



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 11:41 AM UTC

PDB ID : 9DIQ / pdb_00009diq
Title : Crystal structure of Apo-H5 hemagglutinin from the influenza virus A/Texas/37/2024 (H5N1)
Authors : Lin, T.H.; Zhu, Y.; Wilson, I.A.
Deposited on : 2024-09-05
Resolution : 2.69 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

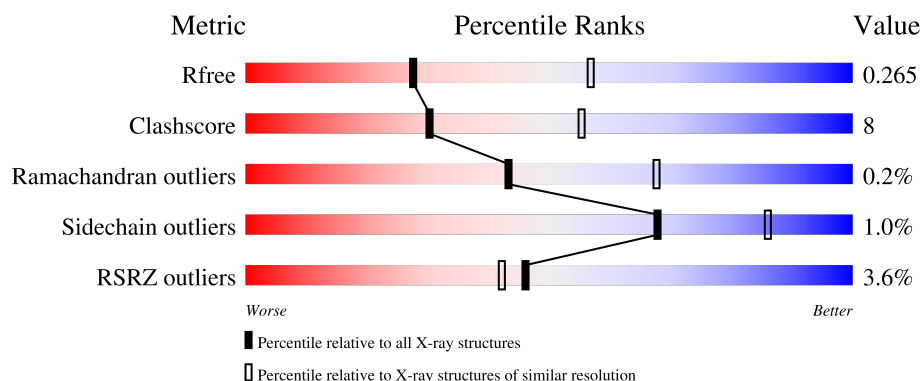
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.69 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	325	3% 76% 22% ..
1	C	325	2% 85% 14% .
1	E	325	% 78% 21% ..
1	G	325	3% 82% 17% .
1	I	325	% 83% 16% .

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Mol	Chain	Length	Quality of chain
1	K	325	
2	B	176	
2	D	176	
2	F	176	
2	H	176	
2	J	176	
2	L	176	
3	M	3	
3	O	3	
3	Q	3	
4	N	2	
4	P	2	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 23750 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 HA1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	I	322	Total	C	N	O	S	0	0	0
			2546	1611	444	476	15			
1	A	323	Total	C	N	O	S	0	0	0
			2553	1616	445	477	15			
1	C	322	Total	C	N	O	S	0	0	0
			2546	1611	444	476	15			
1	E	323	Total	C	N	O	S	0	0	0
			2553	1616	445	477	15			
1	G	322	Total	C	N	O	S	0	0	0
			2546	1611	444	476	15			
1	K	322	Total	C	N	O	S	0	0	0
			2546	1611	444	476	15			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I	7	ALA	-	expression tag	UNP A0A6B7HPT9
I	8	ASP	-	expression tag	UNP A0A6B7HPT9
I	9	PRO	-	expression tag	UNP A0A6B7HPT9
I	10	GLY	-	expression tag	UNP A0A6B7HPT9
I	111	MET	LEU	conflict	UNP A0A6B7HPT9
I	122	GLN	LEU	conflict	UNP A0A6B7HPT9
I	199	ILE	THR	conflict	UNP A0A6B7HPT9
I	214	ALA	VAL	conflict	UNP A0A6B7HPT9
A	7	ALA	-	expression tag	UNP A0A6B7HPT9
A	8	ASP	-	expression tag	UNP A0A6B7HPT9
A	9	PRO	-	expression tag	UNP A0A6B7HPT9
A	10	GLY	-	expression tag	UNP A0A6B7HPT9
A	111	MET	LEU	conflict	UNP A0A6B7HPT9
A	122	GLN	LEU	conflict	UNP A0A6B7HPT9
A	199	ILE	THR	conflict	UNP A0A6B7HPT9
A	214	ALA	VAL	conflict	UNP A0A6B7HPT9
C	7	ALA	-	expression tag	UNP A0A6B7HPT9

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Chain	Residue	Modelled	Actual	Comment	Reference
C	8	ASP	-	expression tag	UNP A0A6B7HPT9
C	9	PRO	-	expression tag	UNP A0A6B7HPT9
C	10	GLY	-	expression tag	UNP A0A6B7HPT9
C	111	MET	LEU	conflict	UNP A0A6B7HPT9
C	122	GLN	LEU	conflict	UNP A0A6B7HPT9
C	199	ILE	THR	conflict	UNP A0A6B7HPT9
C	214	ALA	VAL	conflict	UNP A0A6B7HPT9
E	7	ALA	-	expression tag	UNP A0A6B7HPT9
E	8	ASP	-	expression tag	UNP A0A6B7HPT9
E	9	PRO	-	expression tag	UNP A0A6B7HPT9
E	10	GLY	-	expression tag	UNP A0A6B7HPT9
E	111	MET	LEU	conflict	UNP A0A6B7HPT9
E	122	GLN	LEU	conflict	UNP A0A6B7HPT9
E	199	ILE	THR	conflict	UNP A0A6B7HPT9
E	214	ALA	VAL	conflict	UNP A0A6B7HPT9
G	7	ALA	-	expression tag	UNP A0A6B7HPT9
G	8	ASP	-	expression tag	UNP A0A6B7HPT9
G	9	PRO	-	expression tag	UNP A0A6B7HPT9
G	10	GLY	-	expression tag	UNP A0A6B7HPT9
G	111	MET	LEU	conflict	UNP A0A6B7HPT9
G	122	GLN	LEU	conflict	UNP A0A6B7HPT9
G	199	ILE	THR	conflict	UNP A0A6B7HPT9
G	214	ALA	VAL	conflict	UNP A0A6B7HPT9
K	7	ALA	-	expression tag	UNP A0A6B7HPT9
K	8	ASP	-	expression tag	UNP A0A6B7HPT9
K	9	PRO	-	expression tag	UNP A0A6B7HPT9
K	10	GLY	-	expression tag	UNP A0A6B7HPT9
K	111	MET	LEU	conflict	UNP A0A6B7HPT9
K	122	GLN	LEU	conflict	UNP A0A6B7HPT9
K	199	ILE	THR	conflict	UNP A0A6B7HPT9
K	214	ALA	VAL	conflict	UNP A0A6B7HPT9

- Molecule 2 is a protein called Hemagglutinin HA2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	J	167	Total	C	N	O	S	0	0	0
			1367	847	238	274	8			
2	B	170	Total	C	N	O	S	0	0	0
			1385	859	241	277	8			
2	D	166	Total	C	N	O	S	0	0	0
			1363	845	237	273	8			
2	F	170	Total	C	N	O	S	0	0	0
			1383	858	241	276	8			

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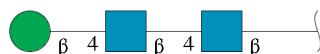
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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	167	Total	C	N	O	S	0	0	0
			1367	847	238	274	8			
2	L	167	Total	C	N	O	S	0	0	0
			1367	847	238	274	8			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
J	175	SER	-	expression tag	UNP A0A6B7HQ27
J	176	GLY	-	expression tag	UNP A0A6B7HQ27
B	175	SER	-	expression tag	UNP A0A6B7HQ27
B	176	GLY	-	expression tag	UNP A0A6B7HQ27
D	175	SER	-	expression tag	UNP A0A6B7HQ27
D	176	GLY	-	expression tag	UNP A0A6B7HQ27
F	175	SER	-	expression tag	UNP A0A6B7HQ27
F	176	GLY	-	expression tag	UNP A0A6B7HQ27
H	175	SER	-	expression tag	UNP A0A6B7HQ27
H	176	GLY	-	expression tag	UNP A0A6B7HQ27
L	175	SER	-	expression tag	UNP A0A6B7HQ27
L	176	GLY	-	expression tag	UNP A0A6B7HQ27

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



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	M	3	Total	C	N	O		0	0	0
			39	22	2	15				
3	O	3	Total	C	N	O		0	0	0
			39	22	2	15				
3	Q	3	Total	C	N	O		0	0	0
			39	22	2	15				

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



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

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	I	5	Total	O	0	0
			5	5		
6	A	6	Total	O	0	0
			6	6		
6	B	1	Total	O	0	0
			1	1		
6	C	7	Total	O	0	0
			7	7		
6	D	1	Total	O	0	0
			1	1		
6	E	7	Total	O	0	0
			7	7		

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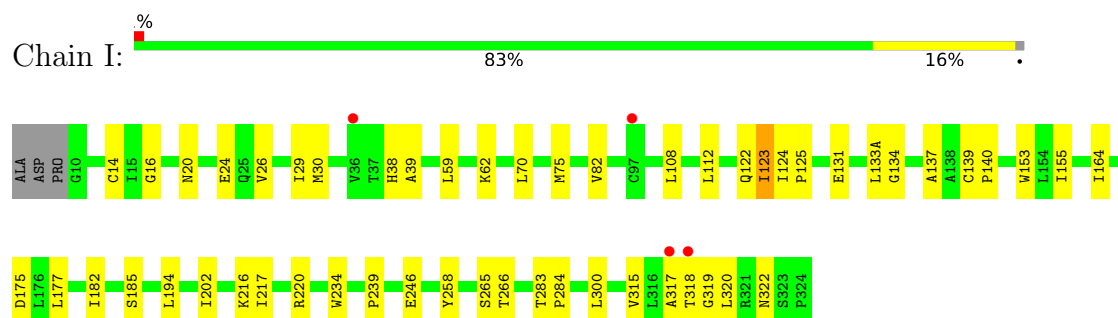
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	F	2	Total 2	O 2	0	0
6	G	4	Total 4	O 4	0	0
6	H	1	Total 1	O 1	0	0
6	K	5	Total 5	O 5	0	0
6	L	2	Total 2	O 2	0	0

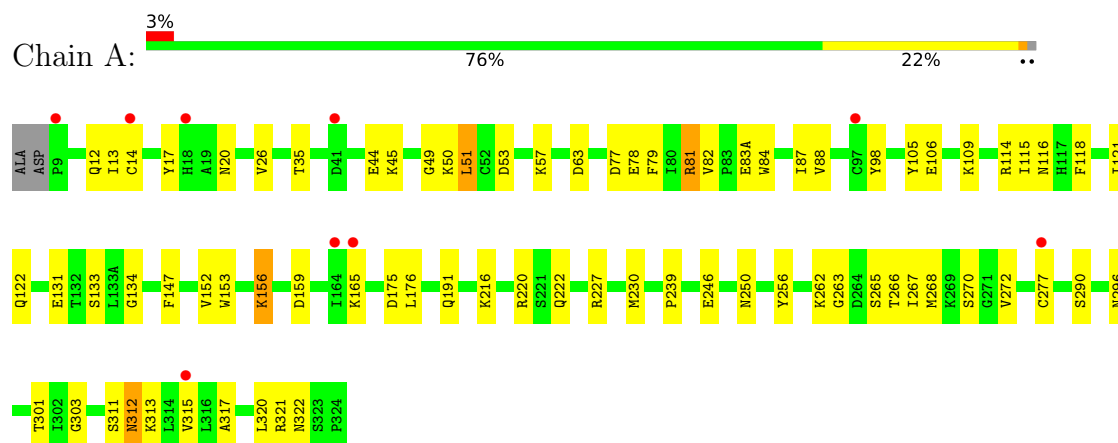
3 Residue-property plots [i](#)

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.

