



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 10, 2022 – 06:07 PM EST

PDB ID : 2BC4
Title : Crystal structure of HLA-DM
Authors : Nicholson, M.J.; Moradi, B.; Seth, N.P.; Xing, X.; Cuny, G.D.; Stein, R.L.;
Wucherpennig, K.W.
Deposited on : 2005-10-18
Resolution : 2.27 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.31.2
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) : 2.31.2

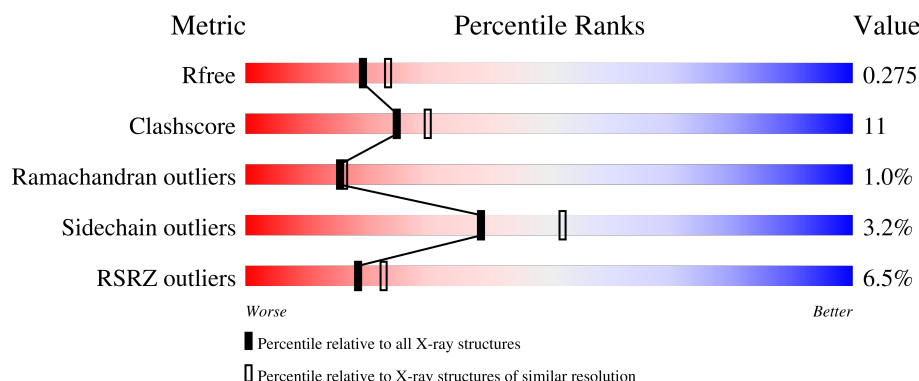
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.27 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	6980 (2.30-2.26)
Clashscore	141614	7711 (2.30-2.26)
Ramachandran outliers	138981	7597 (2.30-2.26)
Sidechain outliers	138945	7598 (2.30-2.26)
RSRZ outliers	127900	6849 (2.30-2.26)

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	211	<div> <div>3%</div> <div>74%</div> <div>13%</div> <div>11%</div> </div>
1	C	211	<div> <div>5%</div> <div>67%</div> <div>24%</div> <div>8%</div> </div>
2	B	211	<div> <div>5%</div> <div>65%</div> <div>24%</div> <div>9%</div> </div>
2	D	211	<div> <div>10%</div> <div>67%</div> <div>25%</div> <div>7%</div> </div>
3	E	4	<div> <div>50%</div> <div>25%</div> <div>25%</div> </div>

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Mol	Chain	Length	Quality of chain
3	F	4	 <div>75% 25%</div>

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
3	NAG	E	1	X	-	-	-
3	BMA	E	4	-	-	-	X
3	NAG	F	1	X	-	-	-
4	CL	C	504	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6667 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 HLA class II histocompatibility antigen, DM alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	188	Total	C	N	O	S	0	0	0
			1508	984	239	279	6			
1	C	195	Total	C	N	O	S	0	0	0
			1570	1024	247	293	6			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	204	ASP	-	expression tag	GB 18765715
A	205	TYR	-	expression tag	GB 18765715
A	206	LYS	-	expression tag	GB 18765715
A	207	ASP	-	expression tag	GB 18765715
A	208	ASP	-	expression tag	GB 18765715
A	209	ASP	-	expression tag	GB 18765715
A	210	ASP	-	expression tag	GB 18765715
A	211	LYS	-	expression tag	GB 18765715
C	204	ASP	-	expression tag	GB 18765715
C	205	TYR	-	expression tag	GB 18765715
C	206	LYS	-	expression tag	GB 18765715
C	207	ASP	-	expression tag	GB 18765715
C	208	ASP	-	expression tag	GB 18765715
C	209	ASP	-	expression tag	GB 18765715
C	210	ASP	-	expression tag	GB 18765715
C	211	LYS	-	expression tag	GB 18765715

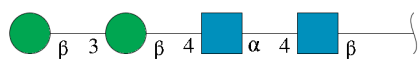
- Molecule 2 is a protein called HLA class II histocompatibility antigen, DM beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	191	Total	C	N	O	S	0	0	0
			1510	959	259	281	11			
2	D	197	Total	C	N	O	S	0	0	0
			1555	987	266	290	12			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	201	LYS	-	expression tag	GB 4504399
B	202	PRO	-	expression tag	GB 4504399
B	203	PRO	-	expression tag	GB 4504399
B	204	THR	-	expression tag	GB 4504399
B	205	PRO	-	expression tag	GB 4504399
B	206	PRO	-	expression tag	GB 4504399
B	207	PRO	-	expression tag	GB 4504399
B	208	GLU	-	expression tag	GB 4504399
B	209	PRO	-	expression tag	GB 4504399
B	210	GLU	-	expression tag	GB 4504399
B	211	THR	-	expression tag	GB 4504399
D	201	LYS	-	expression tag	GB 4504399
D	202	PRO	-	expression tag	GB 4504399
D	203	PRO	-	expression tag	GB 4504399
D	204	THR	-	expression tag	GB 4504399
D	205	PRO	-	expression tag	GB 4504399
D	206	PRO	-	expression tag	GB 4504399
D	207	PRO	-	expression tag	GB 4504399
D	208	GLU	-	expression tag	GB 4504399
D	209	PRO	-	expression tag	GB 4504399
D	210	GLU	-	expression tag	GB 4504399
D	211	THR	-	expression tag	GB 4504399

