



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 16, 2020 – 06:11 PM JST

PDB ID : 7DE5  
Title : Crystal structure of yak lactoperoxidase at 1.55 Å resolution.  
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Deposited on : 2020-11-02  
Resolution : 1.55 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.14.6  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.14.6

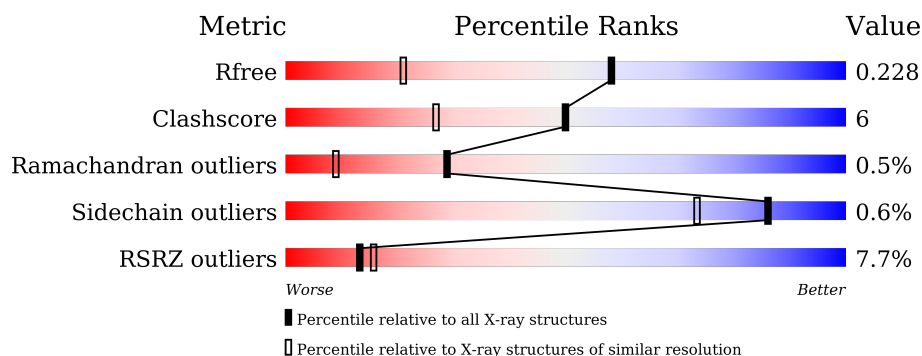
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.55 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	1483 (1.56-1.56)
Clashscore	141614	1529 (1.56-1.56)
Ramachandran outliers	138981	1498 (1.56-1.56)
Sidechain outliers	138945	1495 (1.56-1.56)
RSRZ outliers	127900	1465 (1.56-1.56)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	595	<div> <div>8%</div> <div>89%</div> <div>10%</div> <div>.</div> </div>
2	B	6	<div> <div>83%</div> <div>17%</div> </div>
3	C	2	<div> <div>50%</div> <div>50%</div> </div>

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
6	IOD	A	630	-	-	X	-
6	IOD	A	640	-	-	X	-
6	IOD	A	643	-	-	X	-

## 2 Entry composition [i](#)

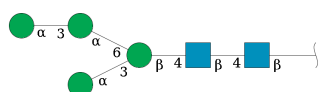
There are 10 unique types of molecules in this entry. The entry contains 5718 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Lactoperoxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	595	Total	C	N	O	S	0	15	0
			4857	3113	852	863	29			

- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



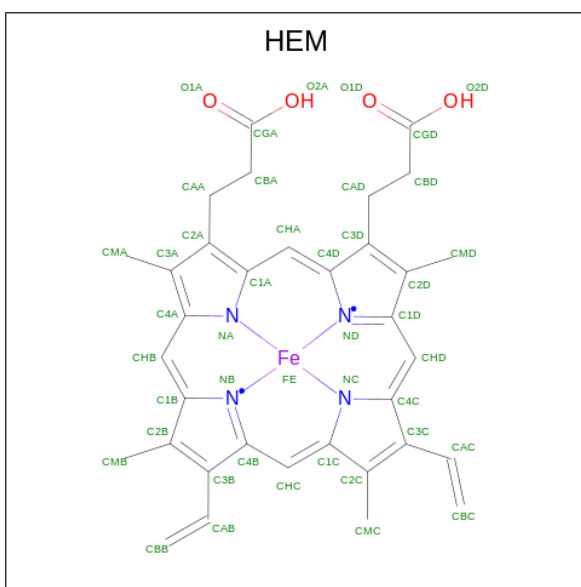
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	6	Total	C	N	O	0	0	0
			72	40	2	30			

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



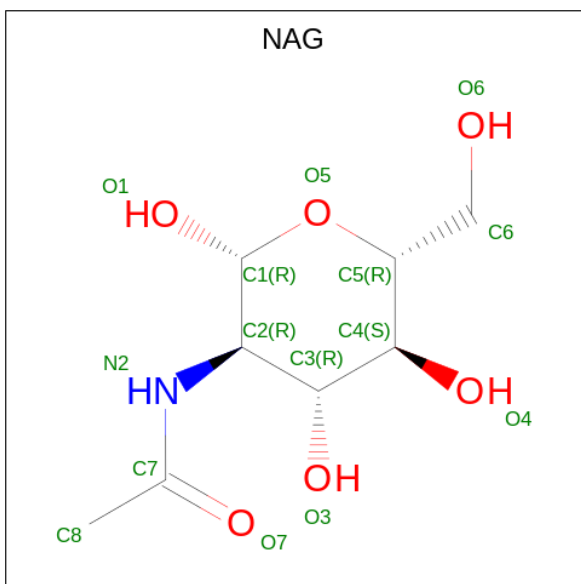
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: C<sub>34</sub>H<sub>32</sub>FeN<sub>4</sub>O<sub>4</sub>) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	Fe	N	O	
			43	34	1	4	4	
							0	0

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

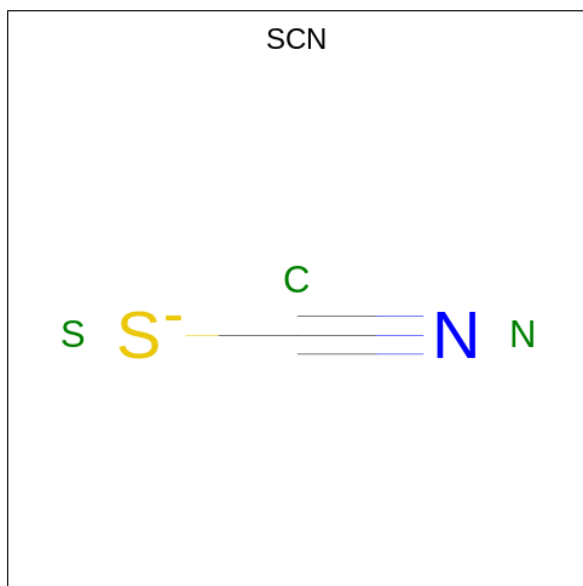


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O		
			14	8	1	5		
							0	0
5	A	1	Total	C	N	O		
			14	8	1	5		
							0	0

- Molecule 6 is IODIDE ION (three-letter code: IOD) (formula: I).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	35	Total I 48 48	0	9

- Molecule 7 is THIOCYANATE ION (three-letter code: SCN) (formula: CNS).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total C N S 3 1 1 1	0	0
7	A	1	Total C N S 3 1 1 1	0	0
7	A	1	Total C N S 3 1 1 1	0	0
7	A	1	Total C N S 3 1 1 1	0	0
7	A	1	Total C N S 3 1 1 1	0	0

- Molecule 8 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total Ca 1 1	0	0

- Molecule 9 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			4	2	2		

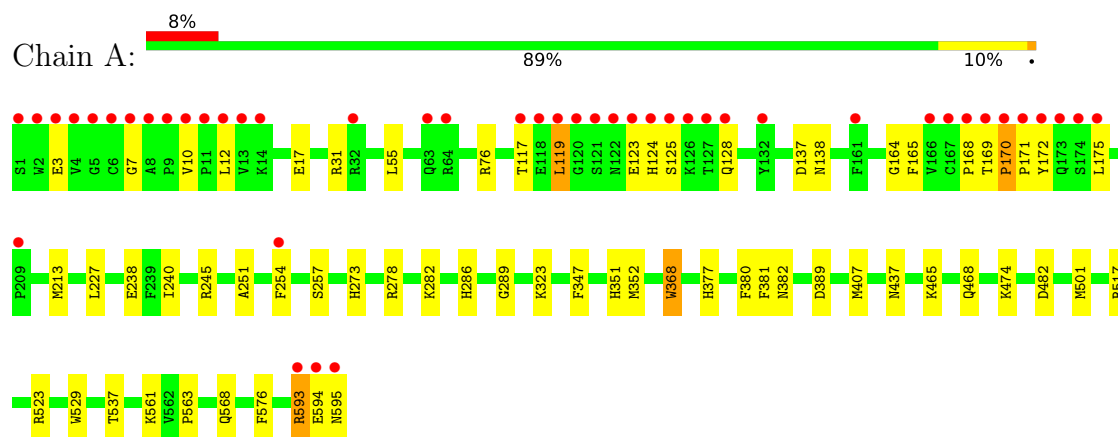
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	622	Total	O	0	0
			622	622		