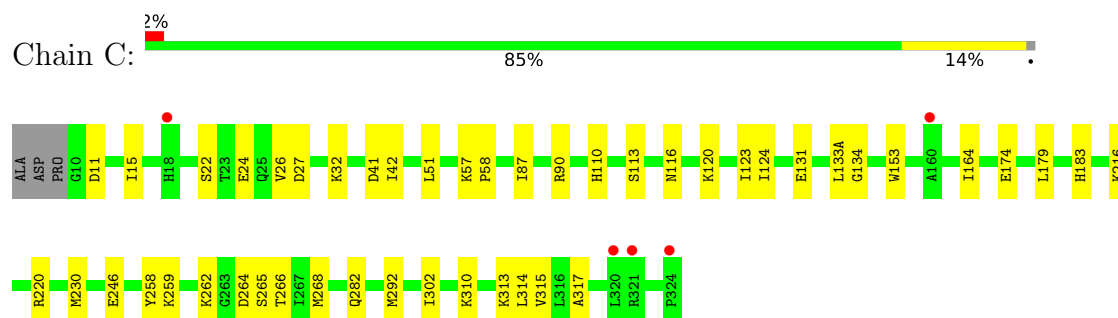
• Molecule 1: Hemagglutinin HA1



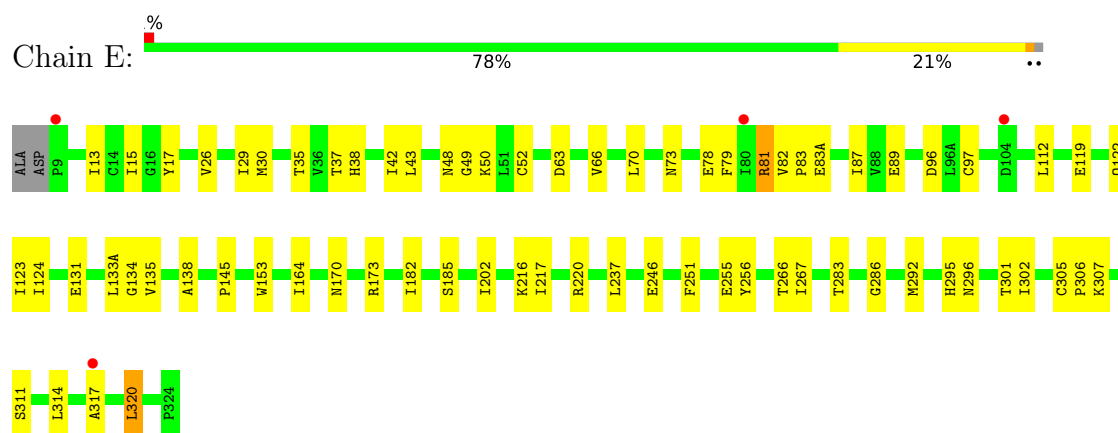
• Molecule 1: Hemagglutinin HA1



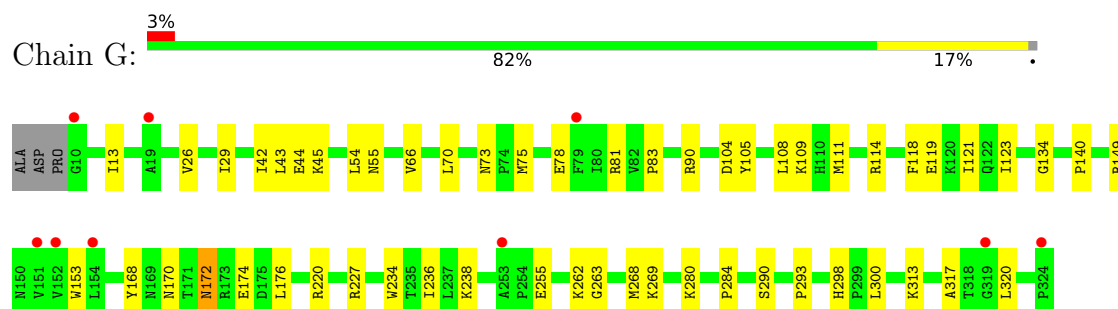
• Molecule 1: Hemagglutinin HA1



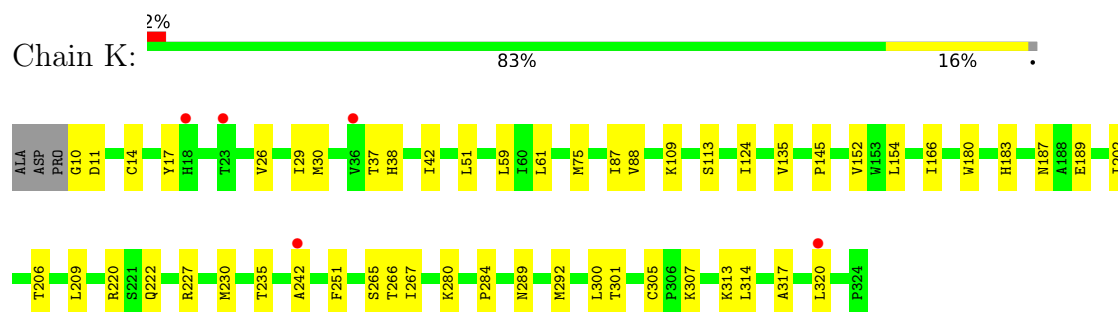
• Molecule 1: Hemagglutinin HA1



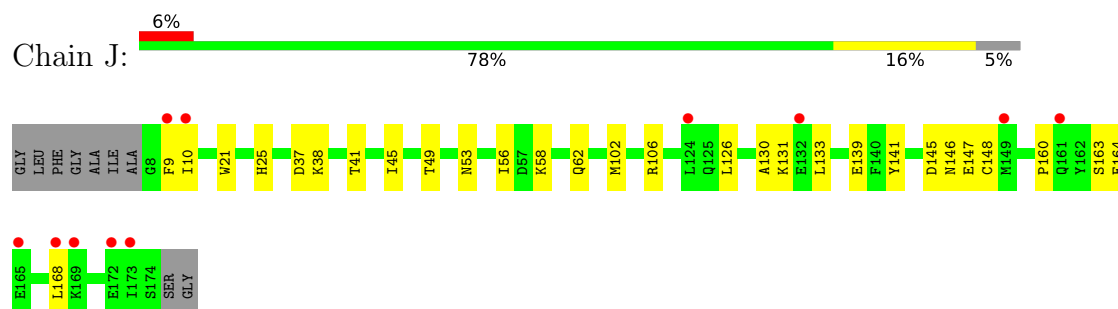
• Molecule 1: Hemagglutinin HA1



• Molecule 1: Hemagglutinin HA1

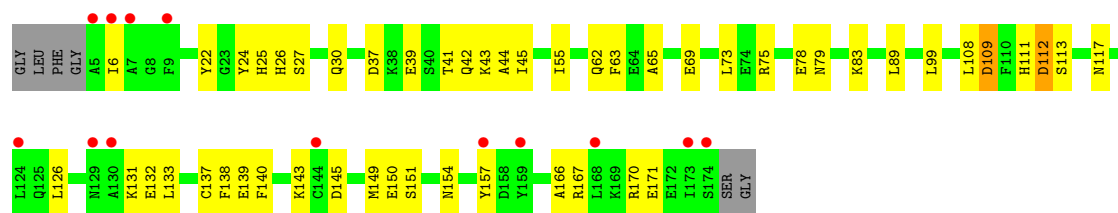


• Molecule 2: Hemagglutinin HA2

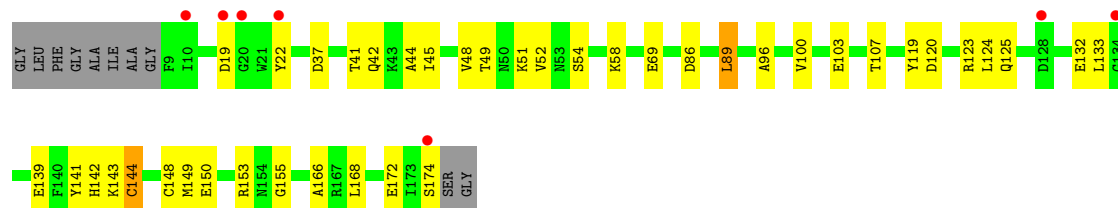


• Molecule 2: Hemagglutinin HA2

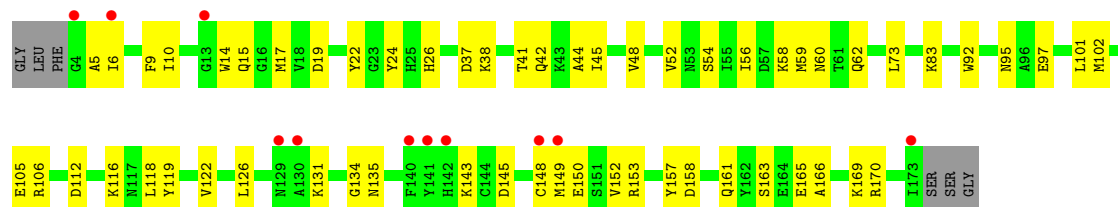




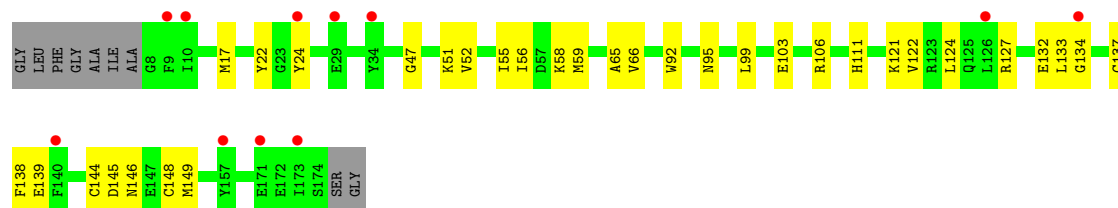
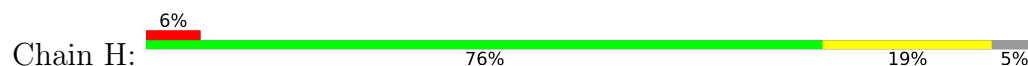
• Molecule 2: Hemagglutinin HA2



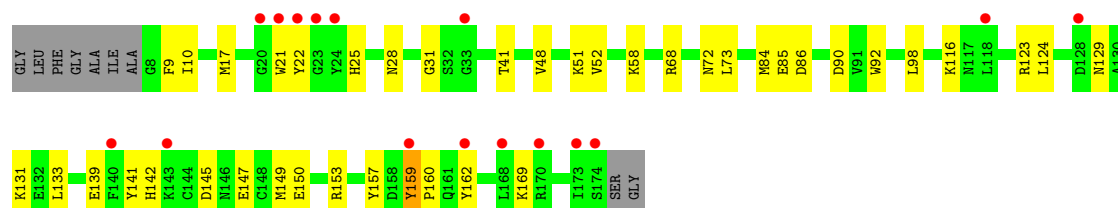
• Molecule 2: Hemagglutinin HA2



• Molecule 2: Hemagglutinin HA2



• Molecule 2: Hemagglutinin HA2



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

Chain M:  67% 33%

MAG1
MAG2
BMA3

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

Chain O:  67% 33%

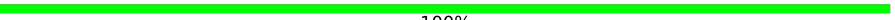
MAG1
MAG2
BMA3

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

Chain Q:  67% 33%

MAG1
MAG2
BMA3

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

Chain N:  100%

MAG1
MAG2

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

Chain P:  100%

MAG1
MAG2

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	130.37Å 72.06Å 206.86Å 90.00° 91.66° 90.00°	Depositor
Resolution (Å)	33.74 – 2.69 33.74 – 2.69	Depositor EDS
% Data completeness (in resolution range)	96.2 (33.74-2.69) 96.3 (33.74-2.69)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.29 (at 2.69Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.210 , 0.265 0.209 , 0.265	Depositor DCC
R_{free} test set	5242 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	65.5	Xtriage
Anisotropy	0.251	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 35.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.027 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	23750	wwPDB-VP
Average B, all atoms (Å ²)	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.19% 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, BMA

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	0.17	0/2618	0.33	0/3559
1	C	0.17	0/2610	0.35	0/3548
1	E	0.14	0/2618	0.32	0/3559
1	G	0.14	0/2610	0.32	0/3548
1	I	0.14	0/2610	0.32	0/3548
1	K	0.14	0/2610	0.31	0/3548
2	B	0.15	0/1411	0.35	0/1897
2	D	0.15	0/1389	0.30	0/1867
2	F	0.14	0/1409	0.31	0/1894
2	H	0.17	0/1393	0.31	0/1872
2	J	0.13	0/1393	0.31	0/1872
2	L	0.15	0/1393	0.36	0/1872
All	All	0.15	0/24064	0.32	0/32584