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



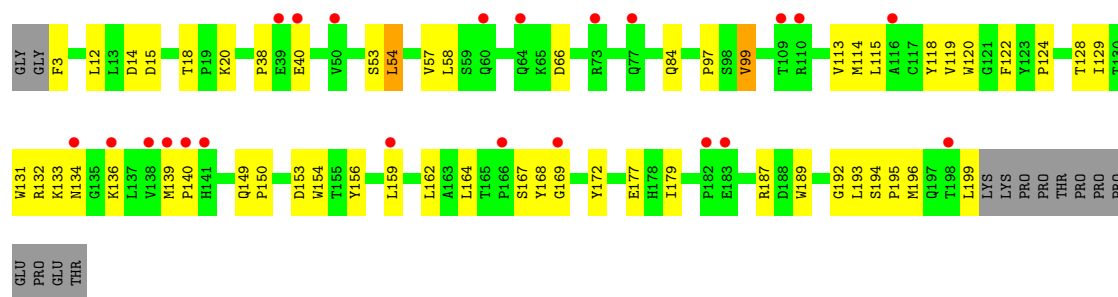
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	E	4	Total	C	N	O	0	0	0
			50	28	2	20			
3	F	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	135	Total 135	O 135	0	0
5	B	96	Total 96	O 96	0	0
5	C	121	Total 121	O 121	0	0
5	D	68	Total 68	O 68	0	0



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

Chain E: 50% 25% 25%



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

Chain F: 75% 25%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	98.07Å 108.42Å 110.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.27 29.40 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.7 (30.00-2.27) 99.1 (29.40-2.20)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.68 (at 2.20Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.229 , 0.268 0.234 , 0.275	Depositor DCC
R_{free} test set	1605 reflections (2.70%)	wwPDB-VP
Wilson B-factor (Å ²)	35.5	Xtriage
Anisotropy	0.190	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.001 for -h,l,k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6667	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 30.98 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.1935e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, CL, NAG, NDG

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	0.40	0/1561	0.80	3/2135 (0.1%)
1	C	0.40	0/1624	0.81	4/2220 (0.2%)
2	B	0.34	0/1556	0.61	0/2131
2	D	0.33	0/1602	0.63	1/2194 (0.0%)
All	All	0.37	0/6343	0.72	8/8680 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	98	ARG	NE-CZ-NH1	-14.63	112.98	120.30
1	C	98	ARG	NE-CZ-NH1	14.50	127.55	120.30
1	C	98	ARG	NE-CZ-NH2	-14.39	113.10	120.30
1	A	98	ARG	NE-CZ-NH2	14.16	127.38	120.30
1	C	98	ARG	CD-NE-CZ	7.57	134.20	123.60
1	A	98	ARG	CD-NE-CZ	7.48	134.07	123.60
2	D	194	SER	N-CA-C	5.60	126.13	111.00
1	C	68	ASP	N-CA-C	5.29	125.29	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1508	0	1417	27	0
1	C	1570	0	1475	43	0
2	B	1510	0	1452	35	0
2	D	1555	0	1499	37	0
3	E	50	0	42	1	0
3	F	50	0	42	2	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	135	0	0	5	0
5	B	96	0	0	1	0
5	C	121	0	0	7	0
5	D	68	0	0	1	0
All	All	6667	0	5927	132	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 11.

