



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 07:01 AM UTC

PDB ID : 6V3R / pdb_00006v3r
Title : Crystal structure of murine cyclooxygenase in complex with a harmaline analog,
4,9-dihydro-3H-pyrido[3,4-b]indole
Authors : Xu, S.; Uddin, M.J.; Banerjee, S.; Marnett, L.J.
Deposited on : 2019-11-26
Resolution : 2.66 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

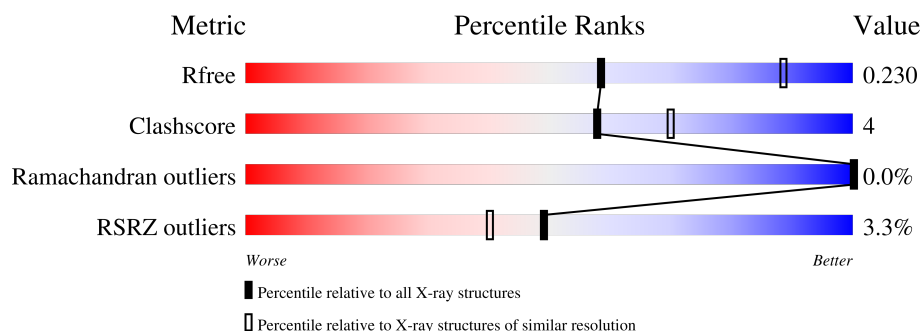
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.66 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	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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	587	2% 82% 12% 6%
1	B	587	4% 84% 10% 6%
1	C	587	3% 84% 10% 6%
1	D	587	3% 83% 11% 6%
2	E	2	100%
2	F	2	100%

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Mol	Chain	Length	Quality of chain
2	G	2	 100%
2	H	2	 50% 50%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 18772 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	552	Total	C	N	O	S	0	3	0
			4499	2899	756	819	25			
1	B	552	Total	C	N	O	S	0	3	0
			4499	2899	756	819	25			
1	C	552	Total	C	N	O	S	0	3	0
			4499	2899	756	819	25			
1	D	552	Total	C	N	O	S	0	3	0
			4499	2899	756	819	25			

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



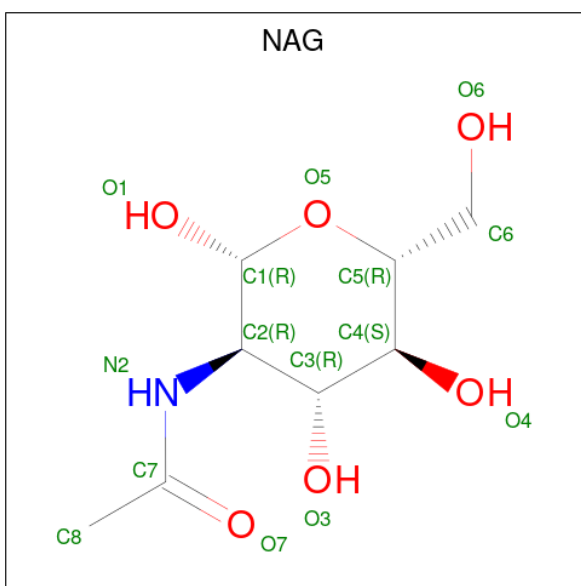
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	H	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: C₃₄H₃₂FeN₄O₄).



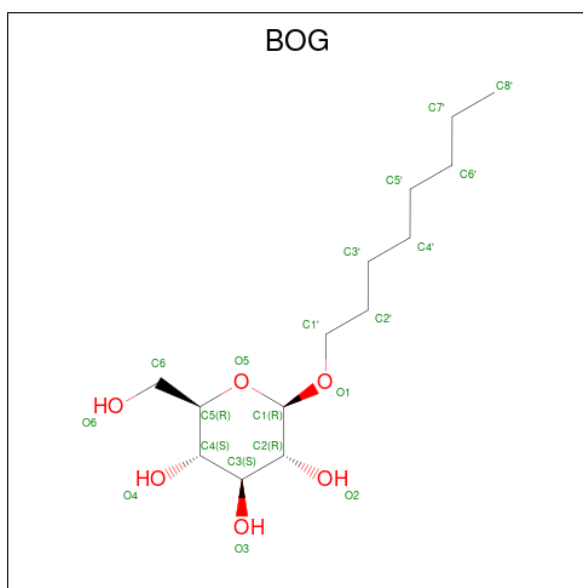
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



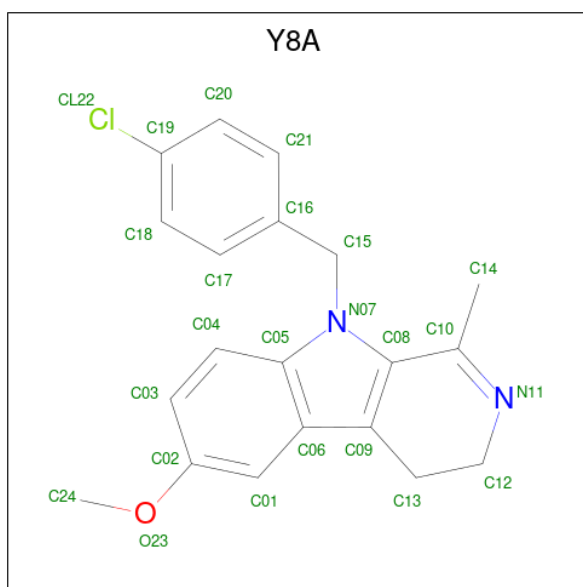
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		
4	C	1	Total	C	N	O	0	0
			14	8	1	5		
4	C	1	Total	C	N	O	0	0
			14	8	1	5		
4	D	1	Total	C	N	O	0	0
			14	8	1	5		
4	D	1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

- Molecule 5 is octyl beta-D-glucopyranoside (CCD ID: BOG) (formula: $C_{14}H_{28}O_6$).



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

- Molecule 6 is 9-[(4-chlorophenyl)methyl]-6-methoxy-1-methyl-4,9-dihydro-3H-beta-carbolin e (CCD ID: Y8A) (formula: $C_{20}H_{19}ClN_2O$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	Cl	N	O	0	0
			24	20	1	2	1		
6	B	1	Total	C	Cl	N	O	0	0
			24	20	1	2	1		
6	C	1	Total	C	Cl	N	O	0	0
			24	20	1	2	1		
6	D	1	Total	C	Cl	N	O	0	0
			24	20	1	2	1		

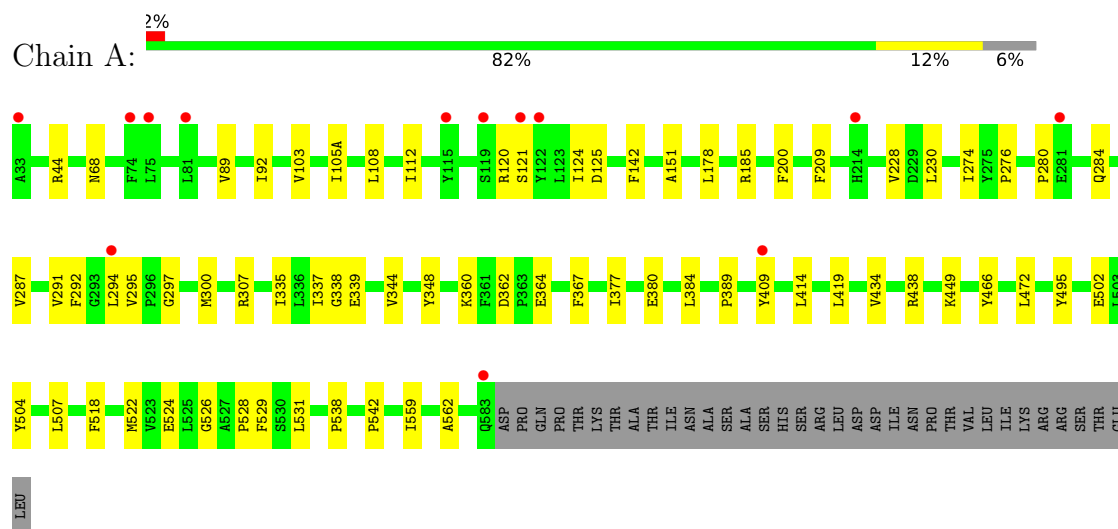
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	56	Total	O	0	0
			56	56		
7	B	55	Total	O	0	0
			55	55		
7	C	57	Total	O	0	0
			57	57		
7	D	76	Total	O	0	0
			76	76		

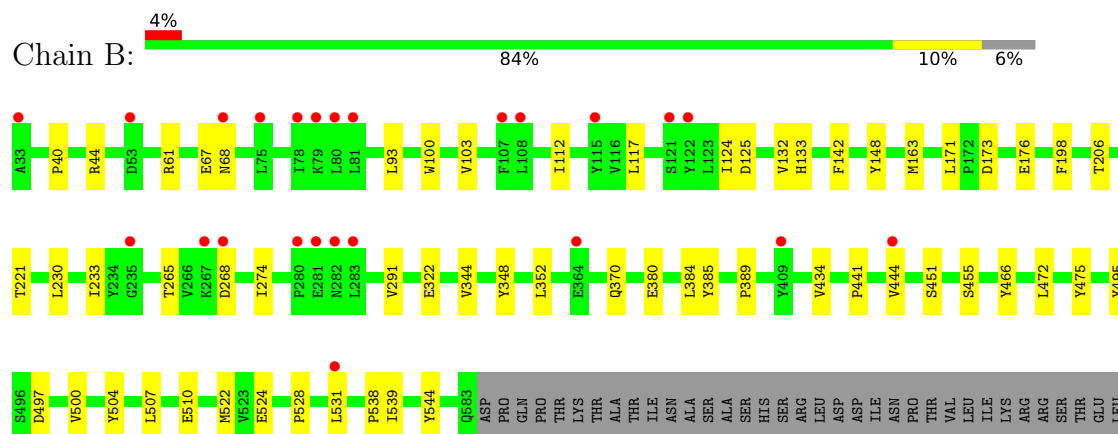
3 Residue-property plots [i](#)

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.

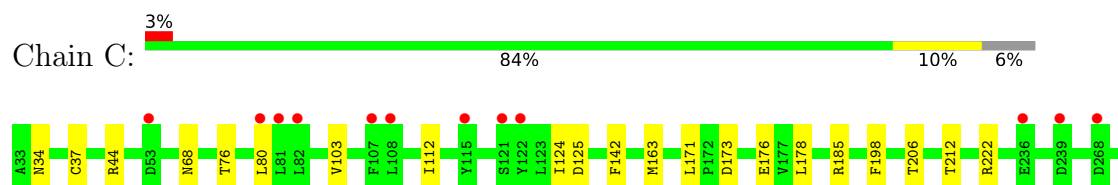
• Molecule 1: Prostaglandin G/H synthase 2



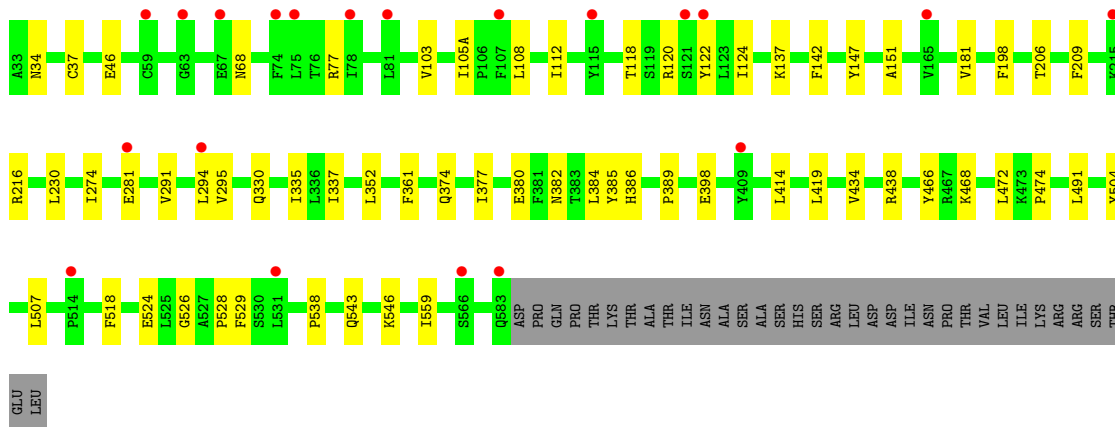
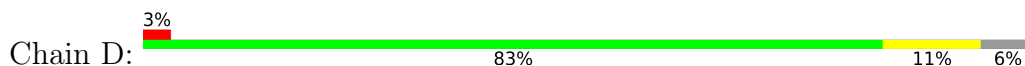
• Molecule 1: Prostaglandin G/H synthase 2



• Molecule 1: Prostaglandin G/H synthase 2



- Molecule 1: Prostaglandin G/H synthase 2



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



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



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	217.15Å 124.38Å 136.54Å 90.00° 123.84° 90.00°	Depositor
Resolution (Å)	102.39 – 2.66 102.39 – 2.66	Depositor EDS
% Data completeness (in resolution range)	98.0 (102.39-2.66) 98.0 (102.39-2.66)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.10 (at 2.65Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.219 , 0.264 (Not available) , 0.230	Depositor DCC
R_{free} test set	2537 reflections (2.94%)	wwPDB-VP
Wilson B-factor (Å ²)	42.6	Xtriage
Anisotropy	0.110	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 38.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for 1/2*h-3/2*k,-1/2*h-1/2*k,-1/2*h +1/2*k-l 0.000 for 1/2*h+3/2*k,1/2*h-1/2*k,-1/2*h- 1/2*k-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	18772	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.24% 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: Y8A, BOG, HEM, NAG

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.11	0/4629	0.31	1/6276 (0.0%)
1	B	0.11	0/4629	0.30	0/6276
1	C	0.10	0/4629	0.29	0/6276
1	D	0.12	0/4629	0.30	0/6276
All	All	0.11	0/18516	0.30	1/25104 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	228	VAL	N-CA-C	5.39	111.97	106.21

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	4499	0	4396	45	0
1	B	4499	0	4396	38	0
1	C	4499	0	4396	36	0
1	D	4499	0	4396	44	0
2	E	28	0	25	0	0
2	F	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	G	28	0	25	0	0
2	H	28	0	25	1	0
3	A	43	0	30	2	0
3	B	43	0	30	2	0
3	C	43	0	30	2	0
3	D	43	0	30	2	0
4	A	28	0	26	0	0
4	B	28	0	26	1	0
4	C	28	0	26	0	0
4	D	28	0	26	0	0
5	A	20	0	28	1	0
5	D	20	0	28	0	0
6	A	24	0	0	2	0
6	B	24	0	0	1	0
6	C	24	0	0	1	0
6	D	24	0	0	2	0
7	A	56	0	0	4	0
7	B	55	0	0	1	0
7	C	57	0	0	1	0
7	D	76	0	0	7	0
All	All	18772	0	17964	164	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 4.

