



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:40 AM UTC

PDB ID : 3ZNM / pdb_00003znm
Title : H5 Haemagglutinin in Complex with Sialyl-Lewis X
Authors : Xiong, X.; Tuzikov, A.; Coombs, P.; Martin, S.R.; Walker, P.A.; Gamblin, S.J.; Bovin, N.; Skehel, J.J.
Deposited on : 2013-02-15
Resolution : 2.40 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

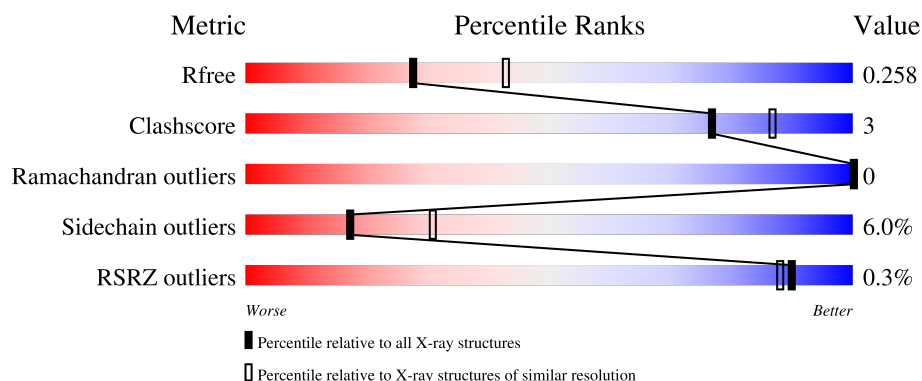
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 Å.

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



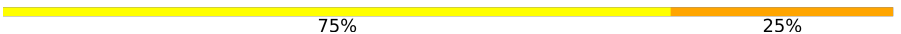
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	326	 87% 11% ..
1	C	326	 88% 10% .
1	E	326	 87% 10% ..
2	B	166	 84% 14% ..
2	D	166	 84% 14% ..

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Mol	Chain	Length	Quality of chain
2	F	166	 85% 13% ..
3	G	4	 75% 25%
3	H	4	 75% 25%
3	I	4	 75% 25%

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
4	NAG	B	1164	X	-	-	-
4	NAG	D	1164	X	-	-	-
4	NAG	F	1164	X	-	-	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 12237 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 HAEMAGGLUTININ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	321	Total	C	N	O	S	0	0	0
			2547	1609	439	484	15			
1	C	321	Total	C	N	O	S	0	0	0
			2547	1609	439	484	15			
1	E	321	Total	C	N	O	S	0	0	0
			2547	1609	439	484	15			

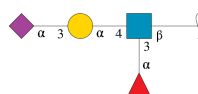
There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	325	THR	ARG	conflict	UNP Q6DQ34
C	325	THR	ARG	conflict	UNP Q6DQ34
E	325	THR	ARG	conflict	UNP Q6DQ34

- Molecule 2 is a protein called HAEMAGGLUTININ.

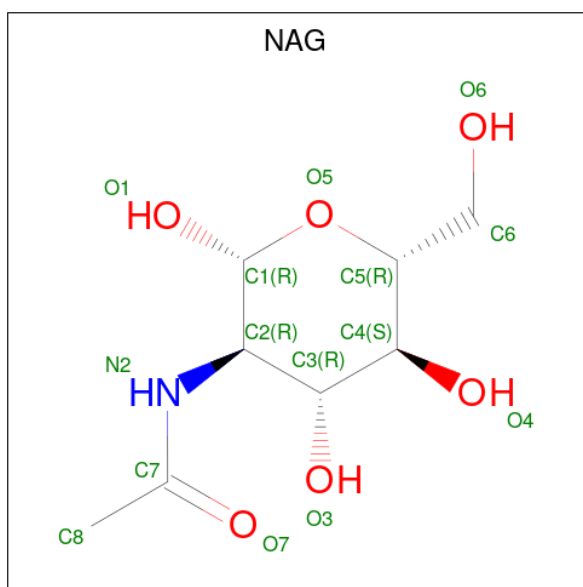
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	163	Total	C	N	O	S	0	0	0
			1318	822	227	261	8			
2	D	163	Total	C	N	O	S	0	0	0
			1318	822	227	261	8			
2	F	163	Total	C	N	O	S	0	0	0
			1318	822	227	261	8			

- Molecule 3 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-3)-alpha-D-galactopyranose-(1-4)-[alpha-L-fucopyranose-(1-3)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	G	4	Total	C	N	O	0	0	0
			56	31	2	23			
3	H	4	Total	C	N	O	0	0	0
			56	31	2	23			
3	I	4	Total	C	N	O	0	0	0
			56	31	2	23			

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



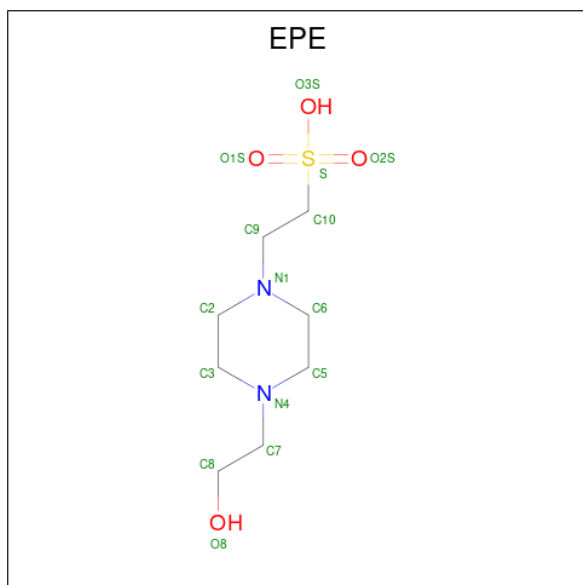
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		
4	C	1	Total	C	N	O	0	0
			14	8	1	5		
4	C	1	Total	C	N	O	0	0
			14	8	1	5		
4	D	1	Total	C	N	O	0	0
			14	8	1	5		
4	E	1	Total	C	N	O	0	0
			14	8	1	5		
4	E	1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

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	F	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: $C_8H_{18}N_2O_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
5	D	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
5	F	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	64	Total	O	0	0
			64	64		
6	B	48	Total	O	0	0
			48	48		
6	C	60	Total	O	0	0
			60	60		
6	D	41	Total	O	0	0
			41	41		
6	E	51	Total	O	0	0
			51	51		

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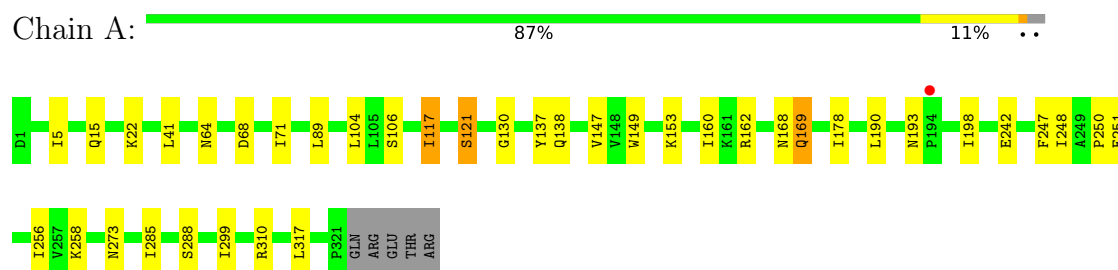
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	F	39	Total	O	0	0
			39	39		

