



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 03:55 PM UTC

PDB ID : 6BL4 / pdb_00006bl4
Title : Crystal Complex of Cyclooxygenase-2 with indomethacin-ethylenediamine-dansyl conjugate
Authors : Xu, S.; Uddin, M.J.; Banerjee, S.; Marnett, L.J.
Deposited on : 2017-11-09
Resolution : 2.22 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

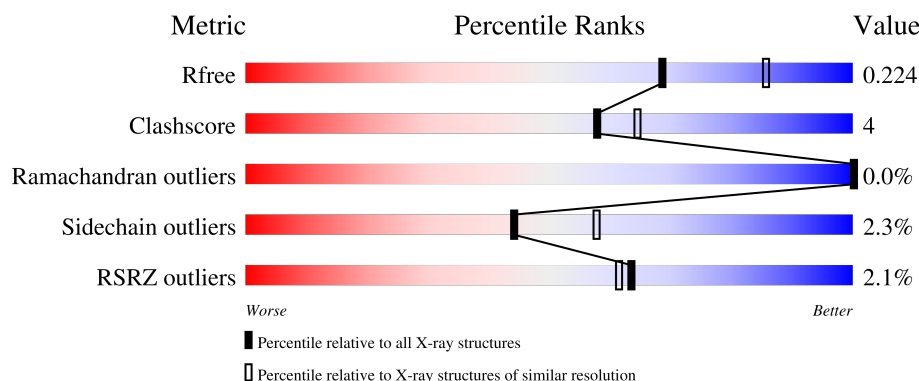
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.22 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	7682 (2.24-2.20)
Clashscore	190562	8402 (2.24-2.20)
Ramachandran outliers	187476	8303 (2.24-2.20)
Sidechain outliers	187428	8304 (2.24-2.20)
RSRZ outliers	180081	7683 (2.24-2.20)

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% 84% 9% 6%
1	B	587	3% 86% 8% 6%
1	C	587	2% 84% 9% 6%
1	D	587	2% 83% 10% 6%
2	E	2	50% 50%

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

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 20002 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.

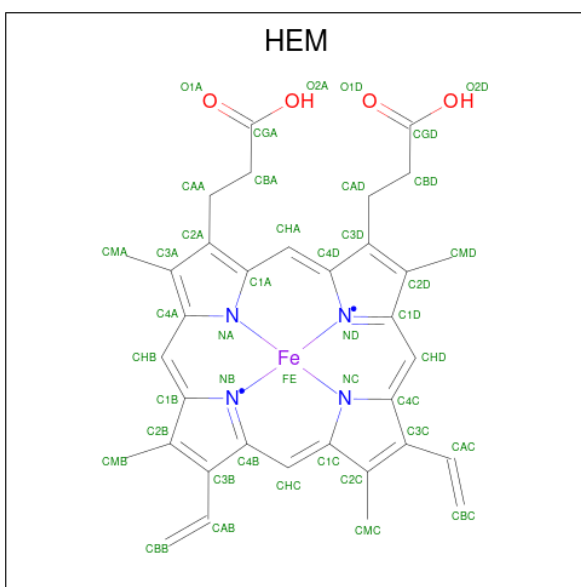
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	552	Total	C	N	O	S	0	0	0
			4474	2885	750	814	25			
1	B	552	Total	C	N	O	S	0	0	0
			4474	2885	750	814	25			
1	C	552	Total	C	N	O	S	0	0	0
			4474	2885	750	814	25			
1	D	552	Total	C	N	O	S	0	0	0
			4474	2885	750	814	25			

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



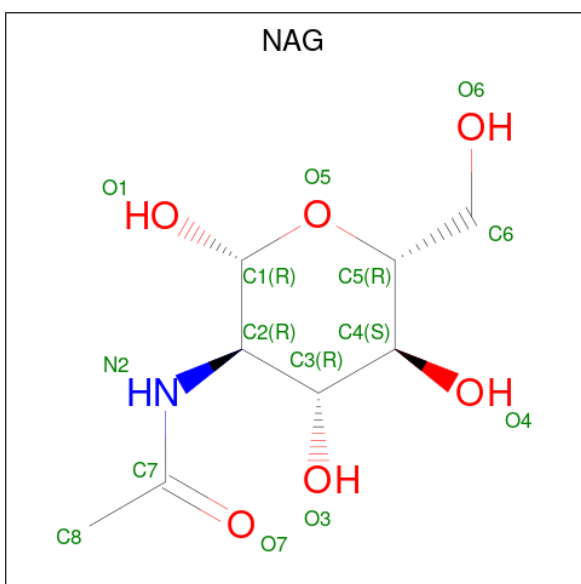
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	H	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



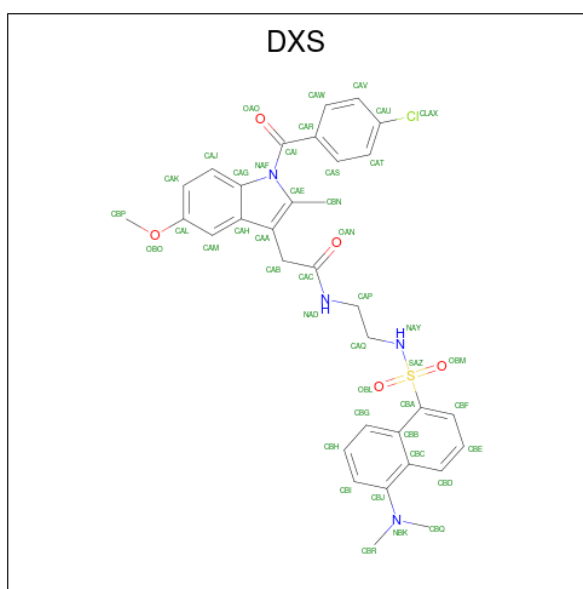
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- 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 14	C 8	N 1	O 5	0	0
4	A	1	Total 14	C 8	N 1	O 5	0	0
4	B	1	Total 14	C 8	N 1	O 5	0	0
4	B	1	Total 14	C 8	N 1	O 5	0	0
4	C	1	Total 14	C 8	N 1	O 5	0	0
4	C	1	Total 14	C 8	N 1	O 5	0	0
4	D	1	Total 14	C 8	N 1	O 5	0	0
4	D	1	Total 14	C 8	N 1	O 5	0	0

- Molecule 5 is 2-[1-(4-chlorobenzene-1-carbonyl)-5-methoxy-2-methyl-1H-indol-3-yl]-N-[2-({[5-(dimethylamino)naphthalen-1-yl]sulfonyl}amino)ethyl]acetamide (CCD ID: DXS) (formula: C₃₃H₃₃ClN₄O₅S).



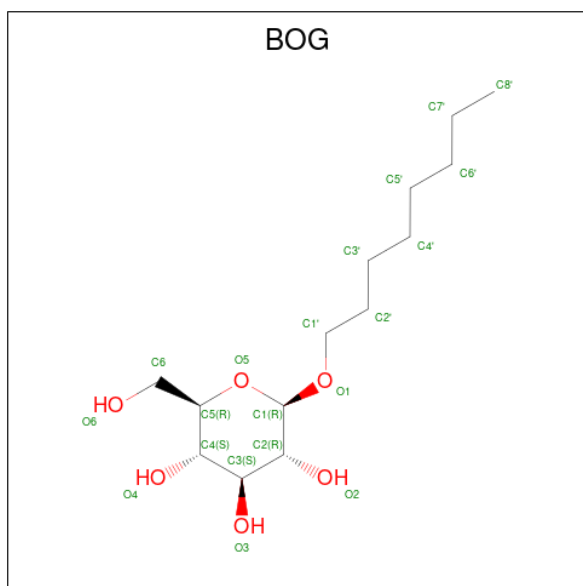
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
5	A	1	Total 44	C 33	Cl 1	N 4	O 5	S 1	0	0
5	B	1	Total 44	C 33	Cl 1	N 4	O 5	S 1	0	0
5	C	1	Total 44	C 33	Cl 1	N 4	O 5	S 1	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	D	1	Total	C	Cl	N	O	S	
			44	33	1	4	5	1	0

- Molecule 6 is octyl beta-D-glucopyranoside (CCD ID: BOG) (formula: $C_{14}H_{28}O_6$).



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

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	395	Total	O		
			395	395	0	0
7	B	324	Total	O		
			324	324	0	0
7	C	373	Total	O		
			373	373	0	0

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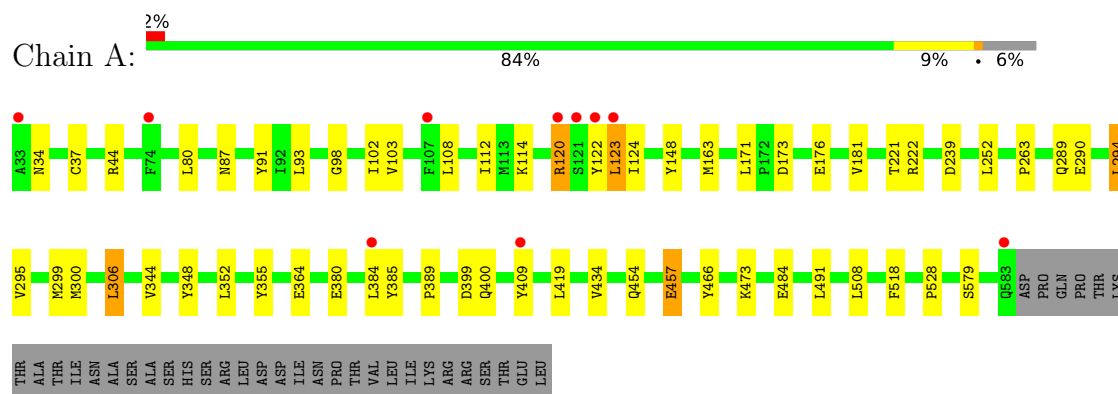
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	D	342	Total	O	0	0
			342	342		

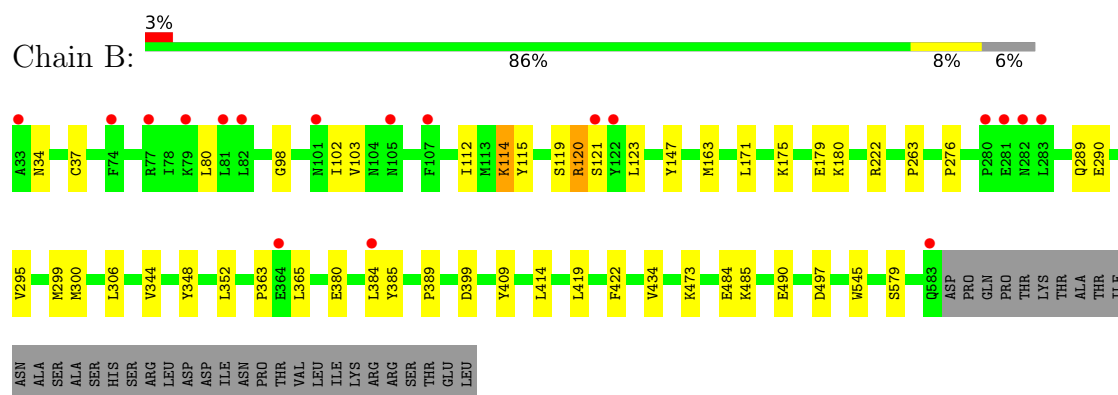
3 Residue-property plots

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

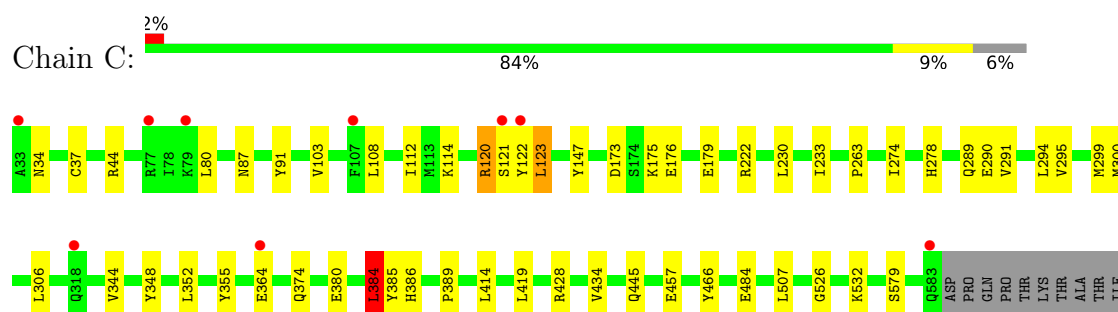
• Molecule 1: Prostaglandin G/H synthase 2



• Molecule 1: Prostaglandin G/H synthase 2



• Molecule 1: Prostaglandin G/H synthase 2



ASN
ALA
SER
ALA
SER
HIS
SER
ARG
LEU
ASP
ASP
ILE
ASN
PRO
THR
VAL
LEU
ILE
LYS
ARG
SER
THR
GLU
LEU

- Molecule 1: Prostaglandin G/H synthase 2

Chain D:  2% 83% 10% 6%

A33 N34 C37 R44 F74 L80 T88 L93 G98 I102 V103 F107 I112 M113 K114 Y115 V116 S119 R120 S121 Y122 L123 I124 M163 L171 F172 D173 E176 K180 F209 R222 L230 R231 H232 L252 P263 Q289 E290

L294 V295 M299 M300 L306 H320 V344 Y348 L352 Y355 Q374 I377 E380 L384 Y385 P389 N396 E401 L414 L419 V434 Y466 K473 E484 E502 P528 K532 P542 S579 V582 Q583 ASP PRO GLN

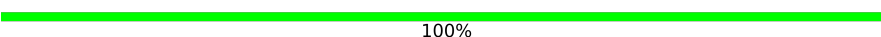
PRO
THR
LYS
THR
ALA
THR
ILE
ASN
ALA
SER
ALA
SER
HIS
SER
ARG
LEU
ASP
ASP
ILE
ASN
PRO
THR
VAL
LEU
ILE
LYS
ARG
SER
THR
GLU
LEU

