



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 3, 2023 – 01:36 PM EDT

PDB ID : 6O9N  
Title : Structural insights on a new fungal aryl-alcohol oxidase  
Authors : Kadowaki, M.A.S.; Polikarpov, I.  
Deposited on : 2019-03-14  
Resolution : 2.60 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 : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.35.1  
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) : 2.35.1

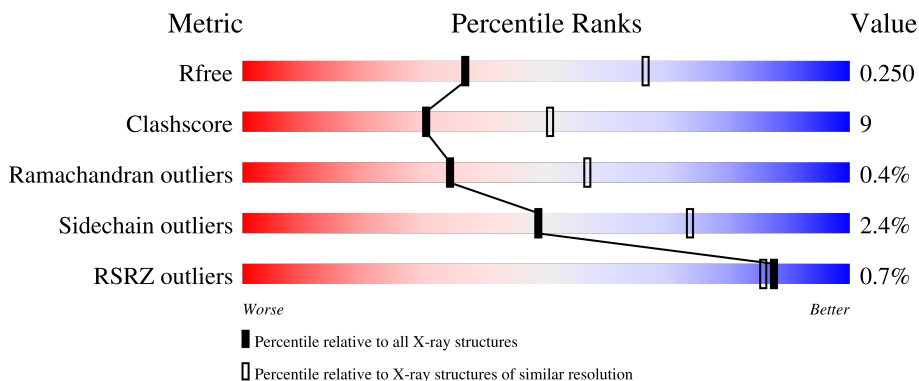
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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	647	 80% 13% • 5%
1	B	647	 79% 14% • 5%
2	C	3	 100%
3	D	2	 50% 50%
3	E	2	 100%

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Mol	Chain	Length	Quality of chain
3	F	2	 100%

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
5	NAG	A	708	-	-	X	-
8	B3P	A	711	-	-	X	-
9	EDO	A	712	-	-	-	X

## 2 Entry composition [i](#)

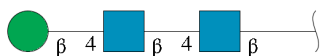
There are 11 unique types of molecules in this entry. The entry contains 10140 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Aryl-alcohol oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	613	Total 4648	C 2958	N 782	O 891	S 17	0	2	0
1	B	613	Total 4651	C 2959	N 783	O 892	S 17	0	2	0

- Molecule 2 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
			Total	C	N	O			
2	C	3	Total 39	C 22	N 2	O 15	0	0	0

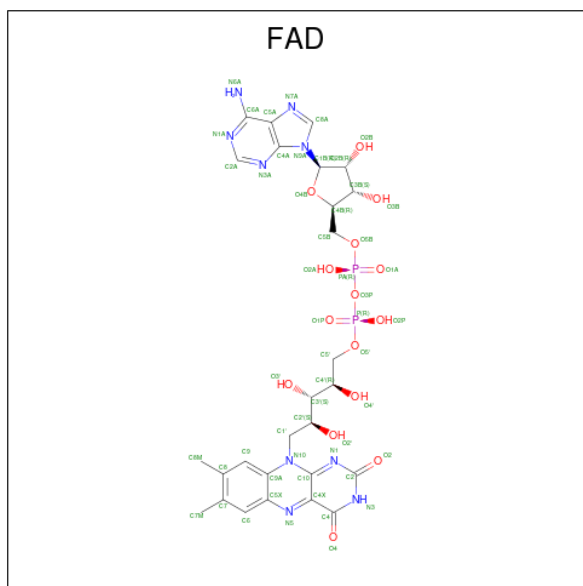
- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



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

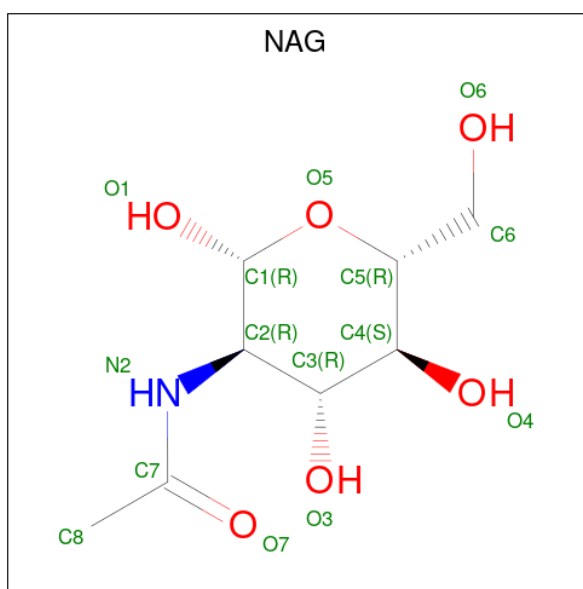
- Molecule 4 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula:

$C_{27}H_{33}N_9O_{15}P_2$ ).



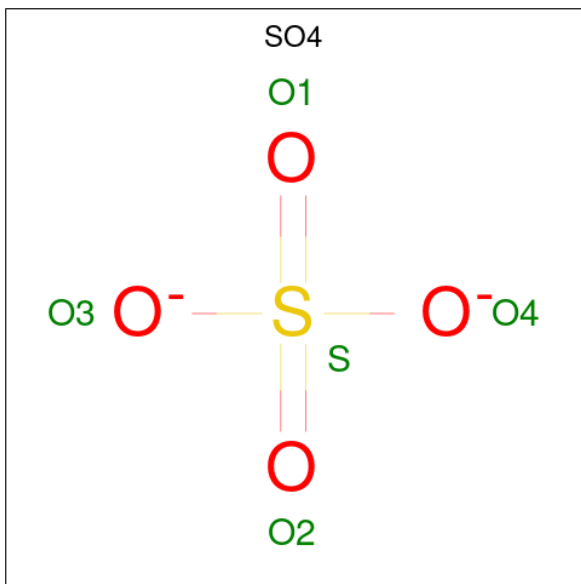
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
4	A	1	Total	C	N	O	P	0	0
			53	27	9	15	2		
4	B	1	Total	C	N	O	P	0	0
			53	27	9	15	2		

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



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

- Molecule 6 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).

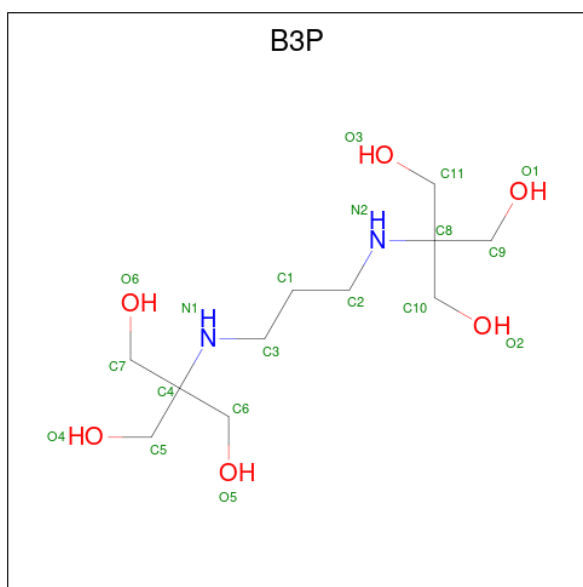


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	O S	0	0
			5	4 1		

- Molecule 7 is CALCIUM ION (three-letter code: CA) (formula: Ca).

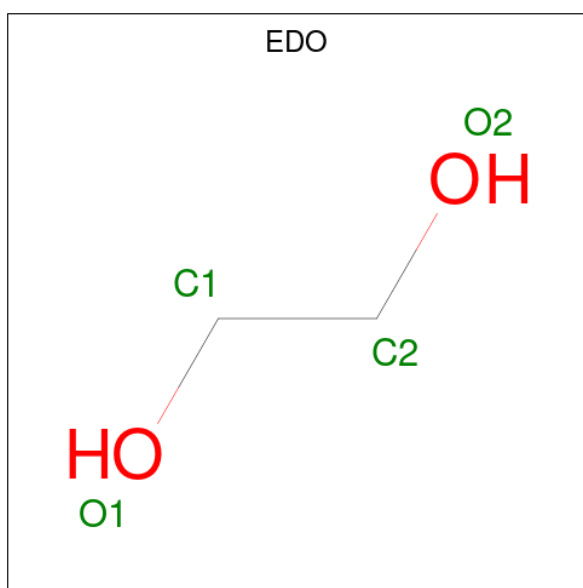
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1	Total	Ca	0	0
			1	1		
7	B	1	Total	Ca	0	0
			1	1		

- Molecule 8 is 2-[3-(2-HYDROXY-1,1-DIHYDROXYMETHYL-ETHYLAMINO)-PROPYL AMINO]-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: B3P) (formula: C<sub>11</sub>H<sub>26</sub>N<sub>2</sub>O<sub>6</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
8	A	1	Total	C	N	O	0	0
			19	11	2	6		
8	B	1	Total	C	N	O	0	0
			19	11	2	6		

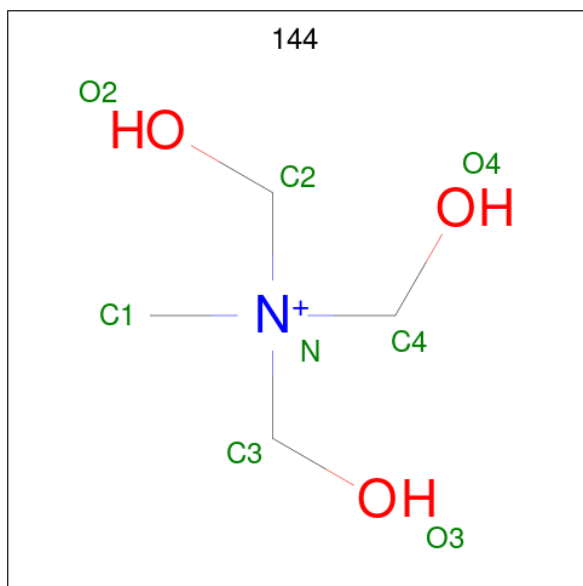
- Molecule 9 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula:  $C_2H_6O_2$ ).



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

- Molecule 10 is TRIS-HYDROXYMETHYL-METHYL-AMMONIUM (three-letter code:

144) (formula: C<sub>4</sub>H<sub>12</sub>NO<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
10	B	1	8	4	1	3	0	0

- Molecule 11 is water.

