



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 07:33 PM UTC

PDB ID : 3QMO / pdb_00003qmo
Title : X-ray crystal structure of NS-398 bound to the cyclooxygenase channel of cyclooxygenase-2
Authors : Vecchio, A.J.; Malkowski, M.G.
Deposited on : 2011-02-04
Resolution : 3.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

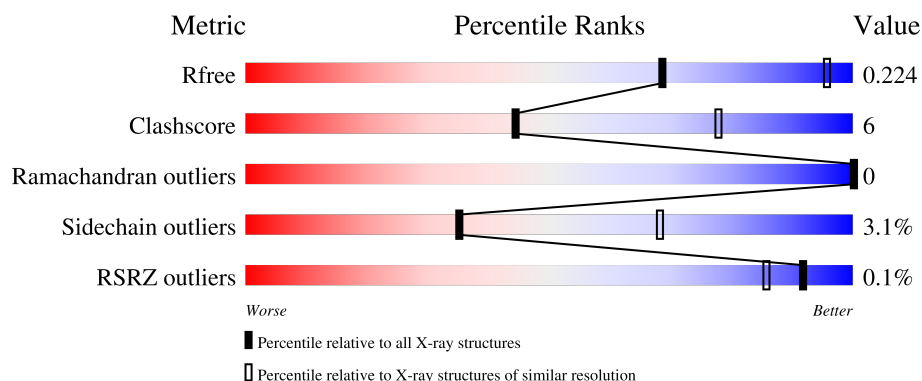
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	610	
1	B	610	
2	C	2	
2	E	2	
3	D	2	

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Mol	Chain	Length	Quality of chain
4	F	3	 33% 67%

2 Entry composition [i](#)

There are 11 unique types of molecules in this entry. The entry contains 9237 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 Prostaglandin G/H synthase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	551	Total	C	N	O	S	0	3	0
			4392	2840	736	791	25			
1	B	551	Total	C	N	O	S	0	0	0
			4379	2838	729	787	25			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	29	HIS	-	expression tag	UNP Q05769
A	30	HIS	-	expression tag	UNP Q05769
A	31	HIS	-	expression tag	UNP Q05769
A	32	HIS	-	expression tag	UNP Q05769
A	33	HIS	-	expression tag	UNP Q05769
A	34	HIS	-	expression tag	UNP Q05769
A	594	ALA	ASN	engineered mutation	UNP Q05769
B	29	HIS	-	expression tag	UNP Q05769
B	30	HIS	-	expression tag	UNP Q05769
B	31	HIS	-	expression tag	UNP Q05769
B	32	HIS	-	expression tag	UNP Q05769
B	33	HIS	-	expression tag	UNP Q05769
B	34	HIS	-	expression tag	UNP Q05769
B	594	ALA	ASN	engineered mutation	UNP Q05769

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



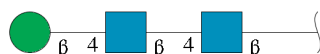
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	E	2	Total	C	N	O	0	0	0
			28	16	2	10			

- 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	D	2	Total	C	N	O	0	0	0
			28	16	2	10			

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



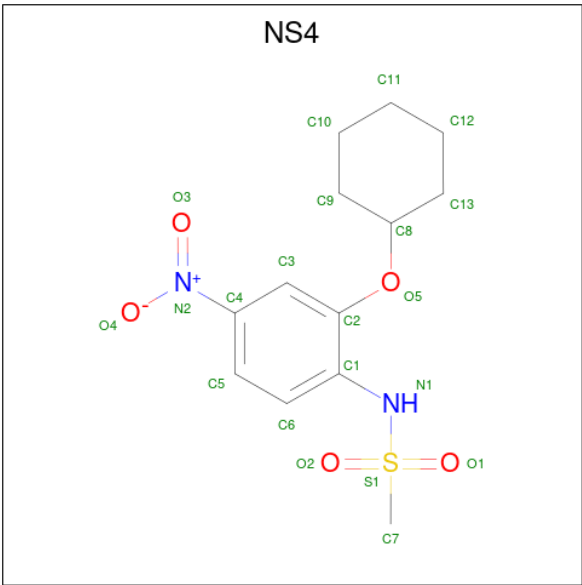
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	F	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



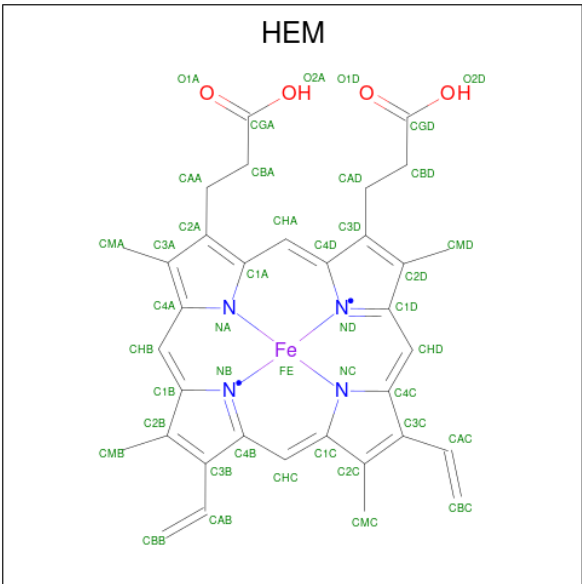
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is N-[2-(cyclohexyloxy)-4-nitrophenyl]methanesulfonamide (CCD ID: NS4) (formula: C₁₃H₁₈N₂O₅S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	N	O	S	0	0
			21	13	2	5	1		
6	B	1	Total	C	N	O	S	0	0
			21	13	2	5	1		

- Molecule 7 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: C₃₄H₃₂FeN₄O₄).



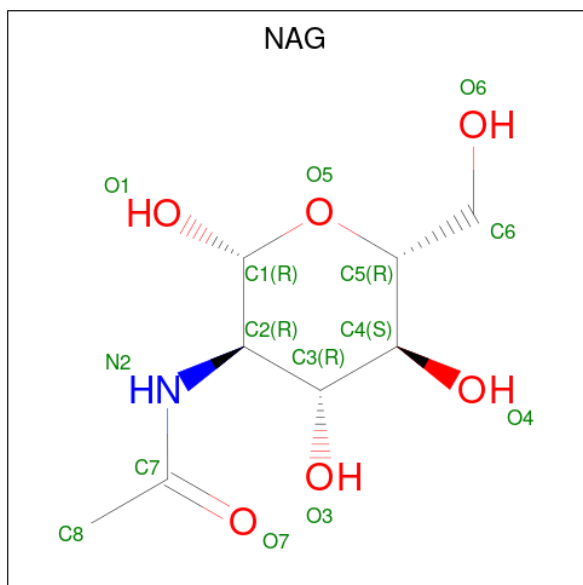
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

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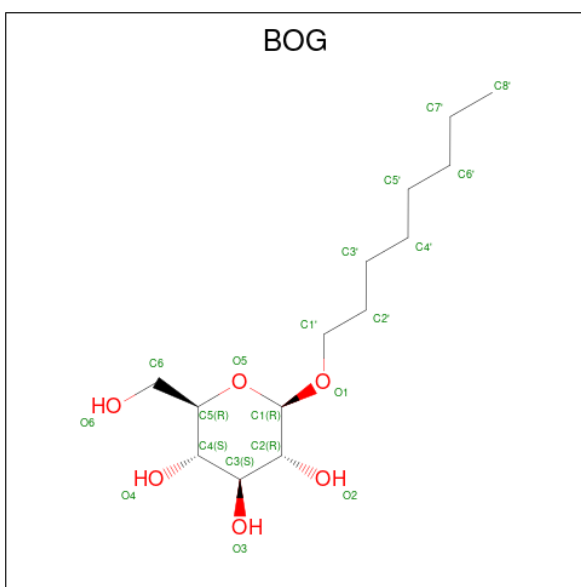
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 8 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



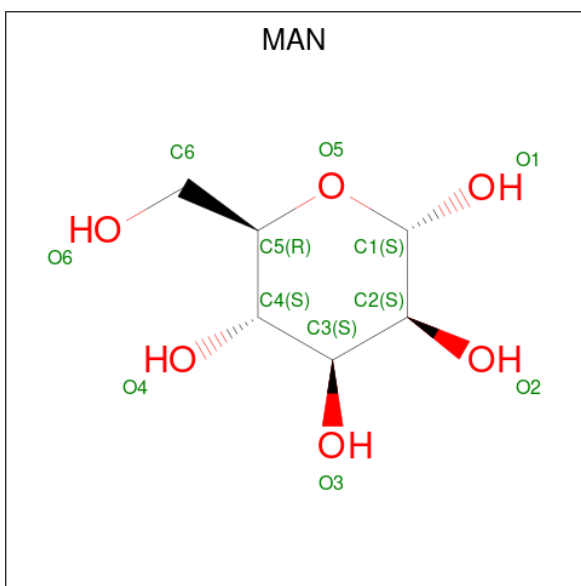
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	A	1	Total	C	N	O	0	0
			14	8	1	5		
8	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 9 is octyl beta-D-glucopyranoside (CCD ID: BOG) (formula: $C_{14}H_{28}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			20	14	6		

- Molecule 10 is alpha-D-mannopyranose (CCD ID: MAN) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	B	1	Total	C	O	0	0
			11	6	5		

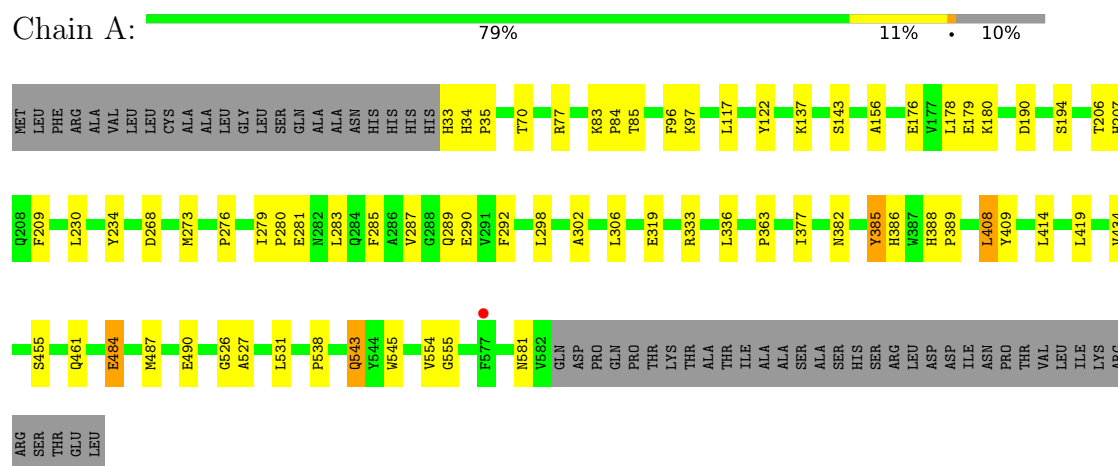
- Molecule 11 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	A	59	Total 59	O 59	0	0
11	B	61	Total 61	O 61	0	0

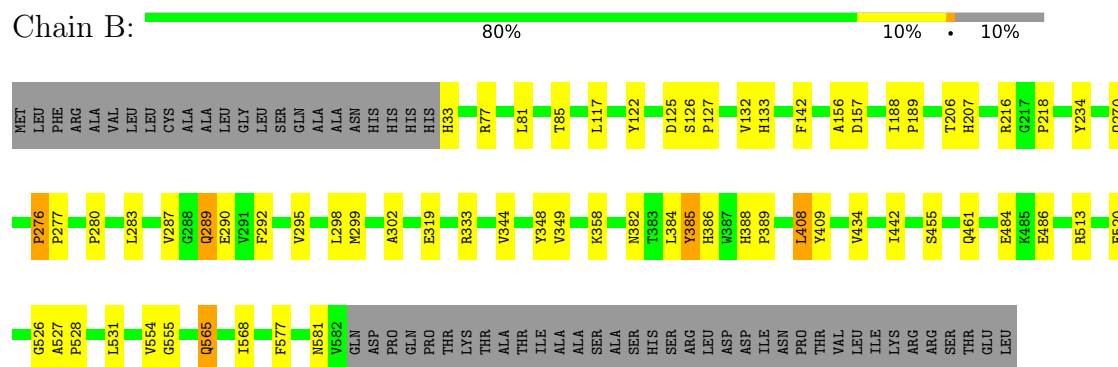
3 Residue-property plots [i](#)

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: Prostaglandin G/H synthase 2