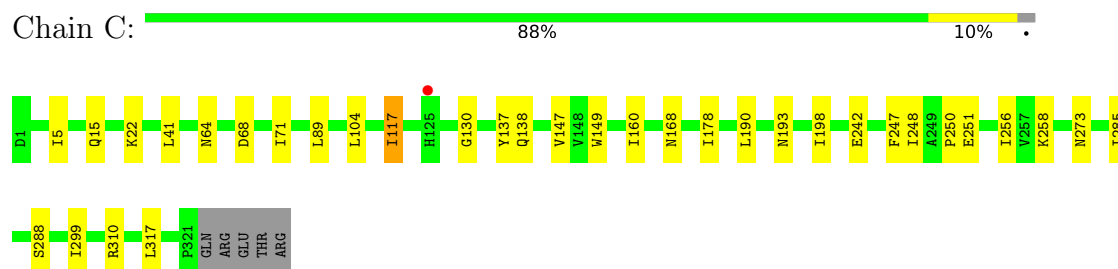
3 Residue-property plots

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.

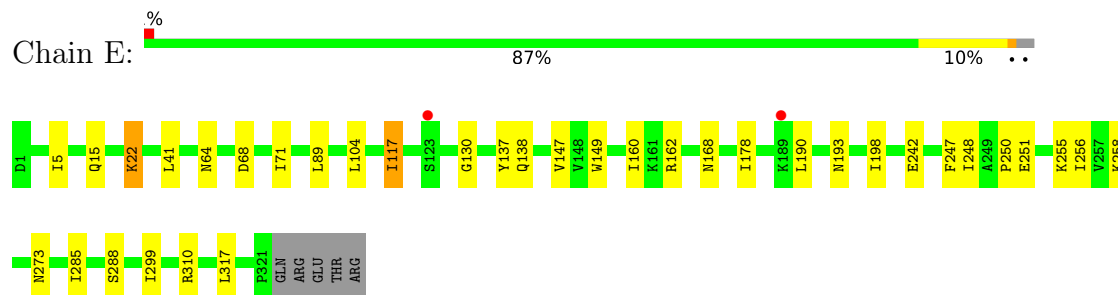
• Molecule 1: HAEMAGGLUTININ



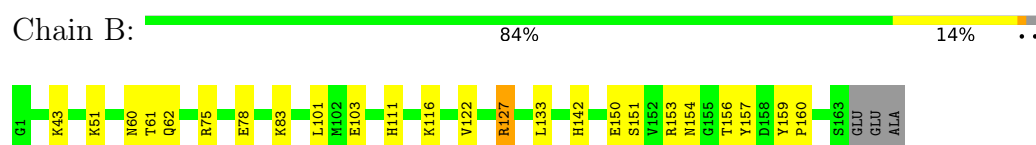
• Molecule 1: HAEMAGGLUTININ



• Molecule 1: HAEMAGGLUTININ



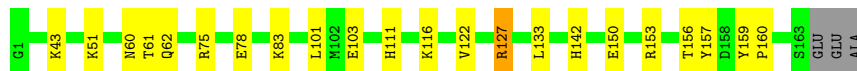
• Molecule 2: HAEMAGGLUTININ



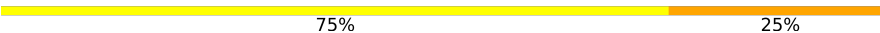
• Molecule 2: HAEMAGGLUTININ

Chain D:  84% 14% ..

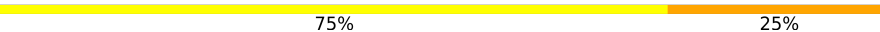
• Molecule 2: HAEMAGGLUTININ

Chain F:  85% 13% ..

• Molecule 3: N-acetyl-alpha-neuraminic acid-(2-3)-alpha-D-galactopyranose-(1-4)-[alpha-L-fucopyranose-(1-3)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:  75% 25%

• Molecule 3: N-acetyl-alpha-neuraminic acid-(2-3)-alpha-D-galactopyranose-(1-4)-[alpha-L-fucopyranose-(1-3)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H:  75% 25%

• Molecule 3: N-acetyl-alpha-neuraminic acid-(2-3)-alpha-D-galactopyranose-(1-4)-[alpha-L-fucopyranose-(1-3)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  75% 25%

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	175.56Å 101.36Å 161.20Å 90.00° 111.29° 90.00°	Depositor
Resolution (Å)	37.15 – 2.40 37.15 – 2.40	Depositor EDS
% Data completeness (in resolution range)	75.6 (37.15-2.40) 75.7 (37.15-2.40)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.69 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.238 , 0.256 0.241 , 0.258	Depositor DCC
R_{free} test set	3882 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	33.4	Xtriage
Anisotropy	0.043	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 16.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.468 for $1/2^*h+3/2^*k, 1/2^*h-1/2^*k, -1/2^*h-1/2^*k-l$ 0.469 for $1/2^*h-3/2^*k, -1/2^*h-1/2^*k, -1/2^*h+1/2^*k-l$	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12237	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.23% 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: EPE, SIA, NAG, GLA, 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	0.40	0/2609	0.64	0/3545
1	C	0.39	0/2609	0.63	0/3545
1	E	0.40	0/2609	0.64	0/3545
2	B	0.43	0/1345	0.73	0/1810
2	D	0.43	0/1345	0.73	0/1810
2	F	0.43	0/1345	0.73	0/1810
All	All	0.41	0/11862	0.67	0/16065