All (164) 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)
3:D:702:HEM:HBC2	3:D:702:HEM:HHD	1.70	0.73
3:B:702:HEM:HHC	3:B:702:HEM:HBB2	1.71	0.72
3:A:701:HEM:HHD	3:A:701:HEM:HBC2	1.72	0.71
1:C:173:ASP:HB3	1:C:176:GLU:HB2	1.70	0.71
3:C:702:HEM:HBB2	3:C:702:HEM:HHC	1.73	0.71
1:C:274:ILE:HD12	1:C:291:VAL:HG12	1.73	0.70
1:D:216:ARG:NH1	2:H:2:NAG:O7	2.26	0.68
3:A:701:HEM:HBB2	3:A:701:HEM:HHC	1.75	0.68
3:D:702:HEM:HBB2	3:D:702:HEM:HHC	1.76	0.67
3:C:702:HEM:HBC2	3:C:702:HEM:HHD	1.77	0.66
3:B:702:HEM:HBC2	3:B:702:HEM:HHD	1.76	0.66
1:A:230:LEU:HD13	1:A:337:ILE:HG12	1.77	0.66
1:D:77:ARG:NH1	7:D:806:HOH:O	2.30	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:274:ILE:HD12	1:D:291:VAL:HG12	1.78	0.64
1:D:294:LEU:HG	1:D:295:VAL:HG22	1.79	0.63
1:A:178:LEU:HD22	1:A:449:LYS:HB2	1.81	0.62
1:D:103:VAL:HG11	1:D:112:ILE:HD12	1.82	0.62
1:B:173:ASP:HB3	1:B:176:GLU:HB2	1.82	0.61
1:A:151:ALA:HB2	1:A:529:PHE:HZ	1.66	0.60
1:B:497:ASP:HB3	1:B:500:VAL:HG23	1.84	0.59
1:A:542:PRO:O	1:B:61:ARG:NH2	2.29	0.59
1:D:103:VAL:HG13	1:D:108:LEU:HB3	1.83	0.59
1:B:274:ILE:HD12	1:B:291:VAL:HG12	1.85	0.58
1:B:68:ASN:N	1:B:68:ASN:OD1	2.34	0.58
1:A:274:ILE:HD12	1:A:291:VAL:HG12	1.86	0.58
1:D:124:ILE:HD11	1:D:528:PRO:HB2	1.84	0.58
1:C:538:PRO:O	1:C:541:SER:OG	2.20	0.58
1:C:389:PRO:HB2	1:C:434:VAL:HA	1.86	0.57
1:C:212:THR:O	1:C:222:ARG:NH2	2.37	0.57
1:D:374:GLN:NE2	7:D:807:HOH:O	2.30	0.57
1:C:103:VAL:HG11	1:C:112:ILE:HD12	1.88	0.56
1:D:209:PHE:HB2	1:D:377:ILE:HG13	1.88	0.56
1:A:103:VAL:HG22	1:A:108:LEU:HD23	1.87	0.56
1:A:294:LEU:HG	1:A:295:VAL:HG13	1.86	0.56
1:A:103:VAL:HG11	1:A:112:ILE:HD12	1.87	0.56
1:B:389:PRO:HB2	1:B:434:VAL:HA	1.87	0.55
1:C:384:LEU:HD13	6:C:701:Y8A:CL22	2.44	0.55
1:B:265:THR:OG1	1:B:268:ASP:OD1	2.22	0.55
1:D:151:ALA:HB2	1:D:529:PHE:HZ	1.71	0.55
1:C:497:ASP:HB3	1:C:500:VAL:HG23	1.88	0.54
1:A:367:PHE:O	1:B:370:GLN:NE2	2.37	0.54
1:C:280:PRO:HD2	1:C:283:LEU:HD12	1.89	0.54
1:D:504:TYR:HA	1:D:507:LEU:HD12	1.90	0.54
1:A:105(A):ILE:HG22	1:A:108:LEU:H	1.73	0.53
5:A:706:BOG:H6'1	1:C:185:ARG:HA	1.91	0.53
1:B:163:MET:HE2	1:B:171:LEU:HD11	1.90	0.53
1:A:121:SER:HB2	1:A:531:LEU:HD12	1.90	0.52
1:D:384:LEU:HD13	6:D:701:Y8A:CL22	2.46	0.52
1:A:384:LEU:HD13	6:A:707:Y8A:CL22	2.47	0.52
1:A:495:TYR:HE2	1:A:502:GLU:HG3	1.75	0.52
1:D:330:GLN:NE2	7:D:808:HOH:O	2.35	0.52
1:B:384:LEU:HB3	1:B:507:LEU:HD13	1.93	0.51
1:B:103:VAL:HG11	1:B:112:ILE:HD12	1.92	0.51
1:A:124:ILE:HD11	1:A:528:PRO:HB2	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:198:PHE:HZ	1:C:352:LEU:HD23	1.76	0.50
1:B:507:LEU:HB3	1:B:522:MET:HE2	1.93	0.50
1:A:504:TYR:HA	1:A:507:LEU:HD12	1.94	0.50
1:B:124:ILE:HD11	1:B:528:PRO:HB2	1.94	0.50
1:B:44:ARG:NH2	1:B:125:ASP:OD1	2.36	0.50
1:C:380:GLU:O	1:C:384:LEU:HG	2.12	0.49
1:A:92:ILE:N	7:A:802:HOH:O	2.44	0.49
1:D:230:LEU:HD13	1:D:337:ILE:HG12	1.94	0.49
1:B:198:PHE:HZ	1:B:352:LEU:HD23	1.77	0.49
1:C:68:ASN:N	1:C:68:ASN:OD1	2.45	0.49
1:D:68:ASN:OD1	1:D:68:ASN:N	2.45	0.49
1:A:389:PRO:HB2	1:A:434:VAL:HA	1.94	0.49
1:D:546:LYS:NZ	7:D:813:HOH:O	2.42	0.49
1:A:380:GLU:HG2	1:A:466:TYR:CE1	2.48	0.48
1:D:147:TYR:HB2	7:D:850:HOH:O	2.12	0.48
1:A:414:LEU:HD11	1:A:419:LEU:HD23	1.96	0.47
1:A:518:PHE:HB3	1:A:522:MET:HB3	1.96	0.47
1:C:475:TYR:OH	1:C:510:GLU:OE2	2.30	0.47
1:A:360:LYS:HE2	1:A:362:ASP:HB2	1.95	0.47
1:D:380:GLU:O	1:D:384:LEU:HG	2.14	0.47
1:C:125:ASP:OD1	1:D:543:GLN:NE2	2.48	0.47
1:D:414:LEU:HD11	1:D:419:LEU:HD23	1.96	0.47
1:B:148:TYR:CZ	1:B:221:THR:HB	2.50	0.47
1:C:380:GLU:HG2	1:C:466:TYR:CE1	2.49	0.47
1:C:142:PHE:CE2	1:D:538:PRO:HG3	2.50	0.47
1:C:530:SER:HA	7:C:825:HOH:O	2.15	0.47
1:C:414:LEU:HD11	1:C:419:LEU:HD23	1.97	0.46
1:D:118:THR:O	1:D:122:TYR:HD1	1.97	0.46
1:D:389:PRO:HB2	1:D:434:VAL:HA	1.97	0.46
1:A:89:VAL:O	7:A:802:HOH:O	2.21	0.46
1:B:380:GLU:O	1:B:384:LEU:HG	2.15	0.46
1:A:185:ARG:HD3	1:A:438:ARG:HD3	1.97	0.46
1:A:287:VAL:HG21	1:A:292:PHE:HB2	1.97	0.46
1:B:132:VAL:HG13	1:B:133:HIS:CD2	2.51	0.46
1:B:539:ILE:HA	1:B:544:TYR:HB3	1.98	0.46
1:B:441:PRO:HG2	1:B:444:VAL:HG22	1.98	0.46
1:C:178:LEU:HD21	1:C:445:GLN:HB2	1.98	0.46
1:D:335:ILE:HA	1:D:559:ILE:HD11	1.98	0.46
1:B:117:LEU:HD22	1:B:531:LEU:HD21	1.98	0.46
1:C:124:ILE:HD11	1:C:528:PRO:HB2	1.98	0.45
1:A:44:ARG:NH2	1:A:125:ASP:OD1	2.39	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:438:ARG:NH1	7:D:822:HOH:O	2.48	0.45
1:A:300:MET:HB3	1:A:300:MET:HE3	1.72	0.45
1:B:206:THR:HG21	1:B:385:TYR:CE2	2.51	0.45
1:D:352:LEU:HD22	1:D:518:PHE:CE2	2.52	0.45
1:B:230:LEU:HG	1:B:233:ILE:HD12	1.98	0.45
1:C:296:PRO:HD2	1:C:407:PHE:CE2	2.51	0.45
1:D:46:GLU:CD	1:D:137:LYS:HZ1	2.25	0.45
1:D:380:GLU:HG2	1:D:466:TYR:CE1	2.52	0.45
1:D:108:LEU:HD13	1:D:108:LEU:HA	1.86	0.45
1:A:338:GLY:HA3	1:A:559:ILE:HD13	2.00	0.44
1:B:40:PRO:HB3	4:B:703:NAG:H83	2.00	0.44
1:C:472:LEU:HD11	1:C:524:GLU:HB2	1.99	0.44
1:D:361:PHE:HA	7:D:816:HOH:O	2.17	0.44
1:C:206:THR:HG21	1:C:385:TYR:CE2	2.53	0.44
1:A:364:GLU:HG2	1:A:367:PHE:CE1	2.53	0.44
1:A:380:GLU:O	1:A:384:LEU:HG	2.17	0.44
1:B:67:GLU:O	7:B:801:HOH:O	2.21	0.44
1:A:538:PRO:HG3	1:B:142:PHE:CE2	2.53	0.44
1:B:495:TYR:O	1:B:497:ASP:N	2.49	0.44
1:B:198:PHE:CZ	1:B:352:LEU:HD23	2.53	0.43
1:C:441:PRO:HG2	1:C:444:VAL:HG22	2.00	0.43
1:B:475:TYR:OH	1:B:510:GLU:OE2	2.36	0.43
1:D:34:ASN:HB3	1:D:37:CYS:SG	2.59	0.43
1:A:68:ASN:OD1	1:A:68:ASN:N	2.51	0.43
1:C:360:LYS:HE2	1:C:362:ASP:HB2	2.00	0.43
1:A:307:ARG:NH1	7:A:817:HOH:O	2.51	0.43
1:C:344:VAL:HA	1:C:348:TYR:HB3	2.00	0.43
1:D:206:THR:HG21	1:D:385:TYR:CE2	2.53	0.43
1:A:280:PRO:O	1:A:284:GLN:HG3	2.19	0.43
1:D:198:PHE:HZ	1:D:352:LEU:HD21	1.83	0.43
1:A:142:PHE:CE2	1:B:538:PRO:HG3	2.54	0.43
1:D:468:LYS:HG3	1:D:474:PRO:HG3	2.00	0.43
1:A:89:VAL:C	7:A:802:HOH:O	2.60	0.42
1:A:294:LEU:HA	1:A:409:TYR:CE1	2.55	0.42
1:C:467:ARG:HB3	1:C:472:LEU:HB2	2.01	0.42
1:D:105(A):ILE:HG22	1:D:108:LEU:H	1.85	0.42
1:B:384:LEU:HD13	6:B:701:Y8A:CL22	2.56	0.42
1:B:504:TYR:HA	1:B:507:LEU:HD12	2.00	0.42
1:C:481:LEU:HD12	1:C:510:GLU:HG3	2.02	0.42
1:A:526:GLY:HA3	6:A:707:Y8A:CL22	2.56	0.42
1:A:339:GLU:HG2	1:A:562:ALA:HB2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:472:LEU:HD21	1:B:524:GLU:HG3	2.02	0.42
1:A:472:LEU:HD21	1:A:524:GLU:HG3	2.00	0.42
1:D:382:ASN:O	1:D:386:HIS:HD2	2.03	0.42
1:C:44:ARG:NH2	1:C:125:ASP:OD1	2.50	0.42
1:D:526:GLY:HA3	6:D:701:Y8A:CL22	2.57	0.42
1:A:344:VAL:HA	1:A:348:TYR:HB3	2.02	0.41
1:D:120:ARG:NH1	1:D:524:GLU:OE2	2.49	0.41
1:B:451:SER:O	1:B:455:SER:OG	2.32	0.41
1:D:472:LEU:HD21	1:D:524:GLU:HG3	2.02	0.41
1:B:344:VAL:HA	1:B:348:TYR:HB3	2.02	0.41
1:D:281:GLU:H	1:D:281:GLU:HG3	1.74	0.41
1:C:323:TRP:HE3	1:C:328:LEU:HD23	1.85	0.41
1:C:538:PRO:HG3	1:D:142:PHE:CZ	2.55	0.41
1:D:181:VAL:HG21	1:D:491:LEU:HD21	2.03	0.41
1:A:335:ILE:HA	1:A:559:ILE:HD11	2.03	0.41
1:B:322:GLU:OE2	1:B:322:GLU:N	2.51	0.41
1:A:209:PHE:HB2	1:A:377:ILE:HG13	2.03	0.41
1:C:76:THR:O	1:C:80:LEU:HB2	2.21	0.41
1:D:377:ILE:HD13	1:D:377:ILE:HA	1.89	0.41
1:A:200:PHE:CD1	1:A:297:GLY:HA3	2.56	0.41
1:B:93:LEU:HD23	1:B:100:TRP:CZ2	2.56	0.41
1:B:380:GLU:HG2	1:B:466:TYR:CE1	2.56	0.41
1:C:34:ASN:HB3	1:C:37:CYS:SG	2.61	0.41
1:C:295:VAL:HG23	1:C:298:LEU:HG	2.03	0.41
1:A:120:ARG:HH12	1:A:524:GLU:CD	2.29	0.40
1:D:137:LYS:HE3	1:D:137:LYS:HB2	1.94	0.40
1:A:276:PRO:HD3	1:A:409:TYR:CD1	2.57	0.40
1:C:163:MET:HE2	1:C:171:LEU:HD11	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	553/587 (94%)	533 (96%)	20 (4%)	0	100	100
1	B	553/587 (94%)	533 (96%)	20 (4%)	0	100	100
1	C	553/587 (94%)	534 (97%)	19 (3%)	0	100	100
1	D	553/587 (94%)	533 (96%)	19 (3%)	1 (0%)	43	60
All	All	2212/2348 (94%)	2133 (96%)	78 (4%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	398	GLU

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

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](#)

8 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	E	1	2,1	14,14,15	0.22	0	17,19,21	0.54	0
2	NAG	E	2	2	14,14,15	0.29	0	17,19,21	0.43	0
2	NAG	F	1	2,1	14,14,15	0.33	0	17,19,21	0.47	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	F	2	2	14,14,15	0.32	0	17,19,21	0.44	0
2	NAG	G	1	2,1	14,14,15	0.29	0	17,19,21	0.49	0
2	NAG	G	2	2	14,14,15	0.28	0	17,19,21	0.48	0
2	NAG	H	1	2,1	14,14,15	0.30	0	17,19,21	0.52	0
2	NAG	H	2	2	14,14,15	0.37	0	17,19,21	0.47	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	E	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	E	2	2	-	0/6/23/26	0/1/1/1
2	NAG	F	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	F	2	2	-	0/6/23/26	0/1/1/1
2	NAG	G	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	G	2	2	-	0/6/23/26	0/1/1/1
2	NAG	H	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	H	2	2	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

