



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:43 AM UTC

PDB ID : 9C0V / pdb_00009c0v
Title : Crystal structure of chimeric hemagglutinin cH5/1 in complex with broad protective antibody 3E1
Authors : Nguyen, T.K.Y.; Wilson, I.A.
Deposited on : 2024-05-27
Resolution : 3.50 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

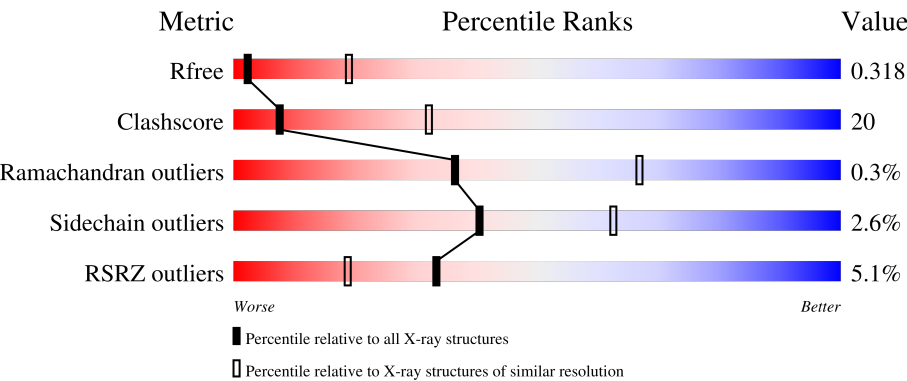
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.50 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	1085 (3.54-3.46)
Clashscore	190562	1140 (3.54-3.46)
Ramachandran outliers	187476	1113 (3.54-3.46)
Sidechain outliers	187428	1114 (3.54-3.46)
RSRZ outliers	180081	1084 (3.54-3.46)

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	541	6% 31% 25% 40%
1	B	541	% 20% 9% 71%
2	H	224	2% 68% 30%
3	L	214	2% 64% 35%
4	C	3	33% 67%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
4	S	3	

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	S	2	X	-	-	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7207 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	323	Total	C	N	O	S	0	0	0
			2552	1615	439	486	12			
1	B	157	Total	C	N	O	S	0	0	0
			1269	790	215	258	6			

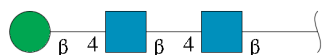
- Molecule 2 is a protein called Antibody 3E1 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	221	Total	C	N	O	S	0	0	0
			1658	1050	278	325	5			

- Molecule 3 is a protein called Antibody 3E1 Fab light chain.

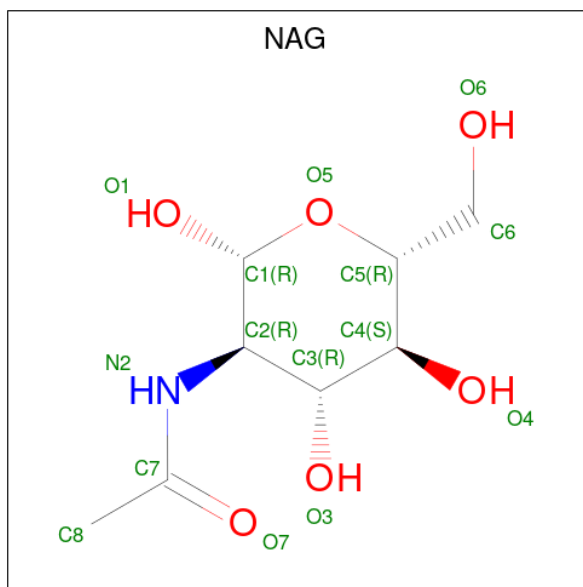
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	212	Total	C	N	O	S	0	0	0
			1636	1029	273	329	5			

- Molecule 4 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.



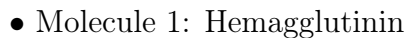
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	S	3	Total	C	N	O	0	0	0
			39	22	2	15			
4	C	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 1: Hemagglutinin



Chain C:

33%

67%

HA01
HA02
HA03

4 Data and refinement statistics

Property	Value	Source
Space group	I 21 3	Depositor
Cell constants a, b, c, α , β , γ	241.49Å 241.49Å 241.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.25 – 3.50 40.25 – 3.50	Depositor EDS
% Data completeness (in resolution range)	97.0 (40.25-3.50) 99.1 (40.25-3.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 3.48Å)	Xtriage
Refinement program	PHENIX (1.19.2_4158: ???)	Depositor
R, R_{free}	0.292 , 0.321 0.290 , 0.318	Depositor DCC
R_{free} test set	1412 reflections (4.77%)	wwPDB-VP
Wilson B-factor (Å ²)	107.5	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 105.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.029 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	7207	wwPDB-VP
Average B, all atoms (Å ²)	121.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.94% 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

5.1 Standard geometry

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

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	0.86	1/2613 (0.0%)	1.37	25/3554 (0.7%)
1	B	0.46	0/1292	0.89	6/1742 (0.3%)
2	H	0.45	0/1700	0.68	1/2323 (0.0%)
3	L	0.46	0/1674	0.86	5/2273 (0.2%)
All	All	0.63	1/7279 (0.0%)	1.05	37/9892 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
2	H	0	1
3	L	0	1
All	All	0	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	324	PRO	N-CD	-6.32	1.39	1.47

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	276	ASN	N-CA-C	30.54	150.12	112.86
1	A	276	ASN	CB-CA-C	-18.61	81.02	111.36
1	B	18	VAL	N-CA-C	-16.67	97.19	113.53
1	A	279	THR	N-CA-C	10.22	124.67	109.95
1	A	277	CYS	N-CA-CB	-9.37	92.52	109.10

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	17	MET	N-CA-C	-9.36	95.84	109.59
3	L	138	ASN	N-CA-C	8.93	123.57	111.24
1	B	58	LYS	N-CA-C	-8.88	95.04	109.96
1	A	89	GLU	N-CA-C	-7.73	97.64	109.85
1	B	58	LYS	CB-CA-C	7.12	121.29	109.53
3	L	61	ARG	N-CA-C	7.12	121.06	112.38
1	A	77	ASP	N-CA-C	-6.78	105.13	113.19
1	A	86	TYR	CB-CA-C	6.71	121.72	110.79
1	A	255	GLU	N-CA-C	-6.39	106.69	114.75
1	A	127	TRP	CB-CA-C	6.26	117.50	109.80
1	A	157	LYS	N-CA-C	6.23	122.65	110.56
1	A	46	LYS	CB-CA-C	6.12	122.34	110.46
1	A	59	LEU	CA-C-N	-6.03	115.33	123.10
1	A	59	LEU	C-N-CA	-6.03	115.33	123.10
3	L	76	SER	N-CA-C	6.00	117.62	111.14
1	A	128	SER	N-CA-C	-5.87	104.34	112.03
1	A	279	THR	CB-CA-C	-5.74	100.52	111.48
1	A	127	TRP	N-CA-C	-5.72	102.17	110.64
1	A	277	CYS	N-CA-C	-5.66	100.18	109.80
2	H	158	VAL	N-CA-CB	5.59	120.63	111.58
1	A	125	PRO	N-CA-CB	-5.42	97.56	103.25
3	L	76	SER	CB-CA-C	-5.40	102.36	110.90
1	A	279	THR	N-CA-CB	-5.36	102.46	110.60
3	L	29	ILE	N-CA-C	-5.34	107.52	112.43
1	B	17	MET	CB-CA-C	5.16	118.00	110.26
1	A	158	ASN	N-CA-C	-5.13	97.97	107.71
1	A	45	ASP	N-CA-C	-5.11	107.59	113.88
1	A	171	THR	N-CA-C	-5.07	107.45	113.38
1	A	58	PRO	CB-CA-C	-5.07	103.20	111.56
1	A	80	ILE	N-CA-C	-5.06	107.46	111.81
1	A	60	ILE	N-CA-C	-5.05	101.20	108.53
1	B	135	ASN	N-CA-C	-5.04	100.66	108.31

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	166	ARG	Sidechain
2	H	50	ARG	Sidechain
3	L	142	ARG	Sidechain

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2552	0	2515	148	0
1	B	1269	0	1191	37	0
2	H	1658	0	1640	62	0
3	L	1636	0	1591	59	0
4	C	39	0	34	1	0
4	S	39	0	34	1	0
5	A	14	0	13	1	0
All	All	7207	0	7018	287	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 20.