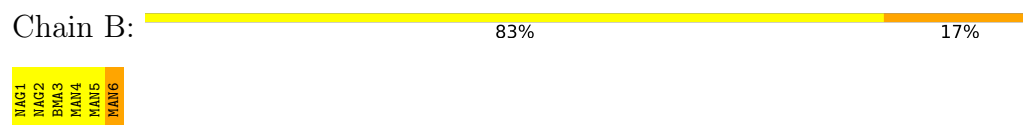
### 3 Residue-property plots

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: Lactoperoxidase



- Molecule 2: alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



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





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	53.91Å 78.98Å 67.82Å 90.00° 92.96° 90.00°	Depositor
Resolution (Å)	43.25 – 1.55 43.25 – 1.55	Depositor EDS
% Data completeness (in resolution range)	98.4 (43.25-1.55) 98.4 (43.25-1.55)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.43 (at 1.55Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, $R_{free}$	0.186 , 0.226 0.195 , 0.228	Depositor DCC
$R_{free}$ test set	1676 reflections (2.07%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	14.0	Xtriage
Anisotropy	0.035	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 43.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.031 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	5718	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.91% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: SCN, NAG, CA, EDO, BMA, HEM, IOD, MAN

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	0.71	0/5033	0.79	2/6826 (0.0%)

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

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	245	ARG	NE-CZ-NH2	-6.37	117.11	120.30
1	A	76	ARG	NE-CZ-NH2	-5.14	117.73	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	165	PHE	Peptide

### 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	4857	0	4837	50	2
2	B	72	0	61	0	1
3	C	28	0	25	0	0
4	A	43	0	30	1	0
5	A	28	0	26	0	0
6	A	48	0	0	13	2
7	A	15	0	0	2	0
8	A	1	0	0	0	0
9	A	4	0	6	0	0
10	A	622	0	0	12	1
All	All	5718	0	4985	58	3

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 (58) 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)
6:A:643:IOD:I	6:A:644:IOD:I	1.53	1.53
6:A:638:IOD:I	10:A:1201:HOH:O	2.21	1.26
1:A:119:LEU:HD12	1:A:138:ASN:HD22	1.33	0.94
1:A:351:HIS:HD1	1:A:437:ASN:HD21	1.15	0.91
1:A:368[A]:TRP:HZ3	1:A:389:ASP:OD1	1.63	0.81
1:A:168:PRO:HB2	1:A:171:PRO:HD2	1.65	0.77
1:A:10:VAL:HG22	1:A:12:LEU:HB2	1.68	0.75
1:A:282:LYS:HE3	10:A:859:HOH:O	1.86	0.74
1:A:238:GLU:OE1	6:A:614:IOD:I	2.79	0.71
1:A:124:HIS:O	1:A:128:GLN:HB2	1.90	0.70
1:A:257:SER:HB2	6:A:636[A]:IOD:I	2.64	0.67
1:A:125:SER:HB3	6:A:620:IOD:I	2.66	0.66
1:A:175:LEU:HG	10:A:716:HOH:O	1.98	0.63
6:A:636[C]:IOD:I	10:A:777:HOH:O	2.86	0.63
1:A:169:THR:HG23	1:A:170:PRO:HD3	1.80	0.61
6:A:610:IOD:I	10:A:1014:HOH:O	2.87	0.60
6:A:635:IOD:I	6:A:640:IOD:I	3.61	0.59
1:A:537:THR:HG21	1:A:593:ARG:NH2	2.18	0.58
1:A:351:HIS:HD1	1:A:437:ASN:ND2	1.95	0.58
1:A:352[A]:MET:SD	1:A:407[A]:MET:SD	3.02	0.57
1:A:169:THR:CG2	1:A:170:PRO:HD3	2.37	0.55
1:A:537:THR:HG21	1:A:593:ARG:HH22	1.72	0.55
1:A:286:HIS:HB2	6:A:612:IOD:I	2.77	0.55
6:A:643:IOD:I	10:A:856:HOH:O	2.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:LEU:O	1:A:172:TYR:OH	2.19	0.54
1:A:561:LYS:HB2	7:A:616:SCN:N	2.23	0.53
4:A:601:HEM:HMC2	4:A:601:HEM:HBC2	1.92	0.52
1:A:3:GLU:N	1:A:3:GLU:OE1	2.43	0.52
1:A:289:GLY:HA3	6:A:631:IOD:I	2.82	0.50
1:A:119:LEU:HD22	1:A:123:GLU:OE2	2.11	0.50
1:A:517:ARG:NE	10:A:704:HOH:O	2.26	0.49
1:A:31:ARG:NH2	10:A:706:HOH:O	2.37	0.48
1:A:377:HIS:HA	1:A:380:PHE:CE2	2.49	0.47
1:A:568:GLN:OE1	10:A:702:HOH:O	2.20	0.47
1:A:368[A]:TRP:CZ3	1:A:389:ASP:OD1	2.55	0.46
1:A:407[A]:MET:HB3	1:A:501:MET:CE	2.45	0.46
1:A:563:PRO:HD3	1:A:576:PHE:CE2	2.51	0.45
1:A:170:PRO:HB2	1:A:171:PRO:HD3	1.98	0.45
1:A:10:VAL:HG13	1:A:10:VAL:O	2.16	0.45
1:A:482:ASP:OD2	10:A:703:HOH:O	2.21	0.44
1:A:117:THR:HG22	1:A:164:GLY:HA2	1.98	0.44
1:A:278:ARG:NH1	10:A:728:HOH:O	2.50	0.44
1:A:468:GLN:HG2	1:A:474:LYS:HA	1.99	0.44
1:A:257:SER:O	1:A:381:PHE:HA	2.18	0.43
1:A:17:GLU:N	7:A:615:SCN:S	2.91	0.43
1:A:7:GLY:O	1:A:10:VAL:N	2.52	0.43
1:A:523:ARG:HG3	1:A:529:TRP:CE2	2.54	0.43
1:A:170:PRO:HD2	1:A:171:PRO:HD3	2.01	0.42
1:A:593:ARG:HD3	1:A:593:ARG:HA	1.74	0.42
1:A:465:LYS:HA	1:A:465:LYS:HD2	1.85	0.42
1:A:594:GLU:O	1:A:595:ASN:HB2	2.20	0.42
1:A:240:ILE:HD11	1:A:382:ASN:HA	2.00	0.41
1:A:170:PRO:CB	1:A:171:PRO:HD3	2.50	0.41
1:A:323:LYS:HA	1:A:323:LYS:HD2	1.75	0.41
6:A:626:IOD:I	10:A:722:HOH:O	2.93	0.41
1:A:254:PHE:CE1	6:A:640:IOD:I	3.45	0.40
1:A:227:LEU:CD2	1:A:251:ALA:HB2	2.51	0.40
1:A:213:MET:HG2	1:A:273:HIS:CD2	2.57	0.40

All (3) 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:323:LYS:NZ	6:A:630:IOD:I[2_645]	1.11	1.09

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:323:LYS:CE	6:A:630:IOD:I[2_645]	1.87	0.33
2:B:6:MAN:O2	10:A:1316:HOH:O[2_555]	2.11	0.09

## 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	608/595 (102%)	585 (96%)	20 (3%)	3 (0%)	29 9

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	137	ASP
1	A	119	LEU
1	A	170	PRO

### 5.3.2 Protein sidechains [i](#)

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	533/518 (103%)	529 (99%)	4 (1%)	81 66