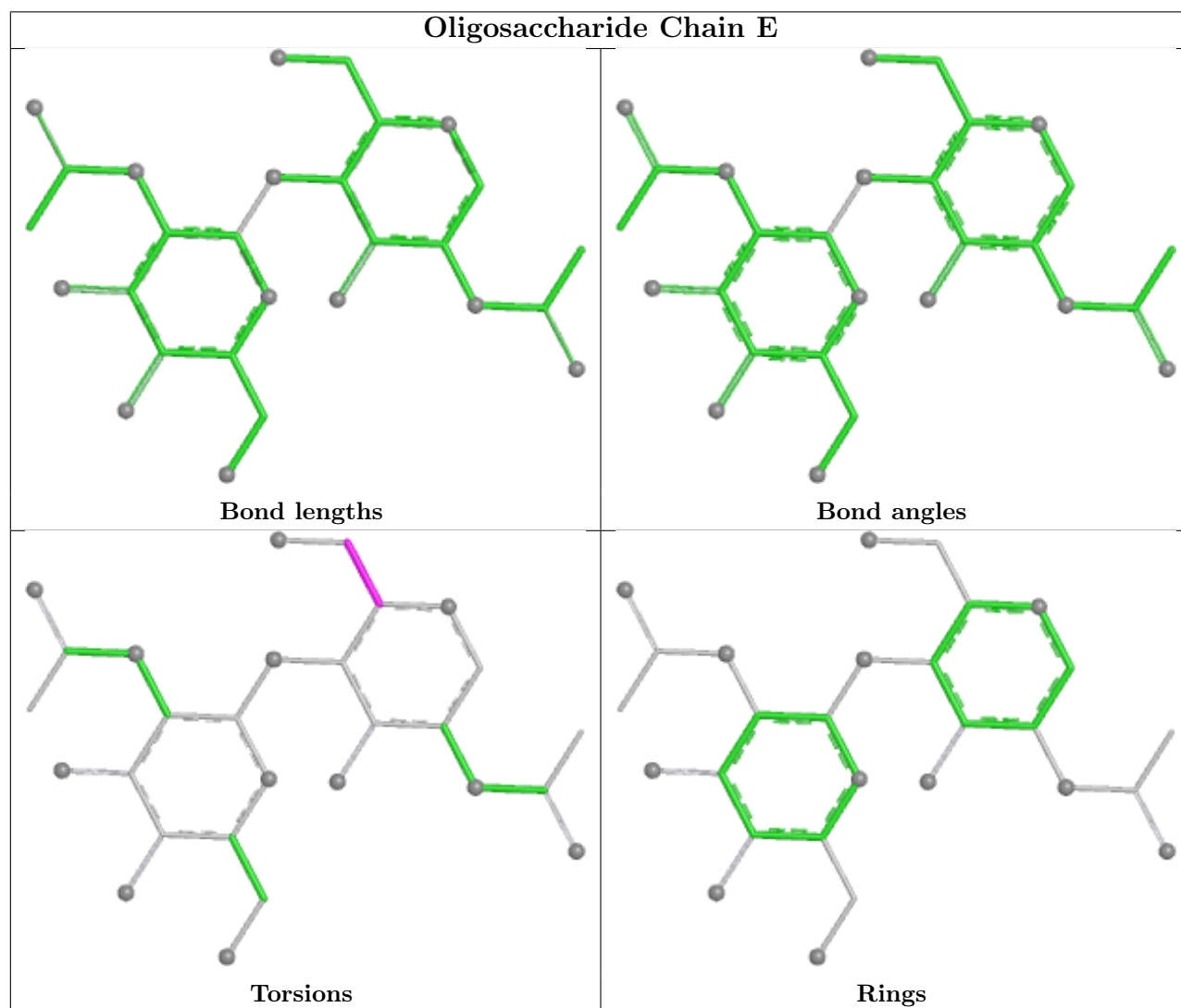
Mol	Chain	Res	Type	Atoms
2	H	2	NAG	O5-C5-C6-O6
2	H	2	NAG	C4-C5-C6-O6
2	E	1	NAG	O5-C5-C6-O6
2	E	1	NAG	C4-C5-C6-O6
2	F	1	NAG	C4-C5-C6-O6
2	F	1	NAG	O5-C5-C6-O6

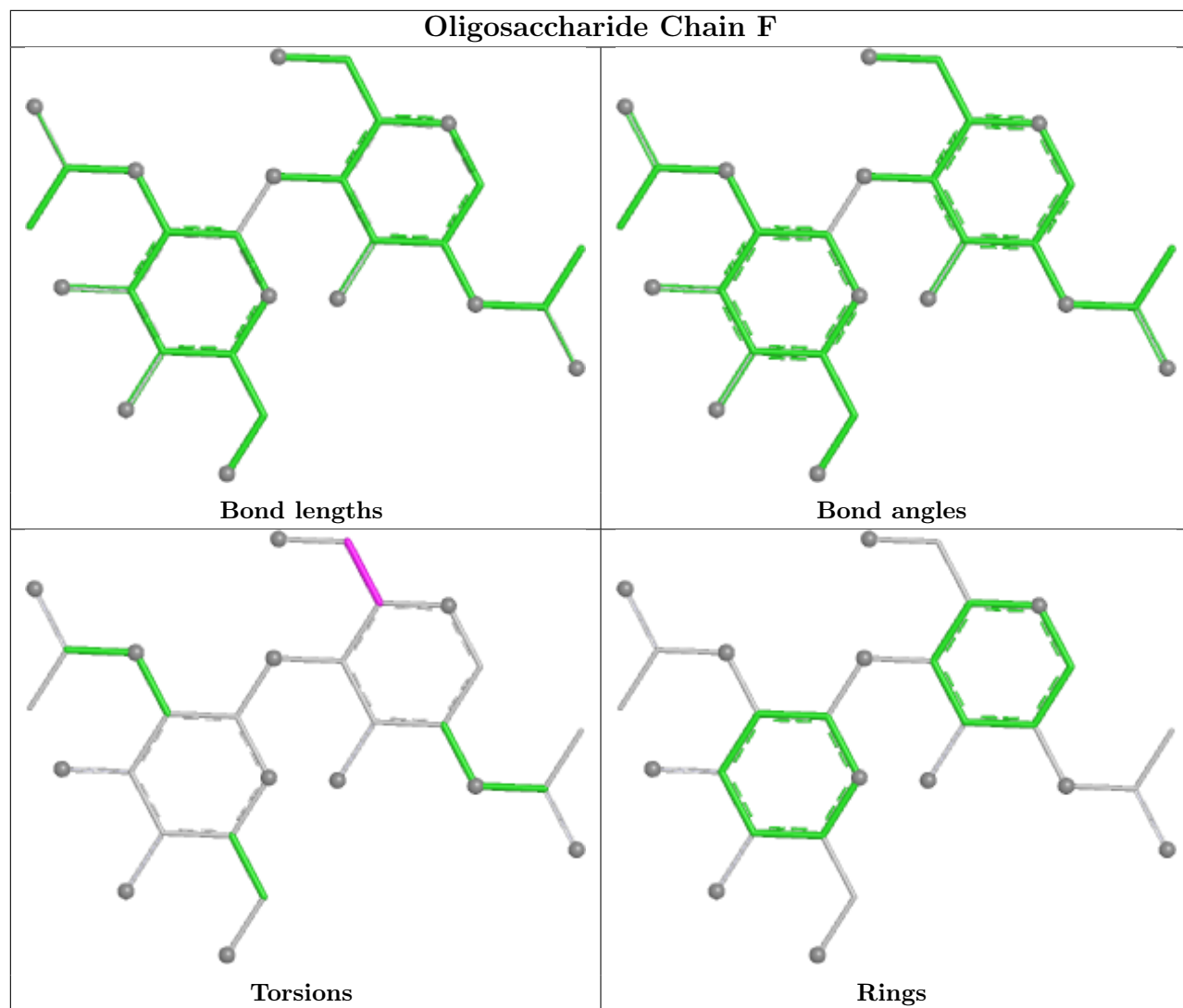
There are no ring outliers.

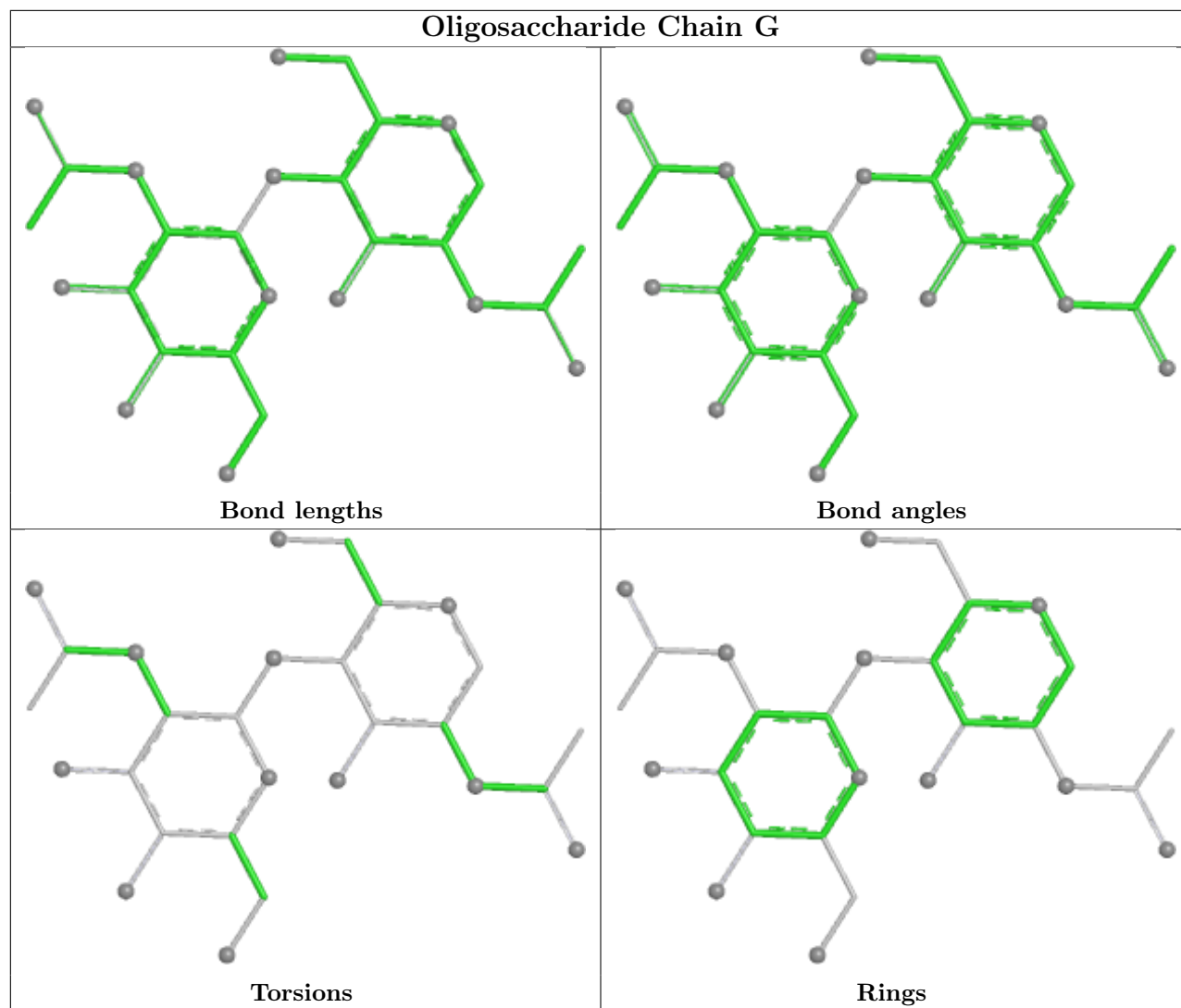
1 monomer is involved in 1 short contact:

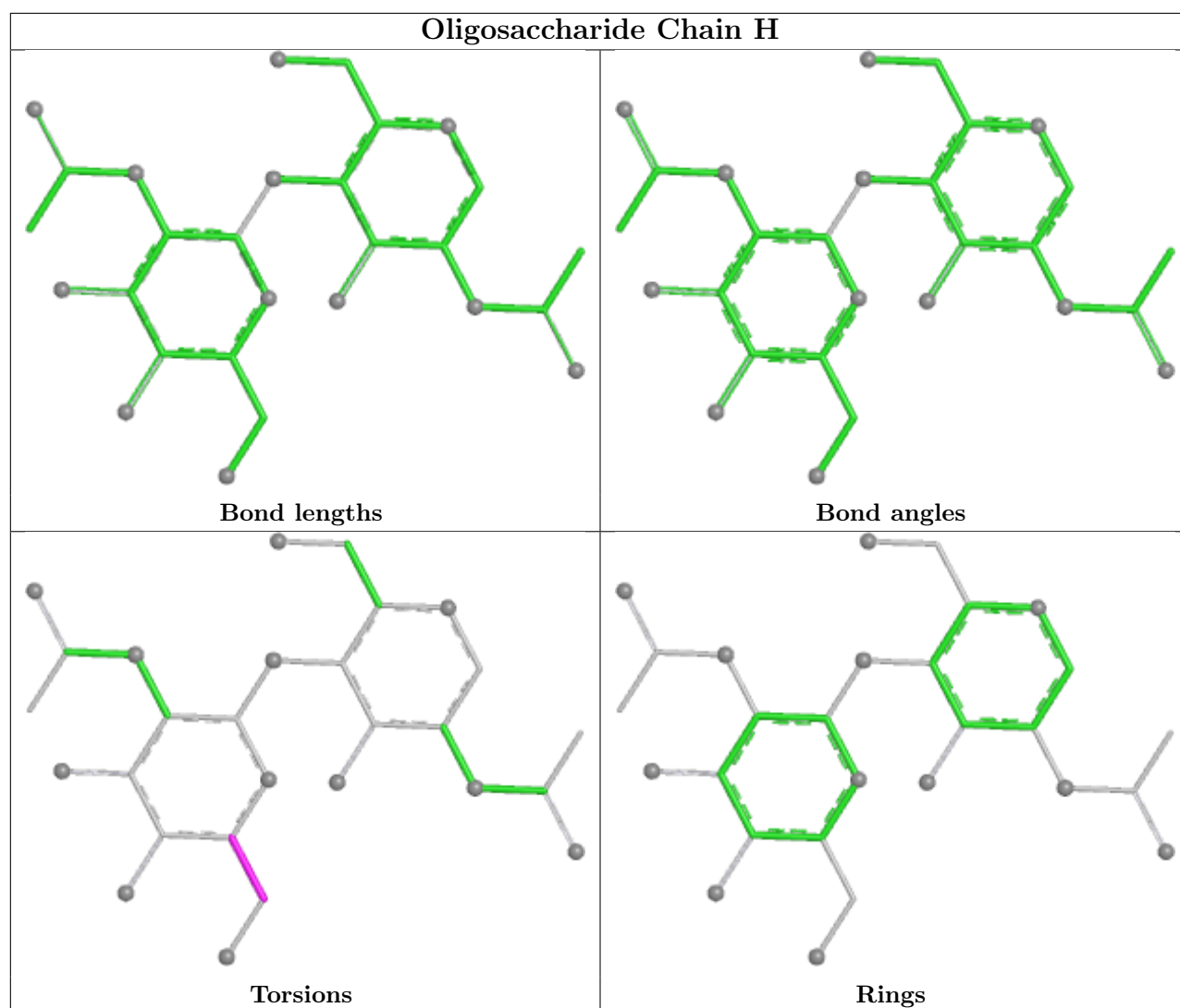
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	H	2	NAG	1	0

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](#)