All (287) 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:130:HIS:CE1	1:A:161:TYR:CE1	2.04	1.46
1:A:130:HIS:CE1	1:A:161:TYR:CD1	2.24	1.26
1:A:122:GLN:HG2	1:A:256:TYR:CE2	1.81	1.14
1:A:130:HIS:NE2	1:A:161:TYR:CE1	2.19	1.10
1:A:87:ILE:HB	1:A:267:ILE:HA	1.34	1.09
1:A:130:HIS:HE1	1:A:161:TYR:CD1	1.67	1.01
1:A:121:ILE:HG13	1:A:123:ILE:HB	1.53	0.91
1:A:130:HIS:HE1	1:A:161:TYR:CE1	1.67	0.90
1:A:122:GLN:HG2	1:A:256:TYR:HE2	1.18	0.88
1:A:121:ILE:HD12	1:A:259:LYS:HE3	1.57	0.86
1:A:130:HIS:NE2	1:A:161:TYR:HE1	1.80	0.79
2:H:116:LEU:HD23	2:H:157:PRO:HD3	1.64	0.78
1:A:121:ILE:CD1	1:A:259:LYS:CE	2.61	0.78
2:H:191:THR:HG21	3:L:137:ASN:ND2	2.00	0.77
1:A:121:ILE:HD12	1:A:259:LYS:CE	2.17	0.75
1:A:121:ILE:HD13	1:A:259:LYS:HE2	1.69	0.74
1:A:201:TYR:H	1:A:248:ASN:HB3	1.50	0.73
1:A:296:ASN:ND2	1:A:311:SER:O	2.23	0.71
1:A:58:PRO:HB3	1:A:86:TYR:CE2	2.26	0.71
2:H:90:THR:HG23	2:H:118:THR:HA	1.71	0.71

