



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:12 AM UTC

PDB ID : 4LKG / pdb_00004lkg
Title : The structure of hemagglutinin from a avian-origin H7N9 influenza virus (A/Shanghai/1/2013) in complex with avian receptor analog 3'SLNLN
Authors : Shi, Y.; Zhang, W.; Wang, F.; Qi, J.; Song, H.; Wu, Y.; Gao, F.; Zhang, Y.; Fan, Z.; Gong, W.; Wang, D.; Shu, Y.; Wang, Y.; Yan, J.; Gao, G.F.
Deposited on : 2013-07-07
Resolution : 2.99 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

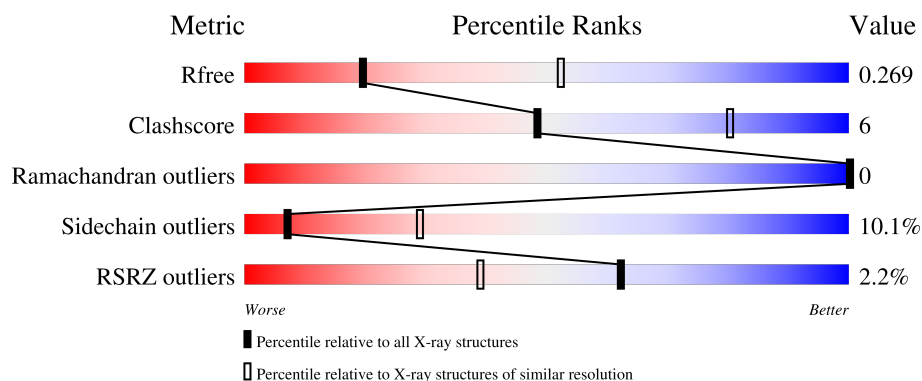
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

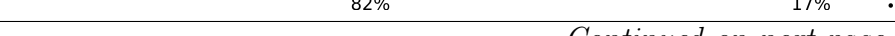
The reported resolution of this entry is 2.99 Å.

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.



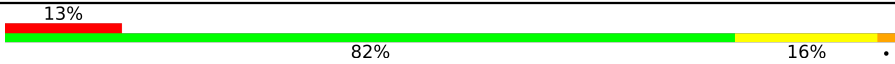
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	316	
1	C	316	
1	E	316	
2	B	170	
2	D	170	

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Mol	Chain	Length	Quality of chain
2	F	170	
3	G	3	
4	H	2	
4	I	2	

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	G	1	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 11531 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.

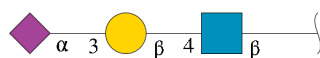
- Molecule 1 is a protein called hemagglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	316	Total	C	N	O	S	0	0	0
			2415	1497	435	468	15			
1	C	316	Total	C	N	O	S	0	0	0
			2415	1497	435	468	15			
1	E	316	Total	C	N	O	S	0	0	0
			2415	1497	435	468	15			

- Molecule 2 is a protein called hemagglutinin.

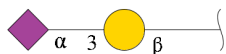
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	170	Total	C	N	O	S	0	0	0
			1378	851	240	280	7			
2	D	170	Total	C	N	O	S	0	0	0
			1378	851	240	280	7			
2	F	170	Total	C	N	O	S	0	0	0
			1378	851	240	280	7			

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



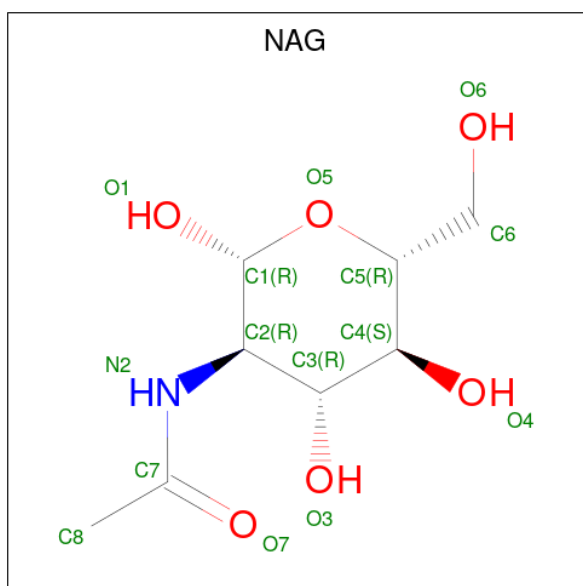
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	G	3	Total	C	N	O	0	0	0
			46	25	2	19			

- Molecule 4 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	H	2	Total	C	N	O	0	0	0
			32	17	1	14			
4	I	2	Total	C	N	O	0	0	0
			32	17	1	14			

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).

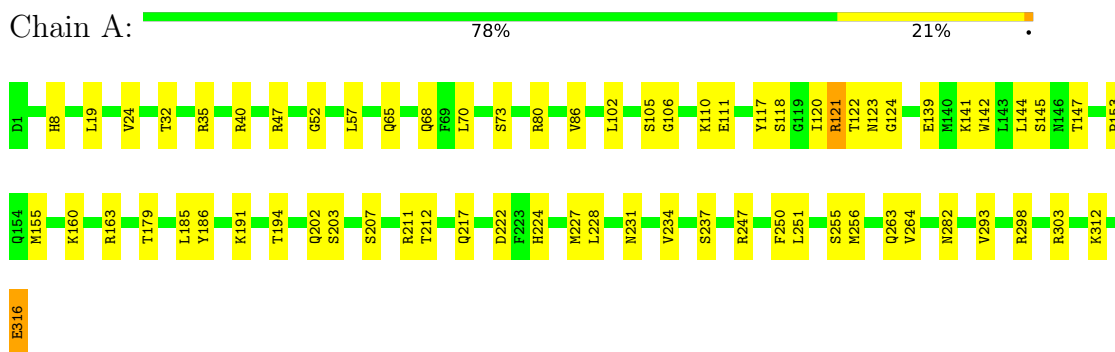


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	N	O	0	0
			14	8	1	5		
5	D	1	Total	C	N	O	0	0
			14	8	1	5		
5	F	1	Total	C	N	O	0	0
			14	8	1	5		

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.