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	2553	0	2503	56	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2546	0	2497	26	0
1	E	2553	0	2503	44	0
1	G	2546	0	2496	38	0
1	I	2546	0	2496	32	0
1	K	2546	0	2498	37	0
2	B	1385	0	1290	42	0
2	D	1363	0	1268	28	0
2	F	1383	0	1288	48	0
2	H	1367	0	1272	27	0
2	J	1367	0	1272	24	0
2	L	1367	0	1272	38	0
3	M	39	0	34	0	0
3	O	39	0	34	0	0
3	Q	39	0	34	2	0
4	N	28	0	25	0	0
4	P	28	0	25	0	0
5	E	14	0	13	0	0
6	A	6	0	0	0	0
6	B	1	0	0	0	0
6	C	7	0	0	0	0
6	D	1	0	0	0	0
6	E	7	0	0	0	0
6	F	2	0	0	0	0
6	G	4	0	0	0	0
6	H	1	0	0	0	0
6	I	5	0	0	0	0
6	K	5	0	0	0	0
6	L	2	0	0	0	0
All	All	23750	0	22820	389	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (389) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:150:GLU:OE1	2:D:153:ARG:NH1	2.13	0.80
1:A:114:ARG:HH21	1:A:263:GLY:HA3	1.46	0.79
1:I:123:ILE:HD12	1:I:124:ILE:HG13	1.68	0.75
2:B:30:GLN:NE2	2:B:145:ASP:OD1	2.20	0.74
2:F:15:GLN:N	2:F:15:GLN:OE1	2.21	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:151:SER:HB3	2:B:157:TYR:HA	1.69	0.73
2:F:26:HIS:HB2	2:F:149:MET:HE3	1.72	0.71
2:F:83:LYS:NZ	2:H:65:ALA:O	2.21	0.71
2:B:75:ARG:NH1	2:B:78:GLU:OE1	2.24	0.71
2:B:131:LYS:HB2	2:B:139:GLU:HB3	1.73	0.71
1:I:216:LYS:O	1:I:220:ARG:NH2	2.24	0.70
1:C:87:ILE:HD12	1:C:113:SER:HA	1.72	0.70
2:F:97:GLU:HB3	2:H:58:LYS:HE3	1.73	0.69
2:D:22:TYR:H	2:D:41:THR:HG22	1.56	0.69
2:D:42:GLN:HA	2:D:45:ILE:HG13	1.74	0.68
2:H:134:GLY:HA2	2:L:124:LEU:HD21	1.75	0.68
1:A:147:PHE:HZ	1:A:230:MET:HE1	1.58	0.68
1:E:283:THR:OG1	1:E:286:GLY:O	2.12	0.68
1:K:301:THR:OG1	1:K:305:CYS:SG	2.50	0.68
2:F:134:GLY:HA2	2:H:124:LEU:HD21	1.77	0.67
1:G:238:LYS:H	1:G:238:LYS:HD2	1.60	0.67
1:A:106:GLU:N	1:A:106:GLU:OE2	2.28	0.66
1:K:280:LYS:NZ	1:K:289:ASN:OD1	2.29	0.65
1:E:131:GLU:HG3	1:E:133(A):LEU:HG	1.77	0.65
2:B:26:HIS:HB2	2:B:149:MET:HE3	1.77	0.65
2:B:166:ALA:O	2:B:170:ARG:N	2.30	0.65
2:B:167:ARG:HG2	2:B:170:ARG:HH22	1.62	0.65
1:A:313:LYS:HG3	1:A:315:VAL:HG23	1.79	0.64
2:B:69:GLU:OE2	2:B:69:GLU:N	2.29	0.64
2:B:132:GLU:HG2	2:B:138:PHE:HE2	1.62	0.63
1:A:134:GLY:HA3	1:A:153:TRP:HB3	1.79	0.63
1:K:11:ASP:OD2	2:L:28:ASN:ND2	2.29	0.63
1:E:185:SER:HB2	1:E:217:ILE:HG12	1.80	0.62
1:A:14:CYS:HB2	2:B:25:HIS:HB3	1.80	0.62
1:G:134:GLY:HA3	1:G:153:TRP:HB3	1.82	0.62
1:I:131:GLU:HG2	1:I:133(A):LEU:HG	1.82	0.62
2:L:149:MET:HE3	2:L:149:MET:H	1.64	0.61
1:A:35:THR:HG22	1:A:322:ASN:HB3	1.82	0.61
1:E:26:VAL:HG21	1:E:317:ALA:HB2	1.82	0.61
2:B:166:ALA:HB3	2:B:170:ARG:HH21	1.66	0.61
1:I:20:ASN:OD1	1:I:322:ASN:ND2	2.34	0.61
1:A:79:PHE:O	1:A:81:ARG:N	2.26	0.61
1:K:42:ILE:HA	1:K:292:MET:HE3	1.82	0.61
1:G:13:ILE:HG22	2:H:138:PHE:HB2	1.83	0.61
1:E:134:GLY:HA3	1:E:153:TRP:HB3	1.85	0.59
2:H:146:ASN:HA	2:H:149:MET:HB2	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:14:CYS:HB2	2:L:25:HIS:HB3	1.84	0.59
1:I:39:ALA:HA	1:I:318:THR:HG22	1.85	0.59
3:Q:1:NAG:H62	3:Q:2:NAG:O5	2.03	0.59
2:L:22:TYR:H	2:L:41:THR:HG22	1.68	0.59
1:E:17:TYR:HB2	1:E:320:LEU:HD22	1.84	0.58
1:A:78:GLU:HG2	1:A:79:PHE:H	1.67	0.58
2:J:102:MET:HE3	2:J:106:ARG:HH22	1.67	0.58
2:J:41:THR:O	2:J:45:ILE:HG12	2.03	0.58
1:G:81:ARG:NH1	1:G:118:PHE:O	2.37	0.58
1:G:29:ILE:HG12	2:L:51:LYS:HD2	1.86	0.57
1:K:183:HIS:CD2	1:K:230:MET:HE2	2.39	0.57
1:A:191:GLN:NE2	1:A:250:ASN:OD1	2.36	0.57
1:G:174:GLU:OE1	1:G:174:GLU:N	2.37	0.57
1:I:26:VAL:HG21	1:I:317:ALA:HB2	1.86	0.57
1:A:26:VAL:HG21	1:A:317:ALA:HB2	1.86	0.57
2:D:51:LYS:NZ	2:D:103:GLU:OE1	2.36	0.57
2:H:51:LYS:O	2:H:55:ILE:HG13	2.04	0.57
1:A:312:ASN:OD1	1:A:312:ASN:N	2.37	0.57
1:C:26:VAL:HG21	1:C:317:ALA:HB2	1.85	0.57
2:F:145:ASP:N	2:F:148:CYS:SG	2.78	0.57
1:K:87:ILE:HG12	1:K:113:SER:HA	1.87	0.57
1:A:87:ILE:HB	1:A:267:ILE:HG12	1.87	0.57
1:C:131:GLU:HG2	1:C:133(A):LEU:HG	1.87	0.56
1:G:123:ILE:HG22	1:G:168:TYR:CZ	2.40	0.56
1:K:206:THR:HG23	1:K:209:LEU:HB3	1.88	0.56
1:I:14:CYS:HB2	2:J:25:HIS:HB3	1.88	0.55
2:F:42:GLN:O	2:F:45:ILE:N	2.39	0.55
2:L:129:ASN:ND2	2:L:157:TYR:OH	2.30	0.55
1:C:134:GLY:HA3	1:C:153:TRP:HB3	1.86	0.55
1:A:216:LYS:O	1:A:220:ARG:NH2	2.40	0.55
1:A:265:SER:OG	1:A:266:THR:N	2.40	0.55
1:C:57:LYS:HG2	1:C:58:PRO:HD2	1.88	0.55
2:L:48:VAL:O	2:L:52:VAL:HG23	2.05	0.55
1:A:116:ASN:HD22	1:A:262:LYS:HG3	1.70	0.55
1:K:220:ARG:O	1:K:227:ARG:NH1	2.39	0.55
2:J:145:ASP:OD1	2:J:146:ASN:N	2.39	0.55
1:E:283:THR:HG22	1:E:301:THR:HG22	1.88	0.55
2:F:22:TYR:H	2:F:41:THR:HG22	1.72	0.55
1:A:81:ARG:HG3	1:A:82:VAL:HG13	1.87	0.55
1:G:114:ARG:NH1	1:G:263:GLY:HA3	2.22	0.55
1:A:156:LYS:HD2	1:A:159:ASP:HA	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:87:ILE:HB	1:K:267:ILE:HG12	1.89	0.55
1:K:109:LYS:HB3	1:K:267:ILE:HD11	1.88	0.54
1:E:63:ASP:N	1:E:63:ASP:OD1	2.41	0.54
2:F:5:ALA:HB1	2:F:10:ILE:HD12	1.90	0.54
1:K:152:VAL:HG22	1:K:154:LEU:HD12	1.89	0.54
2:L:145:ASP:N	2:L:145:ASP:OD1	2.40	0.54
2:F:52:VAL:O	2:F:56:ILE:HG13	2.08	0.54
1:I:75:MET:HE1	1:I:139:CYS:HA	1.89	0.54
2:B:132:GLU:HG2	2:B:138:PHE:CE2	2.43	0.54
1:E:79:PHE:HZ	1:E:119:GLU:HA	1.71	0.54
1:K:307:LYS:HG2	2:L:92:TRP:NE1	2.22	0.54
1:I:185:SER:HB2	1:I:217:ILE:HG12	1.90	0.54
1:A:17:TYR:HE1	2:B:6:ILE:HA	1.73	0.54
1:A:98:TYR:CD1	1:A:230:MET:HG3	2.42	0.53
1:C:22:SER:OG	1:C:24:GLU:OE1	2.24	0.53
2:L:129:ASN:HB3	2:L:162:TYR:CE2	2.43	0.53
2:B:55:ILE:HG12	2:B:99:LEU:HD21	1.91	0.53
2:F:165:GLU:O	2:F:169:LYS:HG2	2.09	0.53
1:G:75:MET:HE1	1:G:140:PRO:HD2	1.91	0.53
2:J:145:ASP:H	2:J:148:CYS:HB3	1.73	0.53
2:H:55:ILE:HG21	2:H:99:LEU:HD21	1.89	0.53
2:L:28:ASN:OD1	2:L:31:GLY:N	2.39	0.53
1:E:266:THR:HG22	1:E:302:ILE:HD11	1.90	0.53
2:H:52:VAL:O	2:H:56:ILE:HG13	2.09	0.53
2:J:37:ASP:OD1	2:J:38:LYS:N	2.42	0.53
1:G:90:ARG:O	1:G:269:LYS:NZ	2.40	0.53
2:D:123:ARG:NH2	2:D:132:GLU:OE1	2.43	0.52
2:F:14:TRP:HE3	2:F:17:MET:HE2	1.73	0.52
1:K:202:ILE:HD11	1:K:251:PHE:HA	1.92	0.52
2:B:22:TYR:HH	2:B:111:HIS:HD1	1.58	0.52
2:J:133:LEU:HD21	2:J:139:GLU:HB2	1.92	0.52
2:H:132:GLU:HG2	2:H:138:PHE:CE1	2.45	0.52
1:I:265:SER:OG	1:I:266:THR:N	2.42	0.52
2:B:150:GLU:OE2	2:B:154:ASN:HB2	2.10	0.52
2:B:37:ASP:O	2:B:41:THR:HG23	2.10	0.52
2:B:41:THR:O	2:B:44:ALA:N	2.43	0.51
2:B:167:ARG:HG2	2:B:170:ARG:NH2	2.25	0.51
1:A:114:ARG:NH2	1:A:263:GLY:HA3	2.21	0.51
2:F:6:ILE:HG12	2:F:112:ASP:OD2	2.10	0.51
1:C:268:MET:HE3	1:C:302:ILE:HD12	1.91	0.51
1:G:220:ARG:O	1:G:227:ARG:NH1	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:149:MET:O	2:L:153:ARG:N	2.33	0.51
1:I:284:PRO:HD3	1:I:300:LEU:O	2.11	0.51
1:C:266:THR:OG1	2:D:69:GLU:OE1	2.16	0.51
1:E:73:ASN:ND2	1:E:96:ASP:O	2.42	0.51
1:E:296:ASN:ND2	1:E:311:SER:O	2.44	0.51
1:K:26:VAL:HG21	1:K:317:ALA:HB2	1.92	0.51
1:A:321:ARG:HE	2:B:108:LEU:HD13	1.74	0.51
1:I:134:GLY:HA3	1:I:153:TRP:HB3	1.93	0.51
1:I:137:ALA:O	1:I:140:PRO:HD3	2.11	0.51
2:J:9:PHE:HB2	2:J:10:ILE:HD12	1.93	0.51
2:B:167:ARG:HG2	2:B:170:ARG:HH12	1.76	0.50
1:G:111:MET:HE1	1:G:236:ILE:HG13	1.92	0.50
1:A:20:ASN:OD1	1:A:322:ASN:ND2	2.43	0.50
1:C:216:LYS:O	1:C:220:ARG:NH2	2.43	0.50
2:L:9:PHE:CD2	2:L:10:ILE:HG13	2.47	0.50
1:A:296:ASN:ND2	1:A:311:SER:O	2.41	0.50
2:F:83:LYS:NZ	2:H:66:VAL:HA	2.26	0.50
2:J:145:ASP:OD1	2:J:147:GLU:N	2.40	0.50
1:A:121:ILE:HD12	1:A:122:GLN:H	1.76	0.50
2:F:38:LYS:HG2	2:F:38:LYS:O	2.10	0.50
2:F:106:ARG:HH11	2:H:106:ARG:NH1	2.09	0.50
2:L:145:ASP:O	2:L:147:GLU:N	2.44	0.50
1:G:78:GLU:HA	1:G:149:ARG:HH22	1.76	0.50
2:L:129:ASN:HD21	2:L:159:TYR:HD1	1.59	0.50
1:G:268:MET:HE3	1:G:284:PRO:HA	1.94	0.50
1:E:307:LYS:HD2	2:F:92:TRP:NE1	2.27	0.49
1:A:165:LYS:HG2	1:A:246:GLU:HG3	1.93	0.49
2:D:133:LEU:HD21	2:D:139:GLU:HB2	1.94	0.49
2:F:150:GLU:OE1	2:F:153:ARG:NH1	2.45	0.49
1:G:293:PRO:HD3	2:H:56:ILE:HG23	1.93	0.49
1:A:115:ILE:HG21	1:A:118:PHE:HB2	1.94	0.49
2:F:60:ASN:ND2	2:L:90:ASP:OD1	2.43	0.49
1:C:11:ASP:OD2	2:D:144:CYS:N	2.36	0.49
1:G:320:LEU:HD23	2:H:111:HIS:HB3	1.95	0.49
1:E:81:ARG:HB3	1:E:83:PRO:HD3	1.95	0.49
2:F:126:LEU:HD23	2:F:157:TYR:CE2	2.48	0.48
1:E:170:ASN:HB2	1:E:237:LEU:HD23	1.94	0.48
2:F:116:LYS:HD2	2:F:116:LYS:O	2.13	0.48
1:I:59:LEU:HD13	1:I:82:VAL:HG21	1.95	0.48
1:A:26:VAL:HG12	1:A:315:VAL:HG12	1.96	0.48
2:D:103:GLU:O	2:D:107:THR:OG1	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:29:ILE:HG23	1:K:30:MET:HG3	1.94	0.48
1:A:81:ARG:HG3	1:A:82:VAL:H	1.77	0.48
1:A:122:GLN:HB2	1:A:256:TYR:CE2	2.49	0.48
1:C:265:SER:OG	1:C:266:THR:N	2.47	0.48
2:D:120:ASP:O	2:D:124:LEU:HG	2.13	0.48
1:K:10:GLY:N	2:L:139:GLU:OE2	2.47	0.48
2:B:113:SER:O	2:B:117:ASN:ND2	2.47	0.48
1:K:187:ASN:OD1	1:K:187:ASN:N	2.47	0.48
1:K:51:LEU:HD21	1:K:88:VAL:HG11	1.95	0.48
2:J:160:PRO:HA	2:J:163:SER:HB2	1.95	0.47
1:C:123:ILE:HG13	1:C:124:ILE:HG13	1.94	0.47
1:E:70:LEU:HD11	1:E:112:LEU:HD11	1.96	0.47
2:H:51:LYS:NZ	2:H:103:GLU:OE1	2.39	0.47
1:C:90:ARG:HD2	1:C:90:ARG:HA	1.59	0.47
1:E:15:ILE:HG13	2:F:119:TYR:HA	1.97	0.47
1:E:30:MET:HG2	2:H:47:GLY:O	2.13	0.47
1:E:42:ILE:HA	1:E:292:MET:SD	2.54	0.47
1:E:216:LYS:O	1:E:220:ARG:NH2	2.46	0.47
1:A:191:GLN:HE22	1:A:250:ASN:CG	2.21	0.47
1:C:120:LYS:HB2	1:C:258:TYR:CE1	2.49	0.47
1:E:43:LEU:HB2	1:E:314:LEU:HD12	1.96	0.47
2:B:79:ASN:O	2:B:83:LYS:HG3	2.15	0.47
2:B:62:GLN:NE2	2:D:86:ASP:HB3	2.29	0.47
1:G:45:LYS:HE2	1:G:313:LYS:HE3	1.96	0.47
1:C:174:GLU:OE1	1:C:259:LYS:HE3	2.14	0.47