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:121:ILE:CD1	1:A:259:LYS:HE3	2.20	0.71
1:A:156:LYS:HD2	1:A:196:GLN:HG3	1.72	0.71
1:A:57:LYS:HG3	1:A:58:PRO:HD2	1.73	0.70
1:A:130:HIS:CE1	1:A:161:TYR:CZ	2.78	0.70
1:A:99:PRO:HB2	1:A:229:ARG:HD3	1.73	0.69
3:L:194:CYS:O	3:L:206:THR:HA	1.92	0.69
1:A:121:ILE:CD1	1:A:259:LYS:HE2	2.22	0.69
1:A:132:ALA:HB3	1:A:155:ILE:HG22	1.75	0.69
3:L:108:ARG:NH1	3:L:170:ASP:O	2.25	0.69
1:A:108:LEU:HD21	1:A:177:LEU:HD21	1.74	0.69
1:A:130:HIS:CD2	1:A:131:GLU:HB3	2.29	0.68
1:A:289:ASN:OD1	5:A:601:NAG:O7	2.11	0.68
3:L:29:ILE:HD13	3:L:90:GLN:HG2	1.76	0.68
2:H:204:CYS:O	2:H:216:ASP:HA	1.94	0.67
1:A:323:VAL:O	1:A:323:VAL:HG23	1.94	0.67
3:L:184:ALA:HB1	3:L:187:GLU:HB2	1.75	0.67
3:L:29:ILE:HG23	3:L:92:ASN:HB2	1.77	0.67
1:A:161:TYR:HE2	1:A:249:GLY:HA2	1.60	0.67
2:H:93:TYR:O	2:H:114:GLY:HA2	1.94	0.66
1:A:151:VAL:HA	1:A:254:PRO:HA	1.77	0.65
1:A:54:LEU:HD22	1:A:57:LYS:HA	1.78	0.65
1:A:161:TYR:CE2	1:A:249:GLY:HA2	2.32	0.64
2:H:51:PHE:HD1	2:H:57:PRO:HB3	1.62	0.64
1:A:121:ILE:HG23	1:A:123:ILE:H	1.62	0.64
3:L:147:GLN:HB2	3:L:195:GLU:HB2	1.81	0.63
3:L:108:ARG:HD3	3:L:140:TYR:HB2	1.79	0.63
1:A:104:ASP:HB2	1:A:234:TRP:HE1	1.64	0.62
2:H:155:PRO:HB2	2:H:210:PRO:HG2	1.81	0.62
2:H:155:PRO:HD2	2:H:208:HIS:NE2	2.14	0.62
2:H:155:PRO:HD2	2:H:208:HIS:CE1	2.35	0.62
1:A:17:TYR:HB2	1:A:320:LEU:HD13	1.80	0.62
1:B:119:TYR:CE1	1:B:136:GLY:HA2	2.35	0.62
2:H:132:LEU:HD22	3:L:118:PHE:HB3	1.82	0.62
2:H:171:VAL:HG22	2:H:190:VAL:HG22	1.83	0.61
1:A:119:GLU:HB3	1:A:259:LYS:HB2	1.82	0.61
1:A:47:HIS:HB3	1:A:288:ILE:HA	1.83	0.60
3:L:80:PRO:HA	3:L:106:ILE:HD13	1.84	0.60
1:A:69:TRP:HE1	1:A:82:VAL:HG12	1.66	0.59
1:A:20:ASN:HD21	1:A:37:THR:HB	1.67	0.59
1:A:176:LEU:HB2	1:A:237:LEU:HB3	1.84	0.59
1:A:83(a):GLU:HA	1:A:117:HIS:ND1	2.17	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:186:ASN:HB3	1:A:219:THR:HG22	1.84	0.59
2:H:154:PHE:CG	2:H:154:PHE:O	2.54	0.59
1:A:121:ILE:HD13	1:A:259:LYS:CE	2.29	0.59
2:H:18:LEU:HD23	2:H:82:LEU:HD12	1.83	0.58
1:A:122:GLN:CG	1:A:256:TYR:CE2	2.74	0.58
1:A:125:PRO:O	1:A:125(a):LYS:HB2	2.03	0.58
3:L:85:ILE:HG12	3:L:103:LYS:HG2	1.85	0.58
1:B:51:LYS:O	1:B:55:VAL:HG23	2.04	0.57
1:A:55(a):GLY:H	1:A:278:ASN:HD21	1.51	0.57
1:A:13:LEU:HD11	1:B:24:TYR:HB3	1.86	0.57
1:A:124:ILE:O	1:A:125:PRO:C	2.48	0.57
1:A:151:VAL:HG12	1:A:252:ILE:HG22	1.85	0.57
3:L:6:GLN:HE21	3:L:99:GLY:HA3	1.67	0.57
1:A:47:HIS:NE2	1:A:289:ASN:HB3	2.20	0.57
1:B:18:VAL:HG13	2:H:100:HIS:HA	1.87	0.57
1:A:83(a):GLU:HA	1:A:117:HIS:CG	2.40	0.56
1:A:126:SER:O	1:A:127:TRP:C	2.48	0.56
1:A:170:ASN:HB2	1:A:172:ASN:HD22	1.69	0.56
1:A:121:ILE:CG1	1:A:123:ILE:HB	2.32	0.56
1:A:182:ILE:O	1:A:230:MET:HA	2.05	0.56
1:A:201:TYR:CD2	1:A:248:ASN:HB2	2.41	0.56
1:B:148:CYS:HG	1:B:162:TYR:HH	1.52	0.56
3:L:6:GLN:HE22	3:L:87:TYR:HA	1.69	0.56
1:A:115:ILE:HA	1:A:262:LYS:HA	1.88	0.56
1:A:179:LEU:O	1:A:254:PRO:HB3	2.04	0.56
1:A:54:LEU:HB2	1:A:278:ASN:HB2	1.85	0.56
1:A:87:ILE:HG13	1:A:267:ILE:HG22	1.86	0.56
3:L:108:ARG:HG2	3:L:109:THR:N	2.20	0.56
1:B:89:LEU:O	1:B:93:THR:HG23	2.05	0.56
1:A:71:LEU:O	1:A:148:PHE:HB3	2.06	0.55
3:L:48:ILE:HG12	3:L:54:LEU:HD23	1.88	0.55
1:A:169:ASN:OD1	1:A:169:ASN:O	2.23	0.55
1:A:172:ASN:O	1:A:239:PRO:HB3	2.07	0.55
1:A:151:VAL:HG12	1:A:152:VAL:N	2.22	0.55
3:L:197:THR:HG22	3:L:204:PRO:HG3	1.87	0.55
1:B:103:GLU:O	1:B:107:THR:HG23	2.06	0.55
1:A:131:GLU:HG3	1:A:154:LEU:HD22	1.89	0.55
2:H:155:PRO:HD2	2:H:208:HIS:HE2	1.72	0.54
1:A:53:ASP:HB3	1:A:277:CYS:C	2.32	0.54
1:A:130:HIS:CD2	1:A:164:ILE:HG21	2.43	0.54
2:H:122:ALA:HB3	2:H:154:PHE:CZ	2.43	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:26:HIS:O	1:B:32:SER:HA	2.08	0.54
1:B:46:ASP:O	1:B:49:THR:OG1	2.24	0.54
3:L:140:TYR:HB3	3:L:141:PRO:HD3	1.90	0.54
1:A:29:VAL:HB	1:B:105:GLU:OE1	2.08	0.54
1:A:54:LEU:HD22	1:A:58:PRO:HD3	1.90	0.54
1:A:139:CYS:HB2	1:A:146:SER:O	2.08	0.53
1:A:47:HIS:HD2	1:A:290:THR:HG23	1.73	0.53
1:A:30:LEU:HG	1:B:105:GLU:OE2	2.09	0.53
2:H:29:ILE:HG23	2:H:34:TRP:NE1	2.24	0.53
2:H:18:LEU:O	2:H:81:LYS:HA	2.09	0.53
2:H:189:VAL:HG21	3:L:135:LEU:HD13	1.89	0.53
3:L:8:PRO:HG2	3:L:11:LEU:HD21	1.90	0.53
2:H:122:ALA:HB3	2:H:154:PHE:CE2	2.43	0.52
2:H:33:TYR:CD2	2:H:52:TYR:HB2	2.43	0.52
2:H:36:TRP:NE1	2:H:80:LEU:HB2	2.23	0.52
2:H:109:ARG:HH21	3:L:55:GLU:HG3	1.75	0.52
1:A:200:THR:HB	1:A:249:GLY:N	2.24	0.52
1:A:74:PRO:HB3	1:A:141:TYR:HB2	1.91	0.52
1:A:175:ASP:OD1	1:A:239:PRO:HD3	2.10	0.52
1:B:84:VAL:HG13	1:B:88:PHE:CE1	2.45	0.52
1:A:96(a):LEU:O	1:A:97:CYS:C	2.54	0.51
1:A:123:ILE:O	1:A:124:ILE:C	2.52	0.51
2:H:85:VAL:HG13	2:H:89:ASP:HB2	1.92	0.51
3:L:125:LEU:HD21	3:L:186:TYR:CE2	2.45	0.51
3:L:186:TYR:HB2	3:L:192:TYR:OH	2.11	0.51
1:A:155:ILE:HD11	1:A:194:LEU:HD22	1.93	0.51
2:H:18:LEU:HB3	2:H:82:LEU:HB2	1.91	0.51
2:H:132:LEU:HD11	2:H:149:LEU:HD21	1.93	0.50
1:A:15:ILE:HG23	1:B:118:LEU:HD23	1.93	0.50
1:A:130:HIS:HE1	1:A:161:TYR:CG	2.25	0.50
1:A:170:ASN:OD1	1:A:239:PRO:HA	2.11	0.50
1:B:102:LEU:O	1:B:106:ARG:HG3	2.11	0.50
3:L:182:SER:HA	3:L:185:ASP:HA	1.93	0.50
2:H:29:ILE:HD12	2:H:76:ASN:C	2.37	0.50
1:A:33:ASN:ND2	4:C:1:NAG:H3	2.27	0.50
1:A:121:ILE:HG21	1:A:124:ILE:HG12	1.94	0.50
2:H:11:LEU:HD21	2:H:154:PHE:CE2	2.47	0.50
1:A:15:ILE:HD11	1:B:122:VAL:HG21	1.94	0.49
1:A:53:ASP:N	1:A:282:GLN:HE22	2.09	0.49
1:A:314:LEU:HD22	1:B:100:VAL:HG21	1.93	0.49
2:H:154:PHE:O	2:H:154:PHE:CD1	2.65	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:129:THR:HA	3:L:182:SER:HB3	1.93	0.49
2:H:14:PRO:O	2:H:15:SER:OG	2.27	0.48
1:A:19:ALA:O	1:B:15:THR:HA	2.13	0.48
1:A:108:LEU:O	1:A:112:LEU:HG	2.12	0.48
1:B:134:GLY:O	1:B:135:ASN:C	2.55	0.48
3:L:186:TYR:O	3:L:192:TYR:OH	2.30	0.48
1:A:54:LEU:C	1:A:278:ASN:HD22	2.21	0.48
2:H:14:PRO:HG3	2:H:119:VAL:HG12	1.95	0.48
2:H:66:ARG:HD2	2:H:83:THR:O	2.13	0.48
1:A:230:MET:HE1	1:A:252:ILE:HG21	1.96	0.48
1:A:274:TYR:CZ	1:A:276:ASN:HA	2.49	0.48
3:L:125:LEU:HD21	3:L:186:TYR:HE2	1.79	0.48
1:A:182:ILE:HD11	1:A:213:LEU:HD13	1.96	0.48