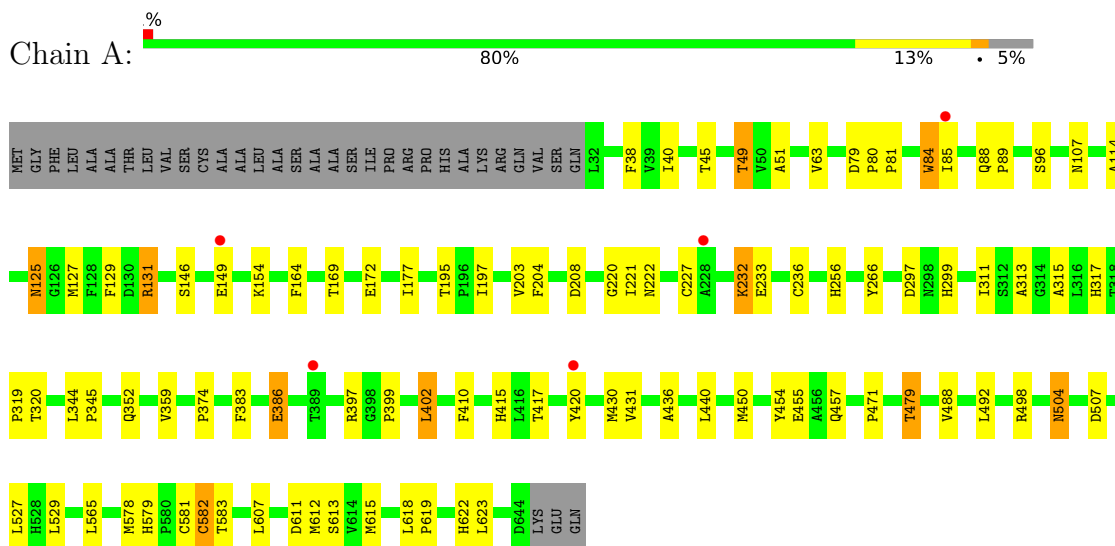
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
11	A	231	231	231	0	0
11	B	282	282	282	0	0



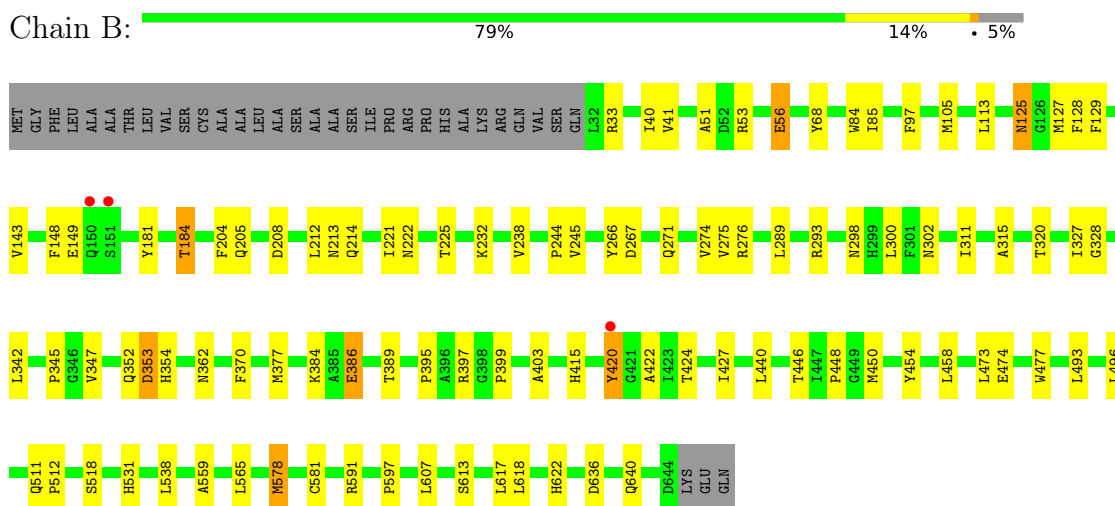
### 3 Residue-property plots i

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Aryl-alcohol oxidase



- Molecule 1: Aryl-alcohol oxidase



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

Chain C:  100%


MAG1  
MAG2  
BMA3

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

Chain D:  50% 50%

MAG1  
MAG2

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

Chain E:  100%

MAG1  
MAG2

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

Chain F:  100%

MAG1  
MAG2

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	63.85Å 109.70Å 220.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	98.19 – 2.60 98.19 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.9 (98.19-2.60) 100.0 (98.19-2.60)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.69 (at 2.58Å)	Xtrriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, $R_{free}$	0.194 , 0.251 0.194 , 0.250	Depositor DCC
$R_{free}$ test set	2458 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	30.4	Xtrriage
Anisotropy	0.056	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 30.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	10140	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.09% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, CA, B3P, FAD, NAG, 144, BMA, EDO

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	0.53	1/4790 (0.0%)	0.66	3/6573 (0.0%)
1	B	0.52	0/4790	0.66	2/6573 (0.0%)
All	All	0.52	1/9580 (0.0%)	0.66	5/13146 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	236	CYS	CB-SG	-5.50	1.72	1.81

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	353	ASP	CB-CG-OD2	-6.56	112.40	118.30
1	B	353	ASP	CB-CG-OD1	5.50	123.25	118.30
1	A	84	TRP	CA-C-N	-5.36	105.41	117.20
1	A	402	LEU	CA-CB-CG	5.24	127.36	115.30
1	A	84	TRP	C-N-CA	5.11	134.47	121.70

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4648	0	4444	96	0
1	B	4651	0	4440	66	0
2	C	39	0	34	0	0
3	D	28	0	25	1	0
3	E	28	0	25	0	0
3	F	28	0	25	3	0
4	A	53	0	31	4	0
4	B	53	0	31	3	0
5	A	28	0	26	9	0
5	B	14	0	13	0	0
6	A	5	0	0	0	0
7	A	1	0	0	0	0
7	B	1	0	0	0	0
8	A	19	0	26	13	0
8	B	19	0	26	7	0
9	A	4	0	6	0	0
10	B	8	0	12	0	0
11	A	231	0	0	3	1
11	B	282	0	0	4	0
All	All	10140	0	9164	174	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

