



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 8, 2020 – 11:42 PM EST

PDB ID : 6OHE
Title : Alpha-L-fucosidase AlfC D200A in complex with Fuca(1,6)GlcNAc
Authors : Klontz, E.H.; Sundberg, E.J.
Deposited on : 2019-04-05
Resolution : 3.14 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 : 2.15.1
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) : 2.15.1

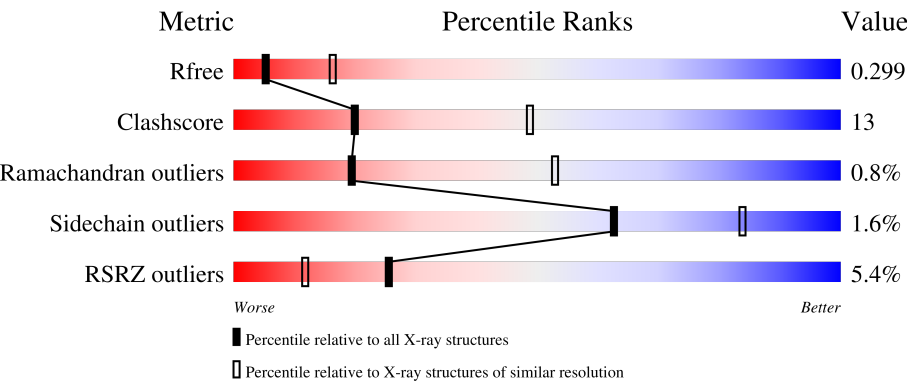
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.14 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	1626 (3.18-3.10)
Clashscore	141614	1735 (3.18-3.10)
Ramachandran outliers	138981	1677 (3.18-3.10)
Sidechain outliers	138945	1677 (3.18-3.10)
RSRZ outliers	127900	1588 (3.18-3.10)

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	345	<div><div>4%</div><div>73%</div><div>21%</div><div>6%</div></div>
1	B	345	<div><div>4%</div><div>68%</div><div>24%</div><div>7%</div></div>
1	C	345	<div><div>5%</div><div>71%</div><div>21%</div><div>7%</div></div>
1	D	345	<div><div>7%</div><div>71%</div><div>20%</div><div>7%</div></div>
2	E	2	<div><div>100%</div></div>

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

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10367 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 AlfC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	325	Total	C	N	O	S	0	0	0
			2586	1651	425	498	12			
1	B	322	Total	C	N	O	S	0	0	0
			2566	1639	421	494	12			
1	C	322	Total	C	N	O	S	0	0	0
			2565	1640	421	492	12			
1	D	320	Total	C	N	O	S	0	0	0
			2550	1629	419	490	12			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	200	ALA	ASP	engineered mutation	UNP K0NB39
A	345	LEU	-	expression tag	UNP K0NB39
B	200	ALA	ASP	engineered mutation	UNP K0NB39
B	345	LEU	-	expression tag	UNP K0NB39
C	200	ALA	ASP	engineered mutation	UNP K0NB39
C	345	LEU	-	expression tag	UNP K0NB39
D	200	ALA	ASP	engineered mutation	UNP K0NB39
D	345	LEU	-	expression tag	UNP K0NB39

- Molecule 2 is an oligosaccharide called alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	2	Total	C	N	O	0	0	0
			25	14	1	10			

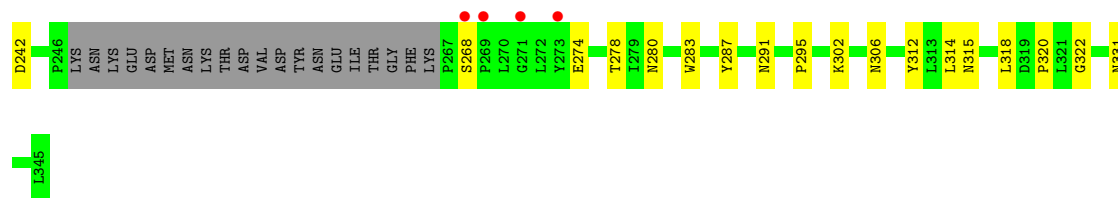
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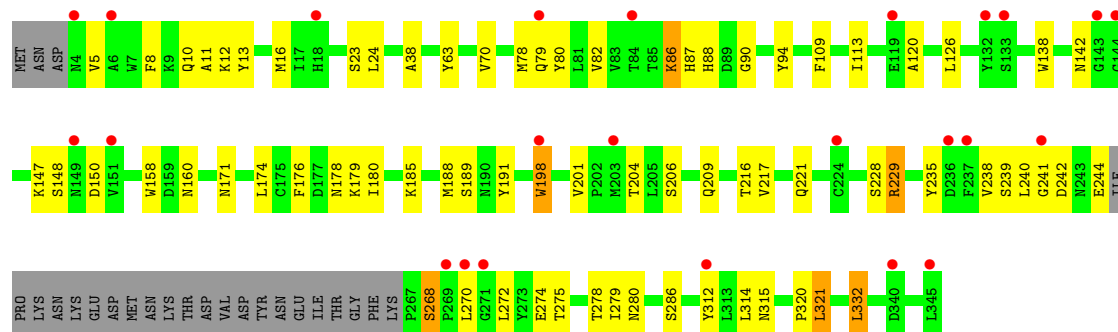
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	F	2	Total	C	N	O	0	0	0
			25	14	1	10			
2	G	2	Total	C	N	O	0	0	0
			25	14	1	10			
2	H	2	Total	C	N	O	0	0	0
			25	14	1	10			

- Molecule 1: AlfC





- Molecule 1: AlfC



- Molecule 2: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 2: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 2: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 2: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	89.02Å 139.25Å 263.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.76 – 3.14 29.76 – 3.14	Depositor EDS
% Data completeness (in resolution range)	99.7 (29.76-3.14) 99.7 (29.76-3.14)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.39 (at 3.11Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.279 , 0.299 0.279 , 0.299	Depositor DCC
R_{free} test set	1445 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	86.3	Xtriage
Anisotropy	0.479	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 46.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	10367	wwPDB-VP
Average B, all atoms (Å ²)	126.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.65% 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, FUC

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	0.36	0/2659	0.65	2/3616 (0.1%)
1	B	0.39	0/2638	0.68	0/3585
1	C	0.35	0/2638	0.65	1/3586 (0.0%)
1	D	0.37	0/2622	0.68	3/3563 (0.1%)
All	All	0.37	0/10557	0.66	6/14350 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	321	LEU	CA-CB-CG	8.72	135.35	115.30
1	D	229	ARG	CA-CB-CG	-6.50	99.10	113.40
1	D	332	LEU	CA-CB-CG	-5.75	102.07	115.30
1	C	240	LEU	CA-CB-CG	-5.73	102.12	115.30
1	A	203	MET	CA-CB-CG	5.57	122.76	113.30
1	A	203	MET	CB-CG-SD	-5.01	97.37	112.40

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	2586	0	2405	54	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2566	0	2391	81	0
1	C	2565	0	2394	65	0
1	D	2550	0	2376	57	0
2	E	25	0	24	1	0
2	F	25	0	24	2	0
2	G	25	0	24	6	0
2	H	25	0	24	1	0
All	All	10367	0	9662	252	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 13.