1:B:111:HIS:O	1:B:115:VAL:HG23	2.14	0.48
1:B:55:VAL:HG22	1:B:103:GLU:OE1	2.13	0.47
2:H:149:LEU:HB3	2:H:151:LYS:HE3	1.95	0.47
2:H:156:GLU:N	2:H:157:PRO:HD2	2.29	0.47
3:L:139:PHE:CZ	3:L:175:LEU:HG	2.48	0.47
3:L:139:PHE:HZ	3:L:175:LEU:HG	1.79	0.47
3:L:24:ARG:HG2	3:L:70:GLU:HG2	1.96	0.47
2:H:35:ILE:HD12	2:H:108:VAL:HG11	1.97	0.47
1:A:87:ILE:CB	1:A:267:ILE:HA	2.25	0.47
2:H:155:PRO:CB	2:H:210:PRO:HG2	2.44	0.47
3:L:37:GLN:HB2	3:L:47:LEU:HD11	1.95	0.47
1:A:117:HIS:HD2	1:A:260(a):VAL:HG11	1.80	0.47
1:B:112:ASP:O	1:B:116:LYS:HG3	2.14	0.47
2:H:36:TRP:HB2	2:H:69:MET:HE2	1.95	0.47
2:H:132:LEU:HD13	3:L:118:PHE:CD2	2.50	0.47
2:H:189:VAL:HG11	3:L:135:LEU:HD22	1.97	0.47
2:H:132:LEU:HB2	2:H:147:GLY:CA	2.45	0.46
2:H:132:LEU:HB2	2:H:147:GLY:HA3	1.96	0.46
3:L:49:TYR:C	3:L:51:ALA:N	2.71	0.46
1:A:108:LEU:HG	1:A:234:TRP:CD2	2.51	0.46
3:L:33:LEU:HD13	3:L:88:CYS:SG	2.56	0.46
2:H:191:THR:HG21	3:L:137:ASN:HD21	1.79	0.46
1:A:130:HIS:O	1:A:157:LYS:HG2	2.15	0.46
2:H:38:ARG:NH2	2:H:46:GLU:OE1	2.49	0.46
3:L:120:PRO:HG3	3:L:186:TYR:CE2	2.51	0.46
2:H:11:LEU:HD21	2:H:154:PHE:HE2	1.82	0.45
2:H:139:THR:O	2:H:140:SER:C	2.58	0.45
3:L:120:PRO:CG	3:L:186:TYR:CE2	2.99	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:80:PRO:HA	3:L:106:ILE:CD1	2.46	0.45
1:A:136:SER:O	1:A:145:SER:HB2	2.16	0.45
3:L:119:PRO:HB3	3:L:209:PHE:CE2	2.52	0.45
1:A:130:HIS:HE2	1:A:161:TYR:HE1	1.51	0.45
3:L:2:ILE:O	3:L:97:THR:HG21	2.16	0.45
3:L:29:ILE:HG22	3:L:32:TRP:HB2	1.98	0.45
2:H:33:TYR:CE2	2:H:52:TYR:HB2	2.52	0.45
2:H:130:PHE:HB3	3:L:121:SER:OG	2.17	0.45
1:A:206:THR:HG22	1:A:243:ILE:HG13	1.98	0.45
1:B:18:VAL:O	2:H:101:ILE:HG13	2.17	0.45
2:H:33:TYR:HB2	2:H:98:GLU:HB3	1.99	0.45
1:A:38:HIS:CD2	1:B:21:TRP:HE1	2.35	0.44
1:A:97:CYS:SG	1:A:148:PHE:CZ	3.11	0.44
1:A:201:TYR:H	1:A:248:ASN:CB	2.27	0.44
1:A:283:THR:N	1:A:286:GLY:O	2.42	0.44
1:A:64:CYS:HB2	1:A:76:CYS:SG	2.57	0.44
1:A:310:LYS:HB2	1:B:93:THR:HG21	1.98	0.44
3:L:124:GLN:O	3:L:127:SER:OG	2.26	0.44
1:A:114:ARG:HE	1:A:262:LYS:HE2	1.82	0.44
2:H:208:HIS:ND1	2:H:210:PRO:HD2	2.33	0.44
3:L:117:ILE:H	3:L:207:LYS:HZ3	1.64	0.44
1:A:130:HIS:HE1	1:A:161:TYR:CZ	2.24	0.44
1:A:105:TYR:CE2	1:A:109:LYS:HE3	2.53	0.44
2:H:37:ILE:HG12	2:H:47:TRP:HA	1.98	0.44
2:H:213:THR:HG22	2:H:215:VAL:HG23	2.00	0.44
1:B:52:VAL:O	1:B:56:ILE:HG13	2.18	0.44
2:H:154:PHE:HB3	2:H:155:PRO:HD3	2.00	0.44
1:B:97:GLU:O	1:B:101:LEU:HG	2.18	0.43
3:L:49:TYR:C	3:L:51:ALA:H	2.25	0.43
2:H:38:ARG:HG2	2:H:38:ARG:HH11	1.83	0.43
1:A:170:ASN:C	1:A:172:ASN:N	2.76	0.43
1:A:281:CYS:SG	1:A:304:LYS:O	2.75	0.43
1:B:141:TYR:O	1:B:169:ASN:ND2	2.49	0.43
1:A:267:ILE:O	1:A:268:MET:C	2.61	0.43
1:B:17:MET:HB3	1:B:17:MET:HE3	1.61	0.43
1:A:151:VAL:CG1	1:A:152:VAL:N	2.82	0.43
1:A:54:LEU:CD2	1:A:58:PRO:HD3	2.47	0.43
1:A:198:PRO:O	1:A:199:THR:C	2.62	0.43
3:L:29:ILE:CD1	3:L:33:LEU:HD21	2.48	0.43
1:A:99:PRO:HB2	1:A:229:ARG:CD	2.44	0.43
1:A:170:ASN:HB2	1:A:172:ASN:ND2	2.34	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:292:LEU:HB3	1:A:293:PRO:HD2	2.01	0.43
1:B:30:GLN:OE1	1:B:146:ASN:N	2.37	0.43
1:B:55:VAL:O	1:B:58:LYS:HG3	2.18	0.43
1:A:74:PRO:HA	1:A:77:ASP:OD2	2.19	0.43
1:A:151:VAL:CG1	1:A:252:ILE:HG22	2.48	0.43
1:A:164:ILE:O	1:A:164:ILE:HD12	2.19	0.43
3:L:49:TYR:HD2	3:L:50:LYS:HG3	1.84	0.43
3:L:184:ALA:CB	3:L:187:GLU:HB2	2.44	0.43
1:A:83(a):GLU:O	1:A:117:HIS:HB2	2.19	0.43
1:A:230:MET:HE3	1:A:252:ILE:HG13	2.01	0.42
1:B:84:VAL:O	1:B:88:PHE:HD1	2.02	0.42
1:B:18:VAL:HG13	2:H:100:HIS:ND1	2.34	0.42
3:L:117:ILE:HG13	3:L:207:LYS:HD3	2.01	0.42
2:H:59:TYR:HB2	2:H:64:ARG:HG2	2.02	0.42
3:L:4:MET:HE3	3:L:23:CYS:SG	2.60	0.42
1:A:170:ASN:C	1:A:172:ASN:H	2.27	0.42
1:A:266:THR:OG1	1:A:302:ILE:HD11	2.20	0.42
1:B:54:SER:O	1:B:57:GLU:HB2	2.20	0.42
1:B:100:VAL:HG12	1:B:104:ASN:HD21	1.84	0.42
2:H:155:PRO:HG2	2:H:210:PRO:HB2	2.01	0.42
3:L:29:ILE:CD1	3:L:90:GLN:HG2	2.47	0.42
1:A:74:PRO:HG2	1:A:139:CYS:O	2.19	0.41
1:A:288:ILE:HD11	1:A:295:GLN:HE22	1.85	0.41
1:A:73:ASN:ND2	1:A:97:CYS:SG	2.93	0.41
1:A:122:GLN:CG	1:A:256:TYR:HE2	2.08	0.41
1:B:61:THR:O	1:B:62:GLN:C	2.63	0.41
2:H:137:LYS:HA	2:H:137:LYS:HD2	1.95	0.41
3:L:110:VAL:HG22	3:L:141:PRO:HG3	2.01	0.41
1:A:65:SER:HB3	1:A:93:PRO:HG3	2.01	0.41
1:A:195:TYR:CE2	1:A:250:ASN:HA	2.56	0.41
1:A:230:MET:CE	1:A:252:ILE:HG21	2.50	0.41
2:H:131:PRO:HD3	2:H:217:LYS:NZ	2.34	0.41
1:A:98:TYR:O	1:A:99:PRO:C	2.63	0.41
1:A:61:LEU:O	1:A:62:ARG:C	2.63	0.41
1:A:75:MET:O	1:A:76:CYS:HB2	2.20	0.41
1:B:26:HIS:HB2	1:B:149:MET:HE3	2.02	0.41
1:A:181:GLY:HA2	1:A:231:GLU:O	2.20	0.41
3:L:49:TYR:CD2	3:L:50:LYS:HG3	2.55	0.41
1:A:96(a):LEU:HD21	1:A:232:PHE:HD2	1.86	0.41
1:A:272:LEU:HD13	1:A:272:LEU:HA	1.86	0.41
1:A:127:TRP:NE1	1:A:166:ARG:HE	2.19	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:165:LYS:HG3	1:A:246:GLU:HG2	2.03	0.41
2:H:153:TYR:CE1	2:H:184:TYR:HB2	2.56	0.41
3:L:122:ASP:HA	3:L:125:LEU:HD12	2.03	0.41
1:A:70:LEU:O	1:A:150:ASN:ND2	2.54	0.41
1:A:130:HIS:CD2	1:A:164:ILE:CG2	3.04	0.41
1:A:168:TYR:CZ	1:A:170:ASN:HA	2.56	0.41
3:L:111:ALA:HB3	3:L:140:TYR:H	1.86	0.41
1:A:124:ILE:HB	1:A:255:GLU:O	2.21	0.40
1:A:61:LEU:HD23	1:A:61:LEU:HA	1.92	0.40
1:A:196:GLN:HE21	1:A:196:GLN:HB3	1.61	0.40
2:H:131:PRO:HD3	2:H:217:LYS:CE	2.52	0.40
3:L:107:LYS:HA	3:L:140:TYR:OH	2.21	0.40
3:L:140:TYR:O	3:L:141:PRO:C	2.64	0.40
3:L:146:VAL:HG22	3:L:196:VAL:HG22	2.03	0.40
1:B:145:ASP:HB2	1:B:147:THR:OG1	2.21	0.40
3:L:161:GLU:HA	3:L:176:SER:O	2.22	0.40
1:A:76:CYS:C	1:A:78:GLU:H	2.28	0.40
1:A:175:ASP:C	1:A:176:LEU:HD12	2.46	0.40
1:A:176:LEU:O	1:A:236:ILE:HA	2.21	0.40
2:H:206:VAL:O	2:H:214:LYS:HA	2.22	0.40
4:S:1:NAG:H5	4:S:2:NAG:H83	2.03	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	321/541 (59%)	275 (86%)	43 (13%)	3 (1%)	14	47
1	B	153/541 (28%)	144 (94%)	9 (6%)	0	100	100
2	H	219/224 (98%)	205 (94%)	14 (6%)	0	100	100
3	L	210/214 (98%)	196 (93%)	14 (7%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	903/1520 (59%)	820 (91%)	80 (9%)	3 (0%)	36 67