18 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	$\# Z > 2$
4	NAG	A	705	1	14,14,15	0.22	0	17,19,21	0.48	0
4	NAG	B	706	1	14,14,15	0.22	0	17,19,21	0.45	0
6	Y8A	D	701	-	25,27,27	1.42	3 (12%)	34,39,39	1.22	3 (8%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	D	703	1	14,14,15	0.34	0	17,19,21	0.40	0
4	NAG	A	702	1	14,14,15	0.44	0	17,19,21	0.36	0
6	Y8A	A	707	-	25,27,27	1.43	4 (16%)	34,39,39	1.15	2 (5%)
4	NAG	D	706	1	14,14,15	0.27	0	17,19,21	0.44	0
6	Y8A	B	701	-	25,27,27	1.44	3 (12%)	34,39,39	1.34	4 (11%)
3	HEM	A	701	-	50,50,50	1.48	6 (12%)	67,82,82	1.27	5 (7%)
3	HEM	C	702	1	50,50,50	1.45	7 (14%)	67,82,82	1.26	6 (8%)
4	NAG	B	703	1	14,14,15	0.34	0	17,19,21	0.58	0
4	NAG	C	706	1	14,14,15	0.25	0	17,19,21	0.44	0
5	BOG	A	706	-	20,20,20	1.23	2 (10%)	25,25,25	0.79	0
5	BOG	D	707	-	20,20,20	1.22	2 (10%)	25,25,25	0.75	0
4	NAG	C	703	1	14,14,15	0.27	0	17,19,21	0.63	0
6	Y8A	C	701	-	25,27,27	1.43	4 (16%)	34,39,39	1.13	2 (5%)
3	HEM	B	702	-	50,50,50	1.52	8 (16%)	67,82,82	1.27	8 (11%)
3	HEM	D	702	1	50,50,50	1.48	7 (14%)	67,82,82	1.26	6 (8%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	A	705	1	-	2/6/23/26	0/1/1/1
4	NAG	B	706	1	-	0/6/23/26	0/1/1/1
6	Y8A	D	701	-	-	2/6/16/16	0/4/4/4
4	NAG	D	703	1	-	3/6/23/26	0/1/1/1
4	NAG	A	702	1	-	0/6/23/26	0/1/1/1
6	Y8A	A	707	-	-	2/6/16/16	0/4/4/4
4	NAG	D	706	1	-	1/6/23/26	0/1/1/1
6	Y8A	B	701	-	-	2/6/16/16	0/4/4/4
3	HEM	A	701	-	-	4/14/54/54	-
3	HEM	C	702	1	-	2/14/54/54	-
4	NAG	B	703	1	-	4/6/23/26	0/1/1/1
4	NAG	C	706	1	-	2/6/23/26	0/1/1/1
5	BOG	A	706	-	-	6/11/31/31	0/1/1/1
5	BOG	D	707	-	-	4/11/31/31	0/1/1/1
4	NAG	C	703	1	-	4/6/23/26	0/1/1/1
6	Y8A	C	701	-	-	2/6/16/16	0/4/4/4

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	HEM	B	702	-	-	2/14/54/54	-
3	HEM	D	702	1	-	2/14/54/54	-

All (46) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	701	Y8A	C13-C09	4.17	1.55	1.50
3	A	701	HEM	FE-NB	4.14	2.07	1.94
3	B	702	HEM	FE-ND	4.10	2.07	1.94
6	C	701	Y8A	C13-C09	4.05	1.55	1.50
6	A	707	Y8A	C13-C09	4.04	1.55	1.50
6	D	701	Y8A	C13-C09	4.00	1.55	1.50
3	B	702	HEM	FE-NB	3.97	2.07	1.94
3	D	702	HEM	FE-NB	3.90	2.06	1.94
3	B	702	HEM	FE-NA	3.84	2.07	1.95
3	C	702	HEM	FE-ND	3.84	2.06	1.94
3	A	701	HEM	FE-ND	3.71	2.06	1.94
3	D	702	HEM	FE-ND	3.66	2.06	1.94
3	C	702	HEM	FE-NA	3.54	2.06	1.95
5	A	706	BOG	O5-C1	3.52	1.50	1.41
5	D	707	BOG	O5-C1	3.48	1.50	1.41
3	C	702	HEM	FE-NB	3.45	2.05	1.94
3	D	702	HEM	FE-NA	3.23	2.05	1.95
3	A	701	HEM	FE-NA	3.21	2.05	1.95
3	D	702	HEM	FE-NC	3.13	2.05	1.95
3	D	702	HEM	CAC-C3C	3.01	1.55	1.47
3	B	702	HEM	CAC-C3C	3.00	1.55	1.47
3	C	702	HEM	CAB-C3B	3.00	1.55	1.47
3	C	702	HEM	CAC-C3C	2.99	1.55	1.47
3	B	702	HEM	CAB-C3B	2.98	1.55	1.47
3	D	702	HEM	CAB-C3B	2.97	1.55	1.47
6	D	701	Y8A	C06-C05	-2.95	1.37	1.41
3	A	701	HEM	FE-NC	2.93	2.04	1.95
6	A	707	Y8A	C06-C05	-2.92	1.37	1.41
3	A	701	HEM	CAB-C3B	2.88	1.55	1.47
6	C	701	Y8A	C06-C05	-2.86	1.37	1.41
3	A	701	HEM	CAC-C3C	2.84	1.55	1.47
6	B	701	Y8A	C06-C05	-2.70	1.37	1.41
3	B	702	HEM	FE-NC	2.57	2.03	1.95
5	D	707	BOG	O1-C1	-2.36	1.36	1.40
5	A	706	BOG	O1-C1	-2.34	1.36	1.40
3	C	702	HEM	FE-NC	2.26	2.02	1.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	701	Y8A	C04-C05	2.11	1.42	1.39
6	D	701	Y8A	C05-N07	-2.08	1.35	1.39
6	A	707	Y8A	C04-C05	2.06	1.42	1.39
6	A	707	Y8A	C05-N07	-2.06	1.35	1.39
6	C	701	Y8A	C05-N07	-2.05	1.35	1.39
3	B	702	HEM	CMD-C2D	2.04	1.55	1.50
6	C	701	Y8A	C04-C05	2.02	1.42	1.39
3	D	702	HEM	CMA-C3A	2.01	1.54	1.50
3	B	702	HEM	CMA-C3A	2.01	1.54	1.50
3	C	702	HEM	CMD-C2D	2.01	1.54	1.50

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	701	Y8A	C12-C13-C09	4.76	115.39	109.14
6	D	701	Y8A	C10-C08-N07	3.38	134.45	128.90
6	B	701	Y8A	C10-C08-N07	3.28	134.29	128.90
6	C	701	Y8A	C12-C13-C09	3.24	113.39	109.14
6	A	707	Y8A	C10-C08-N07	3.21	134.18	128.90
6	D	701	Y8A	C12-C13-C09	3.03	113.11	109.14
3	D	702	HEM	C3B-C2B-C1B	3.03	108.69	106.41
3	C	702	HEM	C3B-C2B-C1B	3.03	108.68	106.41
6	C	701	Y8A	C10-C08-N07	2.97	133.78	128.90
6	A	707	Y8A	C12-C13-C09	2.91	112.95	109.14
3	A	701	HEM	C1B-NB-C4B	2.88	108.62	105.21
3	A	701	HEM	C3B-C4B-NB	-2.86	107.42	109.47
3	A	701	HEM	C3B-C2B-C1B	2.79	108.51	106.41
3	B	702	HEM	C3B-C2B-C1B	2.77	108.49	106.41
3	C	702	HEM	C1B-NB-C4B	2.62	108.31	105.21
3	C	702	HEM	C3B-C4B-NB	-2.50	107.67	109.47
3	B	702	HEM	C4D-ND-C1D	2.50	108.17	105.21
3	D	702	HEM	C1B-NB-C4B	2.49	108.15	105.21
3	D	702	HEM	C3B-C4B-NB	-2.48	107.69	109.47
3	C	702	HEM	C4D-ND-C1D	2.44	108.10	105.21
3	B	702	HEM	C1B-NB-C4B	2.41	108.06	105.21
3	D	702	HEM	C4D-ND-C1D	2.40	108.05	105.21
3	B	702	HEM	C3B-C4B-NB	-2.33	107.79	109.47
3	A	701	HEM	C4D-ND-C1D	2.30	107.93	105.21
3	B	702	HEM	CAA-CBA-CGA	-2.26	107.67	113.67
3	B	702	HEM	CHC-C1C-NC	2.24	126.89	124.45
3	A	701	HEM	CAA-CBA-CGA	-2.22	107.77	113.67
6	B	701	Y8A	C12-N11-C10	2.18	121.44	117.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	702	HEM	C3D-C4D-ND	-2.16	107.80	110.17
3	B	702	HEM	C3D-C4D-ND	-2.16	107.81	110.17
3	B	702	HEM	C2A-C1A-NA	-2.09	107.84	110.15
3	D	702	HEM	C2A-C1A-NA	-2.07	107.86	110.15
3	C	702	HEM	C3D-C4D-ND	-2.04	107.93	110.17
6	B	701	Y8A	C24-O23-C02	-2.04	113.13	117.50
6	D	701	Y8A	C15-N07-C05	-2.03	121.67	124.79
3	C	702	HEM	C2A-C1A-NA	-2.02	107.91	110.15

There are no chirality outliers.

All (44) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	D	707	BOG	C2'-C1'-O1-C1
6	A	707	Y8A	C01-C02-O23-C24
6	B	701	Y8A	C01-C02-O23-C24
6	B	701	Y8A	C03-C02-O23-C24
6	C	701	Y8A	C03-C02-O23-C24
6	D	701	Y8A	C03-C02-O23-C24
6	A	707	Y8A	C03-C02-O23-C24
6	C	701	Y8A	C01-C02-O23-C24
6	D	701	Y8A	C01-C02-O23-C24
4	D	703	NAG	O5-C5-C6-O6
5	D	707	BOG	C4-C5-C6-O6
4	C	703	NAG	O5-C5-C6-O6
4	C	703	NAG	C4-C5-C6-O6
4	B	703	NAG	O5-C5-C6-O6
5	D	707	BOG	O5-C5-C6-O6
4	D	703	NAG	C4-C5-C6-O6
4	C	706	NAG	C4-C5-C6-O6
4	B	703	NAG	C4-C5-C6-O6
4	A	705	NAG	C4-C5-C6-O6
4	C	706	NAG	O5-C5-C6-O6
5	A	706	BOG	O1-C1'-C2'-C3'
5	A	706	BOG	C1'-C2'-C3'-C4'
4	A	705	NAG	O5-C5-C6-O6
5	D	707	BOG	C1'-C2'-C3'-C4'
3	A	701	HEM	C4B-C3B-CAB-CBB
3	A	701	HEM	C4C-C3C-CAC-CBC
3	B	702	HEM	C4B-C3B-CAB-CBB
3	B	702	HEM	C4C-C3C-CAC-CBC
3	C	702	HEM	C4B-C3B-CAB-CBB

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Mol	Chain	Res	Type	Atoms
3	C	702	HEM	C4C-C3C-CAC-CBC
3	D	702	HEM	C4B-C3B-CAB-CBB
3	D	702	HEM	C4C-C3C-CAC-CBC
5	A	706	BOG	C3'-C4'-C5'-C6'
4	B	703	NAG	C3-C2-N2-C7
4	C	703	NAG	C3-C2-N2-C7
5	A	706	BOG	C2'-C3'-C4'-C5'
4	D	706	NAG	C4-C5-C6-O6
4	B	703	NAG	C1-C2-N2-C7
4	C	703	NAG	C1-C2-N2-C7
4	D	703	NAG	C3-C2-N2-C7
3	A	701	HEM	CAA-CBA-CGA-O2A
3	A	701	HEM	CAA-CBA-CGA-O1A
5	A	706	BOG	C2'-C1'-O1-C1
5	A	706	BOG	C2-C1-O1-C1'

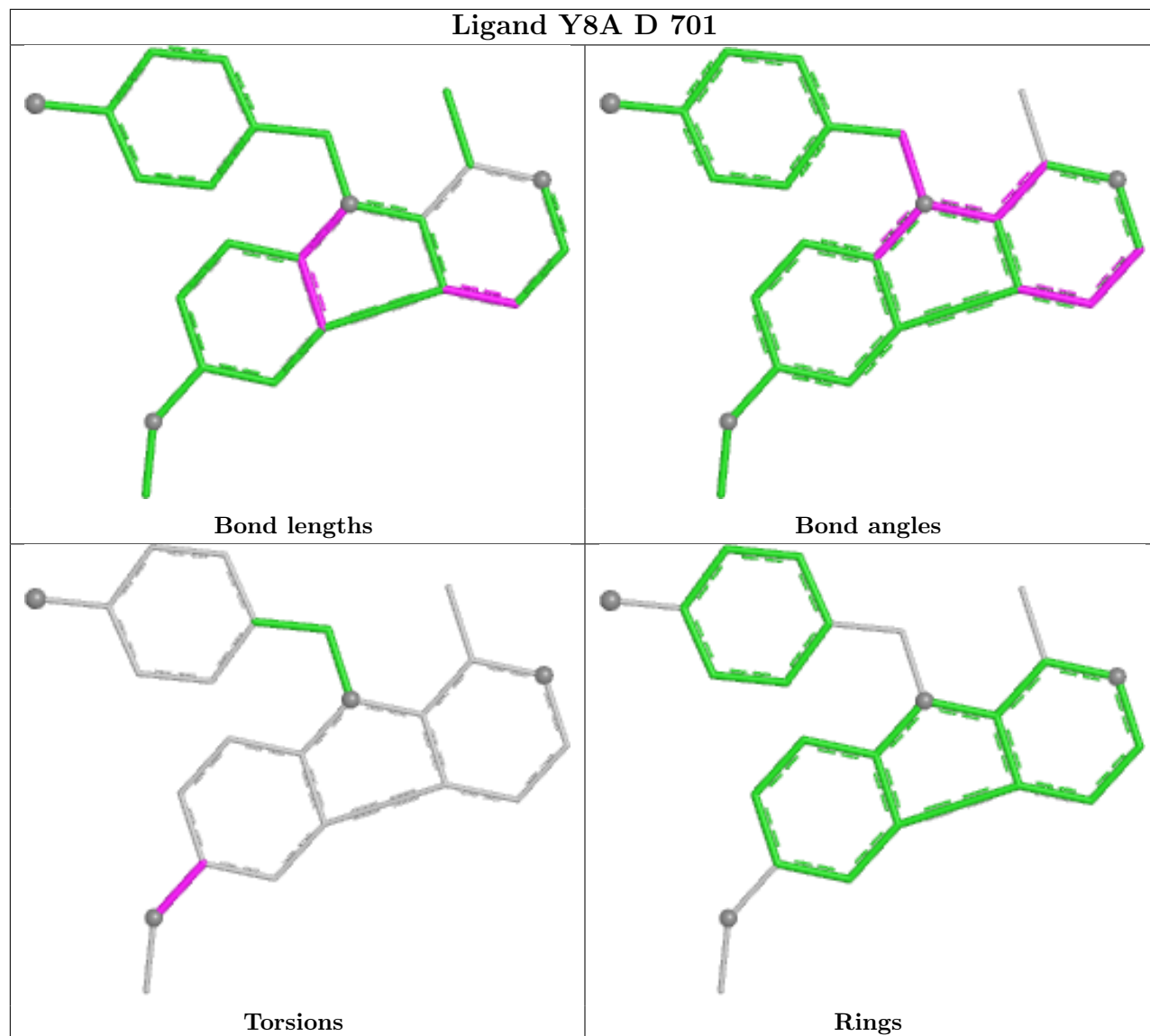
There are no ring outliers.