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

Chain E:  50% 50%

NAG1
NAG2

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

Chain F:  100%

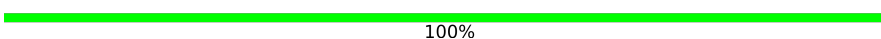
NAG1
NAG2

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

Chain G:  50% 50%

NAG1
NAG2

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

Chain H:  100%

NAG1
NAG2

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	214.79Å 120.41Å 134.42Å 90.00° 123.60° 90.00°	Depositor
Resolution (Å)	111.96 – 2.22 111.96 – 2.22	Depositor EDS
% Data completeness (in resolution range)	98.5 (111.96-2.22) 90.0 (111.96-2.22)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.43 (at 2.22Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.189 , 0.220 0.196 , 0.224	Depositor DCC
R_{free} test set	4150 reflections (3.00%)	wwPDB-VP
Wilson B-factor (Å ²)	24.5	Xtriage
Anisotropy	0.449	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.014 for $1/2^*h-3/2^*k,-1/2^*h-1/2^*k,-1/2^*h+1/2^*k-l$ 0.018 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	0.94	EDS
Total number of atoms	20002	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.59% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: BOG, DXS, NAG, HEM

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	0.30	1/4601 (0.0%)	0.47	0/6239
1	B	0.31	1/4601 (0.0%)	0.48	0/6239
1	C	0.32	1/4601 (0.0%)	0.49	0/6239
1	D	0.34	2/4601 (0.0%)	0.46	0/6239
All	All	0.32	5/18404 (0.0%)	0.48	0/24956

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	384	LEU	C-O	-7.03	1.15	1.24
1	D	121	SER	C-O	-5.91	1.17	1.24
1	B	497	ASP	C-O	-5.84	1.17	1.24
1	A	508	LEU	C-O	-5.52	1.16	1.24
1	D	320	HIS	C-O	-5.34	1.18	1.24

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4474	0	4373	39	0
1	B	4474	0	4373	35	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	4474	0	4373	36	0
1	D	4474	0	4373	39	0
2	E	28	0	25	0	0
2	F	28	0	25	0	0
2	G	28	0	25	0	0
2	H	28	0	25	0	0
3	A	43	0	30	2	0
3	B	43	0	30	2	0
3	C	43	0	30	5	0
3	D	43	0	30	1	0
4	A	28	0	26	0	0
4	B	28	0	26	0	0
4	C	28	0	26	0	0
4	D	28	0	26	0	0
5	A	44	0	0	2	0
5	B	44	0	0	3	0
5	C	44	0	0	3	0
5	D	44	0	0	3	0
6	A	40	0	55	2	0
6	C	20	0	28	3	0
6	D	40	0	56	1	0
7	A	395	0	0	4	0
7	B	324	0	0	3	0
7	C	373	0	0	8	0
7	D	342	0	0	3	0
All	All	20002	0	17955	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)
1:A:120:ARG:HB3	1:A:120:ARG:HH11	1.07	1.12
1:D:120:ARG:HH11	1:D:120:ARG:HB3	1.21	1.04
1:C:120:ARG:O	1:C:122:TYR:N	1.94	1.01
1:D:120:ARG:HB3	1:D:120:ARG:NH1	1.78	0.98
1:B:120:ARG:HH11	1:B:120:ARG:HB3	1.30	0.96
1:D:116:VAL:O	1:D:120:ARG:HD3	1.63	0.96
1:A:120:ARG:HB3	1:A:120:ARG:NH1	1.80	0.95
1:C:263:PRO:HG2	1:C:299:MET:HE1	1.60	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:120:ARG:HH11	1:B:120:ARG:CB	1.92	0.82
1:B:263:PRO:HG2	1:B:299:MET:HE1	1.60	0.81
1:C:445:GLN:OE1	7:C:801:HOH:O	2.02	0.77
1:B:120:ARG:HB3	1:B:120:ARG:NH1	2.00	0.76
3:C:701:HEM:HHC	3:C:701:HEM:HBB2	1.67	0.76
3:B:701:HEM:HHC	3:B:701:HEM:HBB2	1.68	0.75
1:C:384:LEU:HD22	1:C:507:LEU:HD11	1.68	0.74
1:B:180:LYS:HB3	1:B:490:GLU:HG2	1.70	0.73
1:D:263:PRO:HG2	1:D:299:MET:HE1	1.72	0.71
1:A:120:ARG:HD2	1:A:120:ARG:H	1.55	0.71
1:D:119:SER:OG	1:D:120:ARG:HD2	1.90	0.70
1:B:120:ARG:HH11	1:B:120:ARG:CG	2.05	0.70
3:C:701:HEM:HBC2	3:C:701:HEM:HHD	1.75	0.69
1:C:278:HIS:NE2	7:C:809:HOH:O	2.30	0.63
6:C:707:BOG:H8'2	6:C:707:BOG:H4'1	1.82	0.61
1:C:384:LEU:CD2	1:C:507:LEU:HD11	2.31	0.60
1:C:222:ARG:HH12	1:C:290:GLU:CD	2.09	0.60
1:D:222:ARG:HH12	1:D:290:GLU:CD	2.10	0.60
3:A:701:HEM:HBB2	3:A:701:HEM:HHC	1.84	0.59
1:B:222:ARG:HH12	1:B:290:GLU:CD	2.10	0.59
1:A:222:ARG:NE	7:A:803:HOH:O	2.24	0.59
3:D:701:HEM:HBB2	3:D:701:HEM:HHC	1.85	0.59
7:C:928:HOH:O	1:D:542:PRO:HD2	2.01	0.59
1:C:384:LEU:HD22	1:C:507:LEU:CD1	2.33	0.58
1:B:120:ARG:O	1:B:121:SER:C	2.45	0.58
1:A:120:ARG:HD2	1:A:120:ARG:N	2.18	0.57
3:C:701:HEM:HBB2	3:C:701:HEM:CHC	2.29	0.57
1:D:44:ARG:HD3	7:D:1084:HOH:O	2.03	0.57
1:A:222:ARG:HH12	1:A:290:GLU:CD	2.13	0.57
1:B:180:LYS:HD3	1:B:490:GLU:HG3	1.88	0.56
5:C:706:DXS:CBR	5:C:706:DXS:CBD	2.84	0.55
1:A:454:GLN:HA	1:A:457:GLU:HB2	1.89	0.55
1:A:364:GLU:HG3	7:A:988:HOH:O	2.07	0.55
1:A:122:TYR:CE1	1:A:123:LEU:HD13	2.42	0.54
1:A:263:PRO:HG2	1:A:299:MET:HE1	1.88	0.54
1:C:300:MET:HG3	1:C:419:LEU:HD22	1.89	0.54
1:A:163:MET:HE2	1:A:171:LEU:HD11	1.89	0.54
5:A:706:DXS:CBR	5:A:706:DXS:CBD	2.85	0.54
1:A:173:ASP:HB3	1:A:176:GLU:HB2	1.90	0.53
1:D:173:ASP:HB3	1:D:176:GLU:HB2	1.90	0.53
3:C:701:HEM:HHC	3:C:701:HEM:CBB	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:355:TYR:OH	5:C:706:DXS:NAD	2.42	0.53
1:D:115:TYR:O	1:D:119:SER:HB3	2.09	0.52
1:A:344:VAL:HA	1:A:348:TYR:HB3	1.90	0.52
1:B:34:ASN:HB3	1:B:37:CYS:SG	2.49	0.52
1:B:175:LYS:O	1:B:179:GLU:HG3	2.08	0.52
1:B:485:LYS:HG3	7:B:983:HOH:O	2.08	0.52
5:D:706:DXS:CBD	5:D:706:DXS:CBQ	2.85	0.52
1:A:120:ARG:H	1:A:120:ARG:CD	2.23	0.52
1:B:399:ASP:OD1	1:B:399:ASP:N	2.40	0.51
1:A:181:VAL:HG21	1:A:491:LEU:HD21	1.93	0.51
1:C:230:LEU:HG	1:C:233:ILE:HD12	1.93	0.50
1:D:384:LEU:C	1:D:384:LEU:HD12	2.36	0.50
1:D:300:MET:HG3	1:D:419:LEU:HD22	1.94	0.50
1:C:34:ASN:HB3	1:C:37:CYS:SG	2.51	0.50
1:B:300:MET:HG3	1:B:419:LEU:HD22	1.92	0.50
1:C:91:TYR:CD2	6:C:707:BOG:H5'2	2.48	0.49
1:B:222:ARG:NH1	1:B:290:GLU:OE2	2.44	0.49
1:C:122:TYR:CE1	1:C:123:LEU:HD13	2.48	0.49
1:C:428:ARG:HG3	7:C:831:HOH:O	2.12	0.49
1:B:115:TYR:O	1:B:119:SER:HB3	2.12	0.49
1:B:380:GLU:O	1:B:384:LEU:HG	2.12	0.49
1:D:163:MET:HE2	1:D:171:LEU:HD11	1.95	0.49
1:A:34:ASN:HB3	1:A:37:CYS:SG	2.53	0.49
1:D:344:VAL:HA	1:D:348:TYR:HB3	1.94	0.48
1:C:87:ASN:HD22	6:C:707:BOG:H61	1.78	0.48
1:C:380:GLU:HG2	1:C:466:TYR:CE1	2.49	0.48
1:D:180:LYS:HE3	7:D:980:HOH:O	2.14	0.48
1:C:344:VAL:HA	1:C:348:TYR:HB3	1.96	0.48
1:C:120:ARG:C	1:C:122:TYR:N	2.67	0.48
1:D:124:ILE:HD11	1:D:528:PRO:HB2	1.96	0.48
5:A:706:DXS:CAJ	5:A:706:DXS:CAW	2.91	0.48
1:A:44:ARG:HG3	7:A:1110:HOH:O	2.13	0.47
1:C:389:PRO:HB2	1:C:434:VAL:HA	1.95	0.47
1:B:414:LEU:HD11	1:B:419:LEU:HD23	1.95	0.47
1:D:582:VAL:O	1:D:583:GLN:HG3	2.14	0.47
1:B:344:VAL:HA	1:B:348:TYR:HB3	1.96	0.47
1:A:91:TYR:CD2	6:A:707:BOG:H4'2	2.50	0.47
1:C:526:GLY:HA3	5:C:706:DXS:CAT	2.44	0.47
1:D:103:VAL:HG11	1:D:112:ILE:HD12	1.97	0.47
1:D:389:PRO:HB2	1:D:434:VAL:HA	1.96	0.46
1:A:252:LEU:HB2	1:A:306:LEU:HD13	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:252:LEU:HB2	1:D:306:LEU:HD13	1.97	0.46
1:D:414:LEU:HD11	1:D:419:LEU:HD23	1.96	0.46
1:C:386:HIS:CE1	3:C:701:HEM:HAD2	2.51	0.46
1:D:98:GLY:O	1:D:102:ILE:HG12	2.15	0.46
1:A:380:GLU:O	1:A:384:LEU:HG	2.15	0.46
1:C:175:LYS:O	1:C:179:GLU:HG3	2.15	0.46
1:D:473:LYS:NZ	7:D:803:HOH:O	2.29	0.45
1:D:380:GLU:O	1:D:384:LEU:HG	2.17	0.45
1:A:87:ASN:HB2	6:A:707:BOG:H62	1.98	0.45
1:C:44:ARG:HG3	7:C:1121:HOH:O	2.14	0.45
1:B:120:ARG:NH1	1:B:120:ARG:CG	2.72	0.45
1:D:473:LYS:NZ	1:D:473:LYS:HB2	2.31	0.45
1:D:380:GLU:HG2	1:D:466:TYR:CE1	2.52	0.45
1:A:380:GLU:HG2	1:A:466:TYR:CE1	2.51	0.45
1:C:414:LEU:HD11	1:C:419:LEU:HD23	1.99	0.45
1:D:222:ARG:NH1	1:D:290:GLU:OE2	2.46	0.45
1:A:389:PRO:HB2	1:A:434:VAL:HA	1.98	0.44
1:B:473:LYS:HE2	7:B:1095:HOH:O	2.17	0.44
1:C:147:TYR:HB2	7:C:910:HOH:O	2.18	0.44
1:A:103:VAL:HG11	1:A:112:ILE:HD12	1.99	0.44
1:A:222:ARG:NH1	1:A:290:GLU:OE2	2.46	0.44
3:A:701:HEM:HHD	3:A:701:HEM:HBC2	1.99	0.44
1:B:114:LYS:HE3	1:B:365:LEU:O	2.18	0.44
1:B:363:PRO:HG2	1:B:545:TRP:CD2	2.53	0.44
1:A:108:LEU:HD12	1:A:108:LEU:HA	1.82	0.43
1:D:209:PHE:HB2	1:D:377:ILE:HG13	2.01	0.43
1:A:98:GLY:O	1:A:102:ILE:HG12	2.18	0.43
1:C:222:ARG:NH1	1:C:290:GLU:OE2	2.50	0.43
1:D:374:GLN:O	1:D:532:LYS:HE3	2.19	0.43
3:B:701:HEM:HBB2	3:B:701:HEM:CHC	2.40	0.43
1:C:274:ILE:HD12	1:C:291:VAL:HG12	1.99	0.43
1:C:364:GLU:HG3	7:C:824:HOH:O	2.19	0.43
1:A:473:LYS:HB2	1:A:473:LYS:NZ	2.34	0.43
1:D:352:LEU:HD23	1:D:352:LEU:HA	1.91	0.43
1:D:88:THR:HG23	6:D:707:BOG:H4'1	2.00	0.42
1:A:300:MET:HG3	1:A:419:LEU:HD22	2.01	0.42
1:D:294:LEU:HG	1:D:295:VAL:HG22	2.00	0.42
1:B:103:VAL:HG11	1:B:112:ILE:HD12	2.01	0.42
1:A:44:ARG:NH1	1:A:122:TYR:O	2.52	0.42
1:A:93:LEU:HD13	1:A:355:TYR:CZ	2.54	0.42
1:A:399:ASP:OD1	1:A:399:ASP:N	2.46	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:98:GLY:O	1:B:102:ILE:HG12	2.19	0.42
1:B:120:ARG:HH11	1:B:120:ARG:HG2	1.79	0.42
1:B:120:ARG:NH2	5:B:706:DXS:CAP	2.83	0.42
1:B:276:PRO:HD3	1:B:409:TYR:CD1	2.55	0.42
1:A:124:ILE:HD11	1:A:528:PRO:HB2	2.01	0.42
1:B:163:MET:HE2	1:B:171:LEU:HD11	2.02	0.42
1:C:103:VAL:HG11	1:C:112:ILE:HD12	2.01	0.42
1:D:230:LEU:HA	1:D:232:HIS:CE1	2.55	0.42
1:B:414:LEU:HA	1:B:422:PHE:CE1	2.55	0.41
1:D:384:LEU:CD1	5:D:706:DXS:CLAX	3.05	0.41
1:B:473:LYS:HB2	1:B:473:LYS:NZ	2.35	0.41
1:B:147:TYR:HB2	7:B:876:HOH:O	2.19	0.41
1:C:123:LEU:HD12	1:C:123:LEU:HA	1.91	0.41
1:A:294:LEU:HA	1:A:409:TYR:CE1	2.54	0.41
1:C:173:ASP:HB3	1:C:176:GLU:HB2	2.02	0.41
1:A:400:GLN:O	7:A:801:HOH:O	2.22	0.41
1:C:108:LEU:HD12	1:C:108:LEU:HA	1.88	0.41
1:C:352:LEU:HD23	1:C:352:LEU:HA	1.95	0.41
1:B:115:TYR:HE2	5:B:706:DXS:CBA	2.34	0.41
1:C:374:GLN:O	1:C:532:LYS:HE3	2.20	0.41
1:D:384:LEU:HD13	5:D:706:DXS:CLAX	2.58	0.41
1:A:148:TYR:CE1	1:A:221:THR:HB	2.56	0.41
1:A:352:LEU:HA	1:A:352:LEU:HD23	1.88	0.41
1:D:34:ASN:HB3	1:D:37:CYS:SG	2.61	0.41
1:D:93:LEU:HD13	1:D:355:TYR:CZ	2.55	0.41
1:D:163:MET:CE	1:D:502:GLU:HG2	2.50	0.41
1:D:396:ASN:OD1	1:D:401:GLU:HG2	2.21	0.41
1:A:352:LEU:HD22	1:A:518:PHE:CE2	2.56	0.40
1:B:352:LEU:HD23	1:B:352:LEU:HA	1.93	0.40
1:B:389:PRO:HB2	1:B:434:VAL:HA	2.03	0.40
5:B:706:DXS:CBG	5:B:706:DXS:NAY	2.83	0.40
1:A:294:LEU:HA	1:A:409:TYR:CD1	2.57	0.40
1:C:457:GLU:OE1	7:C:802:HOH:O	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	550/587 (94%)	536 (98%)	14 (2%)	0	100	100
1	B	550/587 (94%)	536 (98%)	14 (2%)	0	100	100
1	C	550/587 (94%)	533 (97%)	16 (3%)	1 (0%)	43	50
1	D	550/587 (94%)	536 (98%)	14 (2%)	0	100	100
All	All	2200/2348 (94%)	2141 (97%)	58 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	121	SER