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	125	PRO
1	A	97	CYS
1	A	125(a)	LYS

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	291/475 (61%)	277 (95%)	14 (5%)	23 49
1	B	138/475 (29%)	136 (99%)	2 (1%)	59 71
2	H	190/193 (98%)	189 (100%)	1 (0%)	81 80
3	L	186/188 (99%)	182 (98%)	4 (2%)	45 65
All	All	805/1331 (60%)	784 (97%)	21 (3%)	40 62

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	86	TYR
1	A	125	PRO
1	A	125(a)	LYS
1	A	125(b)	SER
1	A	130	HIS
1	A	135	VAL
1	A	155	ILE
1	A	158	ASN
1	A	177	LEU
1	A	238	LYS
1	A	252	ILE
1	A	272	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	278	ASN
1	A	290	THR
1	B	17	MET
1	B	66	VAL
2	H	50	ARG
3	L	5	THR
3	L	7	SER
3	L	33	LEU
3	L	93	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	ASN
1	A	38	HIS
1	A	103	ASN
1	A	130	HIS
1	A	158	ASN
1	A	172	ASN
1	A	183	HIS
1	A	196	GLN
1	A	250	ASN
1	A	278	ASN
1	A	282	GLN
1	A	295	GLN
1	B	25	HIS
1	B	53	ASN
1	B	114	ASN
2	H	172	HIS
3	L	6	GLN
3	L	137	ASN
3	L	138	ASN
3	L	160	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 ⓘ

6 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	$\# Z > 2$
4	NAG	C	1	4	14,14,15	0.41	0	17,19,21	0.54	0
4	NAG	C	2	4	14,14,15	0.41	0	17,19,21	0.71	1 (5%)
4	BMA	C	3	4	11,11,12	0.19	0	15,15,17	0.53	0
4	NAG	S	1	4,1	14,14,15	0.48	0	17,19,21	0.98	2 (11%)
4	NAG	S	2	4	14,14,15	0.38	0	17,19,21	0.62	0
4	BMA	S	3	4	11,11,12	0.22	0	15,15,17	0.41	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
4	NAG	C	1	4	-	3/6/23/26	0/1/1/1
4	NAG	C	2	4	-	3/6/23/26	0/1/1/1
4	BMA	C	3	4	-	0/2/19/22	0/1/1/1
4	NAG	S	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	S	2	4	1/1/5/7	4/6/23/26	0/1/1/1
4	BMA	S	3	4	-	2/2/19/22	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	S	1	NAG	C4-C3-C2	2.62	114.86	111.02
4	C	2	NAG	O5-C1-C2	-2.29	107.74	111.29
4	S	1	NAG	O5-C1-C2	2.26	114.79	111.29

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	S	2	NAG	C1

All (14) torsion outliers are listed below:

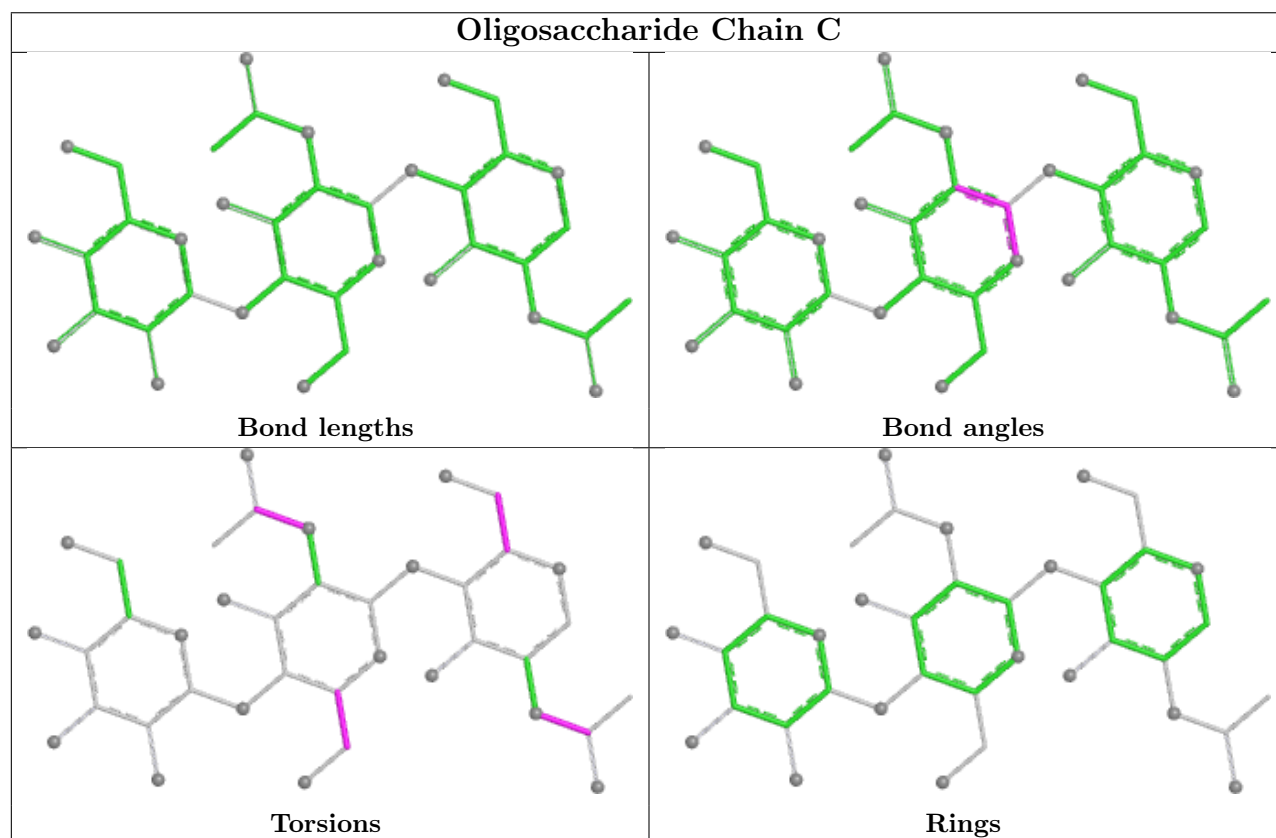
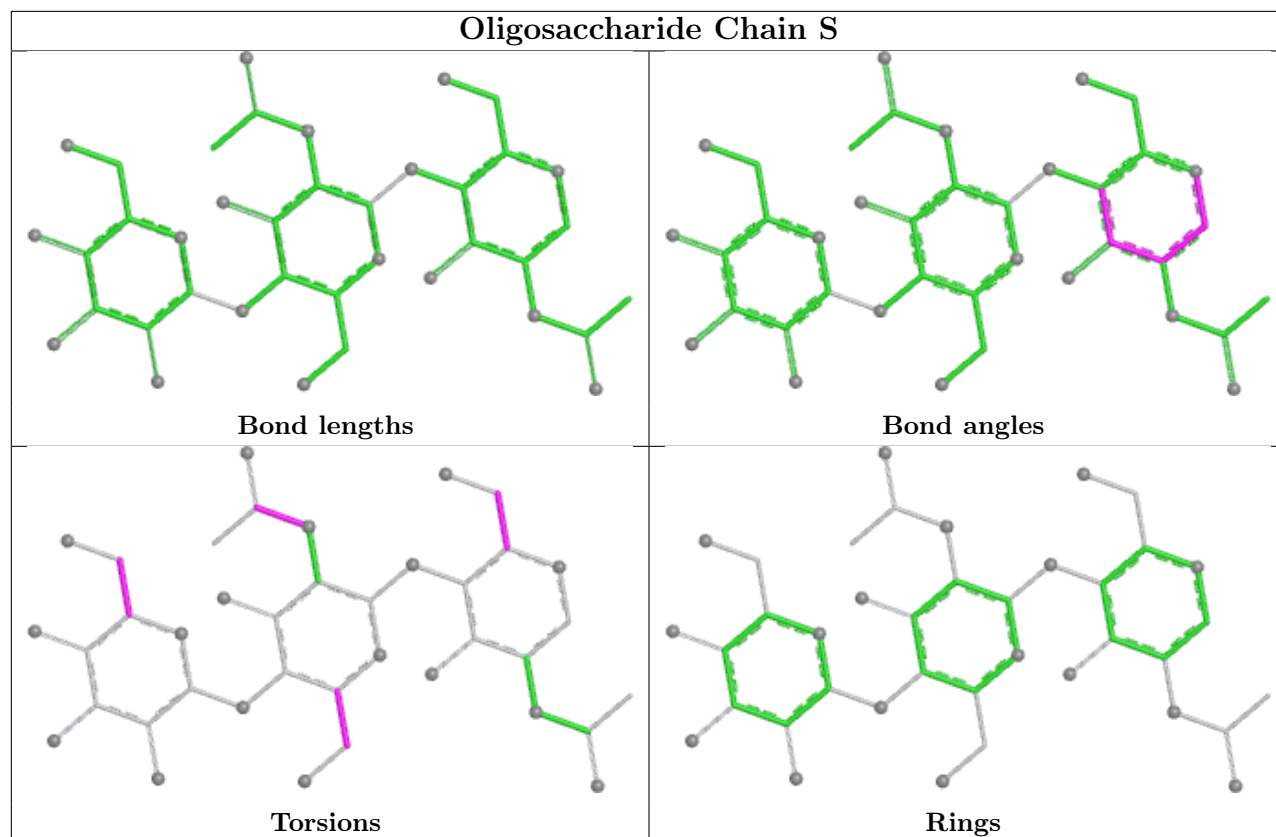
Mol	Chain	Res	Type	Atoms
4	S	2	NAG	C8-C7-N2-C2
4	S	2	NAG	O7-C7-N2-C2
4	C	1	NAG	C8-C7-N2-C2
4	C	1	NAG	O7-C7-N2-C2
4	S	2	NAG	C4-C5-C6-O6
4	C	2	NAG	C8-C7-N2-C2
4	C	2	NAG	O7-C7-N2-C2
4	S	3	BMA	O5-C5-C6-O6
4	S	2	NAG	O5-C5-C6-O6
4	S	1	NAG	C4-C5-C6-O6
4	S	3	BMA	C4-C5-C6-O6
4	S	1	NAG	O5-C5-C6-O6
4	C	1	NAG	O5-C5-C6-O6
4	C	2	NAG	C4-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	S	2	NAG	1	0
4	C	1	NAG	1	0
4	S	1	NAG	1	0

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



5.6 Ligand geometry

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	A	601	-	14,14,15	0.42	0	17,19,21	0.62	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	A	601	-	-	5/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	601	NAG	C8-C7-N2-C2
5	A	601	NAG	O7-C7-N2-C2
5	A	601	NAG	C4-C5-C6-O6
5	A	601	NAG	O5-C5-C6-O6
5	A	601	NAG	C1-C2-N2-C7