All (174) 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:504:ASN:HD21	5:A:708:NAG:C1	1.34	1.38
1:A:581:CYS:SG	1:A:582:CYS:SG	1.42	1.38
1:A:417:THR:HG21	1:A:420:TYR:CE1	1.65	1.31
1:A:504:ASN:ND2	5:A:708:NAG:C1	1.99	1.25
1:A:417:THR:HG21	1:A:420:TYR:HE1	0.96	1.08
1:A:417:THR:CG2	1:A:420:TYR:CE1	2.37	1.07
1:A:417:THR:CB	1:A:420:TYR:CD1	2.39	1.06
1:A:417:THR:HB	1:A:420:TYR:CD1	1.92	1.04
1:A:504:ASN:OD1	5:A:708:NAG:C1	2.11	0.99
1:A:581:CYS:CB	1:A:582:CYS:SG	2.51	0.97
1:A:504:ASN:CG	5:A:708:NAG:C1	2.33	0.97
1:A:417:THR:CG2	1:A:420:TYR:HE1	1.76	0.96
1:A:417:THR:CB	1:A:420:TYR:CE1	2.52	0.93
1:A:146:SER:HA	1:A:149:GLU:HG3	1.57	0.86
1:A:79:ASP:H	1:A:88:GLN:HE22	1.25	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:504:ASN:HD22	1:A:507:ASP:H	1.23	0.83
1:A:417:THR:OG1	1:A:420:TYR:HD1	1.62	0.81
1:A:417:THR:OG1	1:A:420:TYR:CD1	2.34	0.81
8:A:711:B3P:O1	11:A:801:HOH:O	1.98	0.80
1:A:457:GLN:HE22	1:A:498:ARG:HE	1.29	0.78
1:B:345:PRO:HG2	1:B:450:MET:HE3	1.68	0.75
1:A:233:GLU:OE1	8:A:711:B3P:H51	1.88	0.74
1:B:473:LEU:HD13	1:B:531:HIS:CG	2.22	0.74
1:A:415:HIS:CE1	8:A:711:B3P:H52	2.23	0.74
1:A:222:ASN:H	8:A:711:B3P:H31	1.53	0.73
1:B:125:ASN:HB3	1:B:127:MET:H	1.55	0.72
1:B:622:HIS:HB3	4:B:701:FAD:C2	2.21	0.70
1:B:311:ILE:HD11	1:B:607:LEU:HD11	1.72	0.70
1:A:457:GLN:HE21	1:A:498:ARG:HH21	1.39	0.69
1:B:420:TYR:CZ	1:B:422:ALA:HB3	2.27	0.69
1:A:417:THR:CB	1:A:420:TYR:HD1	1.93	0.69
1:A:504:ASN:OD1	5:A:708:NAG:C2	2.40	0.69
1:B:473:LEU:HD13	1:B:531:HIS:HB3	1.75	0.69
1:B:420:TYR:CE1	1:B:422:ALA:HB3	2.28	0.69
1:A:622:HIS:HB3	4:A:701:FAD:C2	2.25	0.66
1:B:474:GLU:OE1	1:B:622:HIS:HE1	1.78	0.66
1:B:289:LEU:HD13	1:B:302:ASN:OD1	1.95	0.66
8:B:709:B3P:O5	8:B:709:B3P:O4	2.14	0.64
1:A:125:ASN:HB3	1:A:127:MET:H	1.61	0.64
1:A:146:SER:HA	1:A:149:GLU:CG	2.27	0.64
1:A:582:CYS:HB3	1:A:615:MET:O	1.97	0.64
1:B:33:ARG:NH1	1:B:267:ASP:OD2	2.32	0.62
1:A:618:LEU:HD12	1:A:619:PRO:HD2	1.81	0.62
1:A:415:HIS:HE1	8:A:711:B3P:H52	1.65	0.61
1:A:208:ASP:OD1	1:A:479:THR:HG22	2.00	0.61
1:B:362:ASN:ND2	11:B:807:HOH:O	2.32	0.61
1:A:195:THR:O	1:A:197:ILE:N	2.32	0.61
1:A:315:ALA:HA	1:A:613:SER:HB3	1.82	0.60
1:B:578:MET:HG2	4:B:701:FAD:HM83	1.83	0.59
1:B:327:ILE:HG23	1:B:342:LEU:HB2	1.84	0.59
1:B:181:TYR:CD1	1:B:205:GLN:HB2	2.37	0.59
1:A:457:GLN:NE2	1:A:498:ARG:HH21	1.99	0.58
1:B:473:LEU:HD13	1:B:531:HIS:CB	2.32	0.58
1:A:311:ILE:HD12	1:A:607:LEU:HD11	1.85	0.58
1:A:529:LEU:HD22	1:A:565:LEU:HG	1.86	0.58
1:A:504:ASN:ND2	1:A:507:ASP:H	1.98	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:A:711:B3P:O4	8:A:711:B3P:O5	2.20	0.56
1:B:458:LEU:HG	1:B:496:LEU:HD23	1.88	0.56
1:A:221:ILE:HA	8:A:711:B3P:H22	1.87	0.56
3:F:2:NAG:H83	3:F:2:NAG:H3	1.87	0.56
8:A:711:B3P:O4	8:A:711:B3P:O6	2.16	0.55
1:A:457:GLN:HE22	1:A:498:ARG:NE	2.03	0.55
1:A:146:SER:CA	1:A:149:GLU:HG3	2.34	0.55
1:B:636:ASP:O	1:B:640:GLN:HG3	2.06	0.55
1:A:204:PHE:CD2	1:A:399:PRO:HA	2.41	0.55
8:A:711:B3P:HN2	8:A:711:B3P:HN1	1.55	0.55
1:A:579:HIS:HB3	1:A:623:LEU:HG	1.90	0.54
1:B:386[A]:GLU:CD	1:B:397:ARG:HH12	2.11	0.54
1:A:504:ASN:OD1	5:A:708:NAG:H2	2.08	0.54
1:A:352:GLN:NE2	1:A:618:LEU:H	2.04	0.53
1:B:68:TYR:O	1:B:271:GLN:HA	2.08	0.53
1:B:352:GLN:NE2	1:B:618:LEU:H	2.06	0.53
1:B:222:ASN:H	8:B:709:B3P:H12	1.74	0.52
1:B:420:TYR:CE1	1:B:422:ALA:CB	2.92	0.52
1:A:79:ASP:H	1:A:88:GLN:NE2	2.00	0.52
1:B:275:VAL:HG12	1:B:276:ARG:HG2	1.91	0.52
1:A:386[A]:GLU:CD	1:A:397:ARG:HH12	2.14	0.52
1:A:177:ILE:HG13	1:A:203:VAL:HG12	1.92	0.52
1:B:149:GLU:HA	1:B:149:GLU:OE1	2.09	0.51
1:A:410:PHE:HB3	1:A:471:PRO:HB3	1.92	0.51
1:A:38:PHE:HB2	1:A:63:VAL:HG22	1.93	0.51
1:A:107:ASN:N	1:A:107:ASN:OD1	2.43	0.51
1:A:457:GLN:NE2	1:A:498:ARG:HE	2.02	0.51
1:A:227:CYS:HB3	1:A:232:LYS:HB2	1.91	0.51
1:A:457:GLN:HE21	1:A:498:ARG:NH2	2.07	0.51
5:A:707:NAG:C6	5:A:707:NAG:H2	2.39	0.51
1:B:415:HIS:HB3	8:B:709:B3P:H51	1.92	0.50
1:B:353:ASP:OD1	1:B:354:HIS:N	2.45	0.50
1:A:221:ILE:HG23	8:A:711:B3P:H11	1.93	0.50
1:B:105:MET:HG3	1:B:518:SER:HB2	1.94	0.50
1:B:352:GLN:HE22	1:B:617:LEU:HA	1.76	0.50
1:A:345:PRO:HG2	1:A:450:MET:HE3	1.93	0.49
1:B:377:MET:HE3	1:B:384:LYS:HG3	1.93	0.49
1:A:504:ASN:ND2	5:A:708:NAG:O5	2.41	0.49
1:A:222:ASN:HB2	8:A:711:B3P:H61	1.93	0.49
1:B:302:ASN:OD1	3:F:1:NAG:O5	2.31	0.49
1:A:220:GLY:O	8:A:711:B3P:O1	2.18	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:204:PHE:CD2	1:B:399:PRO:HA	2.48	0.48
1:A:374:PRO:CD	1:A:479:THR:HG21	2.44	0.48
1:B:320:THR:HG21	1:B:512:PRO:HD3	1.95	0.47
1:B:232:LYS:O	1:B:232:LYS:HG2	2.13	0.47
1:B:143:VAL:O	1:B:591:ARG:NH1	2.46	0.47
1:B:424:THR:HA	1:B:427:ILE:HD12	1.96	0.47
1:B:427:ILE:HG12	1:B:496:LEU:HD21	1.96	0.47
1:B:40:ILE:HD13	1:B:51:ALA:HB2	1.96	0.47
1:A:297:ASP:OD1	1:A:299:HIS:HB3	2.14	0.46
1:A:352:GLN:HE22	1:A:618:LEU:H	1.64	0.46
1:B:559:ALA:HA	1:B:565:LEU:HD13	1.96	0.46
1:B:415:HIS:HE1	8:B:709:B3P:H91	1.81	0.46
1:A:131:ARG:HD2	1:A:164:PHE:CD1	2.50	0.46
4:A:701:FAD:H1'1	4:A:701:FAD:H9	1.77	0.45
1:B:315:ALA:HA	1:B:613:SER:HB3	1.98	0.45
1:A:374:PRO:HD3	1:A:479:THR:HG21	1.99	0.45
1:B:354:HIS:CD2	1:B:493:LEU:HA	2.51	0.45
1:B:181:TYR:CE1	1:B:205:GLN:HB2	2.51	0.45
1:B:311:ILE:CD1	1:B:607:LEU:HD11	2.42	0.45
1:A:578:MET:HG2	4:A:701:FAD:HM83	1.98	0.45
1:A:383:PHE:HD1	3:D:1:NAG:H61	1.81	0.45
8:B:709:B3P:HO4	8:B:709:B3P:HO5	1.57	0.45
1:A:345:PRO:HB2	1:A:450:MET:HE3	1.99	0.45
1:A:504:ASN:HB3	1:A:507:ASP:O	2.17	0.45
1:B:125:ASN:HB3	1:B:127:MET:N	2.29	0.45
1:B:298:ASN:O	1:B:300:LEU:HD12	2.17	0.45
1:A:45:THR:O	1:A:49:THR:HG23	2.17	0.45
1:A:440:LEU:HD21	1:A:454:TYR:CG	2.52	0.45
1:A:611:ASP:HB2	4:A:701:FAD:O2P	2.17	0.45
1:A:204:PHE:HD2	1:A:402:LEU:HD22	1.81	0.44
1:B:184:THR:CG2	1:B:225:THR:O	2.65	0.44
11:B:888:HOH:O	3:F:2:NAG:H2	2.16	0.44
1:B:244:PRO:HG3	11:B:922:HOH:O	2.18	0.44
1:A:492:LEU:HD11	1:A:527:LEU:HD13	1.99	0.44
1:B:415:HIS:CG	8:B:709:B3P:H72	2.53	0.44
1:B:214:GLN:HG2	11:B:1038:HOH:O	2.17	0.44
1:B:328:GLY:HA2	1:B:347:VAL:HB	1.99	0.44
1:A:40:ILE:HD13	1:A:51:ALA:HB2	1.98	0.44
1:A:172:GLU:HG3	11:A:906:HOH:O	2.17	0.44
1:A:344:LEU:HD23	1:A:344:LEU:HA	1.85	0.44
1:A:146:SER:HA	1:A:149:GLU:CD	2.38	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:311:ILE:CD1	1:A:607:LEU:HD11	2.47	0.43
1:B:213:ASN:HB2	1:B:477:TRP:HZ2	1.84	0.43
1:B:446:THR:O	1:B:448:PRO:HD3	2.19	0.43
1:A:417:THR:OG1	1:A:420:TYR:CE1	2.63	0.43
1:B:41:VAL:HG13	1:B:274:VAL:HG21	2.01	0.43
1:A:80:PRO:HA	1:A:81:PRO:HA	1.77	0.43
1:B:148:PHE:CE1	1:B:597:PRO:HD3	2.53	0.43
1:B:97:PHE:CZ	1:B:511:GLN:HB2	2.54	0.42
1:A:345:PRO:CB	1:A:450:MET:HE3	2.49	0.42
1:B:184:THR:HG23	1:B:225:THR:O	2.19	0.42
4:B:701:FAD:H9	4:B:701:FAD:H1'1	1.82	0.42
1:A:45:THR:O	1:A:49:THR:CG2	2.68	0.42
1:B:440:LEU:HD21	1:B:454:TYR:CG	2.54	0.42
1:A:612:MET:CE	1:A:612:MET:HA	2.50	0.42
1:A:208:ASP:OD1	1:A:479:THR:CG2	2.67	0.42
1:A:88:GLN:HA	1:A:89:PRO:HD3	1.96	0.42
1:A:430:MET:HG2	1:A:436:ALA:HB2	2.02	0.42
1:A:233:GLU:OE2	8:A:711:B3P:H12	2.19	0.42
1:B:221:ILE:HA	8:B:709:B3P:N1	2.35	0.41
1:A:49:THR:HB	1:A:256:HIS:NE2	2.36	0.41
1:A:359:VAL:HB	1:A:488:VAL:HG23	2.02	0.41
1:B:53:ARG:HD3	1:B:56:GLU:OE2	2.21	0.41
1:B:245:VAL:HG13	1:B:395:PRO:HB3	2.01	0.41
1:A:431:VAL:HG13	1:A:455:GLU:HG2	2.02	0.41
1:A:457:GLN:NE2	1:A:498:ARG:NH2	2.67	0.41
1:A:125:ASN:HD22	1:A:125:ASN:HA	1.62	0.41
1:A:313:ALA:O	1:A:317:HIS:HB2	2.19	0.41
1:A:319:PRO:HG3	1:A:583:THR:HB	2.02	0.41
5:A:707:NAG:C6	5:A:707:NAG:C2	2.99	0.41
1:A:581:CYS:HB2	1:A:582:CYS:SG	2.53	0.41
1:A:114:ALA:HB1	11:A:952:HOH:O	2.21	0.40
1:B:113:LEU:N	1:B:578:MET:HE3	2.36	0.40
1:B:386[A]:GLU:HA	1:B:389:THR:HG22	2.02	0.40
1:B:128:PHE:CE2	1:B:238:VAL:HG11	2.57	0.40
1:B:208:ASP:HA	1:B:370:PHE:CE1	2.57	0.40
1:B:212:LEU:HD22	1:B:538:LEU:HD11	2.04	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
11:A:880:HOH:O	11:A:880:HOH:O[2_557]	2.17	0.03