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2547	0	2490	16	0
1	C	2547	0	2490	12	0
1	E	2547	0	2490	13	0
2	B	1318	0	1223	11	0
2	D	1318	0	1223	12	0
2	F	1318	0	1223	9	0
3	G	56	0	49	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	H	56	0	49	1	0
3	I	56	0	49	1	0
4	A	28	0	26	0	0
4	B	14	0	13	0	0
4	C	28	0	26	0	0
4	D	14	0	13	0	0
4	E	28	0	26	1	0
4	F	14	0	13	0	0
5	B	15	0	18	0	0
5	D	15	0	18	0	0
5	F	15	0	18	0	0
6	A	64	0	0	1	0
6	B	48	0	0	0	0
6	C	60	0	0	0	0
6	D	41	0	0	0	0
6	E	51	0	0	0	0
6	F	39	0	0	0	0
All	All	12237	0	11457	67	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (67) 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:A:121:SER:OG	1:A:162:ARG:NH2	1.85	1.08
1:A:5:ILE:HD11	2:B:122:VAL:HG21	1.72	0.71
1:C:5:ILE:HD11	2:D:122:VAL:HG21	1.74	0.69
1:E:5:ILE:HD11	2:F:122:VAL:HG21	1.74	0.68
2:D:127:ARG:HG2	2:D:127:ARG:HH11	1.66	0.60
2:F:127:ARG:HH11	2:F:127:ARG:HG2	1.66	0.60
2:B:127:ARG:HH11	2:B:127:ARG:HG2	1.66	0.59
2:D:156:THR:HG22	2:D:156:THR:O	2.03	0.59
2:B:156:THR:O	2:B:156:THR:HG22	2.03	0.57
2:F:156:THR:O	2:F:156:THR:HG22	2.03	0.57
2:F:150:GLU:OE1	2:F:153:ARG:NH1	2.38	0.57
2:D:75:ARG:NH1	2:D:78:GLU:OE1	2.38	0.57
1:A:121:SER:HG	1:A:162:ARG:NH2	2.00	0.56
2:B:150:GLU:OE1	2:B:153:ARG:NH1	2.38	0.56
2:F:75:ARG:NH1	2:F:78:GLU:OE1	2.38	0.56
1:E:68:ASP:O	1:E:71:ILE:HG22	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:150:GLU:OE1	2:D:153:ARG:NH1	2.38	0.55
2:B:75:ARG:NH1	2:B:78:GLU:OE1	2.38	0.55
1:A:68:ASP:O	1:A:71:ILE:HG22	2.06	0.55
1:C:68:ASP:O	1:C:71:ILE:HG22	2.06	0.54
1:A:117:ILE:HB	1:A:250:PRO:O	2.08	0.54
1:E:117:ILE:HB	1:E:250:PRO:O	2.08	0.53
1:C:117:ILE:HB	1:C:250:PRO:O	2.08	0.53
1:A:160:ILE:O	1:A:242:GLU:HA	2.09	0.53
1:C:160:ILE:O	1:C:242:GLU:HA	2.09	0.53
1:E:160:ILE:O	1:E:242:GLU:HA	2.09	0.52
2:D:51:LYS:HE2	2:D:103:GLU:HB3	1.92	0.51
2:B:51:LYS:HE2	2:B:103:GLU:HB3	1.93	0.51
1:A:15:GLN:O	1:A:310:ARG:NH2	2.44	0.50
1:E:15:GLN:O	1:E:310:ARG:NH2	2.44	0.50
1:C:15:GLN:O	1:C:310:ARG:NH2	2.44	0.50
2:F:51:LYS:HE2	2:F:103:GLU:HB3	1.93	0.50
2:D:142:HIS:NE2	2:D:157:TYR:OH	2.46	0.48
2:B:142:HIS:NE2	2:B:157:TYR:OH	2.46	0.47
2:D:159:TYR:HB3	2:D:160:PRO:HD3	1.97	0.47
1:E:190:LEU:HD21	3:I:3:SIA:O10	2.15	0.46
2:B:159:TYR:HB3	2:B:160:PRO:HD3	1.98	0.46
1:C:147:VAL:HB	1:C:248:ILE:HG22	1.97	0.46
1:A:190:LEU:HD21	3:G:3:SIA:O10	2.16	0.46
1:C:190:LEU:HD21	3:H:3:SIA:O10	2.16	0.46
1:E:147:VAL:HB	1:E:248:ILE:HG22	1.97	0.46
1:A:147:VAL:HB	1:A:248:ILE:HG22	1.97	0.45
2:F:159:TYR:HB3	2:F:160:PRO:HD3	1.97	0.45
1:E:130:GLY:HA3	1:E:149:TRP:HB3	1.99	0.44
1:C:130:GLY:HA3	1:C:149:TRP:HB3	1.99	0.44
1:A:106:SER:HB3	6:A:2038:HOH:O	2.17	0.44
1:A:130:GLY:HA3	1:A:149:TRP:HB3	1.99	0.43
1:E:22:LYS:HE3	4:E:1322:NAG:H82	2.00	0.43
1:C:198:ILE:HD13	1:C:247:PHE:HD1	1.85	0.42
2:B:156:THR:O	2:B:156:THR:CG2	2.67	0.42
1:A:169:GLN:HE21	1:A:169:GLN:HB2	1.75	0.42
2:D:156:THR:O	2:D:156:THR:CG2	2.68	0.42
1:A:198:ILE:HD13	1:A:247:PHE:HD1	1.84	0.41
1:E:64:ASN:OD1	1:E:64:ASN:C	2.64	0.41
1:E:198:ILE:HD13	1:E:247:PHE:HD1	1.85	0.41
1:A:137:TYR:CE2	1:A:138:GLN:HG3	2.56	0.41
1:A:317:LEU:HB3	2:B:111:HIS:CG	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:317:LEU:HB3	2:D:111:HIS:CG	2.56	0.41
1:E:137:TYR:CE2	1:E:138:GLN:HG3	2.56	0.41
2:F:142:HIS:NE2	2:F:157:TYR:OH	2.46	0.41
2:D:151:SER:HA	2:D:154:ASN:HB2	2.03	0.41
1:A:64:ASN:OD1	1:A:64:ASN:C	2.64	0.41
1:C:64:ASN:C	1:C:64:ASN:OD1	2.64	0.40
2:B:151:SER:HA	2:B:154:ASN:HB2	2.03	0.40
1:C:137:TYR:CE2	1:C:138:GLN:HG3	2.56	0.40
2:D:127:ARG:HH11	2:D:127:ARG:CG	2.33	0.40
1:E:317:LEU:HB3	2:F:111:HIS:CG	2.56	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	319/326 (98%)	311 (98%)	8 (2%)	0	100	100
1	C	319/326 (98%)	311 (98%)	8 (2%)	0	100	100
1	E	319/326 (98%)	311 (98%)	8 (2%)	0	100	100
2	B	161/166 (97%)	159 (99%)	2 (1%)	0	100	100
2	D	161/166 (97%)	159 (99%)	2 (1%)	0	100	100
2	F	161/166 (97%)	159 (99%)	2 (1%)	0	100	100
All	All	1440/1476 (98%)	1410 (98%)	30 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	289/294 (98%)	271 (94%)	18 (6%)	16	29
1	C	289/294 (98%)	274 (95%)	15 (5%)	21	36
1	E	289/294 (98%)	272 (94%)	17 (6%)	18	31
2	B	139/141 (99%)	130 (94%)	9 (6%)	15	27
2	D	139/141 (99%)	130 (94%)	9 (6%)	15	27
2	F	139/141 (99%)	130 (94%)	9 (6%)	15	27
All	All	1284/1305 (98%)	1207 (94%)	77 (6%)	17	31

All (77) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LYS
1	A	41	LEU
1	A	89	LEU
1	A	104	LEU
1	A	117	ILE
1	A	121	SER
1	A	153	LYS
1	A	168	ASN
1	A	169	GLN
1	A	178	ILE
1	A	193	ASN
1	A	251	GLU
1	A	256	ILE
1	A	258	LYS
1	A	273	ASN
1	A	285	ILE
1	A	288	SER
1	A	299	ILE
2	B	43	LYS
2	B	60	ASN
2	B	61	THR
2	B	62	GLN