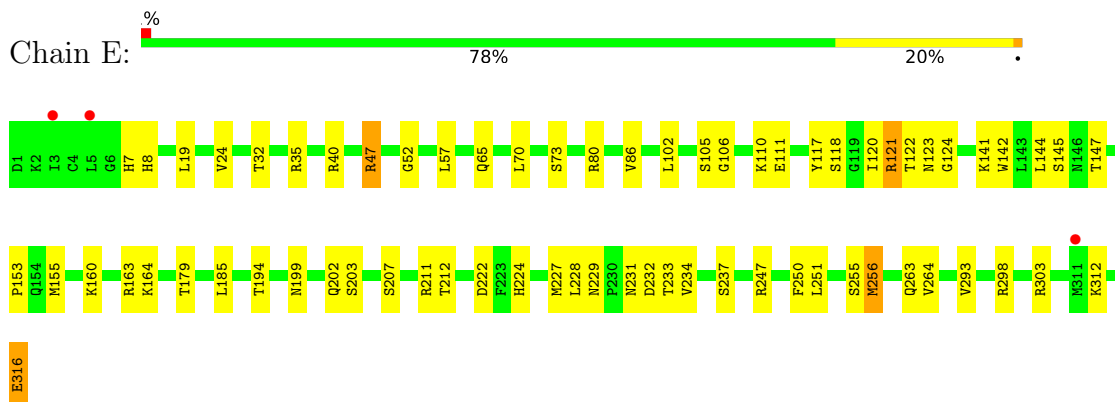
- Molecule 1: hemagglutinin



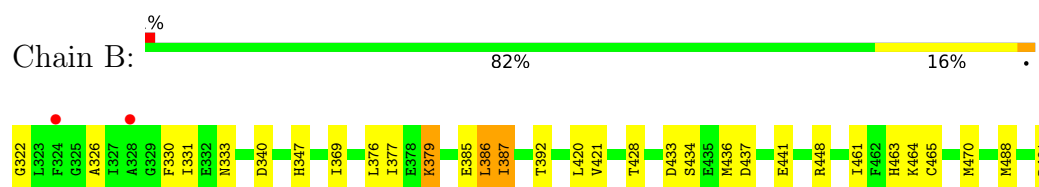
- Molecule 1: hemagglutinin



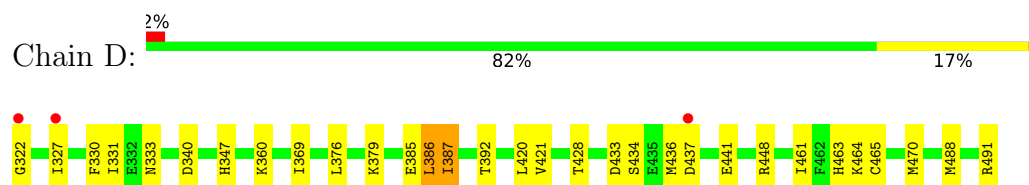
- Molecule 1: hemagglutinin



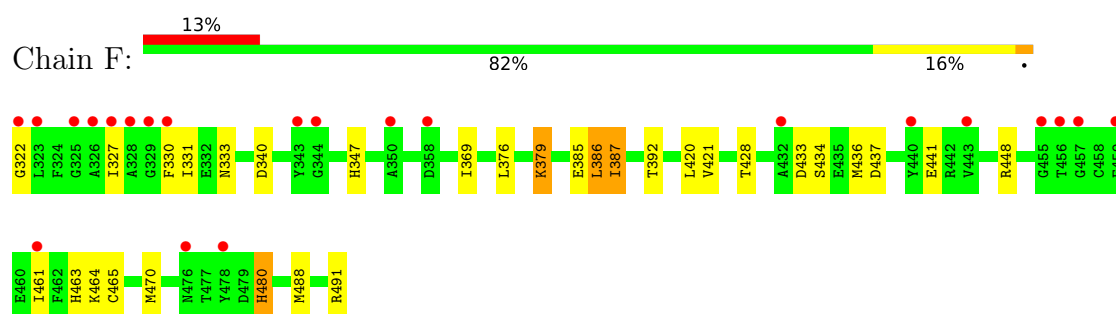
- Molecule 2: hemagglutinin



- Molecule 2: hemagglutinin



- Molecule 2: hemagglutinin



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



- Molecule 4: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose



- Molecule 4: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 3	Depositor
Cell constants a, b, c, α , β , γ	110.76Å 110.76Å 133.23Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	42.59 – 2.99 42.59 – 2.99	Depositor EDS
% Data completeness (in resolution range)	99.7 (42.59-2.99) 99.7 (42.59-2.99)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.21 (at 3.01Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.246 , 0.278 0.238 , 0.269	Depositor DCC
R_{free} test set	1838 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	72.7	Xtriage
Anisotropy	0.168	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 37.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.029 for -h,-k,l 0.109 for h,-h-k,-l 0.037 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	11531	wwPDB-VP
Average B, all atoms (Å ²)	100.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 15.49% 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: GAL, NAG, SIA

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	0.33	0/2460	0.75	0/3323
1	C	0.32	0/2460	0.74	0/3323
1	E	0.31	0/2460	0.75	0/3323
2	B	0.33	0/1402	0.70	1/1889 (0.1%)
2	D	0.32	0/1402	0.70	1/1889 (0.1%)
2	F	0.37	1/1402 (0.1%)	0.70	1/1889 (0.1%)
All	All	0.33	1/11586 (0.0%)	0.73	3/15636 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	480	HIS	CE1-NE2	-5.81	1.26	1.32

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	387	ILE	N-CA-C	-5.22	108.19	113.47
2	B	387	ILE	N-CA-C	-5.11	108.31	113.47
2	D	387	ILE	N-CA-C	-5.10	108.32	113.47

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	2415	0	2366	26	1
1	C	2415	0	2366	36	0
1	E	2415	0	2366	27	1
2	B	1378	0	1278	13	0
2	D	1378	0	1278	14	0
2	F	1378	0	1278	14	0
3	G	46	0	40	14	0
4	H	32	0	28	1	0
4	I	32	0	28	1	0
5	B	14	0	13	0	0
5	D	14	0	13	0	0
5	F	14	0	13	0	0
All	All	11531	0	11067	125	1

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 6.