### 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	613/647 (95%)	579 (94%)	32 (5%)	2 (0%)	41 64
1	B	613/647 (95%)	582 (95%)	28 (5%)	3 (0%)	29 52
All	All	1226/1294 (95%)	1161 (95%)	60 (5%)	5 (0%)	34 57

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	125	ASN
1	B	125	ASN
1	A	84	TRP
1	B	84	TRP
1	B	403	ALA

#### 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	493/516 (96%)	478 (97%)	15 (3%)	41 67
1	B	493/516 (96%)	482 (98%)	11 (2%)	52 76
All	All	986/1032 (96%)	960 (97%)	26 (3%)	49 72

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	THR
1	A	85	ILE
1	A	96	SER
1	A	129	PHE
1	A	131	ARG
1	A	154	LYS
1	A	169	THR
1	A	232	LYS
1	A	266	TYR
1	A	320	THR
1	A	386[A]	GLU
1	A	386[B]	GLU
1	A	479	THR
1	A	504	ASN
1	A	582	CYS
1	B	56	GLU
1	B	85	ILE
1	B	129	PHE
1	B	184	THR
1	B	266	TYR
1	B	293	ARG
1	B	386[A]	GLU
1	B	386[B]	GLU
1	B	420	TYR
1	B	578	MET
1	B	581	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	88	GLN
1	A	125	ASN
1	A	352	GLN
1	A	379	ASN
1	A	457	GLN
1	B	213	ASN
1	B	271	GLN
1	B	352	GLN
1	B	362	ASN
1	B	622	HIS
1	B	640	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

9 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	C	1	1,2	14,14,15	0.90	1 (7%)	17,19,21	1.40	3 (17%)
2	NAG	C	2	2	14,14,15	0.85	1 (7%)	17,19,21	0.48	0
2	BMA	C	3	2	11,11,12	0.95	1 (9%)	15,15,17	0.93	0
3	NAG	D	1	1,3	14,14,15	1.10	2 (14%)	17,19,21	1.16	1 (5%)
3	NAG	D	2	3	14,14,15	0.71	0	17,19,21	0.74	0
3	NAG	E	1	3	14,14,15	1.24	1 (7%)	17,19,21	2.04	1 (5%)
3	NAG	E	2	3	14,14,15	0.88	1 (7%)	17,19,21	1.08	2 (11%)
3	NAG	F	1	1,3	14,14,15	1.20	2 (14%)	17,19,21	1.27	2 (11%)
3	NAG	F	2	3	14,14,15	0.70	1 (7%)	17,19,21	1.53	3 (17%)

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	C	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
2	BMA	C	3	2	-	2/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	D	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	D	2	3	-	2/6/23/26	0/1/1/1
3	NAG	E	1	3	-	2/6/23/26	0/1/1/1
3	NAG	E	2	3	-	2/6/23/26	0/1/1/1
3	NAG	F	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	F	2	3	-	3/6/23/26	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	1	NAG	C1-C2	-3.74	1.46	1.52
3	F	1	NAG	O5-C1	3.37	1.49	1.43
3	E	2	NAG	C1-C2	2.63	1.56	1.52
2	C	2	NAG	O5-C1	-2.52	1.39	1.43
3	D	1	NAG	C1-C2	-2.46	1.48	1.52
2	C	1	NAG	C1-C2	2.45	1.56	1.52
3	F	2	NAG	C1-C2	2.44	1.56	1.52
3	D	1	NAG	O5-C1	-2.36	1.39	1.43
2	C	3	BMA	C4-C3	2.16	1.57	1.52
3	F	1	NAG	C8-C7	-2.03	1.46	1.50

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	1	NAG	C1-O5-C5	7.41	122.24	112.19
3	F	2	NAG	C2-N2-C7	4.08	128.71	122.90
3	D	1	NAG	C1-O5-C5	3.61	117.09	112.19
3	F	2	NAG	C1-O5-C5	3.27	116.63	112.19
3	F	1	NAG	C1-O5-C5	3.23	116.57	112.19
2	C	1	NAG	C1-O5-C5	2.78	115.96	112.19
3	F	2	NAG	C1-C2-N2	2.50	114.76	110.49
3	E	2	NAG	C3-C4-C5	-2.41	105.94	110.24
2	C	1	NAG	O3-C3-C2	-2.39	104.52	109.47
3	E	2	NAG	C1-O5-C5	2.23	115.21	112.19
2	C	1	NAG	C3-C4-C5	-2.12	106.46	110.24
3	F	1	NAG	C3-C4-C5	-2.10	106.49	110.24

There are no chirality outliers.

All (15) torsion outliers are listed below:

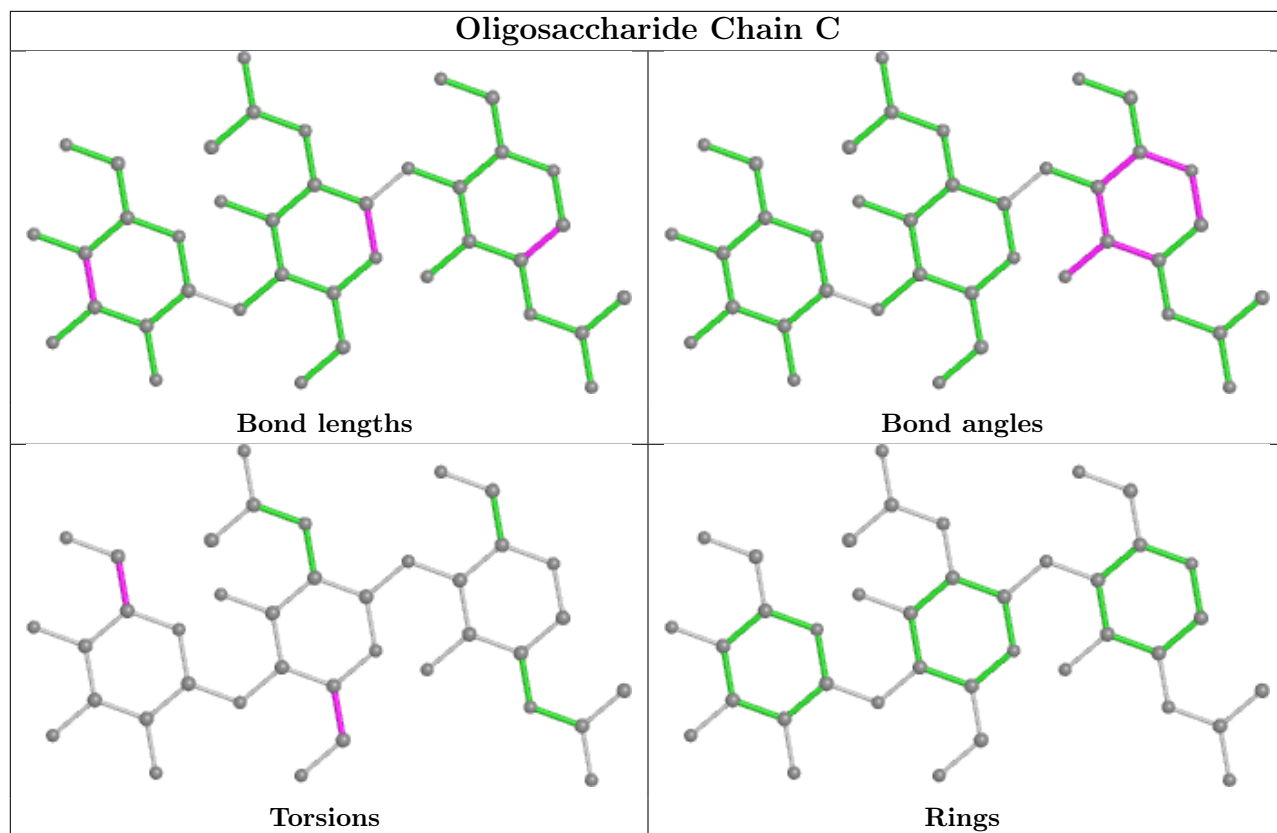
Mol	Chain	Res	Type	Atoms
2	C	3	BMA	O5-C5-C6-O6
2	C	3	BMA	C4-C5-C6-O6
3	D	2	NAG	C4-C5-C6-O6
3	E	2	NAG	O5-C5-C6-O6
3	D	2	NAG	O5-C5-C6-O6
3	E	1	NAG	C8-C7-N2-C2
3	E	1	NAG	O7-C7-N2-C2
3	F	2	NAG	C8-C7-N2-C2
3	F	2	NAG	O7-C7-N2-C2
2	C	2	NAG	C4-C5-C6-O6
3	E	2	NAG	C4-C5-C6-O6
3	D	1	NAG	C4-C5-C6-O6
3	D	1	NAG	O5-C5-C6-O6
2	C	2	NAG	O5-C5-C6-O6
3	F	2	NAG	C3-C2-N2-C7

