



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:54 PM UTC

PDB ID : 9SS7 / pdb\_00009ss7  
Title : Human angiotensin 1-converting enzyme C-domain in complex with rentiapril  
Authors : Gregory, K.S.; Acharya, K.R.  
Deposited on : 2025-09-25  
Resolution : 1.95 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.

The types of validation reports are described at

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

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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

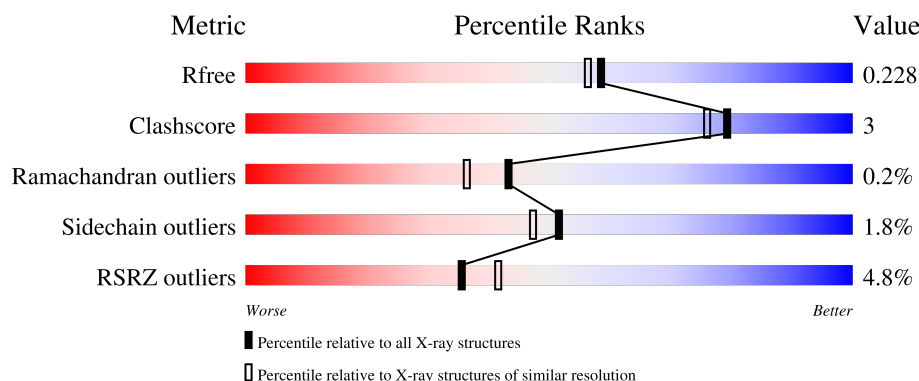
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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  $\geq 3$ , 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	584	<div> <div>5%</div> <div> <div></div> <div>90%</div> <div>9%</div> </div> </div>
2	B	4	<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	MLI	A	708	-	-	X	-

## 2 Entry composition [i](#)

There are 11 unique types of molecules in this entry. The entry contains 5062 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.

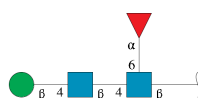
- Molecule 1 is a protein called Angiotensin-converting enzyme, soluble form.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	580	Total	C	N	O	S	0	2	0
			4752	3049	811	868	24			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	64	GLY	GLU	conflict	UNP P12821
A	90	GLN	ASN	conflict	UNP P12821
A	155	GLN	ASN	conflict	UNP P12821
A	337	GLN	ASN	conflict	UNP P12821
A	586	GLN	ASN	conflict	UNP P12821

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	4	Total	C	N	O	0	0	0
			49	28	2	19			

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).

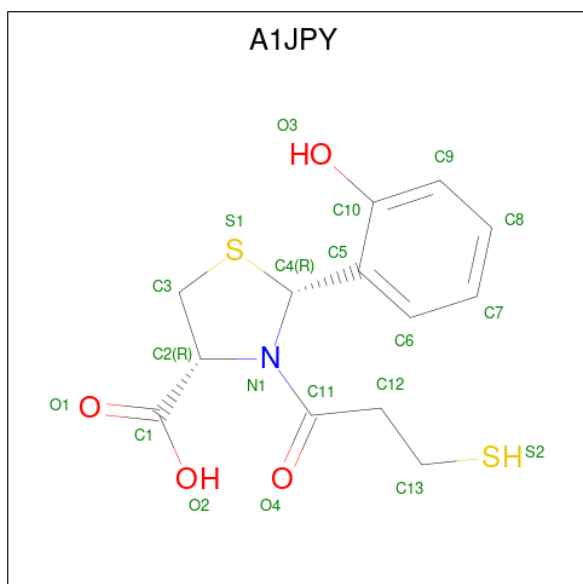


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

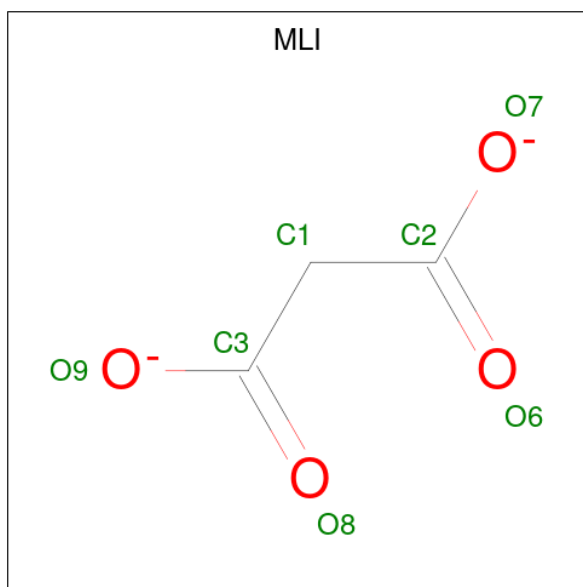
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn	0	0
			1	1		

- Molecule 5 is Rentiapril (CCD ID: A1JPY) (formula: C<sub>13</sub>H<sub>15</sub>NO<sub>4</sub>S<sub>2</sub>) (labeled as "Ligand of Interest" by depositor).



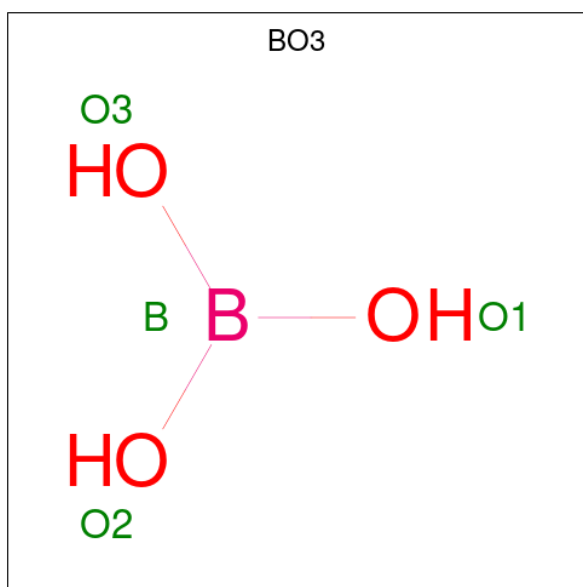
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	S	0	0
			20	13	1	4	2		
5	A	1	Total	C	N	O	S	0	0
			20	13	1	4	2		

- Molecule 6 is MALONATE ION (CCD ID: MLI) (formula:  $C_3H_2O_4$ ).