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	493/525 (94%)	480 (97%)	13 (3%)	40	53
1	B	493/525 (94%)	483 (98%)	10 (2%)	48	63
1	C	493/525 (94%)	481 (98%)	12 (2%)	43	56
1	D	493/525 (94%)	483 (98%)	10 (2%)	48	63
All	All	1972/2100 (94%)	1927 (98%)	45 (2%)	44	58

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	80	LEU
1	A	114	LYS
1	A	120	ARG
1	A	123	LEU
1	A	239	ASP
1	A	289	GLN
1	A	294	LEU
1	A	295	VAL
1	A	306	LEU
1	A	385	TYR
1	A	457	GLU
1	A	484	GLU
1	A	579	SER
1	B	80	LEU
1	B	114	LYS
1	B	120	ARG
1	B	123	LEU
1	B	289	GLN
1	B	295	VAL
1	B	306	LEU
1	B	385	TYR
1	B	484	GLU
1	B	579	SER
1	C	80	LEU
1	C	114	LYS
1	C	120	ARG
1	C	123	LEU
1	C	289	GLN
1	C	294	LEU
1	C	295	VAL
1	C	306	LEU
1	C	384	LEU
1	C	385	TYR
1	C	484	GLU
1	C	579	SER
1	D	80	LEU
1	D	114	LYS
1	D	120	ARG
1	D	123	LEU
1	D	289	GLN
1	D	295	VAL
1	D	306	LEU
1	D	385	TYR

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Mol	Chain	Res	Type
1	D	484	GLU
1	D	579	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	203	GLN
1	A	327	GLN
1	B	39	ASN
1	B	42	GLN
1	B	203	GLN
1	B	282	ASN
1	B	421	GLN
1	B	571	ASN
1	C	39	ASN
1	C	327	GLN
1	C	386	HIS
1	C	461	GLN
1	C	571	ASN
1	D	203	GLN
1	D	270	GLN
1	D	327	GLN
1	D	461	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

8 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond

length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	E	1	2,1	14,14,15	0.39	0	17,19,21	0.65	1 (5%)
2	NAG	E	2	2	14,14,15	0.39	0	17,19,21	0.53	0
2	NAG	F	1	2,1	14,14,15	0.45	0	17,19,21	0.58	0
2	NAG	F	2	2	14,14,15	0.41	0	17,19,21	0.53	0
2	NAG	G	1	2,1	14,14,15	0.41	0	17,19,21	0.58	0
2	NAG	G	2	2	14,14,15	0.60	1 (7%)	17,19,21	0.61	0
2	NAG	H	1	2,1	14,14,15	0.46	0	17,19,21	0.61	0
2	NAG	H	2	2	14,14,15	0.40	0	17,19,21	0.56	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	-	0/6/23/26	0/1/1/1
2	NAG	E	2	2	-	0/6/23/26	0/1/1/1
2	NAG	F	1	2,1	-	0/6/23/26	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	-	0/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	2	NAG	C1-C2	2.09	1.55	1.52

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	1	NAG	C1-O5-C5	2.11	115.01	112.19

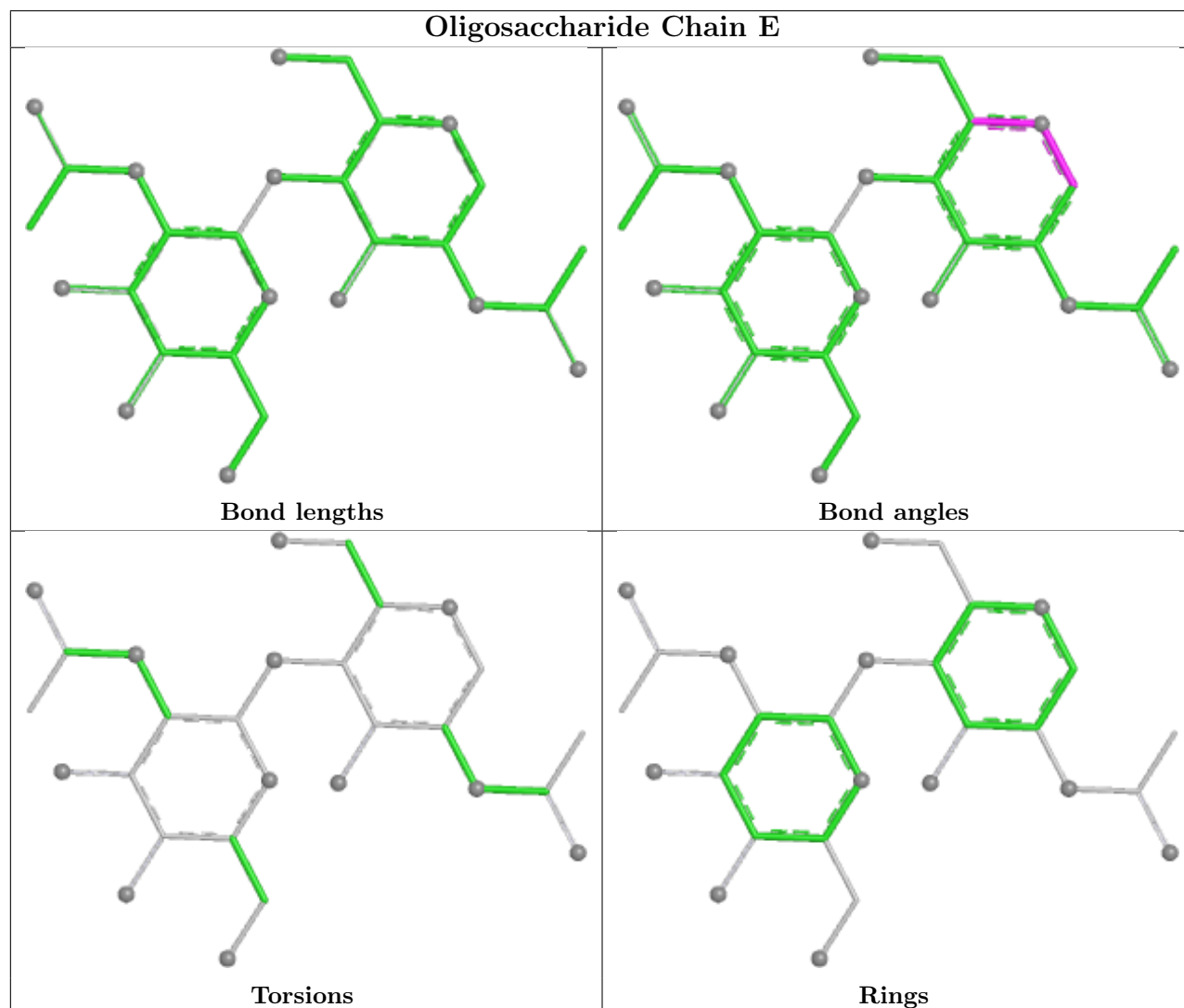
There are no chirality outliers.

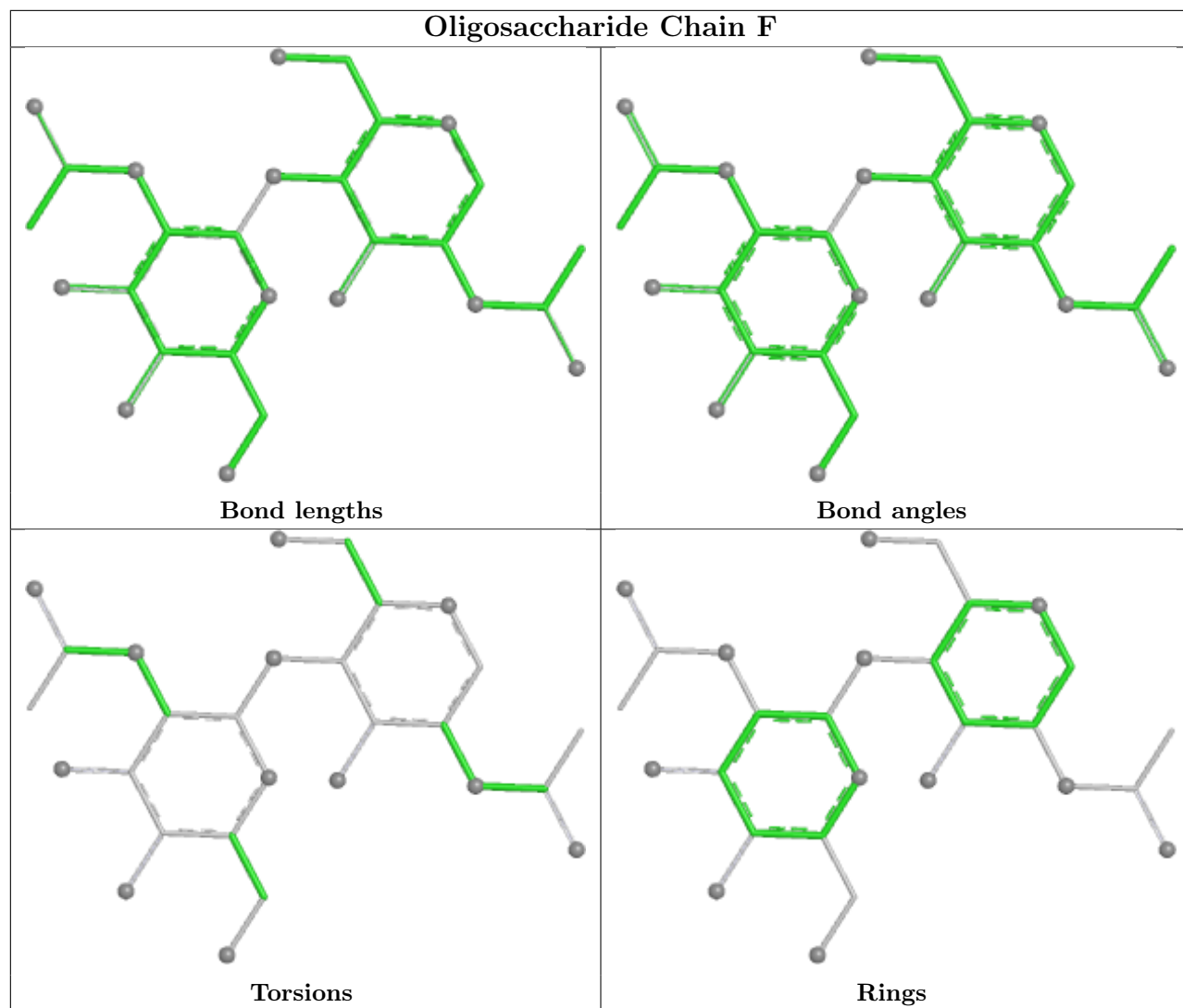
There are no torsion outliers.