All (252) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:19:TRP:HE1	1:C:318:LEU:HD12	1.36	0.90
1:C:228:SER:HB3	1:C:238:VAL:HA	1.59	0.84
1:D:86:LYS:HD3	1:D:90:GLY:HA2	1.60	0.84
1:C:19:TRP:HZ2	1:C:322:GLY:HA2	1.42	0.84
1:C:198:TRP:HH2	1:C:239:SER:HG	1.25	0.83
1:C:86:LYS:NZ	1:C:137:ASP:OD2	2.11	0.83
1:D:5:VAL:HG11	1:D:270:LEU:HD22	1.61	0.82
1:B:138:TRP:O	1:B:160:ASN:ND2	2.12	0.81
1:D:176:PHE:HA	1:D:180:ILE:HD13	1.63	0.80
1:D:63:TYR:HD2	1:D:321:LEU:HD21	1.47	0.80
1:B:24:LEU:HD11	1:B:59:PHE:HD2	1.47	0.79
1:B:228:SER:HB3	1:B:238:VAL:HA	1.65	0.78
1:B:226:ILE:H	1:B:236:ASP:HB2	1.50	0.77
1:C:176:PHE:HA	1:C:180:ILE:HD13	1.68	0.75
1:B:145:GLY:HA2	1:B:204:THR:HG23	1.68	0.74
1:A:176:PHE:HA	1:A:180:ILE:HD13	1.70	0.74
1:B:146:TYR:HE2	1:B:172:PHE:HB2	1.53	0.74
1:B:4:ASN:HB3	1:B:225:LEU:HD11	1.69	0.73
1:A:62:LEU:HD11	1:C:58:ALA:HB2	1.72	0.72
1:D:80:TYR:OH	1:D:274:GLU:OE2	2.06	0.72
1:D:174:LEU:O	1:D:178:ASN:ND2	2.17	0.71
1:D:63:TYR:CD2	1:D:321:LEU:HD21	2.25	0.71
1:B:336:LYS:NZ	1:B:340:ASP:OD1	2.23	0.70
1:B:149:ASN:ND2	1:B:159:ASP:OD2	2.25	0.69
1:B:240:LEU:HD11	1:B:273:TYR:HD1	1.56	0.69
1:B:130:LEU:HD12	1:B:194:ILE:HG21	1.74	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:19:TRP:NE1	1:C:318:LEU:HB2	2.07	0.69
1:C:302:LYS:NZ	1:C:306:ASN:OD1	2.24	0.68
1:B:199:PHE:H	1:B:227:ASN:HD21	1.40	0.68
1:D:238:VAL:HG23	1:D:268:SER:HB2	1.76	0.67
1:B:135:ASP:OD1	1:B:204:THR:OG1	2.13	0.67
1:B:143:GLY:HA2	1:B:172:PHE:HD1	1.57	0.67
1:A:86:LYS:NZ	1:A:137:ASP:OD2	2.28	0.66
1:C:228:SER:CB	1:C:238:VAL:HA	2.26	0.65
1:A:62:LEU:CD1	1:C:58:ALA:HB2	2.26	0.65
1:A:9:LYS:NZ	1:A:270:LEU:O	2.29	0.65
1:B:199:PHE:H	1:B:227:ASN:ND2	1.95	0.65
1:A:287:TYR:CE1	1:C:287:TYR:CE1	2.86	0.64
1:C:278:THR:HG23	1:C:280:ASN:O	1.99	0.62
1:D:23:SER:O	1:D:320:PRO:HG3	1.99	0.62
1:B:145:GLY:CA	1:B:204:THR:HG23	2.30	0.61
1:D:206:SER:N	1:D:209:GLN:OE1	2.29	0.61
1:A:246:PRO:HB2	1:A:308:PHE:HE2	1.66	0.61
1:B:49:ASN:HD21	1:B:160:ASN:HD21	1.47	0.61
1:B:180:ILE:H	1:B:180:ILE:HD12	1.65	0.60
1:A:58:ALA:HB2	1:C:62:LEU:CD1	2.31	0.60
1:C:80:TYR:OH	1:C:274:GLU:OE2	2.07	0.60
1:D:90:GLY:O	1:D:179:LYS:NZ	2.34	0.60
1:D:16:MET:HB3	1:D:315:ASN:HA	1.85	0.59
1:C:19:TRP:HE1	1:C:318:LEU:CD1	2.14	0.59
1:B:24:LEU:HD11	1:B:59:PHE:CD2	2.33	0.58
1:A:80:TYR:OH	1:A:274:GLU:OE2	2.14	0.58
1:D:109:PHE:HE2	1:D:113:ILE:HD11	1.68	0.58
1:A:198:TRP:HZ2	1:A:274:GLU:OE1	1.86	0.58
1:A:278:THR:HG23	1:A:280:ASN:O	2.03	0.58
1:D:314:LEU:HD23	1:D:332:LEU:HD21	1.86	0.57
1:C:200:ALA:HA	1:C:229:ARG:NH1	2.20	0.57
1:C:37:TYR:HE2	2:G:1:NAG:HO1	1.53	0.57
1:C:19:TRP:NE1	1:C:318:LEU:HD12	2.14	0.57
1:B:136:LEU:HA	1:B:144:GLY:HA2	1.87	0.57
1:B:132:TYR:HE1	1:B:180:ILE:HG23	1.71	0.56
1:B:175:CYS:SG	1:B:180:ILE:HD11	2.45	0.56
1:B:278:THR:HG23	1:B:280:ASN:O	2.05	0.56
1:A:198:TRP:CZ2	1:A:274:GLU:OE1	2.59	0.56
1:C:198:TRP:HZ2	1:C:274:GLU:OE1	1.89	0.55
1:C:34:SER:HB3	1:C:41:ILE:HD13	1.89	0.55
1:B:132:TYR:CE1	1:B:180:ILE:HG23	2.41	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:144:GLY:O	1:B:159:ASP:HB3	2.07	0.54
1:C:198:TRP:CZ2	1:C:274:GLU:OE1	2.61	0.54
1:C:87:HIS:HE2	2:G:2:FUC:HO4	1.54	0.54
2:G:1:NAG:O3	2:G:1:NAG:O7	2.23	0.54
1:C:154:ALA:N	2:G:1:NAG:O7	2.40	0.54
1:A:287:TYR:CE1	1:C:287:TYR:HE1	2.26	0.54
1:A:342:ALA:HA	1:A:345:LEU:HD12	1.89	0.54
1:B:142:ASN:HD21	1:B:171:ASN:N	2.06	0.53
1:B:47:ILE:CG2	1:B:52:TYR:HB2	2.38	0.53
1:C:87:HIS:NE2	2:G:2:FUC:O4	2.39	0.53
1:B:94:TYR:HD2	1:B:191:TYR:HH	1.56	0.53
1:A:235:TYR:OH	1:A:267:PRO:O	2.21	0.53
1:B:26:ALA:HB3	1:B:287:TYR:CE2	2.43	0.53
1:D:180:ILE:H	1:D:180:ILE:HD12	1.74	0.53
1:B:94:TYR:HB2	1:B:191:TYR:OH	2.09	0.53
1:B:180:ILE:HD12	1:B:180:ILE:N	2.24	0.53
1:B:89:ASP:HA	1:B:138:TRP:CE3	2.44	0.52
1:C:19:TRP:CZ2	1:C:322:GLY:HA2	2.34	0.52