10 monomers are involved in 16 short contacts:

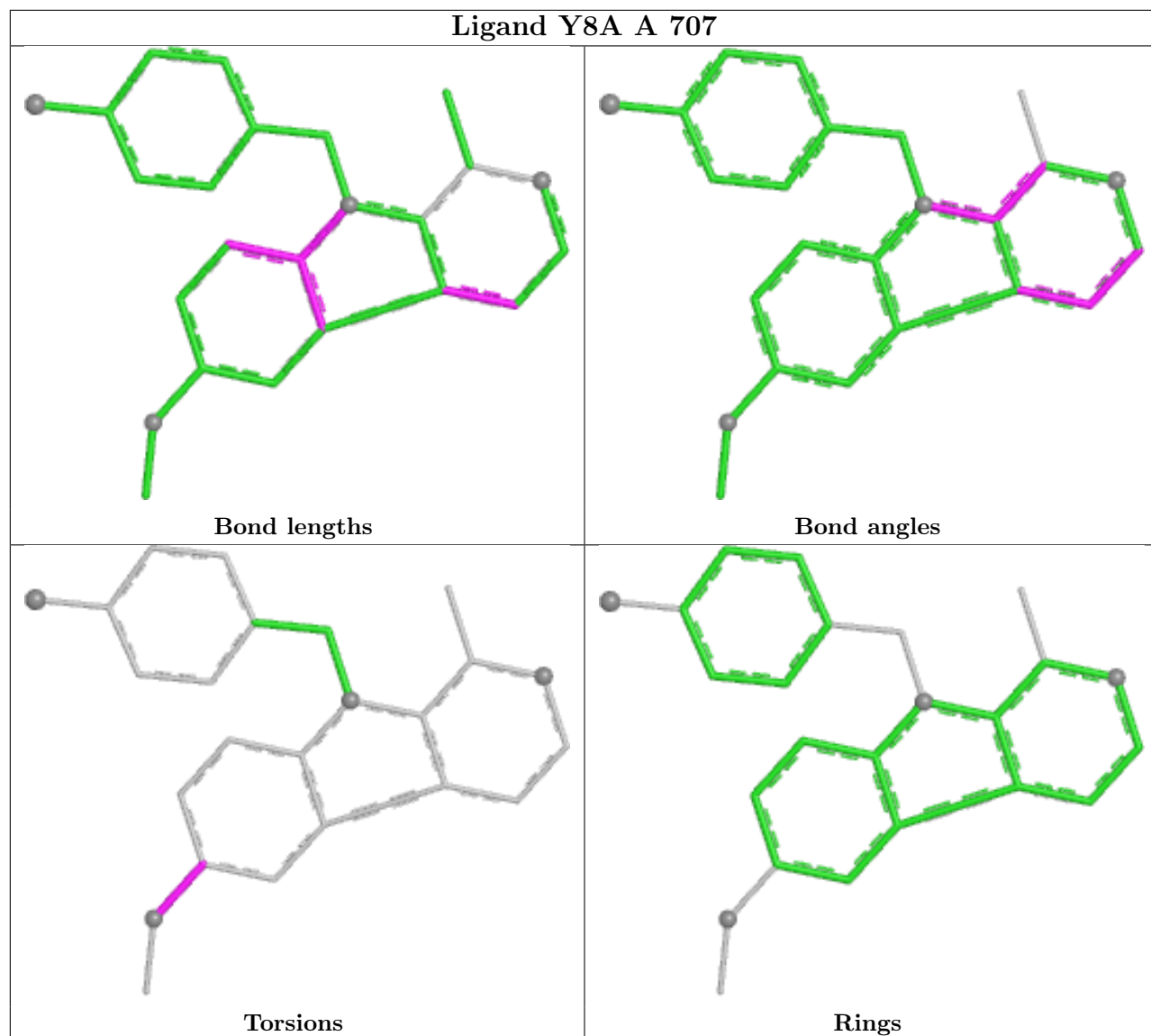
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	701	Y8A	2	0
6	A	707	Y8A	2	0
6	B	701	Y8A	1	0
3	A	701	HEM	2	0
3	C	702	HEM	2	0
4	B	703	NAG	1	0
5	A	706	BOG	1	0
6	C	701	Y8A	1	0
3	B	702	HEM	2	0
3	D	702	HEM	2	0

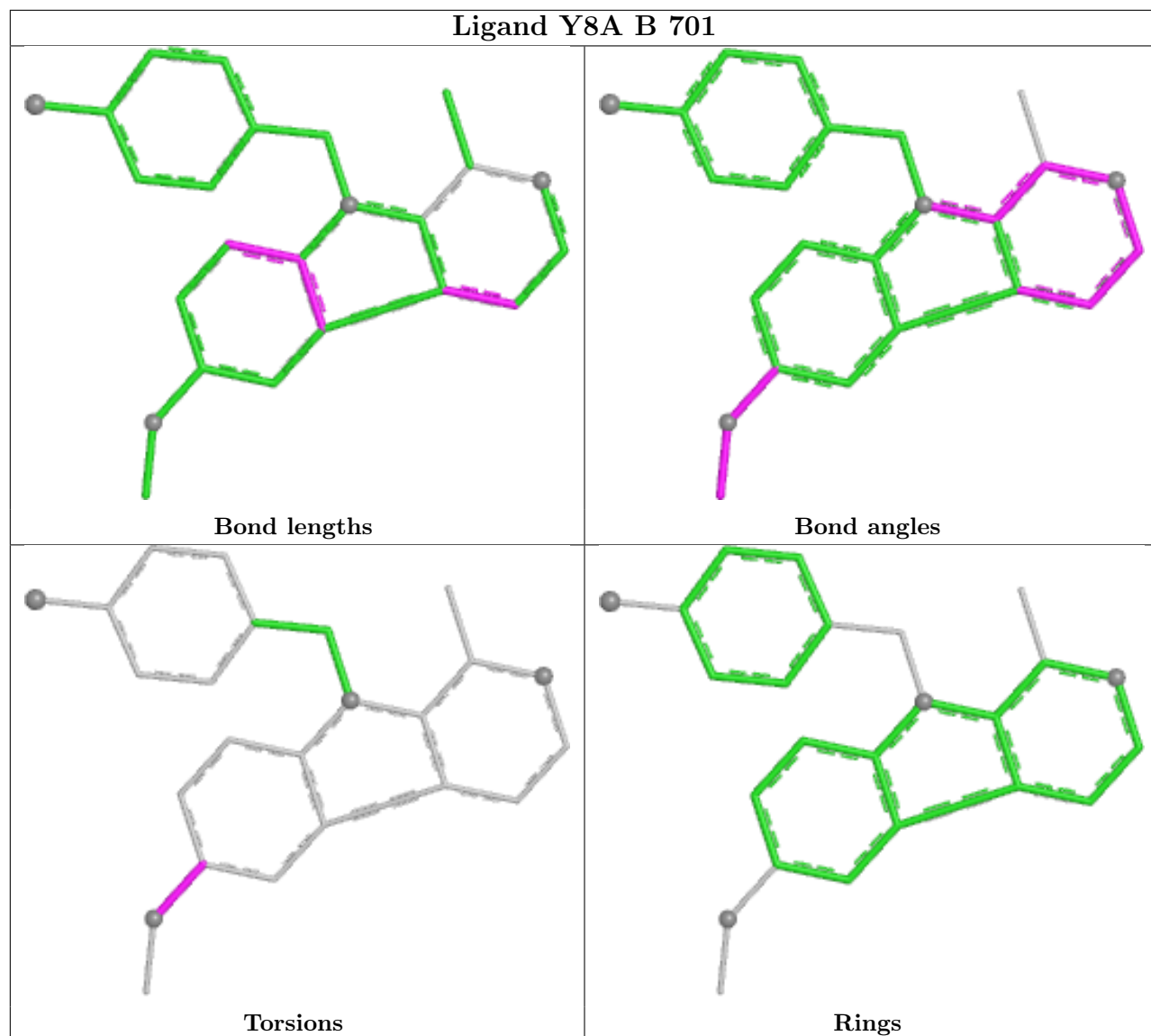
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

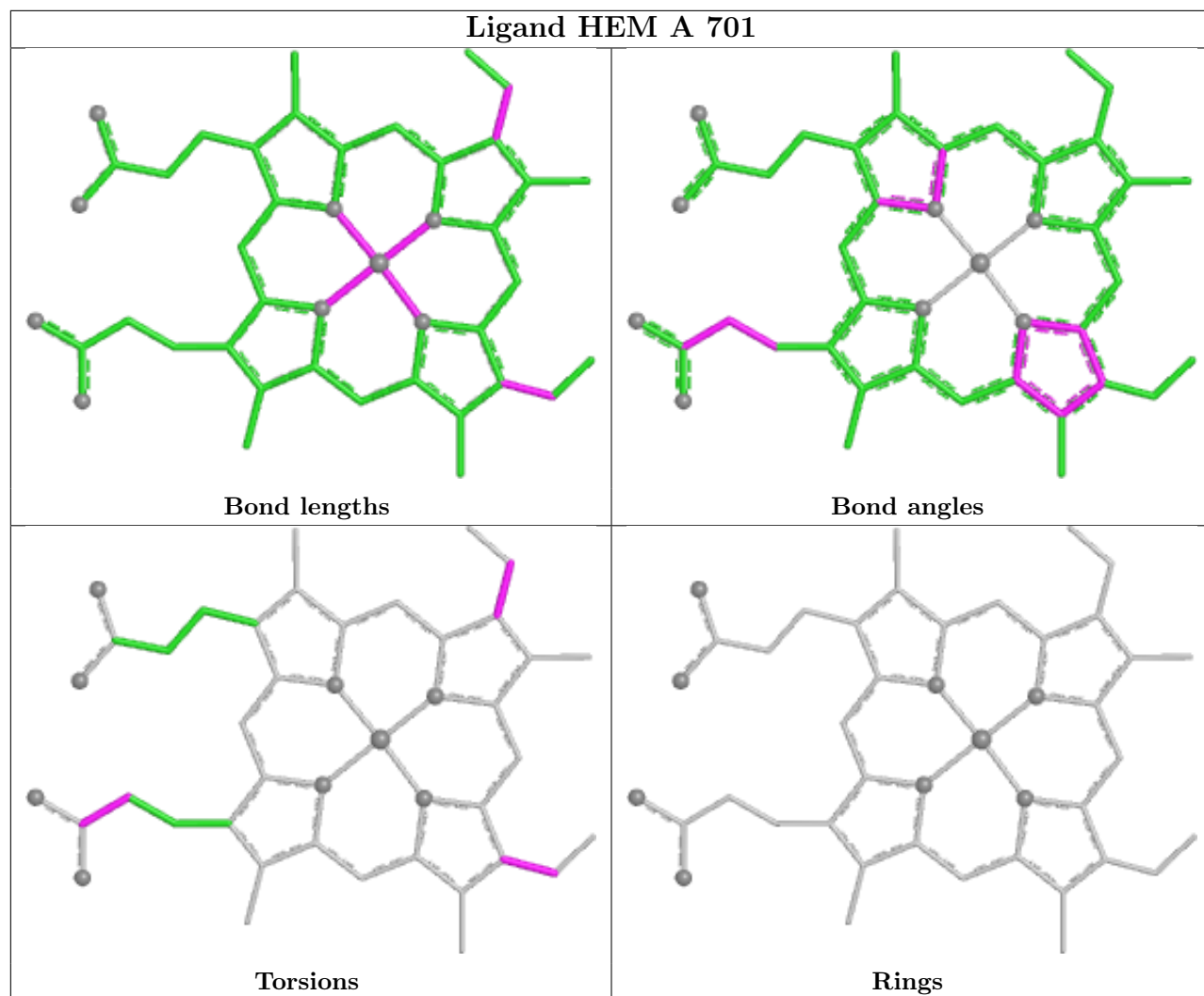
equivalents in the CSD to analyse the geometry.