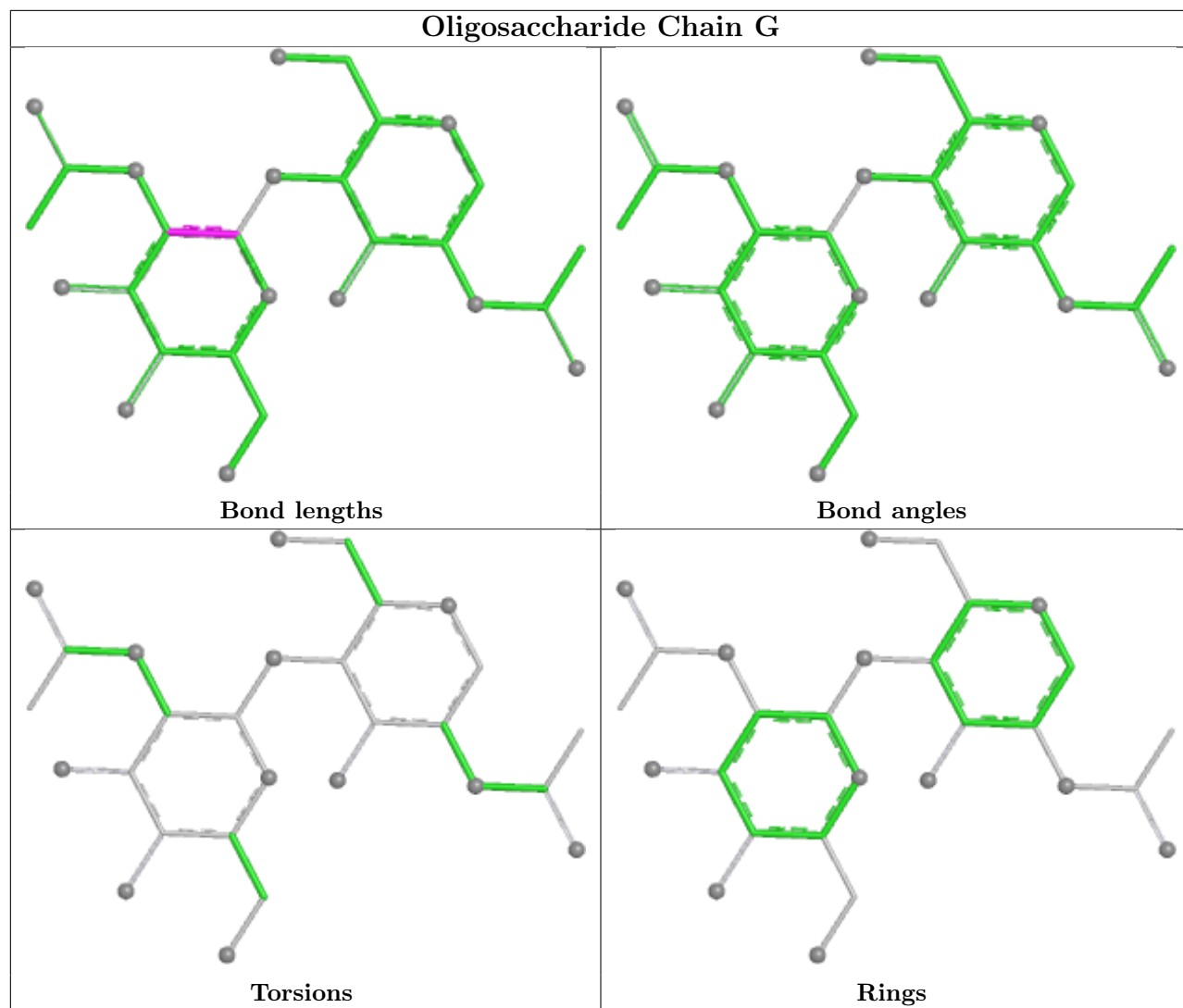
There are no ring outliers.

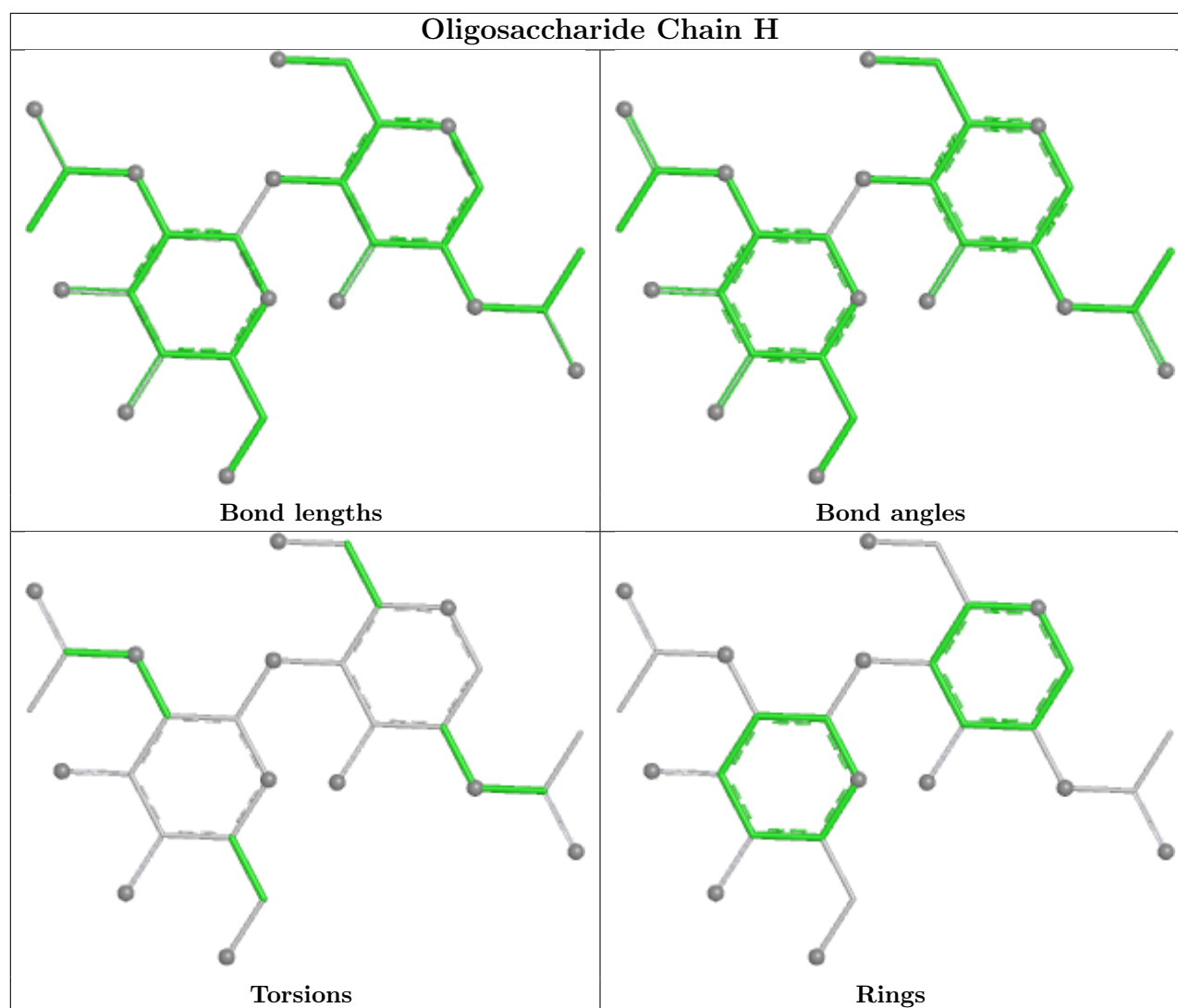
No monomer is involved in short contacts.

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









5.6 Ligand geometry [i](#)

21 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	$\# Z > 2$
3	HEM	D	701	1	50,50,50	1.46	5 (10%)	67,82,82	1.30	6 (8%)
5	DXS	C	706	-	48,48,48	4.39	20 (41%)	69,70,70	1.99	16 (23%)
6	BOG	A	707	-	20,20,20	1.01	1 (5%)	25,25,25	1.19	3 (12%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	BOG	C	707	-	20,20,20	1.09	1 (5%)	25,25,25	1.36	5 (20%)
3	HEM	C	701	1	50,50,50	2.38	22 (44%)	67,82,82	2.01	20 (29%)
6	BOG	D	708	-	20,20,20	1.09	2 (10%)	25,25,25	1.06	2 (8%)
6	BOG	A	708	-	20,20,20	1.06	1 (5%)	25,25,25	1.28	2 (8%)
6	BOG	D	707	-	20,20,20	1.05	1 (5%)	25,25,25	1.59	4 (16%)
4	NAG	D	705	1	14,14,15	0.35	0	17,19,21	0.45	0
5	DXS	A	706	-	48,48,48	4.39	21 (43%)	69,70,70	1.68	9 (13%)
4	NAG	B	705	1	14,14,15	0.31	0	17,19,21	0.46	0
4	NAG	A	705	1	14,14,15	0.40	0	17,19,21	0.45	0
3	HEM	A	701	1	50,50,50	1.43	8 (16%)	67,82,82	1.28	9 (13%)
4	NAG	A	702	1	14,14,15	0.69	1 (7%)	17,19,21	0.78	1 (5%)
4	NAG	C	705	1	14,14,15	0.33	0	17,19,21	0.48	0
4	NAG	C	702	1	14,14,15	0.35	0	17,19,21	0.53	0
4	NAG	D	702	1	14,14,15	0.66	0	17,19,21	0.68	1 (5%)
5	DXS	D	706	-	48,48,48	5.62	19 (39%)	69,70,70	1.89	12 (17%)
4	NAG	B	702	1	14,14,15	0.45	0	17,19,21	0.51	0
5	DXS	B	706	-	48,48,48	3.85	15 (31%)	69,70,70	4.06	29 (42%)
3	HEM	B	701	1	50,50,50	2.49	22 (44%)	67,82,82	2.17	23 (34%)

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
3	HEM	D	701	1	-	2/14/54/54	-
5	DXS	C	706	-	-	4/32/32/32	0/5/5/5
6	BOG	A	707	-	-	4/11/31/31	0/1/1/1
6	BOG	C	707	-	-	6/11/31/31	0/1/1/1
3	HEM	C	701	1	-	1/14/54/54	-
6	BOG	D	708	-	-	9/11/31/31	0/1/1/1
6	BOG	A	708	-	-	4/11/31/31	0/1/1/1
6	BOG	D	707	-	-	4/11/31/31	0/1/1/1
4	NAG	D	705	1	-	0/6/23/26	0/1/1/1
5	DXS	A	706	-	-	7/32/32/32	0/5/5/5
4	NAG	B	705	1	-	0/6/23/26	0/1/1/1
4	NAG	A	705	1	-	0/6/23/26	0/1/1/1
3	HEM	A	701	1	-	1/14/54/54	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	A	702	1	-	2/6/23/26	0/1/1/1
4	NAG	C	705	1	-	0/6/23/26	0/1/1/1
4	NAG	C	702	1	-	2/6/23/26	0/1/1/1
4	NAG	D	702	1	-	2/6/23/26	0/1/1/1
5	DXS	D	706	-	-	3/32/32/32	0/5/5/5
4	NAG	B	702	1	-	2/6/23/26	0/1/1/1
5	DXS	B	706	-	-	14/32/32/32	0/5/5/5
3	HEM	B	701	1	-	3/14/54/54	-