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Mol	Chain	Res	Type
2	B	83	LYS
2	B	101	LEU
2	B	116	LYS
2	B	127	ARG
2	B	133	LEU
1	C	22	LYS
1	C	41	LEU
1	C	89	LEU
1	C	104	LEU
1	C	117	ILE
1	C	168	ASN
1	C	178	ILE
1	C	193	ASN
1	C	251	GLU
1	C	256	ILE
1	C	258	LYS
1	C	273	ASN
1	C	285	ILE
1	C	288	SER
1	C	299	ILE
2	D	43	LYS
2	D	60	ASN
2	D	61	THR
2	D	62	GLN
2	D	83	LYS
2	D	101	LEU
2	D	116	LYS
2	D	127	ARG
2	D	133	LEU
1	E	22	LYS
1	E	41	LEU
1	E	89	LEU
1	E	104	LEU
1	E	117	ILE
1	E	162	ARG
1	E	168	ASN
1	E	178	ILE
1	E	193	ASN
1	E	251	GLU
1	E	255	LYS
1	E	256	ILE
1	E	258	LYS

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Mol	Chain	Res	Type
1	E	273	ASN
1	E	285	ILE
1	E	288	SER
1	E	299	ILE
2	F	43	LYS
2	F	60	ASN
2	F	61	THR
2	F	62	GLN
2	F	83	LYS
2	F	101	LEU
2	F	116	LYS
2	F	127	ARG
2	F	133	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	A	84	ASN
1	A	103	HIS
1	A	169	GLN
2	B	15	GLN
1	C	15	GLN
2	D	15	GLN
2	D	62	GLN
1	E	15	GLN
2	F	15	GLN
2	F	62	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 ⓘ

12 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	G	1	3	15,15,15	0.65	0	21,21,21	1.37	2 (9%)
3	GLA	G	2	3	11,11,12	0.38	0	15,15,17	1.35	2 (13%)
3	SIA	G	3	3	20,20,21	0.55	0	21,28,31	1.07	3 (14%)
3	FUC	G	4	3	10,10,11	0.63	0	14,14,16	1.94	2 (14%)
3	NAG	H	1	3	15,15,15	0.64	0	21,21,21	1.37	2 (9%)
3	GLA	H	2	3	11,11,12	0.29	0	15,15,17	1.21	3 (20%)
3	SIA	H	3	3	20,20,21	0.57	0	21,28,31	1.12	3 (14%)
3	FUC	H	4	3	10,10,11	0.63	0	14,14,16	1.64	2 (14%)
3	NAG	I	1	3	15,15,15	0.61	0	21,21,21	1.30	2 (9%)
3	GLA	I	2	3	11,11,12	0.30	0	15,15,17	1.10	1 (6%)
3	SIA	I	3	3	20,20,21	0.54	0	21,28,31	1.17	3 (14%)
3	FUC	I	4	3	10,10,11	0.65	0	14,14,16	1.87	2 (14%)

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	G	1	3	-	2/6/26/26	0/1/1/1
3	GLA	G	2	3	-	0/2/19/22	0/1/1/1
3	SIA	G	3	3	-	0/18/34/38	0/1/1/1
3	FUC	G	4	3	-	-	0/1/1/1
3	NAG	H	1	3	-	2/6/26/26	0/1/1/1
3	GLA	H	2	3	-	0/2/19/22	0/1/1/1
3	SIA	H	3	3	-	0/18/34/38	0/1/1/1
3	FUC	H	4	3	-	-	0/1/1/1
3	NAG	I	1	3	-	2/6/26/26	0/1/1/1
3	GLA	I	2	3	-	0/2/19/22	0/1/1/1
3	SIA	I	3	3	-	1/18/34/38	0/1/1/1
3	FUC	I	4	3	-	-	0/1/1/1

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	4	FUC	C1-C2-C3	5.44	117.56	109.64
3	G	4	FUC	C1-C2-C3	5.36	117.44	109.64
3	H	1	NAG	C1-C2-N2	-4.72	105.25	110.73
3	G	1	NAG	C1-C2-N2	-4.60	105.40	110.73
3	I	1	NAG	C1-C2-N2	-4.60	105.40	110.73
3	H	4	FUC	C1-C2-C3	4.28	115.88	109.64
3	G	4	FUC	C2-C3-C4	3.62	117.22	110.86
3	H	4	FUC	C2-C3-C4	3.25	116.57	110.86
3	G	2	GLA	C1-C2-C3	3.22	114.33	109.64
3	I	4	FUC	C2-C3-C4	3.10	116.32	110.86
3	H	3	SIA	C4-C5-N5	-2.93	104.67	110.44
3	I	3	SIA	C4-C5-N5	-2.79	104.94	110.44
3	G	1	NAG	O4-C4-C3	2.77	116.91	110.38
3	I	3	SIA	O6-C2-C1	2.66	112.75	107.72
3	H	1	NAG	O4-C4-C3	2.61	116.53	110.38
3	H	2	GLA	C1-C2-C3	2.57	113.39	109.64
3	I	1	NAG	O4-C4-C3	2.56	116.41	110.38
3	G	3	SIA	C4-C5-N5	-2.55	105.42	110.44
3	I	2	GLA	C1-C2-C3	2.20	112.85	109.64
3	H	3	SIA	O6-C2-C1	2.17	111.81	107.72
3	G	3	SIA	O6-C2-C1	2.13	111.73	107.72
3	H	3	SIA	O1B-C1-C2	2.09	118.15	112.71
3	I	3	SIA	O1B-C1-C2	2.08	118.12	112.71
3	G	2	GLA	C1-O5-C5	2.06	114.95	112.19
3	G	3	SIA	O1B-C1-C2	2.06	118.07	112.71
3	H	2	GLA	O5-C5-C6	2.04	111.64	107.66
3	H	2	GLA	C1-O5-C5	2.03	114.90	112.19

There are no chirality outliers.

All (7) torsion outliers are listed below:

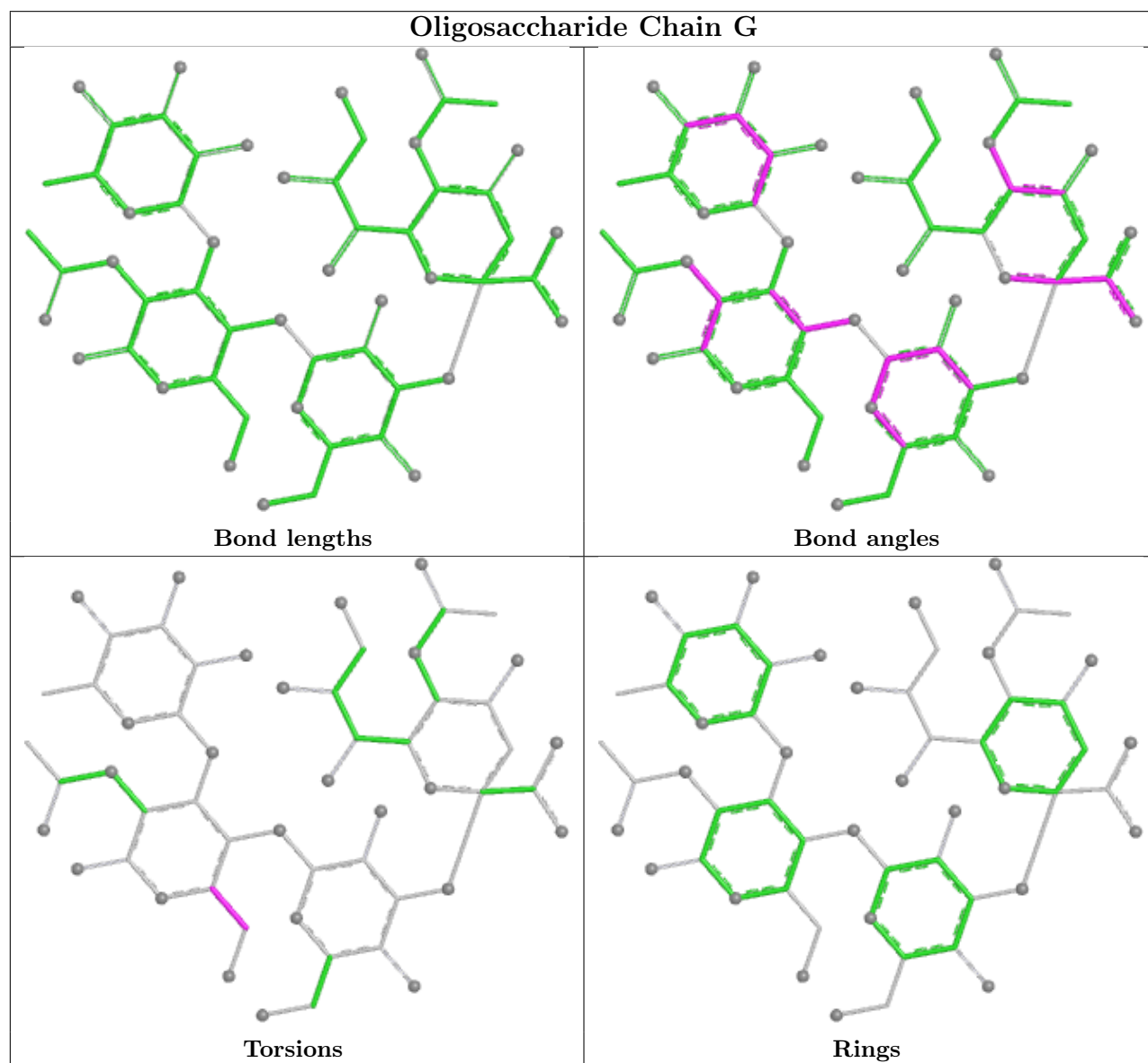
Mol	Chain	Res	Type	Atoms
3	G	1	NAG	C4-C5-C6-O6
3	H	1	NAG	C4-C5-C6-O6
3	H	1	NAG	O5-C5-C6-O6
3	G	1	NAG	O5-C5-C6-O6
3	I	1	NAG	C4-C5-C6-O6
3	I	1	NAG	O5-C5-C6-O6
3	I	3	SIA	C6-C7-C8-O8

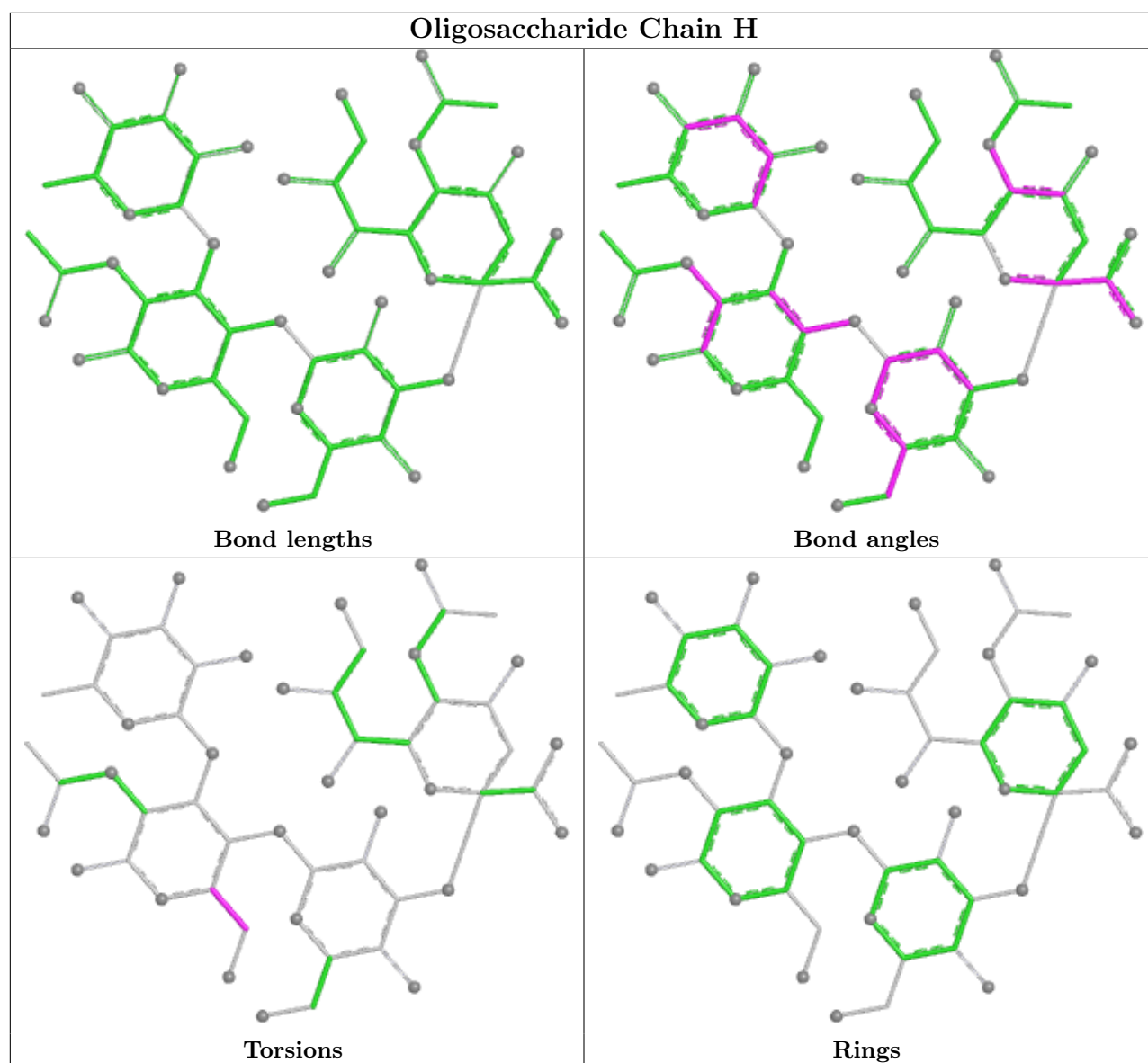
There are no ring outliers.