All (132) 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:B:124:PRO:HD2	2:B:179:ILE:HD11	1.39	1.04
1:A:15:ASN:HB3	2:B:12:LEU:HD22	1.69	0.75
1:C:15:ASN:HB3	2:D:12:LEU:HD22	1.70	0.74
1:C:20:HIS:HD2	1:C:35:GLU:OE2	1.75	0.69
1:A:20:HIS:HD2	1:A:35:GLU:OE2	1.76	0.67
2:B:139:MET:N	2:B:140:PRO:HD3	2.11	0.66
1:A:133:VAL:HG21	1:A:162:SER:OG	2.00	0.61
1:C:39:GLU:HA	5:C:618:HOH:O	2.01	0.60
5:C:510:HOH:O	3:F:1:NAG:H62	2.01	0.60
1:C:130:MET:HE1	1:C:184:ARG:HH21	1.66	0.60
2:B:167:SER:HB2	2:B:170:ASP:OD2	2.02	0.60
1:C:133:VAL:HG21	1:C:162:SER:OG	2.01	0.59
2:D:167:SER:O	2:D:169:GLY:N	2.33	0.59
1:C:81:TRP:NE1	1:C:85:GLN:HG3	2.18	0.58
5:A:600:HOH:O	2:B:82:HIS:HE1	1.88	0.56
1:C:90:LEU:HB3	1:C:94:ILE:CD1	2.35	0.56
1:C:73:LEU:O	1:C:77:GLU:HG3	2.05	0.56
1:C:130:MET:CE	1:C:184:ARG:HH21	2.19	0.55
1:C:16:HIS:HE1	2:D:15:ASP:OD1	1.89	0.55
1:A:16:HIS:HE1	2:B:15:ASP:OD1	1.90	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:94:ILE:H	1:C:94:ILE:HD13	1.72	0.55
2:D:139:MET:N	2:D:140:PRO:HD3	2.22	0.54
1:A:130:MET:CE	1:A:184:ARG:HH21	2.20	0.54
2:B:114:MET:HE3	2:B:163:ALA:HB2	1.89	0.54
1:A:82:MET:O	1:A:86:ILE:O	2.24	0.54
1:C:55:ARG:HG3	1:C:56:LEU:HD13	1.91	0.53
1:A:182:ILE:O	1:A:182:ILE:HG22	2.09	0.53
1:A:19:LEU:O	1:A:35:GLU:HA	2.09	0.53
1:C:201:LEU:HD13	1:C:201:LEU:O	2.09	0.53
2:B:54:LEU:HD22	2:B:58:LEU:HG	1.91	0.53
2:D:54:LEU:HD22	2:D:58:LEU:HG	1.92	0.52
1:C:19:LEU:O	1:C:35:GLU:HA	2.09	0.52
2:D:97:PRO:HB3	2:D:122:PHE:HB3	1.91	0.52
1:A:55:ARG:HG3	1:A:56:LEU:HD13	1.92	0.51
2:B:97:PRO:HB3	2:B:122:PHE:HB3	1.93	0.51
1:A:73:LEU:O	1:A:77:GLU:HG3	2.09	0.51
2:B:51:LEU:HA	5:B:559:HOH:O	2.11	0.51
1:A:20:HIS:HE1	5:A:527:HOH:O	1.93	0.51
2:D:179:ILE:HG23	5:D:546:HOH:O	2.10	0.51
1:A:181:GLU:OE1	1:A:181:GLU:HA	2.11	0.50
2:B:101:VAL:HG23	2:B:187:ARG:HB3	1.94	0.50
1:C:181:GLU:HA	1:C:181:GLU:OE1	2.11	0.50
1:C:182:ILE:HG22	1:C:182:ILE:O	2.11	0.50
1:C:20:HIS:HE1	5:C:528:HOH:O	1.94	0.50
2:D:113:VAL:HB	2:D:164:LEU:HD12	1.94	0.50
2:B:128:THR:OG1	2:B:177:GLU:HB2	2.12	0.49
2:D:167:SER:C	2:D:169:GLY:H	2.15	0.49
2:B:164:LEU:HD22	2:B:172:TYR:CZ	2.47	0.49
2:D:119:VAL:HG21	2:D:129:ILE:HD11	1.94	0.49
1:C:149:PHE:O	1:C:160:ALA:HA	2.13	0.49
2:D:153:ASP:O	2:D:154:TRP:HB2	2.12	0.49
2:D:113:VAL:HG12	2:D:114:MET:N	2.28	0.48
2:D:128:THR:OG1	2:D:177:GLU:HB2	2.13	0.48
2:B:150:PRO:HB3	2:B:156:TYR:CE2	2.49	0.47
2:B:14:ASP:OD2	2:B:18:THR:HB	2.14	0.47
2:B:186:LEU:O	2:B:187:ARG:HD2	2.13	0.47
2:D:14:ASP:OD2	2:D:18:THR:HB	2.14	0.47
2:B:111:GLU:HB2	2:B:163:ALA:HB1	1.97	0.47
1:C:28:SER:O	1:C:130:MET:HG2	2.14	0.47
1:C:146:GLY:HA3	5:C:592:HOH:O	2.15	0.47
2:D:192:GLY:O	2:D:195:PRO:HD2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:104:THR:CG2	2:B:114:MET:HG2	2.45	0.46
2:B:153:ASP:O	2:B:154:TRP:HB2	2.15	0.46
1:C:16:HIS:HD2	1:C:38:ASP:OD2	1.98	0.46
2:D:150:PRO:HB3	2:D:156:TYR:CE2	2.50	0.46
2:B:137:LEU:HD12	2:B:138:VAL:H	1.81	0.46
1:A:16:HIS:HD2	1:A:38:ASP:OD2	1.98	0.46
2:D:131:TRP:CD2	2:D:162:LEU:HD22	2.51	0.46