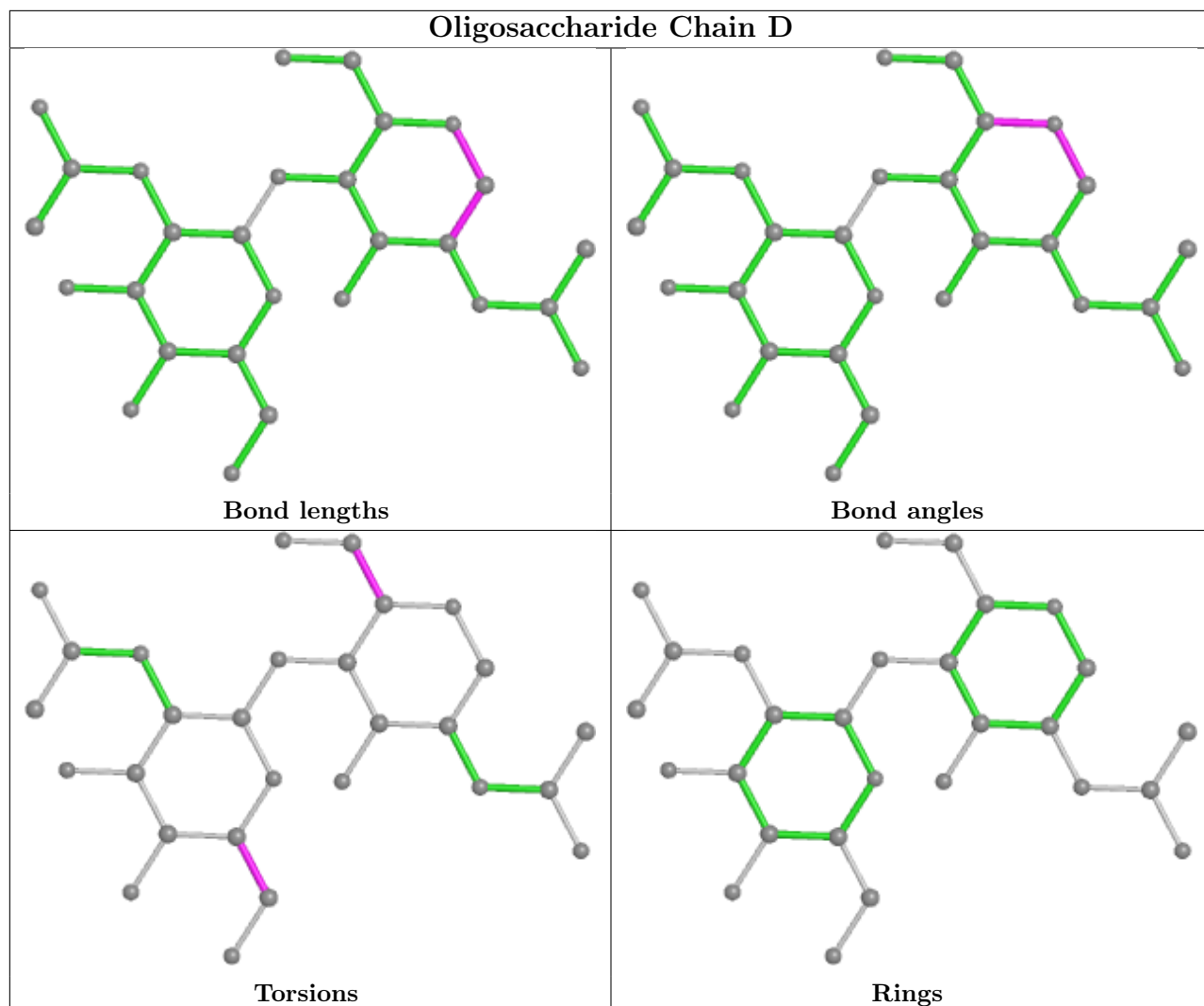
There are no ring outliers.

3 monomers are involved in 4 short contacts:

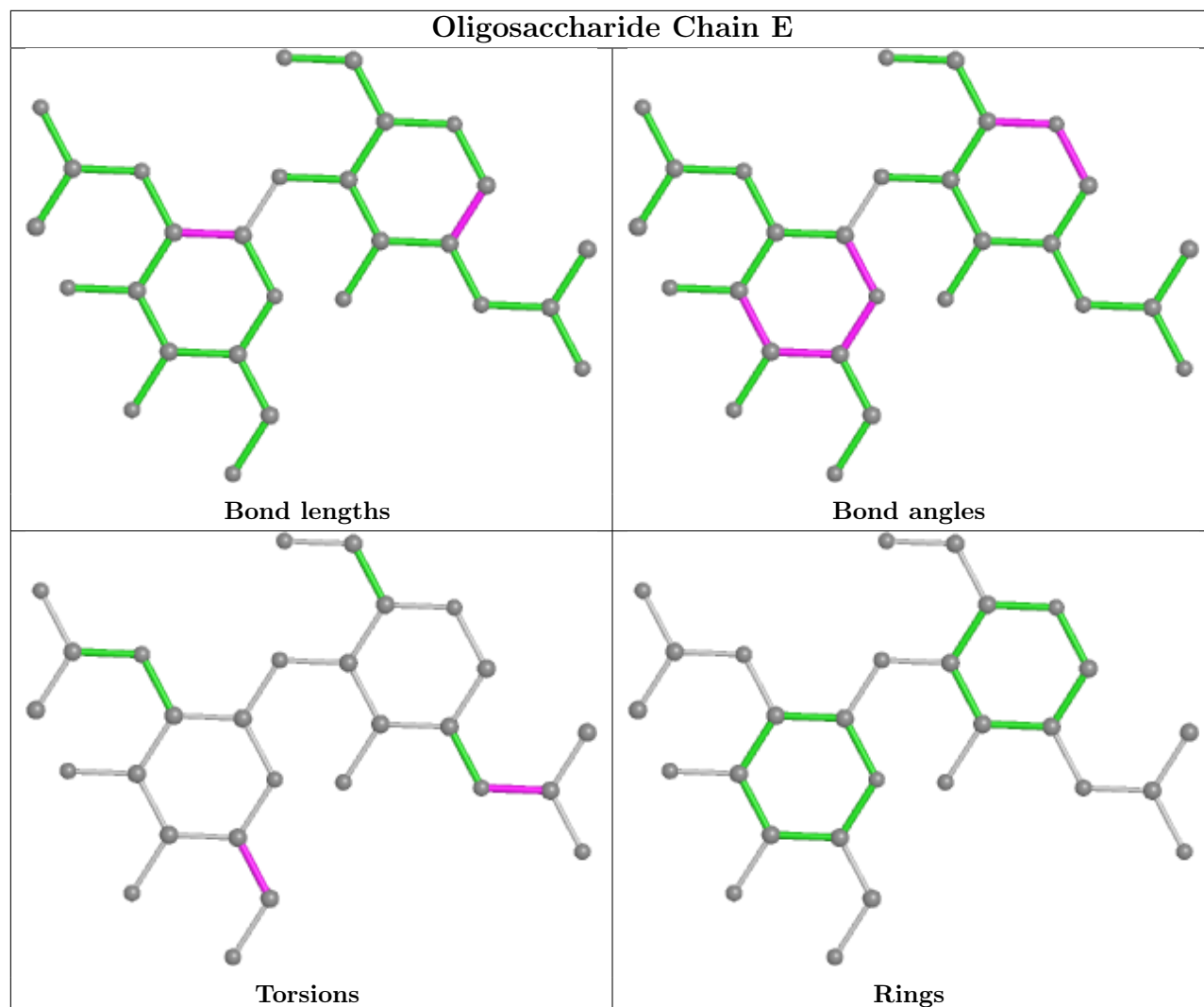
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	2	NAG	2	0
3	F	1	NAG	1	0
3	D	1	NAG	1	0

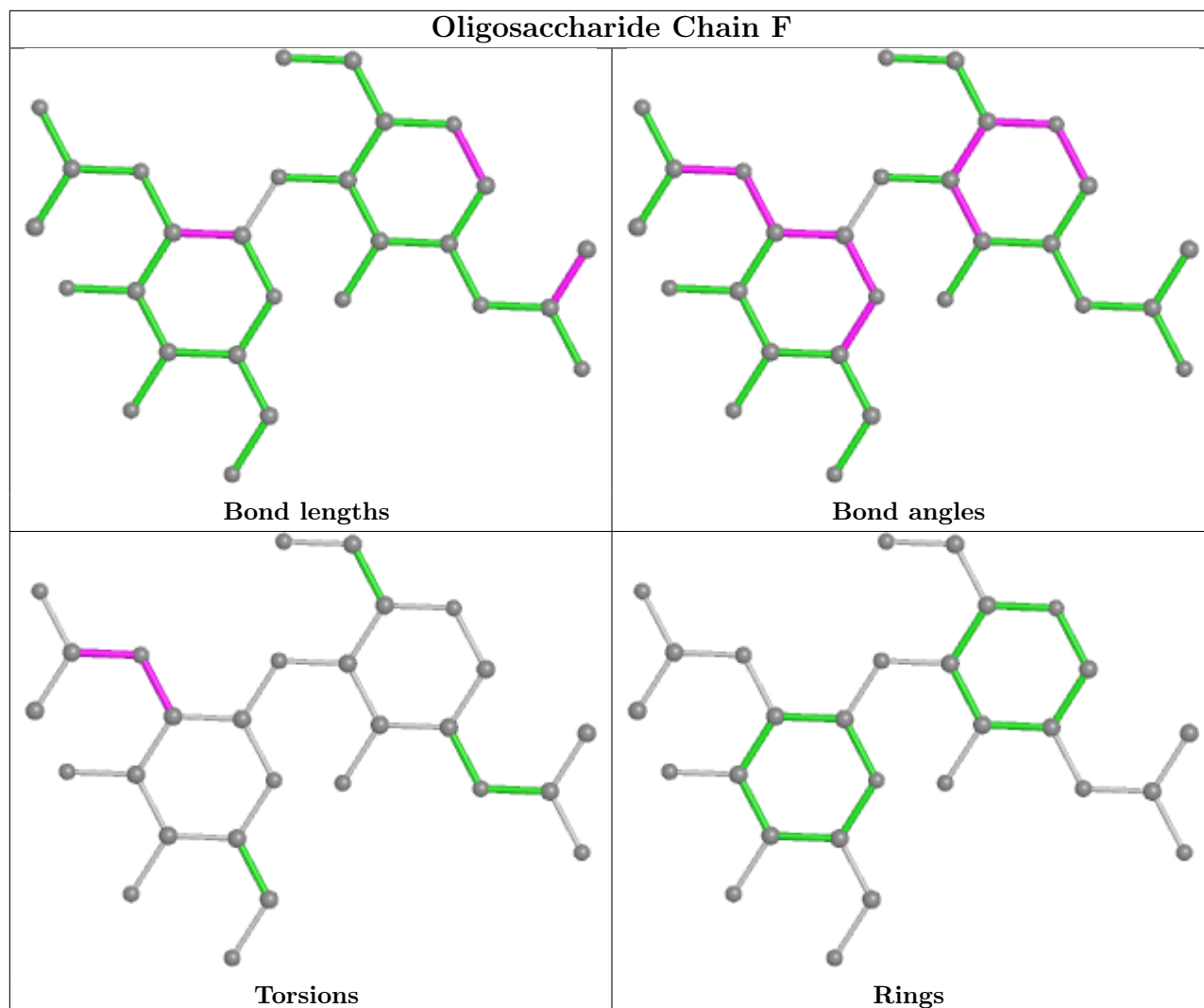
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











## 5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 2 are monoatomic - leaving 10 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
5	NAG	A	707	1	14,14,15	0.94	1 (7%)	17,19,21	1.86	5 (29%)
5	NAG	A	708	-	14,14,15	1.45	1 (7%)	17,19,21	1.30	1 (5%)
4	FAD	A	701	-	53,58,58	0.54	0	68,89,89	0.66	1 (1%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
10	144	B	708	-	1,7,7	1.22	0	3,9,9	0.09	0
9	EDO	A	712	-	3,3,3	0.75	0	2,2,2	0.22	0
8	B3P	A	711	-	18,18,18	1.31	3 (16%)	21,23,23	1.75	7 (33%)
5	NAG	B	706	1	14,14,15	0.76	1 (7%)	17,19,21	0.88	0
6	SO4	A	709	-	4,4,4	0.18	0	6,6,6	0.26	0
4	FAD	B	701	-	53,58,58	0.46	0	68,89,89	0.71	2 (2%)
8	B3P	B	709	-	18,18,18	0.85	0	21,23,23	1.10	1 (4%)