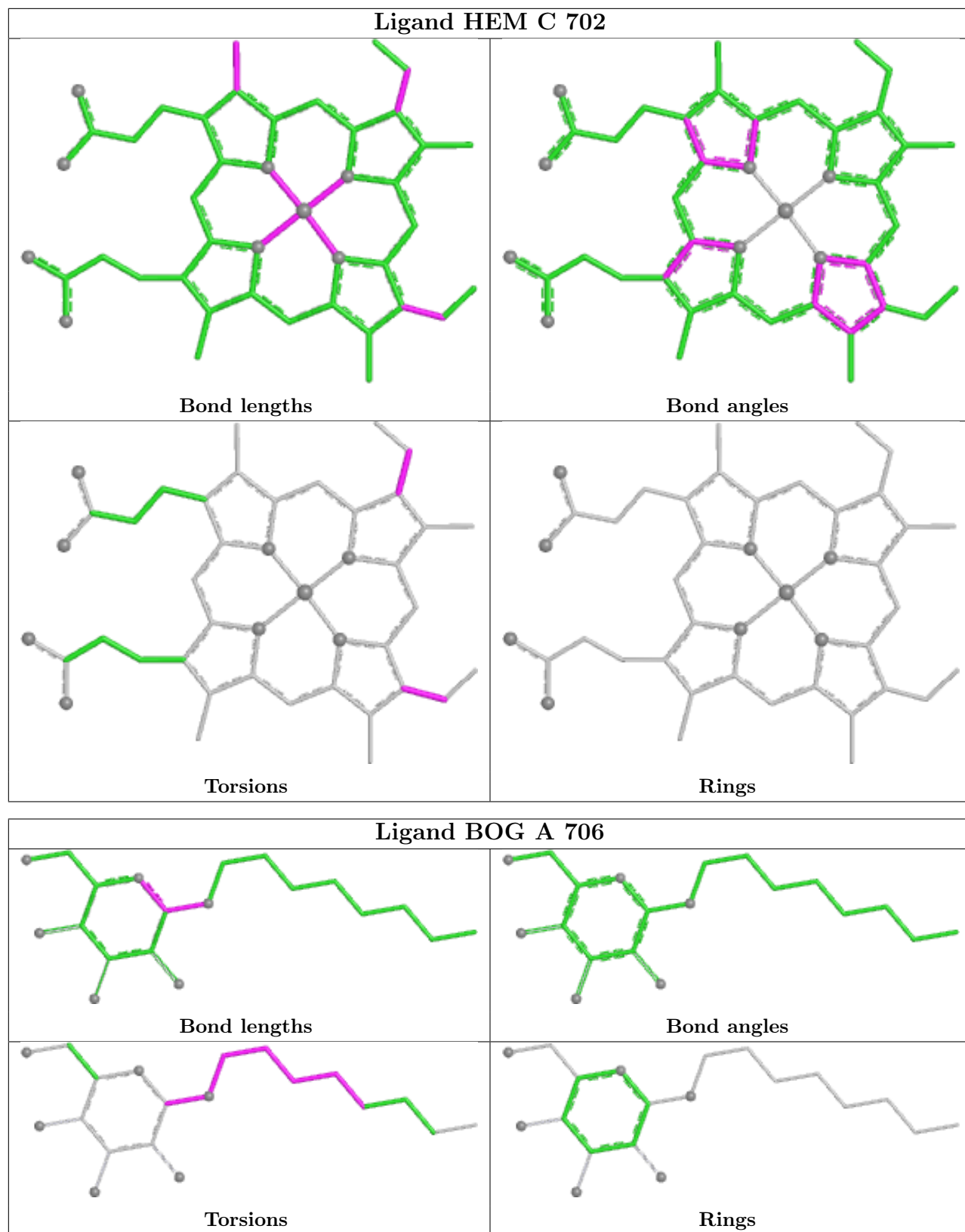


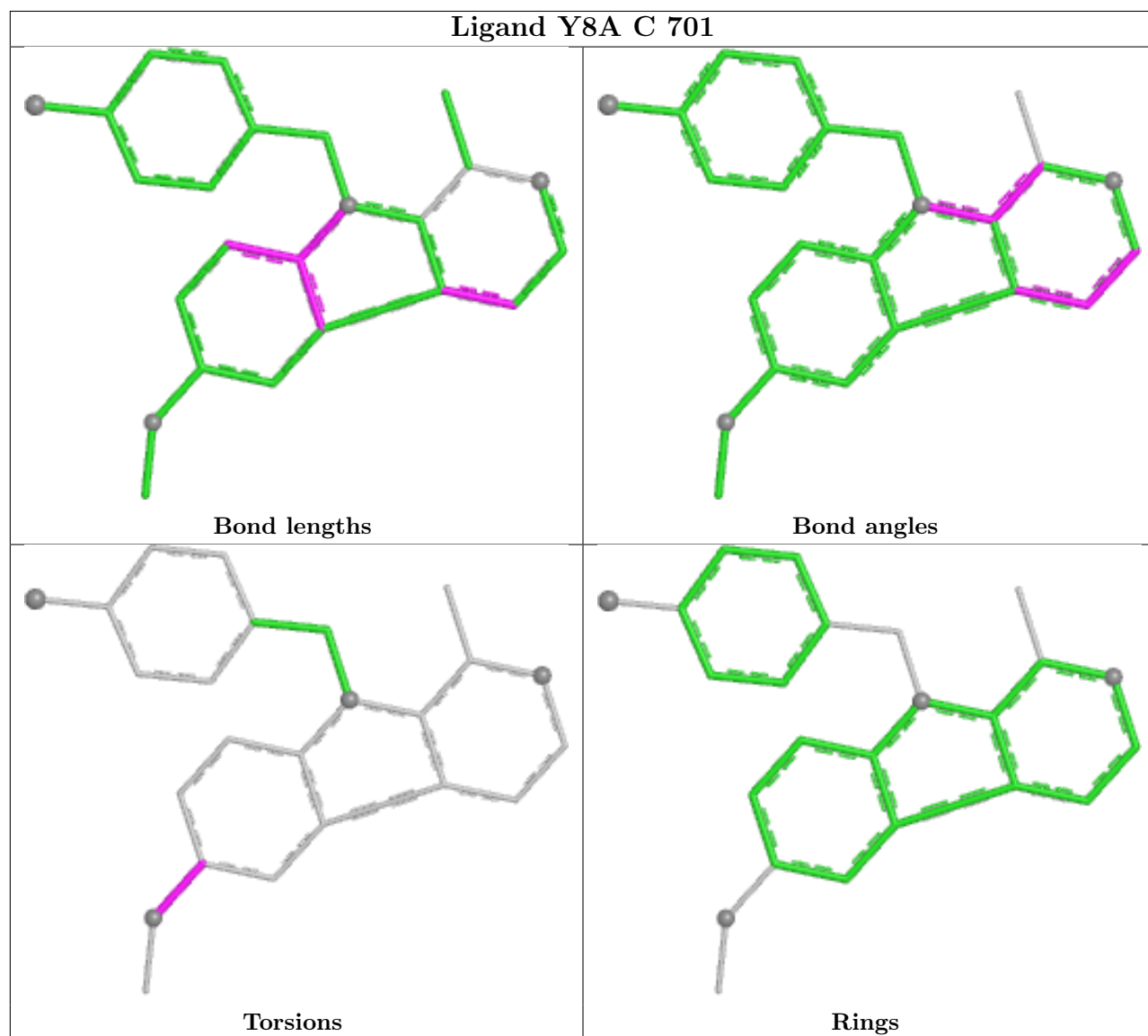
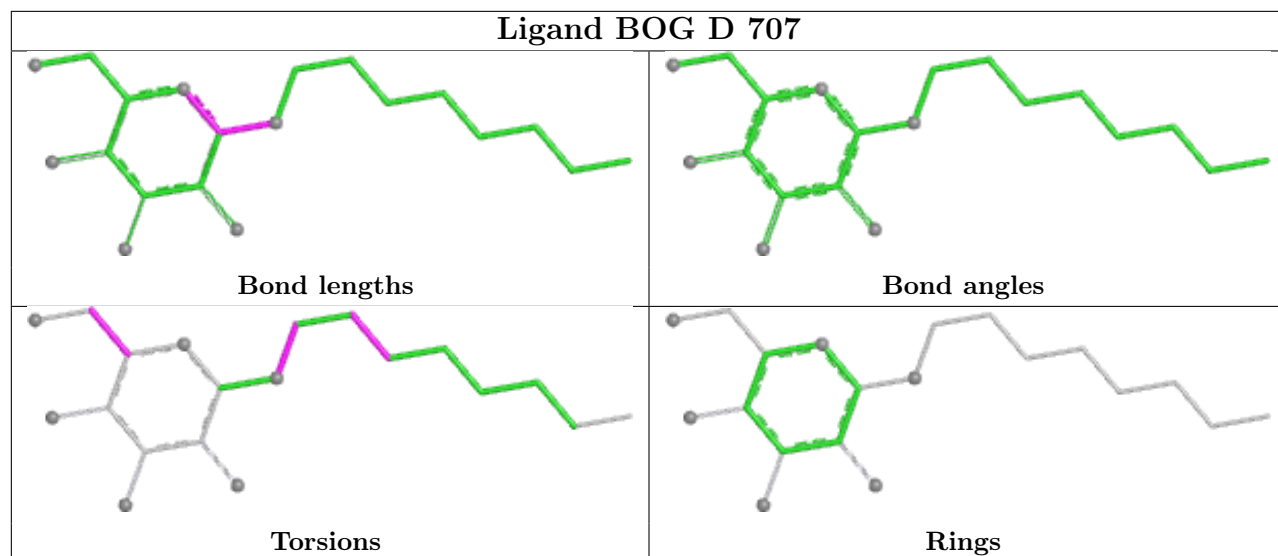
Ligand Y8A A 707

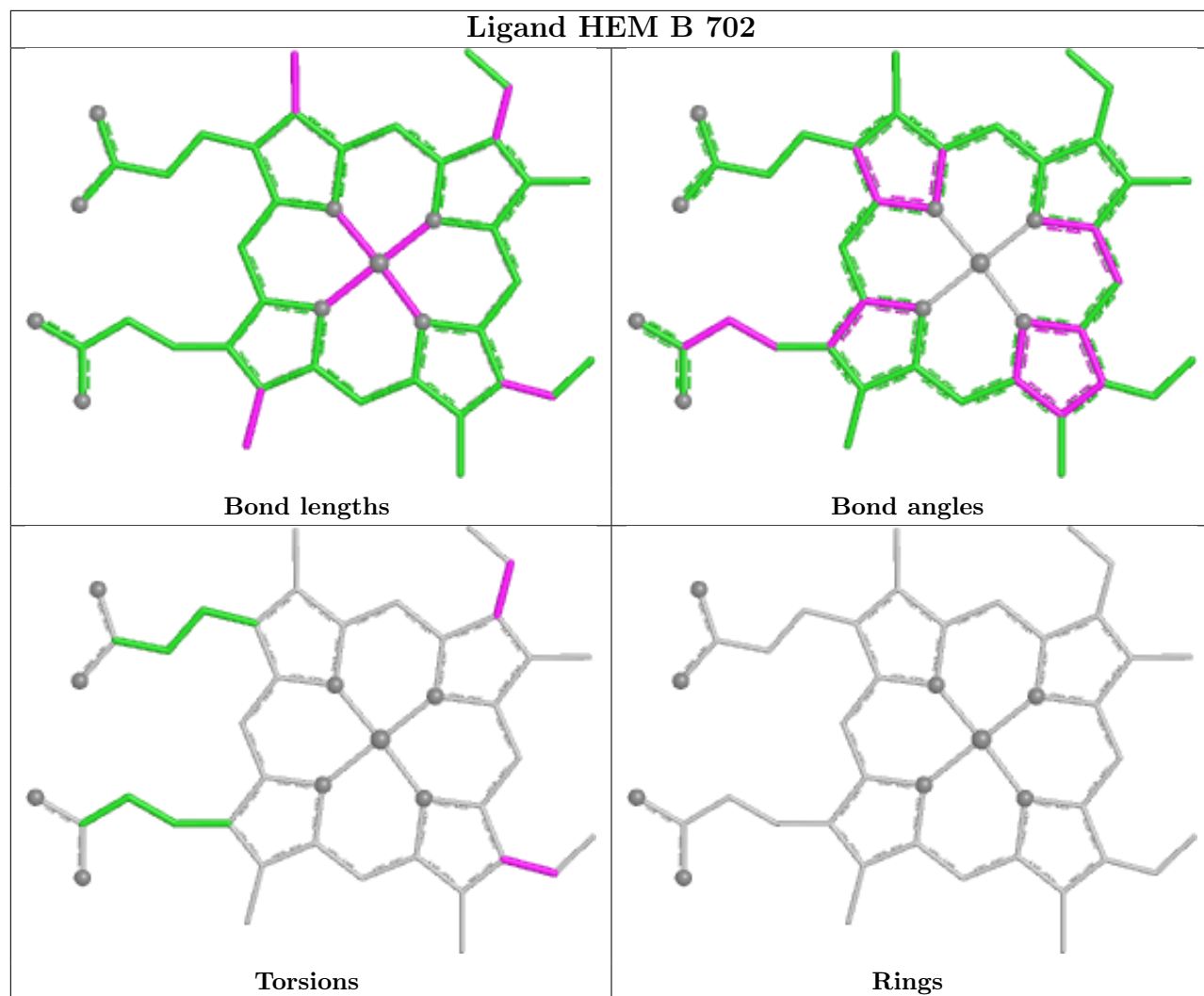


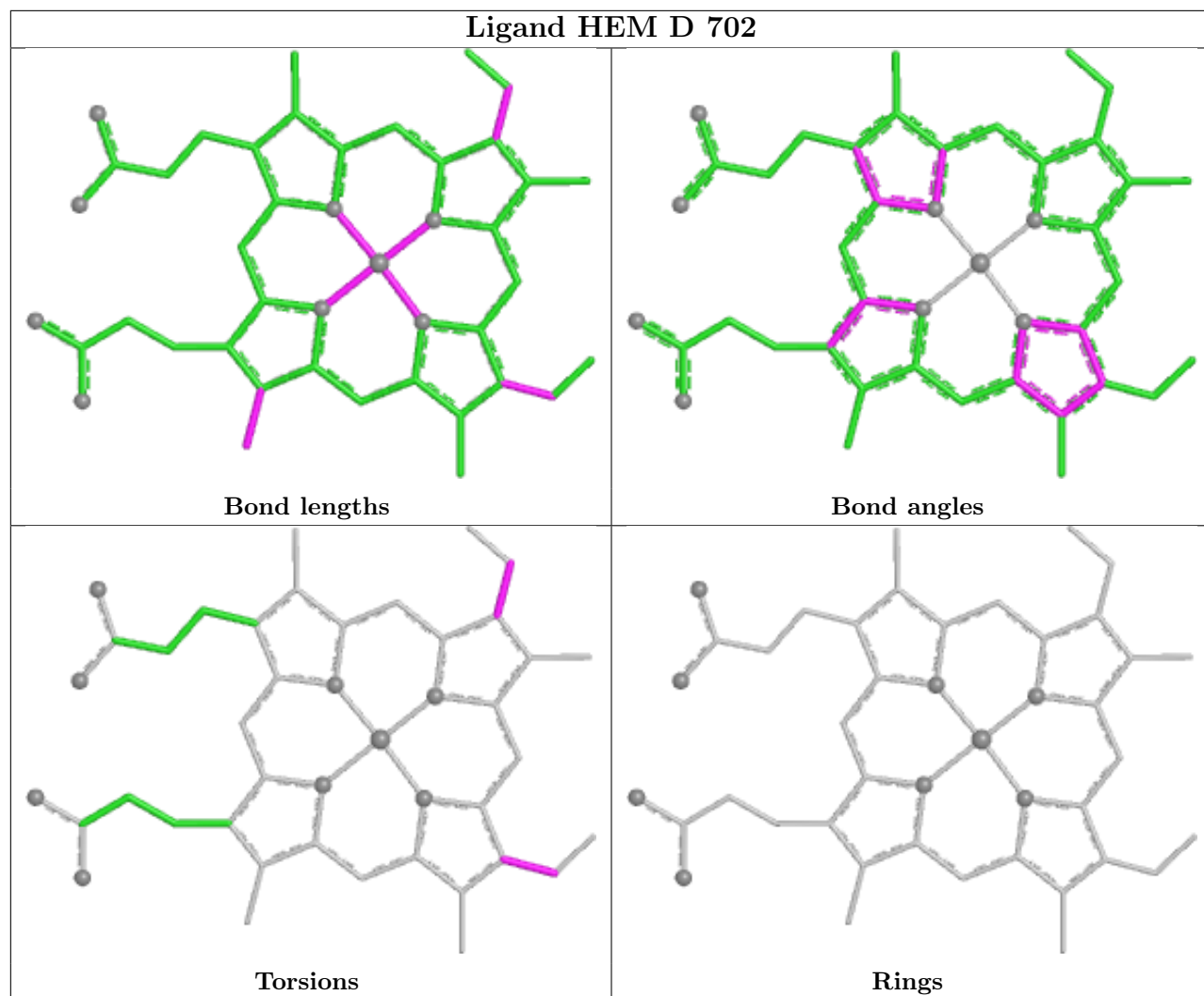












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	552/587 (94%)	0.31	13 (2%) 59 53	24, 47, 70, 87	3 (0%)
1	B	552/587 (94%)	0.37	24 (4%) 40 32	21, 47, 72, 90	3 (0%)
1	C	552/587 (94%)	0.34	15 (2%) 56 49	21, 46, 70, 84	3 (0%)
1	D	552/587 (94%)	0.30	20 (3%) 46 38	24, 48, 68, 84	3 (0%)
All	All	2208/2348 (94%)	0.33	72 (3%) 49 40	21, 47, 70, 90	12 (0%)