2:D:164:LEU:HD22	2:D:172:TYR:CZ	2.51	0.46
1:A:130:MET:HB3	1:A:181:GLU:HB2	1.98	0.46
2:D:14:ASP:OD2	2:D:20:LYS:HE2	2.16	0.46
1:A:149:PHE:O	1:A:160:ALA:HA	2.15	0.45
2:D:99:VAL:HA	2:D:118:TYR:O	2.17	0.45
2:B:139:MET:N	2:B:140:PRO:CD	2.80	0.45
5:A:509:HOH:O	3:E:1:NAG:H62	2.16	0.45
1:A:28:SER:O	1:A:130:MET:HG2	2.15	0.45
1:A:105:VAL:HA	1:A:120:VAL:O	2.17	0.45
2:B:56:ASN:O	2:B:60:GLN:HG3	2.17	0.45
2:D:54:LEU:CD2	2:D:58:LEU:HG	2.47	0.44
1:C:66:GLN:O	1:C:66:GLN:HG3	2.17	0.44
1:C:200:ASP:HB3	1:C:203:GLU:HB3	2.00	0.44
1:A:61:ASP:HB2	5:A:551:HOH:O	2.16	0.44
1:A:130:MET:HE3	1:A:184:ARG:HH21	1.81	0.44
2:D:133:LYS:O	2:D:134:ASN:HB2	2.16	0.44
2:D:124:PRO:HD2	2:D:179:ILE:HD11	2.00	0.44
2:D:124:PRO:CD	2:D:179:ILE:HD11	2.48	0.44
1:A:197:LEU:HD11	2:B:118:TYR:CE1	2.53	0.43
1:C:90:LEU:HB3	1:C:94:ILE:HD12	2.00	0.43
2:B:45:PRO:HG2	2:B:56:ASN:OD1	2.17	0.43
1:C:90:LEU:O	1:C:93:LYS:HB2	2.18	0.43
1:C:104:GLU:OE2	2:D:149:GLN:NE2	2.52	0.43
1:C:119:LEU:HD13	1:C:166:PHE:CE2	2.54	0.43
1:A:104:GLU:OE2	2:B:149:GLN:NE2	2.52	0.43
1:C:130:MET:HB3	1:C:181:GLU:HB2	2.00	0.43
2:D:53:SER:O	2:D:57:VAL:HG23	2.18	0.43
1:C:137:HIS:HE1	1:C:172:ASP:OD1	2.02	0.43
1:A:119:LEU:HD13	1:A:166:PHE:CE2	2.54	0.43
5:C:519:HOH:O	2:D:120:TRP:HB2	2.18	0.43
2:B:14:ASP:OD2	2:B:20:LYS:HE2	2.18	0.43
2:B:99:VAL:HA	2:B:118:TYR:O	2.18	0.43
1:C:138:HIS:O	1:C:139:SER:HB2	2.19	0.43
1:C:57:PRO:HD2	5:C:540:HOH:O	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:202:LEU:HD11	2:D:196:MET:CE	2.49	0.42
1:C:86:ILE:HG22	1:C:90:LEU:HD23	2.01	0.42
2:B:44:ALA:HA	2:B:45:PRO:HD3	1.92	0.42
1:C:105:VAL:HA	1:C:120:VAL:O	2.19	0.42
1:A:130:MET:HE1	1:A:184:ARG:HH21	1.83	0.42
5:C:600:HOH:O	2:D:199:LEU:HB3	2.19	0.42
1:C:16:HIS:CE1	2:D:15:ASP:OD1	2.71	0.42
1:C:119:LEU:HD13	1:C:166:PHE:CZ	2.55	0.42
1:C:87:GLY:H	1:C:90:LEU:HD23	1.85	0.41
2:D:12:LEU:HD13	3:F:1:NAG:H61	2.01	0.41
1:C:76:LYS:O	1:C:80:GLU:HG3	2.19	0.41
1:C:115:LYS:HA	1:C:116:PRO:HD3	1.98	0.41
1:A:16:HIS:CE1	2:B:15:ASP:OD1	2.72	0.41
1:C:197:LEU:HD12	1:C:197:LEU:N	2.35	0.41
5:A:518:HOH:O	2:B:120:TRP:HB2	2.19	0.41
1:C:94:ILE:HD11	2:D:3:PHE:CZ	2.55	0.41
2:D:115:LEU:HD21	2:D:189:TRP:CE2	2.55	0.41
2:D:132:ARG:HA	2:D:136:LYS:O	2.21	0.41
1:A:119:LEU:HD13	1:A:166:PHE:CZ	2.56	0.41
2:D:119:VAL:HG21	2:D:129:ILE:CD1	2.50	0.41
2:D:164:LEU:HD12	2:D:164:LEU:C	2.41	0.41
1:A:15:ASN:OD1	2:B:20:LYS:HG3	2.21	0.41
2:B:69:MET:O	2:B:73:ARG:HG3	2.21	0.41
2:B:40:GLU:CG	2:B:42:LYS:HE2	2.51	0.41
2:D:38:PRO:C	2:D:40:GLU:H	2.24	0.41
1:C:90:LEU:HB3	1:C:94:ILE:HD11	2.03	0.40
2:B:25:CYS:HB3	2:B:32:LEU:HD11	2.03	0.40
2:B:164:LEU:HD12	2:B:164:LEU:C	2.41	0.40
1:A:145:PHE:CD1	1:A:165:ASN:ND2	2.89	0.40
1:C:115:LYS:O	1:C:168:PRO:HD2	2.21	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	186/211 (88%)	179 (96%)	5 (3%)	2 (1%)	14	14
1	C	193/211 (92%)	182 (94%)	7 (4%)	4 (2%)	7	5
2	B	189/211 (90%)	182 (96%)	6 (3%)	1 (0%)	29	34
2	D	195/211 (92%)	182 (93%)	12 (6%)	1 (0%)	29	34
All	All	763/844 (90%)	725 (95%)	30 (4%)	8 (1%)	15	16