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	707	1	-	2/6/23/26	0/1/1/1
5	NAG	A	708	-	-	2/6/23/26	0/1/1/1
4	FAD	A	701	-	-	4/30/50/50	0/6/6/6
10	144	B	708	-	-	0/0/9/9	-
9	EDO	A	712	-	-	1/1/1/1	-
8	B3P	A	711	-	-	7/28/28/28	-
5	NAG	B	706	1	-	2/6/23/26	0/1/1/1
4	FAD	B	701	-	-	4/30/50/50	0/6/6/6
8	B3P	B	709	-	-	14/28/28/28	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	708	NAG	O5-C1	4.84	1.51	1.43
8	A	711	B3P	C9-C8	2.92	1.57	1.53
8	A	711	B3P	C5-C4	-2.47	1.50	1.53
8	A	711	B3P	C3-N1	2.37	1.49	1.46
5	A	707	NAG	C1-C2	-2.12	1.49	1.52
5	B	706	NAG	C1-C2	2.07	1.55	1.52

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	707	NAG	C1-O5-C5	4.86	118.78	112.19
5	A	708	NAG	C1-O5-C5	4.52	118.32	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	711	B3P	C11-C8-C9	-3.61	102.41	110.04
5	A	707	NAG	O4-C4-C5	-3.56	100.47	109.30
8	A	711	B3P	C7-C4-C5	-3.18	103.33	110.04
8	A	711	B3P	C2-N2-C8	2.57	119.72	116.08
4	B	701	FAD	P-O3P-PA	-2.52	124.19	132.83
8	B	709	B3P	C2-N2-C8	-2.44	112.61	116.08
4	B	701	FAD	C5A-C6A-N6A	2.42	124.03	120.35
8	A	711	B3P	C6-C4-C5	-2.31	105.16	110.04
4	A	701	FAD	C5A-C6A-N6A	2.24	123.75	120.35
5	A	707	NAG	C1-C2-N2	2.24	114.31	110.49
8	A	711	B3P	C10-C8-N2	2.23	115.72	109.03
5	A	707	NAG	O5-C5-C6	2.22	110.68	107.20
8	A	711	B3P	O1-C9-C8	2.21	116.11	111.63
8	A	711	B3P	O3-C11-C8	-2.13	107.33	111.63
5	A	707	NAG	O4-C4-C3	-2.08	105.54	110.35

There are no chirality outliers.

All (36) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	701	FAD	O4B-C4B-C5B-O5B
4	A	701	FAD	N10-C1'-C2'-O2'
4	B	701	FAD	O4B-C4B-C5B-O5B
4	B	701	FAD	O4'-C4'-C5'-O5'
8	A	711	B3P	C9-C8-N2-C2
8	A	711	B3P	C10-C8-N2-C2
8	A	711	B3P	C11-C8-N2-C2
8	A	711	B3P	C10-C8-C9-O1
8	B	709	B3P	C5-C4-N1-C3
8	B	709	B3P	C6-C4-N1-C3
8	B	709	B3P	O3-C11-C8-N2
8	B	709	B3P	O3-C11-C8-C9
8	B	709	B3P	O3-C11-C8-C10
5	A	707	NAG	O5-C5-C6-O6
5	B	706	NAG	O5-C5-C6-O6
4	B	701	FAD	C3B-C4B-C5B-O5B
5	A	707	NAG	C4-C5-C6-O6
5	A	708	NAG	O5-C5-C6-O6
5	B	706	NAG	C4-C5-C6-O6
4	A	701	FAD	C3B-C4B-C5B-O5B
8	B	709	B3P	C3-C1-C2-N2
5	A	708	NAG	C4-C5-C6-O6

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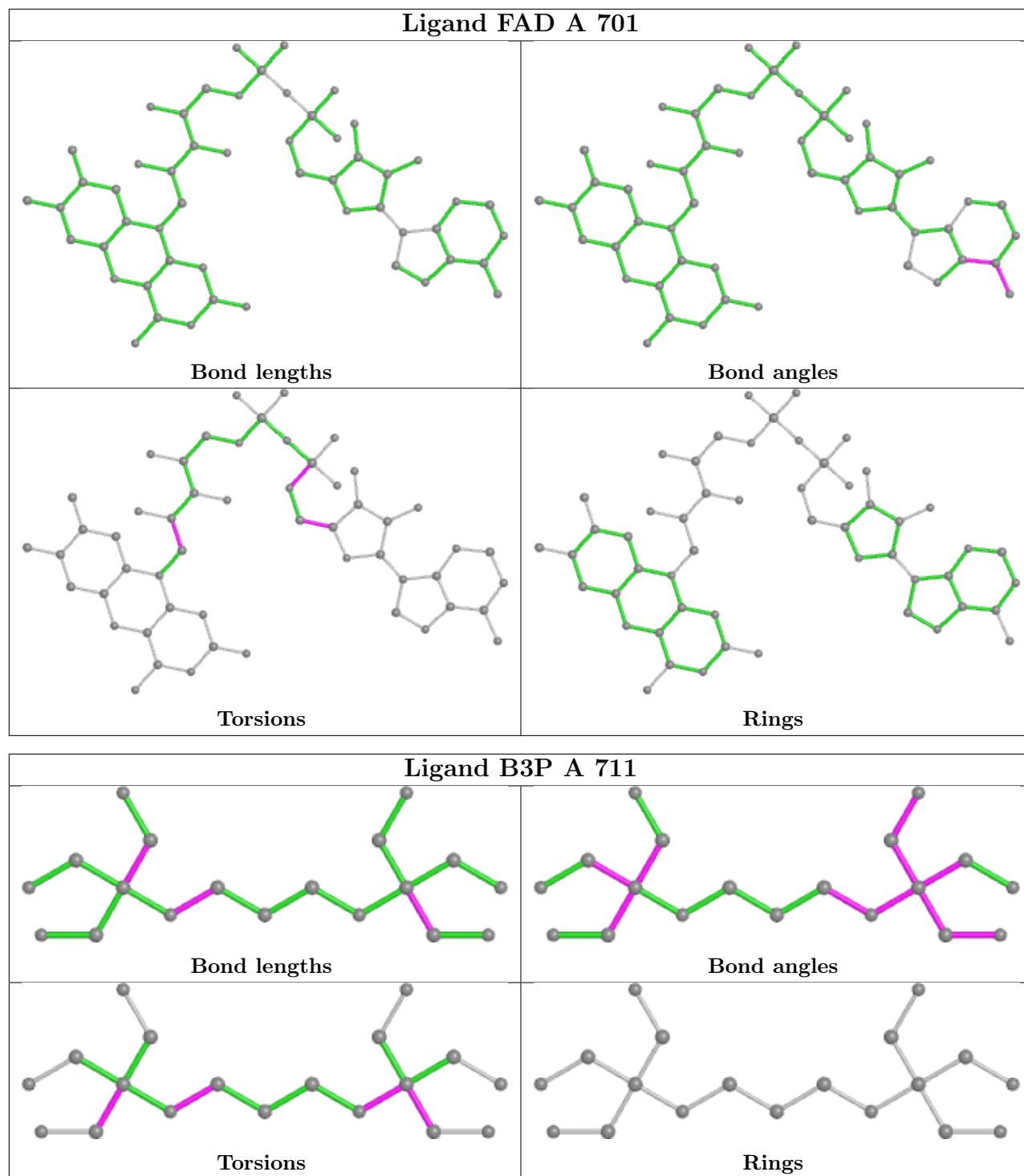
Mol	Chain	Res	Type	Atoms
8	A	711	B3P	C1-C3-N1-C4
8	A	711	B3P	N1-C4-C7-O6
8	B	709	B3P	N1-C4-C7-O6
9	A	712	EDO	O1-C1-C2-O2
8	B	709	B3P	O2-C10-C8-C11
8	B	709	B3P	O2-C10-C8-C9
8	B	709	B3P	C7-C4-N1-C3
8	B	709	B3P	C2-C1-C3-N1
8	B	709	B3P	C5-C4-C7-O6
8	A	711	B3P	C11-C8-C9-O1
8	B	709	B3P	O2-C10-C8-N2
4	A	701	FAD	C5B-O5B-PA-O3P
4	B	701	FAD	C5B-O5B-PA-O1A
8	B	709	B3P	C9-C8-N2-C2