1:D:185:LYS:O	1:D:189:SER:OG	2.16	0.52
1:D:198:TRP:HZ2	1:D:274:GLU:OE1	1.91	0.52
1:B:29:TYR:OH	1:B:47:ILE:HD11	2.10	0.52
1:D:24:LEU:HG	1:D:320:PRO:HB3	1.91	0.52
1:D:87:HIS:HD1	1:D:88:HIS:H	1.57	0.52
1:B:228:SER:CB	1:B:238:VAL:HA	2.37	0.51
1:C:198:TRP:HH2	1:C:239:SER:OG	1.89	0.51
1:C:291:ASN:O	1:C:291:ASN:ND2	2.43	0.51
1:B:29:TYR:CZ	1:B:47:ILE:HD11	2.45	0.51
1:C:180:ILE:HD12	1:C:180:ILE:H	1.75	0.51
1:D:321:LEU:O	1:D:321:LEU:HD23	2.11	0.51
1:A:4:ASN:ND2	1:A:223:ASN:O	2.44	0.51
1:C:7:TRP:CE2	1:C:127:LYS:HE3	2.45	0.51
1:D:78:MET:O	1:D:126:LEU:HD21	2.11	0.51
1:A:246:PRO:HB2	1:A:308:PHE:CE2	2.44	0.50
1:A:242:ASP:HA	1:A:245:ILE:HD12	1.92	0.50
1:D:109:PHE:CE2	1:D:113:ILE:HD11	2.47	0.50
1:D:278:THR:HG23	1:D:280:ASN:O	2.12	0.50
1:A:24:LEU:HD11	1:A:59:PHE:HD2	1.77	0.50
1:C:136:LEU:HD11	1:C:158:TRP:HE3	1.77	0.50
1:D:88:HIS:NE2	2:H:2:FUC:H2	2.27	0.50
1:B:142:ASN:HD21	1:B:171:ASN:H	1.58	0.50
1:D:82:VAL:HG11	1:D:198:TRP:CD1	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:LYS:NZ	1:A:124:ALA:O	2.36	0.49
1:C:10:GLN:O	1:C:12:LYS:HG3	2.12	0.49
1:B:49:ASN:HD21	1:B:160:ASN:ND2	2.10	0.49
1:C:24:LEU:HA	1:C:320:PRO:HG3	1.94	0.49
1:B:240:LEU:HD11	1:B:273:TYR:CD1	2.42	0.49
1:B:146:TYR:CE2	1:B:172:PHE:HB2	2.42	0.48
1:A:58:ALA:HB2	1:C:62:LEU:HD13	1.95	0.48
1:B:148:SER:OG	1:B:203:MET:HB3	2.12	0.48
1:A:111:ARG:NH2	1:A:116:GLU:OE1	2.46	0.48
1:C:214:TYR:HB2	1:C:230:LEU:HD21	1.95	0.48
1:C:89:ASP:HA	1:C:138:TRP:CE3	2.48	0.48
1:D:240:LEU:HB2	1:D:275:THR:OG1	2.13	0.48
1:D:38:ALA:HB2	1:D:286:SER:HB2	1.95	0.48
1:A:148:SER:O	1:A:158:TRP:CD1	2.67	0.48
1:B:227:ASN:OD1	1:B:229:ARG:HB2	2.13	0.48
1:B:228:SER:HA	1:B:235:TYR:CG	2.48	0.48
1:C:7:TRP:CD2	1:C:127:LYS:HE3	2.49	0.47
1:B:84:THR:HG22	1:B:131:TYR:CD1	2.48	0.47
1:A:119:GLU:HG2	1:A:123:LYS:HE3	1.96	0.47
1:A:316:VAL:HG11	1:A:324:VAL:HG22	1.95	0.47
1:B:138:TRP:HA	1:B:160:ASN:HB2	1.95	0.47
1:B:39:GLU:CD	1:B:283:TRP:HB2	2.35	0.47
1:D:198:TRP:HH2	1:D:239:SER:HG	1.61	0.47
1:A:175:CYS:SG	1:A:180:ILE:HD11	2.53	0.47
1:C:185:LYS:HG2	1:C:216:THR:HG21	1.97	0.47
1:C:227:ASN:O	1:C:229:ARG:N	2.45	0.47
1:C:228:SER:N	1:C:237:PHE:O	2.45	0.47
1:A:58:ALA:HB2	1:C:62:LEU:HD11	1.96	0.47
1:C:175:CYS:SG	1:C:180:ILE:HD11	2.55	0.47
2:F:1:NAG:O3	2:F:1:NAG:O7	2.28	0.47
1:D:241:GLY:N	1:D:244:GLU:OE2	2.48	0.47
1:B:84:THR:HA	1:B:131:TYR:HB3	1.96	0.46
1:D:185:LYS:HG2	1:D:216:THR:HG21	1.96	0.46
1:B:218:ARG:NH2	1:B:236:ASP:OD2	2.28	0.46
1:A:153:THR:HG22	1:A:158:TRP:CE2	2.51	0.46
1:A:180:ILE:H	1:A:180:ILE:HD12	1.80	0.46
1:A:336:LYS:NZ	1:A:340:ASP:OD1	2.49	0.46
1:B:130:LEU:CD1	1:B:194:ILE:HG21	2.44	0.46
1:B:40:TRP:HE1	2:F:2:FUC:HO2	1.61	0.46
1:B:49:ASN:OD1	1:B:139:HIS:HA	2.15	0.46
1:C:312:TYR:OH	1:C:314:LEU:HD13	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:240:LEU:HB2	1:A:275:THR:OG1	2.16	0.46
1:B:5:VAL:HA	1:B:225:LEU:HD21	1.97	0.46
1:B:24:LEU:HA	1:B:320:PRO:HG3	1.98	0.45
1:D:147:LYS:HB2	1:D:204:THR:HG22	1.97	0.45
1:B:109:PHE:HE2	1:B:113:ILE:HD11	1.80	0.45
1:B:47:ILE:HG21	1:B:52:TYR:HB2	1.99	0.45
1:C:184:ILE:HG21	1:C:213:ILE:HG23	1.97	0.45
1:C:60:ASN:O	1:C:62:LEU:HD12	2.17	0.45
1:C:87:HIS:CE1	2:G:2:FUC:HO4	2.34	0.45
1:A:22:TYR:CG	1:A:39:GLU:HA	2.52	0.45
1:A:24:LEU:HD11	1:A:59:PHE:CD2	2.52	0.45
1:C:97:LYS:N	1:C:186:GLU:OE1	2.46	0.45
1:B:201:VAL:O	1:B:201:VAL:HG23	2.17	0.45
1:C:153:THR:HG22	1:C:158:TRP:CE2	2.52	0.45
1:D:241:GLY:O	1:D:244:GLU:N	2.35	0.44
1:B:136:LEU:HD23	1:B:144:GLY:HA2	2.00	0.44
1:C:295:PRO:HD3	1:C:331:ASN:ND2	2.32	0.44
1:D:8:PHE:HB3	1:D:272:LEU:HD22	1.99	0.44
1:A:94:TYR:HB2	1:A:191:TYR:OH	2.17	0.44
1:B:56:ALA:HB1	1:B:107:THR:HG22	1.99	0.44
1:D:8:PHE:HD2	1:D:272:LEU:HB3	1.83	0.44
1:A:287:TYR:O	1:A:290:GLN:NE2	2.49	0.44
1:C:240:LEU:HD23	1:C:240:LEU:HA	1.78	0.44