All (125) 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:267:ASP:HB3	3:G:1:NAG:H82	1.32	1.12
1:C:267:ASP:HB3	3:G:1:NAG:C8	1.96	0.96
1:A:217:GLN:HE22	3:G:2:GAL:H3	1.33	0.93
1:C:40:ARG:HG2	1:C:263:GLN:HG3	1.53	0.88
1:C:267:ASP:HB3	3:G:1:NAG:C7	2.08	0.84
1:C:267:ASP:CB	3:G:1:NAG:H82	2.13	0.77
1:C:267:ASP:C	3:G:1:NAG:H82	2.15	0.71
2:F:448:ARG:HD3	2:F:480:HIS:CD2	2.26	0.70
1:E:144:LEU:HD11	1:E:185:LEU:HD22	1.75	0.69
1:A:121:ARG:NH1	1:A:145:SER:O	2.26	0.68
1:A:144:LEU:HD11	1:A:185:LEU:HD22	1.76	0.68
1:E:121:ARG:HB3	1:E:144:LEU:HB3	1.75	0.68
1:A:121:ARG:HB3	1:A:144:LEU:HB3	1.74	0.68
1:C:121:ARG:HB3	1:C:144:LEU:HB3	1.74	0.68
1:E:121:ARG:NH1	1:E:145:SER:O	2.26	0.68
1:C:144:LEU:HD11	1:C:185:LEU:HD22	1.76	0.66
1:C:121:ARG:NH1	1:C:145:SER:O	2.27	0.66
1:E:70:LEU:O	1:E:110:LYS:NZ	2.30	0.64
1:C:70:LEU:O	1:C:110:LYS:NZ	2.31	0.64
1:A:70:LEU:O	1:A:110:LYS:NZ	2.32	0.63
4:H:2:SIA:H6	4:H:2:SIA:O1B	1.99	0.63
1:E:316:GLU:HA	2:F:333:ASN:HD22	1.65	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:448:ARG:HD3	2:F:480:HIS:NE2	2.17	0.59
3:G:1:NAG:H62	3:G:2:GAL:O2	2.02	0.59
1:C:316:GLU:HA	2:D:333:ASN:HD22	1.68	0.59
1:A:316:GLU:HA	2:B:333:ASN:HD22	1.68	0.58
4:I:2:SIA:H6	4:I:2:SIA:O1B	2.03	0.58
1:A:222:ASP:HB3	1:A:224:HIS:CE1	2.40	0.57
1:E:222:ASP:HB3	1:E:224:HIS:CE1	2.40	0.57
1:C:222:ASP:HB3	1:C:224:HIS:CE1	2.40	0.56
1:A:40:ARG:HG2	1:A:263:GLN:HG3	1.87	0.56
1:E:40:ARG:HG2	1:E:263:GLN:HG3	1.87	0.56
1:C:111:GLU:OE2	1:C:163:ARG:NH1	2.40	0.55
1:C:160:LYS:HE3	1:C:231:ASN:HA	1.89	0.55
1:E:111:GLU:OE2	1:E:163:ARG:NH1	2.38	0.55
1:A:111:GLU:OE2	1:A:163:ARG:NH1	2.39	0.54
1:C:267:ASP:HB3	3:G:1:NAG:N2	2.21	0.54
1:C:267:ASP:O	3:G:1:NAG:H82	2.09	0.53
1:A:160:LYS:HE3	1:A:231:ASN:HA	1.90	0.53
1:E:160:LYS:HE3	1:E:231:ASN:HA	1.90	0.53
2:D:437:ASP:O	2:D:441:GLU:HG2	2.10	0.52
2:F:437:ASP:O	2:F:441:GLU:HG2	2.10	0.52
2:B:437:ASP:O	2:B:441:GLU:HG2	2.09	0.51
1:C:207:SER:O	1:C:211:ARG:NH2	2.42	0.51
3:G:3:SIA:O1B	3:G:3:SIA:H6	2.09	0.51
1:E:293:VAL:HG11	2:F:386:LEU:HD13	1.93	0.50
2:F:461:ILE:HG22	2:F:463:HIS:H	1.77	0.49
1:C:293:VAL:HG11	2:D:386:LEU:HD13	1.95	0.49
2:F:376:LEU:HD22	2:F:420:LEU:HD21	1.95	0.48
1:A:217:GLN:HE22	3:G:2:GAL:C3	2.14	0.48
1:A:207:SER:O	1:A:211:ARG:NH2	2.45	0.48
2:D:376:LEU:HD22	2:D:420:LEU:HD21	1.96	0.48
1:A:153:PRO:O	1:A:155:MET:HG3	2.14	0.48
2:B:376:LEU:HD22	2:B:420:LEU:HD21	1.95	0.48
1:C:153:PRO:O	1:C:155:MET:HG3	2.14	0.48
1:A:47:ARG:NH1	1:A:73:SER:HB3	2.29	0.47
1:C:163:ARG:HD3	1:C:250:PHE:CE2	2.50	0.47
2:D:461:ILE:HG22	2:D:463:HIS:H	1.78	0.47
2:B:461:ILE:HG22	2:B:463:HIS:H	1.78	0.47
1:E:124:GLY:HA3	1:E:142:TRP:HB3	1.97	0.47
1:E:207:SER:O	1:E:211:ARG:NH2	2.44	0.47
1:E:153:PRO:O	1:E:155:MET:HG3	2.15	0.46
1:E:35:ARG:NH1	1:E:303:ARG:O	2.47	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:163:ARG:HD3	1:A:250:PHE:CE2	2.50	0.46
1:A:35:ARG:NH1	1:A:303:ARG:O	2.48	0.46
1:E:106:GLY:HA2	1:E:255:SER:HB3	1.98	0.46
2:D:327:ILE:N	2:D:433:ASP:OD1	2.42	0.46
1:E:163:ARG:HD3	1:E:250:PHE:CE2	2.50	0.46
1:E:47:ARG:NH1	1:E:73:SER:HB3	2.30	0.45
1:A:106:GLY:HA2	1:A:255:SER:HB3	1.98	0.45
1:C:35:ARG:NH1	1:C:303:ARG:O	2.49	0.45
1:C:106:GLY:HA2	1:C:255:SER:HB3	1.99	0.45
2:F:327:ILE:N	2:F:433:ASP:OD1	2.42	0.45
1:E:7:HIS:CD2	2:F:327:ILE:HG12	2.51	0.45
2:D:330:PHE:CD1	2:D:331:ILE:HG13	2.52	0.45
2:D:448:ARG:HE	2:D:448:ARG:HB3	1.61	0.45
1:E:163:ARG:HD3	1:E:250:PHE:CZ	2.52	0.45
1:C:47:ARG:NH1	1:C:73:SER:HB3	2.32	0.45
1:A:293:VAL:HG11	2:B:386:LEU:HD13	1.97	0.45
2:F:330:PHE:CD1	2:F:331:ILE:HG13	2.52	0.45
2:F:347:HIS:HB2	2:F:470:MET:HE3	1.99	0.45
1:C:117:TYR:HB3	1:C:120:ILE:HD11	1.99	0.44
1:C:267:ASP:CB	3:G:1:NAG:HN2	2.30	0.44
1:A:124:GLY:HA3	1:A:142:TRP:HB3	1.99	0.44
1:C:124:GLY:HA3	1:C:142:TRP:HB3	1.98	0.44
2:B:322:GLY:HA3	2:B:433:ASP:OD2	2.18	0.44
2:D:322:GLY:HA3	2:D:433:ASP:OD2	2.18	0.43
1:A:105:SER:HB2	1:A:251:LEU:HD22	2.00	0.43
1:A:117:TYR:HB3	1:A:120:ILE:HD11	2.00	0.43
1:C:256:MET:HB3	1:C:256:MET:HE3	1.78	0.43
2:F:379:LYS:HE3	2:F:379:LYS:HB2	1.82	0.43
1:A:163:ARG:HD3	1:A:250:PHE:CZ	2.53	0.43
2:B:330:PHE:CD1	2:B:331:ILE:HG13	2.52	0.43
1:E:117:TYR:HB3	1:E:120:ILE:HD11	2.00	0.43
2:D:347:HIS:HB2	2:D:470:MET:HE3	2.00	0.43
1:C:105:SER:HB2	1:C:251:LEU:HD22	2.01	0.43
1:C:267:ASP:HB3	3:G:1:NAG:HN2	1.84	0.43
2:D:369:ILE:HD11	2:D:428:THR:HG23	2.01	0.43
2:B:448:ARG:HE	2:B:448:ARG:HB3	1.61	0.42
1:E:52:GLY:HA2	1:E:80:ARG:HG3	2.00	0.42
2:F:322:GLY:HA3	2:F:433:ASP:OD2	2.19	0.42
1:C:52:GLY:HA2	1:C:80:ARG:HG3	2.01	0.42
1:C:163:ARG:HD3	1:C:250:PHE:CZ	2.53	0.42
1:E:199:ASN:OD1	1:E:229:ASN:ND2	2.51	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:186:TYR:O	1:A:191:LYS:NZ	2.44	0.42
1:A:282:ASN:HB3	2:B:377:ILE:HG23	2.02	0.42
1:E:105:SER:HB2	1:E:251:LEU:HD22	2.02	0.42
2:B:347:HIS:HB2	2:B:470:MET:HE3	2.01	0.41
1:C:199:ASN:OD1	1:C:229:ASN:ND2	2.49	0.41
2:F:369:ILE:HD11	2:F:428:THR:HG23	2.02	0.41
1:C:57:LEU:HD23	1:C:102:LEU:HD12	2.03	0.41
2:B:369:ILE:HD11	2:B:428:THR:HG23	2.01	0.41
1:C:267:ASP:CA	3:G:1:NAG:H82	2.51	0.41
1:A:52:GLY:HA2	1:A:80:ARG:HG3	2.02	0.41
1:C:35:ARG:HD3	1:C:303:ARG:HG2	2.03	0.41
1:E:232:ASP:OD1	1:E:233:THR:N	2.50	0.41
1:E:256:MET:HB3	1:E:256:MET:HE3	1.78	0.41
2:B:379:LYS:HE3	2:B:379:LYS:HB2	1.82	0.41
1:E:35:ARG:HD3	1:E:303:ARG:HG2	2.02	0.41
1:E:57:LEU:HD23	1:E:102:LEU:HD12	2.03	0.41
1:A:57:LEU:HD23	1:A:102:LEU:HD12	2.03	0.41
2:D:360:LYS:HB2	2:D:360:LYS:HE3	1.86	0.41
2:B:326:ALA:HB3	2:B:433:ASP:OD1	2.22	0.40
2:D:331:ILE:HG13	2:D:331:ILE:H	1.74	0.40
1:C:7:HIS:CD2	2:D:327:ILE:HG12	2.57	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:GLU:OE1	1:E:164:LYS:NZ[2_545]	2.11	0.09