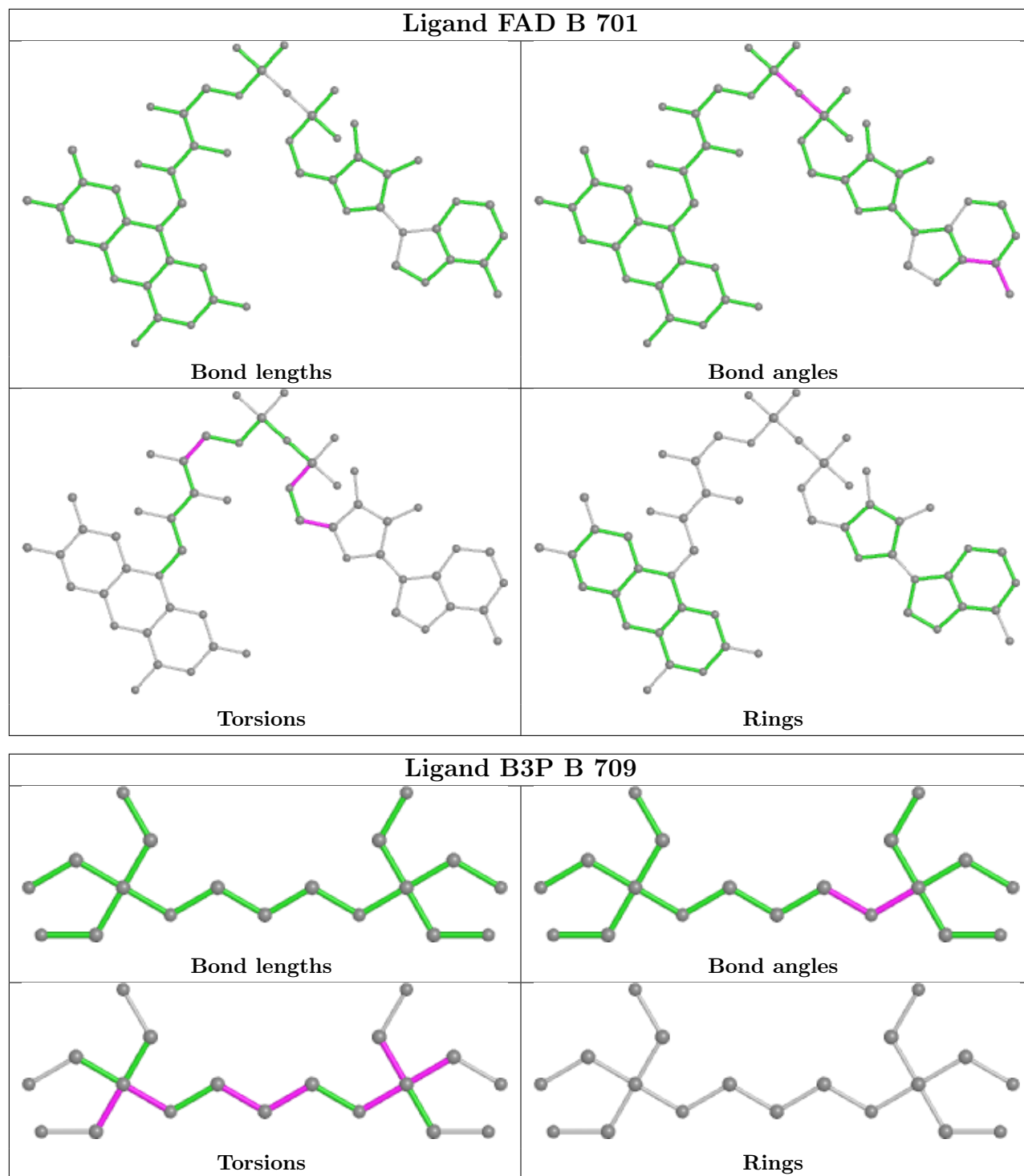
There are no ring outliers.

6 monomers are involved in 36 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	707	NAG	2	0
5	A	708	NAG	7	0
4	A	701	FAD	4	0
8	A	711	B3P	13	0
4	B	701	FAD	3	0
8	B	709	B3P	7	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

There are no chain breaks in this entry.



## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	613/647 (94%)	-0.02	5 (0%) 86 84	18, 27, 46, 96	0
1	B	613/647 (94%)	-0.06	3 (0%) 91 89	19, 25, 42, 74	0
All	All	1226/1294 (94%)	-0.04	8 (0%) 87 86	18, 26, 44, 96	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	420	TYR	5.2
1	B	420	TYR	4.6
1	A	389	THR	2.9
1	A	228	ALA	2.7
1	B	150	GLN	2.6
1	B	151	SER	2.3
1	A	149	GLU	2.3
1	A	85	ILE	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.

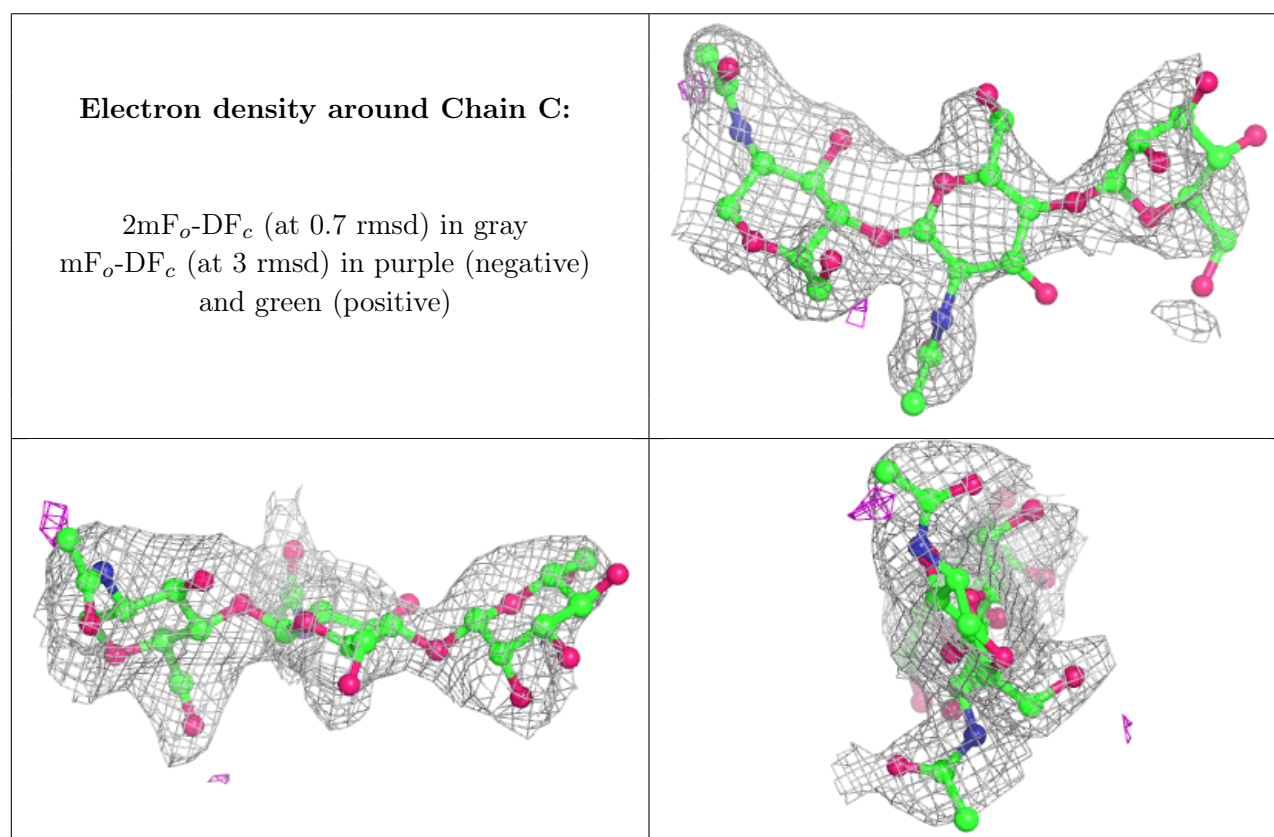
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	BMA	C	3	11/12	0.68	0.24	74,75,79,80	0

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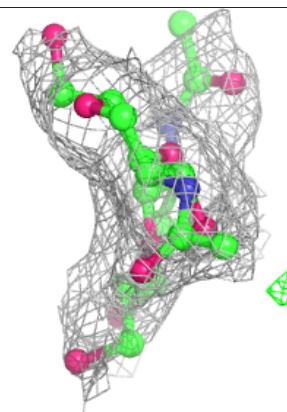
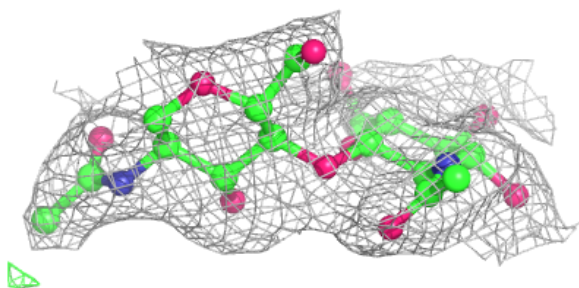
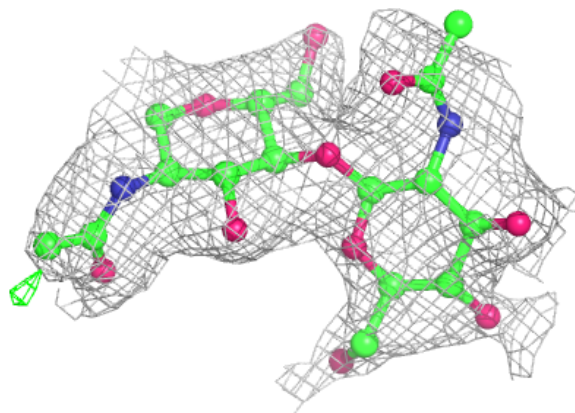
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	NAG	F	2	14/15	0.80	0.18	57,67,70,71	0
3	NAG	E	2	14/15	0.81	0.22	45,59,69,69	0
2	NAG	C	2	14/15	0.82	0.22	59,65,73,74	0
3	NAG	D	2	14/15	0.83	0.21	55,61,71,71	0
2	NAG	C	1	14/15	0.89	0.18	31,44,50,59	0
3	NAG	E	1	14/15	0.89	0.15	34,40,50,53	0
3	NAG	F	1	14/15	0.93	0.13	25,38,49,49	0
3	NAG	D	1	14/15	0.93	0.13	25,37,48,53	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.



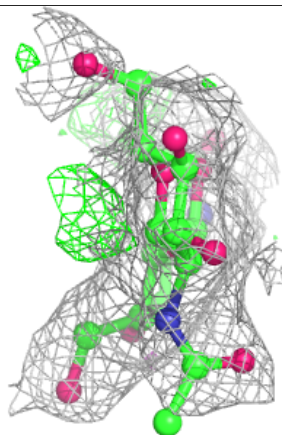
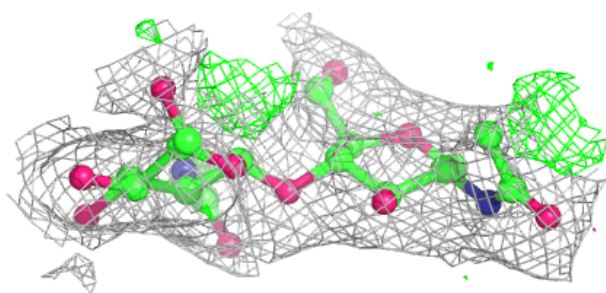
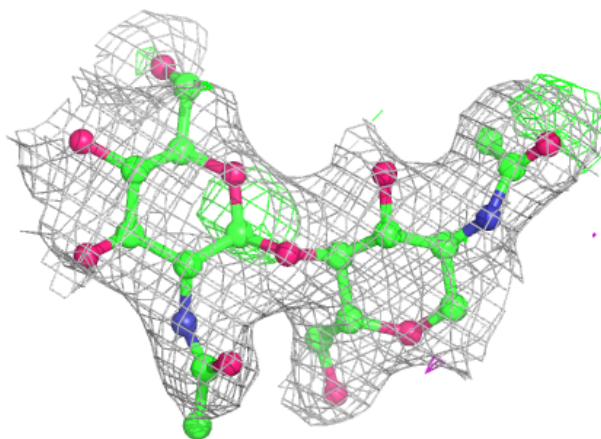
**Electron density around Chain D:**