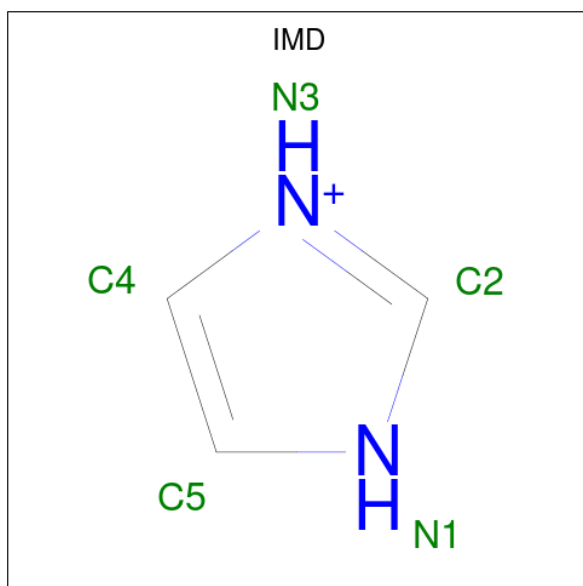
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			7	3	4		
6	A	1	Total	C	O	0	0
			7	3	4		

- Molecule 7 is BORIC ACID (CCD ID: BO3) (formula:  $BH_3O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	B	O	0	0
			4	1	3		

- Molecule 8 is IMIDAZOLE (CCD ID: IMD) (formula:  $C_3H_5N_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	N	0	0
			5	3	2		

- Molecule 9 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula:  $C_2H_6O_2$ ).



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

- Molecule 10 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	2	Total	Cl	0	0
			2	2		

- Molecule 11 is water.

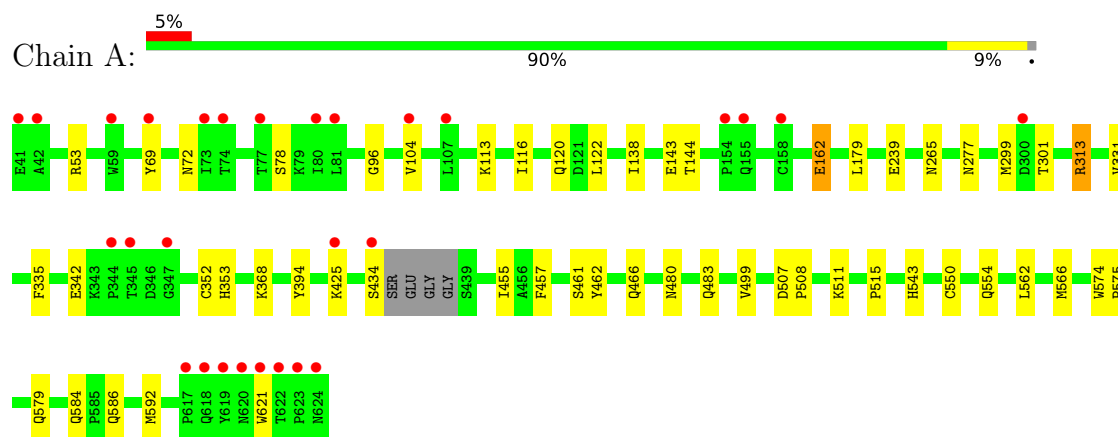
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	177	Total	O	0	0
			177	177		



### 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: Angiotensin-converting enzyme, soluble form



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



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	56.73Å 84.94Å 134.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	71.86 – 1.95 71.86 – 1.95	Depositor EDS
% Data completeness (in resolution range)	100.0 (71.86-1.95) 99.9 (71.86-1.95)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.30 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.8.0431 (refmacat 0.4.105)	Depositor
R, $R_{free}$	0.186 , 0.223 0.194 , 0.228	Depositor DCC
$R_{free}$ test set	2438 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	27.0	Xtriage
Anisotropy	0.115	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 31.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5062	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.56% 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: NAG, A1JPY, BMA, MLI, ZN, EDO, IMD, BO3, CL, FUC

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	0.59	0/4892	1.07	6/6651 (0.1%)

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	2

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	301	THR	CA-CB-OG1	-7.56	98.26	109.60
1	A	143	GLU	N-CA-CB	7.28	120.94	110.16
1	A	162	GLU	CB-CG-CD	5.34	121.68	112.60
1	A	579	GLN	CB-CA-C	-5.33	101.95	110.79
1	A	331	VAL	N-CA-CB	5.26	115.41	109.99
1	A	584	GLN	CB-CA-C	5.07	117.83	109.97

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	313	ARG	Sidechain
1	A	53	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	4752	0	4590	26	0
2	B	49	0	43	0	0
3	A	14	0	13	0	0
4	A	1	0	0	0	0
5	A	40	0	0	1	0
6	A	14	0	4	2	0
7	A	4	0	3	0	0
8	A	5	0	5	0	0
9	A	4	0	6	0	0
10	A	2	0	0	0	0
11	A	177	0	0	3	0
All	All	5062	0	4664	26	0