1:E:301:THR:HB	1:E:305:CYS:SG	2.55	0.47
2:F:19:ASP:OD1	2:F:19:ASP:N	2.42	0.47
2:F:158:ASP:OD2	2:F:161:GLN:HG3	2.15	0.47
1:E:37:THR:HG1	1:E:38:HIS:CE1	2.25	0.47
2:F:62:GLN:NE2	2:L:86:ASP:HB3	2.29	0.47
2:F:131:LYS:NZ	2:H:127:ARG:HH12	2.12	0.47
1:C:110:HIS:O	1:C:113:SER:OG	2.28	0.46
2:B:133:LEU:HD21	2:B:139:GLU:HB2	1.97	0.46
1:I:16:GLY:HA2	2:J:10:ILE:HG21	1.96	0.46
1:E:307:LYS:HD2	2:F:92:TRP:CE2	2.51	0.46
2:F:106:ARG:HE	2:F:106:ARG:HB3	1.59	0.46
2:H:59:MET:HE1	2:H:92:TRP:CH2	2.50	0.46
2:H:121:LYS:HE2	2:H:122:VAL:CG1	2.45	0.46
1:E:122:GLN:NE2	1:E:255:GLU:OE2	2.48	0.46
1:G:262:LYS:HA	1:G:262:LYS:HD3	1.74	0.46
2:L:68:ARG:NH2	2:L:85:GLU:OE2	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:ASP:OD1	1:A:63:ASP:N	2.47	0.46
1:C:41:ASP:OD1	1:C:313:LYS:HD3	2.16	0.46
1:K:189:GLU:OE2	1:K:189:GLU:N	2.49	0.46
2:D:37:ASP:O	2:D:41:THR:HG23	2.16	0.46
1:G:44:GLU:OE2	1:G:290:SER:HB2	2.16	0.46
2:H:133:LEU:HD12	2:H:137:CYS:HB3	1.98	0.46
1:K:313:LYS:O	1:K:314:LEU:HD23	2.16	0.46
2:L:72:ASN:OD1	2:L:72:ASN:N	2.48	0.46
1:A:88:VAL:HG22	1:A:268:MET:HE2	1.98	0.45
2:B:39:GLU:O	2:B:43:LYS:HG2	2.16	0.45
2:F:5:ALA:HA	2:F:9:PHE:HB3	1.99	0.45
1:I:70:LEU:HD11	1:I:112:LEU:HD11	1.98	0.45
2:J:53:ASN:HA	2:J:56:ILE:HD12	1.97	0.45
1:C:116:ASN:HD22	1:C:262:LYS:HE2	1.81	0.45
1:E:164:ILE:O	1:E:246:GLU:HA	2.16	0.45
2:F:44:ALA:O	2:F:48:VAL:HG22	2.16	0.45
1:K:280:LYS:H	1:K:280:LYS:HD2	1.81	0.45
1:A:14:CYS:HA	2:B:137:CYS:HA	1.97	0.45
1:G:26:VAL:HG11	1:G:317:ALA:HB2	1.98	0.45
2:F:73:LEU:HD23	2:F:73:LEU:HA	1.85	0.45
1:K:242:ALA:HB3	3:Q:1:NAG:H82	1.99	0.45
2:D:49:THR:HA	2:D:52:VAL:HG12	1.99	0.45
1:E:78:GLU:OE1	1:E:78:GLU:N	2.48	0.45
1:E:87:ILE:HB	1:E:267:ILE:HG12	1.99	0.45
1:E:295:HIS:HB3	1:E:306:PRO:HG2	1.99	0.45
1:K:265:SER:OG	1:K:266:THR:N	2.49	0.45
1:G:81:ARG:NE	1:G:119:GLU:HG3	2.31	0.45
1:K:320:LEU:HD11	2:L:21:TRP:HB3	1.99	0.45
1:E:29:ILE:HD12	2:F:101:LEU:HD23	1.99	0.45
1:G:73:ASN:OD1	1:G:75:MET:N	2.44	0.45
1:G:170:ASN:OD1	1:G:172:ASN:ND2	2.50	0.45
1:K:222:GLN:HG3	1:K:227:ARG:HG3	1.98	0.45
1:E:255:GLU:HG2	1:E:256:TYR:CD2	2.52	0.45
2:L:141:TYR:HB3	2:L:169:LYS:HB3	1.98	0.45
1:I:164:ILE:O	1:I:246:GLU:HA	2.16	0.45
2:J:126:LEU:HD11	2:J:130:ALA:H	1.82	0.45
1:E:82:VAL:O	1:E:83(A):GLU:N	2.50	0.45
2:F:148:CYS:O	2:F:152:VAL:HG23	2.18	0.44
1:I:24:GLU:OE2	1:I:39:ALA:HB3	2.17	0.44
1:A:53:ASP:OD2	1:A:57:LYS:HA	2.18	0.44
1:A:83(A):GLU:HB2	1:A:116:ASN:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:314:LEU:HD12	2:D:100:VAL:HG21	1.98	0.44
1:A:51:LEU:HD13	1:A:272:VAL:CG2	2.47	0.44
2:J:131:LYS:HG2	2:J:141:TYR:HD2	1.82	0.44
1:G:105:TYR:CZ	1:G:109:LYS:HD2	2.52	0.44
1:C:183:HIS:HB2	1:C:230:MET:HE2	1.99	0.44
1:C:310:LYS:HB2	2:D:89:LEU:HD11	1.98	0.44
1:G:176:LEU:HD12	1:G:176:LEU:HA	1.85	0.44
1:G:313:LYS:HD3	1:G:313:LYS:HA	1.75	0.44
2:L:17:MET:HG3	2:L:17:MET:O	2.15	0.44
1:I:155:ILE:HD12	1:I:194:LEU:HD22	2.00	0.44
2:D:142:HIS:ND1	2:D:143:LYS:O	2.51	0.44
1:G:108:LEU:HD13	1:G:234:TRP:CD2	2.53	0.44
2:L:73:LEU:HD23	2:L:73:LEU:HA	1.84	0.44
1:I:38:HIS:HB2	1:I:318:THR:HG23	2.00	0.44
1:A:320:LEU:HD21	2:B:22:TYR:HE1	1.82	0.44
2:D:44:ALA:O	2:D:48:VAL:HG22	2.17	0.44
2:H:121:LYS:CE	2:H:122:VAL:HG12	2.47	0.44
2:H:144:CYS:SG	2:H:148:CYS:HB3	2.58	0.44
1:I:122:GLN:NE2	1:I:125:PRO:HB3	2.33	0.44
1:A:175:ASP:OD1	1:A:239:PRO:HD3	2.18	0.44
1:E:48:ASN:ND2	1:E:52:CYS:SG	2.86	0.44
1:E:182:ILE:HB	1:E:202:ILE:HD12	2.00	0.44
1:G:43:LEU:HD22	1:G:45:LYS:HG3	1.99	0.44
2:L:131:LYS:HD2	2:L:133:LEU:HD23	2.00	0.44
2:J:58:LYS:HD3	2:J:58:LYS:HA	1.69	0.43
2:B:109:ASP:HA	2:B:112:ASP:HB2	1.99	0.43
2:D:141:TYR:HB3	2:D:166:ALA:HA	2.00	0.43
1:A:49:GLY:O	1:A:50:LYS:HD3	2.17	0.43
1:C:164:ILE:O	1:C:246:GLU:HA	2.18	0.43
1:A:222:GLN:HG3	1:A:227:ARG:HG3	1.99	0.43
2:B:126:LEU:HD13	2:B:140:PHE:CE1	2.53	0.43
2:F:59:MET:HE1	2:F:92:TRP:CZ3	2.54	0.43
1:K:75:MET:HE2	1:K:75:MET:HB3	1.75	0.43
1:I:29:ILE:HG12	2:D:51:LYS:HG3	2.01	0.43
1:I:182:ILE:HB	1:I:202:ILE:HD12	2.00	0.43
2:J:126:LEU:HD11	2:J:130:ALA:HB3	2.01	0.43
1:A:131:GLU:OE2	1:A:133:SER:OG	2.37	0.43
1:A:147:PHE:CZ	1:A:230:MET:HE1	2.45	0.43
1:A:303:GLY:HA2	2:B:63:PHE:CE2	2.54	0.43
2:L:149:MET:H	2:L:149:MET:CE	2.32	0.43
1:A:266:THR:HG22	2:B:65:ALA:HB1	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:149:MET:O	2:D:153:ARG:HG3	2.18	0.43
1:E:202:ILE:HD11	1:E:251:PHE:HA	2.00	0.43
2:L:133:LEU:HD21	2:L:139:GLU:HB2	2.01	0.43
2:L:142:HIS:CG	2:L:162:TYR:CE1	3.06	0.43
1:I:177:LEU:HB3	1:I:258:TYR:HB2	2.00	0.43
2:J:126:LEU:HD11	2:J:130:ALA:N	2.33	0.43
2:F:102:MET:O	2:F:106:ARG:HG2	2.19	0.43
1:G:284:PRO:HG2	1:G:298:HIS:CE1	2.54	0.43
2:D:125:GLN:HE22	2:D:155:GLY:HA2	1.84	0.43
2:D:144:CYS:SG	2:D:148:CYS:HB3	2.59	0.43
1:G:121:ILE:C	1:G:121:ILE:HD12	2.44	0.43
1:A:105:TYR:CZ	1:A:109:LYS:HD2	2.54	0.42
1:E:173:ARG:N	1:E:173:ARG:HD2	2.34	0.42
2:F:135:ASN:OD1	2:F:135:ASN:N	2.51	0.42
1:K:17:TYR:HB3	2:L:22:TYR:HE1	1.84	0.42
2:J:102:MET:O	2:J:106:ARG:HG3	2.18	0.42
1:E:15:ILE:HD11	2:F:122:VAL:HG21	2.01	0.42
1:G:284:PRO:HD3	1:G:300:LEU:O	2.19	0.42
2:J:164:GLU:O	2:J:168:LEU:HG	2.18	0.42
2:B:41:THR:OG1	2:B:42:GLN:N	2.53	0.42
2:L:159:TYR:HB3	2:L:160:PRO:HD3	2.00	0.42
2:J:102:MET:HB3	2:J:106:ARG:NH1	2.34	0.42
1:A:320:LEU:HD23	2:B:111:HIS:HB3	2.01	0.42
1:K:59:LEU:HD21	1:K:61:LEU:HD21	2.01	0.42
2:B:167:ARG:NH2	2:D:174:SER:HA	2.34	0.42
1:C:27:ASP:OD1	1:C:32:LYS:HA	2.19	0.42
1:I:26:VAL:HG11	1:I:317:ALA:HB2	2.00	0.42
1:A:78:GLU:HG2	1:A:79:PHE:N	2.34	0.42
1:E:13:ILE:HD12	1:E:13:ILE:HA	1.84	0.42
1:A:12:GLN:HA	2:B:138:PHE:O	2.19	0.42
1:C:51:LEU:HA	1:C:282:GLN:NE2	2.35	0.42
2:L:58:LYS:HA	2:L:58:LYS:HD3	1.80	0.42
1:E:66:VAL:HG23	1:E:89:GLU:OE2	2.20	0.42
2:F:37:ASP:OD2	2:F:118:LEU:HD21	2.20	0.42
1:A:81:ARG:HG3	1:A:82:VAL:N	2.35	0.42
2:D:168:LEU:O	2:D:172:GLU:HB2	2.19	0.42
1:G:104:ASP:HB2	1:G:234:TRP:HE1	1.84	0.42
2:B:73:LEU:HD23	2:B:73:LEU:HA	1.81	0.41
2:F:95:ASN:HD21	2:H:95:ASN:HD21	1.68	0.41
1:K:135:VAL:HG13	1:K:145:PRO:C	2.45	0.41
1:E:135:VAL:CG1	1:E:145:PRO:HB2	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:143:LYS:HD3	2:F:143:LYS:HA	1.90	0.41
2:H:17:MET:HE1	2:H:22:TYR:C	2.45	0.41
1:E:97:CYS:HB2	1:E:138:ALA:O	2.20	0.41
2:F:24:TYR:OH	2:F:37:ASP:OD2	2.31	0.41
2:F:118:LEU:HA	2:F:118:LEU:HD23	1.70	0.41
1:I:108:LEU:HD13	1:I:234:TRP:CD2	2.55	0.41
1:A:268:MET:HE3	1:A:270:SER:HB3	2.03	0.41
1:G:70:LEU:HD12	1:G:70:LEU:HA	1.86	0.41
1:A:44:GLU:HG3	1:A:290:SER:CB	2.51	0.41
1:A:82:VAL:HG23	1:A:84:TRP:HE3	1.85	0.41
1:E:123:ILE:HG13	1:E:124:ILE:HG13	2.02	0.41
1:K:180:TRP:HZ3	1:K:235:THR:HG22	1.86	0.41
2:J:45:ILE:O	2:J:49:THR:OG1	2.24	0.41
1:G:54:LEU:HG	1:G:55:ASN:ND2	2.35	0.41
1:G:81:ARG:NH1	1:G:83:PRO:HA	2.34	0.41
2:L:116:LYS:NZ	2:L:123:ARG:HH22	2.19	0.41
2:F:166:ALA:O	2:F:170:ARG:HG3	2.21	0.41
1:G:66:VAL:HG21	1:G:109:LYS:HG2	2.01	0.41
1:K:37:THR:OG1	1:K:38:HIS:N	2.53	0.41
2:D:144:CYS:HB3	2:D:149:MET:HE2	2.02	0.41
1:E:30:MET:HB2	2:F:105:GLU:OE1	2.21	0.41
1:G:280:LYS:HE3	1:G:280:LYS:HB3	1.59	0.41
1:A:176:LEU:HD12	1:A:176:LEU:HA	1.90	0.41
2:B:41:THR:O	2:B:45:ILE:HD12	2.20	0.41
2:F:131:LYS:NZ	2:H:127:ARG:HH22	2.18	0.41
2:F:131:LYS:HB2	2:F:131:LYS:HE2	1.85	0.41
1:K:17:TYR:HB3	2:L:22:TYR:CE1	2.55	0.41
1:I:24:GLU:HG2	1:I:39:ALA:HB3	2.01	0.41
1:A:44:GLU:OE2	1:A:45:LYS:N	2.54	0.41
2:B:43:LYS:HG2	2:B:43:LYS:H	1.61	0.41
2:F:24:TYR:CE1	2:F:153:ARG:HB2	2.56	0.41
1:K:124:ILE:HD13	1:K:166:ILE:HD13	2.03	0.41
1:K:284:PRO:HD3	1:K:300:LEU:O	2.21	0.41
1:I:320:LEU:HD11	2:J:21:TRP:CG	2.56	0.40
1:I:175:ASP:OD1	1:I:239:PRO:HD3	2.20	0.40
2:J:62:GLN:NE2	2:B:89:LEU:HD22	2.35	0.40
2:D:96:ALA:O	2:D:100:VAL:HG23	2.20	0.40
1:E:49:GLY:O	1:E:50:LYS:HD3	2.22	0.40
1:I:26:VAL:HG12	1:I:315:VAL:HG12	2.03	0.40
2:J:102:MET:HE2	2:J:106:ARG:HH12	1.86	0.40
1:K:59:LEU:HD23	1:K:87:ILE:HD12	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:84:MET:HE2	2:L:84:MET:HB3	1.84	0.40
2:L:84:MET:HE3	2:L:85:GLU:HG3	2.04	0.40
1:I:30:MET:HE2	1:I:30:MET:HB3	1.95	0.40
2:B:170:ARG:HD2	2:B:171:GLU:OE2	2.22	0.40
2:D:54:SER:O	2:D:58:LYS:HB2	2.21	0.40
2:H:145:ASP:O	2:H:149:MET:N	2.46	0.40
2:L:149:MET:SD	2:L:150:GLU:N	2.95	0.40
1:I:38:HIS:H	1:I:319:GLY:HA3	1.86	0.40
1:A:13:ILE:HD11	2:B:24:TYR:HB3	2.03	0.40
1:C:15:ILE:HG13	2:D:119:TYR:HA	2.03	0.40
1:C:42:ILE:HA	1:C:292:MET:SD	2.62	0.40
2:F:54:SER:O	2:F:58:LYS:HB2	2.21	0.40
1:G:42:ILE:H	1:G:42:ILE:HG13	1.71	0.40
2:H:133:LEU:HD21	2:H:139:GLU:HB2	2.02	0.40
1:K:14:CYS:SG	2:L:25:HIS:ND1	2.78	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	321/325 (99%)	306 (95%)	13 (4%)	2 (1%)	21	44
1	C	320/325 (98%)	311 (97%)	9 (3%)	0	100	100
1	E	321/325 (99%)	305 (95%)	15 (5%)	1 (0%)	36	60
1	G	320/325 (98%)	313 (98%)	7 (2%)	0	100	100
1	I	320/325 (98%)	306 (96%)	14 (4%)	0	100	100
1	K	320/325 (98%)	310 (97%)	10 (3%)	0	100	100
2	B	168/176 (96%)	161 (96%)	6 (4%)	1 (1%)	21	44
2	D	164/176 (93%)	161 (98%)	3 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	F	168/176 (96%)	160 (95%)	8 (5%)	0	100	100
2	H	165/176 (94%)	160 (97%)	5 (3%)	0	100	100
2	J	165/176 (94%)	156 (94%)	9 (6%)	0	100	100
2	L	165/176 (94%)	152 (92%)	12 (7%)	1 (1%)	21	44
All	All	2917/3006 (97%)	2801 (96%)	111 (4%)	5 (0%)	43	68