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	347	PHE

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Mol	Chain	Res	Type
1	A	368[A]	TRP
1	A	368[B]	TRP
1	A	593	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	138	ASN
1	A	545	GLN
1	A	568	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 ⓘ

8 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	NAG	B	1	1,2	14,14,15	0.65	0	17,19,21	1.52	3 (17%)
2	NAG	B	2	2	14,14,15	0.45	0	17,19,21	1.24	2 (11%)
2	BMA	B	3	2	11,11,12	0.56	0	15,15,17	1.21	1 (6%)
2	MAN	B	4	2	11,11,12	0.56	0	15,15,17	1.10	1 (6%)
2	MAN	B	5	2	11,11,12	0.58	0	15,15,17	1.07	1 (6%)
2	MAN	B	6	2	11,11,12	0.51	0	15,15,17	1.52	3 (20%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	NAG	C	1	1,3	14,14,15	0.66	0	17,19,21	0.84	0
3	NAG	C	2	3	14,14,15	0.64	0	17,19,21	1.83	5 (29%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	B	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	2	2	-	0/6/23/26	0/1/1/1
2	BMA	B	3	2	-	1/2/19/22	0/1/1/1
2	MAN	B	4	2	-	0/2/19/22	0/1/1/1
2	MAN	B	5	2	-	0/2/19/22	0/1/1/1
2	MAN	B	6	2	-	0/2/19/22	0/1/1/1
3	NAG	C	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	C	2	3	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	2	NAG	C1-O5-C5	4.79	118.68	112.19
2	B	6	MAN	C2-C3-C4	-3.54	104.77	110.89
2	B	3	BMA	C1-O5-C5	3.39	116.78	112.19
2	B	2	NAG	C1-O5-C5	3.37	116.76	112.19
2	B	1	NAG	C1-O5-C5	-3.16	107.91	112.19
2	B	1	NAG	C1-C2-N2	2.83	115.32	110.49
3	C	2	NAG	C3-C4-C5	2.76	115.16	110.24
3	C	2	NAG	O5-C5-C6	2.75	111.52	107.20
2	B	6	MAN	O5-C1-C2	-2.65	106.67	110.77
2	B	1	NAG	C6-C5-C4	-2.61	106.90	113.00
2	B	2	NAG	C1-C2-N2	-2.41	106.37	110.49
3	C	2	NAG	O5-C1-C2	-2.24	107.75	111.29
2	B	4	MAN	O4-C4-C5	2.22	114.81	109.30
2	B	5	MAN	C1-C2-C3	2.11	112.26	109.67
2	B	6	MAN	O3-C3-C4	2.03	115.04	110.35
3	C	2	NAG	O4-C4-C3	-2.01	105.70	110.35

There are no chirality outliers.

All (1) torsion outliers are listed below:

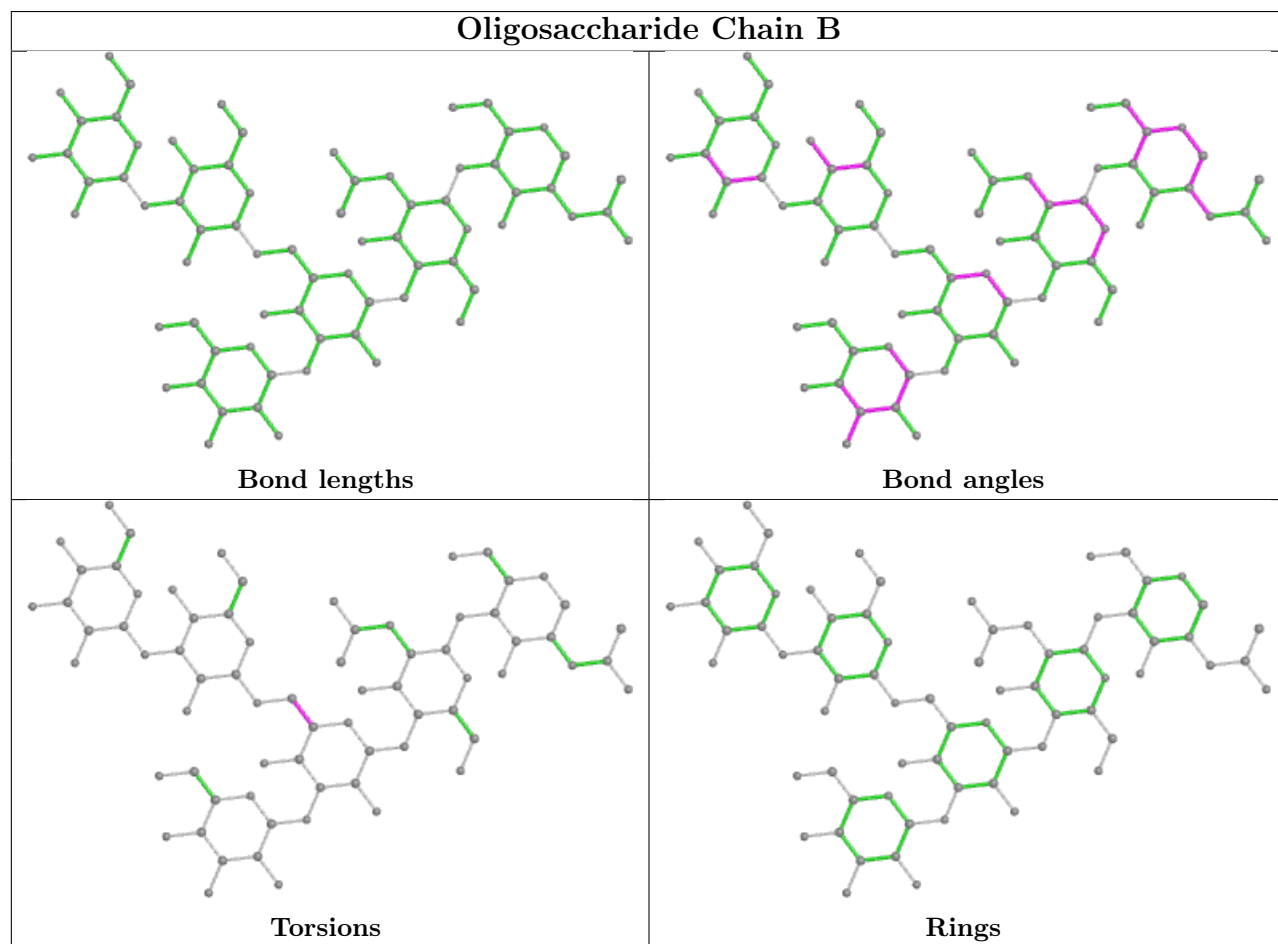
Mol	Chain	Res	Type	Atoms
2	B	3	BMA	C4-C5-C6-O6

There are no ring outliers.