1:C:39:GLU:CD	1:C:283:TRP:HB2	2.38	0.44
1:B:150:ASP:OD1	1:B:151:VAL:HG23	2.18	0.44
1:D:279:ILE:HD11	1:D:332:LEU:HD11	1.99	0.44
1:D:5:VAL:HG22	1:D:270:LEU:HD13	2.00	0.44
1:C:145:GLY:HA2	1:C:203:MET:HE2	1.99	0.44
1:C:19:TRP:CE2	1:C:318:LEU:HB2	2.53	0.44
1:A:134:GLN:HB3	1:A:199:PHE:CD1	2.52	0.44
1:B:142:ASN:CB	1:B:175:CYS:HB2	2.48	0.44
1:D:201:VAL:HG22	1:D:229:ARG:HH11	1.82	0.44
1:A:92:ALA:HB3	1:A:102:ASN:HB3	2.00	0.43
1:D:94:TYR:HB2	1:D:191:TYR:OH	2.18	0.43
1:A:126:LEU:HD23	1:A:126:LEU:HA	1.78	0.43
1:B:131:TYR:CD2	1:B:198:TRP:CE3	3.06	0.43
1:A:78:MET:O	1:A:126:LEU:HD21	2.19	0.43
1:D:11:ALA:HA	1:D:79:GLN:OE1	2.18	0.43
1:B:92:ALA:HB3	1:B:102:ASN:HB3	1.99	0.43
1:B:131:TYR:CE2	1:B:198:TRP:CE3	3.06	0.43
1:B:10:GLN:O	1:B:12:LYS:HG3	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:43:SER:HA	1:A:162:TRP:CH2	2.54	0.43
1:B:312:TYR:OH	1:B:314:LEU:HD13	2.18	0.43
1:C:145:GLY:HA2	1:C:203:MET:CE	2.49	0.43
1:C:49:ASN:OD1	1:C:160:ASN:ND2	2.52	0.43
1:D:10:GLN:O	1:D:12:LYS:HG3	2.19	0.43
1:B:35:SER:HB3	1:B:44:LYS:HE2	2.00	0.43
1:D:70:VAL:HG21	1:D:120:ALA:HB1	2.01	0.43
1:A:99:ASP:OD2	1:A:179:LYS:HE3	2.19	0.42
1:D:188:MET:SD	1:D:221:GLN:HG3	2.59	0.42
1:A:199:PHE:O	1:A:229:ARG:HD3	2.19	0.42
1:B:109:PHE:CE2	1:B:113:ILE:HD11	2.54	0.42
1:A:245:ILE:HA	1:A:246:PRO:HD3	1.72	0.42
1:B:136:LEU:HD22	1:B:158:TRP:O	2.19	0.42
1:D:113:ILE:HD13	1:D:113:ILE:HA	1.85	0.42
1:B:39:GLU:OE1	1:B:283:TRP:HB2	2.19	0.42
1:D:228:SER:HA	1:D:235:TYR:CB	2.49	0.42
1:D:312:TYR:OH	1:D:314:LEU:HD13	2.19	0.42
1:C:142:ASN:HB3	1:C:171:ASN:O	2.19	0.42
1:A:42:GLN:HB2	1:A:52:TYR:CE1	2.55	0.42
1:B:94:TYR:OH	1:B:186:GLU:OE2	2.32	0.42
1:D:198:TRP:CZ2	1:D:274:GLU:OE1	2.71	0.42
1:D:228:SER:HB3	1:D:238:VAL:HA	2.01	0.42
1:B:131:TYR:CE2	1:B:198:TRP:CZ3	3.08	0.42
1:C:220:LEU:HA	1:C:220:LEU:HD23	1.92	0.42
1:C:56:ALA:HB1	1:C:107:THR:HG22	2.02	0.42
1:D:142:ASN:HB3	1:D:171:ASN:O	2.19	0.42
1:A:86:LYS:HE2	1:A:132:TYR:CZ	2.55	0.42
1:B:93:MET:O	1:B:114:ILE:HG12	2.20	0.42
1:C:39:GLU:OE1	1:C:283:TRP:HB2	2.20	0.42
1:A:203:MET:HB3	1:A:203:MET:HE3	1.57	0.41
1:B:130:LEU:HD22	1:B:187:ILE:HB	2.01	0.41
1:D:138:TRP:O	1:D:160:ASN:ND2	2.40	0.41
1:A:342:ALA:O	1:A:345:LEU:HB2	2.20	0.41
1:B:142:ASN:HB2	1:B:175:CYS:HB2	2.02	0.41
1:B:342:ALA:HA	1:B:345:LEU:HD12	2.01	0.41
1:B:47:ILE:HD13	1:B:48:PRO:HD2	2.01	0.41
1:C:89:ASP:OD1	1:C:90:GLY:N	2.53	0.41
1:D:126:LEU:HD23	1:D:126:LEU:HA	1.77	0.41
1:D:24:LEU:HA	1:D:320:PRO:HG3	2.02	0.41
1:B:160:ASN:HA	1:B:162:TRP:CZ3	2.56	0.41
1:A:152:GLU:O	2:E:1:NAG:H83	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:26:ALA:HB3	1:B:287:TYR:CD2	2.55	0.41
1:D:148:SER:O	1:D:158:TRP:CD1	2.74	0.41
1:A:229:ARG:HH11	1:A:239:SER:HB3	1.86	0.41
1:B:131:TYR:HE2	1:B:198:TRP:CZ3	2.37	0.41
1:A:285:PHE:HE2	1:A:287:TYR:CE1	2.39	0.41
1:D:188:MET:HG3	1:D:217:VAL:HG22	2.02	0.41
1:B:132:TYR:OH	1:B:180:ILE:HA	2.20	0.41
1:D:201:VAL:HG22	1:D:229:ARG:NH1	2.36	0.41
1:D:8:PHE:CD2	1:D:272:LEU:HB3	2.56	0.41
1:B:240:LEU:HD23	1:B:244:GLU:HB3	2.03	0.41
1:A:136:LEU:HD11	1:A:158:TRP:HE3	1.86	0.40
1:D:150:ASP:N	1:D:150:ASP:OD1	2.53	0.40
1:C:126:LEU:HD23	1:C:126:LEU:HA	1.86	0.40
1:C:218:ARG:NH2	1:C:236:ASP:OD2	2.50	0.40
1:D:63:TYR:HB2	1:D:321:LEU:CD2	2.51	0.40
1:A:158:TRP:CE3	1:A:203:MET:HE1	2.56	0.40
1:A:156:THR:OG1	1:A:157:THR:N	2.53	0.40
1:B:67:LYS:O	1:B:70:VAL:HG22	2.22	0.40
1:A:97:LYS:N	1:A:186:GLU:OE1	2.51	0.40
1:C:16:MET:HG2	1:C:315:ASN:HB2	2.04	0.40
1:C:7:TRP:CZ3	1:C:8:PHE:HD1	2.38	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	321/345 (93%)	301 (94%)	18 (6%)	2 (1%)	25 59
1	B	318/345 (92%)	297 (93%)	18 (6%)	3 (1%)	17 50
1	C	318/345 (92%)	298 (94%)	17 (5%)	3 (1%)	17 50
1	D	316/345 (92%)	295 (93%)	19 (6%)	2 (1%)	25 59