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	168	TYR
1	A	182	ILE
1	C	182	ILE
1	C	13	LEU
1	C	181	GLU
1	A	181	GLU
2	B	168	TYR
1	C	69	ALA

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	167/189 (88%)	165 (99%)	2 (1%)	71	82
1	C	174/189 (92%)	170 (98%)	4 (2%)	50	65
2	B	169/187 (90%)	160 (95%)	9 (5%)	22	29
2	D	175/187 (94%)	168 (96%)	7 (4%)	31	42
All	All	685/752 (91%)	663 (97%)	22 (3%)	39	52

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	98	ARG
1	A	181	GLU
2	B	54	LEU
2	B	66	ASP
2	B	76	LEU
2	B	84	GLN
2	B	99	VAL
2	B	114	MET
2	B	145	HIS
2	B	159	LEU
2	B	179	ILE
1	C	84	GLN
1	C	94	ILE
1	C	98	ARG
1	C	181	GLU
2	D	54	LEU
2	D	66	ASP
2	D	84	GLN
2	D	99	VAL
2	D	159	LEU
2	D	187	ARG
2	D	193	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	HIS
1	A	20	HIS
1	A	50	ASN
1	A	125	ASN
1	A	137	HIS
2	B	78	ASN
2	B	82	HIS
2	B	108	ASN
2	B	157	GLN
1	C	16	HIS
1	C	20	HIS
1	C	125	ASN
1	C	137	HIS
2	D	70	GLN
2	D	78	ASN
2	D	108	ASN
2	D	157	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 ⓘ

8 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	3,1	14,14,15	0.77	0	17,19,21	0.73	1 (5%)
3	NDG	E	2	3	14,14,15	0.73	0	17,19,21	0.74	1 (5%)
3	BMA	E	3	3	11,11,12	0.71	0	15,15,17	0.39	0
3	BMA	E	4	3	11,11,12	0.53	0	15,15,17	0.60	0
3	NAG	F	1	3,1	14,14,15	0.70	0	17,19,21	0.75	1 (5%)
3	NDG	F	2	3	14,14,15	0.71	0	17,19,21	0.72	0
3	BMA	F	3	3	11,11,12	0.73	0	15,15,17	0.40	0
3	BMA	F	4	3	11,11,12	0.51	0	15,15,17	0.61	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	3,1	1/1/5/7	1/6/23/26	0/1/1/1
3	NDG	E	2	3	-	4/6/23/26	0/1/1/1
3	BMA	E	3	3	-	0/2/19/22	0/1/1/1
3	BMA	E	4	3	-	0/2/19/22	0/1/1/1
3	NAG	F	1	3,1	1/1/5/7	1/6/23/26	0/1/1/1
3	NDG	F	2	3	-	4/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	BMA	F	3	3	-	0/2/19/22	0/1/1/1
3	BMA	F	4	3	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	1	NAG	C2-N2-C7	-2.14	119.85	122.90
3	F	1	NAG	C2-N2-C7	-2.10	119.91	122.90
3	E	2	NDG	C2-N2-C7	-2.04	120.00	122.90

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	E	1	NAG	C1
3	F	1	NAG	C1

All (10) torsion outliers are listed below:

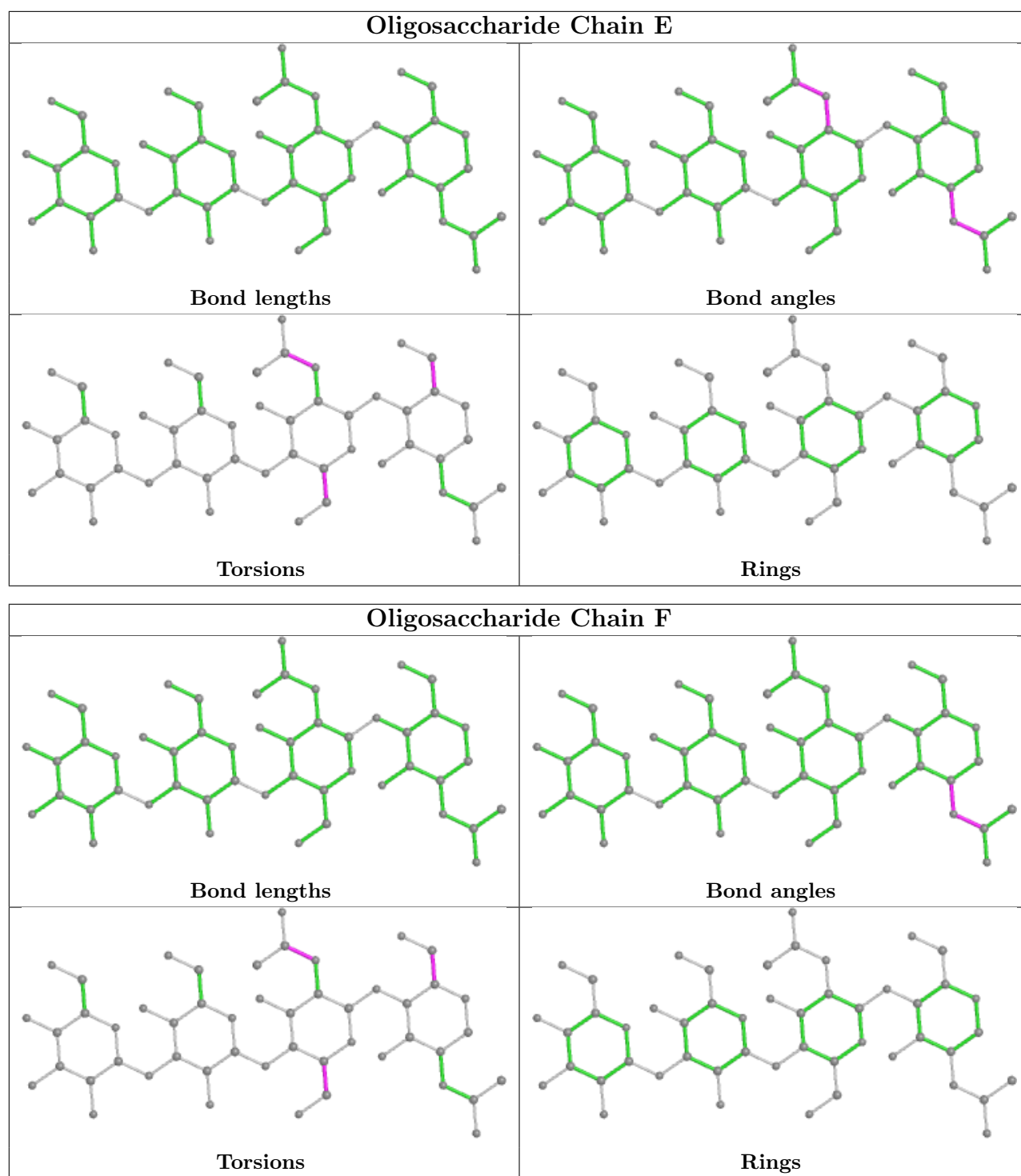
Mol	Chain	Res	Type	Atoms
3	E	2	NDG	C8-C7-N2-C2
3	E	2	NDG	O7-C7-N2-C2
3	F	2	NDG	C8-C7-N2-C2
3	F	2	NDG	O7-C7-N2-C2
3	E	2	NDG	O5-C5-C6-O6
3	F	2	NDG	O5-C5-C6-O6
3	E	2	NDG	C4-C5-C6-O6
3	F	2	NDG	C4-C5-C6-O6
3	F	1	NAG	O5-C5-C6-O6
3	E	1	NAG	O5-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	1	NAG	1	0
3	F	1	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](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

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

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	188/211 (89%)	0.14	7 (3%) 41 47	20, 31, 52, 67	0
1	C	195/211 (92%)	0.11	10 (5%) 28 33	20, 34, 60, 70	0
2	B	191/211 (90%)	0.21	11 (5%) 23 28	22, 38, 58, 72	0
2	D	197/211 (93%)	0.72	22 (11%) 5 6	22, 46, 68, 74	0
All	All	771/844 (91%)	0.30	50 (6%) 18 23	20, 37, 63, 74	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	141	HIS	6.0
2	B	141	HIS	5.7
2	D	136	LYS	5.2
2	D	110	ARG	4.1
2	B	139	MET	4.0
2	D	60	GLN	3.9
2	D	198	THR	3.8
2	D	139	MET	3.7
2	D	140	PRO	3.7
2	D	40	GLU	3.5
1	C	12	ASP	3.3
2	B	39	GLU	3.1
1	A	182	ILE	3.0
1	A	189	ALA	2.9
1	A	105	VAL	2.9
2	B	140	PRO	2.8
2	B	92	ASN	2.8
2	D	183	GLU	2.8
2	B	110	ARG	2.7
2	D	64	GLN	2.7
1	A	13	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
2	D	134	ASN	2.6
2	D	39	GLU	2.6
2	D	169	GLY	2.6
2	D	73	ARG	2.5
1	C	188	ILE	2.5
2	D	77	GLN	2.5
1	C	189	ALA	2.5
1	C	13	LEU	2.4
2	B	145	HIS	2.4
1	A	188	ILE	2.4
1	C	93	LYS	2.4
2	D	116	ALA	2.3
1	C	66	GLN	2.3
1	C	105	VAL	2.3
1	C	205	TYR	2.3
1	C	145	PHE	2.3
2	D	182	PRO	2.3
2	B	158	THR	2.2
2	D	50	VAL	2.2
2	D	138	VAL	2.2
2	D	159	LEU	2.2
1	C	116	PRO	2.2
1	A	199	SER	2.2
2	D	109	THR	2.1
1	A	190	TYR	2.1
2	B	22	PHE	2.1
2	B	181	ALA	2.1
2	D	166	PRO	2.1
2	B	136	LYS	2.0