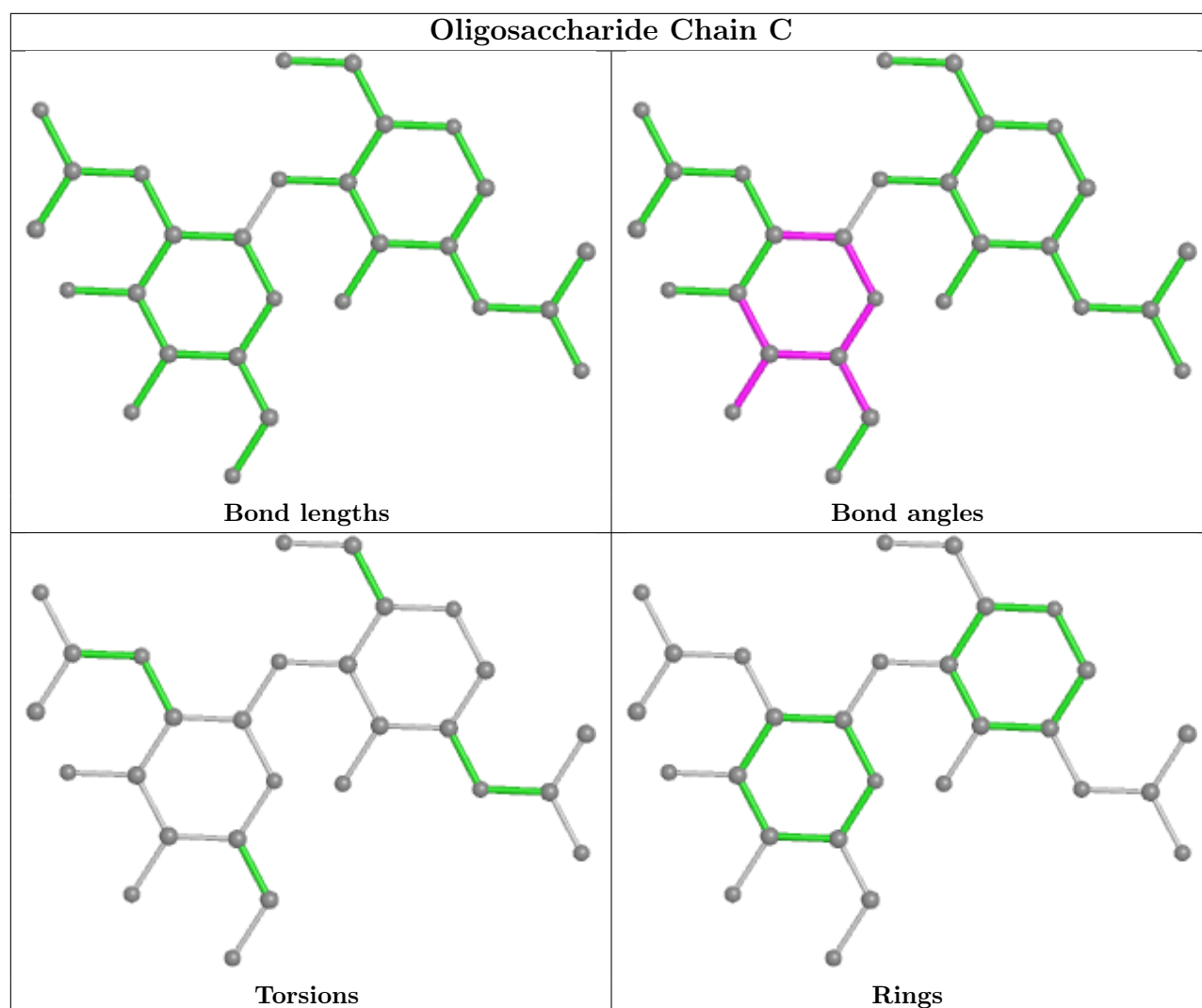
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	6	MAN	0	1

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 ⓘ

Of 58 ligands modelled in this entry, 49 are monoatomic - leaving 9 for Mogul analysis.

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	HEM	A	601	1,10	27,50,50	1.05	2 (7%)	17,82,82	1.40	3 (17%)
5	NAG	A	603	1	14,14,15	0.79	0	17,19,21	1.33	2 (11%)
7	SCN	A	617	-	1,2,2	0.96	0	0,1,1	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	NAG	A	602	1	14,14,15	0.53	0	17,19,21	1.45	2 (11%)
7	SCN	A	615	-	1,2,2	2.52	1 (100%)	0,1,1	0.00	-
9	EDO	A	642	-	3,3,3	0.10	0	2,2,2	0.63	0
7	SCN	A	641	-	1,2,2	0.10	0	0,1,1	0.00	-
7	SCN	A	618	-	1,2,2	1.51	0	0,1,1	0.00	-
7	SCN	A	616	-	1,2,2	0.52	0	0,1,1	0.00	-

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	HEM	A	601	1,10	-	0/6/54/54	-
9	EDO	A	642	-	-	0/1/1/1	-
5	NAG	A	603	1	-	2/6/23/26	0/1/1/1
5	NAG	A	602	1	-	0/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	601	HEM	C4A-NA	2.59	1.41	1.36
7	A	615	SCN	C-N	2.52	1.23	1.15
4	A	601	HEM	C4D-C3D	2.27	1.47	1.42

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	603	NAG	C1-O5-C5	3.46	116.88	112.19
5	A	602	NAG	O5-C1-C2	-3.03	106.50	111.29
4	A	601	HEM	C1D-C2D-C3D	-2.51	105.25	107.00
4	A	601	HEM	C4C-C3C-C2C	-2.18	105.38	106.90
4	A	601	HEM	CBD-CAD-C3D	-2.16	108.50	112.48
5	A	603	NAG	C4-C3-C2	-2.11	107.92	111.02
5	A	602	NAG	C1-C2-N2	-2.11	106.89	110.49

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	603	NAG	C8-C7-N2-C2

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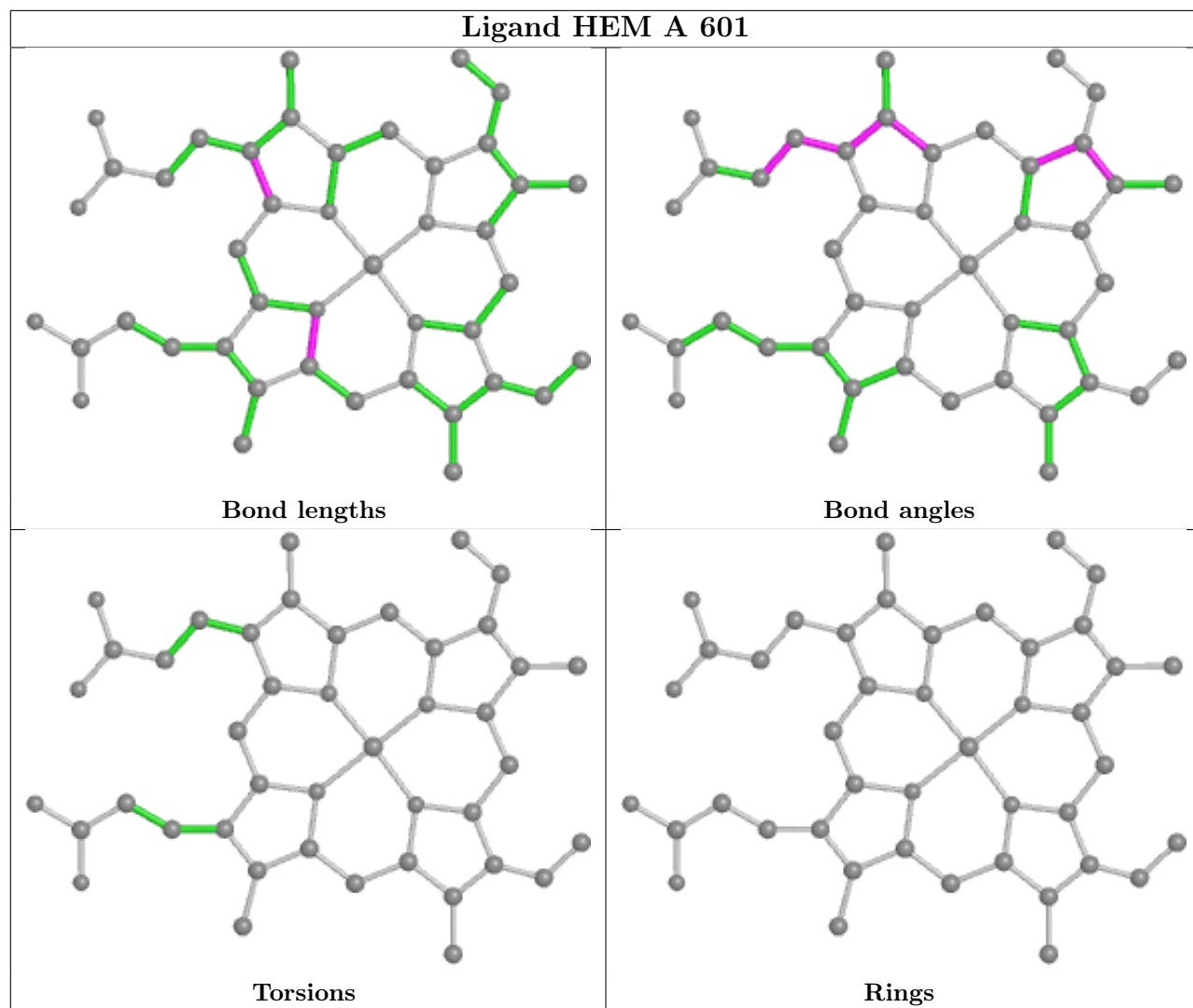
Mol	Chain	Res	Type	Atoms
5	A	603	NAG	O7-C7-N2-C2

