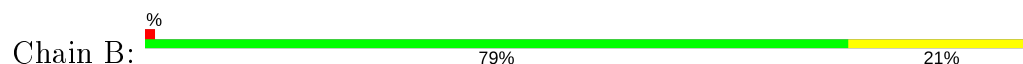
- Molecule 1: Acid ceramidase subunit alpha



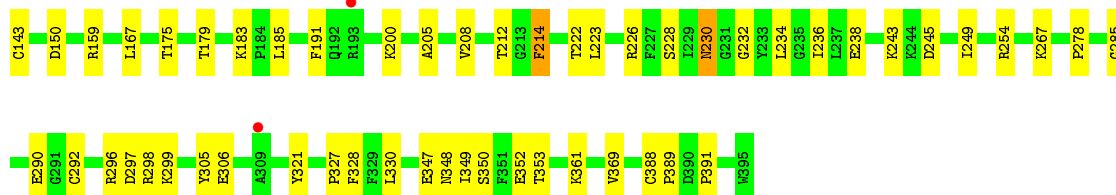
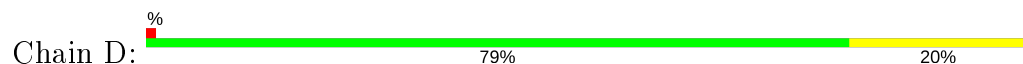
- Molecule 1: Acid ceramidase subunit alpha



- Molecule 2: Acid ceramidase subunit beta



- Molecule 2: Acid ceramidase subunit beta



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:  33% 33% 33%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:  67% 33%



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

Chain F:  100%



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

Chain H:  100%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	153.72Å 68.65Å 98.39Å 90.00° 120.73° 90.00°	Depositor
Resolution (Å)	48.65 – 2.74 48.65 – 2.74	Depositor EDS
% Data completeness (in resolution range)	99.7 (48.65-2.74) 94.4 (48.65-2.74)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.41 (at 2.73Å)	Xtriage
Refinement program	PHENIX 1.11.1 _2575	Depositor
R, R_{free}	0.214 , 0.262 0.214 , 0.262	Depositor DCC
R_{free} test set	2000 reflections (8.60%)	wwPDB-VP
Wilson B-factor (Å ²)	49.7	Xtriage
Anisotropy	0.362	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 52.1	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	0.92	EDS
Total number of atoms	6208	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.54% 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: NAG, JRY, SO4

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	0.39	0/939	0.48	0/1281
1	C	0.27	0/935	0.43	0/1277
2	B	0.38	1/2081 (0.0%)	0.51	0/2821
2	D	0.35	1/2063 (0.0%)	0.50	0/2802
All	All	0.36	2/6018 (0.0%)	0.49	0/8181

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	143	CYS	CB-SG	-10.12	1.65	1.82
2	D	143	CYS	CB-SG	-6.85	1.70	1.82

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	910	0	904	16	0
1	C	906	0	893	15	0
2	B	2033	0	2000	43	0
2	D	2015	0	1952	40	0
3	E	42	0	37	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	G	42	0	37	3	0
4	F	28	0	25	0	0
4	H	28	0	25	0	0
5	A	15	0	0	1	0
5	B	15	0	0	3	0
5	C	15	0	0	0	0
5	D	20	0	0	2	0
6	B	14	0	13	0	0
6	D	28	0	26	0	0
7	B	9	0	0	0	0
7	D	9	0	0	0	0
8	A	13	0	0	0	0
8	B	29	0	0	1	0
8	C	12	0	0	0	0
8	D	25	0	0	0	0
All	All	6208	0	5912	97	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 8.