• Molecule 1: Prostaglandin G/H synthase 2



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



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

Chain E:  100%

MAG1
BDG2

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

Chain D:  50% 50%

MAG1
MAG2

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

Chain F:  33% 67%

MAG1
MAG2
BMA3

4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	120.43Å 131.21Å 179.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.99 – 3.00 19.99 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (19.99-3.00) 99.4 (19.99-3.00)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 2.98Å)	Xtriage
Refinement program	REFMAC 5.5.0088	Depositor
R, R_{free}	0.176 , 0.225 0.175 , 0.224	Depositor DCC
R_{free} test set	1452 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	47.0	Xtriage
Anisotropy	0.143	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 48.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9237	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.36% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, NS4, HEM, NAG, GOL, MAN, NDG, BOG

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	0.63	0/4527	0.88	2/6158 (0.0%)
1	B	0.62	0/4508	0.89	3/6130 (0.0%)
All	All	0.62	0/9035	0.89	5/12288 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	276	PRO	CA-C-N	5.44	125.52	119.32
1	B	276	PRO	C-N-CA	5.44	125.52	119.32
1	B	276	PRO	O-C-N	5.18	123.59	121.15
1	A	276	PRO	CA-C-N	5.01	125.04	119.32
1	A	276	PRO	C-N-CA	5.01	125.04	119.32

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	4392	0	4171	52	0
1	B	4379	0	4172	49	0
2	C	28	0	24	1	0
2	E	28	0	24	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	28	0	24	1	0
4	F	39	0	34	4	0
5	A	18	0	24	3	0
5	B	18	0	24	3	0
6	A	21	0	18	5	0
6	B	21	0	18	6	0
7	A	43	0	30	2	0
7	B	43	0	30	3	0
8	A	14	0	13	0	0
8	B	14	0	13	0	0
9	A	20	0	28	1	0
10	B	11	0	10	0	0
11	A	59	0	0	7	0
11	B	61	0	0	6	0
All	All	9237	0	8657	107	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 6.

All (107) 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:137[B]:LYS:H	1:A:137[B]:LYS:CD	1.51	1.18
1:A:137[B]:LYS:N	1:A:137[B]:LYS:HD2	1.50	1.14
7:A:705:HEM:HHA	7:A:705:HEM:HBA1	1.43	0.95
1:B:156:ALA:H	5:B:703:GOL:H32	1.33	0.91
1:A:543[A]:GLN:HA	1:A:543[A]:GLN:OE1	1.71	0.90
1:A:137[B]:LYS:H	1:A:137[B]:LYS:HD2	0.74	0.89
1:A:137[B]:LYS:HD3	11:B:830:HOH:O	1.84	0.77
6:A:704:NS4:O2	6:A:704:NS4:H6	1.90	0.71
1:A:319:GLU:HG3	1:A:554:VAL:HG11	1.73	0.69
1:B:319:GLU:HG3	1:B:554:VAL:HG11	1.75	0.67
1:B:526:GLY:HA3	6:B:705:NS4:H12	1.77	0.66
7:B:706:HEM:HBC2	7:B:706:HEM:HHD	1.77	0.66
1:B:295:VAL:HG21	7:B:706:HEM:HBB2	1.79	0.64
1:A:179:GLU:HB3	9:A:707:BOG:H3'1	1.81	0.63
1:A:156:ALA:H	5:A:701:GOL:H32	1.64	0.63
1:A:287:VAL:HG11	1:A:302:ALA:HB1	1.80	0.62
1:B:531:LEU:HD11	6:B:705:NS4:H7	1.82	0.61
11:A:852:HOH:O	4:F:3:BMA:H61	2.01	0.60
1:B:216:ARG:NH1	4:F:2:NAG:H83	2.16	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:389:PRO:HB2	1:A:434:VAL:HA	1.83	0.60
1:A:268:ASP:CG	11:A:840:HOH:O	2.45	0.60
1:B:289:GLN:HG2	1:B:292:PHE:CE1	2.38	0.58
1:B:287:VAL:HG11	1:B:302:ALA:HB1	1.86	0.57
1:A:382:ASN:O	1:A:386:HIS:HD2	1.88	0.57
1:A:77:ARG:NH1	11:A:802:HOH:O	2.38	0.57
1:A:289:GLN:HG2	1:A:292:PHE:CE1	2.40	0.56
1:A:527:ALA:HB1	6:A:704:NS4:H7B	1.88	0.55
1:A:414:LEU:HD11	1:A:419:LEU:HD12	1.89	0.54
1:B:389:PRO:HB2	1:B:434:VAL:HA	1.90	0.53
1:B:218:PRO:HD2	11:B:855:HOH:O	2.10	0.52
1:B:77:ARG:NH1	11:B:803:HOH:O	2.43	0.51
1:A:143:SER:O	5:A:703:GOL:H12	2.10	0.51
6:B:705:NS4:H6	6:B:705:NS4:O2	2.10	0.51
1:A:234:TYR:CE2	1:A:333:ARG:HG3	2.47	0.50
1:B:382:ASN:O	1:B:386:HIS:HD2	1.95	0.49
1:B:531:LEU:HD21	6:B:705:NS4:H7A	1.94	0.49
1:B:234:TYR:CE2	1:B:333:ARG:HG3	2.48	0.48
1:A:137[B]:LYS:CD	11:B:830:HOH:O	2.54	0.48
1:B:565:GLN:OE1	1:B:577:PHE:HA	2.13	0.48
1:B:292:PHE:O	1:B:299:MET:HE2	2.13	0.48
1:B:531:LEU:HD11	6:B:705:NS4:C7	2.44	0.48
1:A:531:LEU:HD11	6:A:704:NS4:C7	2.45	0.47
1:A:137[B]:LYS:CD	1:A:137[B]:LYS:N	2.29	0.47
1:A:554:VAL:HG13	1:A:555:GLY:N	2.30	0.47
1:A:96:PHE:O	1:A:97:LYS:C	2.57	0.46
5:A:702:GOL:H32	7:A:705:HEM:NA	2.30	0.46
1:B:554:VAL:HG13	1:B:555:GLY:N	2.29	0.46
1:A:408:LEU:O	1:A:409:TYR:HB2	2.14	0.46
1:A:117:LEU:HD23	1:A:531:LEU:HD13	1.98	0.46
1:B:207:HIS:O	1:B:289:GLN:NE2	2.49	0.46
11:A:838:HOH:O	3:D:1:NAG:H82	2.15	0.45
1:B:388:HIS:N	1:B:389:PRO:CD	2.80	0.45
1:A:382:ASN:O	1:A:386:HIS:CD2	2.68	0.45
1:A:543[A]:GLN:HG3	1:B:125:ASP:OD1	2.16	0.45
1:B:206:THR:HG21	1:B:385:TYR:CE2	2.51	0.45
1:B:216:ARG:HG2	4:F:2:NAG:H81	1.98	0.45
1:B:513:ARG:HH21	1:B:520:GLU:HG3	1.81	0.45
1:A:268:ASP:CB	11:A:841:HOH:O	2.64	0.45
1:B:280:PRO:HG2	1:B:283:LEU:HD12	1.99	0.45
1:A:33:HIS:ND1	1:A:33:HIS:C	2.75	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:531:LEU:HD11	6:A:704:NS4:H7	1.98	0.44
1:B:527:ALA:HB1	6:B:705:NS4:H7B	1.99	0.44
1:B:117:LEU:HD23	1:B:531:LEU:HD13	1.99	0.44
1:B:358:LYS:HE2	1:B:358:LYS:HB3	1.76	0.44
1:B:289:GLN:HG2	1:B:292:PHE:CZ	2.52	0.44
1:A:292:PHE:CD1	1:A:298:LEU:HD13	2.53	0.44
1:B:384:LEU:HD12	1:B:384:LEU:C	2.44	0.43
2:C:2:NDG:H2	2:C:2:NDG:H8C1	1.84	0.43
1:A:190:ASP:OD2	1:A:194:SER:OG	2.34	0.43
1:A:137[B]:LYS:NZ	11:A:801:HOH:O	2.34	0.43
1:B:270:GLN:HA	1:B:270:GLN:HE21	1.84	0.43
1:B:527:ALA:HB3	1:B:528:PRO:HD3	2.01	0.43
1:A:280:PRO:HG2	1:A:283:LEU:HD12	2.00	0.43
1:A:279:ILE:HA	1:A:280:PRO:HD2	1.84	0.43
1:A:363:PRO:HG2	1:A:545:TRP:CD2	2.54	0.42
1:A:538:PRO:HG3	1:B:142:PHE:CE2	2.54	0.42
1:B:156:ALA:N	5:B:703:GOL:H32	2.15	0.42
1:B:126:SER:HA	1:B:127:PRO:C	2.44	0.42
5:B:704:GOL:H31	7:B:706:HEM:C1A	2.54	0.42
1:A:77:ARG:CZ	11:A:802:HOH:O	2.67	0.42
1:B:77:ARG:NH2	11:B:810:HOH:O	2.53	0.42
1:A:34:HIS:HA	1:A:35:PRO:HD3	1.96	0.42
1:B:344:VAL:HA	1:B:348:TYR:HB3	2.01	0.42
1:B:216:ARG:HH11	4:F:2:NAG:H83	1.83	0.42
1:A:484:GLU:CD	1:A:487:MET:H	2.27	0.41
1:A:83:LYS:HA	1:A:84:PRO:HD3	1.96	0.41
1:A:209:PHE:HB2	1:A:377:ILE:HG13	2.02	0.41
1:A:230:LEU:CD2	1:A:336:LEU:HB3	2.50	0.41
1:A:306:LEU:C	1:A:306:LEU:HD23	2.45	0.41
1:A:388:HIS:N	1:A:389:PRO:CD	2.84	0.41
1:B:132:VAL:HG13	1:B:133:HIS:CD2	2.55	0.41
1:B:388:HIS:N	1:B:389:PRO:HD2	2.35	0.41
1:A:273:MET:HE2	1:A:285:PHE:O	2.20	0.41
1:B:333:ARG:NH1	11:B:807:HOH:O	2.50	0.41
1:A:207:HIS:O	1:A:289:GLN:NE2	2.54	0.41
1:B:276:PRO:HA	1:B:277:PRO:HD3	1.91	0.41
1:B:33:HIS:ND1	1:B:33:HIS:C	2.79	0.41
1:A:538:PRO:HG3	1:B:142:PHE:CZ	2.56	0.40
1:B:344:VAL:O	1:B:349:VAL:HG23	2.21	0.40
1:B:442:ILE:H	1:B:442:ILE:HD12	1.85	0.40
1:A:526:GLY:HA3	6:A:704:NS4:H12	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:292:PHE:CD1	1:B:298:LEU:HD13	2.56	0.40
1:B:408:LEU:O	1:B:409:TYR:HB2	2.21	0.40
1:B:188:ILE:HA	1:B:189:PRO:HD2	1.87	0.40
1:A:176:GLU:HG2	1:A:180:LYS:HD2	2.03	0.40
1:A:180:LYS:HE3	1:A:490:GLU:OE1	2.22	0.40
1:A:206:THR:HG21	1:A:385:TYR:CE2	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	552/610 (90%)	530 (96%)	22 (4%)	0	100	100
1	B	549/610 (90%)	529 (96%)	20 (4%)	0	100	100
All	All	1101/1220 (90%)	1059 (96%)	42 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	461/542 (85%)	447 (97%)	14 (3%)	36	69
1	B	460/542 (85%)	445 (97%)	15 (3%)	33	67

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	921/1084 (85%)	892 (97%)	29 (3%)	35 68

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	70	THR
1	A	85	THR
1	A	122	TYR
1	A	178	LEU
1	A	281	GLU
1	A	290	GLU
1	A	385	TYR
1	A	408	LEU
1	A	455	SER
1	A	461	GLN
1	A	484	GLU
1	A	543[A]	GLN
1	A	543[B]	GLN
1	A	581	ASN
1	B	81	LEU
1	B	85	THR
1	B	122	TYR
1	B	157	ASP
1	B	289	GLN
1	B	290	GLU
1	B	385	TYR
1	B	408	LEU
1	B	455	SER
1	B	461	GLN
1	B	484	GLU
1	B	486	GLU
1	B	565	GLN
1	B	568	ILE
1	B	581	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	GLN
1	A	270	GLN
1	A	284	GLN

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Mol	Chain	Res	Type
1	B	42	GLN
1	B	207	HIS
1	B	270	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 ⓘ