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	314/316 (99%)	292 (93%)	22 (7%)	0	100	100
1	C	314/316 (99%)	291 (93%)	23 (7%)	0	100	100
1	E	314/316 (99%)	291 (93%)	23 (7%)	0	100	100
2	B	168/170 (99%)	159 (95%)	9 (5%)	0	100	100
2	D	168/170 (99%)	159 (95%)	9 (5%)	0	100	100
2	F	168/170 (99%)	159 (95%)	9 (5%)	0	100	100
All	All	1446/1458 (99%)	1351 (93%)	95 (7%)	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	265/265 (100%)	237 (89%)	28 (11%)	6	27
1	C	265/265 (100%)	236 (89%)	29 (11%)	6	26
1	E	265/265 (100%)	237 (89%)	28 (11%)	6	27
2	B	145/145 (100%)	132 (91%)	13 (9%)	9	34
2	D	145/145 (100%)	132 (91%)	13 (9%)	9	34
2	F	145/145 (100%)	132 (91%)	13 (9%)	9	34
All	All	1230/1230 (100%)	1106 (90%)	124 (10%)	7	29

All (124) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	HIS
1	A	19	LEU
1	A	24	VAL
1	A	32	THR
1	A	65	GLN
1	A	68	GLN
1	A	86	VAL

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Mol	Chain	Res	Type
1	A	118	SER
1	A	121	ARG
1	A	122	THR
1	A	123	ASN
1	A	141	LYS
1	A	147	THR
1	A	179	THR
1	A	194	THR
1	A	202	GLN
1	A	203	SER
1	A	212	THR
1	A	227	MET
1	A	228	LEU
1	A	234	VAL
1	A	237	SER
1	A	247	ARG
1	A	256	MET
1	A	264	VAL
1	A	298	ARG
1	A	312	LYS
1	A	316	GLU
2	B	340	ASP
2	B	379	LYS
2	B	385	GLU
2	B	386	LEU
2	B	387	ILE
2	B	392	THR
2	B	421	VAL
2	B	434	SER
2	B	436	MET
2	B	464	LYS
2	B	465	CYS
2	B	488	MET
2	B	491	ARG
1	C	8	HIS
1	C	19	LEU
1	C	24	VAL
1	C	32	THR
1	C	47	ARG
1	C	65	GLN
1	C	68	GLN
1	C	86	VAL

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Mol	Chain	Res	Type
1	C	118	SER
1	C	121	ARG
1	C	122	THR
1	C	123	ASN
1	C	141	LYS
1	C	147	THR
1	C	179	THR
1	C	194	THR
1	C	202	GLN
1	C	203	SER
1	C	212	THR
1	C	227	MET
1	C	228	LEU
1	C	234	VAL
1	C	237	SER
1	C	247	ARG
1	C	256	MET
1	C	264	VAL
1	C	298	ARG
1	C	312	LYS
1	C	316	GLU
2	D	340	ASP
2	D	379	LYS
2	D	385	GLU
2	D	386	LEU
2	D	387	ILE
2	D	392	THR
2	D	421	VAL
2	D	434	SER
2	D	436	MET
2	D	464	LYS
2	D	465	CYS
2	D	488	MET
2	D	491	ARG
1	E	8	HIS
1	E	19	LEU
1	E	24	VAL
1	E	32	THR
1	E	47	ARG
1	E	65	GLN
1	E	86	VAL
1	E	118	SER