All (97) 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:D:230:ASN:OD1	2:D:234:LEU:HD12	1.62	0.99
2:B:229:ILE:HD12	2:B:230:ASN:H	1.30	0.94
1:C:36:TYR:HB3	2:D:352:GLU:HG2	1.49	0.94
2:B:352:GLU:N	2:B:352:GLU:OE2	2.03	0.89
2:B:334:ARG:NH1	8:B:501:HOH:O	2.06	0.86
2:B:229:ILE:HD12	2:B:230:ASN:N	1.98	0.77
1:C:140:PHE:HE2	2:D:232:GLY:H	1.33	0.76
1:A:36:TYR:CD2	2:B:352:GLU:HG3	2.26	0.70
2:D:298:ARG:NH1	5:D:412:SO4:O1	2.25	0.70
1:A:95:GLN:OE1	5:A:203:SO4:O4	2.09	0.69
1:A:140:PHE:HE2	2:B:232:GLY:H	1.41	0.67
2:B:274:LYS:NZ	5:B:409:SO4:O4	2.29	0.66
2:B:327:PRO:HG3	2:B:334:ARG:NH2	2.11	0.64
2:D:230:ASN:OD1	2:D:234:LEU:CD1	2.43	0.64
2:B:359:SER:OG	2:B:386:ARG:NH2	2.30	0.64
1:C:53:ASN:HB3	1:C:56:LEU:HD13	1.80	0.64
1:A:36:TYR:HD2	2:B:352:GLU:HG3	1.61	0.64
1:C:80:LEU:HD23	2:D:167:LEU:HD21	1.78	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:215:LYS:HD3	2:B:218:LEU:HD12	1.81	0.62
2:D:228:SER:HB2	2:D:278:PRO:HG3	1.81	0.62
1:A:27:TRP:CE3	1:A:27:TRP:HA	2.33	0.61
1:A:91:GLY:O	1:A:95:GLN:HG2	2.01	0.61
2:D:226:ARG:NH1	2:D:296:ARG:O	2.34	0.60
1:A:42:THR:HG23	2:B:386:ARG:HH21	1.68	0.58
2:D:223:LEU:HD11	2:D:249:ILE:CG2	2.33	0.58
2:B:234:LEU:HD12	2:B:234:LEU:O	2.04	0.58
2:B:298:ARG:NH1	5:B:408:SO4:O2	2.37	0.58
2:D:223:LEU:C	2:D:223:LEU:HD23	2.25	0.57
2:B:224:ASN:HB2	2:B:280:TYR:HB2	1.86	0.57
2:B:149:GLU:OE2	2:B:289:GLY:N	2.35	0.56
2:B:274:LYS:NZ	5:B:409:SO4:S	2.79	0.56
2:B:212:THR:HG23	2:B:222:THR:HG22	1.87	0.55
2:B:228:SER:HB2	2:B:278:PRO:HG3	1.89	0.55
1:C:66:MET:HE1	2:D:205:ALA:HB1	1.88	0.54
1:A:84:ILE:HD13	1:A:93:VAL:HG11	1.90	0.54
2:B:328:PHE:CE2	2:B:330:LEU:HB2	2.42	0.54
2:D:347:GLU:HG3	2:D:348:ASN:OD1	2.08	0.54
1:A:79:SER:HB3	2:B:167:LEU:HD23	1.89	0.54
2:B:351:PHE:CE1	2:B:375:ASP:HB2	2.43	0.54
2:D:285:GLY:HA3	2:D:290:GLU:HB3	1.90	0.54
2:B:210:MET:O	2:B:253:THR:OG1	2.19	0.53
2:B:232:GLY:O	2:B:236:ILE:HG13	2.09	0.52
2:D:175:THR:HG21	3:G:1:NAG:H82	1.91	0.52
2:D:243:LYS:HA	2:D:243:LYS:HE2	1.92	0.51
2:B:223:LEU:HD12	2:B:281:PHE:CE1	2.46	0.51
2:B:167:LEU:HD22	2:B:178:ILE:HG12	1.93	0.51
1:C:140:PHE:HE2	2:D:232:GLY:N	2.05	0.51
2:D:245:ASP:N	2:D:245:ASP:OD1	2.44	0.51
1:C:79:SER:HB3	2:D:167:LEU:HD13	1.91	0.50
2:D:299:LYS:NZ	5:D:410:SO4:O1	2.36	0.50
2:D:232:GLY:O	2:D:236:ILE:HG13	2.12	0.50
2:B:167:LEU:HB2	2:B:178:ILE:HG12	1.93	0.49
2:B:361:LYS:NZ	2:B:391:PRO:O	2.41	0.49
2:D:234:LEU:O	2:D:238:GLU:HG3	2.12	0.49
2:D:185:LEU:HD22	2:D:205:ALA:HB3	1.94	0.49
2:D:350:SER:N	2:D:353:THR:OG1	2.46	0.48
2:B:266:ALA:O	2:B:270:LEU:HD12	2.13	0.48
2:D:230:ASN:OD1	2:D:234:LEU:HB2	2.13	0.48
1:A:42:THR:CG2	2:B:386:ARG:HH21	2.26	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:114:GLU:HA	1:C:117:LYS:HB2	1.95	0.47
2:B:245:ASP:N	2:B:245:ASP:OD2	2.39	0.46
1:C:54:LEU:HD21	2:D:191:PHE:HD2	1.81	0.46
2:D:212:THR:HG23	2:D:222:THR:HG22	1.98	0.46
2:D:267:LYS:NZ	2:D:306:GLU:OE1	2.33	0.46
2:B:351:PHE:HE1	2:B:375:ASP:HB2	1.79	0.45
2:B:150:ASP:OD1	2:B:154:HIS:N	2.44	0.45
1:C:44:ARG:NH2	3:G:1:NAG:H83	2.31	0.45
2:D:361:LYS:NZ	2:D:391:PRO:O	2.34	0.45
1:A:118:GLY:O	1:A:122:VAL:HG23	2.18	0.44
2:D:150:ASP:HB3	2:D:349:ILE:HG22	2.00	0.44
2:B:272:LYS:HA	2:B:272:LYS:HD3	1.81	0.43
1:C:96:VAL:HG13	1:C:100:LYS:HD2	2.00	0.43
2:D:200:LYS:HB2	2:D:214:PHE:HE1	1.82	0.43
1:A:75:VAL:HG21	3:E:3:NAG:H5	1.99	0.43
3:E:3:NAG:O7	3:E:3:NAG:C3	2.67	0.43
2:B:234:LEU:HD12	2:B:234:LEU:C	2.39	0.43
1:A:112:PHE:HE1	2:B:254:ARG:HG2	1.84	0.42
2:D:285:GLY:N	2:D:290:GLU:O	2.51	0.42
2:B:175:THR:HG21	3:E:1:NAG:H82	2.02	0.42
1:C:100:LYS:O	1:C:104:LEU:HB2	2.20	0.42
2:D:321:TYR:CD2	2:D:327:PRO:HB3	2.54	0.42
2:B:349:ILE:HD12	2:B:349:ILE:HA	1.90	0.42
3:G:3:NAG:O3	3:G:3:NAG:H82	2.19	0.42
2:B:185:LEU:HD22	2:B:205:ALA:HB3	2.01	0.41
2:D:292:CYS:HA	2:D:305:TYR:O	2.19	0.41
1:C:134:ASN:OD1	2:D:208:VAL:HG22	2.21	0.41
2:D:183:LYS:HG3	2:D:369:VAL:HG13	2.02	0.41
1:C:41:PRO:HB2	2:D:389:PRO:HG2	2.03	0.41
2:B:229:ILE:CD1	2:B:230:ASN:N	2.77	0.41
2:B:270:LEU:HB3	2:B:281:PHE:CD2	2.55	0.41
1:A:97:VAL:O	1:A:101:LEU:HB3	2.21	0.41
2:D:328:PHE:CE2	2:D:330:LEU:HB2	2.56	0.41
1:A:134:ASN:OD1	2:B:208:VAL:HG22	2.21	0.40
2:D:296:ARG:HD2	2:D:296:ARG:HA	1.90	0.40
1:A:47:VAL:HG12	2:B:385:LEU:HG	2.02	0.40
1:C:112:PHE:HE1	2:D:254:ARG:HG2	1.87	0.40
2:D:179:THR:HG23	2:D:369:VAL:HG22	2.03	0.40

There are no symmetry-related clashes.