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	2,1	14,14,15	0.64	0	17,19,21	1.39	3 (17%)
2	NDG	C	2	2	14,14,15	0.78	1 (7%)	17,19,21	2.32	8 (47%)
3	NAG	D	1	1,3	14,14,15	0.71	0	17,19,21	1.45	3 (17%)
3	NAG	D	2	3	14,14,15	0.55	0	17,19,21	1.36	4 (23%)
2	NAG	E	1	2,1	14,14,15	0.62	0	17,19,21	0.70	0
2	NDG	E	2	2	14,14,15	0.54	0	17,19,21	0.89	0
4	NAG	F	1	1,4	14,14,15	1.90	1 (7%)	17,19,21	2.59	6 (35%)
4	NAG	F	2	4	14,14,15	0.69	0	17,19,21	2.58	5 (29%)
4	BMA	F	3	4	11,11,12	0.82	0	15,15,17	1.84	3 (20%)

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	2,1	-	0/6/23/26	0/1/1/1
2	NDG	C	2	2	-	4/6/23/26	0/1/1/1
3	NAG	D	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	D	2	3	-	4/6/23/26	0/1/1/1
2	NAG	E	1	2,1	-	2/6/23/26	0/1/1/1
2	NDG	E	2	2	-	0/6/23/26	0/1/1/1
4	NAG	F	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	F	2	4	-	4/6/23/26	0/1/1/1
4	BMA	F	3	4	-	2/2/19/22	1/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	1	NAG	O5-C1	-6.45	1.32	1.43
2	C	2	NDG	C1-C2	2.47	1.55	1.52

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	1	NAG	C1-O5-C5	-7.78	101.76	112.19
2	C	2	NDG	C3-C4-C5	-5.08	101.03	110.23
4	F	3	BMA	C1-O5-C5	4.89	118.74	112.19
4	F	2	NAG	O4-C4-C5	4.79	121.12	109.32
4	F	2	NAG	O5-C5-C6	4.58	116.58	107.66
4	F	2	NAG	C4-C3-C2	-4.48	104.45	111.02
4	F	2	NAG	C3-C4-C5	-4.34	102.37	110.23
4	F	2	NAG	O4-C4-C3	4.03	119.87	110.38
4	F	1	NAG	O5-C1-C2	-3.91	105.23	111.29
4	F	1	NAG	O5-C5-C6	3.85	115.15	107.66
2	C	2	NDG	O5-C1-C2	3.68	116.98	111.29
2	C	2	NDG	C1-O5-C5	3.40	116.74	112.19
3	D	1	NAG	C2-N2-C7	3.34	127.37	122.90
2	C	1	NAG	C3-C4-C5	-3.09	104.63	110.23
4	F	3	BMA	C1-C2-C3	2.94	113.93	109.64
4	F	3	BMA	O2-C2-C1	2.90	115.86	109.22
3	D	1	NAG	C1-C2-N2	-2.87	105.91	110.43
2	C	2	NDG	O4-C4-C5	2.75	116.11	109.32
3	D	2	NAG	C1-O5-C5	2.75	115.88	112.19
4	F	1	NAG	O4-C4-C5	2.72	116.01	109.32
2	C	2	NDG	C8-C7-N2	2.53	120.32	116.12
2	C	1	NAG	O5-C5-C6	2.47	112.47	107.66
2	C	1	NAG	O4-C4-C3	2.42	116.08	110.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1	NAG	O7-C7-N2	2.40	126.23	121.98
3	D	2	NAG	C2-N2-C7	-2.40	119.69	122.90
2	C	2	NDG	O5-C5-C6	2.39	112.32	107.66
4	F	1	NAG	C2-N2-C7	-2.28	119.84	122.90
3	D	2	NAG	O5-C1-C2	-2.16	107.95	111.29
2	C	2	NDG	O5-C5-C4	-2.16	105.58	110.83
2	C	2	NDG	C6-C5-C4	2.07	118.09	113.02
3	D	2	NAG	O4-C4-C5	2.04	114.34	109.32
4	F	1	NAG	C3-C4-C5	-2.03	106.55	110.23

There are no chirality outliers.

All (16) torsion outliers are listed below:

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

All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	F	3	BMA	C1-C2-C3-C4-C5-O5

4 monomers are involved in 6 short contacts:

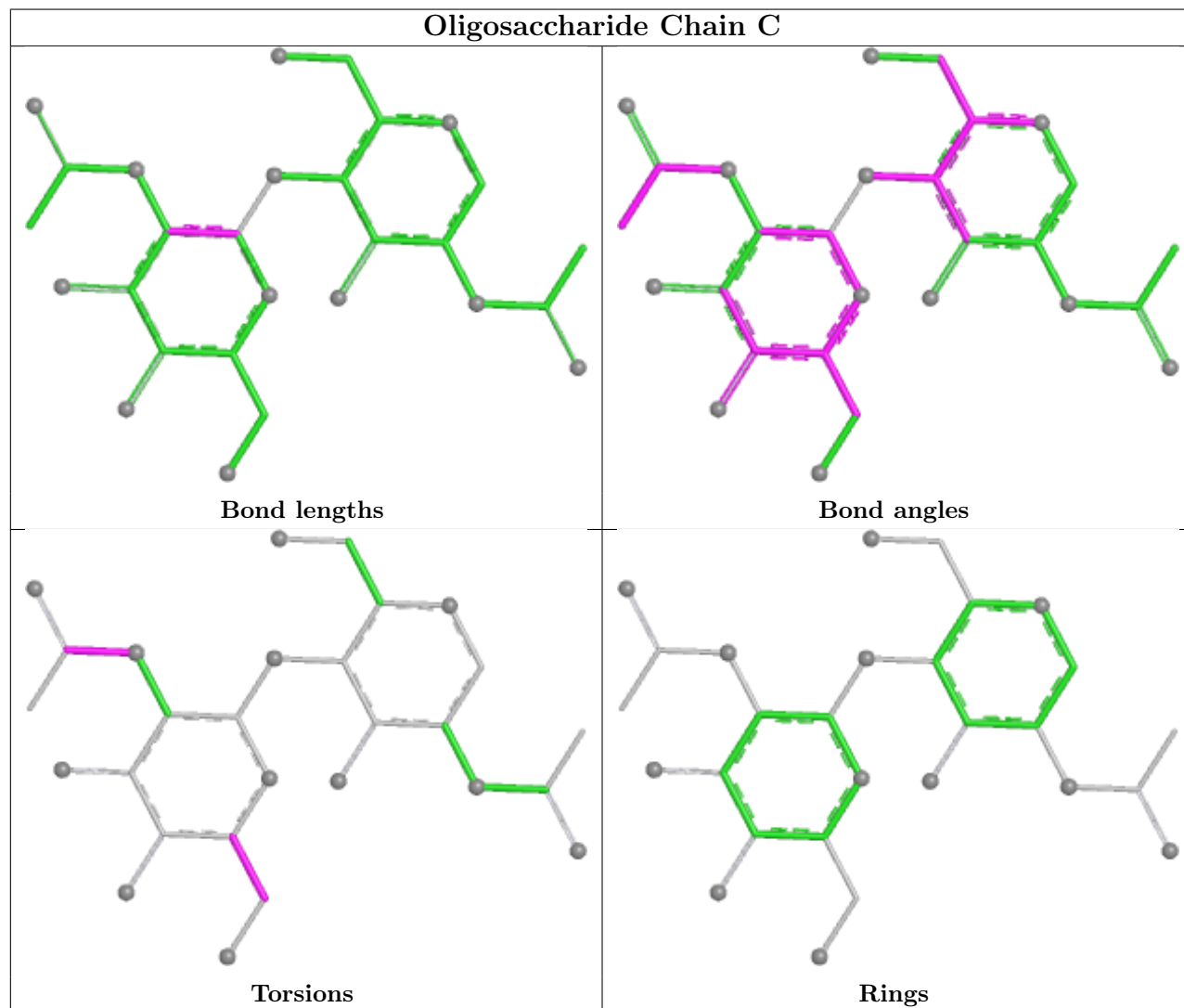
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	1	NAG	1	0
4	F	2	NAG	3	0