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Mol	Chain	Res	Type
1	E	121	ARG
1	E	122	THR
1	E	123	ASN
1	E	141	LYS
1	E	147	THR
1	E	179	THR
1	E	194	THR
1	E	202	GLN
1	E	203	SER
1	E	212	THR
1	E	227	MET
1	E	228	LEU
1	E	234	VAL
1	E	237	SER
1	E	247	ARG
1	E	256	MET
1	E	264	VAL
1	E	298	ARG
1	E	312	LYS
1	E	316	GLU
2	F	340	ASP
2	F	379	LYS
2	F	385	GLU
2	F	386	LEU
2	F	387	ILE
2	F	392	THR
2	F	421	VAL
2	F	434	SER
2	F	436	MET
2	F	464	LYS
2	F	465	CYS
2	F	488	MET
2	F	491	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	217	GLN
1	A	224	HIS
2	B	383	GLN
2	B	480	HIS
2	B	490	ASN

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Mol	Chain	Res	Type
1	C	224	HIS
2	D	480	HIS
1	E	224	HIS
2	F	333	ASN

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 ⓘ

7 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	0.48	0	21,21,21	1.32	3 (14%)
3	GAL	G	2	3	11,11,12	0.64	0	15,15,17	0.90	0
3	SIA	G	3	3	20,20,21	0.55	0	21,28,31	0.81	1 (4%)
4	GAL	H	1	4	12,12,12	0.50	0	17,17,17	0.54	0
4	SIA	H	2	4	20,20,21	0.56	0	21,28,31	0.81	1 (4%)
4	GAL	I	1	4	12,12,12	0.46	0	17,17,17	0.52	0
4	SIA	I	2	4	20,20,21	0.56	0	21,28,31	0.82	2 (9%)

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	GAL	G	2	3	-	0/2/19/22	0/1/1/1
3	SIA	G	3	3	-	4/18/34/38	0/1/1/1
4	GAL	H	1	4	-	2/2/22/22	0/1/1/1
4	SIA	H	2	4	-	0/18/34/38	0/1/1/1
4	GAL	I	1	4	-	2/2/22/22	0/1/1/1
4	SIA	I	2	4	-	3/18/34/38	0/1/1/1

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	1	NAG	C4-C3-C2	4.02	116.26	110.40
3	G	1	NAG	C1-C2-C3	2.32	113.71	110.54
3	G	3	SIA	O1B-C1-C2	2.22	118.50	112.71
4	I	2	SIA	O1B-C1-C2	2.22	118.49	112.71
4	H	2	SIA	O1B-C1-C2	2.20	118.44	112.71
3	G	1	NAG	C3-C2-N2	-2.02	106.90	110.62
4	I	2	SIA	O1A-C1-C2	-2.01	118.51	122.85

There are no chirality outliers.

All (13) torsion outliers are listed below:

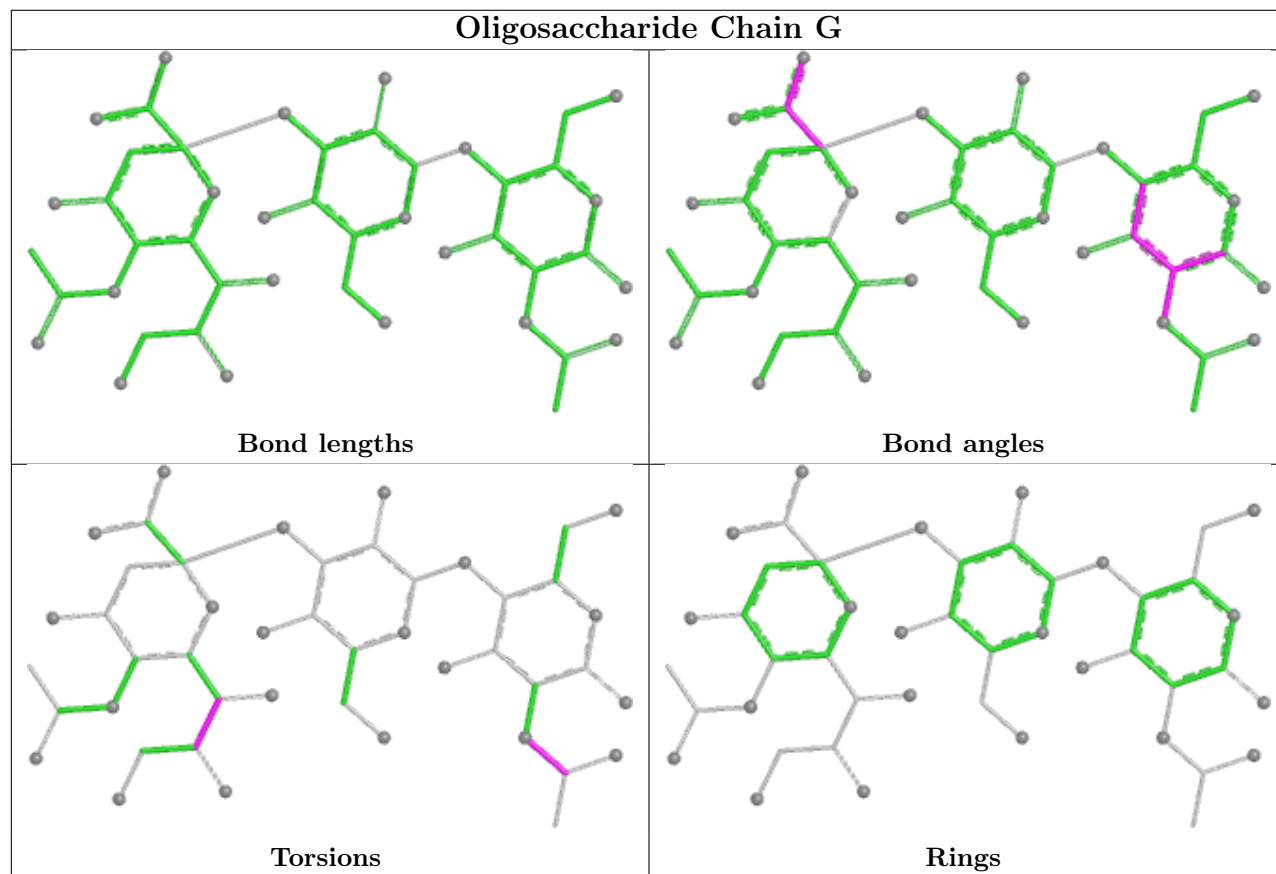
Mol	Chain	Res	Type	Atoms
4	H	1	GAL	O5-C5-C6-O6
3	G	1	NAG	C8-C7-N2-C2
4	H	1	GAL	C4-C5-C6-O6
3	G	1	NAG	O7-C7-N2-C2
3	G	3	SIA	O7-C7-C8-C9
3	G	3	SIA	C6-C7-C8-O8
4	I	1	GAL	O5-C5-C6-O6
3	G	3	SIA	C6-C7-C8-C9
3	G	3	SIA	O7-C7-C8-O8
4	I	2	SIA	C6-C7-C8-O8
4	I	1	GAL	C4-C5-C6-O6
4	I	2	SIA	O7-C7-C8-C9
4	I	2	SIA	O7-C7-C8-O8