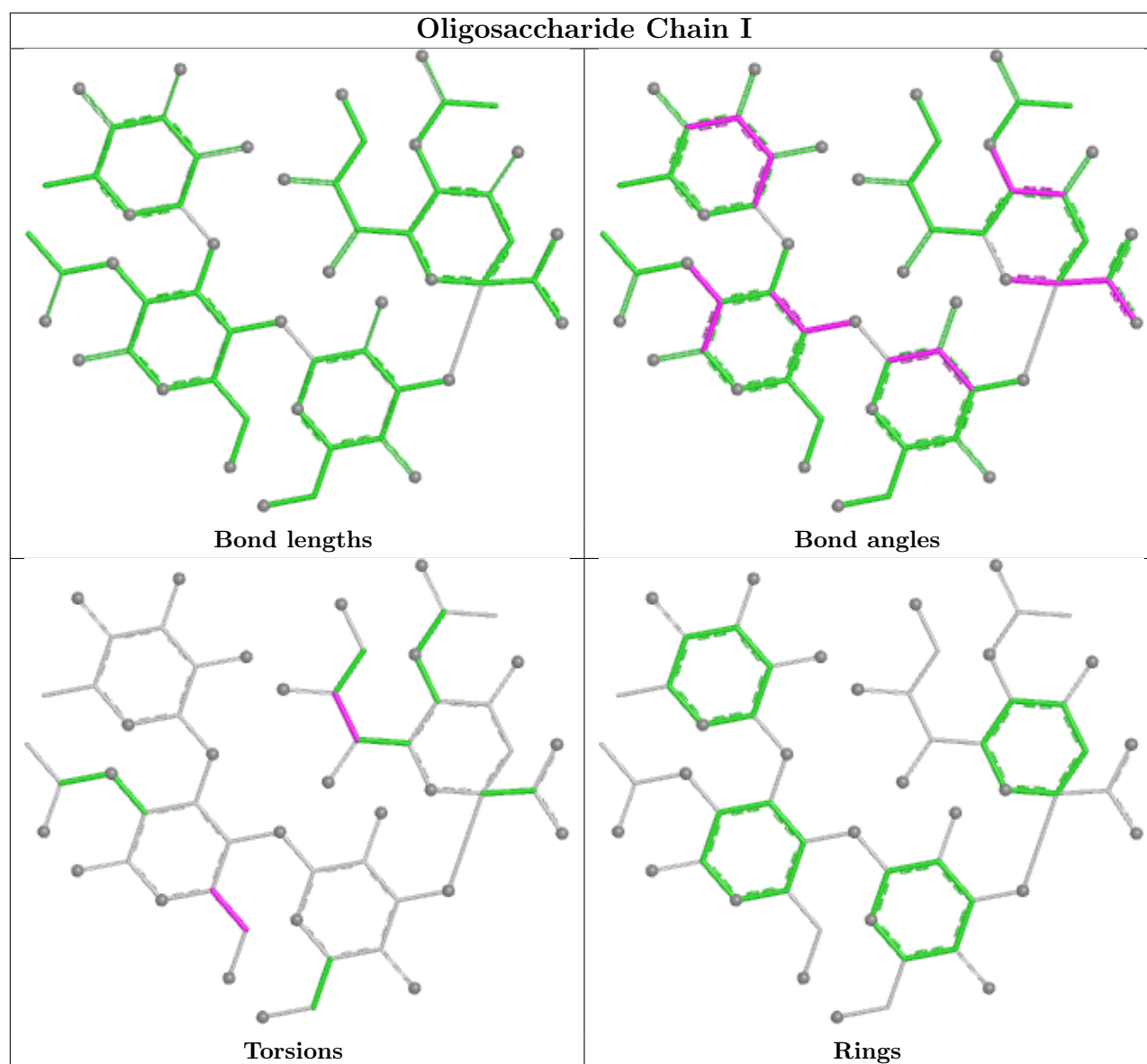
3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	3	SIA	1	0
3	I	3	SIA	1	0
3	G	3	SIA	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](#)

12 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	EPE	B	1165	-	15,15,15	2.00	1 (6%)	19,20,20	2.68	11 (57%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	EPE	F	1165	-	15,15,15	2.04	1 (6%)	19,20,20	2.70	9 (47%)
4	NAG	E	1323	1	14,14,15	0.60	0	17,19,21	1.58	3 (17%)
4	NAG	A	1323	1	14,14,15	0.66	0	17,19,21	1.53	3 (17%)
4	NAG	F	1164	2	14,14,15	0.53	0	17,19,21	1.29	1 (5%)
4	NAG	A	1322	1	14,14,15	0.47	0	17,19,21	0.87	0
4	NAG	E	1322	1	14,14,15	0.46	0	17,19,21	1.17	2 (11%)
4	NAG	C	1322	1	14,14,15	0.51	0	17,19,21	0.98	1 (5%)
4	NAG	C	1323	1	14,14,15	0.66	0	17,19,21	1.55	3 (17%)
4	NAG	B	1164	2	14,14,15	0.44	0	17,19,21	0.84	0
4	NAG	D	1164	2	14,14,15	0.46	0	17,19,21	1.57	2 (11%)
5	EPE	D	1165	-	15,15,15	2.02	1 (6%)	19,20,20	2.64	7 (36%)

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
5	EPE	B	1165	-	-	6/9/19/19	0/1/1/1
5	EPE	F	1165	-	-	4/9/19/19	0/1/1/1
4	NAG	E	1323	1	-	1/6/23/26	0/1/1/1
4	NAG	A	1323	1	-	1/6/23/26	0/1/1/1
4	NAG	C	1323	1	-	1/6/23/26	0/1/1/1
4	NAG	F	1164	2	1/1/5/7	0/6/23/26	0/1/1/1
4	NAG	A	1322	1	-	2/6/23/26	0/1/1/1
4	NAG	E	1322	1	-	2/6/23/26	0/1/1/1
4	NAG	C	1322	1	-	2/6/23/26	0/1/1/1
4	NAG	B	1164	2	1/1/5/7	0/6/23/26	0/1/1/1
4	NAG	D	1164	2	1/1/5/7	0/6/23/26	0/1/1/1
5	EPE	D	1165	-	-	8/9/19/19	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	F	1165	EPE	C10-S	-7.55	1.67	1.77
5	D	1165	EPE	C10-S	-7.48	1.67	1.77
5	B	1165	EPE	C10-S	-7.34	1.67	1.77