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

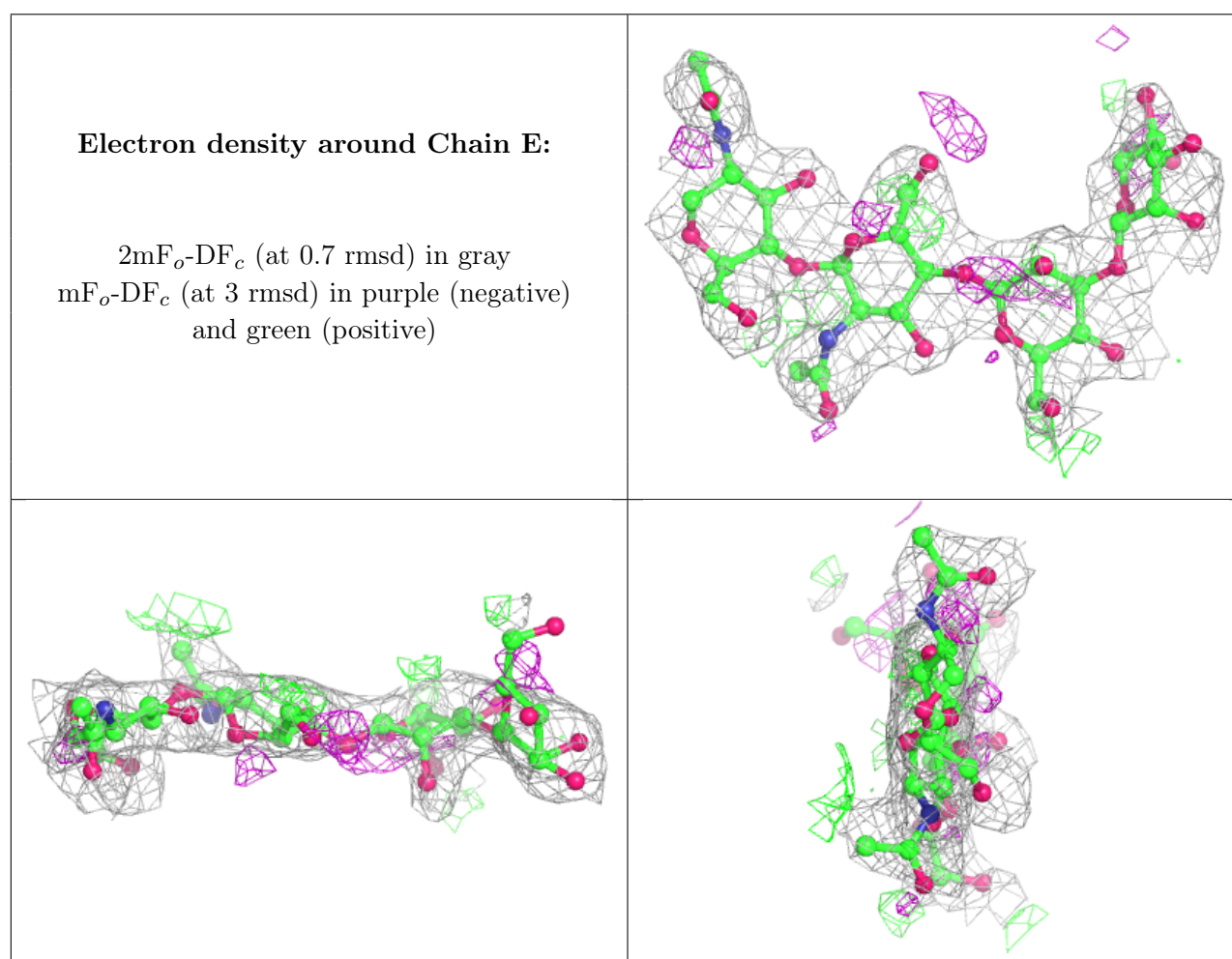
There are no non-standard protein/DNA/RNA residues in this entry.