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	282	ASN	5.4
1	B	122	TYR	5.1
1	B	121	SER	4.6
1	C	122	TYR	4.3
1	C	81	LEU	4.2
1	A	122	TYR	4.1
1	A	409	TYR	3.8
1	A	75	LEU	3.6
1	B	280	PRO	3.5
1	B	282	ASN	3.5
1	B	281	GLU	3.5
1	A	294	LEU	3.4
1	D	107	PHE	3.4
1	B	33	ALA	3.3
1	D	121	SER	3.2
1	D	409	TYR	3.1
1	D	583	GLN	3.0
1	B	283	LEU	3.0
1	A	583	GLN	3.0
1	B	107	PHE	3.0
1	C	82	LEU	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	33	ALA	3.0
1	D	122	TYR	2.9
1	A	81	LEU	2.9
1	C	115	TYR	2.8
1	D	115	TYR	2.8
1	C	239	ASP	2.8
1	B	531	LEU	2.7
1	A	281	GLU	2.7
1	C	352	LEU	2.7
1	C	268	ASP	2.7
1	B	81	LEU	2.7
1	B	75	LEU	2.6
1	A	121	SER	2.6
1	B	115	TYR	2.6
1	C	53	ASP	2.6
1	C	107	PHE	2.5
1	B	108	LEU	2.5
1	D	514	PRO	2.4
1	B	78	ILE	2.3
1	C	108	LEU	2.3
1	C	121	SER	2.3
1	B	68	ASN	2.3
1	B	444	VAL	2.3
1	D	81	LEU	2.3
1	D	294	LEU	2.3
1	B	364	GLU	2.3
1	D	78	ILE	2.2
1	B	268	ASP	2.2
1	B	79	LYS	2.2
1	B	235	GLY	2.2
1	D	63	GLY	2.2
1	C	80	LEU	2.2
1	D	67	GLU	2.2
1	D	531	LEU	2.2
1	A	115	TYR	2.1
1	B	267	LYS	2.1
1	B	80	LEU	2.1
1	B	53	ASP	2.1
1	D	566	SER	2.1
1	C	236	GLU	2.1
1	D	74	PHE	2.1
1	A	214	HIS	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	59	CYS	2.1
1	A	74	PHE	2.1
1	C	416	GLU	2.0
1	A	119	SER	2.0
1	D	215	LYS	2.0
1	D	75	LEU	2.0
1	D	281	GLU	2.0
1	D	165	VAL	2.0
1	B	409	TYR	2.0

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

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

6.3 Carbohydrates [i](#)

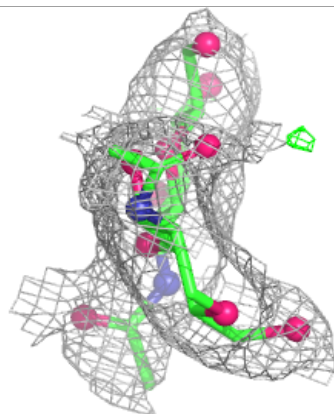
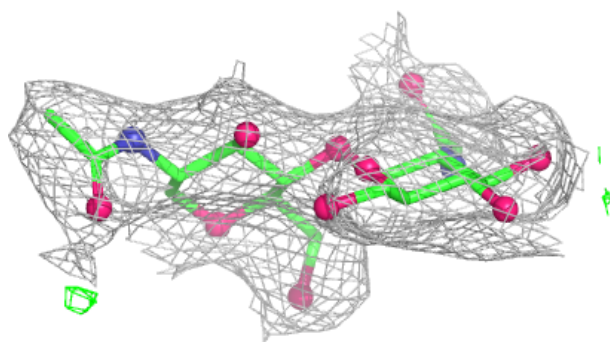
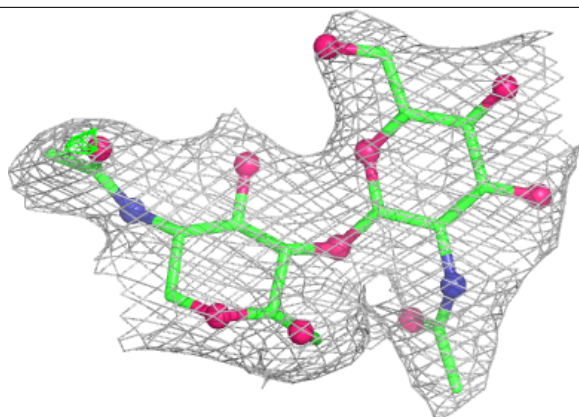
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	G	2	14/15	0.79	0.16	53,61,69,71	0
2	NAG	F	2	14/15	0.83	0.14	41,66,70,73	0
2	NAG	H	2	14/15	0.83	0.14	61,66,71,73	0
2	NAG	E	2	14/15	0.84	0.14	58,67,73,74	0
2	NAG	H	1	14/15	0.90	0.10	35,43,50,57	0
2	NAG	F	1	14/15	0.91	0.10	34,45,54,59	0
2	NAG	E	1	14/15	0.91	0.11	27,41,48,50	0
2	NAG	G	1	14/15	0.91	0.10	27,37,49,54	0

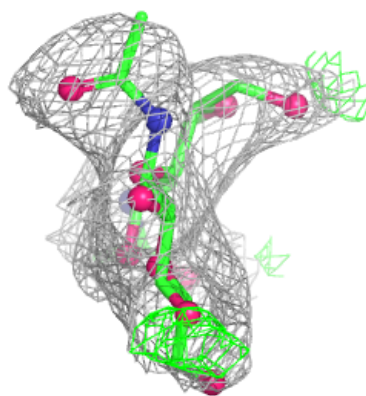
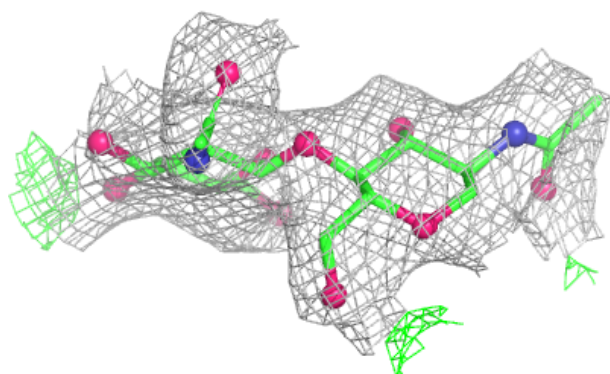
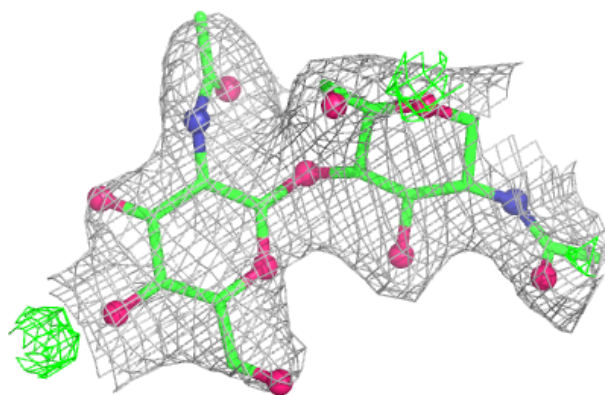
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

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)

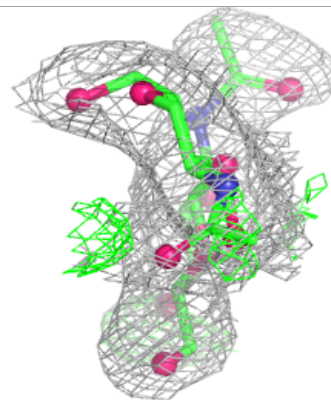
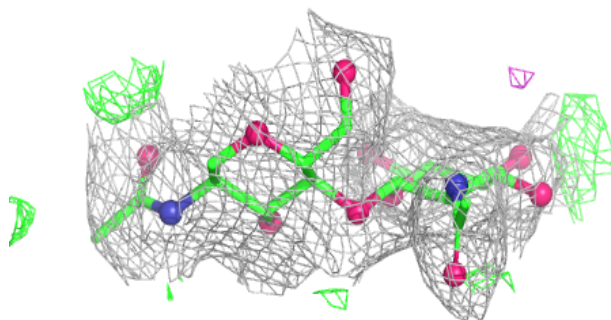
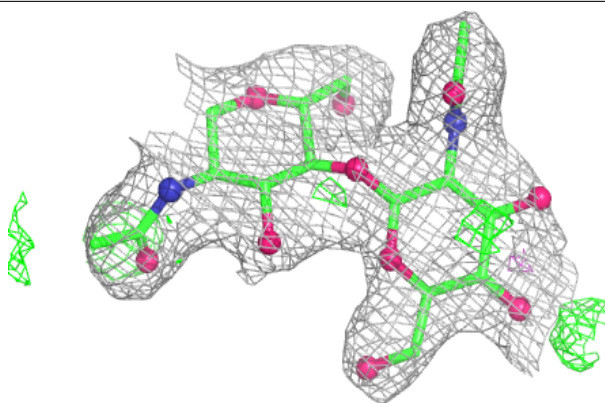
**Electron density around Chain F:**

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

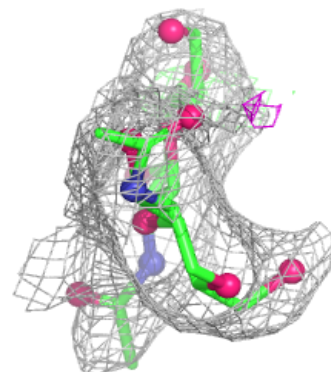
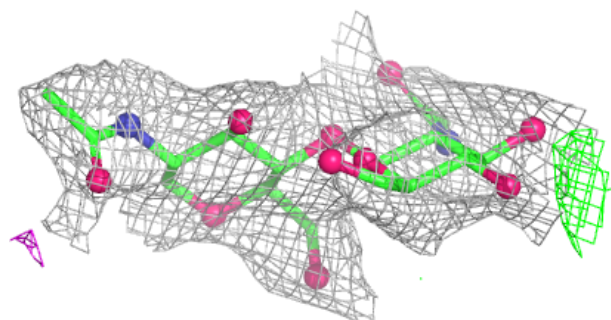
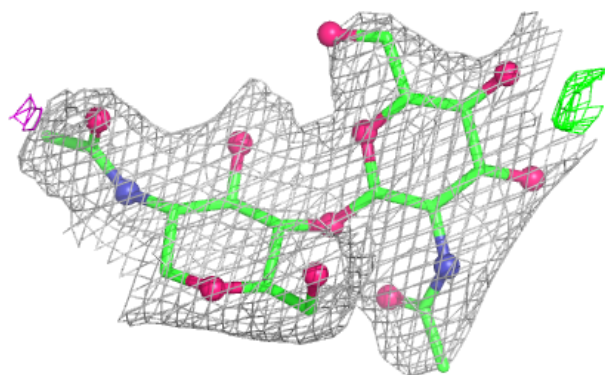


Electron density around Chain G:

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

**Electron density around Chain H:**

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



6.4 Ligands ⓘ

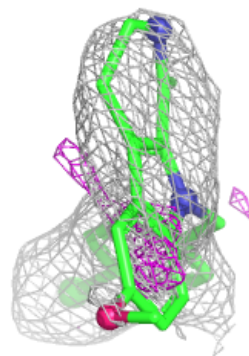
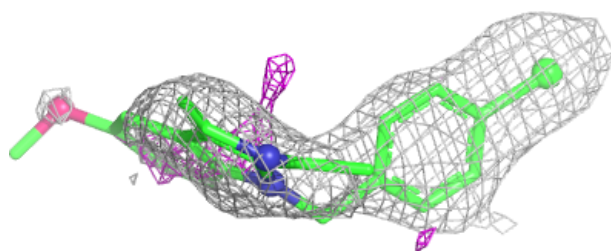
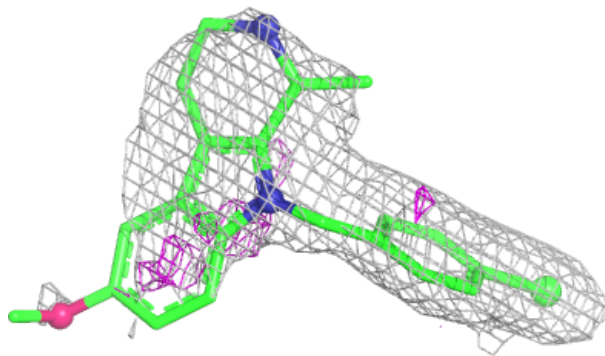
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	NAG	C	703	14/15	0.63	0.18	56,73,80,82	0
4	NAG	B	703	14/15	0.69	0.17	63,73,78,80	0
4	NAG	A	702	14/15	0.77	0.17	49,67,76,76	0
4	NAG	D	703	14/15	0.80	0.16	55,73,79,83	0
4	NAG	A	705	14/15	0.81	0.14	50,66,71,80	0
4	NAG	D	706	14/15	0.83	0.13	50,60,71,75	0
4	NAG	C	706	14/15	0.86	0.12	50,63,67,71	0
6	Y8A	B	701	24/24	0.86	0.21	65,72,80,93	0
6	Y8A	C	701	24/24	0.87	0.22	67,78,92,127	0
4	NAG	B	706	14/15	0.88	0.12	58,68,72,77	0
6	Y8A	D	701	24/24	0.88	0.20	72,81,88,113	0
6	Y8A	A	707	24/24	0.89	0.18	67,73,81,97	0
5	BOG	A	706	20/20	0.92	0.13	39,44,50,52	0
5	BOG	D	707	20/20	0.93	0.11	34,44,55,58	0
3	HEM	B	702	43/43	0.94	0.12	34,45,74,87	0
3	HEM	D	702	43/43	0.94	0.12	24,44,68,79	0
3	HEM	A	701	43/43	0.95	0.17	36,58,76,84	0
3	HEM	C	702	43/43	0.96	0.10	29,42,64,82	0

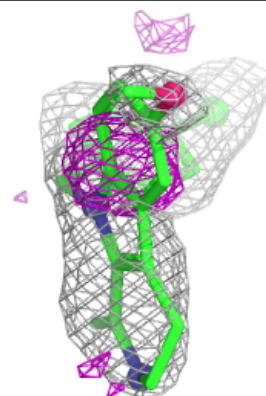
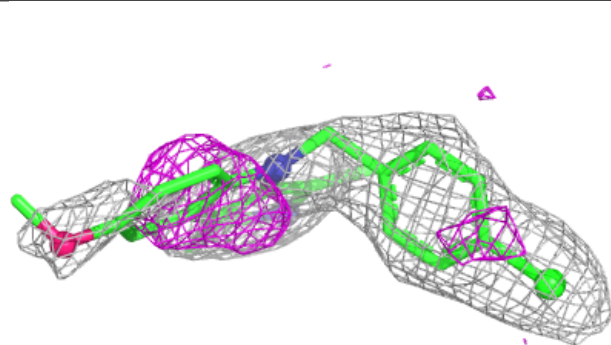
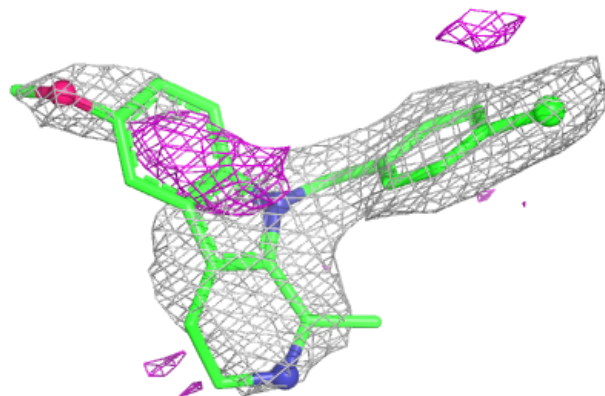
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around Y8A 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)