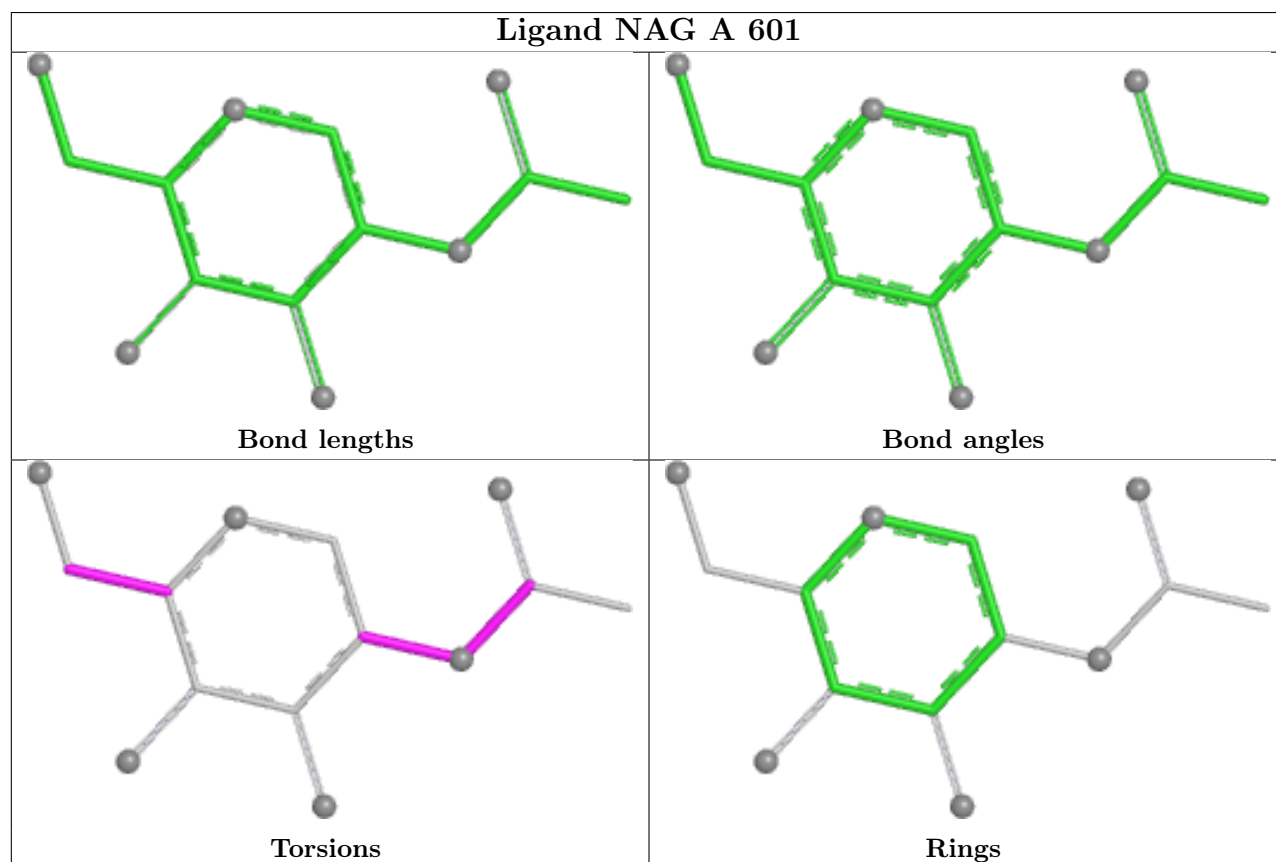
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	601	NAG	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

bond angles, torsion angles, and ring geometry for 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	323/541 (59%)	0.91	31 (9%) 13 9	61, 144, 185, 218	0
1	B	157/541 (29%)	0.43	7 (4%) 38 20	62, 119, 195, 231	0
2	H	221/224 (98%)	0.31	4 (1%) 67 42	62, 100, 132, 224	0
3	L	212/214 (99%)	0.35	5 (2%) 59 35	60, 101, 139, 165	0
All	All	913/1520 (60%)	0.55	47 (5%) 33 19	60, 113, 181, 231	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	60	ILE	4.7
1	A	131	GLU	4.7
1	A	55	ASP	4.6
1	B	66	VAL	4.3
1	A	270	SER	4.1
1	A	153	TRP	4.0
3	L	3	GLN	3.7
1	A	125	PRO	3.5
1	A	84	TRP	3.5
1	A	123	ILE	3.5
1	B	60	ASN	3.4
1	A	52	CYS	3.3
1	A	140	PRO	3.2
1	A	125(b)	SER	3.1
1	A	143	GLY	3.0
1	B	65	ALA	2.9
1	A	47	HIS	2.8
1	A	161	TYR	2.8
1	B	77	ILE	2.7
1	A	135	VAL	2.7
3	L	90	GLN	2.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	141	TYR	2.7
1	B	128	ASN	2.7
1	A	45	ASP	2.6
1	A	139	CYS	2.6
1	A	142	GLN	2.6
1	A	49	GLY	2.6
1	A	132	ALA	2.5
3	L	212	GLY	2.4
3	L	155	GLN	2.3
2	H	155	PRO	2.3
1	A	199	THR	2.3
1	A	261	LYS	2.3
3	L	50	LYS	2.3
1	A	48	ASN	2.3
1	A	144	LYS	2.3
1	A	133(a)	LEU	2.2
1	B	20	GLY	2.2
1	A	295	GLN	2.2
2	H	5	GLN	2.2
2	H	156	GLU	2.2
1	B	132	GLU	2.1
1	A	133	SER	2.0
1	A	51	LEU	2.0
1	A	122	GLN	2.0
1	A	86	TYR	2.0
2	H	140	SER	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.

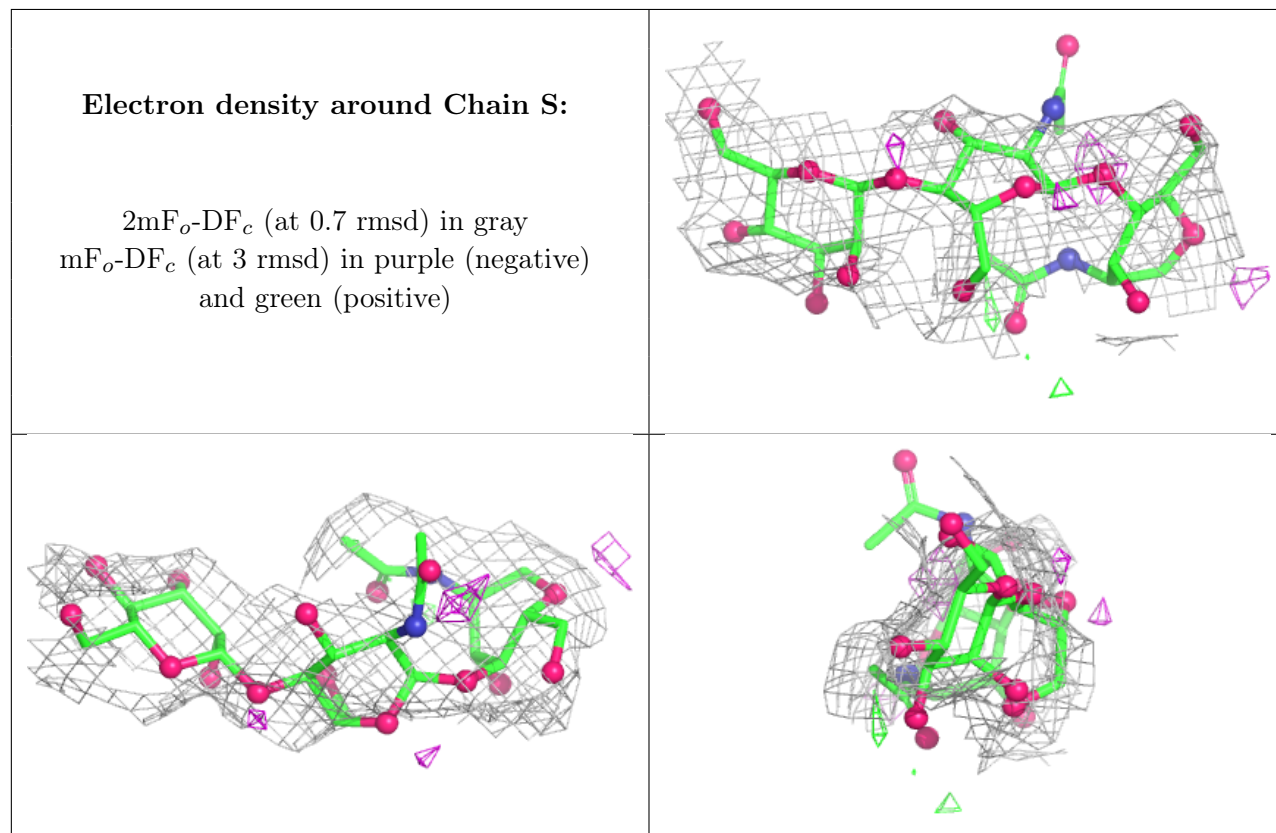
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	NAG	S	1	14/15	-	-	161,187,209,225	0
4	NAG	S	2	14/15	-	-	193,211,214,215	0