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	112/131 (86%)	110 (98%)	2 (2%)	0	100	100
1	C	112/131 (86%)	111 (99%)	1 (1%)	0	100	100
2	B	251/253 (99%)	241 (96%)	10 (4%)	0	100	100
2	D	251/253 (99%)	244 (97%)	7 (3%)	0	100	100
All	All	726/768 (94%)	706 (97%)	20 (3%)	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	101/118 (86%)	101 (100%)	0	100	100
1	C	100/118 (85%)	100 (100%)	0	100	100
2	B	222/224 (99%)	215 (97%)	7 (3%)	39	59
2	D	218/224 (97%)	213 (98%)	5 (2%)	50	70
All	All	641/684 (94%)	629 (98%)	12 (2%)	57	74

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

Mol	Chain	Res	Type
2	B	159	ARG
2	B	214	PHE

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Mol	Chain	Res	Type
2	B	234	LEU
2	B	297	ASP
2	B	326	HIS
2	B	387	ASP
2	B	388	CYS
2	D	159	ARG
2	D	214	PHE
2	D	230	ASN
2	D	297	ASP
2	D	388	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	287	GLN
2	B	318	GLN
2	B	346	GLN
2	B	380	GLN
1	C	53	ASN
1	C	107	ASN
2	D	287	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 ⓘ

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	$\# Z > 2$
3	NAG	E	1	3,2	14,14,15	0.36	0	17,19,21	0.62	0
3	NAG	E	2	3	14,14,15	0.23	0	17,19,21	0.63	0
3	NAG	E	3	3	14,14,15	0.48	0	17,19,21	1.22	2 (11%)
4	NAG	F	1	2,4	14,14,15	0.54	0	17,19,21	0.55	0
4	NAG	F	2	4	14,14,15	0.24	0	17,19,21	0.53	0
3	NAG	G	1	3,2	14,14,15	0.34	0	17,19,21	0.54	0
3	NAG	G	2	3	14,14,15	0.28	0	17,19,21	0.82	1 (5%)
3	NAG	G	3	3	14,14,15	0.70	1 (7%)	17,19,21	0.57	0
4	NAG	H	1	2,4	14,14,15	0.44	0	17,19,21	0.50	0
4	NAG	H	2	4	14,14,15	0.21	0	17,19,21	0.58	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
3	NAG	E	1	3,2	-	2/6/23/26	0/1/1/1
3	NAG	E	2	3	-	2/6/23/26	0/1/1/1
3	NAG	E	3	3	-	3/6/23/26	0/1/1/1
4	NAG	F	1	2,4	-	2/6/23/26	0/1/1/1
4	NAG	F	2	4	-	0/6/23/26	0/1/1/1
3	NAG	G	1	3,2	-	2/6/23/26	0/1/1/1
3	NAG	G	2	3	-	2/6/23/26	0/1/1/1
3	NAG	G	3	3	-	4/6/23/26	0/1/1/1
4	NAG	H	1	2,4	-	2/6/23/26	0/1/1/1
4	NAG	H	2	4	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	3	NAG	C1-C2	2.32	1.55	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	E	3	NAG	C2-N2-C7	4.02	128.62	122.90
3	E	3	NAG	C1-O5-C5	2.08	115.01	112.19
3	G	2	NAG	C1-O5-C5	2.00	114.90	112.19

There are no chirality outliers.

All (21) torsion outliers are listed below:

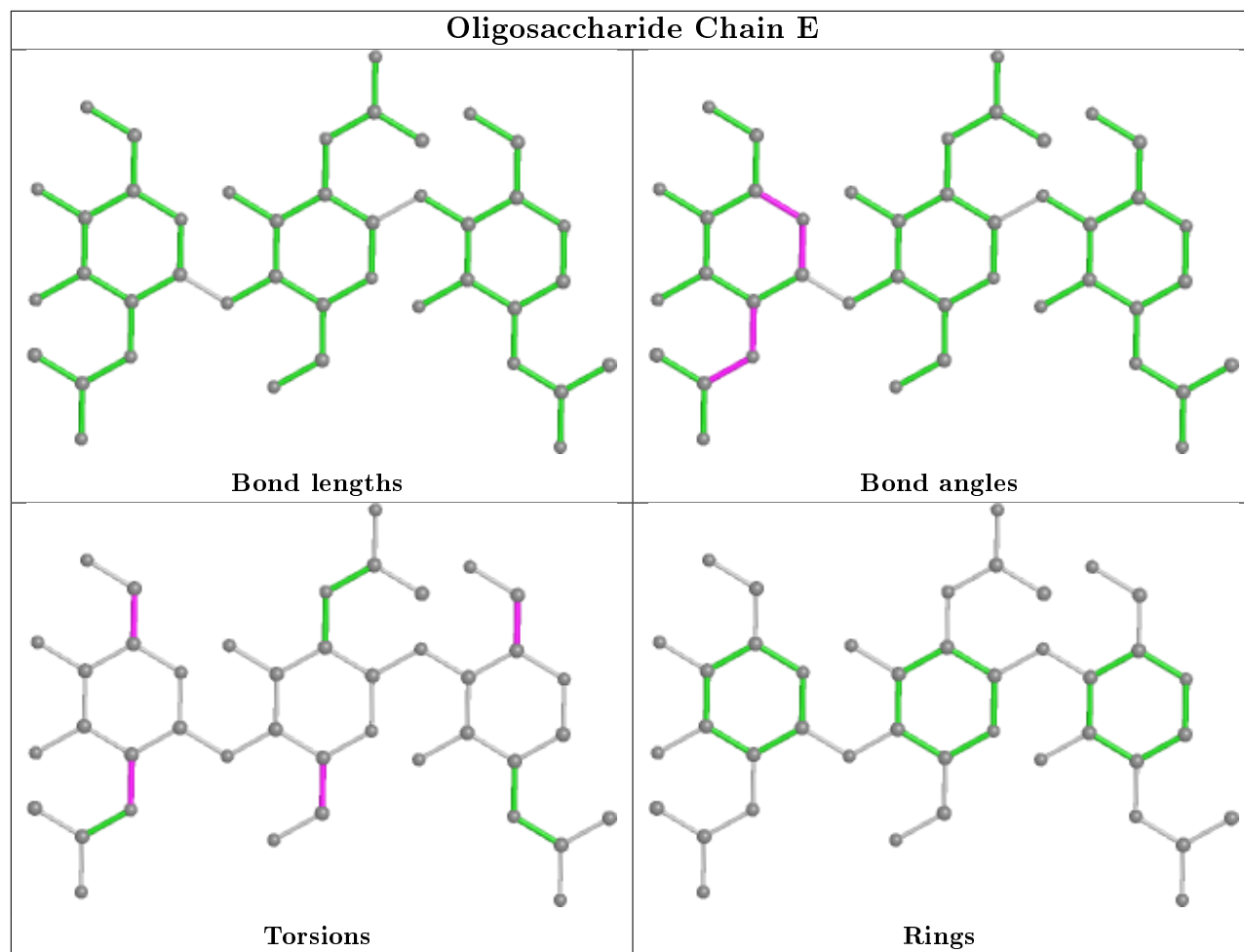
Mol	Chain	Res	Type	Atoms
3	E	3	NAG	C3-C2-N2-C7
4	H	1	NAG	O5-C5-C6-O6
3	G	3	NAG	O5-C5-C6-O6
4	F	1	NAG	O5-C5-C6-O6
3	E	3	NAG	C4-C5-C6-O6
4	H	1	NAG	C4-C5-C6-O6
3	G	3	NAG	C8-C7-N2-C2
3	G	3	NAG	O7-C7-N2-C2
3	E	2	NAG	O5-C5-C6-O6
4	F	1	NAG	C4-C5-C6-O6
3	G	2	NAG	O5-C5-C6-O6
3	E	2	NAG	C4-C5-C6-O6
3	G	3	NAG	C4-C5-C6-O6
3	E	3	NAG	O5-C5-C6-O6
3	E	1	NAG	O5-C5-C6-O6
3	E	1	NAG	C4-C5-C6-O6
3	G	1	NAG	C4-C5-C6-O6
3	G	2	NAG	C4-C5-C6-O6
3	G	1	NAG	O5-C5-C6-O6
4	H	2	NAG	O5-C5-C6-O6
4	H	2	NAG	C4-C5-C6-O6