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	81	ARG
2	L	159	TYR
1	A	81	ARG
2	B	143	LYS
1	A	77	ASP

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	286/287 (100%)	280 (98%)	6 (2%)	47	75
1	C	285/287 (99%)	282 (99%)	3 (1%)	65	85
1	E	286/287 (100%)	284 (99%)	2 (1%)	76	90
1	G	285/287 (99%)	283 (99%)	2 (1%)	76	90
1	I	285/287 (99%)	282 (99%)	3 (1%)	65	85
1	K	285/287 (99%)	285 (100%)	0	100	100
2	B	147/150 (98%)	144 (98%)	3 (2%)	48	76
2	D	146/150 (97%)	143 (98%)	3 (2%)	47	75
2	F	146/150 (97%)	145 (99%)	1 (1%)	76	90
2	H	146/150 (97%)	145 (99%)	1 (1%)	76	90
2	J	146/150 (97%)	146 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	L	146/150 (97%)	145 (99%)	1 (1%)	76	90
All	All	2589/2622 (99%)	2564 (99%)	25 (1%)	68	86

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	I	62	LYS
1	I	123	ILE
1	I	283	THR
1	A	51	LEU
1	A	152	VAL
1	A	156	LYS
1	A	277	CYS
1	A	301	THR
1	A	312	ASN
2	B	27	SER
2	B	109	ASP
2	B	112	ASP
1	C	179	LEU
1	C	264	ASP
1	C	315	VAL
2	D	19	ASP
2	D	89	LEU
2	D	144	CYS
1	E	35	THR
1	E	320	LEU
2	F	163	SER
1	G	172	ASN
1	G	255	GLU
2	H	24	TYR
2	L	98	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	I	25	GLN
1	I	122	GLN
1	I	158	ASN
1	I	172	ASN
2	J	50	ASN
2	J	62	GLN

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Mol	Chain	Res	Type
2	J	125	GLN
1	A	55	ASN
1	A	116	ASN
1	A	170	ASN
1	A	172	ASN
2	B	53	ASN
2	B	117	ASN
2	B	125	GLN
1	C	12	GLN
1	C	21	ASN
1	C	40	GLN
1	C	116	ASN
1	C	172	ASN
1	C	222	GLN
1	C	282	GLN
2	D	62	GLN
2	D	154	ASN
2	F	50	ASN
1	G	117	HIS
1	G	122	GLN
1	G	224	ASN
2	H	95	ASN
2	H	129	ASN
2	H	161	GLN
1	K	25	GLN
1	K	172	ASN
2	L	129	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 ⓘ

13 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	M	1	3,1	14,14,15	0.26	0	17,19,21	0.52	0
3	NAG	M	2	3	14,14,15	0.22	0	17,19,21	0.46	0
3	BMA	M	3	3	11,11,12	0.52	0	15,15,17	0.89	1 (6%)
4	NAG	N	1	4,1	14,14,15	0.26	0	17,19,21	0.53	0
4	NAG	N	2	4	14,14,15	0.27	0	17,19,21	0.41	0
3	NAG	O	1	3,1	14,14,15	0.33	0	17,19,21	0.48	0
3	NAG	O	2	3	14,14,15	0.21	0	17,19,21	0.45	0
3	BMA	O	3	3	11,11,12	0.60	0	15,15,17	0.89	1 (6%)
4	NAG	P	1	4,1	14,14,15	0.25	0	17,19,21	0.51	0
4	NAG	P	2	4	14,14,15	0.29	0	17,19,21	0.44	0
3	NAG	Q	1	3,1	14,14,15	1.04	1 (7%)	17,19,21	1.05	2 (11%)
3	NAG	Q	2	3	14,14,15	0.49	0	17,19,21	0.62	0
3	BMA	Q	3	3	11,11,12	0.58	0	15,15,17	0.89	1 (6%)

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	M	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	M	2	3	-	4/6/23/26	0/1/1/1
3	BMA	M	3	3	-	0/2/19/22	0/1/1/1
4	NAG	N	1	4,1	-	4/6/23/26	0/1/1/1
4	NAG	N	2	4	-	2/6/23/26	0/1/1/1
3	NAG	O	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	O	2	3	-	2/6/23/26	0/1/1/1
3	BMA	O	3	3	-	0/2/19/22	0/1/1/1
4	NAG	P	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	P	2	4	-	2/6/23/26	0/1/1/1
3	NAG	Q	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	Q	2	3	-	1/6/23/26	0/1/1/1
3	BMA	Q	3	3	-	2/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	Q	1	NAG	O5-C1	-3.69	1.37	1.43

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	M	3	BMA	C1-O5-C5	2.28	115.24	112.19
3	Q	1	NAG	C3-C4-C5	2.24	114.30	110.23
3	O	3	BMA	C1-O5-C5	2.23	115.18	112.19
3	Q	3	BMA	C1-O5-C5	2.18	115.11	112.19
3	Q	1	NAG	O4-C4-C3	-2.15	105.31	110.38

There are no chirality outliers.