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

All (26) 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:353:HIS:HE1	6:A:708:MLI:O9	1.59	0.84
1:A:554:GLN:NE2	11:A:802:HOH:O	2.21	0.71
1:A:353:HIS:CE1	6:A:708:MLI:O9	2.46	0.66
1:A:299:MET:HE1	1:A:434:SER:OG	1.99	0.62
1:A:353:HIS:HD2	11:A:809:HOH:O	1.83	0.62
1:A:96:GLY:HA3	1:A:122:LEU:CD2	2.32	0.59
1:A:239:GLU:OE1	11:A:801:HOH:O	2.17	0.57
1:A:277:ASN:ND2	5:A:707:A1JPY:S1	2.80	0.54
1:A:96:GLY:HA3	1:A:122:LEU:HD22	1.89	0.54
1:A:511:LYS:O	1:A:515:PRO:HD2	2.08	0.53
1:A:574:TRP:N	1:A:575:PRO:CD	2.74	0.51
1:A:562:LEU:HB3	1:A:566:MET:HE2	1.94	0.50
1:A:69:TYR:HH	1:A:144:THR:HG1	1.61	0.47
1:A:507:ASP:N	1:A:508:PRO:CD	2.77	0.47
1:A:457:PHE:CE2	1:A:461:SER:HB3	2.50	0.47
1:A:116:ILE:O	1:A:120:GLN:HG3	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:104:VAL:HG13	1:A:113:LYS:HG2	1.99	0.44
1:A:299:MET:HE3	1:A:299:MET:HB2	1.75	0.44
1:A:179:LEU:HD11	1:A:499:VAL:HG23	2.01	0.43
1:A:455:ILE:HD13	1:A:592:MET:HE2	2.00	0.43
1:A:462:TYR:O	1:A:466:GLN:HG2	2.19	0.43
1:A:480:ASN:HA	1:A:483:GLN:OE1	2.18	0.42
1:A:335:PHE:CD1	1:A:335:PHE:C	2.96	0.42
1:A:265:ASN:HD22	1:A:621:TRP:CG	2.37	0.42
1:A:543:HIS:CD2	1:A:550:CYS:HB2	2.55	0.41
1:A:353:HIS:O	1:A:368:LYS:NZ	2.54	0.41

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	578/584 (99%)	567 (98%)	10 (2%)	1 (0%)	43	36

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	72	ASN

### 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	509/509 (100%)	500 (98%)	9 (2%)	51	47

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	78	SER
1	A	138	ILE
1	A	162	GLU
1	A	313	ARG
1	A	342	GLU
1	A	352	CYS
1	A	394	TYR
1	A	425	LYS
1	A	586	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	106	GLN
1	A	160	GLN
1	A	167	ASN
1	A	353	HIS
1	A	618	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 ⓘ

4 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	2,1	14,14,15	0.32	0	17,19,21	1.16	2 (11%)
2	NAG	B	2	2	14,14,15	0.36	0	17,19,21	0.87	0
2	BMA	B	3	2	11,11,12	0.49	0	15,15,17	1.16	2 (13%)
2	FUC	B	4	2	10,10,11	0.37	0	14,14,16	0.38	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
2	NAG	B	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	B	2	2	-	2/6/23/26	0/1/1/1
2	BMA	B	3	2	-	1/2/19/22	0/1/1/1
2	FUC	B	4	2	-	-	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	NAG	C1-C2-N2	3.60	116.11	110.43
2	B	3	BMA	C1-O5-C5	2.61	115.68	112.19
2	B	3	BMA	C2-C3-C4	-2.24	106.92	110.86
2	B	1	NAG	O4-C4-C3	-2.12	105.38	110.38

There are no chirality outliers.

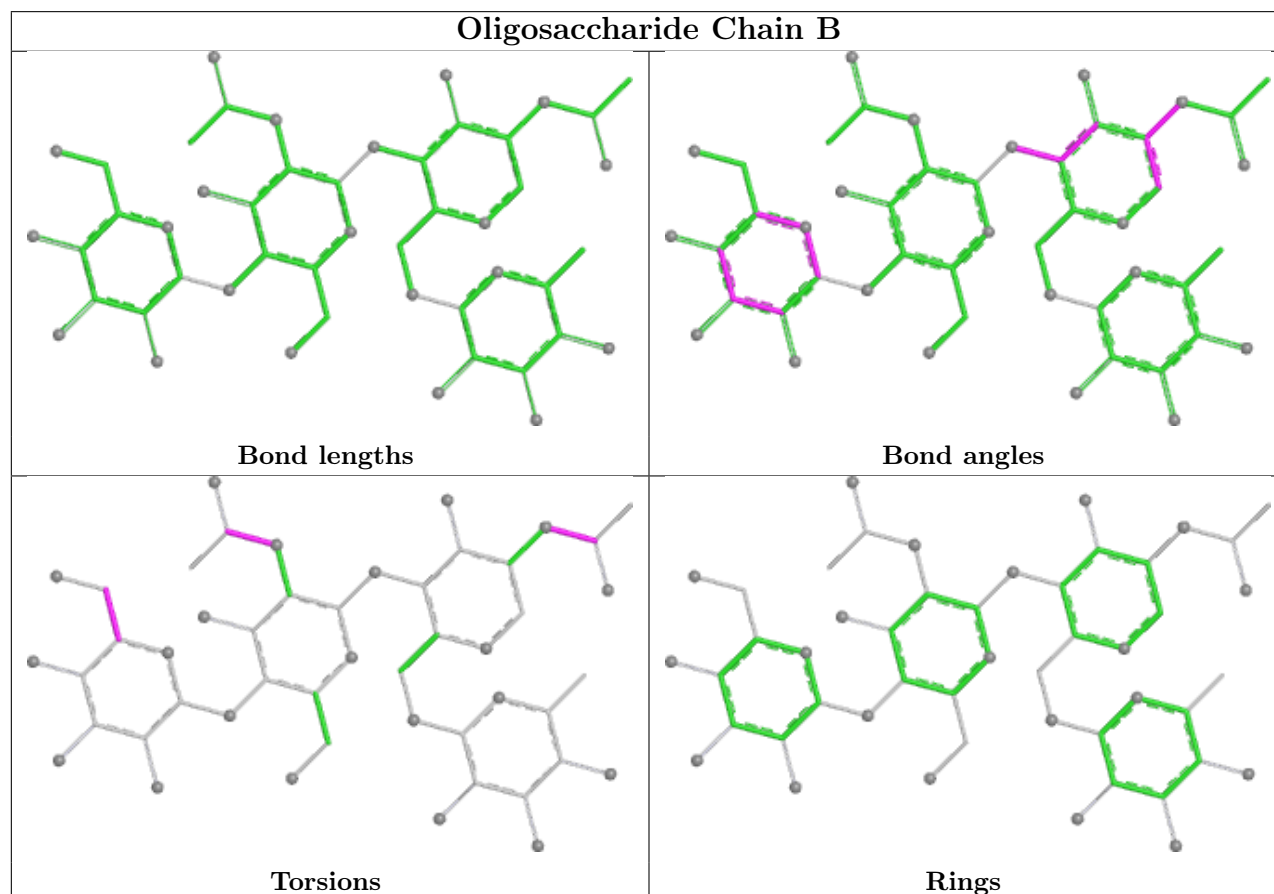
All (5) torsion outliers are listed below:

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

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 oligosaccharide.



## 5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 3 are monoatomic - leaving 8 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$
7	BO3	A	705	-	3,3,3	0.21	0	3,3,3	0.85	0
6	MLI	A	708	-	6,6,6	1.61	1 (16%)	7,7,7	0.98	0
3	NAG	A	701	1	14,14,15	0.47	0	17,19,21	1.64	3 (17%)
8	IMD	A	706	-	5,5,5	0.42	0	5,5,5	0.48	0



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
9	EDO	A	709	-	3,3,3	0.33	0	2,2,2	0.41	0
5	A1JPY	A	703	5	20,21,21	0.49	0	24,29,29	1.29	4 (16%)
6	MLI	A	704	4	6,6,6	1.72	2 (33%)	7,7,7	1.26	0
5	A1JPY	A	707	5	20,21,21	0.56	0	24,29,29	1.40	3 (12%)

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
6	MLI	A	708	-	-	2/4/4/4	-
3	NAG	A	701	1	-	4/6/23/26	0/1/1/1
8	IMD	A	706	-	-	-	0/1/1/1
9	EDO	A	709	-	-	1/1/1/1	-
5	A1JPY	A	703	5	-	1/15/28/28	0/2/2/2
6	MLI	A	704	4	-	0/4/4/4	-
5	A1JPY	A	707	5	-	4/15/28/28	0/2/2/2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	708	MLI	O6-C2	2.48	1.30	1.22
6	A	704	MLI	C1-C2	2.46	1.54	1.51
6	A	704	MLI	O9-C3	-2.02	1.24	1.30

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	701	NAG	C1-C2-N2	4.87	118.10	110.43
5	A	707	A1JPY	C5-C4-S1	4.55	117.76	111.28
5	A	703	A1JPY	C5-C4-N1	2.78	117.09	113.94
3	A	701	NAG	O5-C1-C2	2.51	115.17	111.29
5	A	707	A1JPY	O1-C1-C2	-2.48	114.83	122.44
5	A	703	A1JPY	C3-C2-C1	-2.39	107.03	111.28
3	A	701	NAG	C2-N2-C7	2.36	126.06	122.90
5	A	707	A1JPY	C2-C3-S1	2.28	108.35	105.12
5	A	703	A1JPY	O1-C1-C2	-2.22	115.62	122.44
5	A	703	A1JPY	C5-C4-S1	-2.20	108.15	111.28

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	707	A1JPY	S1-C4-C5-C6
5	A	707	A1JPY	C11-C12-C13-S2
5	A	707	A1JPY	O1-C1-C2-C3
3	A	701	NAG	C8-C7-N2-C2
3	A	701	NAG	O7-C7-N2-C2
3	A	701	NAG	O5-C5-C6-O6
3	A	701	NAG	C4-C5-C6-O6
9	A	709	EDO	O1-C1-C2-O2
5	A	707	A1JPY	O2-C1-C2-C3
6	A	708	MLI	C2-C1-C3-O8
6	A	708	MLI	C2-C1-C3-O9
5	A	703	A1JPY	C11-C12-C13-S2