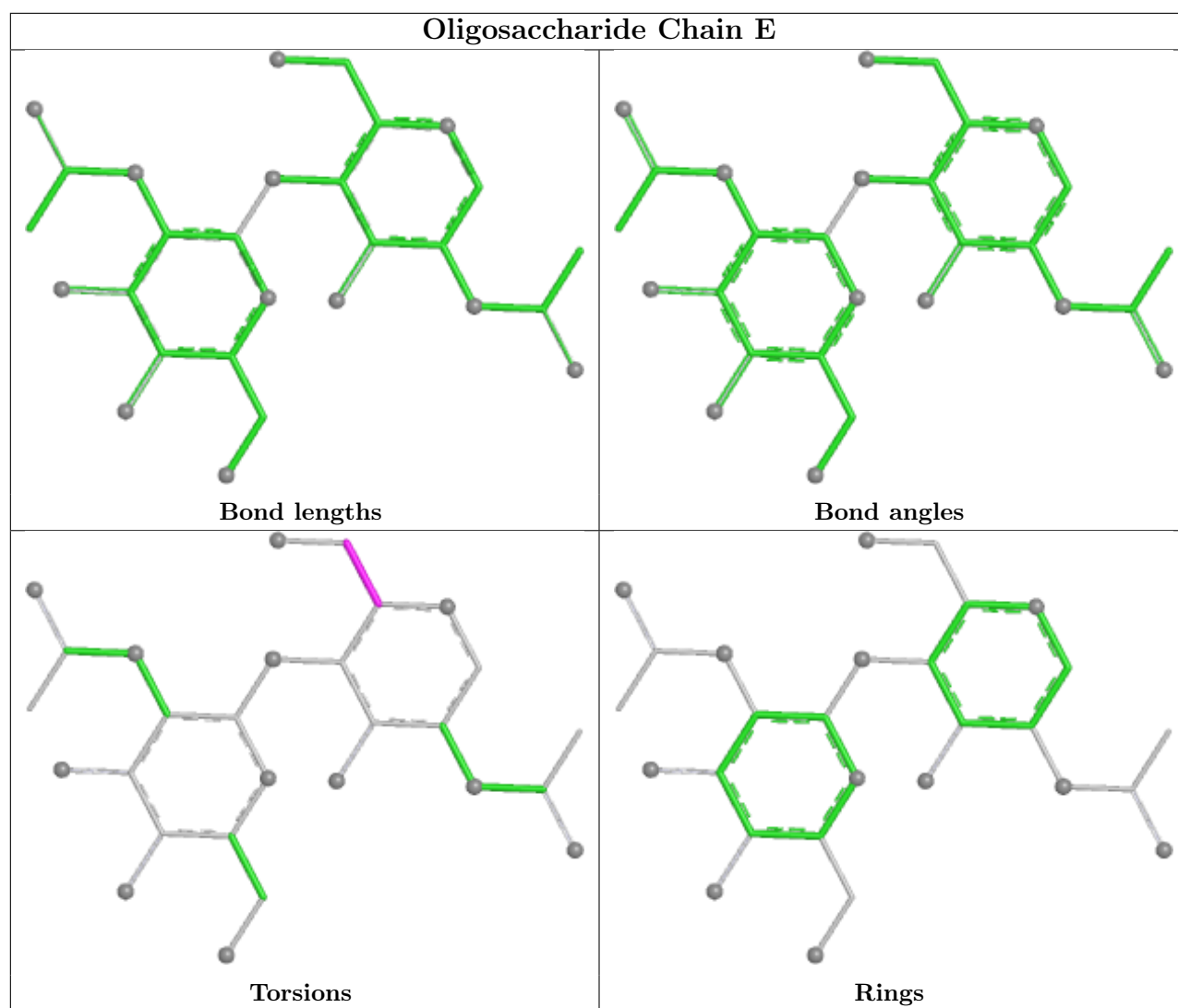
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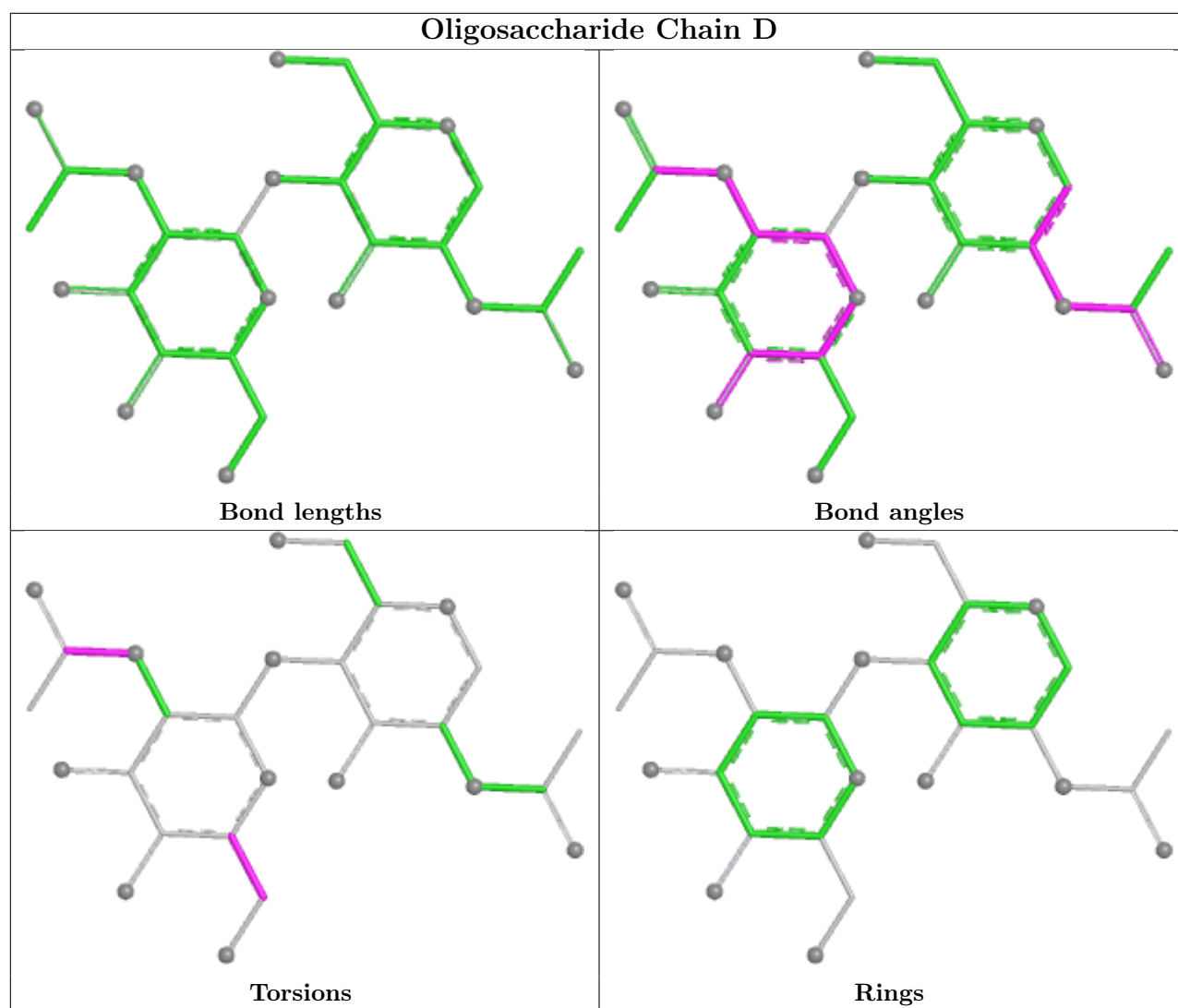
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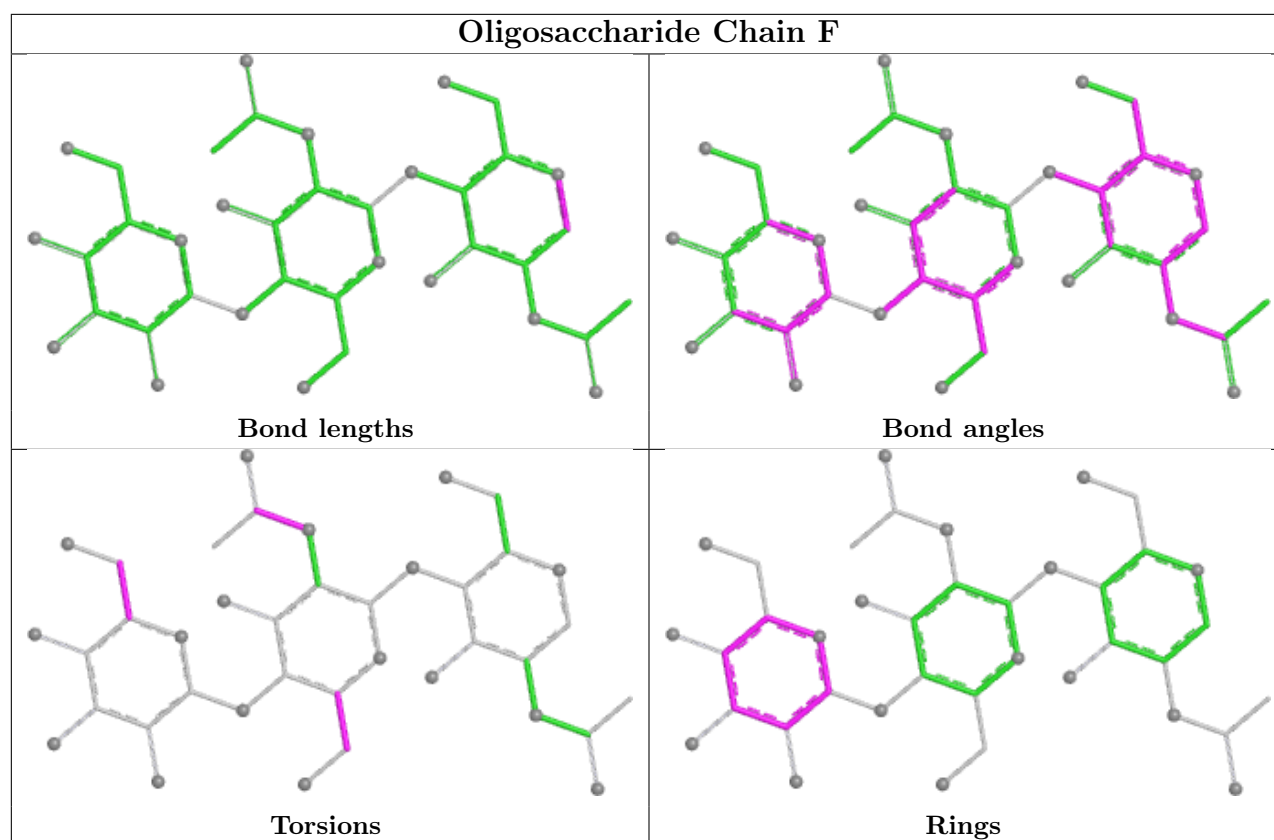
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	2	NDG	1	0
4	F	3	BMA	1	0

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](#)

14 ligands 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	$\# Z > 2$
6	NS4	A	704	-	22,22,22	3.50	5 (22%)	29,31,31	2.31	6 (20%)
5	GOL	A	701	-	5,5,5	0.64	0	5,5,5	0.40	0
5	GOL	B	704	-	5,5,5	0.58	0	5,5,5	0.29	0
5	GOL	B	702	-	5,5,5	0.56	0	5,5,5	0.42	0
7	HEM	A	705	1	50,50,50	1.80	9 (18%)	67,82,82	1.22	3 (4%)
8	NAG	B	707	1	14,14,15	0.76	0	17,19,21	1.97	4 (23%)
10	MAN	B	701	-	11,11,12	0.84	0	15,15,17	1.93	3 (20%)
5	GOL	B	703	-	5,5,5	0.30	0	5,5,5	0.35	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	NAG	A	706	1	14,14,15	1.50	2 (14%)	17,19,21	1.25	2 (11%)
9	BOG	A	707	-	20,20,20	1.19	2 (10%)	25,25,25	1.02	1 (4%)
5	GOL	A	702	-	5,5,5	0.44	0	5,5,5	0.44	0
5	GOL	A	703	-	5,5,5	0.33	0	5,5,5	0.50	0
6	NS4	B	705	-	22,22,22	3.46	5 (22%)	29,31,31	1.91	4 (13%)
7	HEM	B	706	1	50,50,50	1.87	8 (16%)	67,82,82	1.14	2 (2%)

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	NS4	A	704	-	-	5/11/21/21	0/2/2/2
5	GOL	A	701	-	-	2/4/4/4	-
5	GOL	B	704	-	-	4/4/4/4	-
5	GOL	B	702	-	-	2/4/4/4	-
7	HEM	A	705	1	-	8/14/54/54	-
8	NAG	B	707	1	-	1/6/23/26	0/1/1/1
10	MAN	B	701	-	-	2/2/19/22	0/1/1/1
5	GOL	B	703	-	-	0/4/4/4	-
8	NAG	A	706	1	-	2/6/23/26	0/1/1/1
9	BOG	A	707	-	-	6/11/31/31	0/1/1/1
5	GOL	A	702	-	-	4/4/4/4	-
5	GOL	A	703	-	-	4/4/4/4	-
6	NS4	B	705	-	-	5/11/21/21	0/2/2/2
7	HEM	B	706	1	-	5/14/54/54	-

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	704	NS4	O3-N2	12.59	1.44	1.22
6	B	705	NS4	O3-N2	11.96	1.43	1.22
7	B	706	HEM	C3D-C2D	7.61	1.53	1.36
7	A	705	HEM	C3D-C2D	7.46	1.52	1.36
6	B	705	NS4	C1-N1	-6.43	1.32	1.42
6	A	704	NS4	C1-N1	-5.85	1.33	1.42
7	B	706	HEM	FE-NC	5.49	2.13	1.95
6	A	704	NS4	C4-N2	-5.42	1.32	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	705	NS4	C4-N2	-5.39	1.32	1.45
6	B	705	NS4	O2-S1	4.67	1.51	1.43
6	B	705	NS4	O1-S1	4.55	1.51	1.43
6	A	704	NS4	O2-S1	4.47	1.50	1.43
8	A	706	NAG	O5-C1	-4.43	1.36	1.43
7	A	705	HEM	FE-NB	4.32	2.08	1.94
6	A	704	NS4	O1-S1	3.92	1.49	1.43
9	A	707	BOG	O1-C1	3.90	1.46	1.40
7	B	706	HEM	FE-NB	3.68	2.06	1.94
7	B	706	HEM	FE-ND	3.65	2.06	1.94
7	A	705	HEM	FE-NA	3.44	2.06	1.95
7	A	705	HEM	CAC-C3C	3.09	1.55	1.47
8	A	706	NAG	C1-C2	2.95	1.56	1.52
7	B	706	HEM	CAC-C3C	2.89	1.55	1.47
7	A	705	HEM	FE-ND	2.72	2.03	1.94
7	A	705	HEM	CAB-C3B	2.68	1.54	1.47
7	A	705	HEM	FE-NC	2.56	2.03	1.95
7	B	706	HEM	CAB-C3B	2.41	1.53	1.47
9	A	707	BOG	O5-C1	2.37	1.47	1.41
7	B	706	HEM	CMB-C2B	2.25	1.55	1.50
7	A	705	HEM	CMB-C2B	2.21	1.55	1.50
7	A	705	HEM	CMA-C3A	2.09	1.55	1.50
7	B	706	HEM	CMA-C3A	2.01	1.54	1.50

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	704	NS4	O2-S1-O1	-8.64	107.28	118.87
6	B	705	NS4	O2-S1-O1	-7.07	109.39	118.87
8	B	707	NAG	C4-C3-C2	5.48	119.05	111.02
10	B	701	MAN	C1-C2-C3	5.32	117.39	109.64
7	A	705	HEM	C4D-ND-C1D	4.84	110.94	105.21
6	A	704	NS4	O5-C2-C1	4.80	119.94	114.51
7	B	706	HEM	C4D-ND-C1D	4.71	110.78	105.21
6	B	705	NS4	O5-C2-C1	4.61	119.73	114.51
6	A	704	NS4	O5-C8-C9	-4.37	98.11	108.31
8	A	706	NAG	O5-C1-C2	-3.76	105.47	111.29
6	B	705	NS4	O5-C8-C13	3.60	116.71	108.31
8	B	707	NAG	C2-N2-C7	3.42	127.49	122.90
10	B	701	MAN	C2-C3-C4	3.27	116.61	110.86
6	A	704	NS4	C7-S1-N1	3.18	110.15	106.56
8	B	707	NAG	C1-O5-C5	2.66	115.75	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	705	HEM	CAD-C3D-C4D	2.60	129.22	124.70
8	B	707	NAG	C3-C4-C5	2.46	114.69	110.23
8	A	706	NAG	O5-C5-C6	2.29	112.12	107.66
9	A	707	BOG	C1'-O1-C1	-2.27	109.81	113.68
6	B	705	NS4	O3-N2-C4	2.16	121.79	118.82
10	B	701	MAN	O5-C5-C6	2.16	111.86	107.66
7	A	705	HEM	CAA-C2A-C1A	2.14	129.12	124.94
6	A	704	NS4	O5-C8-C13	2.11	113.24	108.31
6	A	704	NS4	C13-C8-C9	2.09	115.94	111.67
7	B	706	HEM	CBD-CAD-C3D	-2.00	107.00	112.53

There are no chirality outliers.