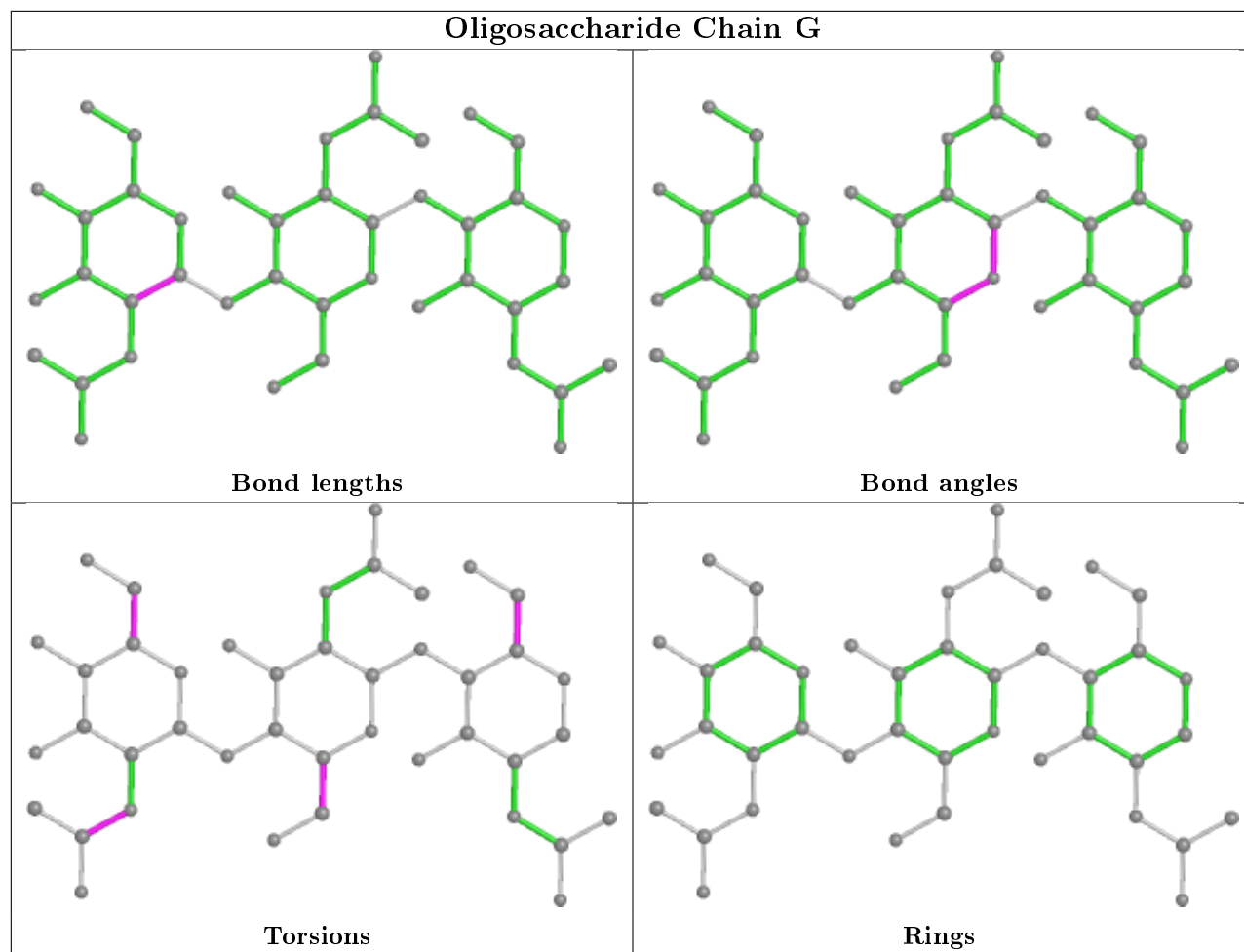
There are no ring outliers.