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1273/1380 (92%)	1191 (94%)	72 (6%)	10 (1%)	19	53

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	242	ASP
1	B	172	PHE
1	B	242	ASP
1	C	228	SER
1	C	242	ASP
1	D	242	ASP
1	A	228	SER
1	B	268	SER
1	C	268	SER
1	D	268	SER

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	267/292 (91%)	265 (99%)	2 (1%)	84	93
1	B	265/292 (91%)	257 (97%)	8 (3%)	41	70
1	C	265/292 (91%)	261 (98%)	4 (2%)	65	84
1	D	263/292 (90%)	260 (99%)	3 (1%)	73	88
All	All	1060/1168 (91%)	1043 (98%)	17 (2%)	62	84

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	TYR
1	A	301	TYR
1	B	10	GLN
1	B	13	TYR
1	B	142	ASN

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Mol	Chain	Res	Type
1	B	150	ASP
1	B	171	ASN
1	B	227	ASN
1	B	236	ASP
1	B	243	ASN
1	C	10	GLN
1	C	13	TYR
1	C	19	TRP
1	C	168	ASP
1	D	13	TYR
1	D	86	LYS
1	D	198	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	49	ASN
1	B	79	GLN
1	B	142	ASN
1	B	227	ASN
1	C	88	HIS
1	D	306	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

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	$\# Z > 2$
2	NAG	E	1	2	15,15,15	0.29	0	21,21,21	0.34	0
2	FUC	E	2	2	10,10,11	0.98	1 (10%)	14,14,16	0.99	0
2	NAG	F	1	2	15,15,15	0.33	0	21,21,21	0.31	0
2	FUC	F	2	2	10,10,11	1.15	1 (10%)	14,14,16	0.88	0
2	NAG	G	1	2	15,15,15	0.13	0	21,21,21	0.38	0
2	FUC	G	2	2	10,10,11	0.85	1 (10%)	14,14,16	1.03	1 (7%)
2	NAG	H	1	2	15,15,15	0.41	0	21,21,21	0.45	0
2	FUC	H	2	2	10,10,11	0.53	0	14,14,16	0.83	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	E	1	2	-	2/6/26/26	0/1/1/1
2	FUC	E	2	2	-	-	0/1/1/1
2	NAG	F	1	2	-	2/6/26/26	0/1/1/1
2	FUC	F	2	2	-	-	0/1/1/1
2	NAG	G	1	2	-	2/6/26/26	0/1/1/1
2	FUC	G	2	2	-	-	0/1/1/1
2	NAG	H	1	2	-	2/6/26/26	0/1/1/1
2	FUC	H	2	2	-	-	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	2	FUC	O5-C1	-2.64	1.39	1.43
2	G	2	FUC	O5-C1	-2.48	1.39	1.43
2	E	2	FUC	O5-C1	-2.36	1.39	1.43

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	2	FUC	O2-C2-C1	2.61	114.48	109.15

There are no chirality outliers.

All (8) torsion outliers are listed below:

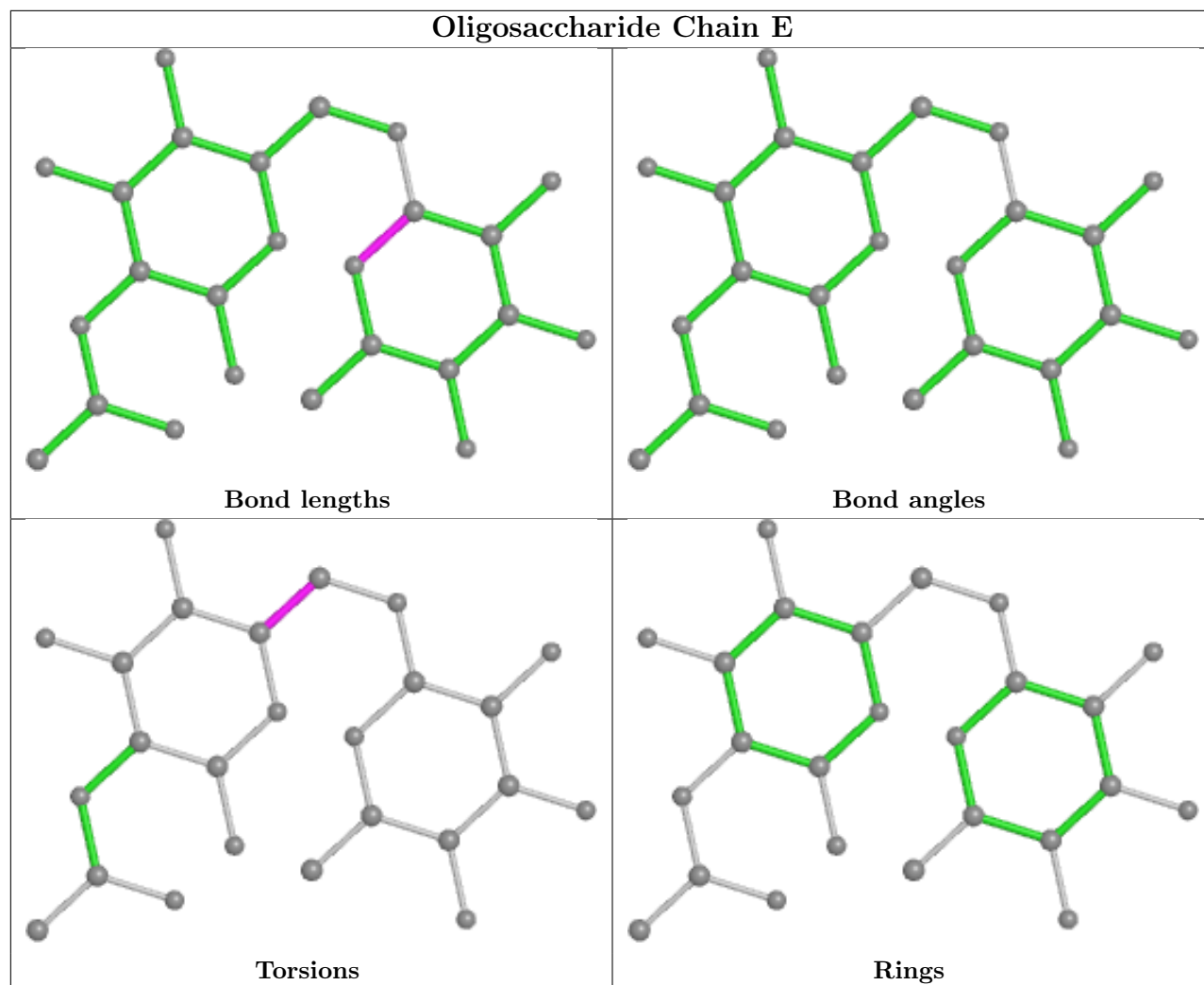
Mol	Chain	Res	Type	Atoms
2	G	1	NAG	C1-C2-N2-C7
2	F	1	NAG	C1-C2-N2-C7
2	H	1	NAG	C1-C2-N2-C7
2	E	1	NAG	C4-C5-C6-O6
2	E	1	NAG	O5-C5-C6-O6
2	G	1	NAG	C3-C2-N2-C7
2	F	1	NAG	C3-C2-N2-C7
2	H	1	NAG	C3-C2-N2-C7