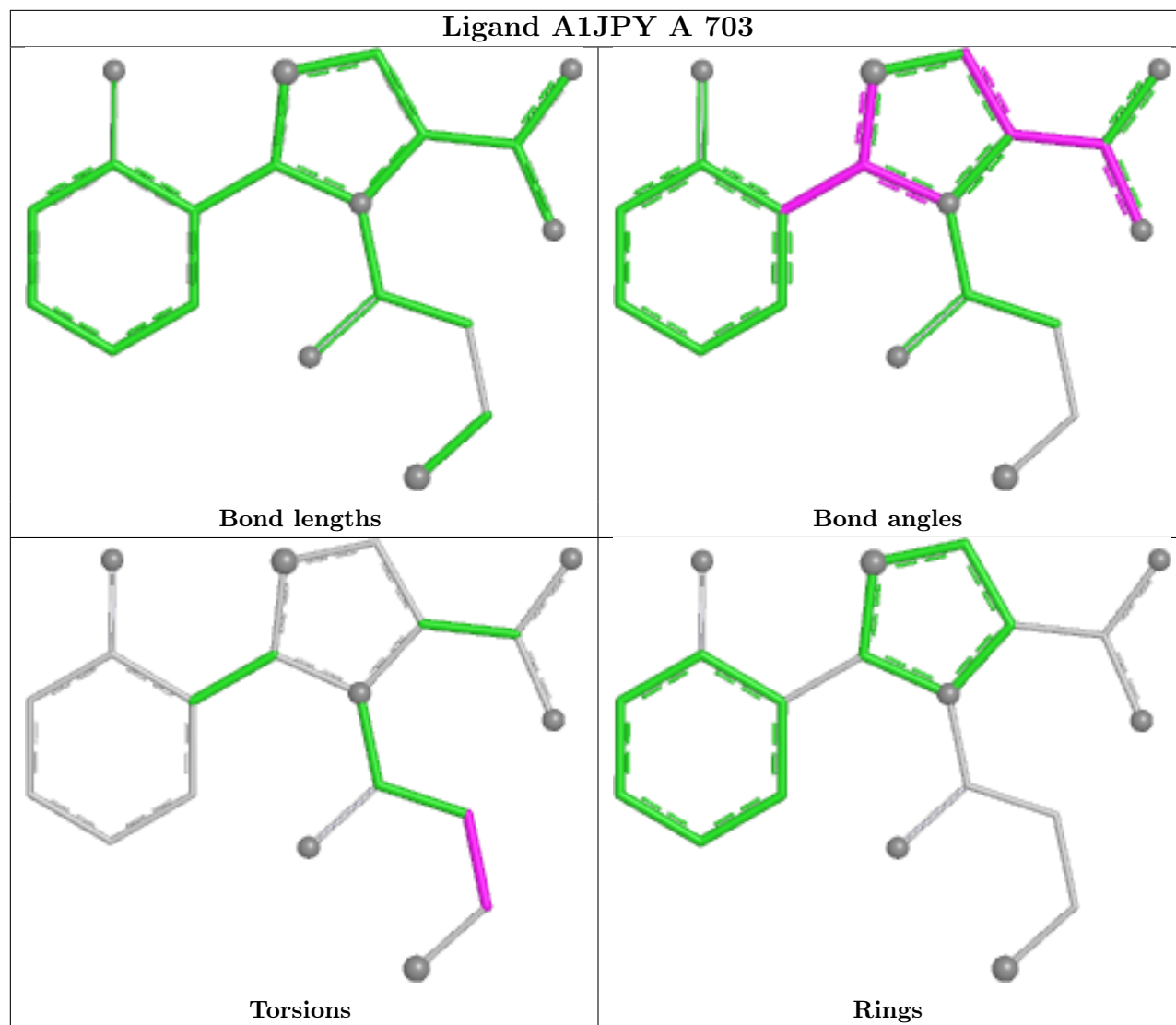
There are no ring outliers.

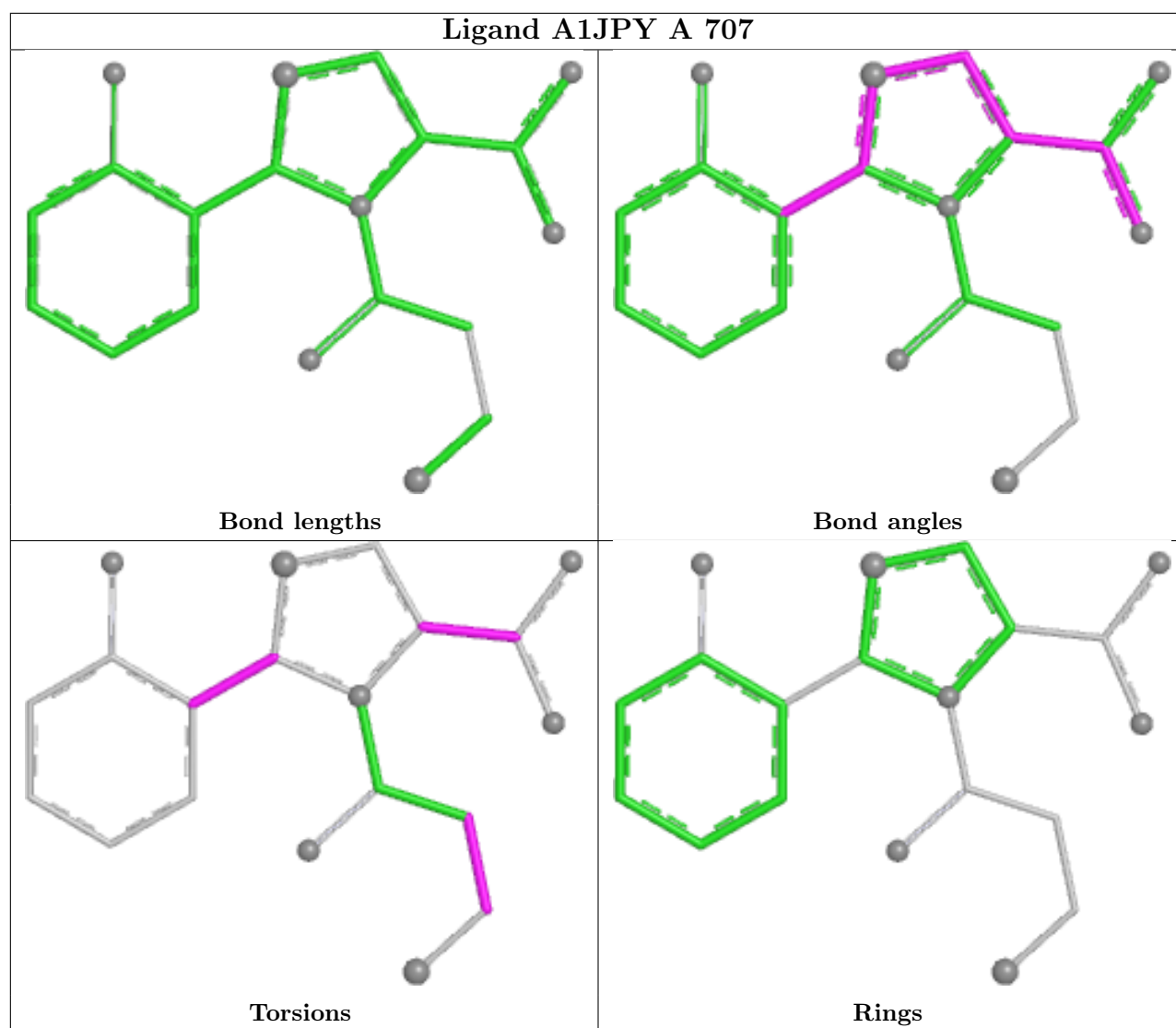
2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	708	MLI	2	0
5	A	707	A1JPY	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.