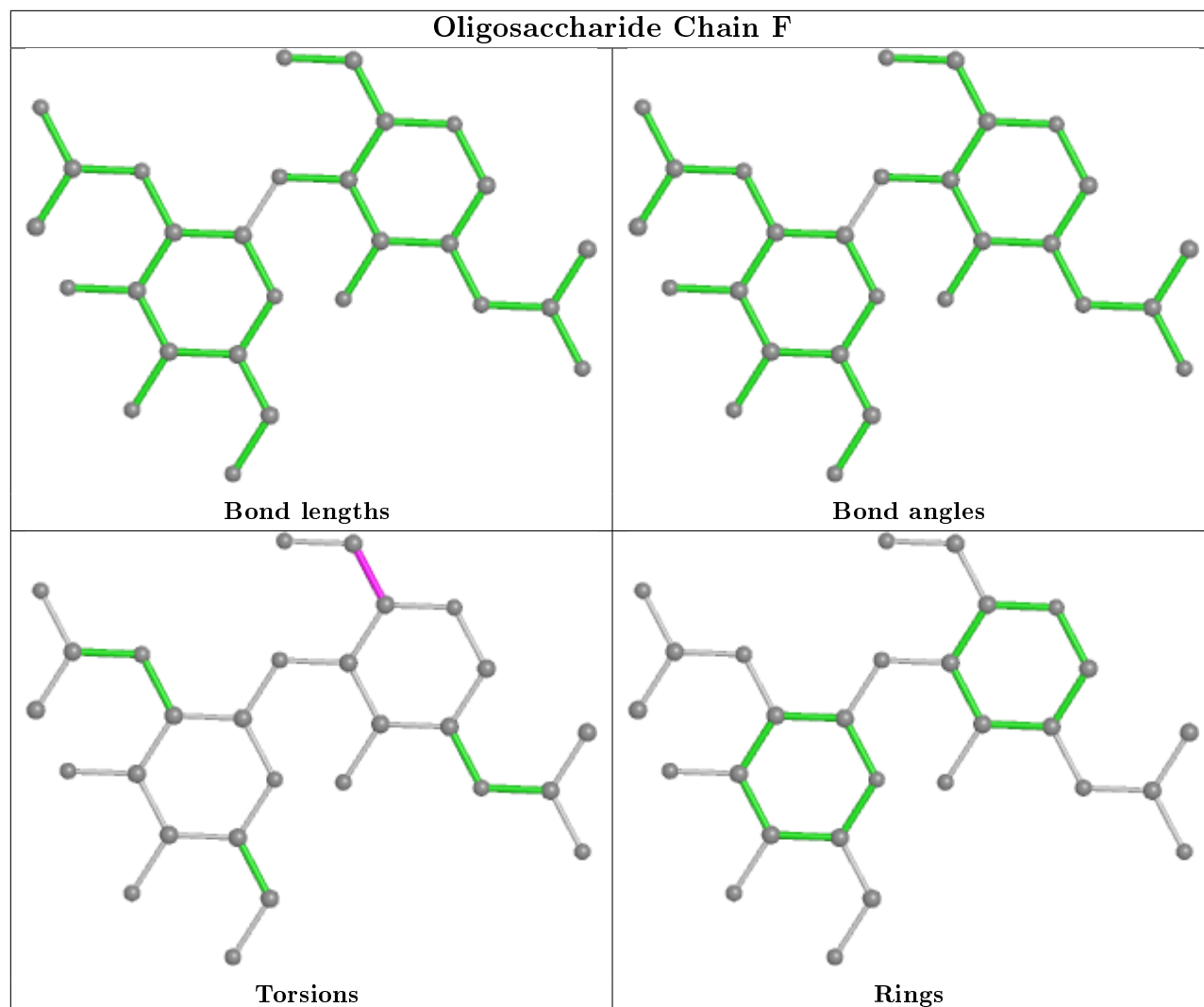
4 monomers are involved in 6 short contacts:

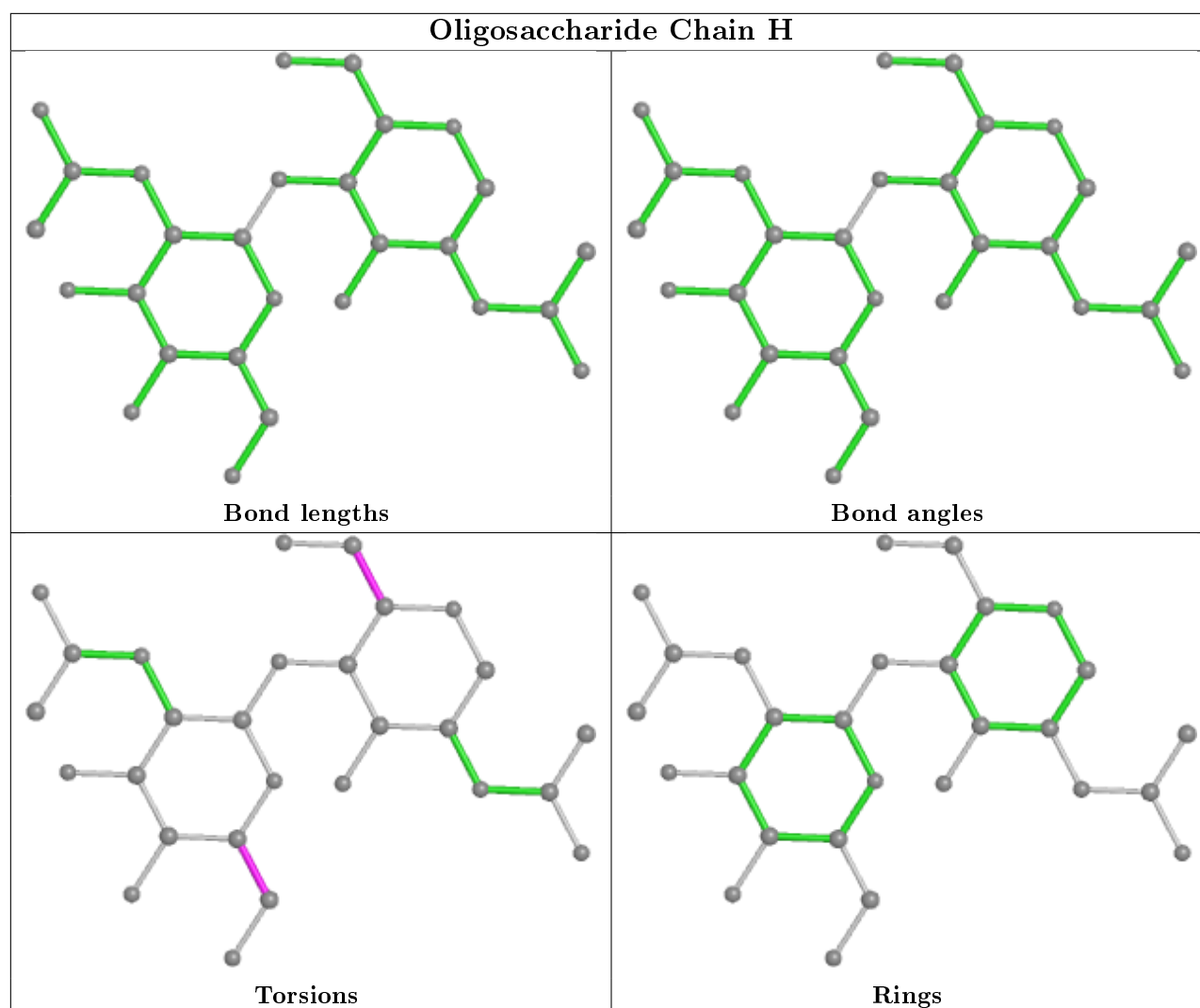
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	1	NAG	1	0
3	G	1	NAG	2	0
3	G	3	NAG	1	0
3	E	3	NAG	2	0