All (50) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	701	GOL	O1-C1-C2-C3
5	A	702	GOL	C1-C2-C3-O3
5	A	702	GOL	O2-C2-C3-O3
5	A	703	GOL	O1-C1-C2-C3
5	A	703	GOL	C1-C2-C3-O3
5	B	702	GOL	O1-C1-C2-C3
5	B	704	GOL	O1-C1-C2-C3
6	A	704	NS4	C1-N1-S1-C7
6	A	704	NS4	C1-C2-O5-C8
6	B	705	NS4	C1-C2-O5-C8
6	B	705	NS4	C13-C8-O5-C2
7	A	705	HEM	C1A-C2A-CAA-CBA
7	A	705	HEM	C2B-C3B-CAB-CBB
7	A	705	HEM	C4B-C3B-CAB-CBB
7	B	706	HEM	C2B-C3B-CAB-CBB
7	B	706	HEM	C4B-C3B-CAB-CBB
10	B	701	MAN	O5-C5-C6-O6
10	B	701	MAN	C4-C5-C6-O6
7	A	705	HEM	C3A-C2A-CAA-CBA
7	A	705	HEM	C4D-C3D-CAD-CBD
8	A	706	NAG	C4-C5-C6-O6
9	A	707	BOG	C2-C1-O1-C1'
7	A	705	HEM	C2D-C3D-CAD-CBD
7	A	705	HEM	C2A-CAA-CBA-CGA
5	A	702	GOL	O1-C1-C2-C3
9	A	707	BOG	O5-C1-O1-C1'
5	A	701	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
5	A	703	GOL	O1-C1-C2-O2
5	B	702	GOL	O1-C1-C2-O2
6	B	705	NS4	C1-N1-S1-C7
6	A	704	NS4	C3-C2-O5-C8
6	A	704	NS4	C13-C8-O5-C2
7	A	705	HEM	C3D-CAD-CBD-CGD
5	A	703	GOL	O2-C2-C3-O3
5	B	704	GOL	O1-C1-C2-O2
8	B	707	NAG	O5-C5-C6-O6
8	A	706	NAG	O5-C5-C6-O6
5	B	704	GOL	O2-C2-C3-O3
9	A	707	BOG	C5'-C6'-C7'-C8'
9	A	707	BOG	C4-C5-C6-O6
5	A	702	GOL	O1-C1-C2-O2
6	A	704	NS4	C1-N1-S1-O1
5	B	704	GOL	C1-C2-C3-O3
9	A	707	BOG	C1'-C2'-C3'-C4'
6	B	705	NS4	C9-C8-O5-C2
9	A	707	BOG	O5-C5-C6-O6
6	B	705	NS4	C3-C2-O5-C8
7	B	706	HEM	CAA-CBA-CGA-O2A
7	B	706	HEM	CAA-CBA-CGA-O1A
7	B	706	HEM	C4C-C3C-CAC-CBC

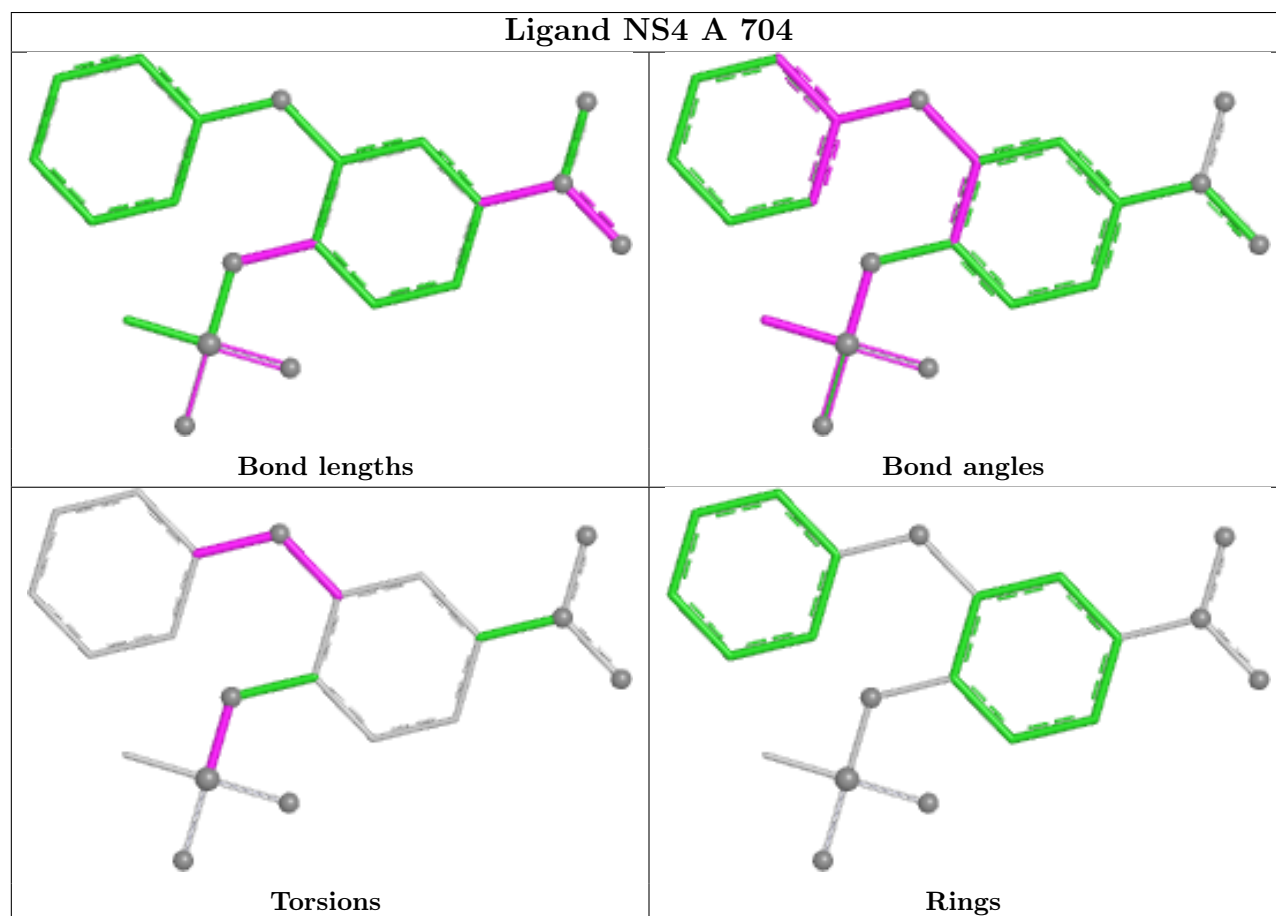
There are no ring outliers.

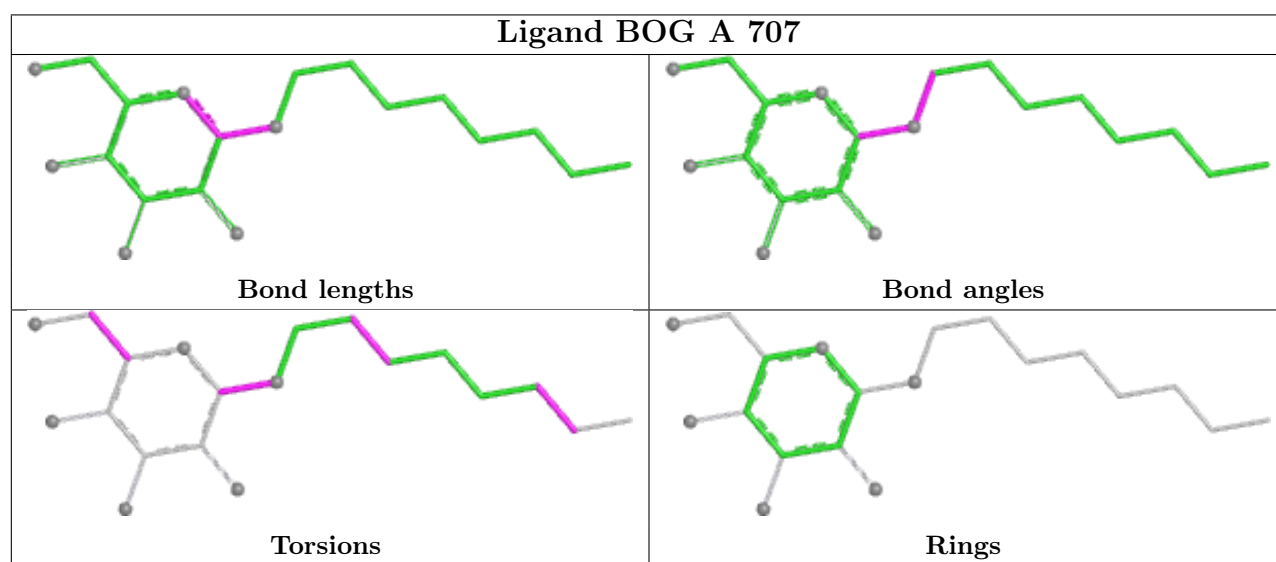
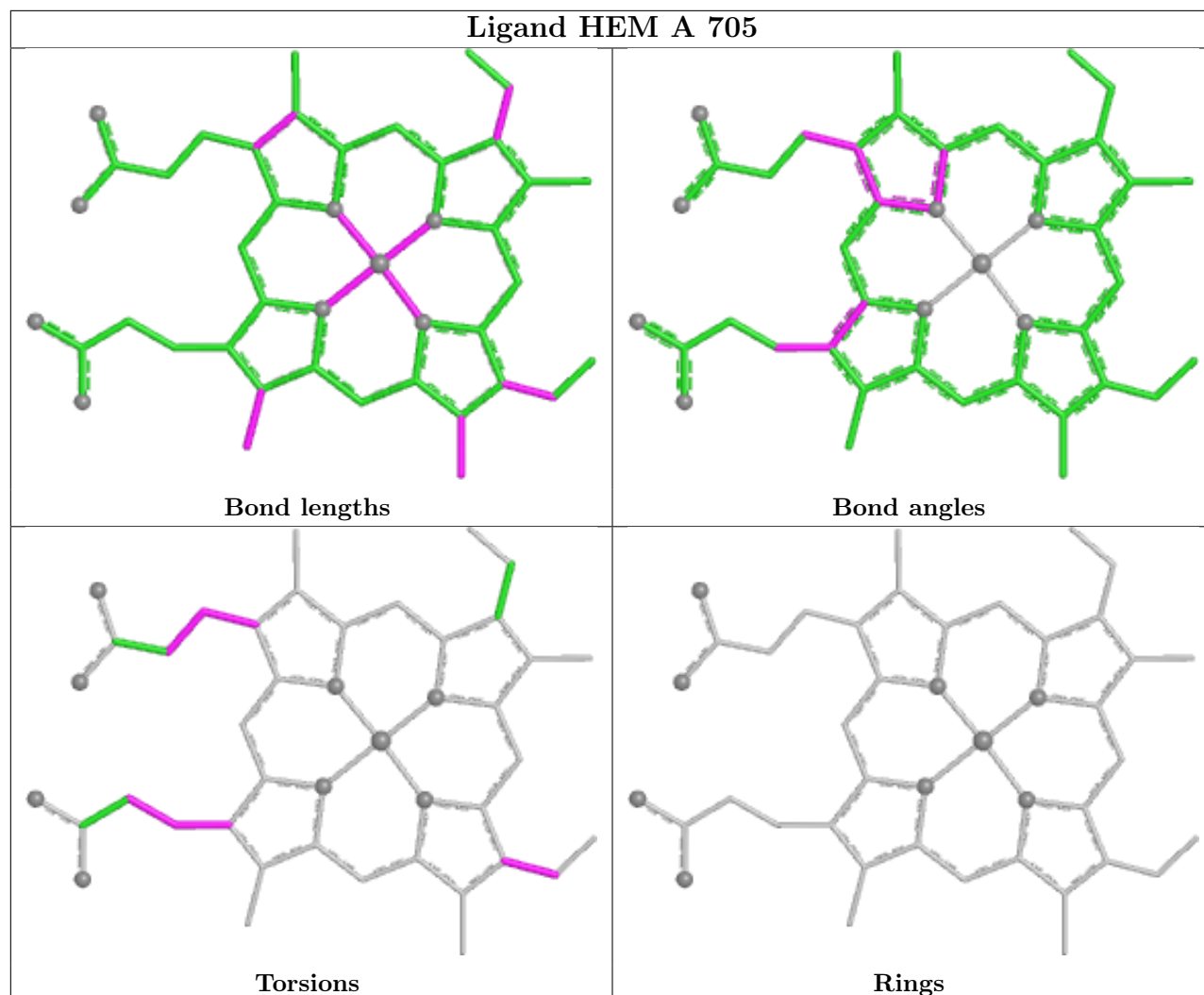
10 monomers are involved in 21 short contacts:

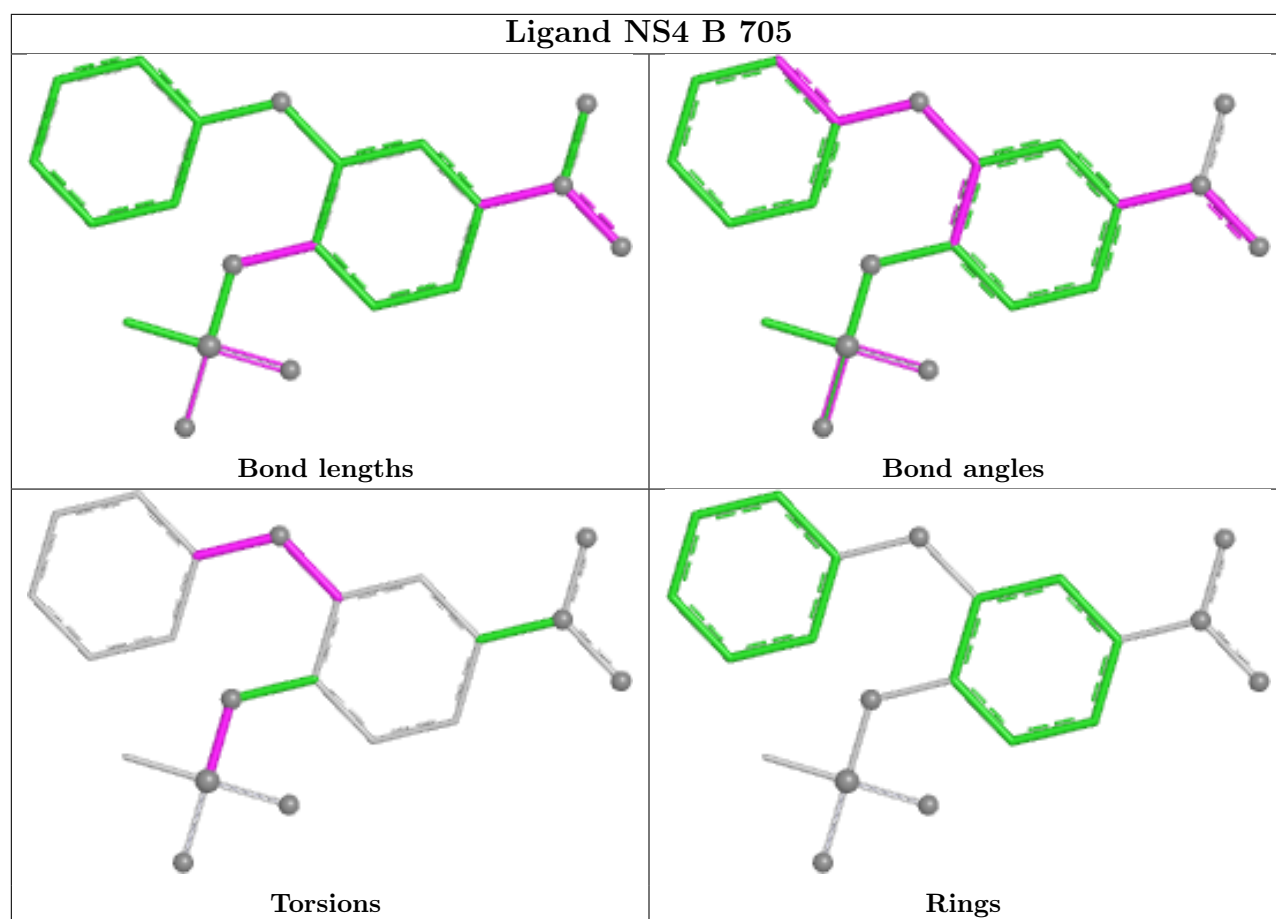
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	704	NS4	5	0
5	A	701	GOL	1	0
5	B	704	GOL	1	0
7	A	705	HEM	2	0
5	B	703	GOL	2	0
9	A	707	BOG	1	0
5	A	702	GOL	1	0
5	A	703	GOL	1	0
6	B	705	NS4	6	0
7	B	706	HEM	3	0

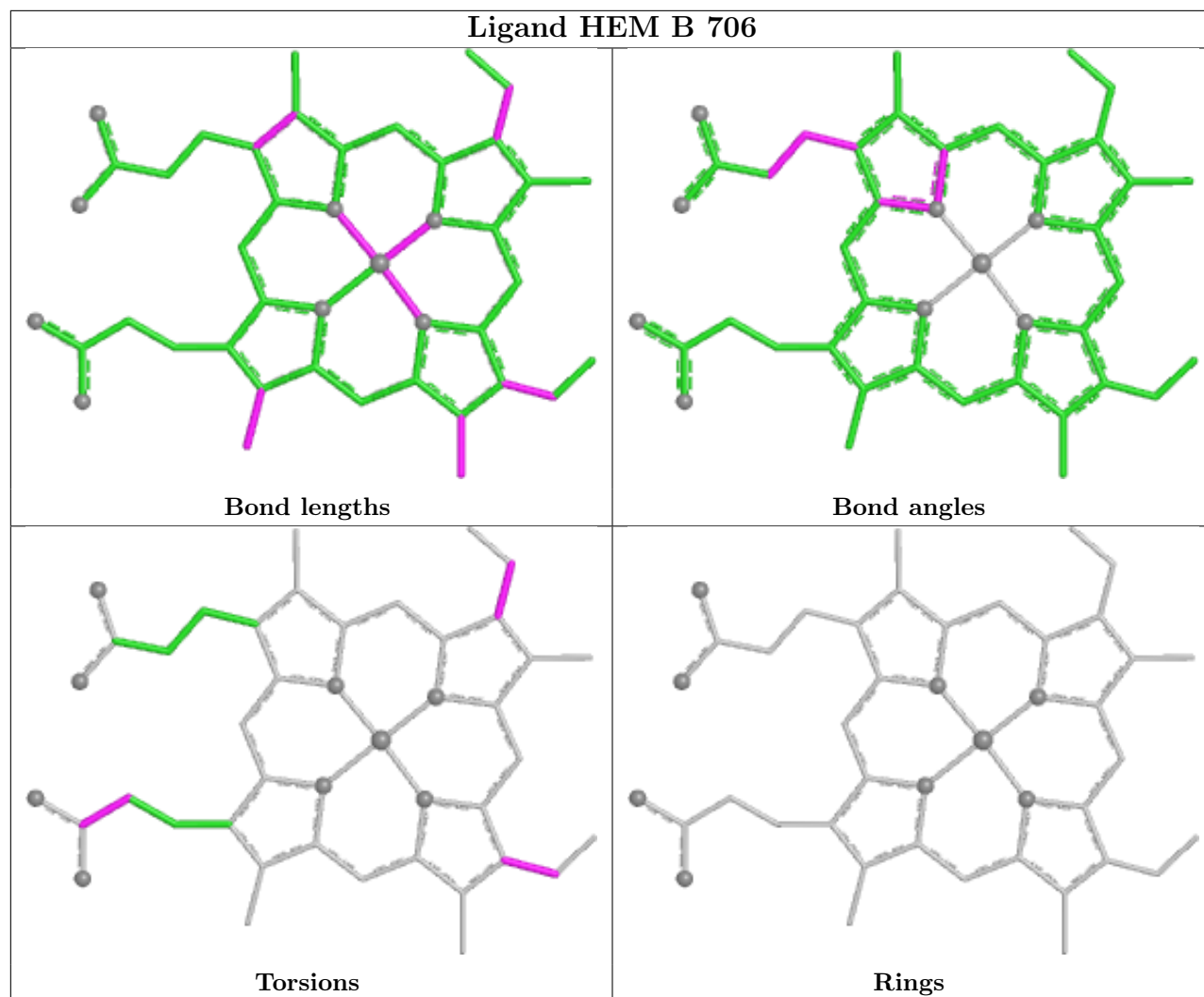
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

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 [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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	551/610 (90%)	-0.63	1 (0%) 91 83	10, 33, 51, 60	3 (0%)
1	B	551/610 (90%)	-0.65	0 100 100	17, 33, 51, 60	0
All	All	1102/1220 (90%)	-0.64	1 (0%) 92 86	10, 33, 51, 60	3 (0%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	577	PHE	4.0

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

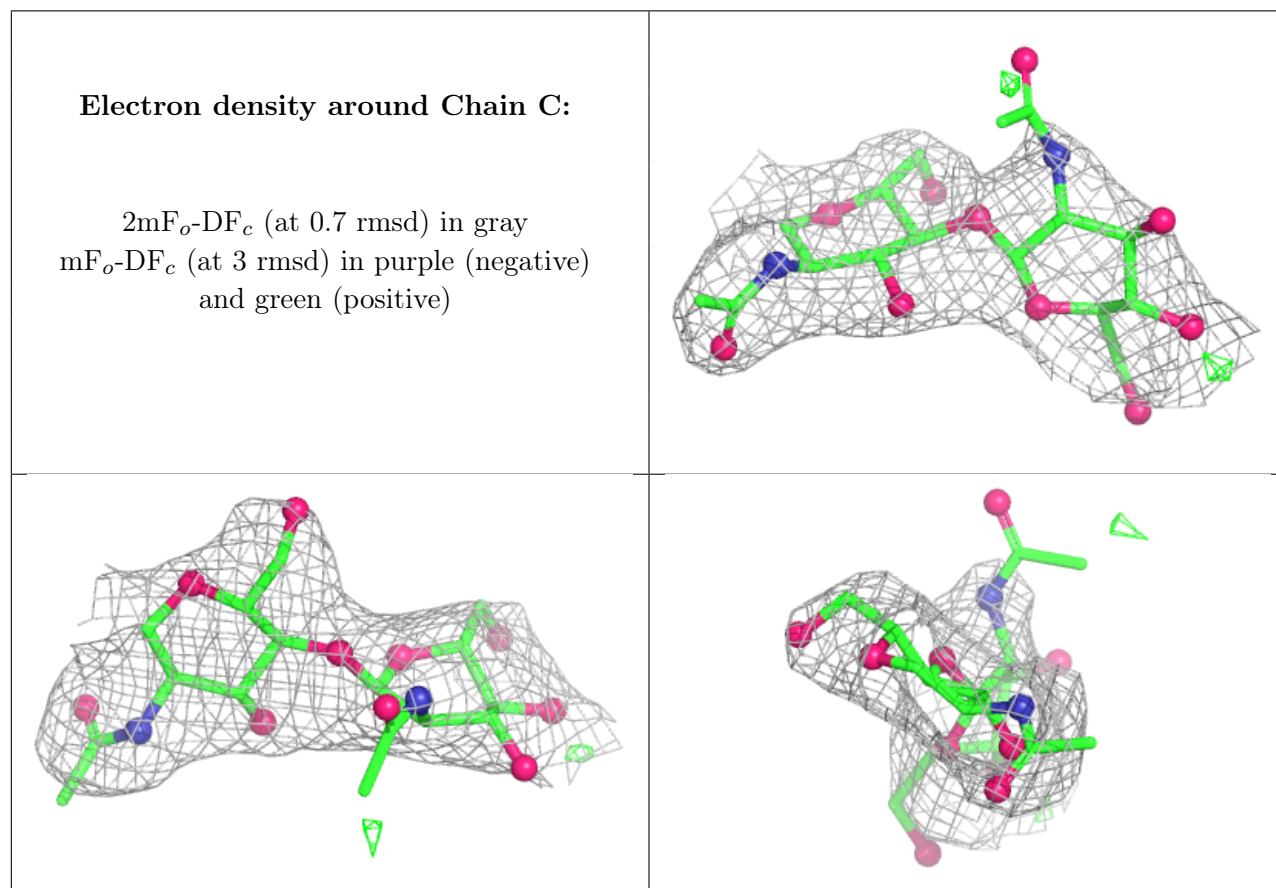
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