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

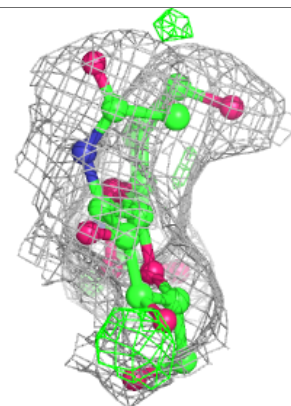
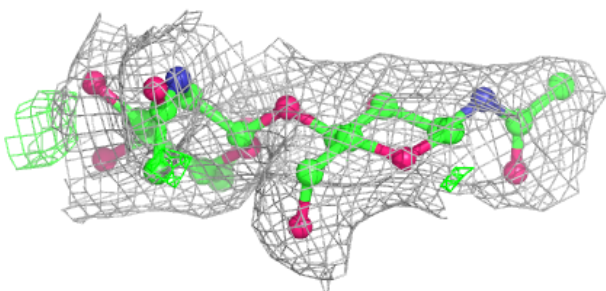
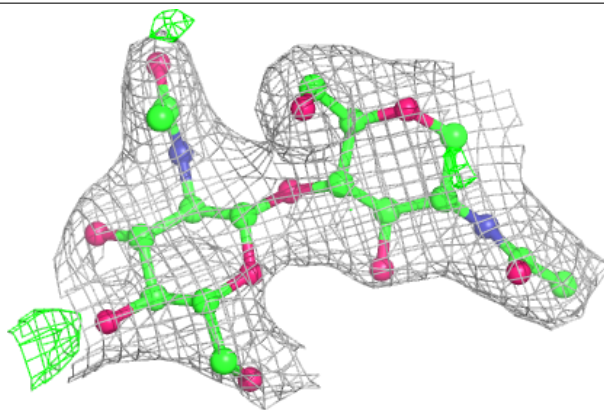


**Electron density around Chain E:**

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

**Electron density around Chain F:**

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



## 6.4 Ligands

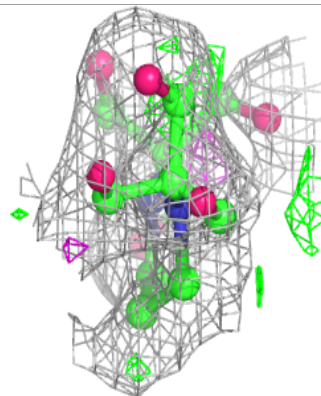
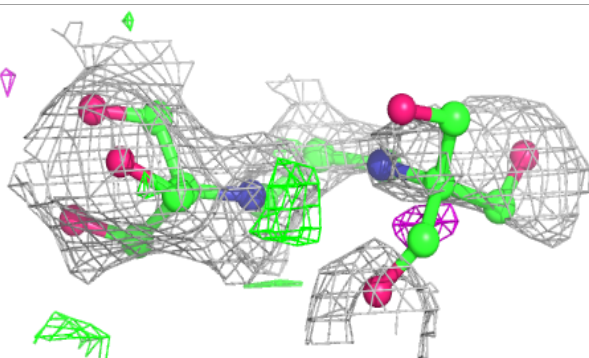
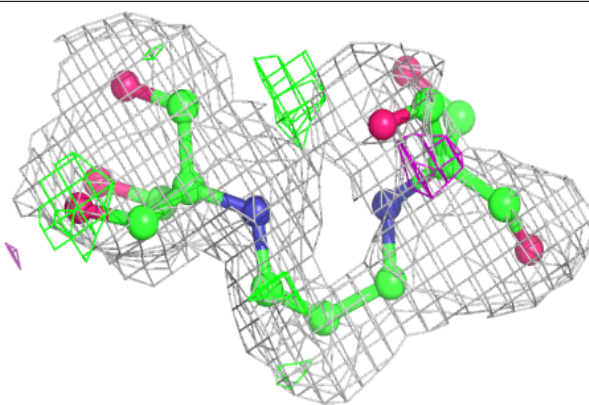
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<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
5	NAG	A	708	14/15	0.73	0.40	61,74,88,89	0
8	B3P	A	711	19/19	0.73	0.29	39,66,79,83	0
8	B3P	B	709	19/19	0.74	0.29	50,57,76,79	0
9	EDO	A	712	4/4	0.75	0.42	26,27,29,31	0
5	NAG	B	706	14/15	0.79	0.27	56,66,72,76	0
5	NAG	A	707	14/15	0.82	0.17	60,68,79,80	0
6	SO4	A	709	5/5	0.87	0.15	78,84,86,86	0
10	144	B	708	8/8	0.89	0.14	54,61,64,68	0
4	FAD	A	701	53/53	0.96	0.14	18,23,30,37	0
4	FAD	B	701	53/53	0.96	0.15	18,23,32,36	0
7	CA	B	707	1/1	0.98	0.11	20,20,20,20	0
7	CA	A	710	1/1	0.99	0.07	19,19,19,19	0

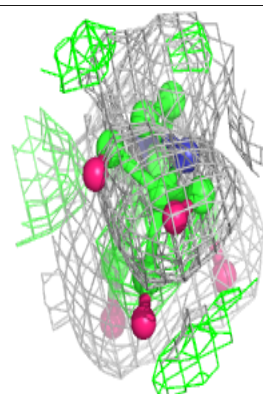
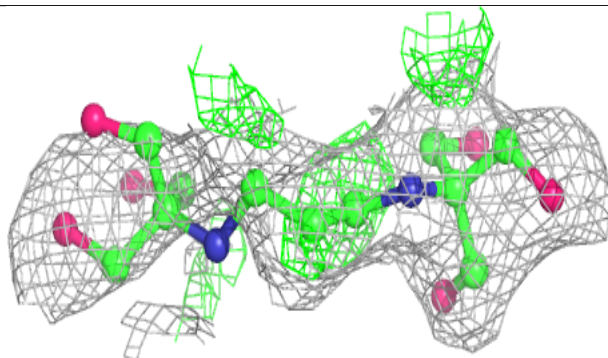
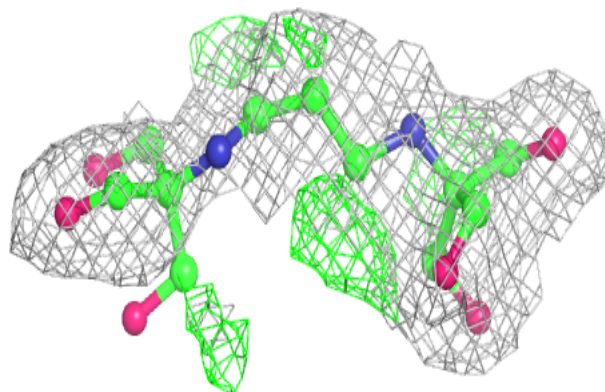
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around B3P A 711:**

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

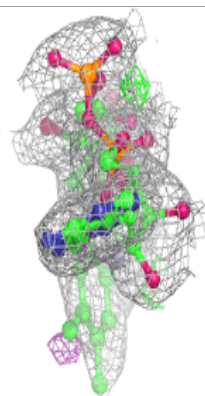
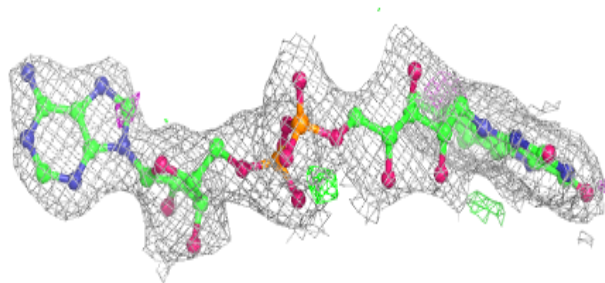
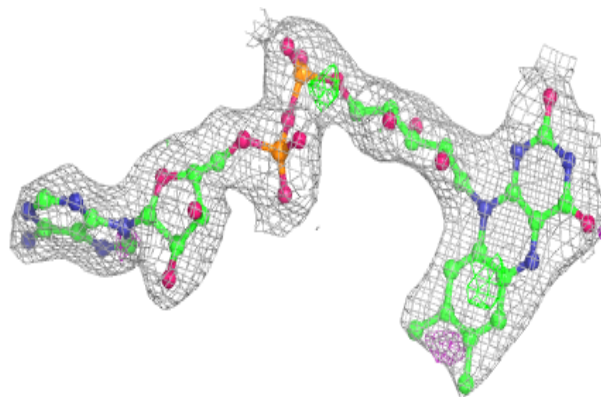
**Electron density around B3P B 709:**

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

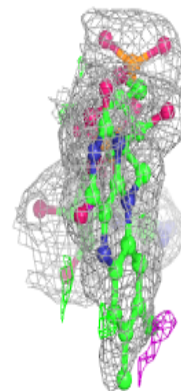
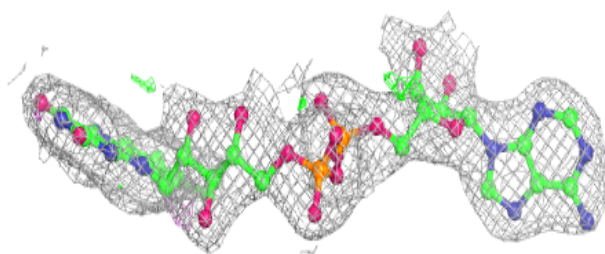
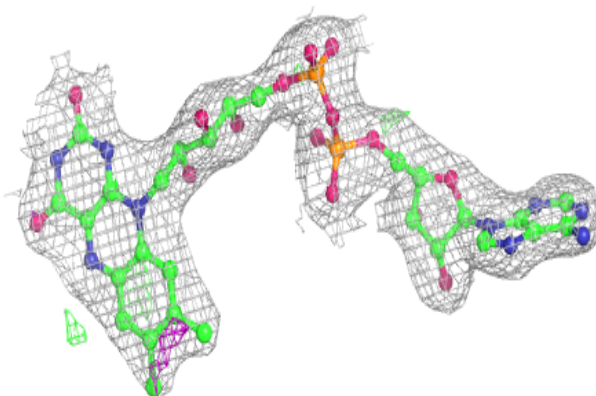


**Electron density around FAD A 701:**

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

**Electron density around FAD B 701:**

$2mF_o-DF_c$  (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.