All (139) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	706	DXS	SAZ-NAY	-24.55	1.25	1.61
5	C	706	DXS	SAZ-NAY	-21.31	1.29	1.61
5	D	706	DXS	CBA-SAZ	-20.94	1.55	1.77
5	A	706	DXS	SAZ-NAY	-16.28	1.37	1.61
5	A	706	DXS	CBA-SAZ	-13.90	1.63	1.77
5	B	706	DXS	SAZ-NAY	13.08	1.81	1.61
5	B	706	DXS	CBA-SAZ	10.77	1.89	1.77
5	B	706	DXS	OBM-SAZ	10.16	1.55	1.43
5	B	706	DXS	OBL-SAZ	9.90	1.55	1.43
5	C	706	DXS	CAH-CAA	-8.26	1.29	1.44
5	D	706	DXS	CAU-CLAX	-8.02	1.55	1.74
5	A	706	DXS	CAH-CAA	-7.86	1.29	1.44
5	C	706	DXS	CAU-CLAX	-7.57	1.56	1.74
5	C	706	DXS	CAG-NAF	-7.45	1.28	1.41
5	D	706	DXS	CAH-CAA	-7.33	1.30	1.44
5	B	706	DXS	CAH-CAA	-7.18	1.31	1.44
5	A	706	DXS	CAH-CAG	-7.03	1.31	1.41
5	D	706	DXS	CAG-NAF	-6.96	1.29	1.41
5	A	706	DXS	CAU-CLAX	-6.75	1.58	1.74
5	C	706	DXS	CAH-CAG	-6.72	1.32	1.41
5	A	706	DXS	CAG-NAF	-6.72	1.29	1.41
5	B	706	DXS	CAR-CAI	-6.40	1.39	1.50
5	D	706	DXS	CAH-CAG	-6.26	1.32	1.41
5	A	706	DXS	CAE-NAF	-6.19	1.27	1.40
5	C	706	DXS	CAR-CAI	-6.07	1.39	1.50
5	A	706	DXS	CAR-CAI	-5.98	1.39	1.50
5	D	706	DXS	CAR-CAI	-5.92	1.39	1.50
3	B	701	HEM	C1B-NB	-5.86	1.30	1.40
5	D	706	DXS	CAJ-CAG	-5.60	1.30	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	706	DXS	CAE-NAF	-5.49	1.29	1.40
5	D	706	DXS	CAE-NAF	-5.42	1.29	1.40
5	C	706	DXS	CAM-CAH	-5.40	1.31	1.39
5	C	706	DXS	CAJ-CAG	-5.34	1.31	1.39
5	D	706	DXS	OBM-SAZ	5.18	1.49	1.43
5	A	706	DXS	CAJ-CAG	-5.17	1.31	1.39
3	D	701	HEM	FE-NB	5.00	2.10	1.94
5	D	706	DXS	OBL-SAZ	4.98	1.49	1.43
3	C	701	HEM	FE-NB	4.92	2.10	1.94
3	C	701	HEM	C1B-NB	-4.92	1.31	1.40
3	C	701	HEM	C4D-ND	-4.89	1.31	1.40
3	B	701	HEM	C4D-ND	-4.73	1.32	1.40
3	C	701	HEM	C3C-C4C	-4.66	1.37	1.46
3	B	701	HEM	C1C-C2C	-4.63	1.36	1.45
5	D	706	DXS	CAM-CAH	-4.63	1.32	1.39
5	B	706	DXS	CAB-CAC	4.59	1.58	1.51
5	A	706	DXS	CAM-CAH	-4.58	1.32	1.39
3	B	701	HEM	C1B-C2B	-4.48	1.35	1.44
3	C	701	HEM	C1C-C2C	-4.47	1.36	1.45
5	B	706	DXS	CAE-NAF	-4.40	1.31	1.40
3	C	701	HEM	C1A-C2A	-4.34	1.35	1.44
5	B	706	DXS	CAH-CAG	-4.34	1.35	1.41
3	A	701	HEM	FE-NB	4.26	2.08	1.94
5	A	706	DXS	OBM-SAZ	4.26	1.48	1.43
5	A	706	DXS	CBJ-NBK	-4.26	1.31	1.43
3	B	701	HEM	C4B-NB	-4.23	1.30	1.38
5	D	706	DXS	CBJ-NBK	-4.19	1.31	1.43
5	C	706	DXS	CBJ-NBK	-4.19	1.31	1.43
3	B	701	HEM	FE-NB	3.95	2.07	1.94
3	B	701	HEM	C1C-NC	-3.91	1.32	1.39
5	D	706	DXS	CBC-CBB	-3.81	1.36	1.43
3	B	701	HEM	C1A-C2A	-3.79	1.37	1.44
3	D	701	HEM	FE-ND	3.76	2.06	1.94
5	B	706	DXS	CAM-CAH	-3.72	1.34	1.39
3	C	701	HEM	C1B-C2B	-3.70	1.37	1.44
3	B	701	HEM	C3C-C4C	-3.64	1.39	1.46
3	C	701	HEM	O2D-CGD	-3.54	1.19	1.30
3	B	701	HEM	C2A-C3A	-3.47	1.29	1.38
3	B	701	HEM	C1D-ND	-3.46	1.32	1.38
5	A	706	DXS	CAE-CAA	-3.45	1.29	1.36
5	D	706	DXS	CAE-CAA	-3.44	1.29	1.36
3	A	701	HEM	FE-ND	3.42	2.05	1.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	706	DXS	CBA-CBB	-3.39	1.37	1.43
3	C	701	HEM	C1D-ND	-3.38	1.32	1.38
5	C	706	DXS	CAE-CAA	-3.34	1.29	1.36
5	D	706	DXS	CAI-NAF	-3.31	1.33	1.41
3	B	701	HEM	C3B-C2B	-3.30	1.30	1.37
3	C	701	HEM	C4B-NB	-3.28	1.32	1.38
5	A	706	DXS	OBL-SAZ	3.26	1.47	1.43
5	C	706	DXS	OBL-SAZ	3.26	1.47	1.43
3	B	701	HEM	O2D-CGD	-3.25	1.20	1.30
5	C	706	DXS	OBO-CAL	-3.24	1.31	1.37
5	A	706	DXS	CBC-CBB	-3.23	1.37	1.43
3	C	701	HEM	FE-ND	-3.16	1.85	1.94
5	B	706	DXS	CAJ-CAG	-3.15	1.34	1.39
5	D	706	DXS	CBA-CBB	-3.14	1.37	1.43
3	B	701	HEM	C4A-NA	-3.09	1.33	1.39
3	B	701	HEM	C4C-NC	-3.07	1.33	1.39
5	A	706	DXS	CAI-NAF	-2.96	1.34	1.41
3	B	701	HEM	FE-NC	2.92	2.04	1.95
3	B	701	HEM	C3D-C2D	-2.90	1.30	1.36
5	C	706	DXS	CAI-NAF	-2.88	1.34	1.41
3	A	701	HEM	CAC-C3C	2.88	1.55	1.47
3	C	701	HEM	FE-NC	2.88	2.04	1.95
3	C	701	HEM	C2A-C3A	-2.84	1.31	1.38
3	C	701	HEM	C4A-NA	-2.83	1.34	1.39
3	C	701	HEM	C3B-C2B	-2.81	1.31	1.37
3	D	701	HEM	CAB-C3B	2.80	1.54	1.47
3	D	701	HEM	CAC-C3C	2.80	1.54	1.47
3	C	701	HEM	C4C-NC	-2.79	1.34	1.39
3	A	701	HEM	CAB-C3B	2.79	1.54	1.47
3	B	701	HEM	O2A-CGA	-2.75	1.21	1.30
3	B	701	HEM	FE-ND	-2.75	1.86	1.94
5	C	706	DXS	CBC-CBB	-2.65	1.38	1.43
3	C	701	HEM	C4A-C3A	-2.62	1.37	1.43
6	A	708	BOG	O2-C2	-2.53	1.36	1.43
3	B	701	HEM	C4A-C3A	-2.52	1.37	1.43
5	C	706	DXS	OBO-CBP	-2.49	1.35	1.42
5	B	706	DXS	CBN-CAE	2.47	1.53	1.49
3	C	701	HEM	C1A-NA	-2.47	1.35	1.39
5	C	706	DXS	CBG-CBB	-2.46	1.37	1.42
3	A	701	HEM	FE-NA	2.45	2.03	1.95
5	A	706	DXS	CBG-CBB	-2.45	1.37	1.42
5	C	706	DXS	CAB-CAA	2.44	1.56	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	D	708	BOG	O2-C2	-2.43	1.36	1.43
3	C	701	HEM	C3D-C2D	-2.38	1.31	1.36
3	B	701	HEM	C3C-C2C	-2.35	1.32	1.37
5	A	706	DXS	CBJ-CBC	-2.34	1.37	1.42
6	A	707	BOG	O2-C2	-2.34	1.37	1.43
6	D	707	BOG	O2-C2	-2.32	1.37	1.43
6	C	707	BOG	O2-C2	-2.31	1.37	1.43
5	A	706	DXS	CAW-CAR	-2.30	1.35	1.39
3	B	701	HEM	CHB-C4A	-2.29	1.34	1.39
3	A	701	HEM	FE-NC	2.26	2.02	1.95
5	A	706	DXS	CBD-CBC	-2.25	1.37	1.42
4	A	702	NAG	O5-C1	2.22	1.47	1.43
3	C	701	HEM	CHA-C1A	-2.21	1.34	1.39
5	B	706	DXS	CBJ-NBK	-2.20	1.37	1.43
5	B	706	DXS	CAE-CAA	-2.18	1.31	1.36
3	A	701	HEM	C2A-C3A	-2.15	1.33	1.38
5	C	706	DXS	CBD-CBC	-2.12	1.38	1.42
3	D	701	HEM	C2A-C3A	-2.11	1.33	1.38
5	B	706	DXS	CAI-NAF	-2.11	1.36	1.41
3	C	701	HEM	O2A-CGA	-2.11	1.23	1.30
6	D	708	BOG	O5-C5	-2.10	1.39	1.44
5	D	706	DXS	OBO-CAL	-2.09	1.33	1.37
5	C	706	DXS	CAB-CAC	2.08	1.54	1.51
5	D	706	DXS	CAS-CAR	-2.08	1.36	1.39
3	C	701	HEM	C4D-C3D	-2.05	1.41	1.45
3	A	701	HEM	CMA-C3A	2.02	1.54	1.50