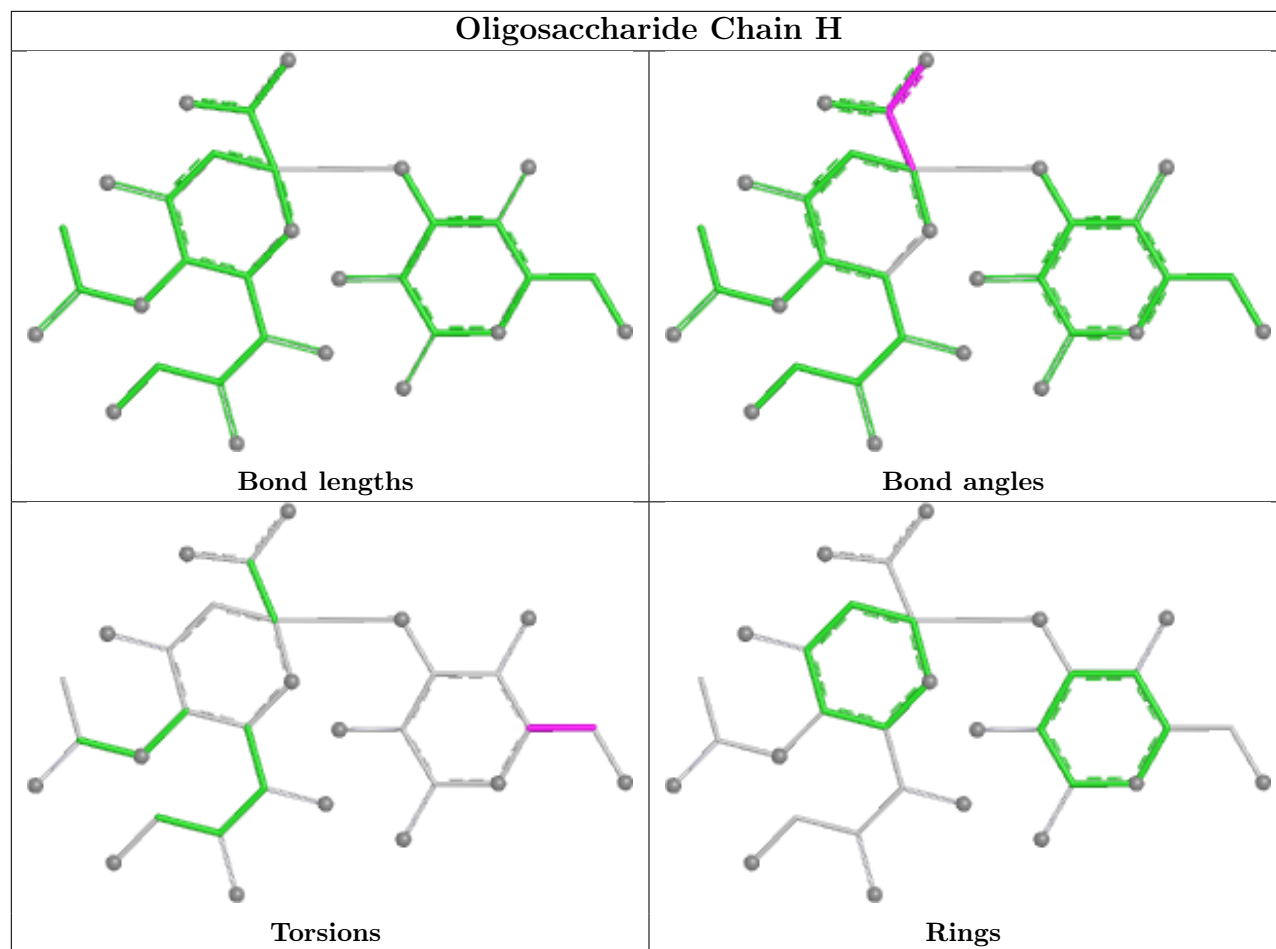
There are no ring outliers.