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	1165	EPE	C5-N4-C3	5.32	120.30	108.84
4	D	1164	NAG	C1-O5-C5	5.29	119.28	112.19
5	D	1165	EPE	C5-N4-C3	5.24	120.12	108.84
5	B	1165	EPE	C5-N4-C3	5.09	119.81	108.84
5	D	1165	EPE	C9-N1-C2	5.00	124.57	111.24
5	D	1165	EPE	C6-N1-C2	4.59	118.73	108.84
5	F	1165	EPE	C6-N1-C2	4.53	118.59	108.84
5	B	1165	EPE	C6-N1-C2	4.45	118.43	108.84
4	F	1164	NAG	C1-O5-C5	4.12	117.71	112.19
5	F	1165	EPE	C9-N1-C2	3.91	121.67	111.24
5	B	1165	EPE	C9-N1-C6	3.86	121.52	111.24
5	B	1165	EPE	C7-N4-C5	3.80	121.36	111.24
5	F	1165	EPE	C5-C6-N1	-3.80	103.00	110.65
5	D	1165	EPE	C7-N4-C3	3.79	121.33	111.24
5	F	1165	EPE	C7-N4-C3	3.59	120.80	111.24
4	E	1323	NAG	C1-C2-N2	3.32	115.67	110.43
5	B	1165	EPE	C9-N1-C2	3.31	120.05	111.24
4	E	1322	NAG	C1-O5-C5	3.27	116.56	112.19
5	B	1165	EPE	C3-C2-N1	-3.25	104.10	110.65
4	C	1323	NAG	C1-C2-N2	3.19	115.46	110.43
5	F	1165	EPE	C9-N1-C6	3.18	119.72	111.24
5	D	1165	EPE	O1S-S-C10	3.12	111.44	106.73
4	C	1323	NAG	C2-N2-C7	3.05	126.98	122.90
4	A	1323	NAG	C1-C2-N2	2.93	115.05	110.43
5	F	1165	EPE	C7-N4-C5	2.87	118.90	111.24
5	B	1165	EPE	C7-N4-C3	2.85	118.82	111.24
4	A	1323	NAG	C2-N2-C7	2.84	126.71	122.90
5	B	1165	EPE	O1S-S-C10	2.82	110.98	106.73
4	E	1323	NAG	C2-N2-C7	2.80	126.65	122.90
5	D	1165	EPE	C7-N4-C5	2.74	118.55	111.24
5	F	1165	EPE	O1S-S-C10	2.33	110.25	106.73
4	E	1322	NAG	O5-C5-C6	2.32	112.18	107.66
4	A	1323	NAG	O7-C7-C8	-2.31	117.94	122.05
4	D	1164	NAG	O5-C1-C2	2.21	114.71	111.29
5	F	1165	EPE	O3S-S-C10	2.19	110.29	106.00
4	E	1323	NAG	O7-C7-C8	-2.17	118.19	122.05
4	C	1323	NAG	O7-C7-C8	-2.08	118.35	122.05
5	B	1165	EPE	C8-C7-N4	-2.07	106.19	113.44
5	B	1165	EPE	C2-C3-N4	-2.05	106.51	110.65
5	D	1165	EPE	C9-N1-C6	2.04	116.68	111.24
5	B	1165	EPE	O2S-S-C10	2.01	109.76	106.73
4	C	1322	NAG	C1-O5-C5	2.00	114.87	112.19

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	B	1164	NAG	C1
4	D	1164	NAG	C1
4	F	1164	NAG	C1

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1323	NAG	C1-C2-N2-C7
4	C	1323	NAG	C1-C2-N2-C7
4	E	1323	NAG	C1-C2-N2-C7
5	B	1165	EPE	C10-C9-N1-C6
5	B	1165	EPE	C9-C10-S-O2S
5	D	1165	EPE	C10-C9-N1-C2
5	D	1165	EPE	C9-C10-S-O2S
5	D	1165	EPE	C9-C10-S-O3S
5	F	1165	EPE	S-C10-C9-N1
4	E	1322	NAG	O5-C5-C6-O6
4	E	1322	NAG	C4-C5-C6-O6
5	F	1165	EPE	N4-C7-C8-O8
4	C	1322	NAG	C4-C5-C6-O6
4	C	1322	NAG	O5-C5-C6-O6
5	B	1165	EPE	N4-C7-C8-O8
5	B	1165	EPE	C9-C10-S-O3S
5	D	1165	EPE	S-C10-C9-N1
5	F	1165	EPE	C8-C7-N4-C3
5	D	1165	EPE	C8-C7-N4-C3
5	D	1165	EPE	N4-C7-C8-O8
5	B	1165	EPE	C9-C10-S-O1S
5	D	1165	EPE	C9-C10-S-O1S
5	B	1165	EPE	C8-C7-N4-C5
4	A	1322	NAG	C4-C5-C6-O6
4	A	1322	NAG	O5-C5-C6-O6
5	F	1165	EPE	C10-C9-N1-C6
5	D	1165	EPE	C10-C9-N1-C6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	1322	NAG	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	321/326 (98%)	-0.71	1 (0%) 90 88	25, 55, 82, 101	0
1	C	321/326 (98%)	-0.72	1 (0%) 90 88	23, 55, 81, 102	0
1	E	321/326 (98%)	-0.66	2 (0%) 85 83	21, 55, 80, 97	0
2	B	163/166 (98%)	-1.28	0 100 100	10, 32, 64, 82	0
2	D	163/166 (98%)	-1.29	0 100 100	11, 32, 62, 80	0
2	F	163/166 (98%)	-1.29	0 100 100	12, 35, 57, 75	0
All	All	1452/1476 (98%)	-0.90	4 (0%) 90 88	10, 47, 77, 102	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	194	PRO	3.4
1	E	123	SER	3.0
1	E	189	LYS	2.3
1	C	125	HIS	2.2

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

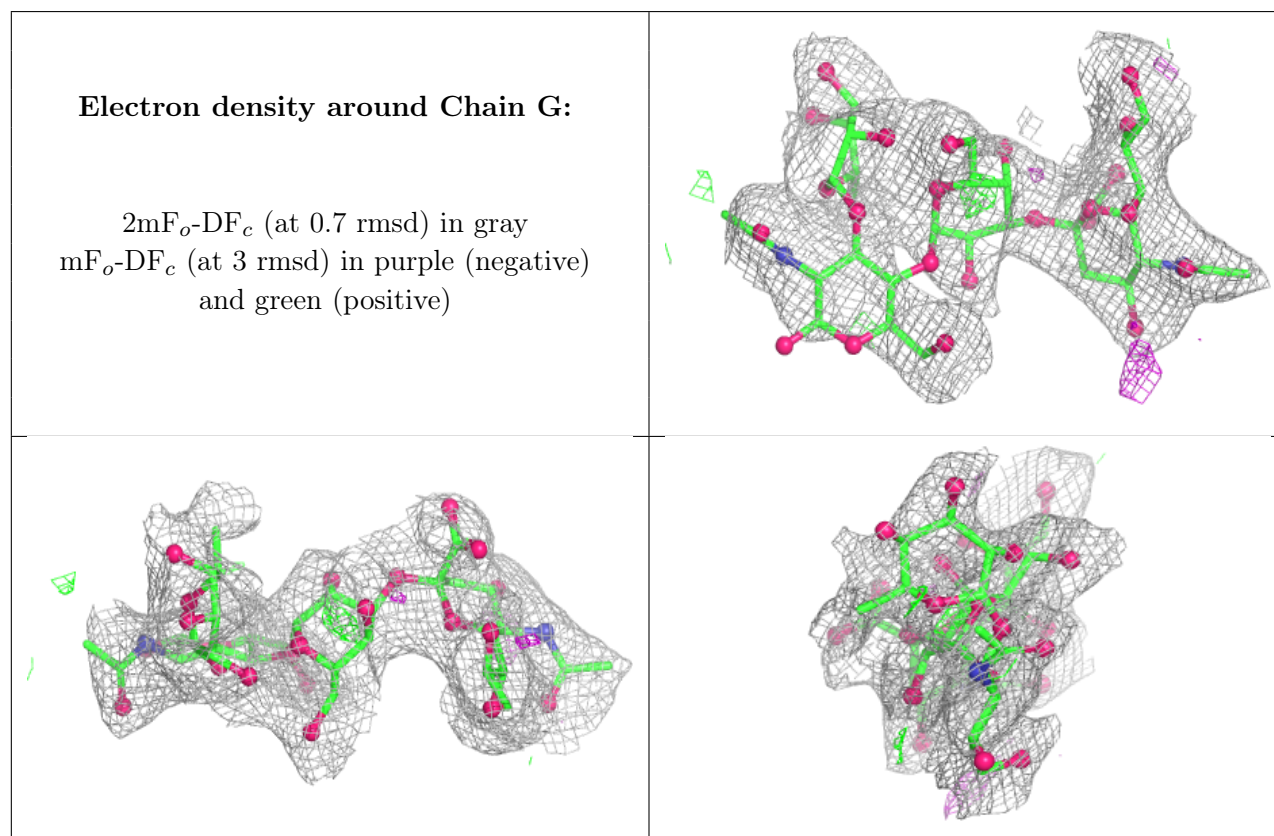
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