## Ligand A1JPY A 703





## 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	580/584 (99%)	0.29	28 (4%) 35 41	10, 30, 56, 83	2 (0%)

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	619	TYR	5.0
1	A	621	TRP	4.0
1	A	618	GLN	3.9
1	A	80	ILE	3.4
1	A	623	PRO	3.3
1	A	158	CYS	3.3
1	A	41	GLU	3.1
1	A	69	TYR	3.0
1	A	155	GLN	3.0
1	A	347	GLY	3.0
1	A	154	PRO	2.9
1	A	622	THR	2.8
1	A	345	THR	2.8
1	A	620	ASN	2.7
1	A	73	ILE	2.6
1	A	624	ASN	2.6
1	A	104	VAL	2.5
1	A	77	THR	2.4
1	A	617	PRO	2.4
1	A	434	SER	2.4
1	A	81	LEU	2.3
1	A	59	TRP	2.2
1	A	42	ALA	2.2
1	A	107	LEU	2.1
1	A	425	LYS	2.1
1	A	300	ASP	2.1
1	A	344	PRO	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	74	THR	2.0

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

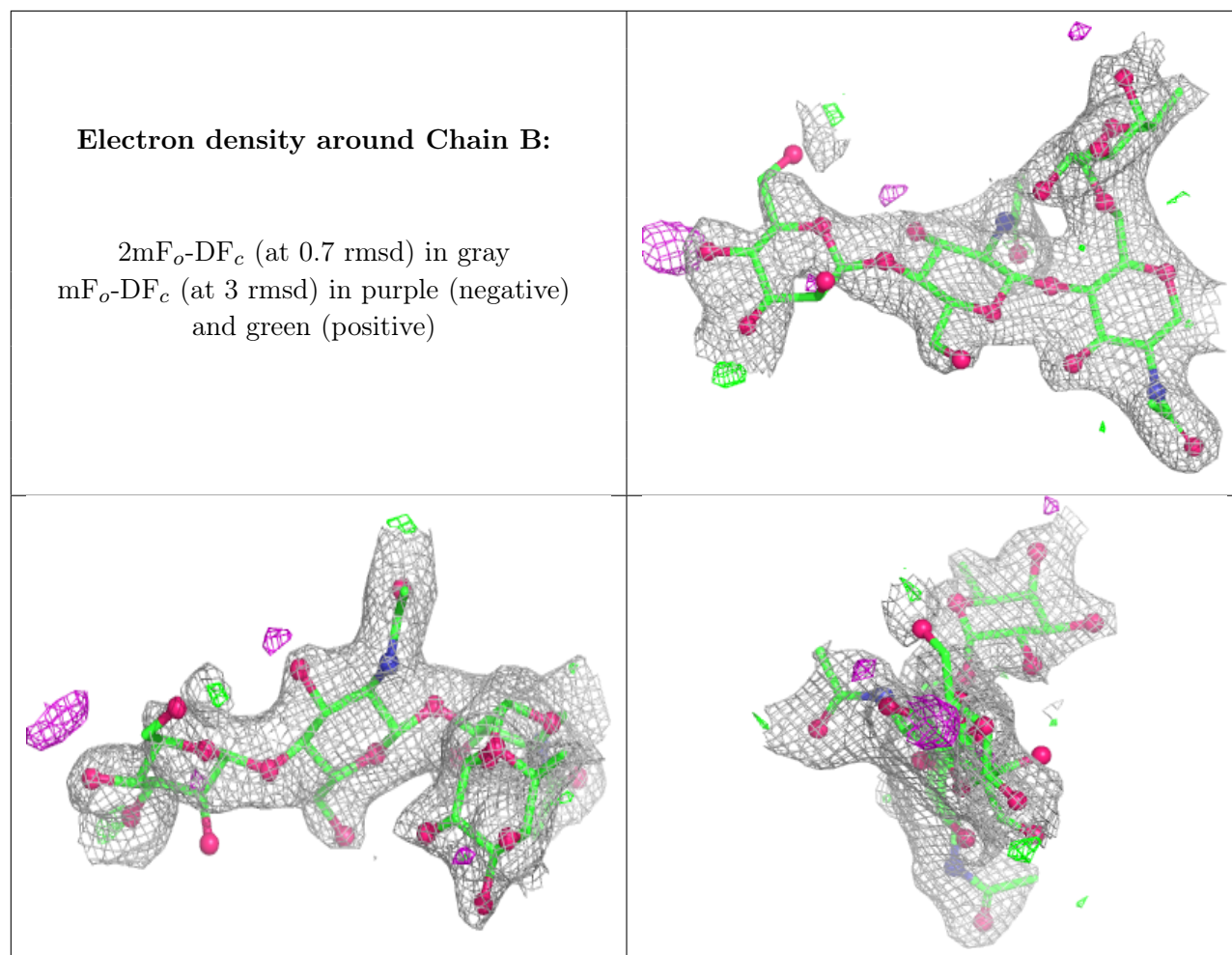
There are no non-standard protein/DNA/RNA residues in this entry.