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

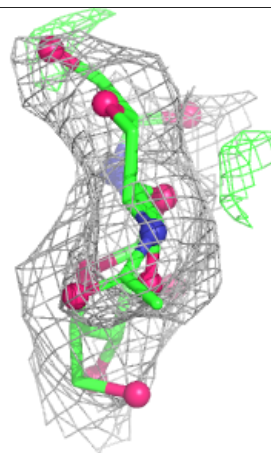
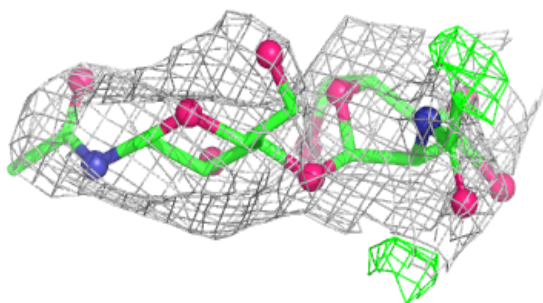
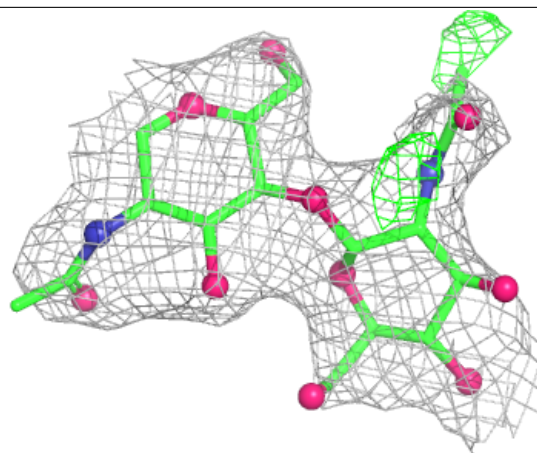
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NDG	C	2	14/15	0.63	0.15	74,77,78,79	0
2	NDG	E	2	14/15	0.73	0.12	72,73,74,74	0
4	BMA	F	3	11/12	0.78	0.13	55,57,58,58	0
3	NAG	D	2	14/15	0.86	0.09	47,50,54,58	0
4	NAG	F	2	14/15	0.91	0.09	43,45,50,52	0
4	NAG	F	1	14/15	0.92	0.09	31,34,39,39	0
2	NAG	C	1	14/15	0.92	0.09	59,61,65,69	0
2	NAG	E	1	14/15	0.92	0.08	60,63,65,69	0
3	NAG	D	1	14/15	0.93	0.09	32,36,38,43	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.



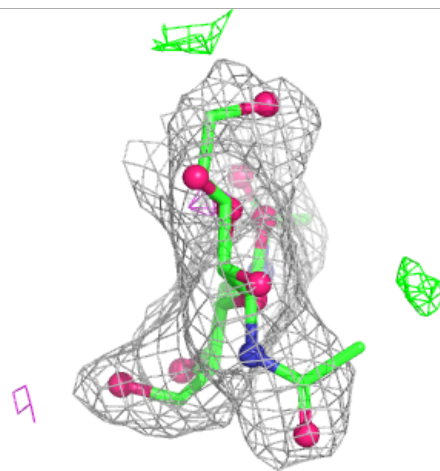
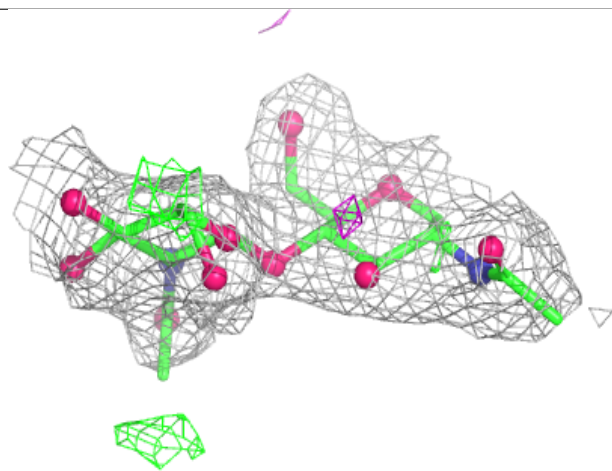
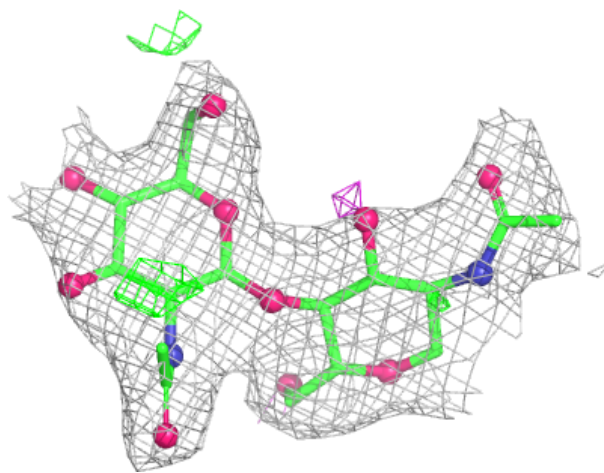
Electron density around Chain E:

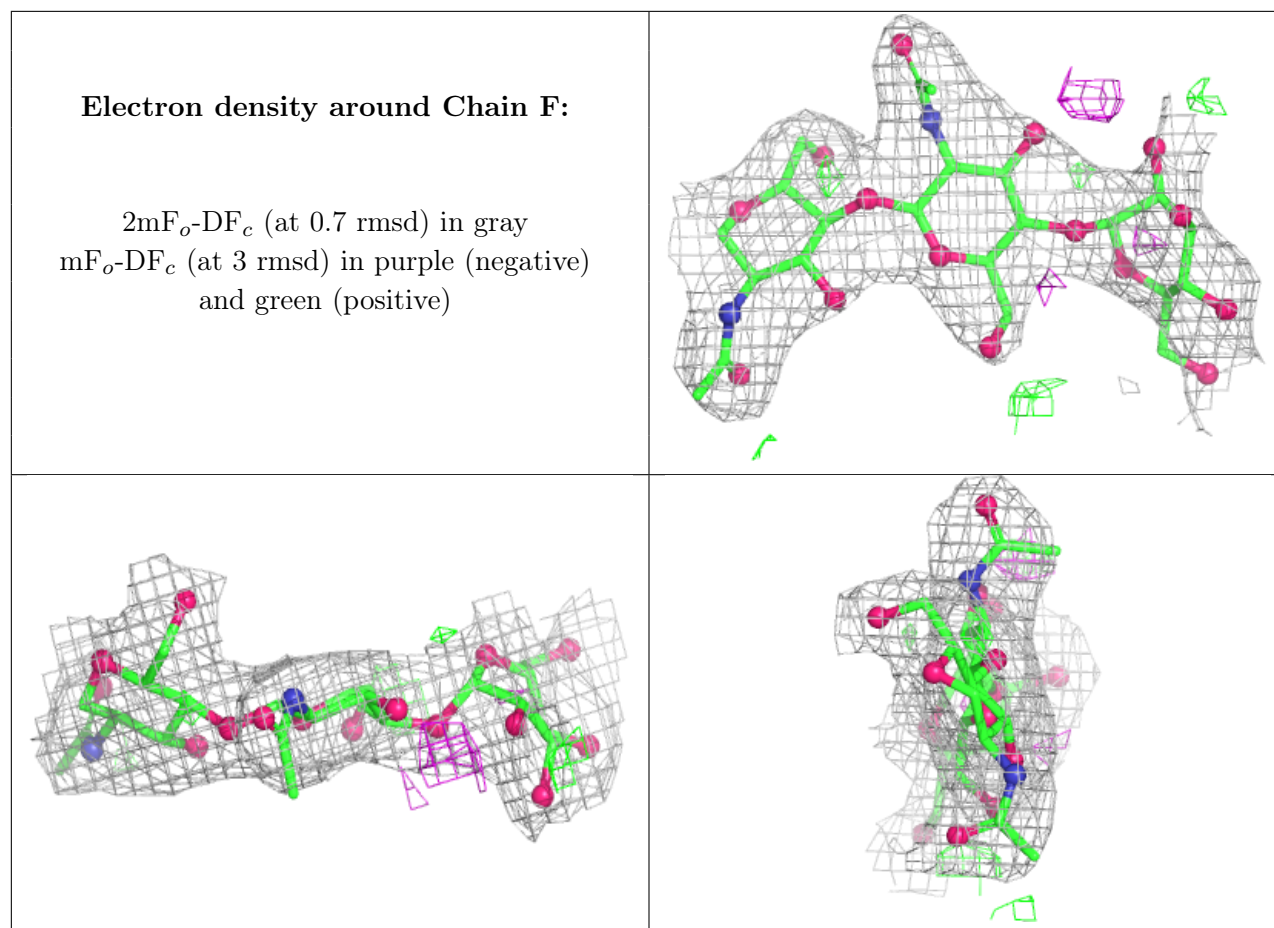
$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain D:

$2mF_o-DF_c$ (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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

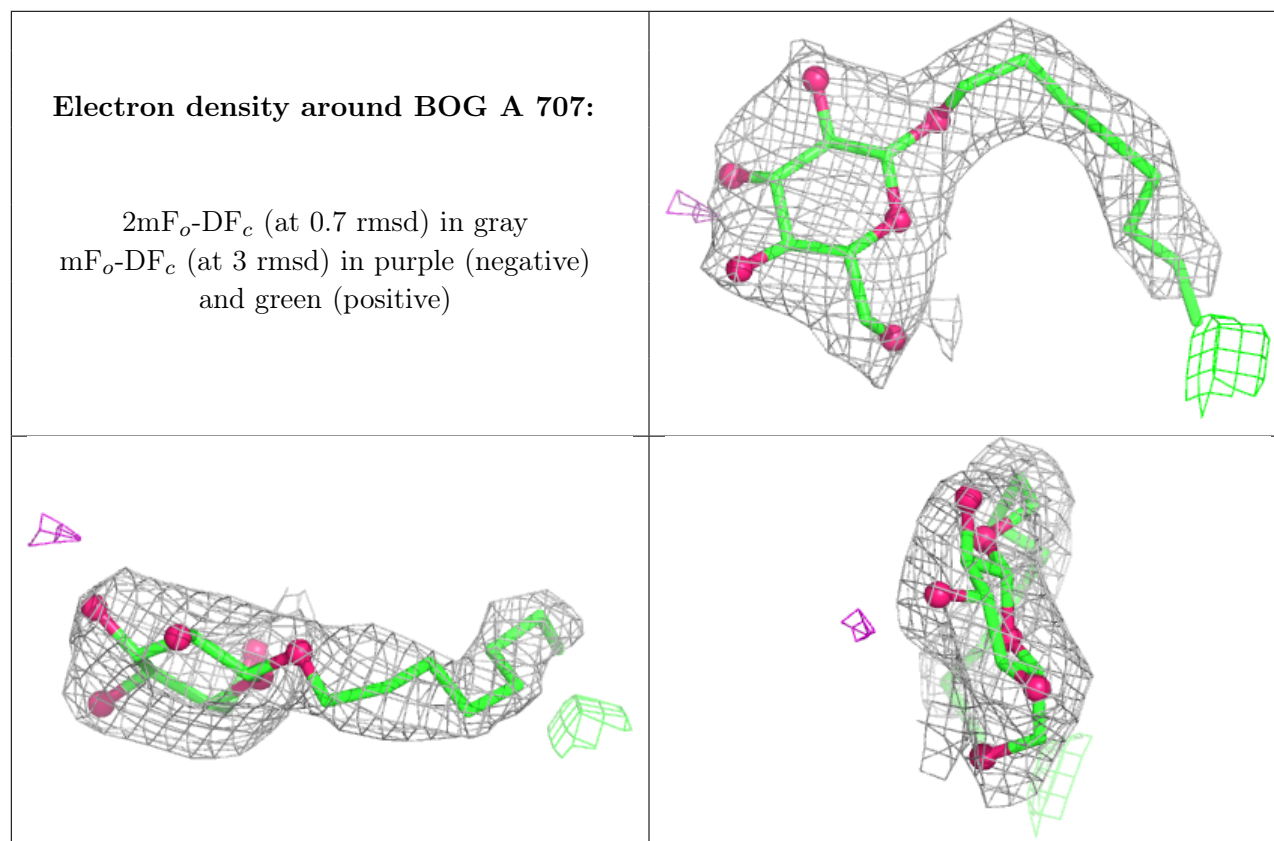
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	NAG	B	707	14/15	0.76	0.13	61,66,67,68	0
10	MAN	B	701	11/12	0.81	0.11	61,63,65,65	0
5	GOL	B	704	6/6	0.88	0.13	57,58,59,59	0
5	GOL	B	702	6/6	0.89	0.12	53,58,58,59	0
8	NAG	A	706	14/15	0.89	0.12	58,59,61,61	0
5	GOL	A	703	6/6	0.92	0.09	49,49,50,50	0
5	GOL	A	702	6/6	0.92	0.12	49,51,51,52	0
9	BOG	A	707	20/20	0.93	0.09	49,51,53,54	0
5	GOL	B	703	6/6	0.93	0.11	54,55,56,56	0
5	GOL	A	701	6/6	0.95	0.09	51,52,52,53	0
6	NS4	B	705	21/21	0.95	0.08	39,42,43,46	0
6	NS4	A	704	21/21	0.96	0.08	29,35,38,41	0