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

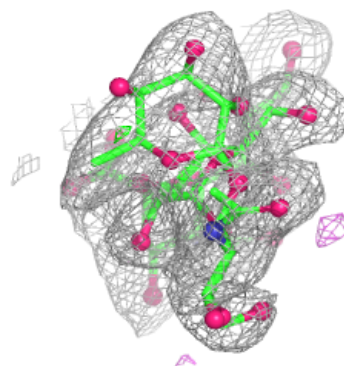
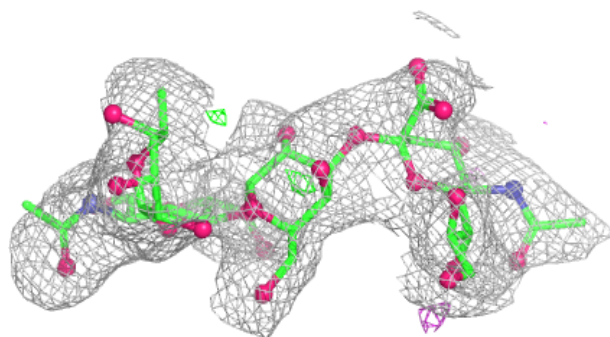
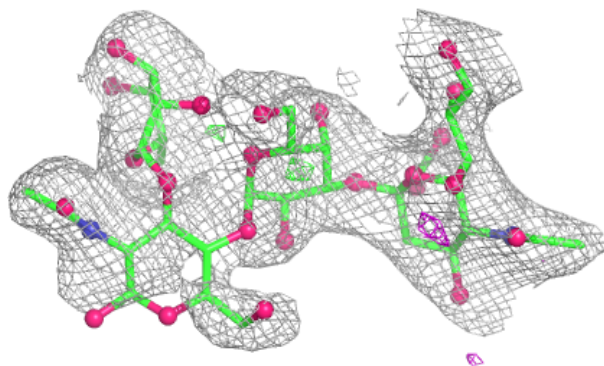
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	G	1	15/15	0.95	0.10	72,78,80,81	0
3	NAG	I	1	15/15	0.95	0.09	72,78,81,81	0
3	FUC	I	4	10/11	0.95	0.07	75,79,80,80	0
3	FUC	H	4	10/11	0.96	0.07	71,74,75,75	0
3	FUC	G	4	10/11	0.96	0.07	69,74,75,75	0
3	NAG	H	1	15/15	0.96	0.08	70,76,79,79	0
3	GLA	I	2	11/12	0.97	0.06	56,62,66,67	0
3	GLA	G	2	11/12	0.97	0.07	56,60,65,66	0
3	GLA	H	2	11/12	0.98	0.05	56,60,65,66	0
3	SIA	H	3	20/21	0.98	0.06	42,46,51,51	0
3	SIA	I	3	20/21	0.99	0.06	44,46,52,52	0
3	SIA	G	3	20/21	0.99	0.05	41,46,52,53	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.

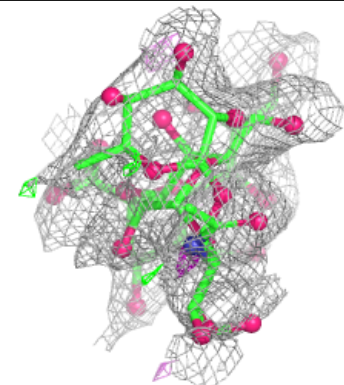
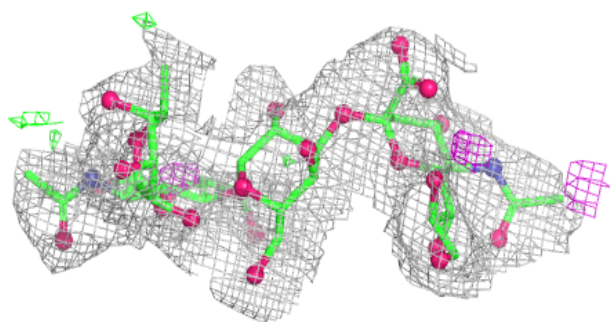
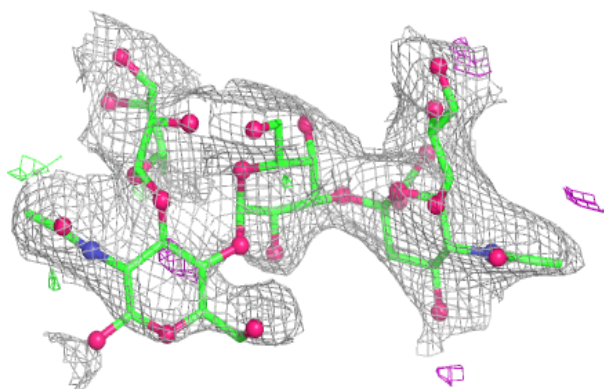


Electron density around Chain H:

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

**Electron density around Chain I:**

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



6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	NAG	B	1164	14/15	0.96	0.06	40,47,49,49	0
4	NAG	A	1322	14/15	0.97	0.05	30,36,37,38	0
4	NAG	D	1164	14/15	0.97	0.05	39,46,47,47	0
4	NAG	E	1322	14/15	0.97	0.04	28,34,35,36	0
4	NAG	F	1164	14/15	0.97	0.05	40,46,47,48	0
4	NAG	C	1323	14/15	0.98	0.04	31,35,39,41	0
4	NAG	E	1323	14/15	0.98	0.04	32,37,42,44	0
4	NAG	C	1322	14/15	0.98	0.04	27,32,34,35	0
4	NAG	A	1323	14/15	0.99	0.04	30,35,41,42	0
5	EPE	B	1165	15/15	0.99	0.04	50,51,61,63	0
5	EPE	D	1165	15/15	0.99	0.06	53,56,65,66	0
5	EPE	F	1165	15/15	0.99	0.06	52,55,64,66	0

6.5 Other polymers [i](#)

There are no such residues in this entry.