All (21) torsion outliers are listed below:

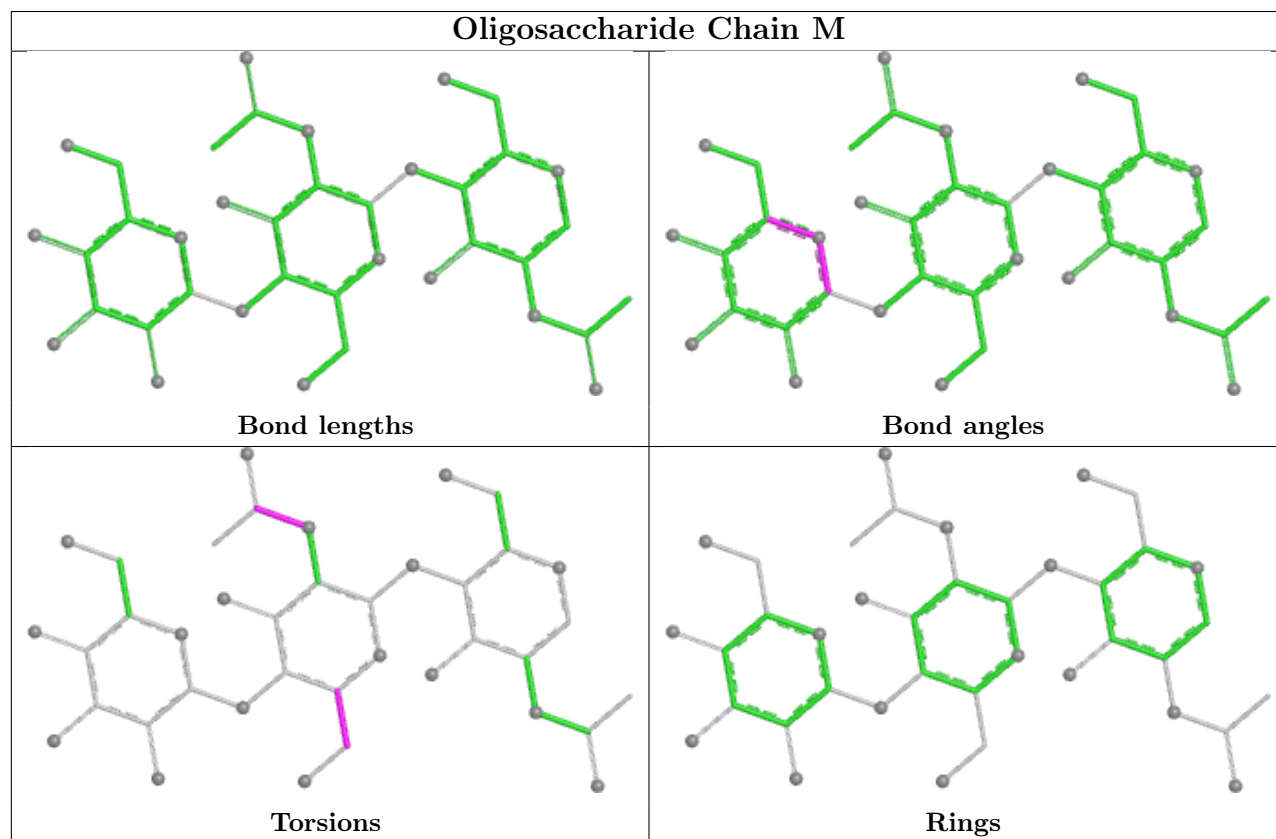
Mol	Chain	Res	Type	Atoms
3	M	2	NAG	O5-C5-C6-O6
3	M	2	NAG	C4-C5-C6-O6
4	N	1	NAG	O5-C5-C6-O6
3	M	2	NAG	C8-C7-N2-C2
3	M	2	NAG	O7-C7-N2-C2
3	O	2	NAG	C8-C7-N2-C2
3	O	2	NAG	O7-C7-N2-C2
4	N	1	NAG	C8-C7-N2-C2
4	N	1	NAG	O7-C7-N2-C2
4	P	1	NAG	C8-C7-N2-C2
4	P	1	NAG	O7-C7-N2-C2
4	P	2	NAG	O5-C5-C6-O6
4	P	2	NAG	C4-C5-C6-O6
3	O	1	NAG	C4-C5-C6-O6
3	O	1	NAG	O5-C5-C6-O6
4	N	1	NAG	C4-C5-C6-O6
3	Q	2	NAG	C3-C2-N2-C7
4	N	2	NAG	C4-C5-C6-O6
3	Q	3	BMA	O5-C5-C6-O6
4	N	2	NAG	O5-C5-C6-O6
3	Q	3	BMA	C4-C5-C6-O6

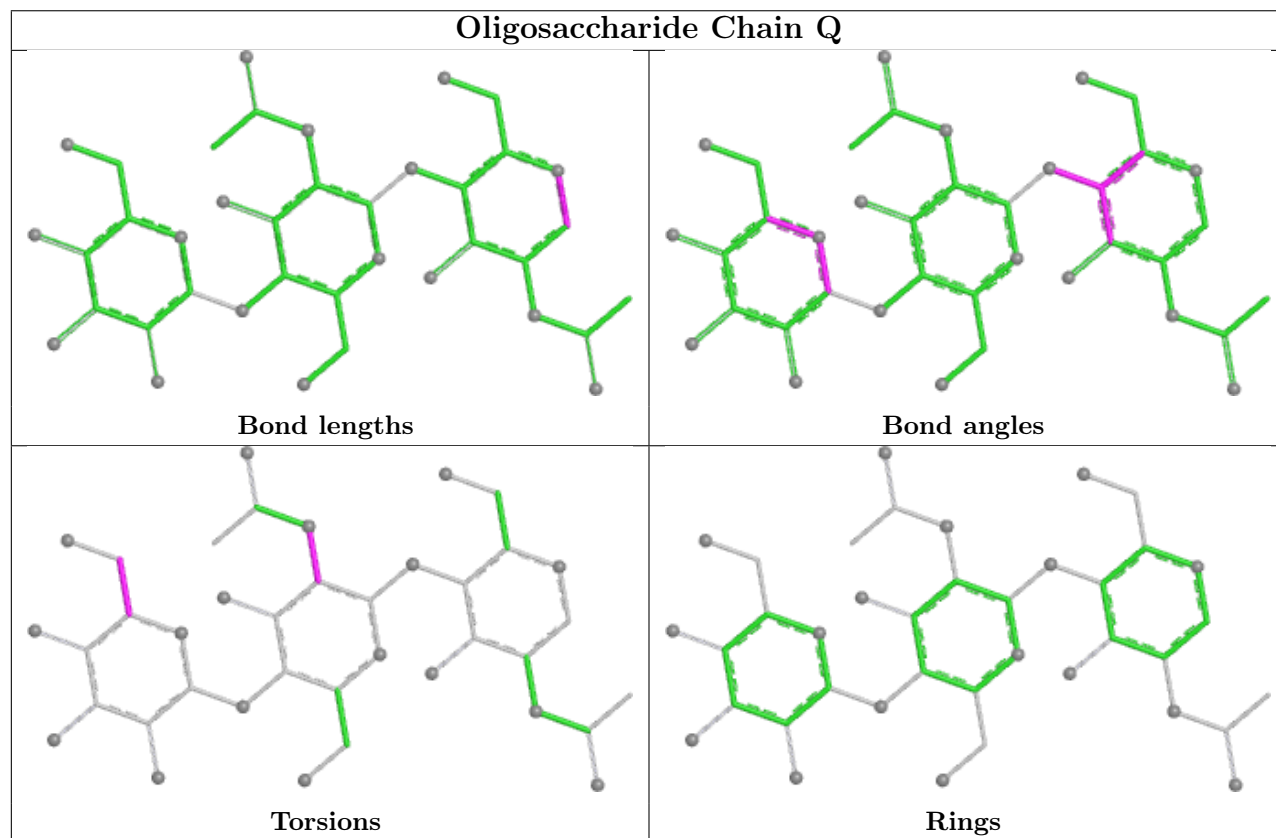
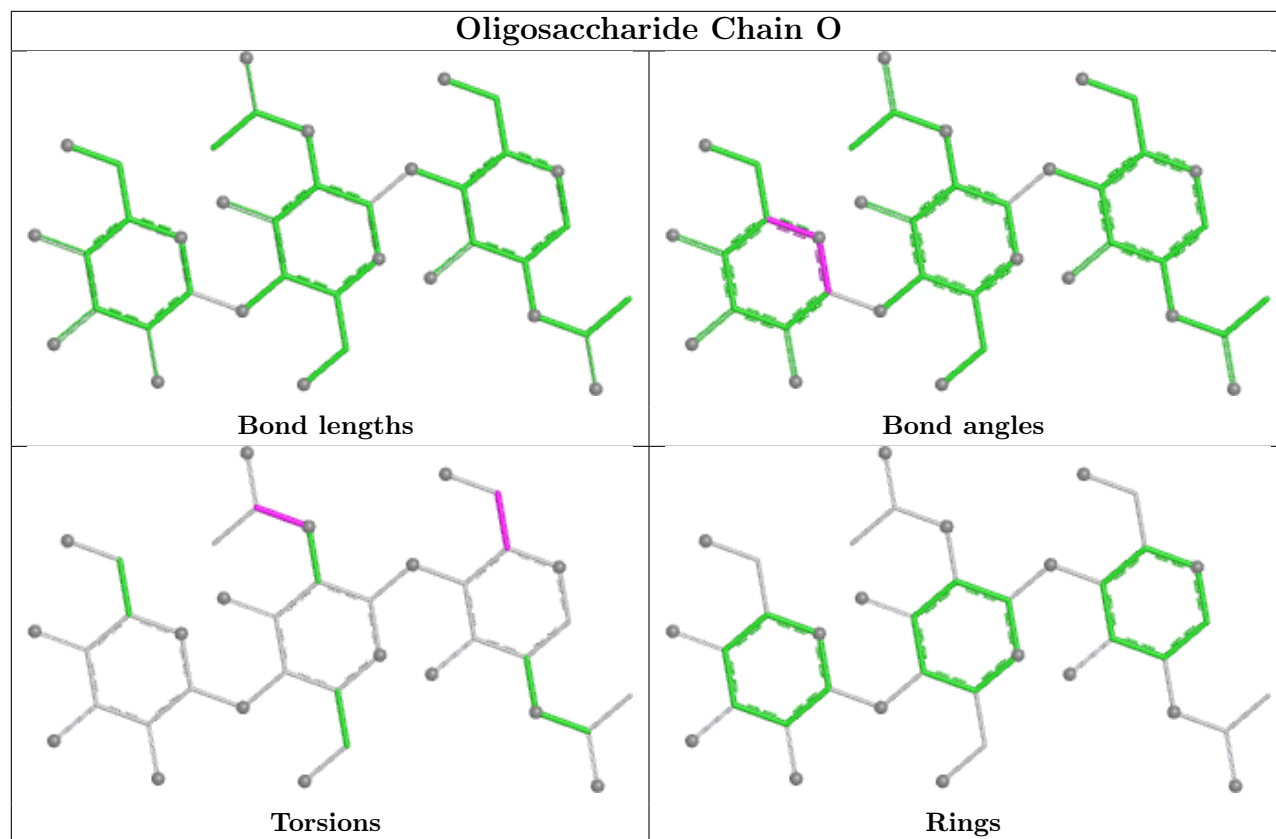
There are no ring outliers.

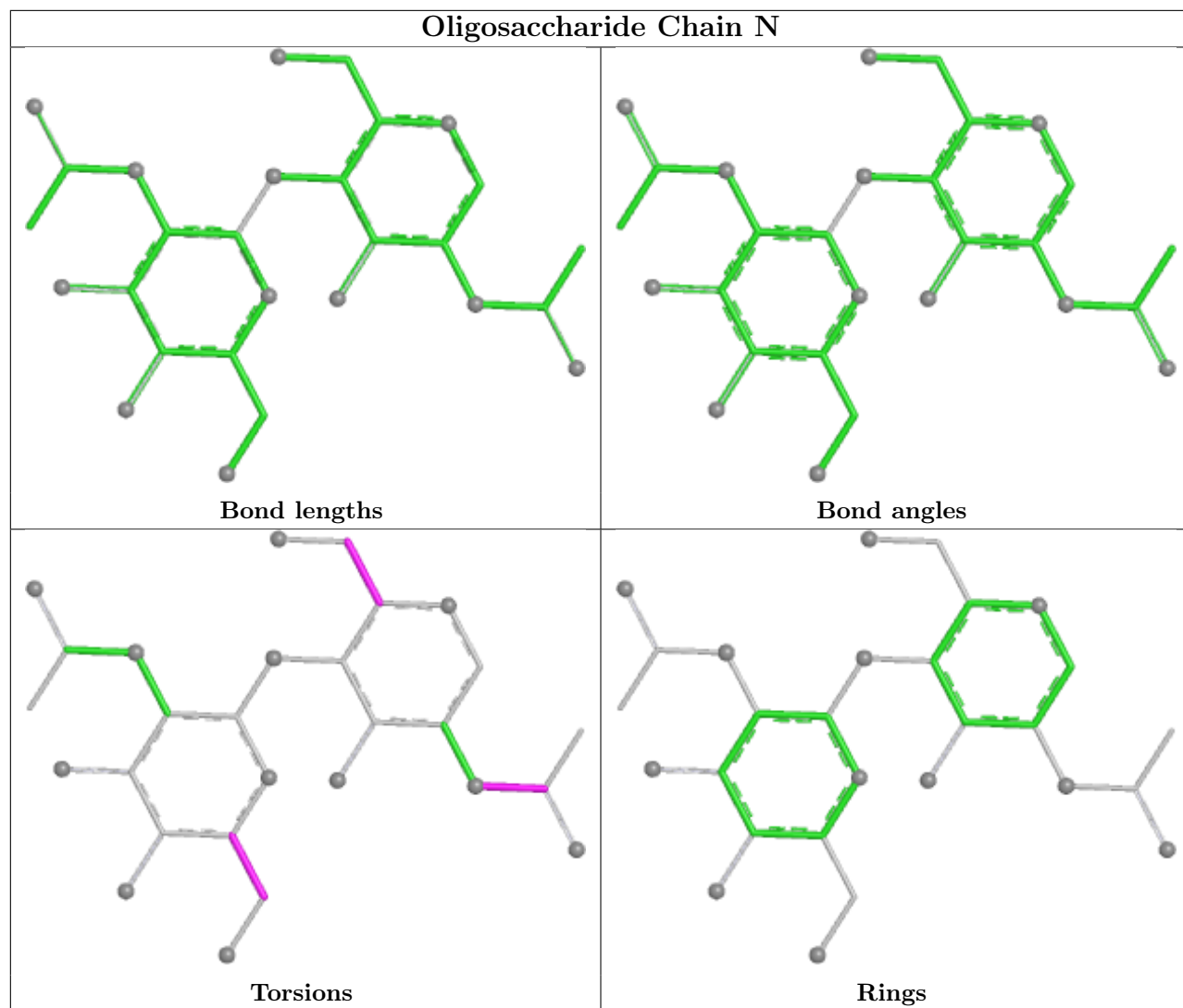
2 monomers are involved in 2 short contacts:

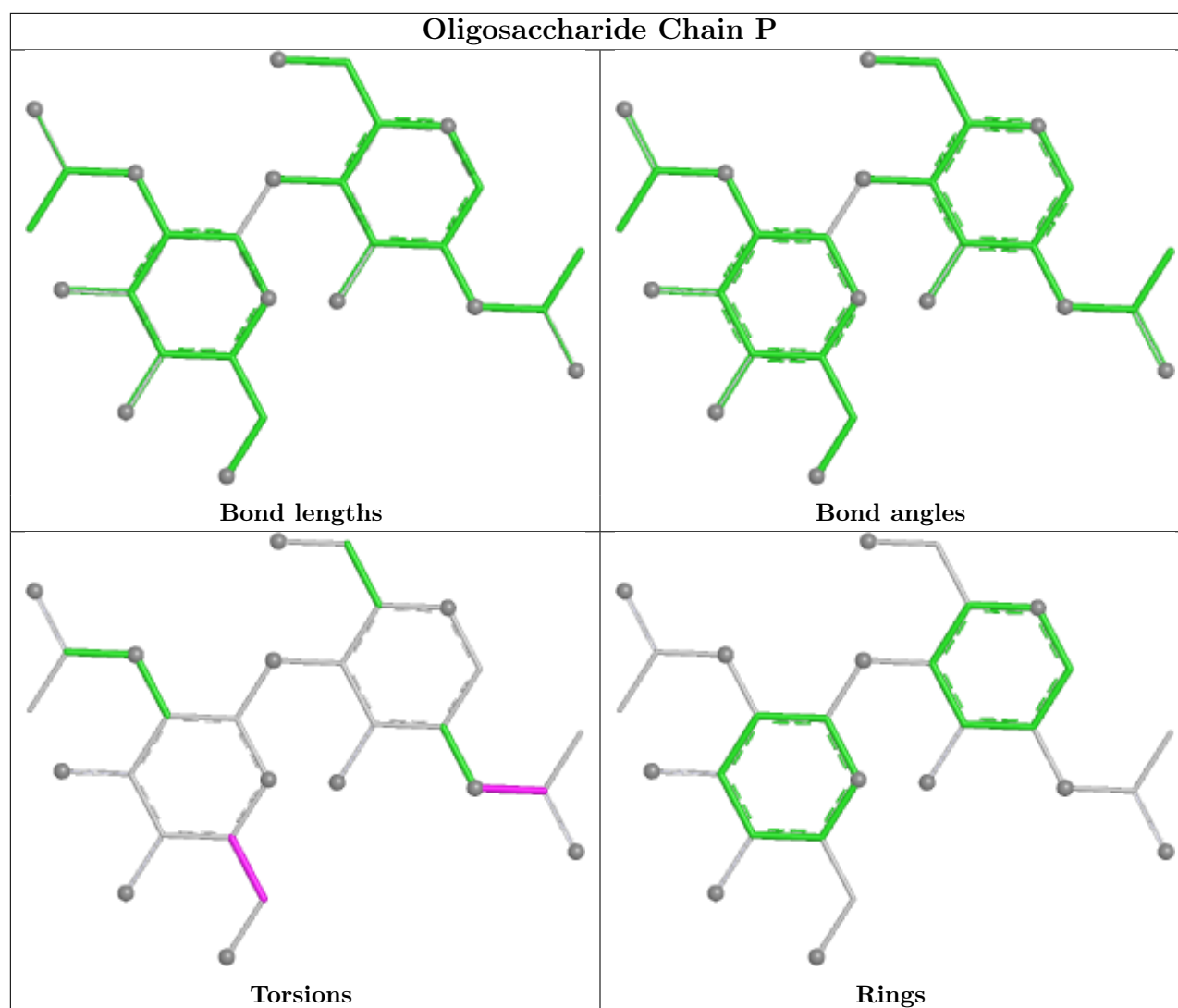
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	Q	2	NAG	1	0
3	Q	1	NAG	2	0

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









5.6 Ligand geometry [i](#)

1 ligand is 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	E	401	1	14,14,15	0.25	0	17,19,21	0.54	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
5	NAG	E	401	1	-	4/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

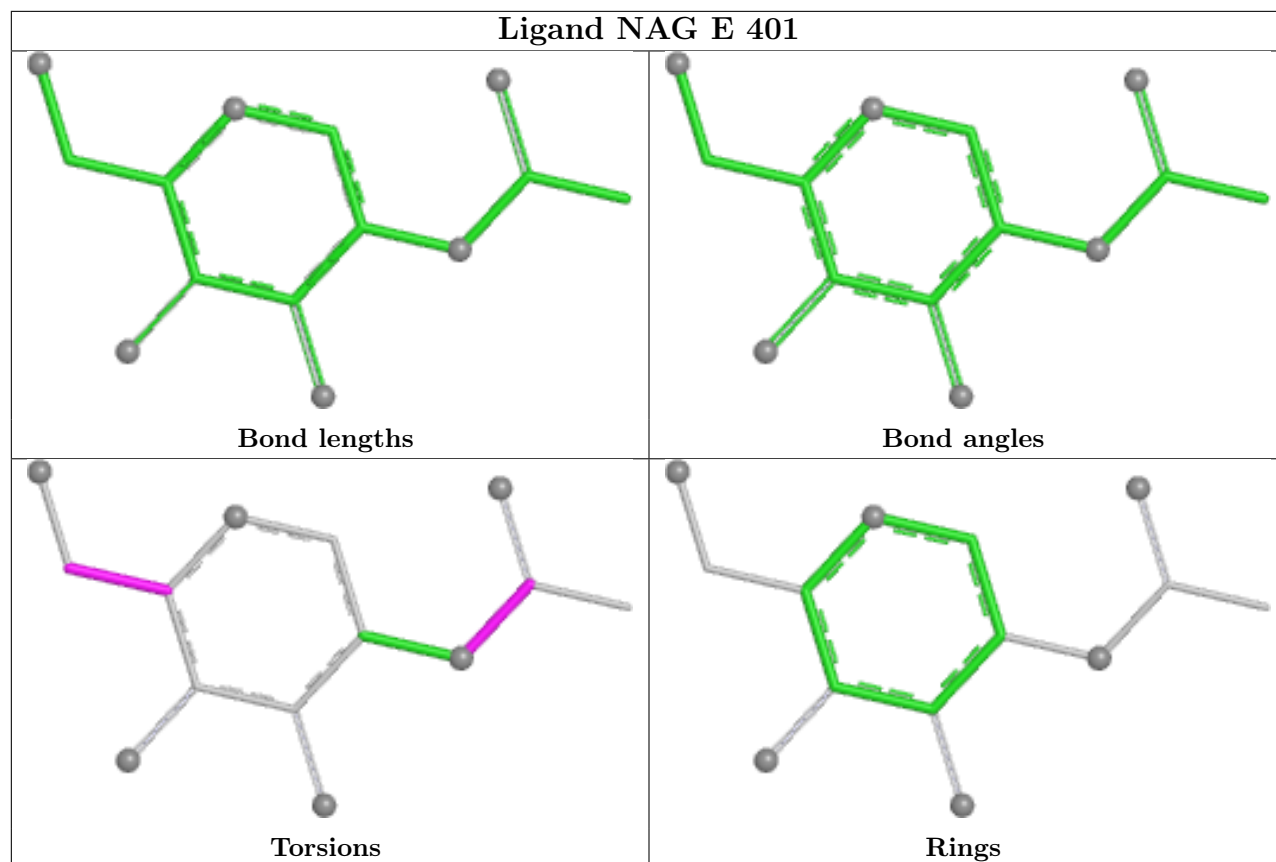
All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	E	401	NAG	C8-C7-N2-C2
5	E	401	NAG	O7-C7-N2-C2
5	E	401	NAG	O5-C5-C6-O6
5	E	401	NAG	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	323/325 (99%)	0.02	9 (2%) 55 51	48, 65, 101, 142	0
1	C	322/325 (99%)	-0.09	5 (1%) 70 68	45, 66, 101, 126	0
1	E	323/325 (99%)	-0.06	4 (1%) 76 75	43, 64, 98, 146	0
1	G	322/325 (99%)	0.20	9 (2%) 55 51	50, 77, 118, 148	0
1	I	322/325 (99%)	-0.05	4 (1%) 76 75	43, 65, 107, 140	0
1	K	322/325 (99%)	0.05	5 (1%) 70 68	49, 70, 115, 163	0
2	B	170/176 (96%)	0.47	13 (7%) 20 17	49, 99, 153, 172	0
2	D	166/176 (94%)	0.36	7 (4%) 40 37	48, 100, 138, 171	0
2	F	170/176 (96%)	0.43	11 (6%) 25 22	59, 93, 135, 210	0
2	H	167/176 (94%)	0.41	11 (6%) 24 21	52, 114, 149, 181	0
2	J	167/176 (94%)	0.45	11 (6%) 24 21	54, 108, 167, 208	0
2	L	167/176 (94%)	0.66	16 (9%) 13 11	57, 122, 174, 193	0
All	All	2941/3006 (97%)	0.17	105 (3%) 46 42	43, 76, 140, 210	0

All (105) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	22	TYR	6.5
2	L	174	SER	4.9
2	B	168	LEU	4.5
2	L	24	TYR	4.4
2	F	140	PHE	4.4
2	L	140	PHE	4.4
2	F	141	TYR	4.3
1	C	320	LEU	4.2
2	D	20	GLY	4.1
2	J	173	ILE	4.0
2	L	168	LEU	3.7

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Mol	Chain	Res	Type	RSRZ
2	F	130	ALA	3.7
1	G	151	VAL	3.7
2	J	149	MET	3.7
2	B	6	ILE	3.6
2	B	159	TYR	3.6
1	C	321	ARG	3.6
2	J	132	GLU	3.5
2	F	142	HIS	3.3
2	L	173	ILE	3.3
1	I	317	ALA	3.3
2	B	174	SER	3.2
2	F	173	ILE	3.2
1	G	253	ALA	3.1
2	J	124	LEU	3.1
2	H	9	PHE	3.1
1	A	9	PRO	3.0
2	F	6	ILE	3.0
2	B	129	ASN	2.9
2	J	9	PHE	2.9
2	F	4	GLY	2.9
2	L	128	ASP	2.9
1	A	165	LYS	2.9
2	F	13	GLY	2.8
2	L	143	LYS	2.8
2	D	10	ILE	2.8
2	L	20	GLY	2.8
2	J	161	GLN	2.8
1	G	10	GLY	2.8
2	H	24	TYR	2.8
1	G	152	VAL	2.7
2	H	171	GLU	2.7
2	D	174	SER	2.7
1	I	36	VAL	2.7
2	J	168	LEU	2.7
2	B	124	LEU	2.7
2	D	128	ASP	2.7
2	B	130	ALA	2.6
2	F	148	CYS	2.6
2	B	9	PHE	2.6
1	G	154	LEU	2.6
2	H	126	LEU	2.6
2	J	169	LYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	G	319	GLY	2.6
2	L	170	ARG	2.5
2	H	34	TYR	2.5
1	G	19	ALA	2.5
2	B	173	ILE	2.5
1	K	320	LEU	2.5
2	H	29	GLU	2.5
1	C	324	PRO	2.5
1	A	18	HIS	2.4
1	K	18	HIS	2.4
1	A	14	CYS	2.4
1	K	36	VAL	2.4
1	C	160	ALA	2.4
2	D	22	TYR	2.4
1	K	242	ALA	2.4
2	L	162	TYR	2.3
1	A	41	ASP	2.3
1	E	317	ALA	2.3
2	H	173	ILE	2.3
1	A	97	CYS	2.3
2	L	118	LEU	2.3
1	E	80	ILE	2.3
1	C	18	HIS	2.3
2	H	134	GLY	2.3
2	L	21	TRP	2.3
2	B	144	CYS	2.2
1	G	324	PRO	2.2
2	J	165	GLU	2.2
2	L	159	TYR	2.2
1	E	9	PRO	2.2
2	D	19	ASP	2.2
2	J	10	ILE	2.2
2	B	5	ALA	2.1
1	A	277	CYS	2.1
2	D	134	GLY	2.1
1	E	104	ASP	2.1
1	I	318	THR	2.1
1	K	23	THR	2.1
2	B	157	TYR	2.1
1	A	164	ILE	2.1
2	H	10	ILE	2.1
1	G	79	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
2	J	172	GLU	2.1
2	F	149	MET	2.1
1	I	97	CYS	2.1
2	H	140	PHE	2.1
2	B	7	ALA	2.1
2	F	129	ASN	2.1
1	A	315	VAL	2.1
2	H	157	TYR	2.0
2	L	23	GLY	2.0
2	L	33	GLY	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](#)