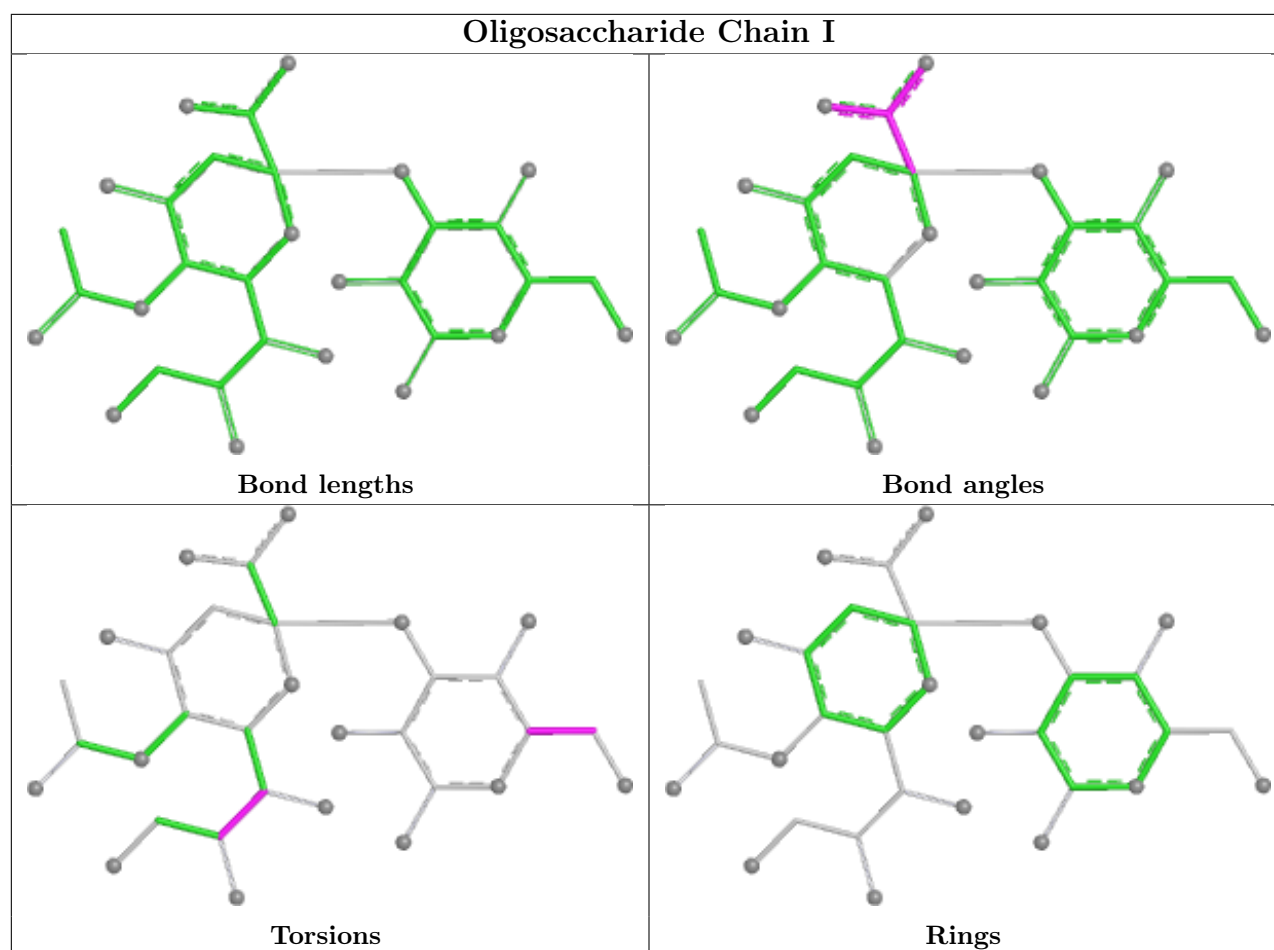
5 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	H	2	SIA	1	0
4	I	2	SIA	1	0
3	G	1	NAG	11	0
3	G	2	GAL	3	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 ⓘ

3 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	NAG	B	601	2	14,14,15	0.49	0	17,19,21	0.93	1 (5%)
5	NAG	D	601	2	14,14,15	0.52	0	17,19,21	0.62	0
5	NAG	F	601	2	14,14,15	0.47	0	17,19,21	0.86	1 (5%)

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	NAG	B	601	2	-	0/6/23/26	0/1/1/1
5	NAG	D	601	2	-	0/6/23/26	0/1/1/1
5	NAG	F	601	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	601	NAG	C1-O5-C5	3.22	116.50	112.19
5	F	601	NAG	C1-O5-C5	2.54	115.59	112.19

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 [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	316/316 (100%)	-0.82	0 100 100	43, 73, 111, 140	0
1	C	316/316 (100%)	-0.53	2 (0%) 85 69	46, 91, 145, 232	0
1	E	316/316 (100%)	-0.38	3 (0%) 81 61	59, 101, 169, 236	0
2	B	170/170 (100%)	-0.71	2 (1%) 76 55	39, 80, 148, 277	0
2	D	170/170 (100%)	-0.45	3 (1%) 67 44	41, 99, 156, 229	0
2	F	170/170 (100%)	0.51	22 (12%) 7 5	55, 158, 221, 306	0
All	All	1458/1458 (100%)	-0.45	32 (2%) 62 39	39, 91, 179, 306	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	327	ILE	8.4
2	F	326	ALA	6.9
2	F	330	PHE	5.4
2	F	455	GLY	4.0
2	F	457	GLY	3.9
2	F	322	GLY	3.7
2	F	329	GLY	3.6
2	F	328	ALA	3.2
2	F	350	ALA	3.2
1	E	311	MET	3.2
2	F	459	PHE	3.2
2	B	324	PHE	3.1
1	C	123	ASN	2.9
1	E	3	ILE	2.9
2	F	358	ASP	2.8
2	F	432	ALA	2.8
2	F	344	GLY	2.8
2	F	343	TYR	2.7
1	E	5	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
2	F	443	VAL	2.6
2	F	461	ILE	2.5
2	D	437	ASP	2.5
2	F	440	TYR	2.3
1	C	190	ASN	2.2
2	F	456	THR	2.2
2	F	325	GLY	2.2
2	F	323	LEU	2.1
2	D	327	ILE	2.1
2	F	476	ASN	2.1
2	D	322	GLY	2.1
2	F	478	TYR	2.1
2	B	328	ALA	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](#)

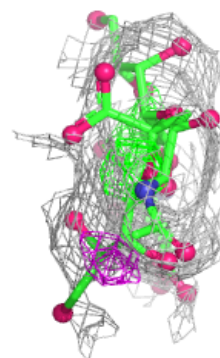
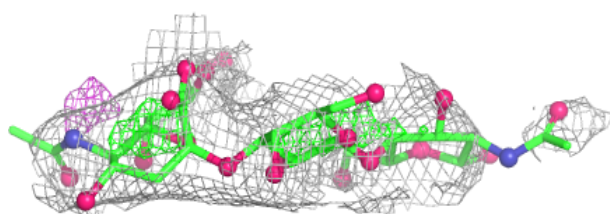
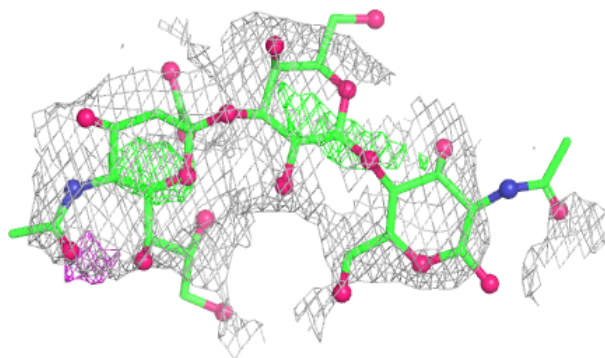
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GAL	H	1	12/12	0.93	0.08	125,148,178,196	0
4	GAL	I	1	12/12	0.93	0.07	123,142,151,156	0
4	SIA	H	2	20/21	0.96	0.09	92,102,113,115	0
3	NAG	G	1	15/15	0.96	0.09	133,154,172,175	0
4	SIA	I	2	20/21	0.96	0.10	130,136,145,147	0
3	SIA	G	3	20/21	0.97	0.10	90,106,115,115	0
3	GAL	G	2	11/12	0.98	0.09	84,89,95,106	0