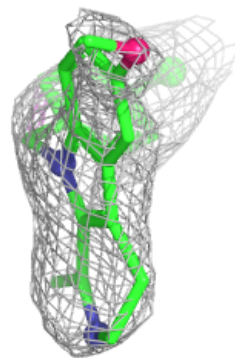
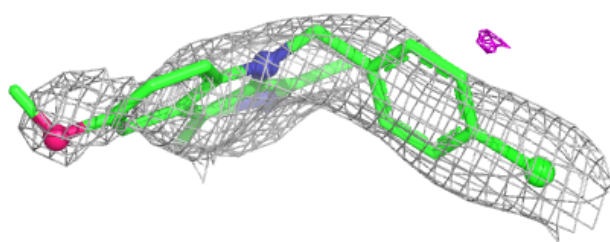
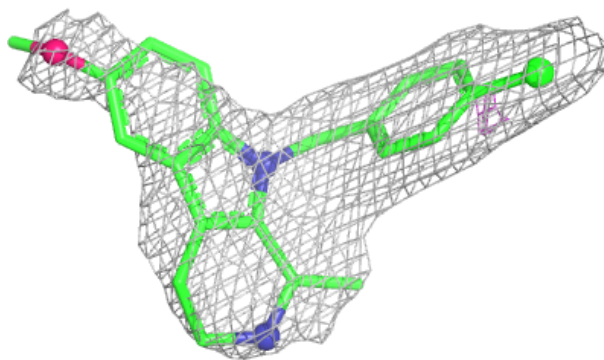
**Electron density around Y8A C 701:**

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

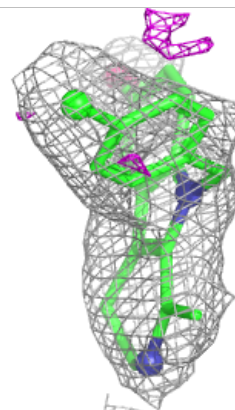
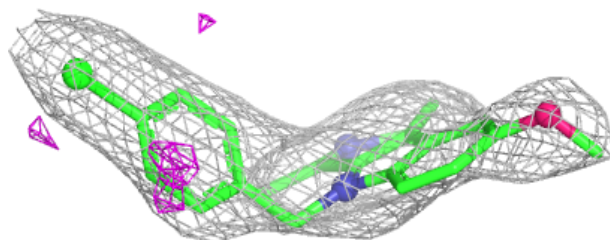
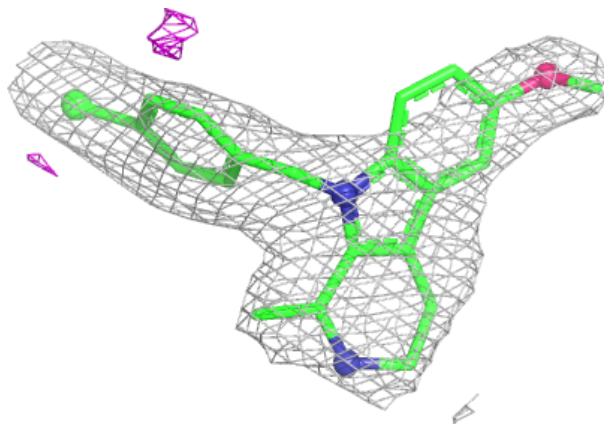


Electron density around Y8A D 701:

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

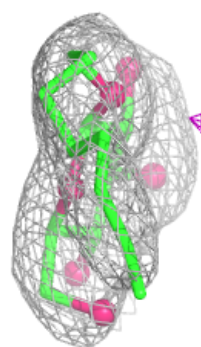
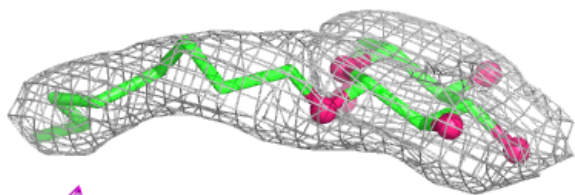
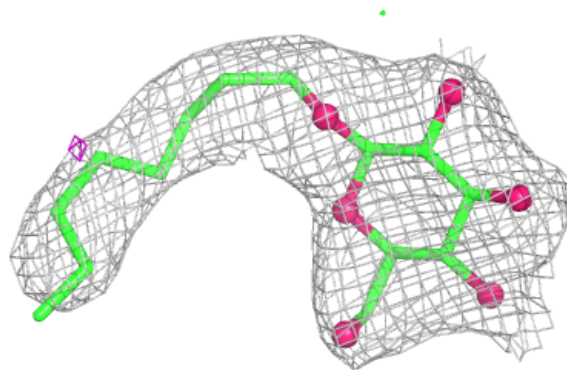
**Electron density around Y8A A 707:**

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

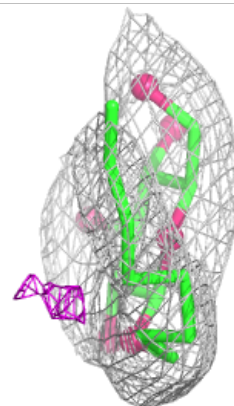
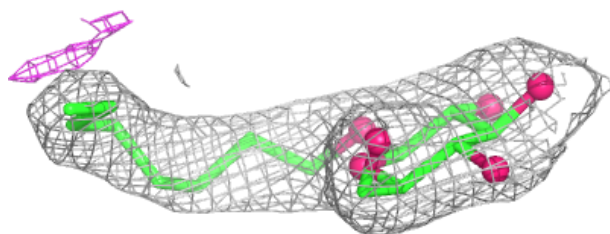
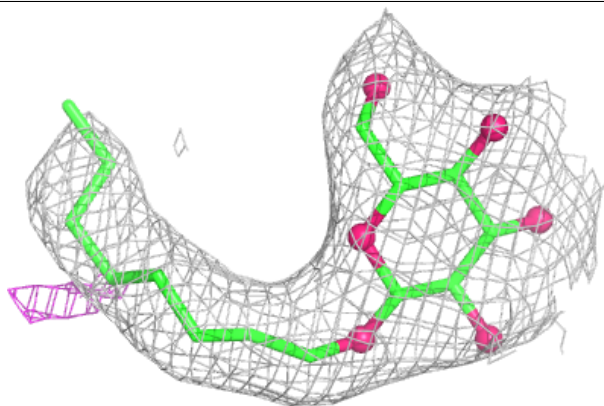


Electron density around BOG A 706:

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

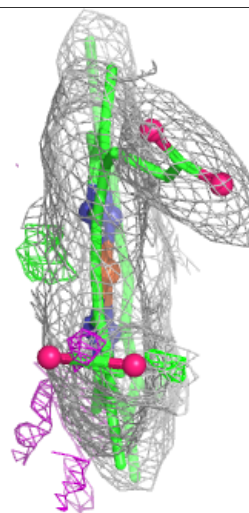
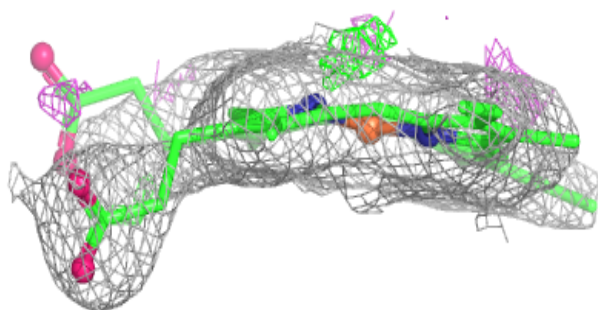
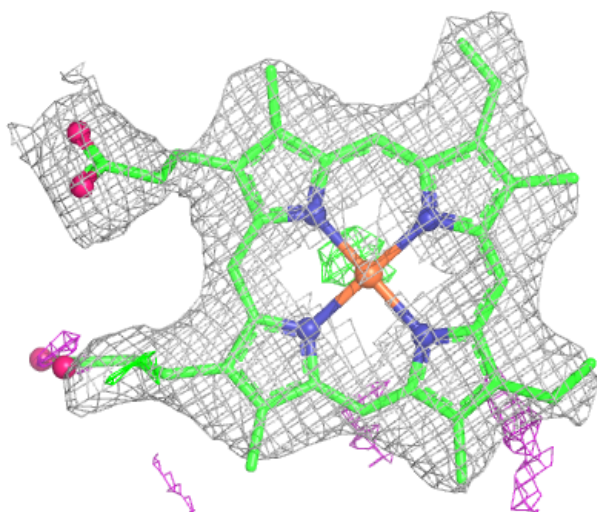
**Electron density around BOG D 707:**

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



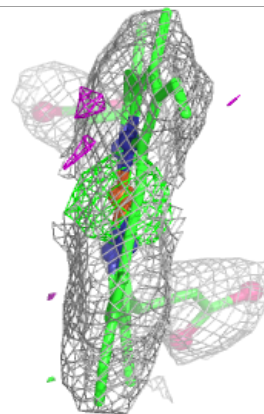
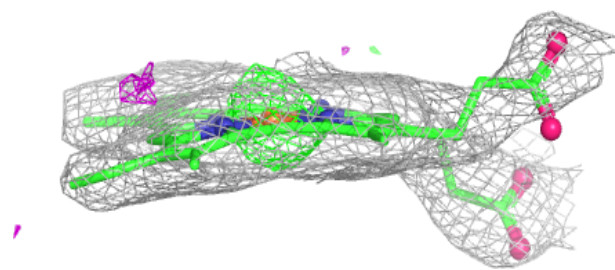
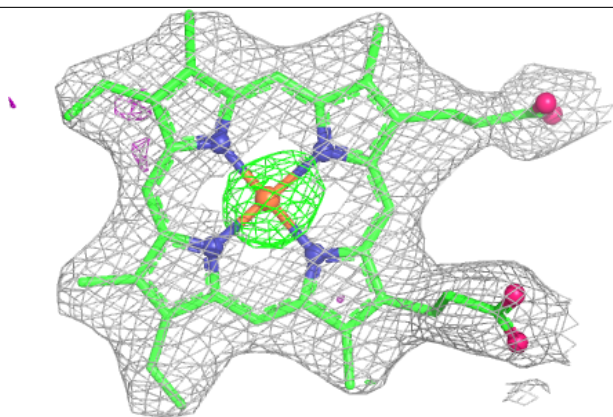
Electron density around HEM B 702:

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



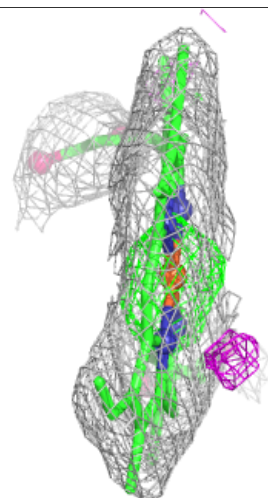
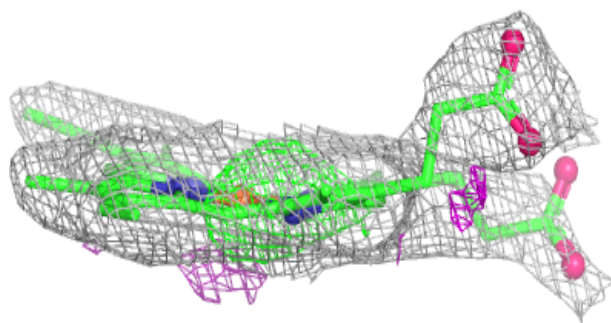
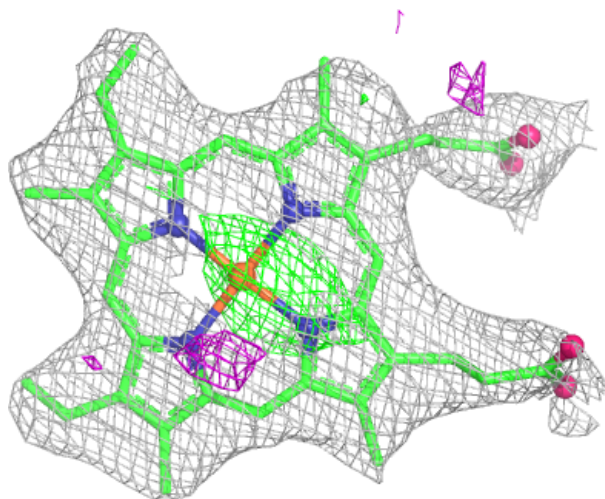
Electron density around HEM D 702:

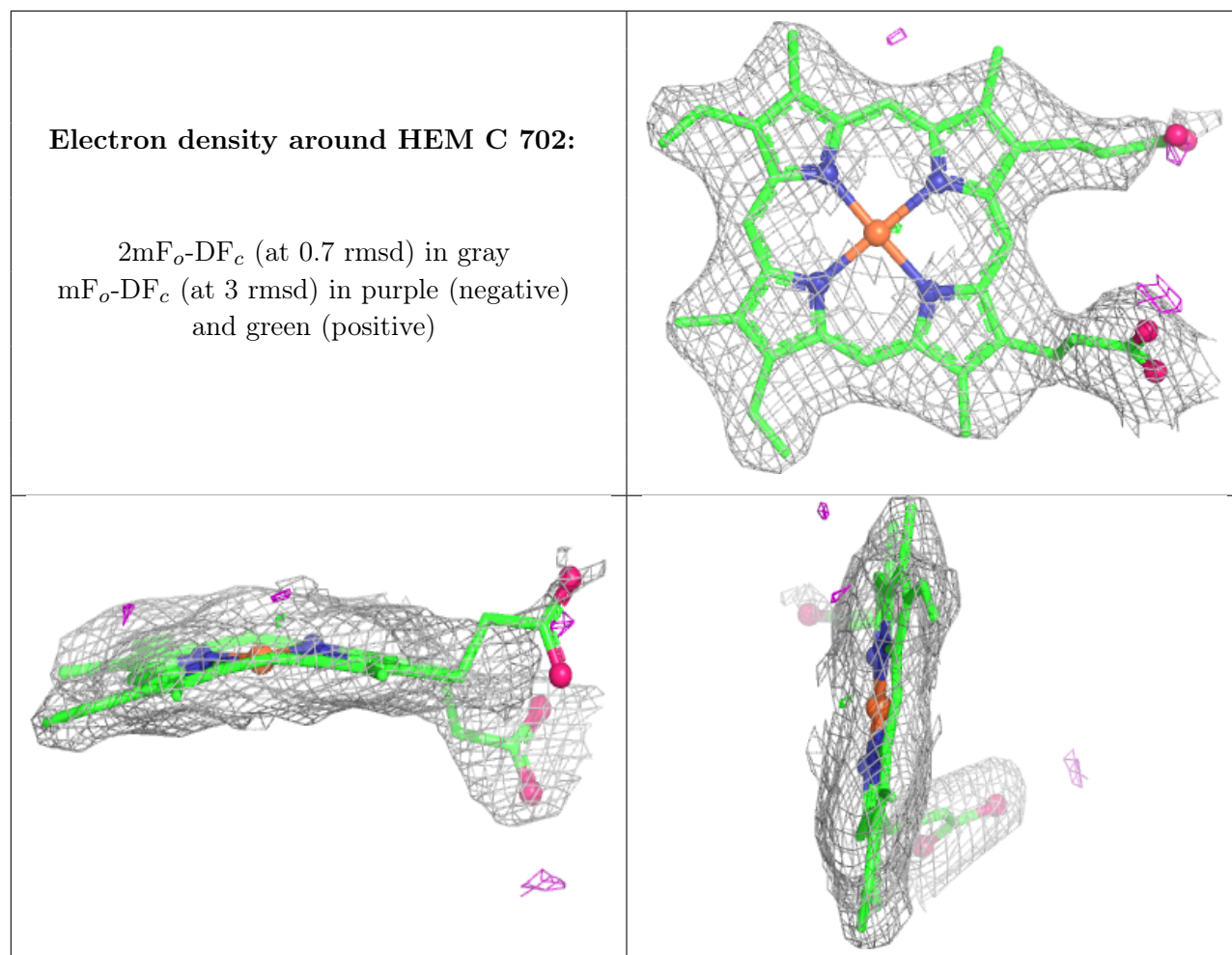
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM A 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 ⓘ

There are no such residues in this entry.