## 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, 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
2	BMA	B	3	11/12	0.64	0.16	58,71,78,84	0
2	NAG	B	2	14/15	0.81	0.12	43,55,60,64	0
2	FUC	B	4	10/11	0.88	0.11	43,49,53,53	0
2	NAG	B	1	14/15	0.89	0.10	39,47,53,55	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, 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	A	701	14/15	0.59	0.19	67,76,83,84	0
5	A1JPY	A	707	20/20	0.77	0.21	50,64,96,107	0
7	BO3	A	705	4/4	0.82	0.13	41,42,48,55	0
9	EDO	A	709	4/4	0.82	0.17	44,52,52,54	0
6	MLI	A	708	7/7	0.88	0.11	33,39,48,51	0
6	MLI	A	704	7/7	0.92	0.08	28,30,40,40	0
5	A1JPY	A	703	20/20	0.93	0.09	21,29,39,49	0
8	IMD	A	706	5/5	0.95	0.09	36,36,38,39	0
10	CL	A	711	1/1	0.99	0.02	22,22,22,22	0
10	CL	A	710	1/1	1.00	0.03	22,22,22,22	0

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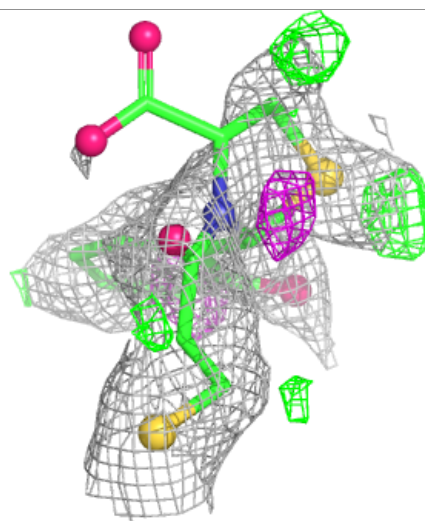
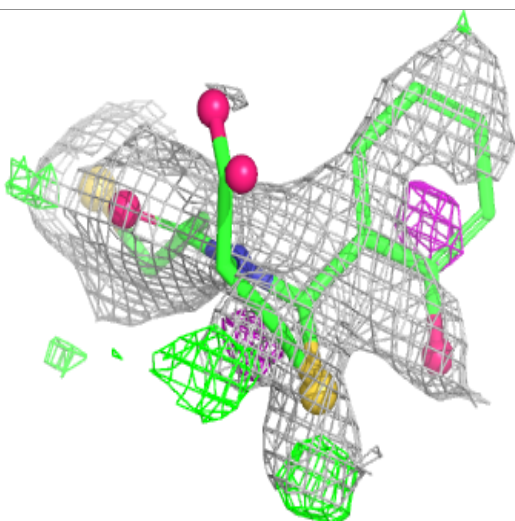
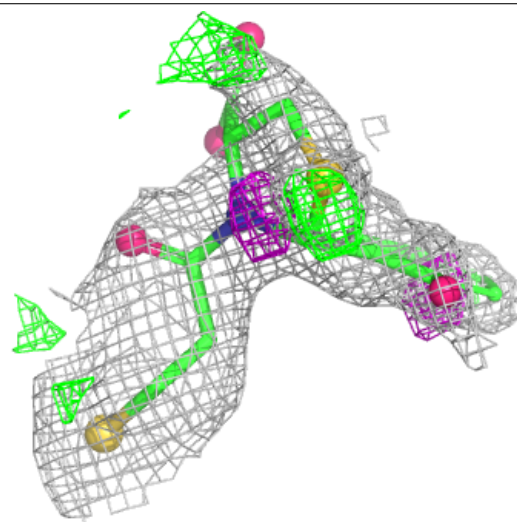
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	ZN	A	702	1/1	1.00	0.01	24,24,24,24	0