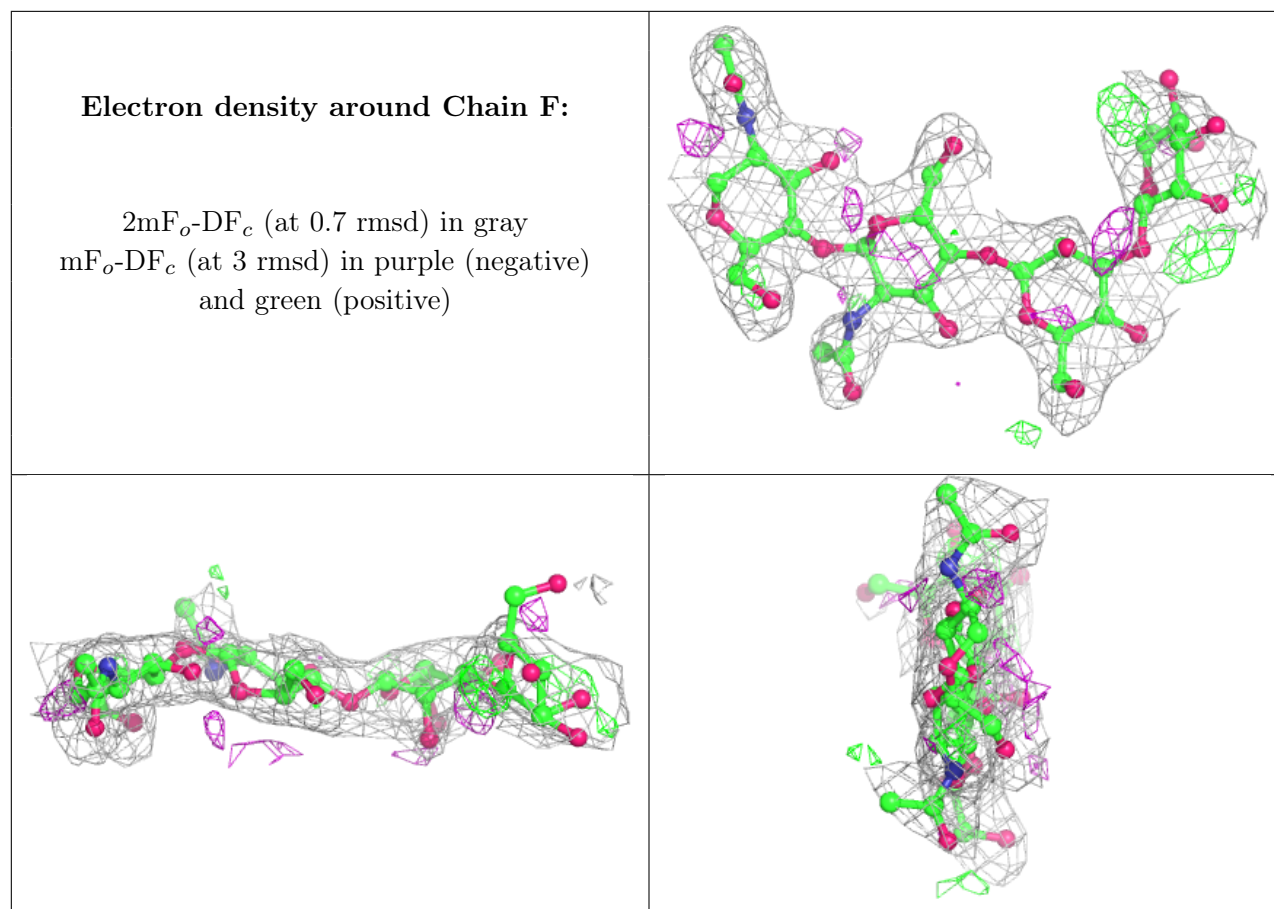
6.3 Carbohydrates ⓘ

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	BMA	E	3	11/12	0.56	0.30	70,71,73,76	0
3	BMA	F	4	11/12	0.63	0.33	77,79,81,81	0
3	BMA	E	4	11/12	0.73	0.41	78,79,81,81	0
3	NDG	E	2	14/15	0.77	0.21	56,59,62,67	0
3	BMA	F	3	11/12	0.82	0.14	67,69,72,75	0
3	NDG	F	2	14/15	0.84	0.19	55,59,61,63	0
3	NAG	E	1	14/15	0.86	0.15	40,43,46,52	0
3	NAG	F	1	14/15	0.92	0.14	39,42,45,50	0

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





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	CL	C	504	1/1	0.79	0.43	71,71,71,71	0
4	CL	A	503	1/1	0.88	0.50	68,68,68,68	0
4	CL	B	501	1/1	0.91	0.22	57,57,57,57	0
4	CL	D	502	1/1	0.96	0.28	51,51,51,51	0

6.5 Other polymers [i](#)

There are no such residues in this entry.