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](#)

18 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	$\# Z > 2$
5	SO4	C	203	-	4,4,4	0.17	0	6,6,6	0.09	0
5	SO4	B	409	-	4,4,4	0.15	0	6,6,6	0.06	0
6	NAG	D	404	2	14,14,15	0.30	0	17,19,21	0.35	0
5	SO4	D	412	-	4,4,4	0.14	0	6,6,6	0.06	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	SO4	B	408	-	4,4,4	0.14	0	6,6,6	0.10	0
5	SO4	C	201	-	4,4,4	0.14	0	6,6,6	0.05	0
5	SO4	D	411	-	4,4,4	0.13	0	6,6,6	0.05	0
6	NAG	D	407	2	14,14,15	1.05	1 (7%)	17,19,21	0.92	1 (5%)
5	SO4	A	202	-	4,4,4	0.14	0	6,6,6	0.16	0
5	SO4	C	202	-	4,4,4	0.13	0	6,6,6	0.07	0
6	NAG	B	406	2	14,14,15	0.28	0	17,19,21	0.56	0
5	SO4	A	201	-	4,4,4	0.14	0	6,6,6	0.05	0
7	JRY	D	408	2	8,8,9	1.85	1 (12%)	6,7,9	6.12	4 (66%)
7	JRY	B	407	2	8,8,9	1.99	3 (37%)	6,7,9	7.49	4 (66%)
5	SO4	D	410	-	4,4,4	0.14	0	6,6,6	0.04	0
5	SO4	A	203	-	4,4,4	0.39	0	6,6,6	0.29	0
5	SO4	B	410	-	4,4,4	0.13	0	6,6,6	0.09	0
5	SO4	D	409	-	4,4,4	0.13	0	6,6,6	0.10	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
6	NAG	B	406	2	-	1/6/23/26	0/1/1/1
6	NAG	D	407	2	-	2/6/23/26	0/1/1/1
6	NAG	D	404	2	-	2/6/23/26	0/1/1/1
7	JRY	D	408	2	-	4/6/6/7	-
7	JRY	B	407	2	-	5/6/6/7	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	D	408	JRY	C07-N08	4.43	1.42	1.33
7	B	407	JRY	C07-N08	4.37	1.42	1.33
6	D	407	NAG	C1-C2	3.58	1.57	1.52
7	B	407	JRY	C09-N08	-2.53	1.40	1.46
7	B	407	JRY	O15-C07	-2.43	1.15	1.22

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	407	JRY	O15-C07-N08	-16.79	107.93	124.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	D	408	JRY	O15-C07-N08	-13.24	111.52	124.89
7	B	407	JRY	C10-C09-N08	6.39	136.17	112.84
7	D	408	JRY	C10-C09-N08	5.40	132.55	112.84
7	D	408	JRY	C11-C10-C09	3.71	131.14	113.56
7	B	407	JRY	C11-C10-C09	2.72	126.44	113.56
6	D	407	NAG	C4-C3-C2	2.54	114.73	111.02
7	B	407	JRY	C12-C11-C10	-2.48	101.86	114.42
7	D	408	JRY	C12-C11-C10	-2.45	101.97	114.42

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	D	404	NAG	O5-C5-C6-O6
7	D	408	JRY	N08-C09-C10-C11
6	D	407	NAG	O5-C5-C6-O6
7	D	408	JRY	C10-C11-C12-C13
7	D	408	JRY	C10-C09-N08-C07
7	B	407	JRY	C10-C09-N08-C07
6	D	404	NAG	C4-C5-C6-O6
7	B	407	JRY	C09-C10-C11-C12
7	D	408	JRY	O15-C07-N08-C09
7	B	407	JRY	O15-C07-N08-C09
7	B	407	JRY	C10-C11-C12-C13
6	D	407	NAG	C4-C5-C6-O6
6	B	406	NAG	C3-C2-N2-C7
7	B	407	JRY	N08-C09-C10-C11

There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	409	SO4	2	0
5	D	412	SO4	1	0
5	B	408	SO4	1	0
5	D	410	SO4	1	0
5	A	203	SO4	1	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	114/131 (87%)	0.09	1 (0%) 84 88	38, 52, 71, 84	0
1	C	114/131 (87%)	-0.02	0 100 100	36, 53, 69, 80	0
2	B	253/253 (100%)	0.02	3 (1%) 79 83	32, 46, 67, 82	0
2	D	253/253 (100%)	0.10	2 (0%) 86 89	34, 47, 68, 82	0
All	All	734/768 (95%)	0.05	6 (0%) 86 89	32, 49, 69, 84	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	309	ALA	2.9
1	A	60	LYS	2.9
2	B	310	LYS	2.7
2	D	193	ARG	2.5
2	B	343	ARG	2.3
2	B	345	SER	2.1

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.