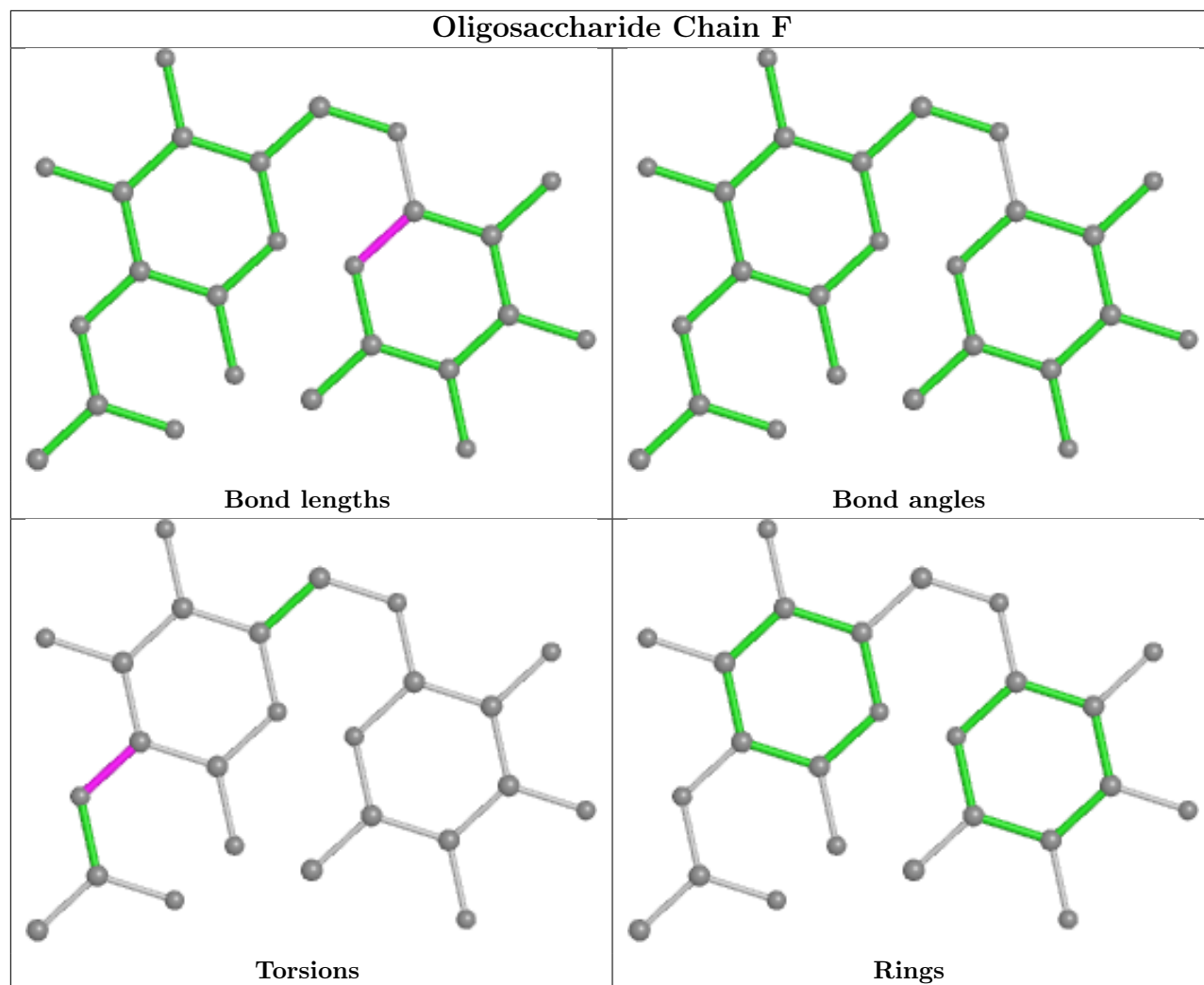
There are no ring outliers.

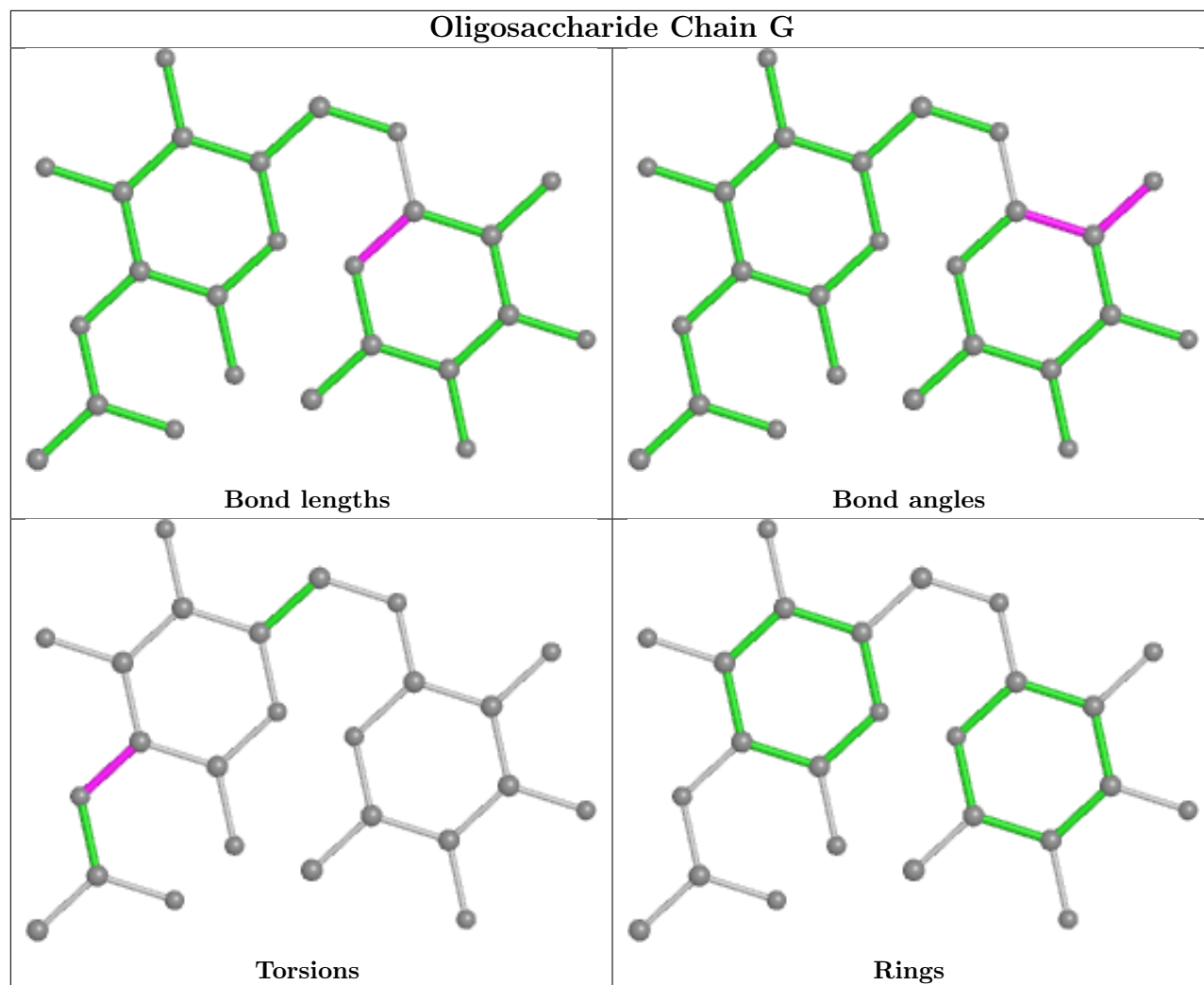
6 monomers are involved in 10 short contacts:

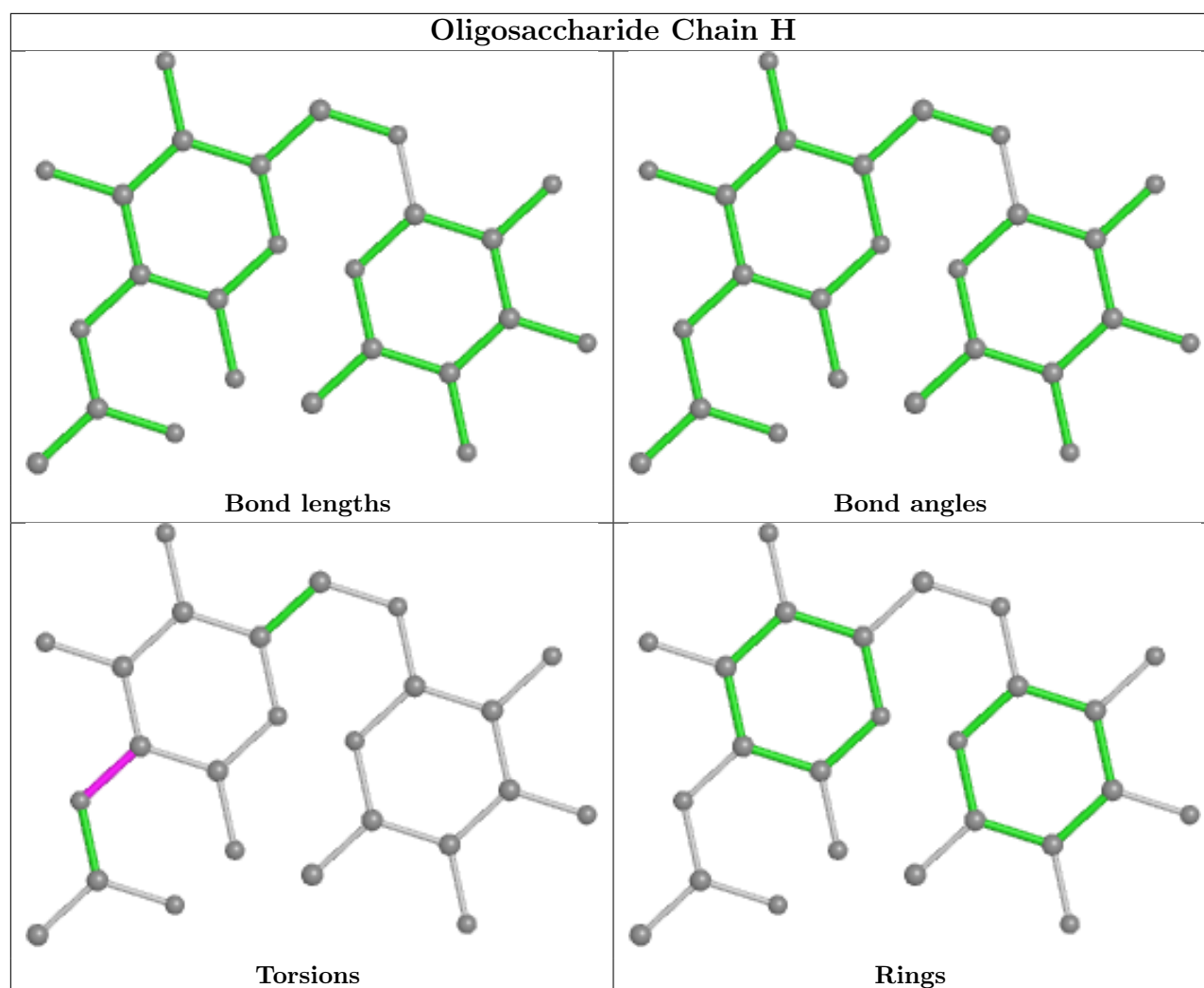
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	2	FUC	1	0
2	G	2	FUC	3	0
2	H	2	FUC	1	0
2	G	1	NAG	3	0
2	F	1	NAG	1	0
2	E	1	NAG	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.









5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	325/345 (94%)	0.18	13 (4%) 38 20	71, 116, 149, 165	0
1	B	322/345 (93%)	0.22	15 (4%) 31 15	68, 106, 155, 171	0
1	C	322/345 (93%)	0.38	18 (5%) 24 11	92, 142, 174, 183	0
1	D	320/345 (92%)	0.44	24 (7%) 14 5	78, 144, 193, 207	0
All	All	1289/1380 (93%)	0.30	70 (5%) 25 12	68, 128, 175, 207	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	271	GLY	5.7
1	D	236	ASP	4.6
1	D	84	THR	4.3
1	D	4	ASN	4.2
1	D	6	ALA	4.2
1	B	267	PRO	4.0
1	A	133	SER	3.6
1	C	273	TYR	3.6
1	D	270	LEU	3.5
1	D	340	ASP	3.4
1	D	224	CYS	3.4
1	B	160	ASN	3.4
1	D	345	LEU	3.3
1	D	269	PRO	3.3
1	D	18	HIS	3.3
1	A	100	ALA	3.2
1	A	224	CYS	3.1
1	D	144	GLY	3.1
1	C	87	HIS	3.0
1	A	20	GLY	3.0
1	C	99	ASP	2.9

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Mol	Chain	Res	Type	RSRZ
1	C	18	HIS	2.9
1	D	149	ASN	2.9
1	C	224	CYS	2.9
1	D	119	GLU	2.9
1	D	271	GLY	2.9
1	B	174	LEU	2.8
1	B	245	ILE	2.8
1	D	133	SER	2.7
1	D	237	PHE	2.7
1	C	189	SER	2.7
1	B	169	GLN	2.7
1	B	345	LEU	2.7
1	D	203	MET	2.7
1	C	15	MET	2.6
1	D	143	GLY	2.6
1	C	238	VAL	2.6
1	B	3	ASP	2.6
1	B	151	VAL	2.6
1	C	268	SER	2.5
1	D	79	GLN	2.4
1	A	88	HIS	2.4
1	C	20	GLY	2.4
1	A	23	SER	2.4
1	B	269	PRO	2.4
1	A	18	HIS	2.4
1	D	198	TRP	2.4
1	B	234	LYS	2.4
1	A	84	THR	2.4
1	A	242	ASP	2.3
1	A	17	ILE	2.3
1	B	271	GLY	2.3
1	C	82	VAL	2.3
1	C	166	ASP	2.2
1	C	89	ASP	2.2
1	C	269	PRO	2.2
1	D	241	GLY	2.2
1	B	146	TYR	2.2
1	A	166	ASP	2.1
1	A	45	PHE	2.1
1	B	87	HIS	2.1
1	B	131	TYR	2.1
1	A	89	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	312	TYR	2.0
1	C	88	HIS	2.0
1	D	151	VAL	2.0
1	C	177	ASP	2.0
1	B	161	SER	2.0
1	C	17	ILE	2.0
1	D	132	TYR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