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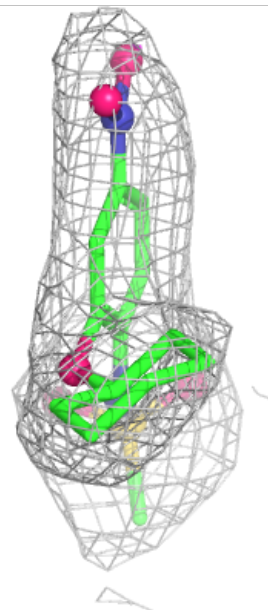
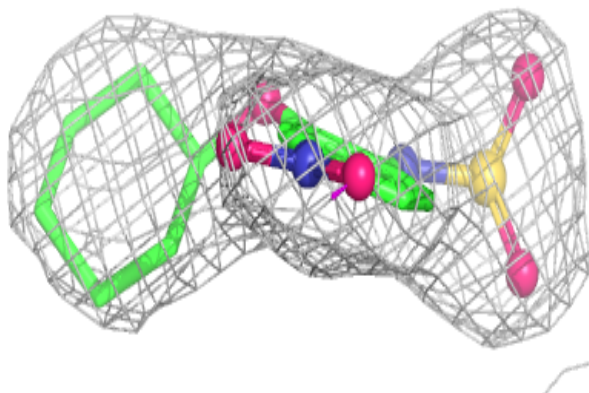
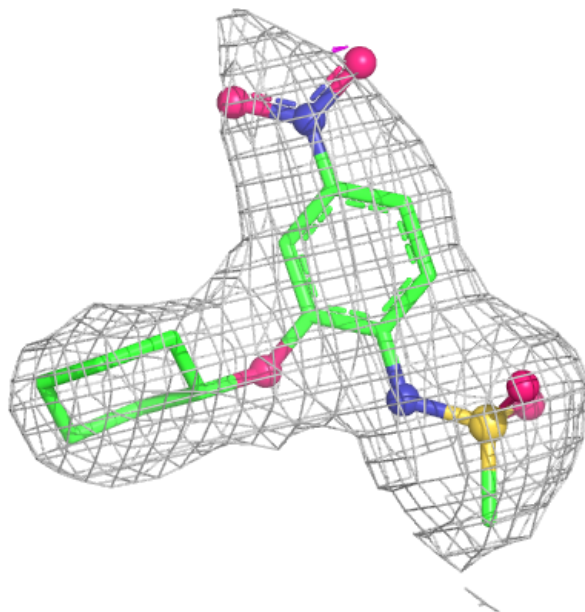
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	HEM	A	705	43/43	0.98	0.06	20,26,41,47	0
7	HEM	B	706	43/43	0.98	0.07	24,29,43,49	0

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.



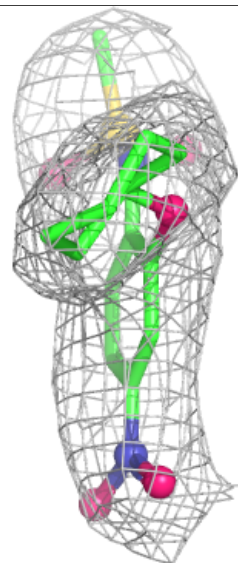
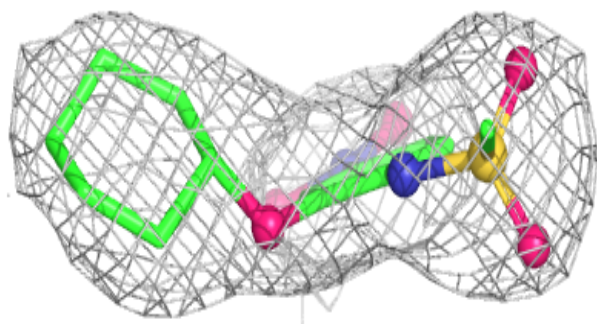
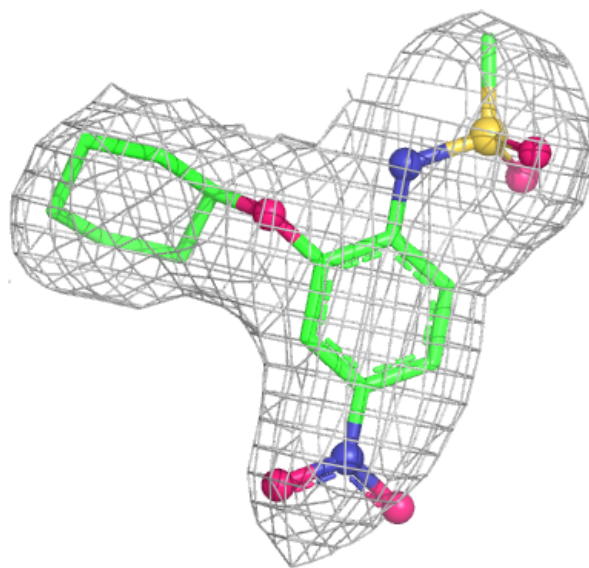
Electron density around NS4 B 705:

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



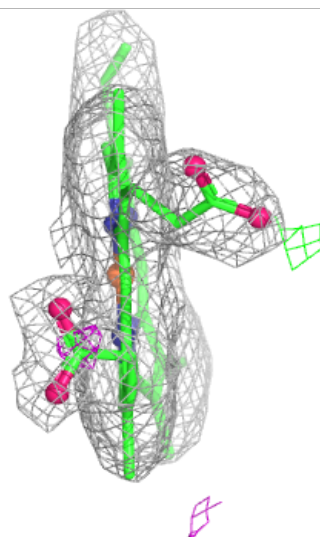
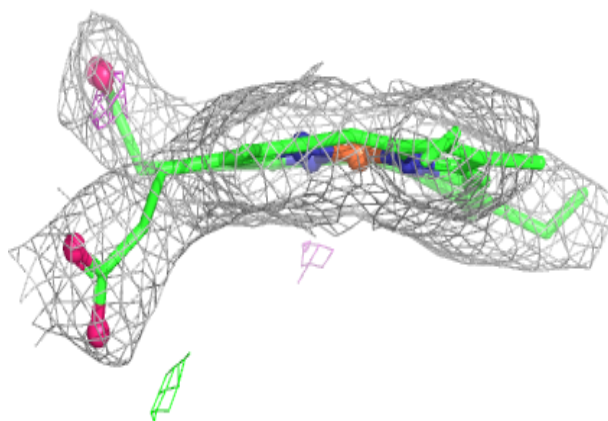
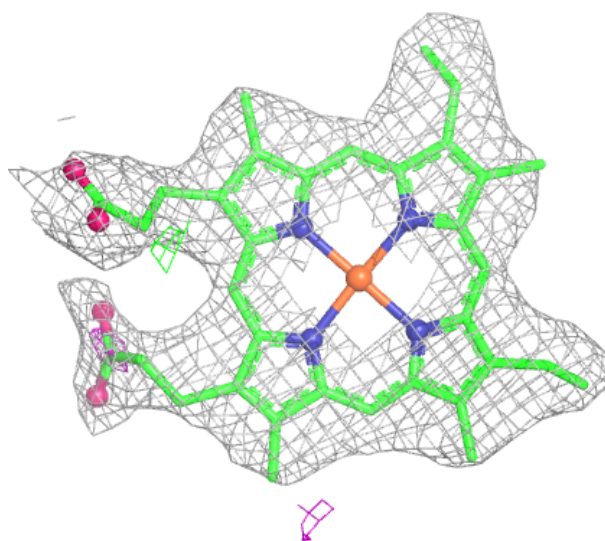
Electron density around NS4 A 704:

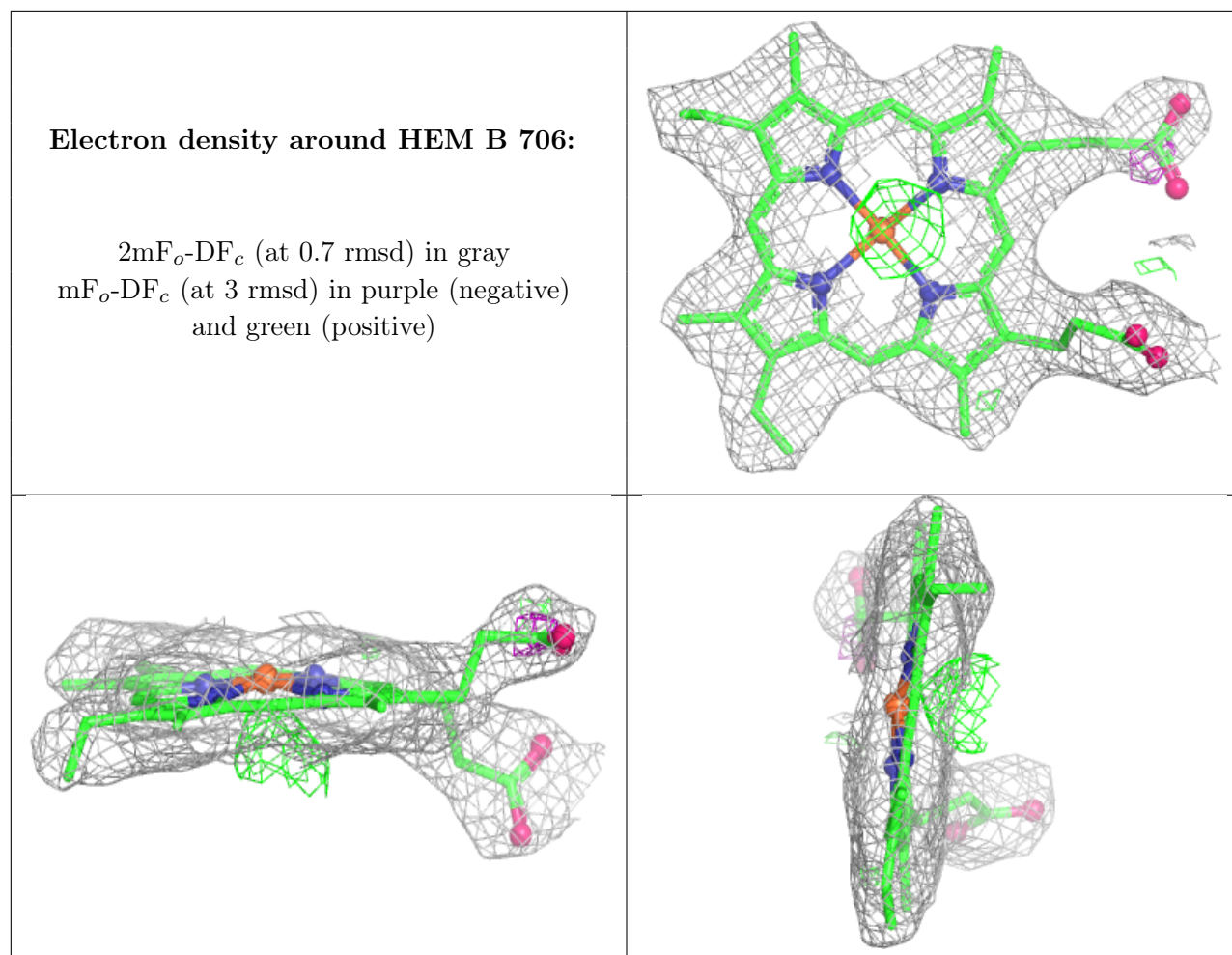
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM A 705:

$2mF_o-DF_c$ (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.