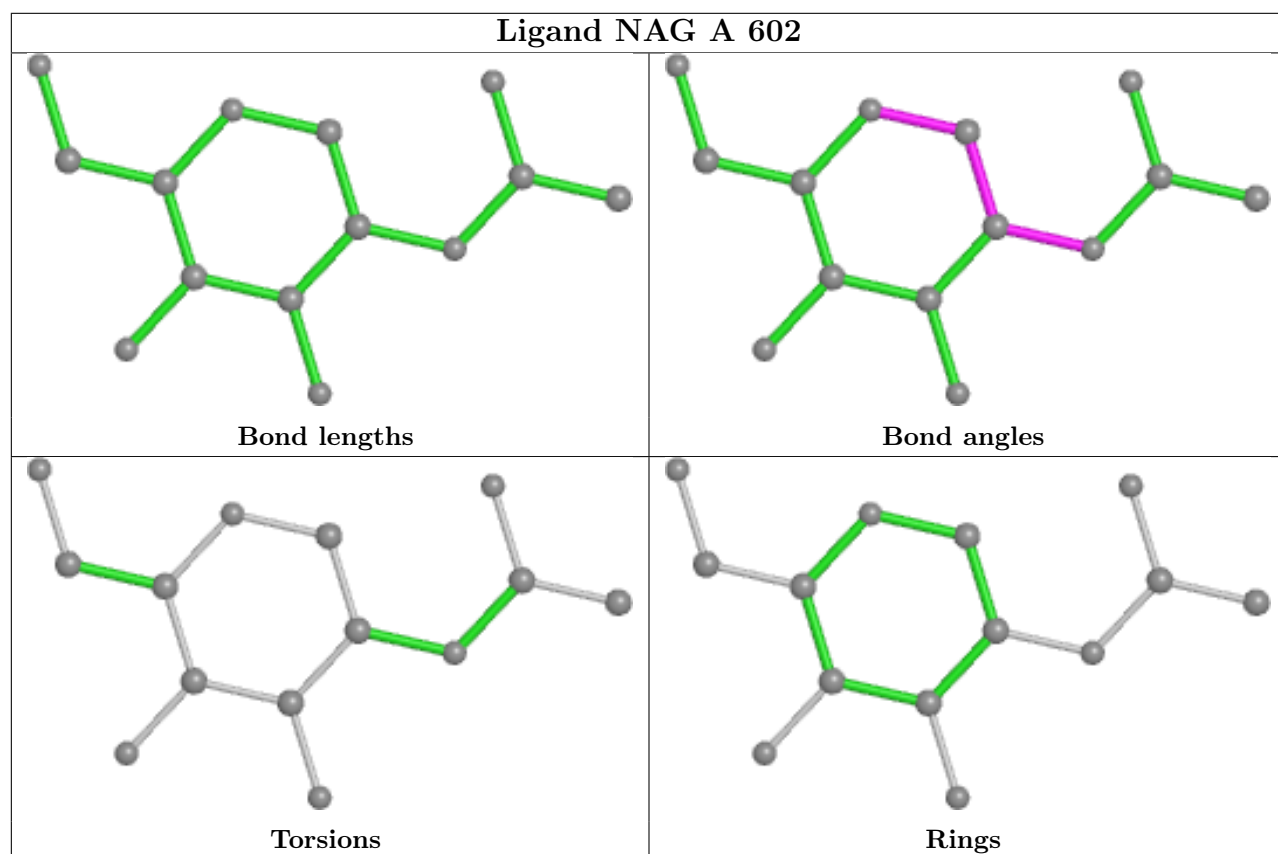
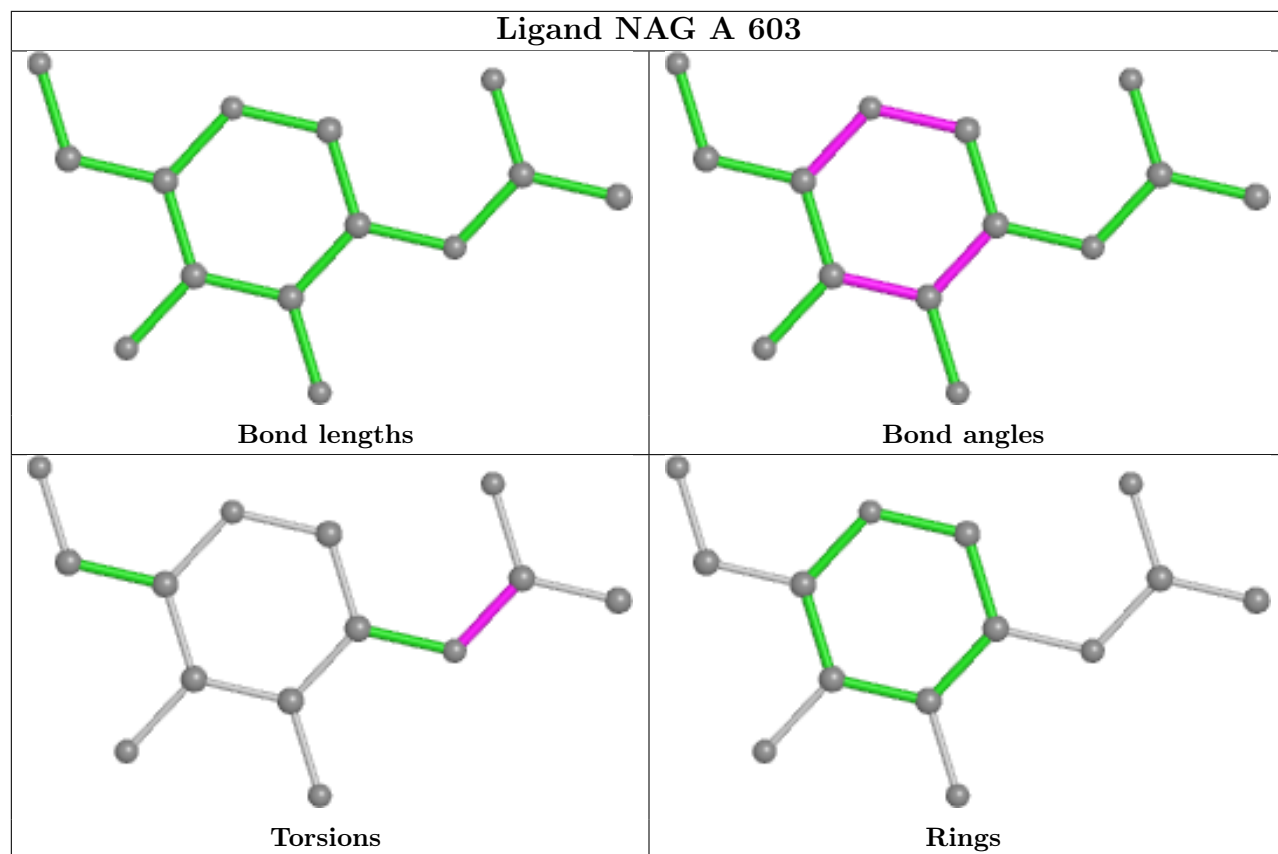
There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	601	HEM	1	0
7	A	615	SCN	1	0
7	A	616	SCN	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	595/595 (100%)	0.55	46 (7%) 13 16	8, 15, 60, 126	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	TRP	38.4
1	A	1	SER	27.2
1	A	120	GLY	21.3
1	A	174	SER	18.9
1	A	122	ASN	17.0
1	A	121	SER	15.0
1	A	119	LEU	14.6
1	A	11	PRO	14.4
1	A	12	LEU	13.8
1	A	7	GLY	12.5
1	A	124	HIS	11.3
1	A	172	TYR	11.1
1	A	171	PRO	9.2
1	A	173	GLN	8.9
1	A	13	VAL	8.8
1	A	123	GLU	8.5
1	A	8	ALA	7.4
1	A	170	PRO	7.4
1	A	169	THR	7.1
1	A	10	VAL	7.0
1	A	9	PRO	6.9
1	A	594	GLU	5.8
1	A	595	ASN	5.5
1	A	5	GLY	5.4
1	A	3	GLU	5.4
1	A	128	GLN	5.2
1	A	175	LEU	5.0