All (142) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	706	DXS	OBL-SAZ-NAY	-15.72	82.55	107.03
5	B	706	DXS	CBA-SAZ-NAY	-14.49	80.02	106.78
5	B	706	DXS	OBM-SAZ-NAY	-11.67	88.86	107.03
5	B	706	DXS	CAG-NAF-CAE	-11.09	98.76	107.80
5	D	706	DXS	OBM-SAZ-OBL	-9.43	108.06	119.52
5	C	706	DXS	OBM-SAZ-OBL	-8.59	109.08	119.52
5	B	706	DXS	OBL-SAZ-CBA	8.53	122.95	108.09
3	B	701	HEM	C3B-C4B-NB	-7.83	103.84	109.47
5	A	706	DXS	OBM-SAZ-OBL	-7.51	110.39	119.52
5	B	706	DXS	CAJ-CAG-CAH	-5.74	115.54	121.95
6	D	707	BOG	C1'-O1-C1	5.66	123.35	113.68
3	C	701	HEM	C3B-C4B-NB	-5.56	105.48	109.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	706	DXS	CBN-CAE-CAA	-5.48	119.49	128.07
3	B	701	HEM	C1B-NB-C4B	5.10	111.24	105.21
5	B	706	DXS	CBC-CBJ-NBK	4.99	126.00	118.37
3	B	701	HEM	CHD-C1D-ND	4.92	129.72	124.42
3	C	701	HEM	CHC-C4B-NB	4.78	129.57	124.42
5	B	706	DXS	CBI-CBJ-NBK	-4.77	114.53	121.61
5	C	706	DXS	CAV-CAU-CAT	-4.77	115.33	121.24
5	B	706	DXS	CAA-CAE-NAF	4.75	113.36	108.85
3	C	701	HEM	O2D-CGD-O1D	-4.70	111.25	123.33
5	B	706	DXS	CAP-CAQ-NAY	4.49	126.18	111.80
5	D	706	DXS	CAV-CAU-CAT	-4.27	115.94	121.24
3	B	701	HEM	CHC-C4B-NB	4.15	128.88	124.42
5	A	706	DXS	CAV-CAU-CAT	-4.06	116.21	121.24
5	B	706	DXS	OBM-SAZ-CBA	4.05	115.16	108.09
5	D	706	DXS	OBM-SAZ-NAY	3.84	113.00	107.03
3	B	701	HEM	O2D-CGD-O1D	-3.83	113.47	123.33
3	B	701	HEM	CHB-C1B-NB	3.77	129.03	124.37
6	A	708	BOG	C1'-O1-C1	3.72	120.03	113.68
5	B	706	DXS	CAM-CAH-CAG	3.61	124.12	119.82
3	C	701	HEM	CHB-C1B-NB	3.60	128.82	124.37
6	A	708	BOG	C1-O5-C5	-3.57	106.74	113.72
5	C	706	DXS	CAS-CAT-CAU	3.51	122.77	119.24
3	C	701	HEM	C1B-NB-C4B	3.49	109.34	105.21
5	B	706	DXS	CAT-CAS-CAR	3.40	124.43	120.80
5	C	706	DXS	CAG-NAF-CAE	-3.40	105.03	107.80
5	C	706	DXS	CAR-CAI-NAF	3.35	124.36	117.92
5	C	706	DXS	CAW-CAV-CAU	3.35	122.61	119.24
5	B	706	DXS	CBR-NBK-CBJ	3.32	124.29	114.19
3	B	701	HEM	CHD-C1D-C2D	-3.31	119.81	125.03
3	C	701	HEM	CHC-C1C-NC	3.30	128.04	124.45
3	C	701	HEM	CHA-C4D-C3D	-3.28	119.18	125.23
3	C	701	HEM	CHA-C4D-ND	3.27	128.41	124.37
5	D	706	DXS	CAS-CAT-CAU	3.23	122.49	119.24
5	B	706	DXS	CAS-CAR-CAI	-3.21	112.05	120.28
5	D	706	DXS	CAV-CAU-CLAX	3.18	124.05	119.36
5	C	706	DXS	CBN-CAE-CAA	-3.17	123.10	128.07
6	A	707	BOG	C1'-O1-C1	3.17	119.10	113.68
5	C	706	DXS	OBM-SAZ-NAY	3.15	111.94	107.03
5	B	706	DXS	CAR-CAI-NAF	3.14	123.95	117.92
5	C	706	DXS	CBA-SAZ-NAY	3.13	112.57	106.78
6	C	707	BOG	C6-C5-C4	-3.12	105.36	113.02
5	B	706	DXS	CAQ-CAP-NAD	-3.08	101.97	111.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	706	DXS	CAG-NAF-CAE	-3.07	105.30	107.80
3	C	701	HEM	CHD-C1D-ND	3.02	127.67	124.42
5	A	706	DXS	CBG-CBB-CBA	-3.00	119.08	123.58
5	A	706	DXS	OBM-SAZ-NAY	2.98	111.68	107.03
5	B	706	DXS	CAW-CAR-CAS	-2.97	114.78	118.57
3	C	701	HEM	CAD-CBD-CGD	2.97	121.55	113.67
6	D	707	BOG	O1-C1-C2	2.96	112.77	108.27
5	B	706	DXS	OAO-CAI-CAR	-2.95	114.44	120.29
5	A	706	DXS	CAR-CAI-NAF	2.95	123.59	117.92
5	A	706	DXS	CAT-CAU-CLAX	2.94	123.70	119.36
6	D	707	BOG	O5-C5-C4	2.93	114.98	109.70
5	C	706	DXS	CAA-CAE-NAF	2.91	111.61	108.85
3	D	701	HEM	CHC-C1C-NC	2.88	127.59	124.45
3	B	701	HEM	C4C-CHD-C1D	-2.87	119.92	126.02
3	B	701	HEM	O2A-CGA-CBA	2.87	123.06	114.00
5	B	706	DXS	CAH-CAG-NAF	2.85	110.88	107.69
3	B	701	HEM	CAC-C3C-C4C	2.82	131.56	124.82
3	A	701	HEM	CBA-CAA-C2A	-2.76	104.91	112.53
3	B	701	HEM	O2D-CGD-CBD	2.75	122.68	114.00
6	C	707	BOG	C1'-O1-C1	2.74	118.36	113.68
5	B	706	DXS	CAW-CAR-CAI	2.73	127.28	120.28
3	B	701	HEM	CHA-C4D-ND	2.72	127.73	124.37
3	C	701	HEM	CHD-C4C-NC	2.70	127.40	124.45
3	C	701	HEM	CMD-C2D-C1D	2.70	129.25	125.03
3	D	701	HEM	CAA-CBA-CGA	-2.68	106.57	113.67
3	D	701	HEM	C2A-C1A-NA	-2.66	107.20	110.15
3	C	701	HEM	CAC-C3C-C4C	2.64	131.13	124.82
3	A	701	HEM	C3B-C4B-NB	-2.61	107.59	109.47
6	A	707	BOG	O5-C5-C4	2.59	114.36	109.70
3	B	701	HEM	CHA-C4D-C3D	-2.59	120.46	125.23
6	D	708	BOG	C1'-O1-C1	2.59	118.10	113.68
5	A	706	DXS	CAW-CAV-CAU	2.58	121.84	119.24
4	A	702	NAG	C1-O5-C5	2.58	115.65	112.19
5	C	706	DXS	CBB-CBA-SAZ	2.57	124.68	121.66
5	D	706	DXS	CAR-CAI-NAF	2.56	122.84	117.92
3	A	701	HEM	CHC-C1C-NC	2.53	127.21	124.45
5	D	706	DXS	CAA-CAE-NAF	2.51	111.23	108.85
3	B	701	HEM	CBA-CAA-C2A	-2.50	105.63	112.53
5	A	706	DXS	OBM-SAZ-CBA	2.49	112.44	108.09
5	B	706	DXS	CBE-CBD-CBC	2.48	124.27	120.91
5	B	706	DXS	CAK-CAJ-CAG	2.48	124.08	119.10
3	C	701	HEM	O2D-CGD-CBD	2.46	121.78	114.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	701	HEM	C1C-CHC-C4B	-2.46	120.80	126.02
3	C	701	HEM	C4C-CHD-C1D	-2.45	120.81	126.02
3	B	701	HEM	CBD-CAD-C3D	-2.45	105.76	112.53
3	B	701	HEM	CAB-C3B-C2B	-2.43	120.54	128.43
5	C	706	DXS	CAT-CAU-CLAX	2.42	122.93	119.36
5	B	706	DXS	CBG-CBB-CBA	-2.41	119.96	123.58
3	C	701	HEM	CAC-C3C-C2C	-2.40	120.63	128.43
5	D	706	DXS	CAW-CAR-CAS	-2.40	115.52	118.57
6	C	707	BOG	C4'-C3'-C2'	-2.30	102.73	114.37
3	C	701	HEM	CHB-C1B-C2B	-2.27	120.50	126.95
3	A	701	HEM	C4D-ND-C1D	2.26	107.89	105.21
6	D	707	BOG	O5-C1-O1	-2.25	104.73	110.04
3	D	701	HEM	C4D-ND-C1D	2.25	107.87	105.21
5	A	706	DXS	CBN-CAE-NAF	2.24	125.42	122.07
3	D	701	HEM	CBA-CAA-C2A	-2.24	106.35	112.53
3	C	701	HEM	O2A-CGA-CBA	2.23	121.05	114.00
5	B	706	DXS	CAG-CAH-CAA	-2.23	105.34	107.38
5	B	706	DXS	CAK-CAL-CAM	-2.22	117.52	120.50
3	A	701	HEM	C2A-C1A-NA	-2.22	107.69	110.15
5	B	706	DXS	CBD-CBC-CBB	-2.21	115.68	118.42
3	A	701	HEM	CAA-CBA-CGA	-2.18	107.87	113.67
3	A	701	HEM	C3B-C2B-C1B	2.17	108.04	106.41
5	C	706	DXS	CBR-NBK-CBJ	2.17	120.78	114.19
3	B	701	HEM	O2A-CGA-O1A	-2.16	117.77	123.33
5	C	706	DXS	CAH-CAA-CAE	-2.15	106.25	108.13
5	C	706	DXS	CBN-CAE-NAF	2.15	125.28	122.07
3	B	701	HEM	C4A-CHB-C1B	-2.14	121.20	126.25
5	B	706	DXS	CBA-CBB-CBC	2.14	120.82	117.97
3	B	701	HEM	CAD-CBD-CGD	2.14	119.34	113.67
6	C	707	BOG	C1-O5-C5	-2.14	109.54	113.72
6	C	707	BOG	O5-C5-C4	2.14	113.55	109.70
3	B	701	HEM	CAC-C3C-C2C	-2.13	121.52	128.43
3	A	701	HEM	C3C-C2C-C1C	2.12	109.06	107.05
5	D	706	DXS	CAW-CAV-CAU	2.11	121.37	119.24
6	D	708	BOG	O5-C5-C6	-2.11	101.22	106.44
3	B	701	HEM	CHC-C1C-NC	2.10	126.74	124.45
5	D	706	DXS	CBR-NBK-CBQ	2.10	122.94	116.18
5	D	706	DXS	CAH-CAG-NAF	2.08	110.02	107.69
3	B	701	HEM	CHB-C1B-C2B	-2.07	121.05	126.95
3	C	701	HEM	CBB-CAB-C3B	-2.06	117.22	127.53
3	A	701	HEM	C1B-NB-C4B	2.06	107.65	105.21
3	B	701	HEM	C1A-CHA-C4D	-2.04	121.45	126.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	706	DXS	CBF-CBA-CBB	-2.03	118.68	120.95
6	A	707	BOG	C4-C3-C2	-2.02	107.28	110.83
4	D	702	NAG	C1-O5-C5	2.00	114.87	112.19
3	D	701	HEM	C4C-NC-C1C	2.00	109.08	105.82

There are no chirality outliers.

All (70) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	706	DXS	CBB-CBA-SAZ-OBL
5	B	706	DXS	CBB-CBA-SAZ-OBM
5	B	706	DXS	CBF-CBA-SAZ-OBL
5	B	706	DXS	CBF-CBA-SAZ-OBM
5	B	706	DXS	CAQ-NAY-SAZ-CBA
6	A	708	BOG	O5-C1-O1-C1'
6	A	708	BOG	C2'-C1'-O1-C1
6	C	707	BOG	C2'-C1'-O1-C1
6	D	707	BOG	O5-C1-O1-C1'
5	B	706	DXS	CAQ-NAY-SAZ-OBM
6	D	708	BOG	C4-C5-C6-O6
6	D	708	BOG	O5-C5-C6-O6
6	A	707	BOG	O5-C5-C6-O6
4	D	702	NAG	O5-C5-C6-O6
5	D	706	DXS	CAQ-NAY-SAZ-OBL
4	A	702	NAG	O5-C5-C6-O6
4	B	702	NAG	O5-C5-C6-O6
4	C	702	NAG	O5-C5-C6-O6
6	D	707	BOG	O5-C5-C6-O6
4	D	702	NAG	C4-C5-C6-O6
6	C	707	BOG	O1-C1'-C2'-C3'
5	A	706	DXS	CAQ-NAY-SAZ-OBL
5	B	706	DXS	CAR-CAI-NAF-CAE
6	D	707	BOG	C2-C1-O1-C1'
6	D	708	BOG	C2-C1-O1-C1'
6	D	708	BOG	O5-C1-O1-C1'
5	D	706	DXS	CAQ-NAY-SAZ-OBM
6	D	708	BOG	C2'-C1'-O1-C1
5	B	706	DXS	CBC-CBJ-NBK-CBR
6	A	707	BOG	C4-C5-C6-O6
4	A	702	NAG	C4-C5-C6-O6
5	B	706	DXS	OAO-CAI-CAR-CAW
6	C	707	BOG	C1'-C2'-C3'-C4'

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Mol	Chain	Res	Type	Atoms
5	C	706	DXS	CAQ-NAY-SAZ-OBL
6	C	707	BOG	C2'-C3'-C4'-C5'
6	A	708	BOG	C2'-C3'-C4'-C5'
5	A	706	DXS	CAQ-NAY-SAZ-OBM
5	D	706	DXS	CAQ-NAY-SAZ-CBA
6	D	708	BOG	C1'-C2'-C3'-C4'
4	B	702	NAG	C4-C5-C6-O6
6	C	707	BOG	C3'-C4'-C5'-C6'
6	D	708	BOG	C2'-C3'-C4'-C5'
5	A	706	DXS	CAQ-NAY-SAZ-CBA
4	C	702	NAG	C4-C5-C6-O6
6	D	708	BOG	C5'-C6'-C7'-C8'
6	D	708	BOG	C4'-C5'-C6'-C7'
5	B	706	DXS	CBI-CBJ-NBK-CBR
6	C	707	BOG	C5'-C6'-C7'-C8'
5	B	706	DXS	CAP-CAQ-NAY-SAZ
5	A	706	DXS	OAQ-CAI-NAF-CAE
5	B	706	DXS	CBC-CBJ-NBK-CBQ
3	B	701	HEM	C4B-C3B-CAB-CBB
3	C	701	HEM	C4C-C3C-CAC-CBC
6	A	708	BOG	C1'-C2'-C3'-C4'
5	A	706	DXS	OAQ-CAI-CAR-CAW
5	C	706	DXS	CAQ-NAY-SAZ-CBA
6	D	707	BOG	C4-C5-C6-O6
5	C	706	DXS	CAQ-NAY-SAZ-OBM
5	B	706	DXS	NAF-CAI-CAR-CAW
6	A	707	BOG	O1-C1'-C2'-C3'
6	A	707	BOG	O5-C1-O1-C1'
3	B	701	HEM	CAA-CBA-CGA-O2A
5	A	706	DXS	OAQ-CAI-CAR-CAS
5	C	706	DXS	CBC-CBJ-NBK-CBR
3	B	701	HEM	CAA-CBA-CGA-O1A
5	A	706	DXS	NAF-CAI-CAR-CAW
3	D	701	HEM	CAA-CBA-CGA-O2A
5	B	706	DXS	OAQ-CAI-NAF-CAG
3	A	701	HEM	CAA-CBA-CGA-O2A
3	D	701	HEM	CAA-CBA-CGA-O1A

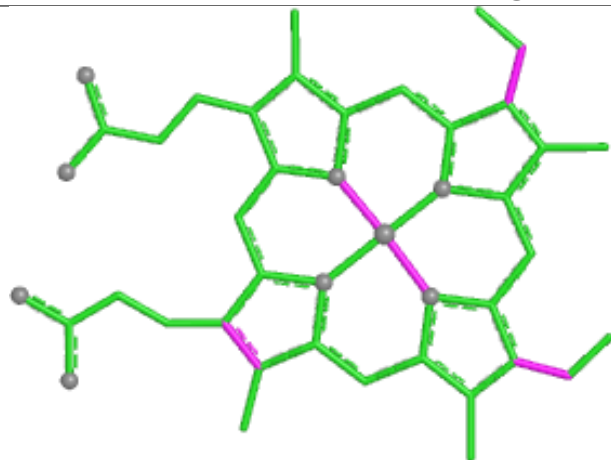
There are no ring outliers.

11 monomers are involved in 27 short contacts:

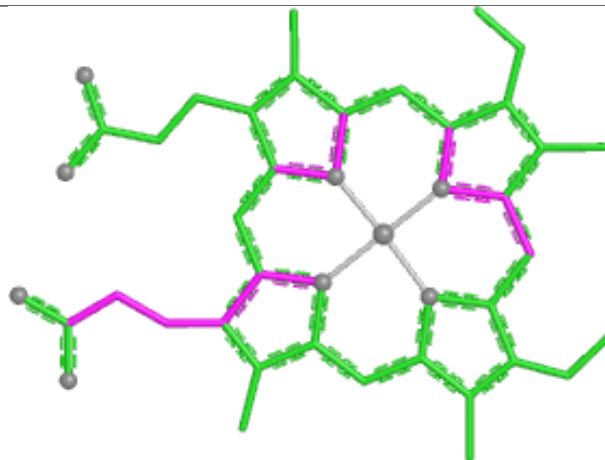
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	701	HEM	1	0
5	C	706	DXS	3	0
6	A	707	BOG	2	0
6	C	707	BOG	3	0
3	C	701	HEM	5	0
6	D	707	BOG	1	0
5	A	706	DXS	2	0
3	A	701	HEM	2	0
5	D	706	DXS	3	0
5	B	706	DXS	3	0
3	B	701	HEM	2	0

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.