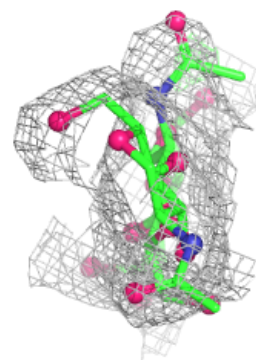
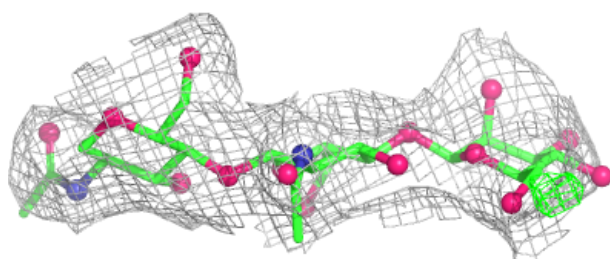
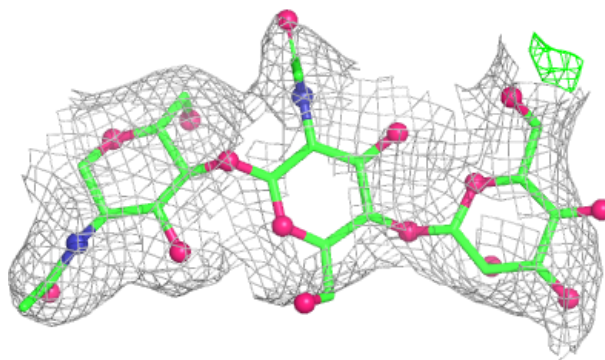
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	BMA	O	3	11/12	0.41	0.13	108,122,131,132	0
3	BMA	Q	3	11/12	0.49	0.12	115,134,147,147	0
3	BMA	M	3	11/12	0.52	0.11	104,125,141,143	0
4	NAG	P	2	14/15	0.63	0.11	95,108,119,130	0
4	NAG	N	2	14/15	0.64	0.11	98,116,126,128	0
3	NAG	Q	2	14/15	0.70	0.13	97,119,127,129	0
3	NAG	O	2	14/15	0.72	0.12	95,114,123,136	0
4	NAG	N	1	14/15	0.80	0.10	59,87,119,121	0
3	NAG	M	2	14/15	0.82	0.09	102,121,134,147	0
4	NAG	P	1	14/15	0.83	0.07	70,89,99,104	0
3	NAG	Q	1	14/15	0.88	0.16	76,88,103,124	0
3	NAG	M	1	14/15	0.91	0.09	64,78,92,109	0
3	NAG	O	1	14/15	0.92	0.12	62,76,88,92	0

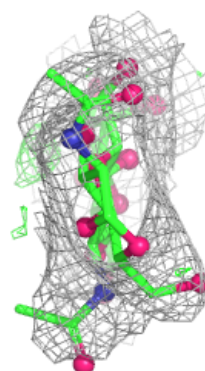
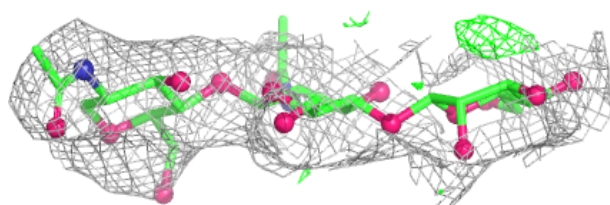
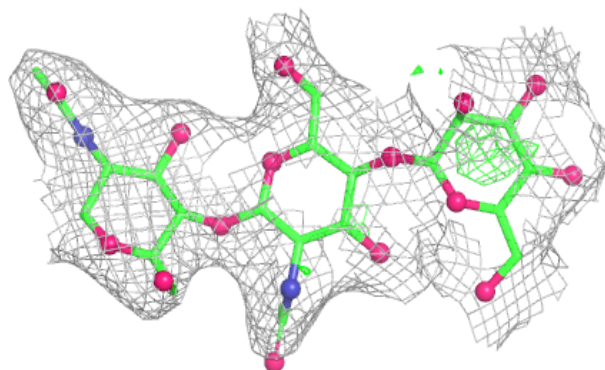
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain M:

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

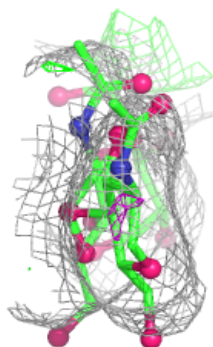
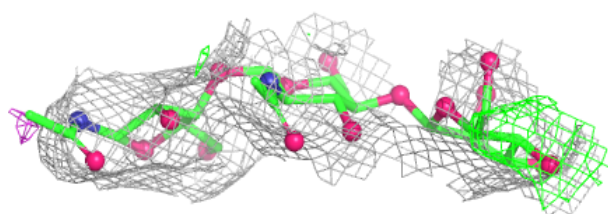
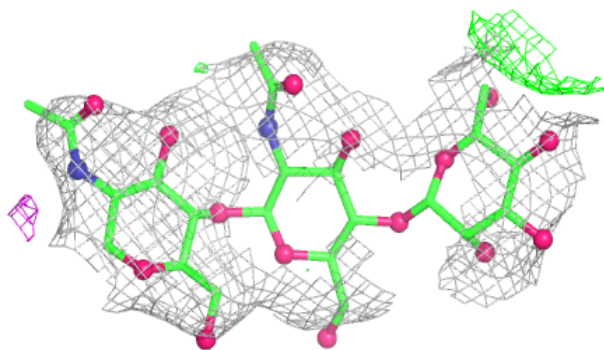
**Electron density around Chain O:**

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

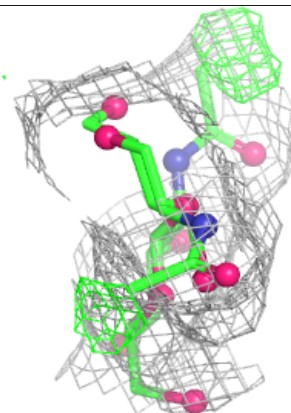
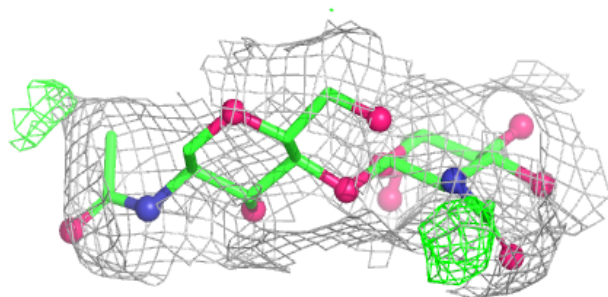
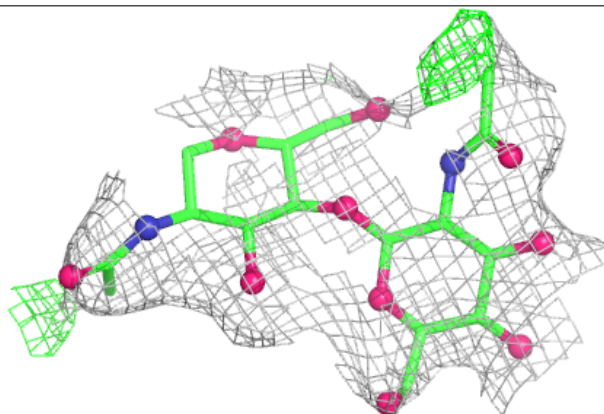


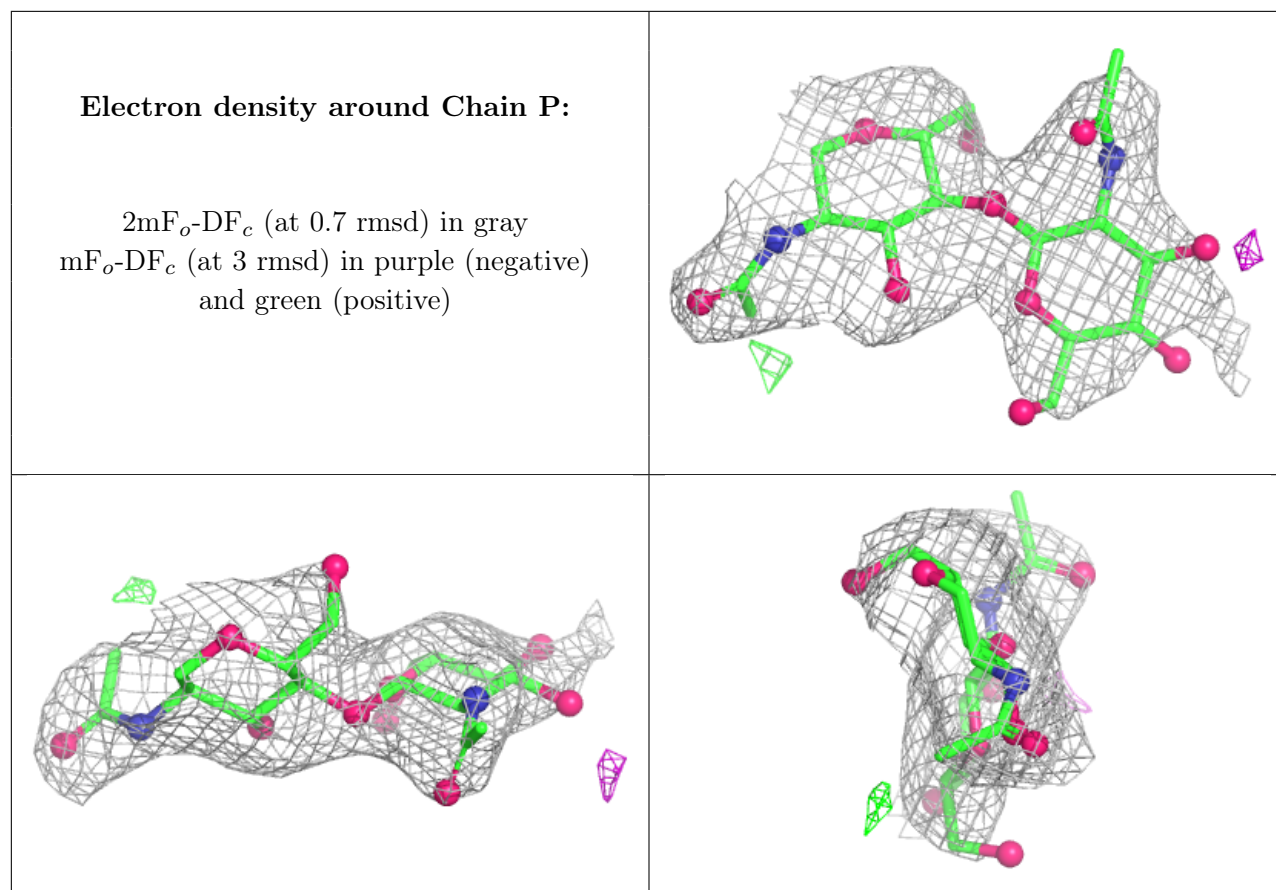
Electron density around Chain Q:

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

**Electron density around Chain N:**

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





6.4 Ligands [i](#)

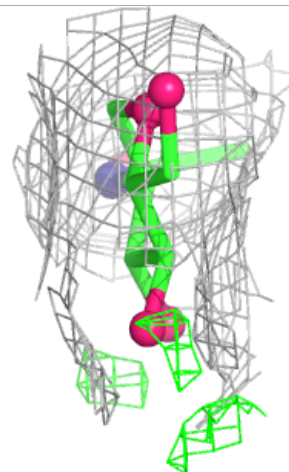
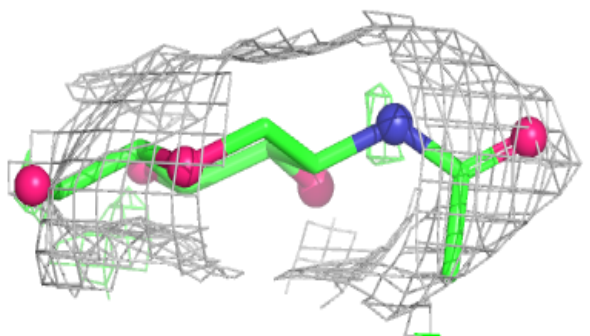
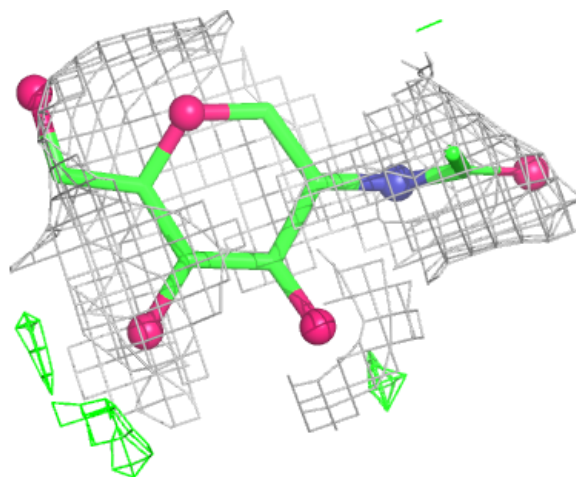
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	E	401	14/15	0.83	0.12	69,80,106,111	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around NAG E 401:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.