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Mol	Chain	Res	Type	RSRZ
1	A	4	VAL	4.9
1	A	166	VAL	4.5
1	A	6	CYS	4.3
1	A	168	PRO	3.9
1	A	127	THR	3.9
1	A	167	CYS	3.9
1	A	161[A]	PHE	3.8
1	A	14	LYS	3.7
1	A	126	LYS	3.6
1	A	118	GLU	3.6
1	A	254	PHE	3.3
1	A	117	THR	3.1
1	A	132	TYR	3.1
1	A	64	ARG	3.0
1	A	593	ARG	2.7
1	A	209	PRO	2.5
1	A	63	GLN	2.4
1	A	32	ARG	2.3
1	A	125	SER	2.1

## 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, 95<sup>th</sup> 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.

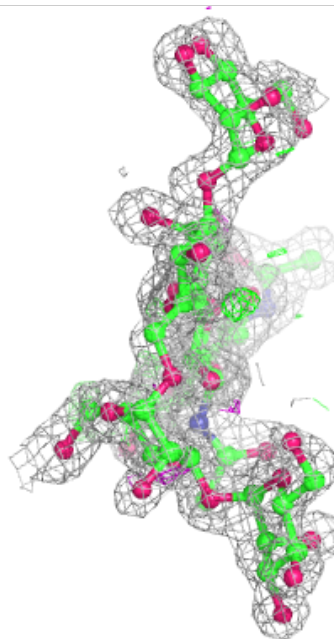
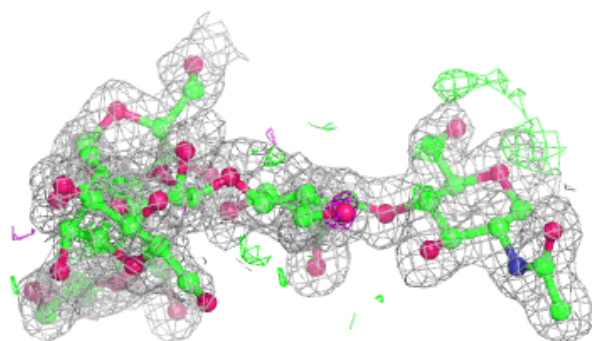
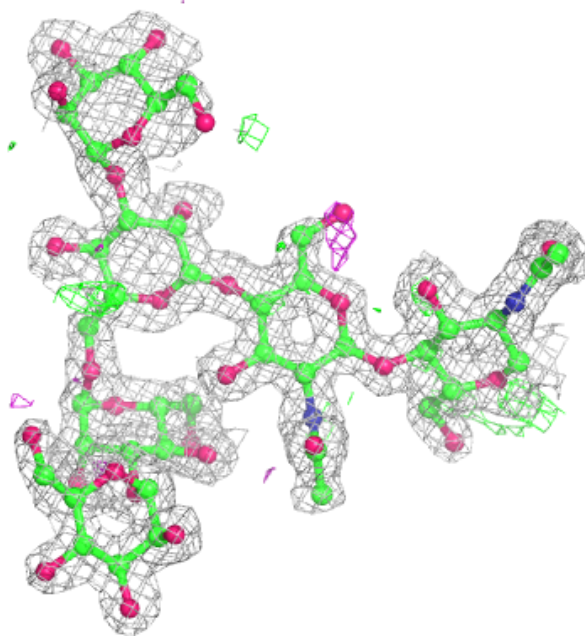
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	NAG	C	2	14/15	0.77	0.21	36,41,47,48	0
2	MAN	B	6	11/12	0.86	0.16	34,37,41,43	0
2	BMA	B	3	11/12	0.87	0.15	24,26,27,28	0
2	NAG	B	2	14/15	0.88	0.15	20,24,28,37	0
2	MAN	B	4	11/12	0.90	0.14	21,23,25,26	0
2	MAN	B	5	11/12	0.90	0.18	25,28,32,33	0
2	NAG	B	1	14/15	0.90	0.09	17,19,24,29	0
3	NAG	C	1	14/15	0.96	0.07	16,19,24,25	0