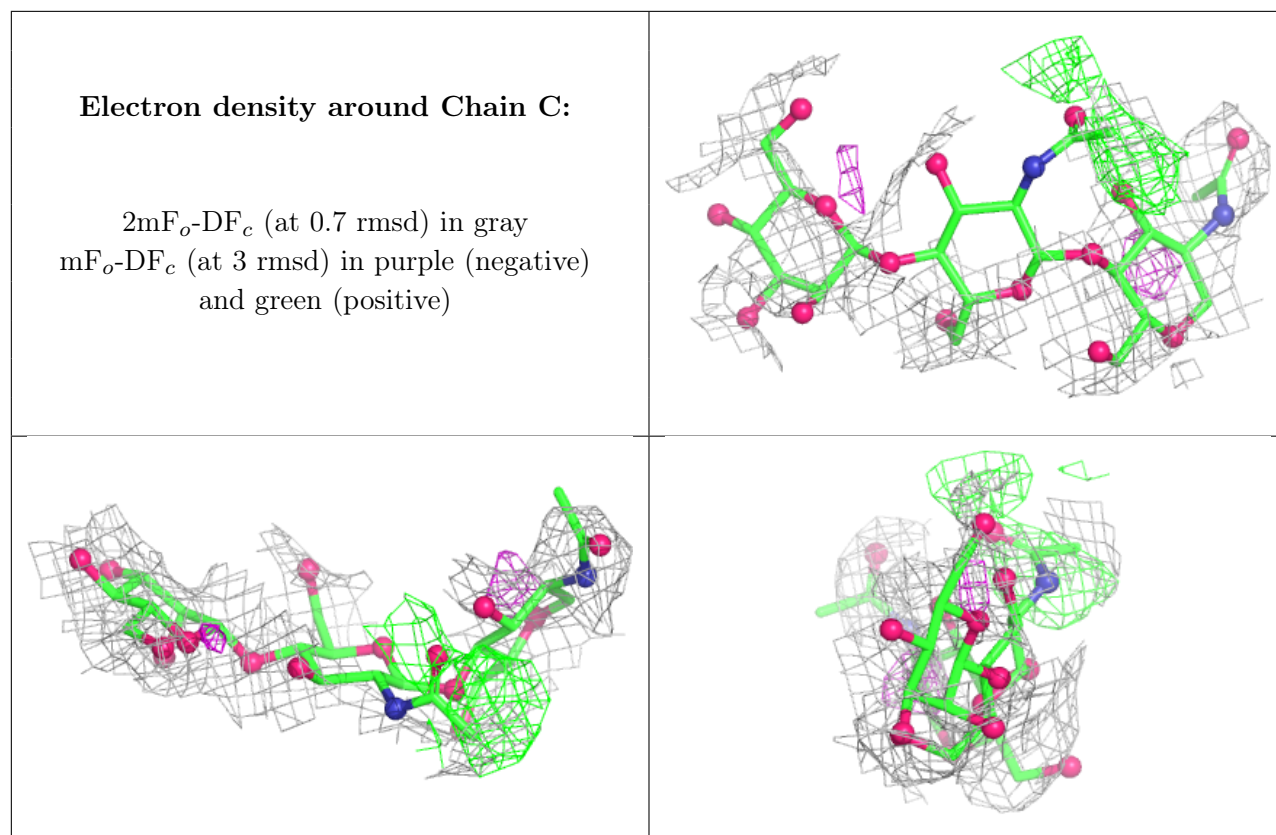
Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	BMA	S	3	11/12	-	-	148,192,201,206	0
4	NAG	C	1	14/15	-	-	178,178,178,178	0
4	NAG	C	2	14/15	-	-	190,190,190,190	0
4	BMA	C	3	11/12	-	-	210,210,210,210	0

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





6.4 Ligands [i](#)

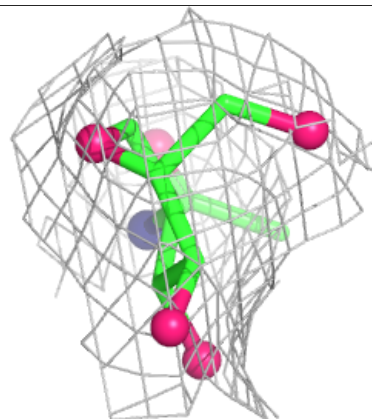
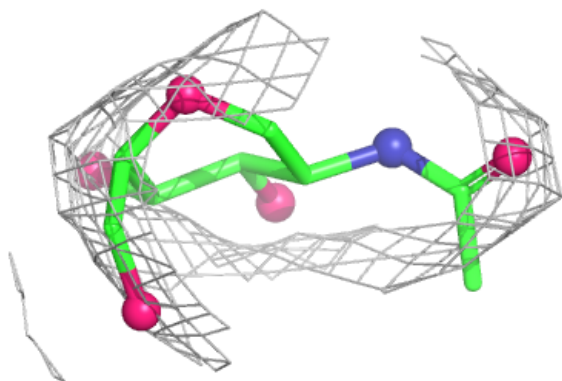
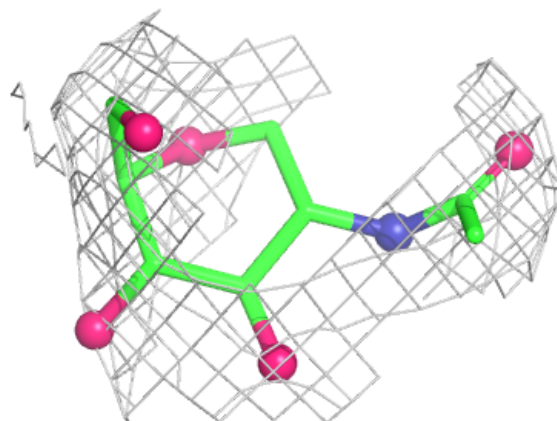
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	NAG	A	601	14/15	0.50	0.14	147,157,163,176	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 A 601:

$2mF_o-DF_c$ (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.