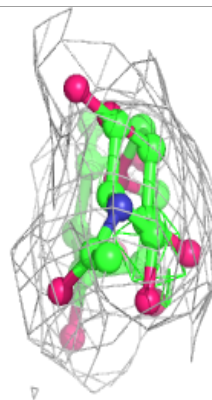
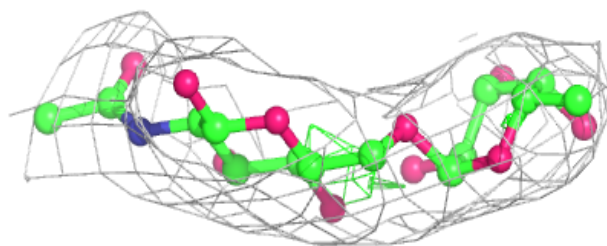
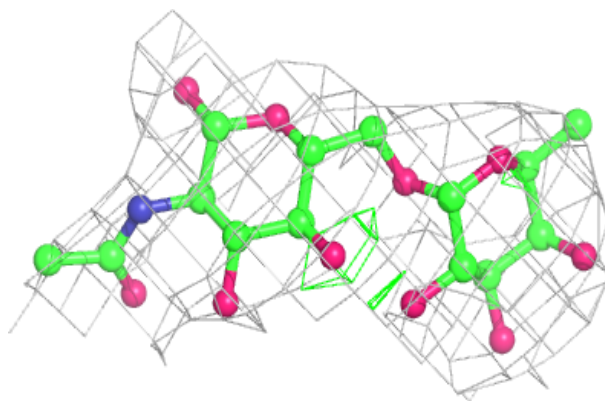
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	E	1	15/15	0.78	0.22	98,98,98,98	0
2	NAG	F	1	15/15	0.86	0.20	106,106,106,106	0
2	NAG	H	1	15/15	0.86	0.24	130,130,130,130	0
2	FUC	G	2	10/11	0.88	0.38	110,110,110,110	0
2	NAG	G	1	15/15	0.90	0.16	111,111,111,111	0
2	FUC	H	2	10/11	0.91	0.29	123,123,123,123	0
2	FUC	F	2	10/11	0.94	0.32	97,97,97,97	0
2	FUC	E	2	10/11	0.97	0.33	94,94,94,94	0

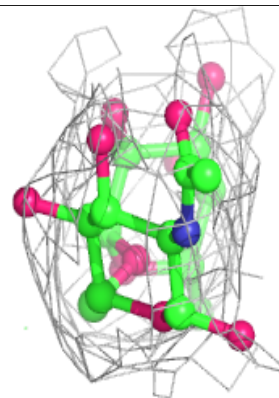
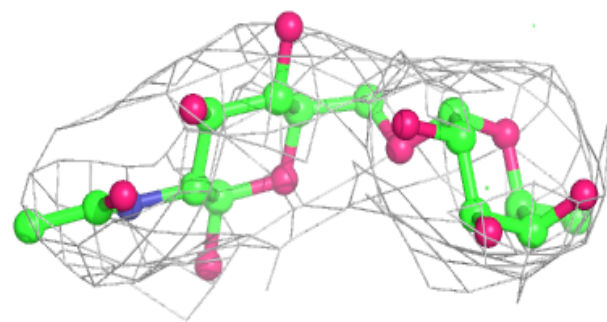
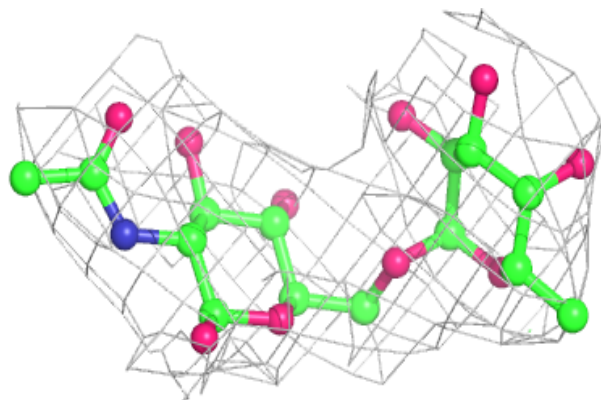
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain E:

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

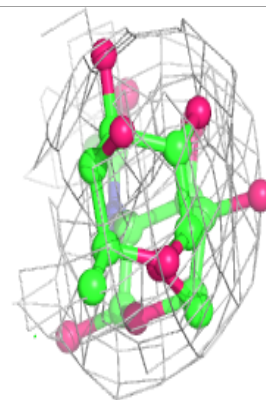
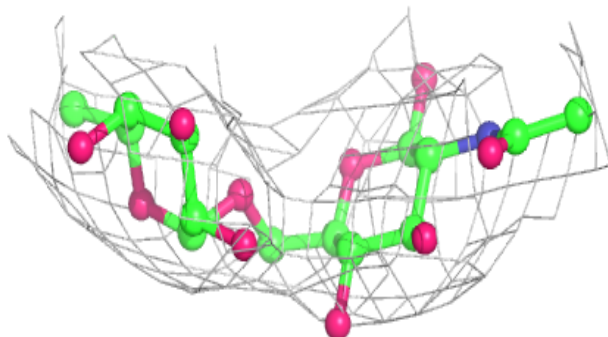
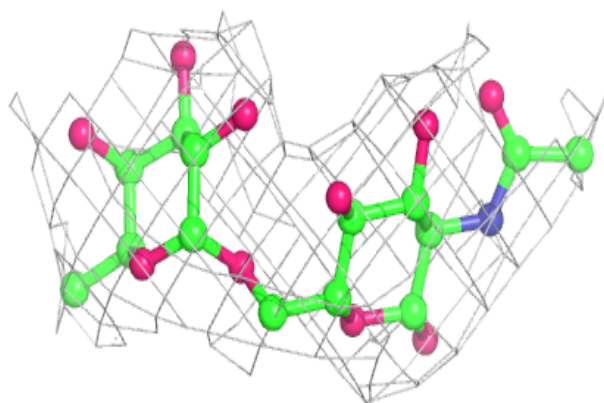
**Electron density around Chain F:**

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

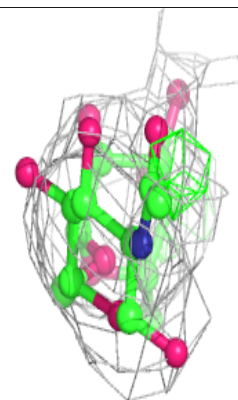
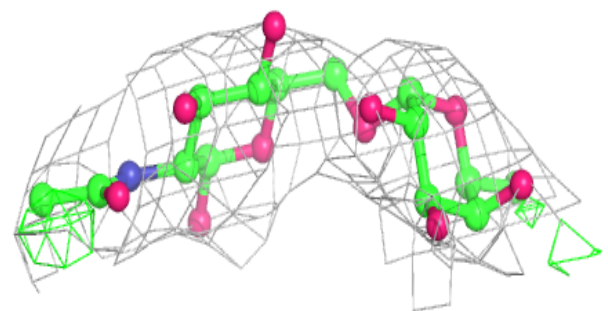
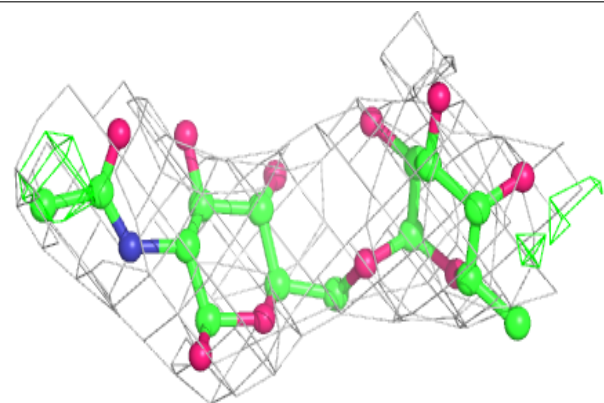


Electron density around Chain G:

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

**Electron density around Chain H:**

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



6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.