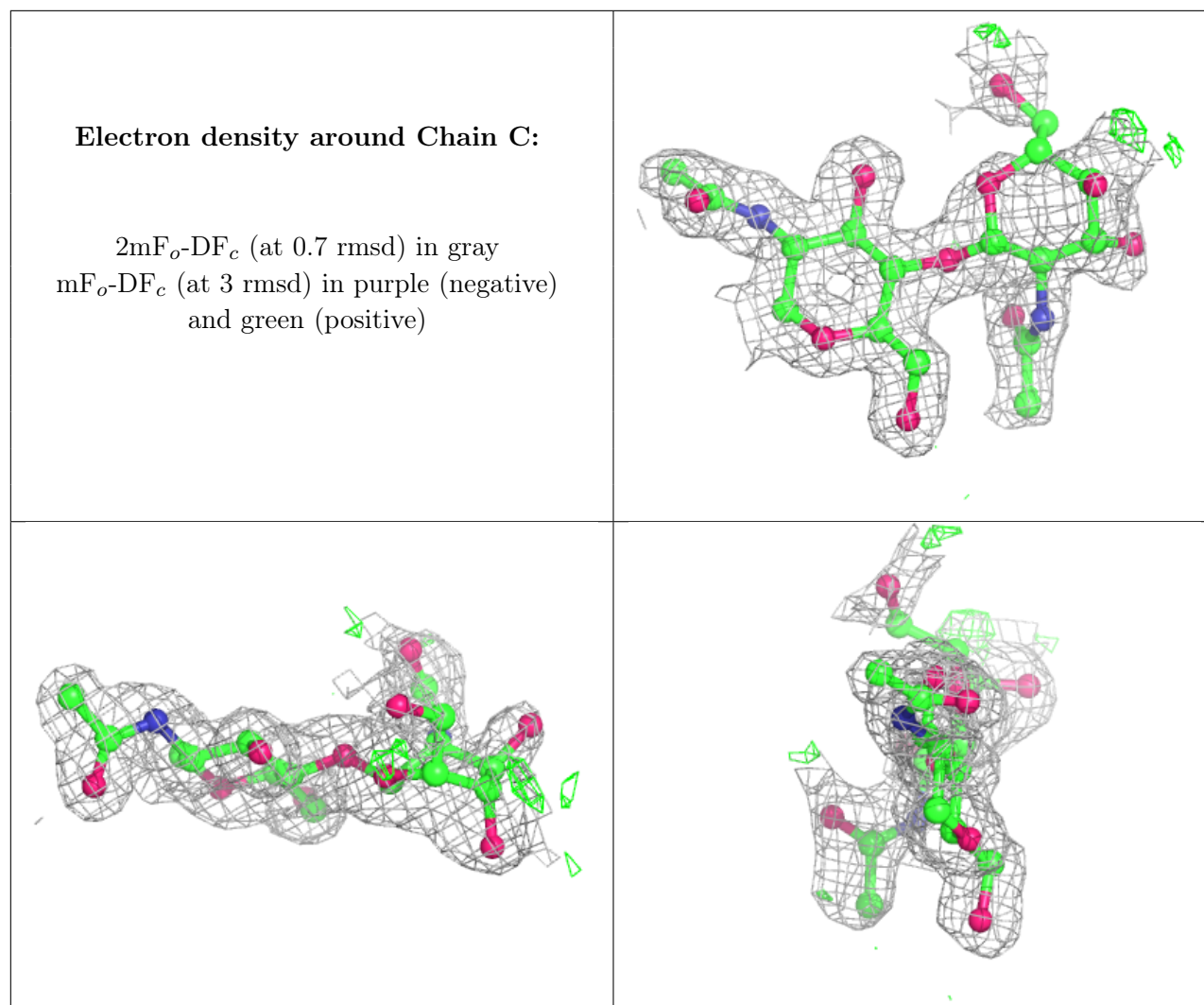


The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

**Electron density around Chain B:**

$2mF_o-DF_c$  (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, 95<sup>th</sup> 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	A	603	14/15	0.79	0.16	27,31,36,36	0
6	IOD	A	632	1/1	0.79	0.14	62,62,62,62	1
5	NAG	A	602	14/15	0.83	0.20	28,33,37,39	0
6	IOD	A	637	1/1	0.84	0.11	55,55,55,55	1
7	SCN	A	617	3/3	0.87	0.14	40,40,45,52	0
9	EDO	A	642	4/4	0.89	0.10	27,28,31,32	0
6	IOD	A	639[B]	1/1	0.91	0.12	61,61,61,61	1
6	IOD	A	639[A]	1/1	0.91	0.12	61,61,61,61	1

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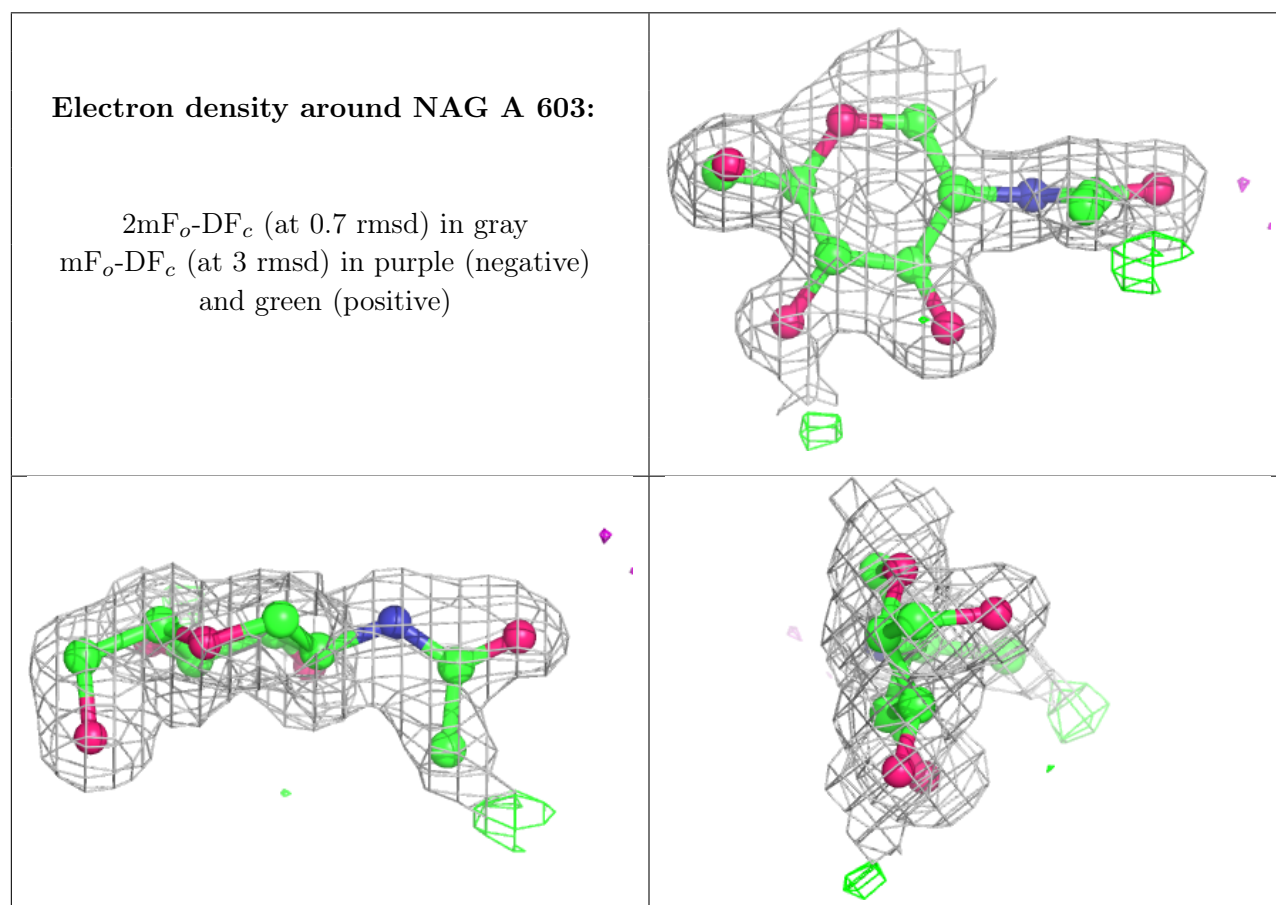
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
6	IOD	A	645	1/1	0.91	0.14	34,34,34,34	1
7	SCN	A	615	3/3	0.92	0.13	29,29,40,45	0
6	IOD	A	638	1/1	0.92	0.07	32,32,32,32	1
6	IOD	A	620	1/1	0.93	0.07	50,50,50,50	1
6	IOD	A	630	1/1	0.93	0.15	63,63,63,63	1
6	IOD	A	631	1/1	0.94	0.07	25,25,25,25	1
6	IOD	A	623	1/1	0.96	0.05	32,32,32,32	1
6	IOD	A	622	1/1	0.96	0.06	31,31,31,31	1
6	IOD	A	640	1/1	0.96	0.06	39,39,39,39	1
6	IOD	A	636[A]	1/1	0.97	0.05	30,30,30,30	1
7	SCN	A	618	3/3	0.97	0.08	26,26,29,35	0
6	IOD	A	636[B]	1/1	0.97	0.05	34,34,34,34	1
6	IOD	A	636[C]	1/1	0.97	0.05	46,46,46,46	1
6	IOD	A	613	1/1	0.97	0.04	23,23,23,23	1
6	IOD	A	621	1/1	0.98	0.04	26,26,26,26	1
7	SCN	A	641	3/3	0.98	0.07	14,14,16,22	0
6	IOD	A	628	1/1	0.98	0.03	27,27,27,27	1
6	IOD	A	635	1/1	0.98	0.04	31,31,31,31	1
6	IOD	A	643	1/1	0.98	0.04	20,20,20,20	1
6	IOD	A	612	1/1	0.98	0.03	22,22,22,22	1
4	HEM	A	601	43/43	0.98	0.10	8,10,12,13	0
7	SCN	A	616	3/3	0.98	0.09	22,22,24,29	0
6	IOD	A	627[B]	1/1	0.98	0.09	37,37,37,37	1
6	IOD	A	627[C]	1/1	0.98	0.09	41,41,41,41	1
6	IOD	A	607	1/1	0.98	0.04	29,29,29,29	1
6	IOD	A	626	1/1	0.98	0.04	21,21,21,21	1
6	IOD	A	633	1/1	0.98	0.03	33,33,33,33	1
6	IOD	A	625[B]	1/1	0.99	0.03	21,21,21,21	1
6	IOD	A	629[B]	1/1	0.99	0.03	25,25,25,25	1
6	IOD	A	608[C]	1/1	0.99	0.03	21,21,21,21	1
6	IOD	A	611[C]	1/1	0.99	0.03	20,20,20,20	1
6	IOD	A	608[B]	1/1	0.99	0.03	22,22,22,22	1
6	IOD	A	624[A]	1/1	0.99	0.03	26,26,26,26	1
6	IOD	A	608[A]	1/1	0.99	0.03	18,18,18,18	1
6	IOD	A	614	1/1	0.99	0.06	42,42,42,42	1
6	IOD	A	625[A]	1/1	0.99	0.03	21,21,21,21	1
6	IOD	A	624[B]	1/1	0.99	0.03	26,26,26,26	1
6	IOD	A	644	1/1	0.99	0.07	29,29,29,29	1
6	IOD	A	634	1/1	0.99	0.03	24,24,24,24	1
6	IOD	A	629[A]	1/1	0.99	0.03	24,24,24,24	1
6	IOD	A	609	1/1	0.99	0.04	19,19,19,19	1
6	IOD	A	611[A]	1/1	0.99	0.03	18,18,18,18	1