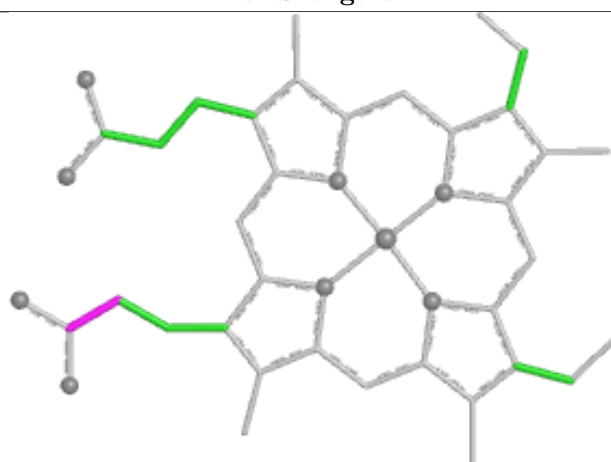
Ligand HEM D 701



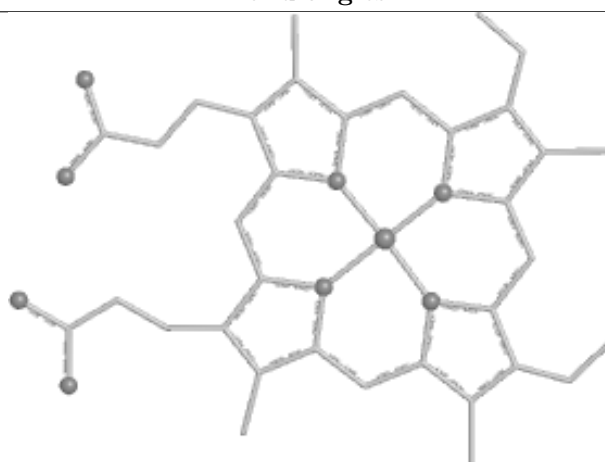
Bond lengths



Bond angles

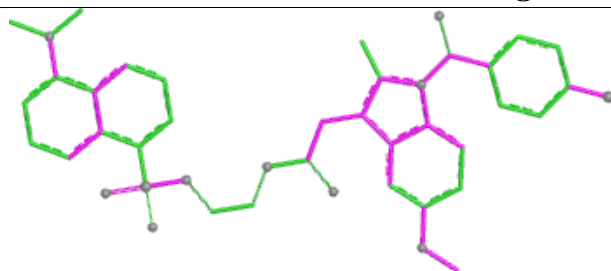


Torsions

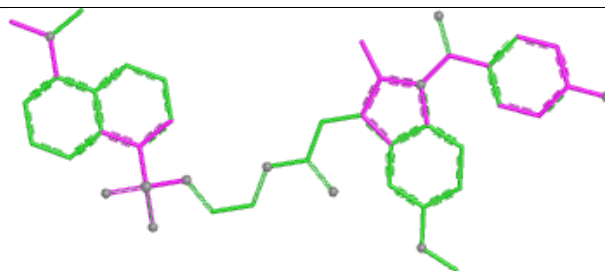


Rings

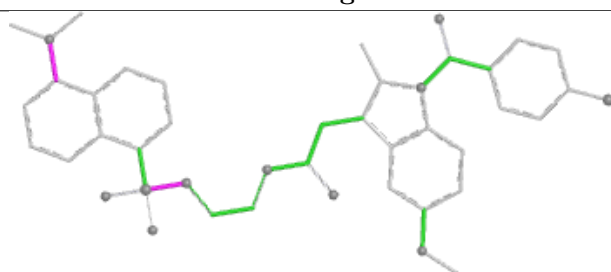
Ligand DXS C 706



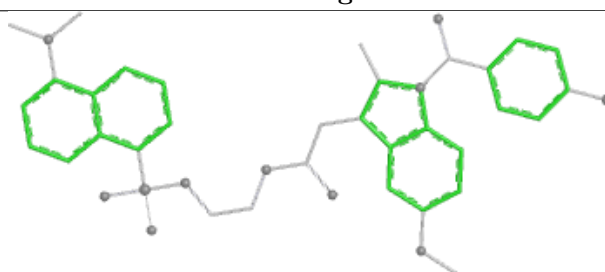
Bond lengths



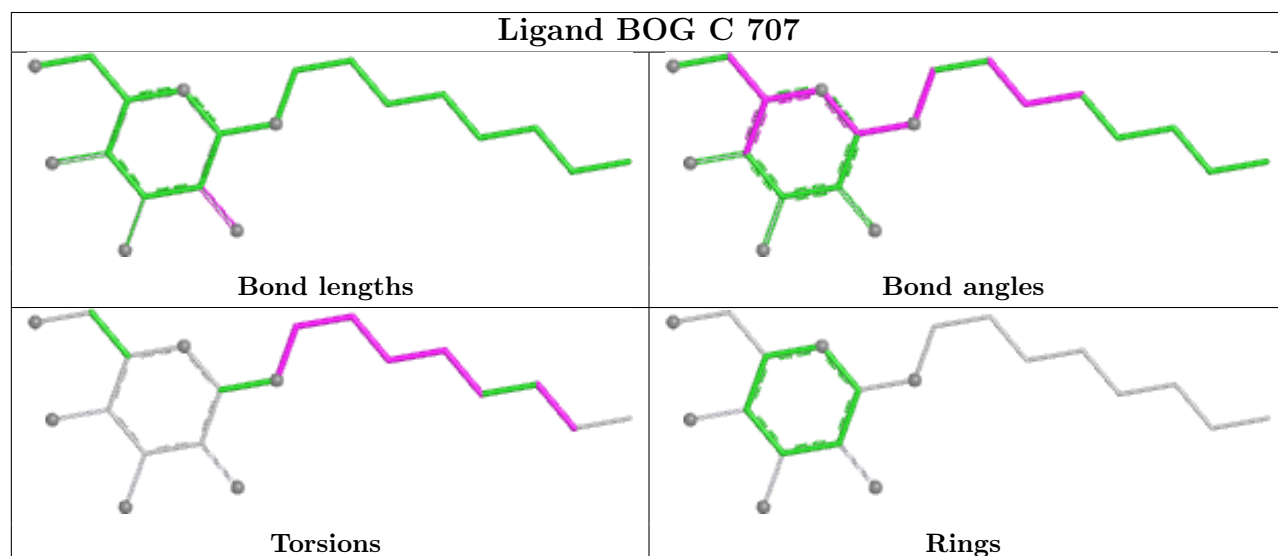
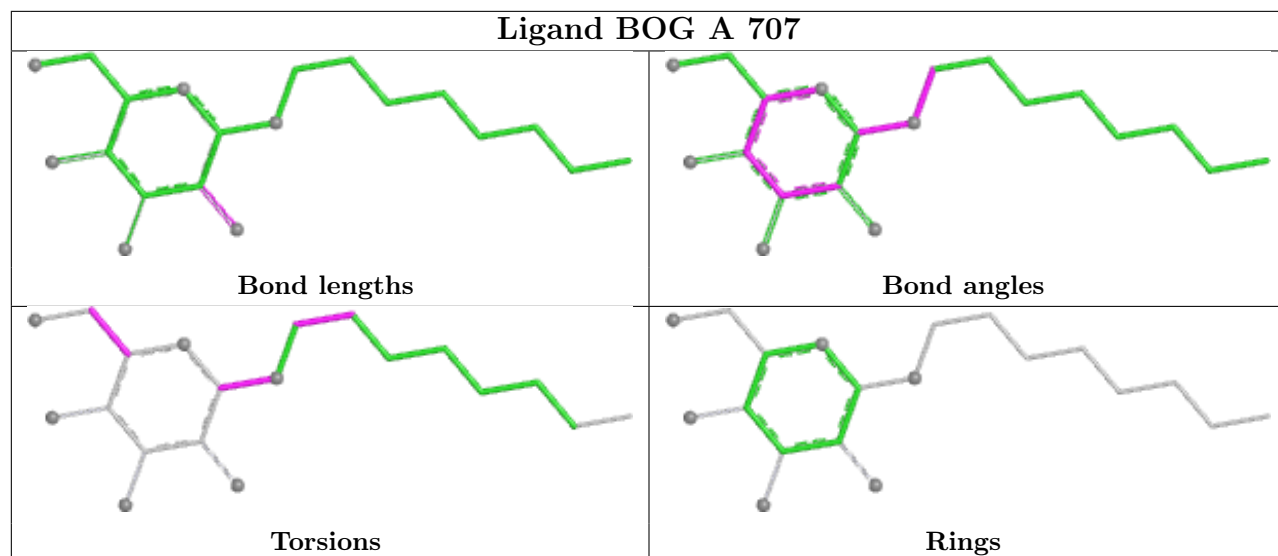
Bond angles

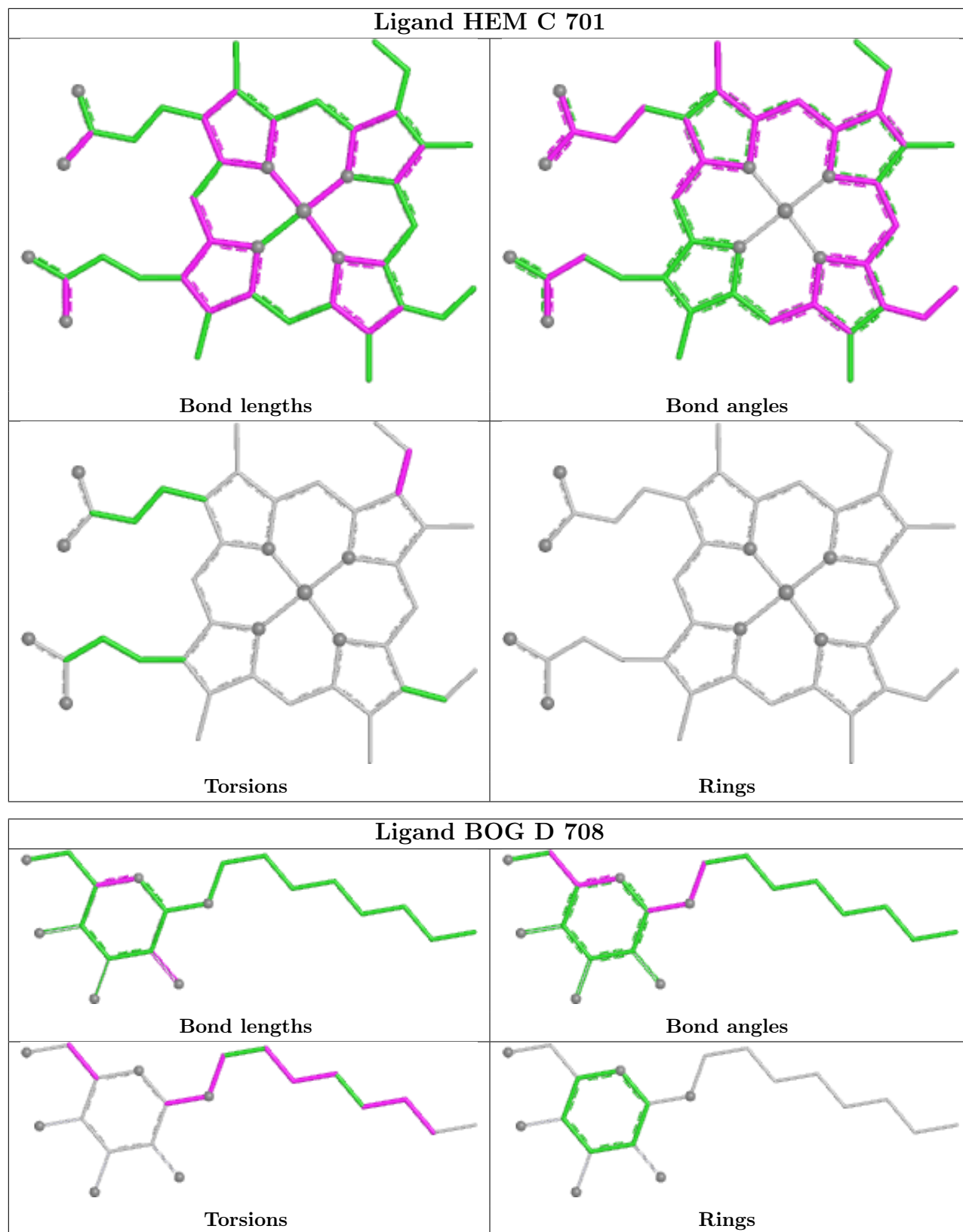


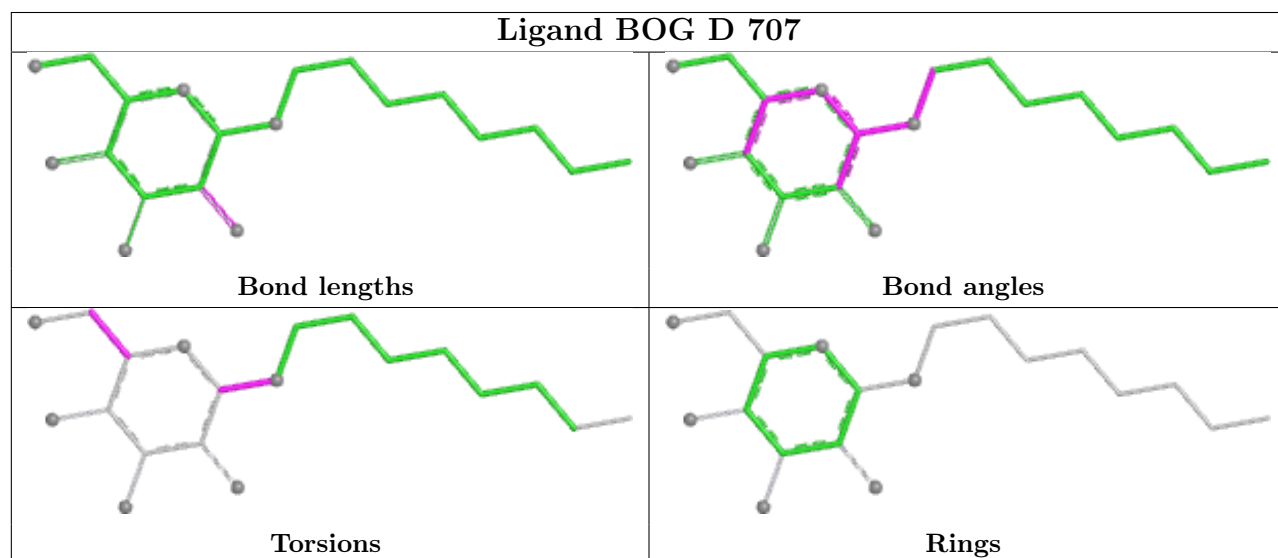
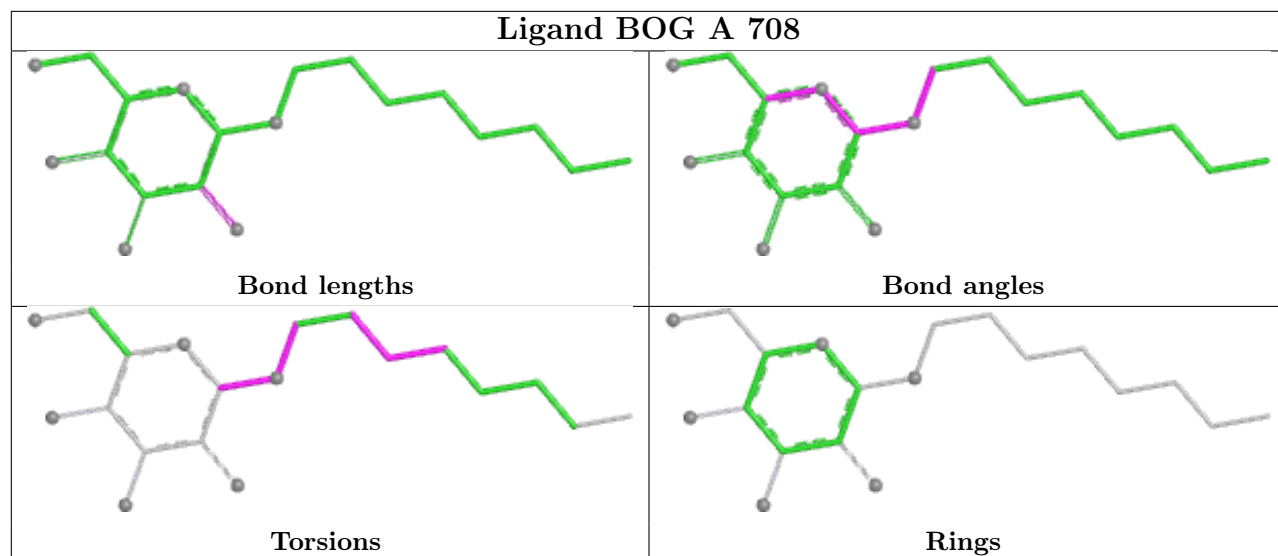
Torsions