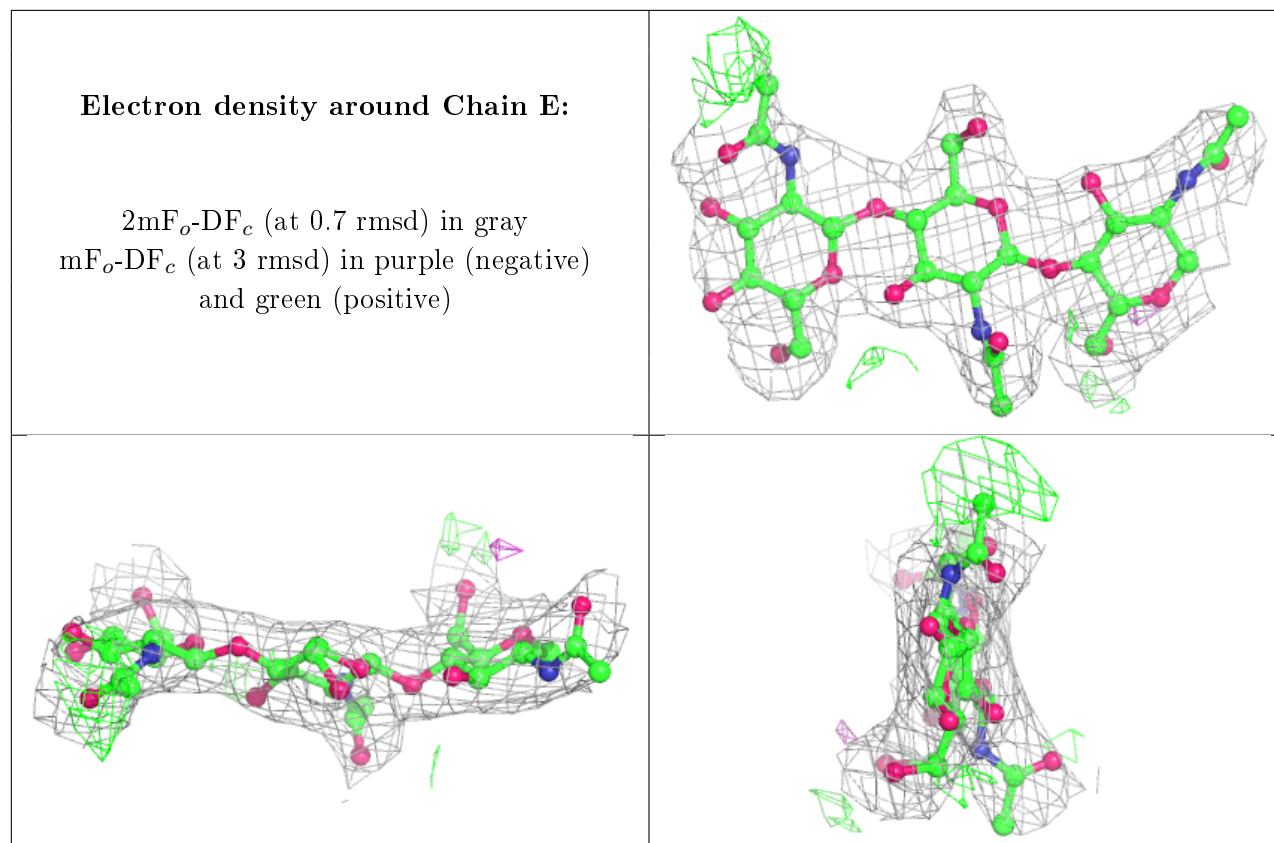
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	NAG	F	2	14/15	0.67	0.25	62,82,84,85	0

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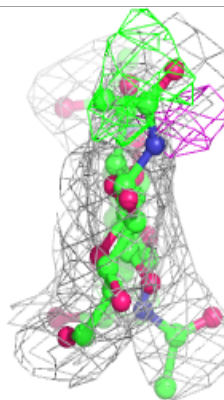
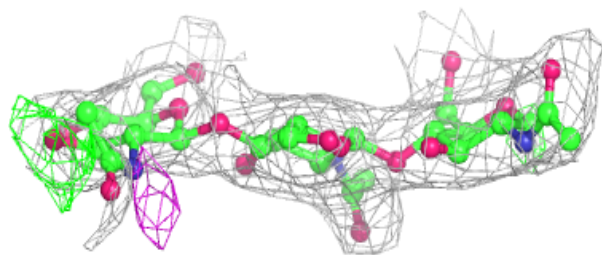
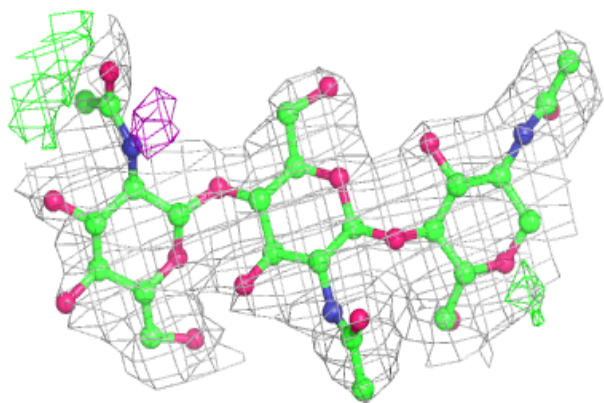
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	NAG	G	3	14/15	0.73	0.27	61,66,76,78	0
4	NAG	H	2	14/15	0.77	0.24	65,80,84,90	0
3	NAG	E	3	14/15	0.80	0.20	56,62,70,71	0
4	NAG	H	1	14/15	0.92	0.12	53,67,72,78	0
4	NAG	F	1	14/15	0.93	0.14	51,71,80,86	0
3	NAG	G	1	14/15	0.94	0.16	34,37,47,49	0
3	NAG	E	1	14/15	0.95	0.15	37,43,48,50	0
3	NAG	E	2	14/15	0.95	0.15	39,49,56,58	0
3	NAG	G	2	14/15	0.95	0.12	39,49,57,63	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.



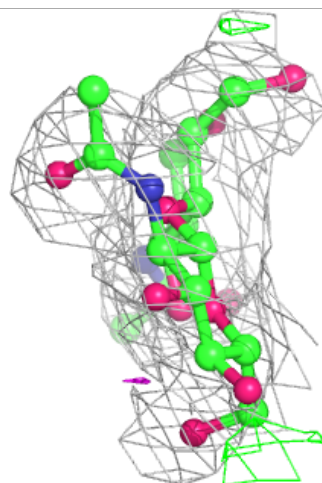
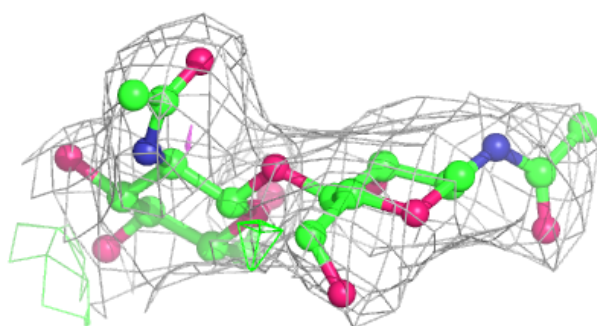
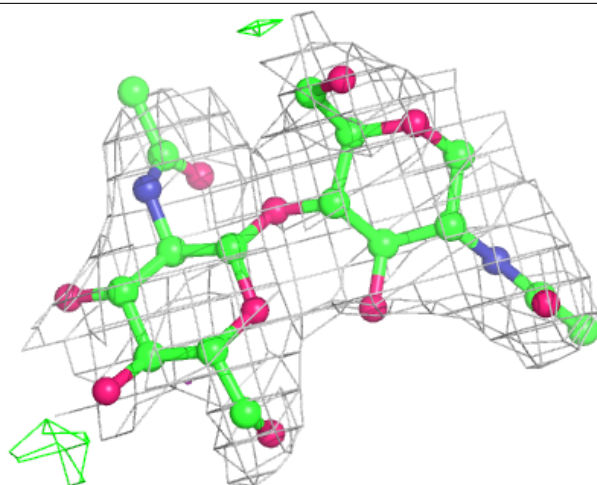
Electron density around Chain G:

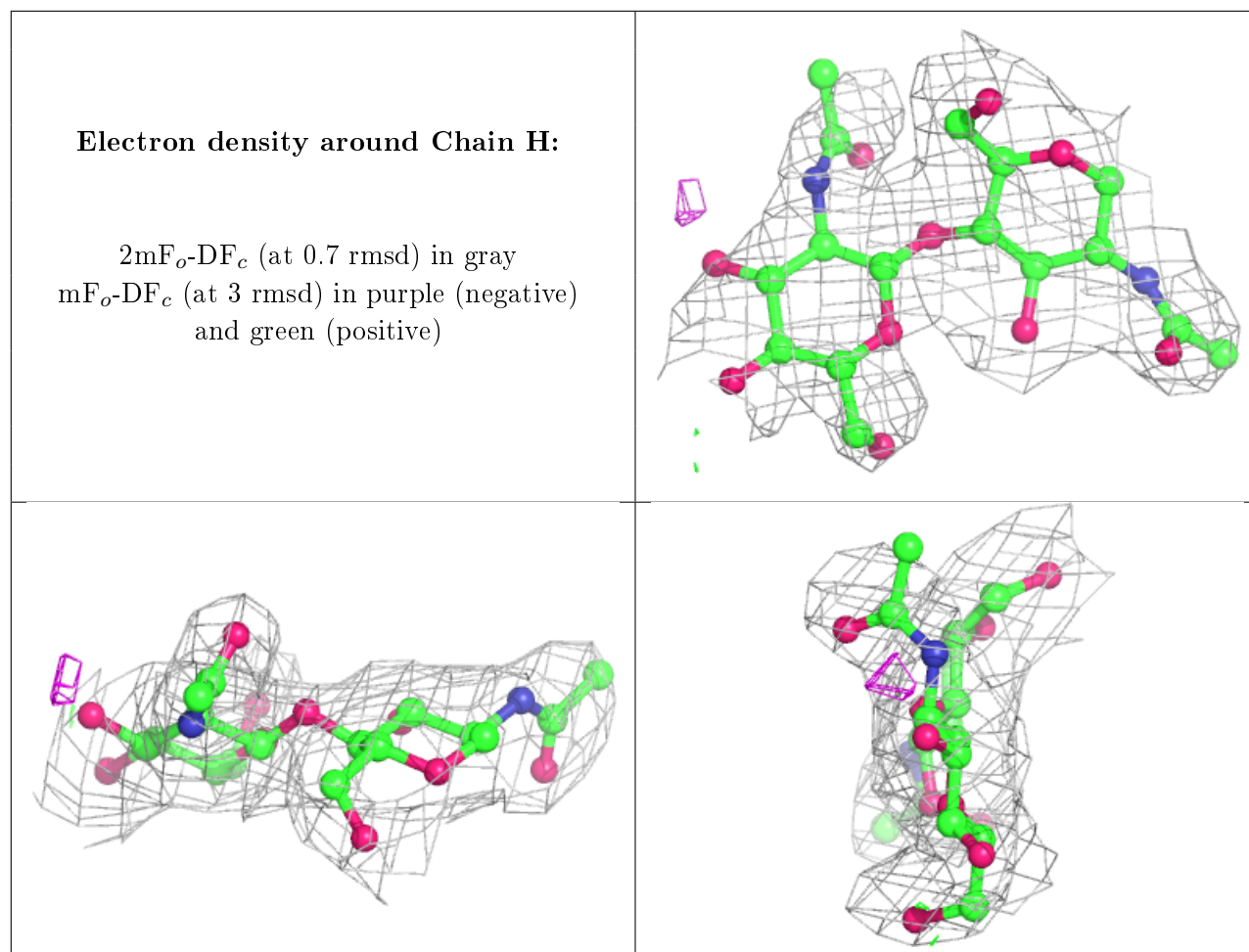
$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain F:

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





6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	SO4	A	201	5/5	0.73	0.30	77,78,88,98	0
5	SO4	D	412	5/5	0.74	0.21	85,85,90,106	0
6	NAG	D	407	14/15	0.75	0.28	70,82,89,91	0
5	SO4	B	410	5/5	0.75	0.26	79,81,85,102	0
5	SO4	D	411	5/5	0.81	0.18	94,95,108,113	0
5	SO4	A	202	5/5	0.81	0.31	80,91,106,108	0
6	NAG	B	406	14/15	0.84	0.17	70,79,84,84	0
6	NAG	D	404	14/15	0.85	0.28	73,83,87,91	0
5	SO4	B	409	5/5	0.85	0.17	92,95,98,112	0
5	SO4	C	201	5/5	0.85	0.14	75,76,84,95	0
5	SO4	D	410	5/5	0.87	0.17	74,79,85,97	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	SO4	B	408	5/5	0.87	0.18	68,72,85,100	0
7	JRY	B	407	9/10	0.88	0.23	43,51,58,59	0
5	SO4	C	202	5/5	0.89	0.20	84,85,90,104	0
5	SO4	D	409	5/5	0.90	0.19	64,65,77,90	0
5	SO4	A	203	5/5	0.91	0.28	66,71,75,86	0
7	JRY	D	408	9/10	0.92	0.20	47,50,56,58	0
5	SO4	C	203	5/5	0.95	0.35	70,70,76,85	0

6.5 Other polymers [i](#)

There are no such residues in this entry.