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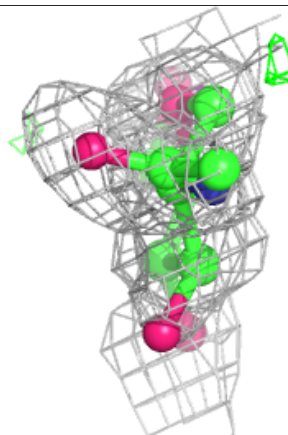
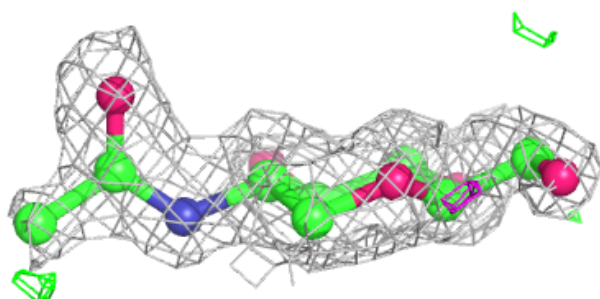
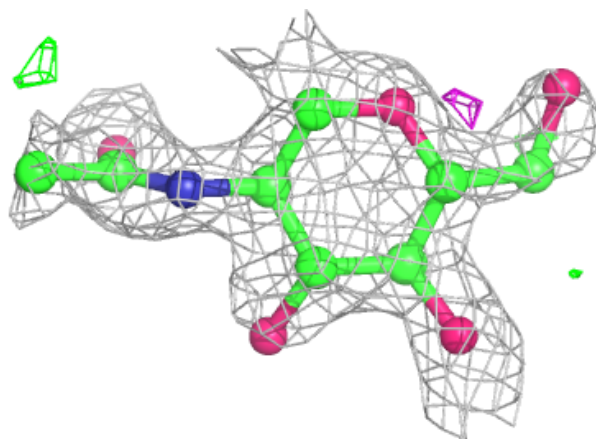
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
6	IOD	A	605	1/1	0.99	0.02	24,24,24,24	1
6	IOD	A	611[B]	1/1	0.99	0.03	22,22,22,22	1
6	IOD	A	606	1/1	0.99	0.04	23,23,23,23	1
6	IOD	A	604[A]	1/1	1.00	0.02	15,15,15,15	1
6	IOD	A	610	1/1	1.00	0.03	19,19,19,19	1
6	IOD	A	604[B]	1/1	1.00	0.02	9,9,9,9	1
8	CA	A	619	1/1	1.00	0.04	9,9,9,9	0
6	IOD	A	604[C]	1/1	1.00	0.02	12,12,12,12	1

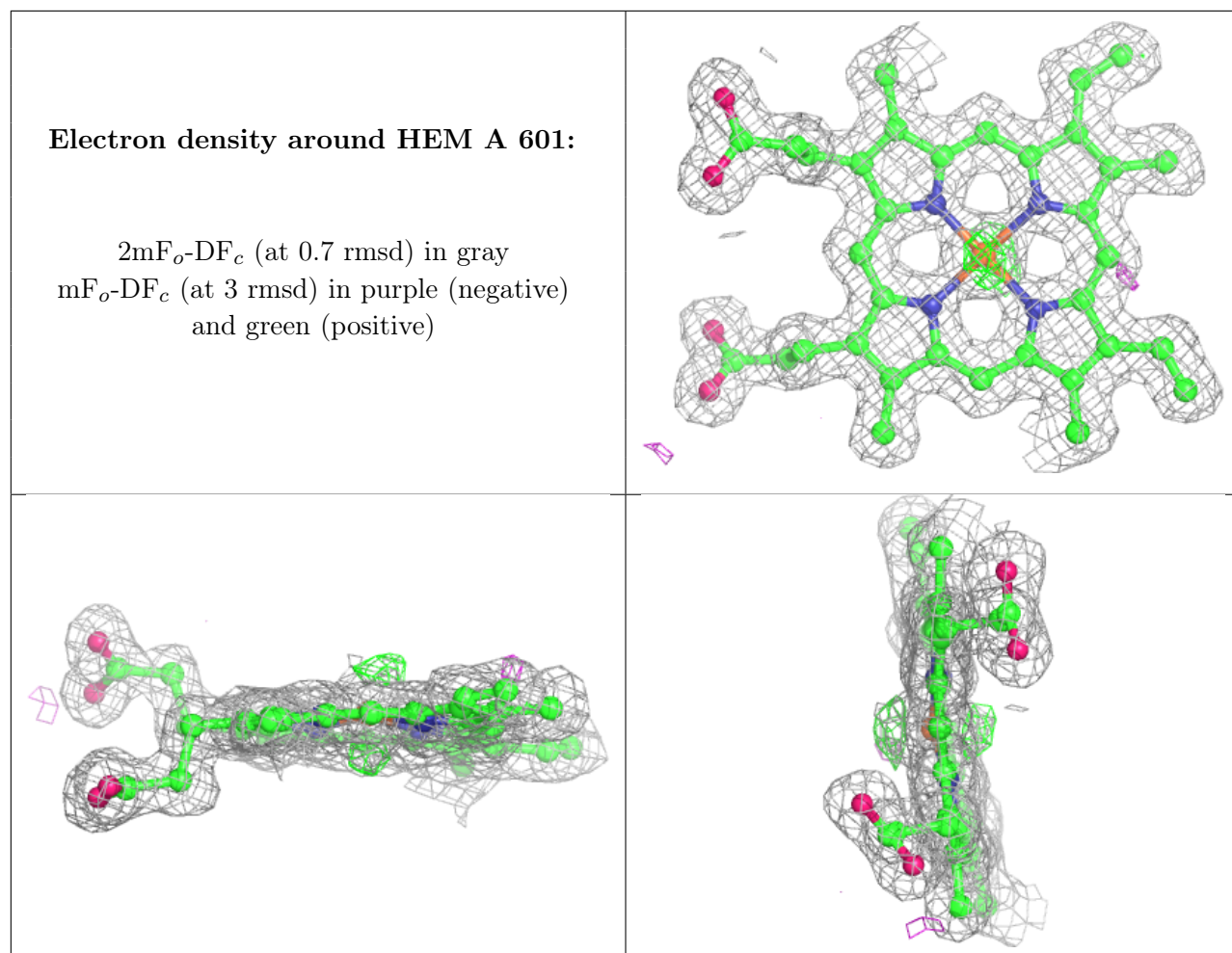
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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



**Electron density around NAG A 602:**

$2mF_o-DF_c$  (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.