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 A1JPY A 707:**

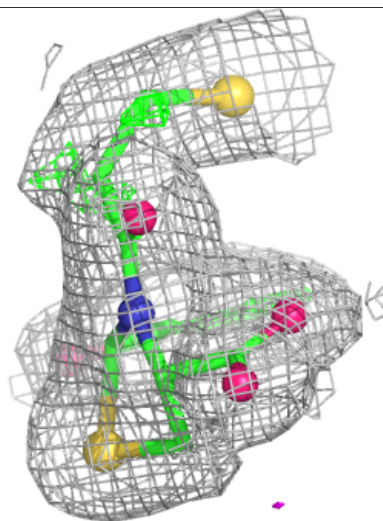
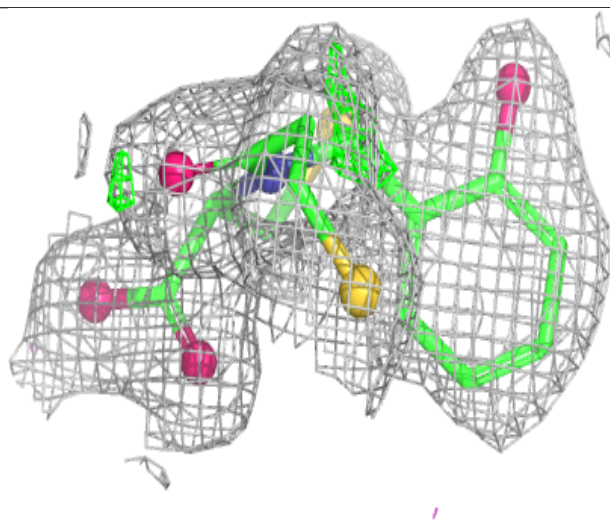
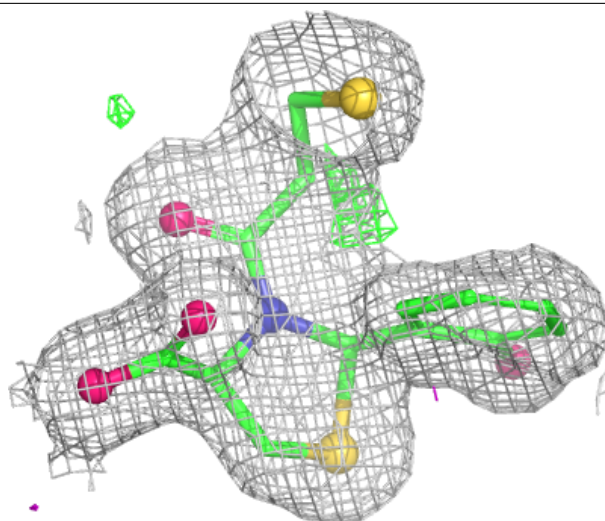
2mF<sub>o</sub>-DF<sub>c</sub> (at 0.7 rmsd) in gray  
mF<sub>o</sub>-DF<sub>c</sub> (at 3 rmsd) in purple (negative)  
and green (positive)





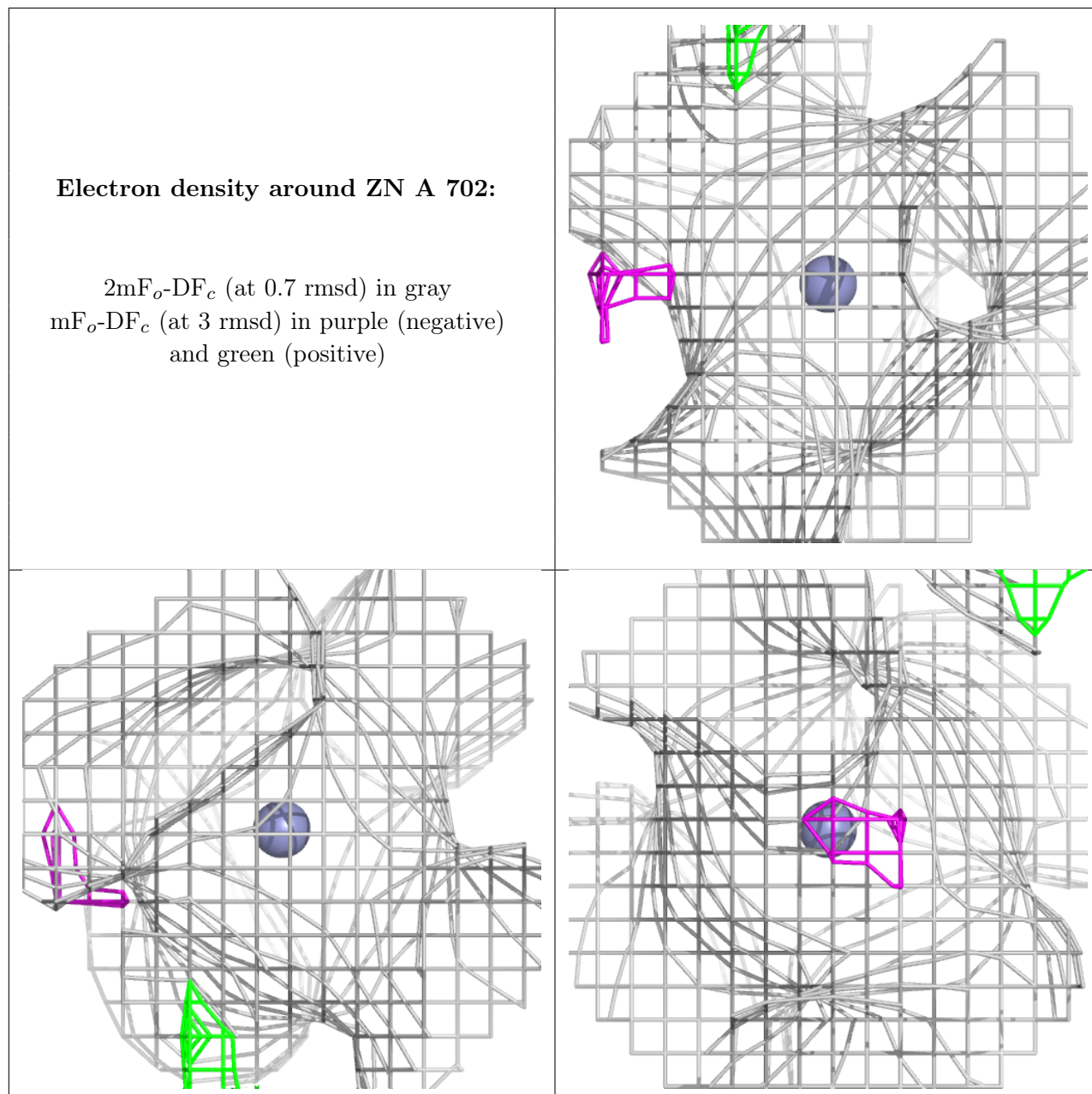
**Electron density around A1JPY A 703:**

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



**Electron density around ZN A 702:**

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



## 6.5 Other polymers ⓘ

There are no such residues in this entry.