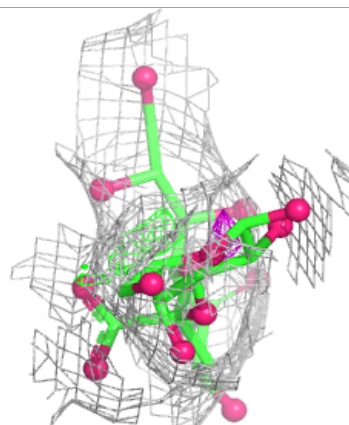
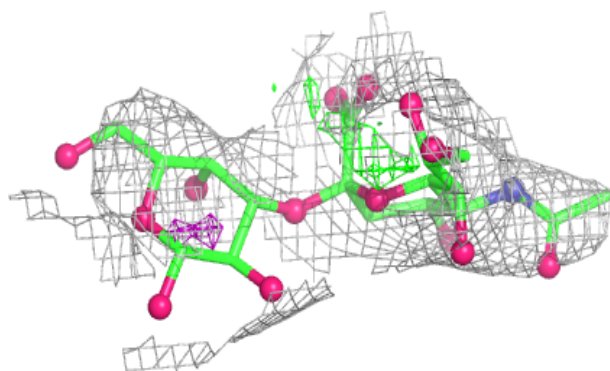
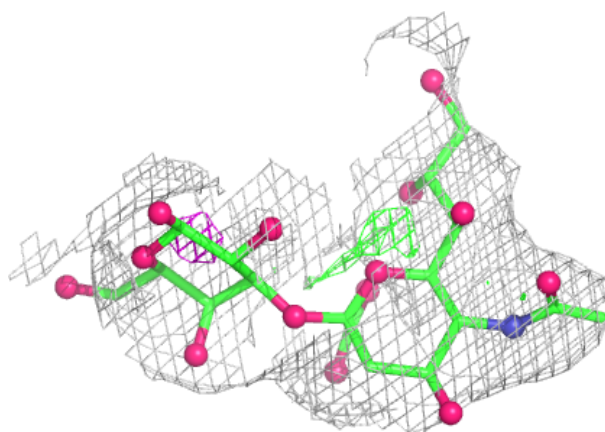
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

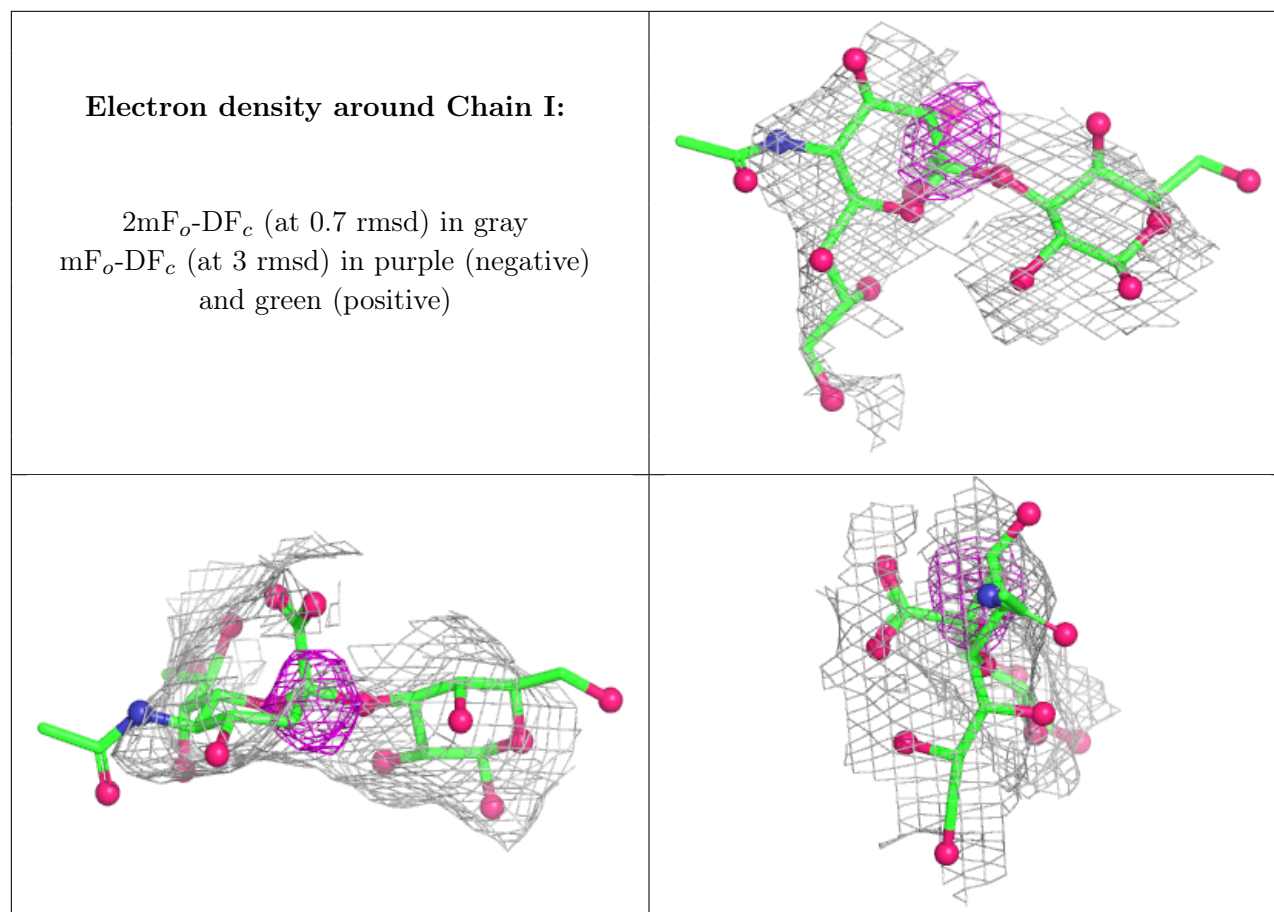
Electron density around Chain G:

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

**Electron density around Chain 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](#)

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
5	NAG	D	601	14/15	0.96	0.08	83,90,107,114	0
5	NAG	F	601	14/15	0.97	0.10	96,120,140,140	0
5	NAG	B	601	14/15	0.98	0.06	70,84,88,89	0

6.5 Other polymers [i](#)

There are no such residues in this entry.