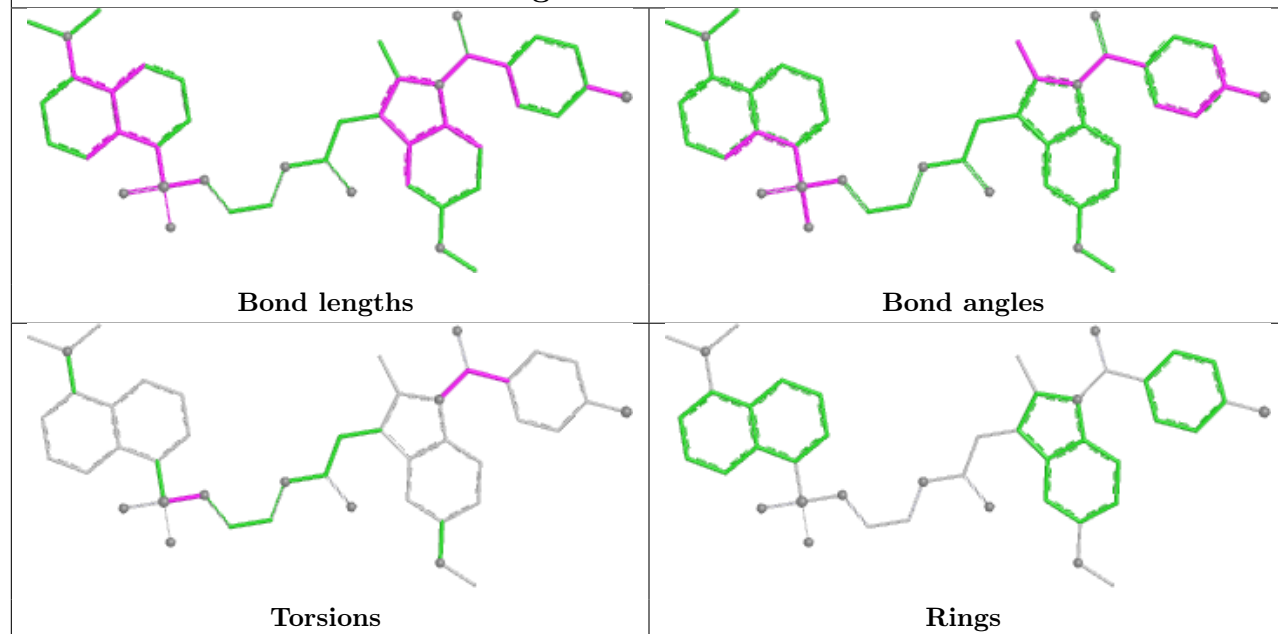
Rings



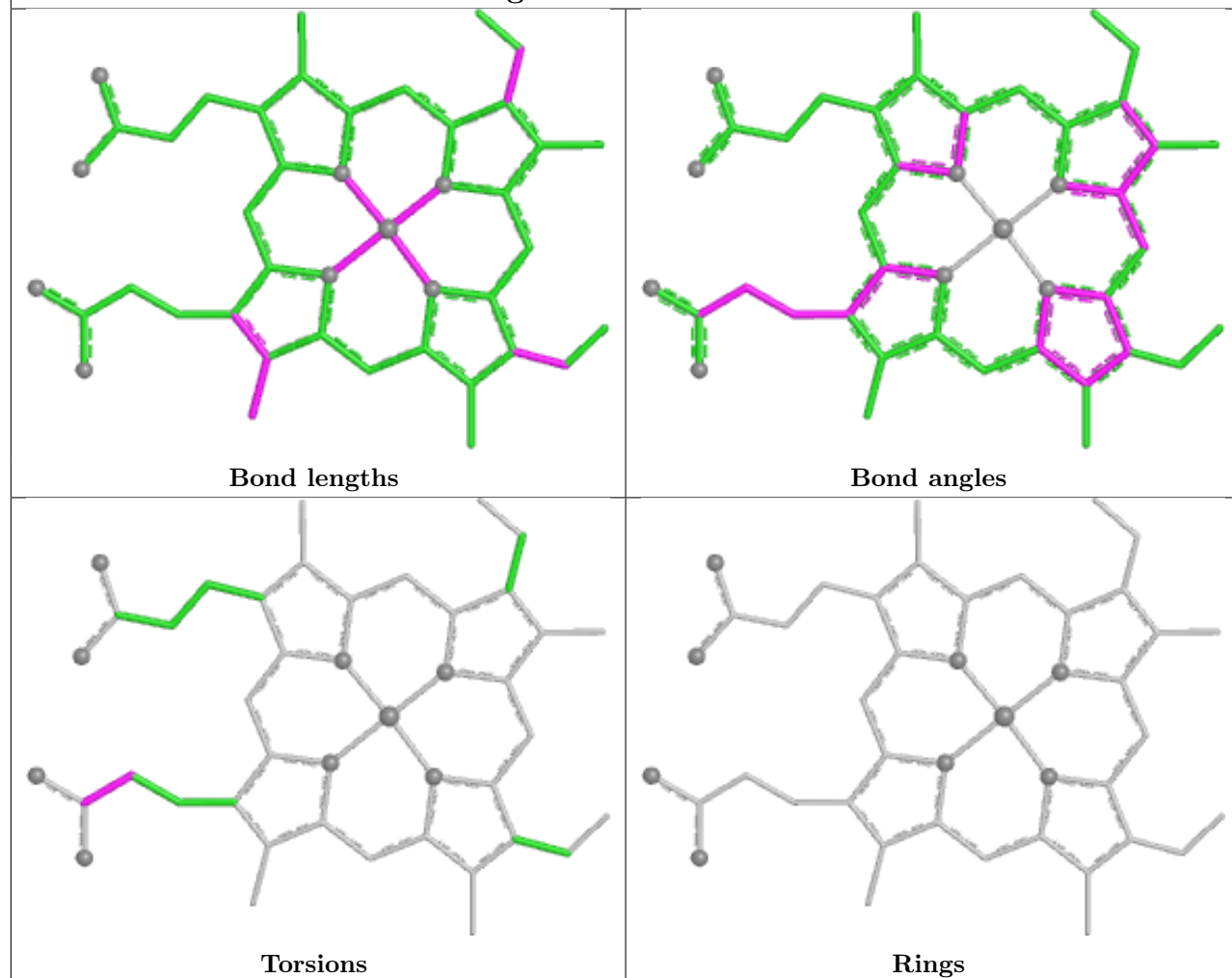




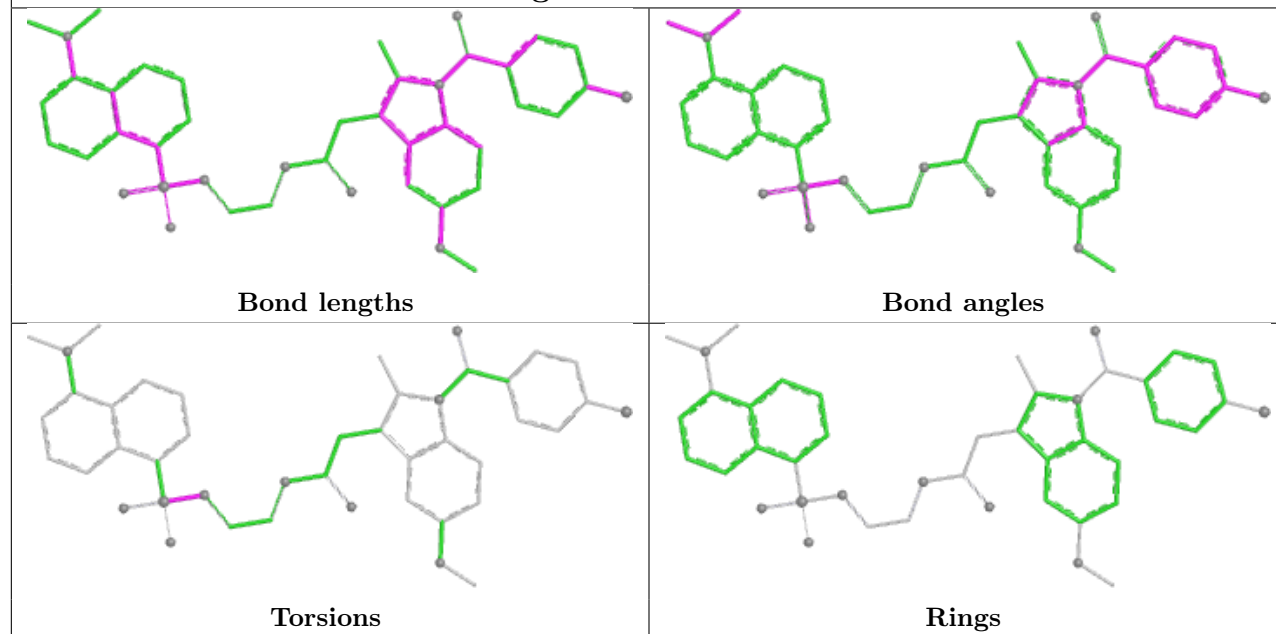
Ligand DXS A 706



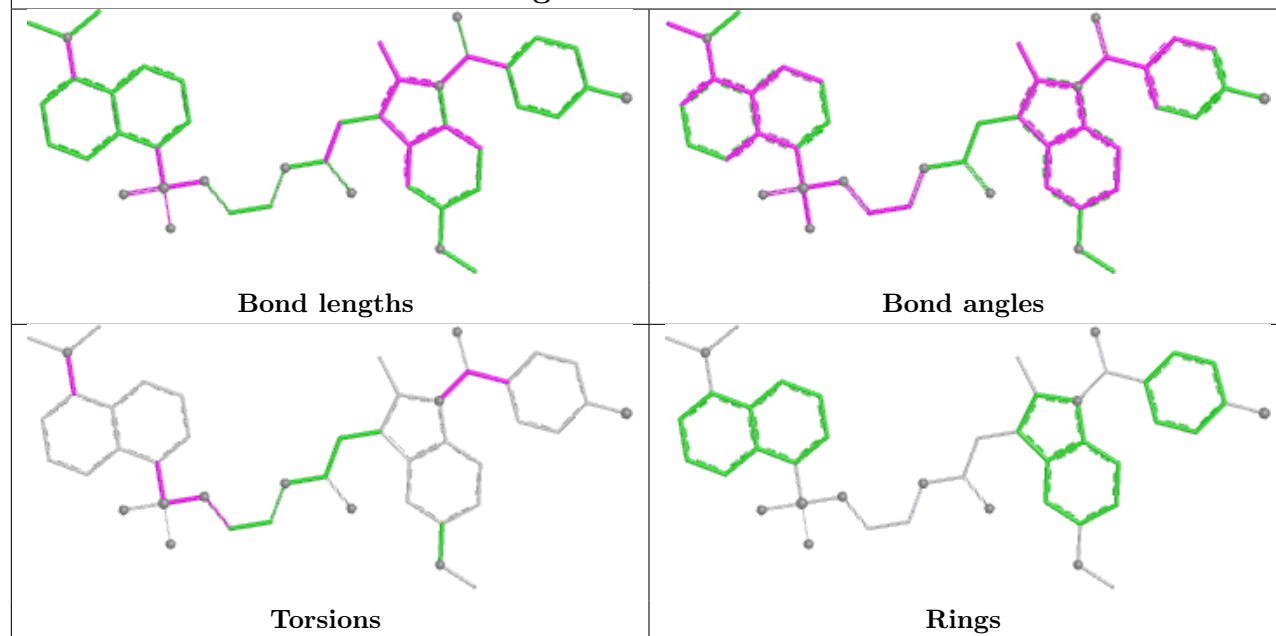
Ligand HEM A 701

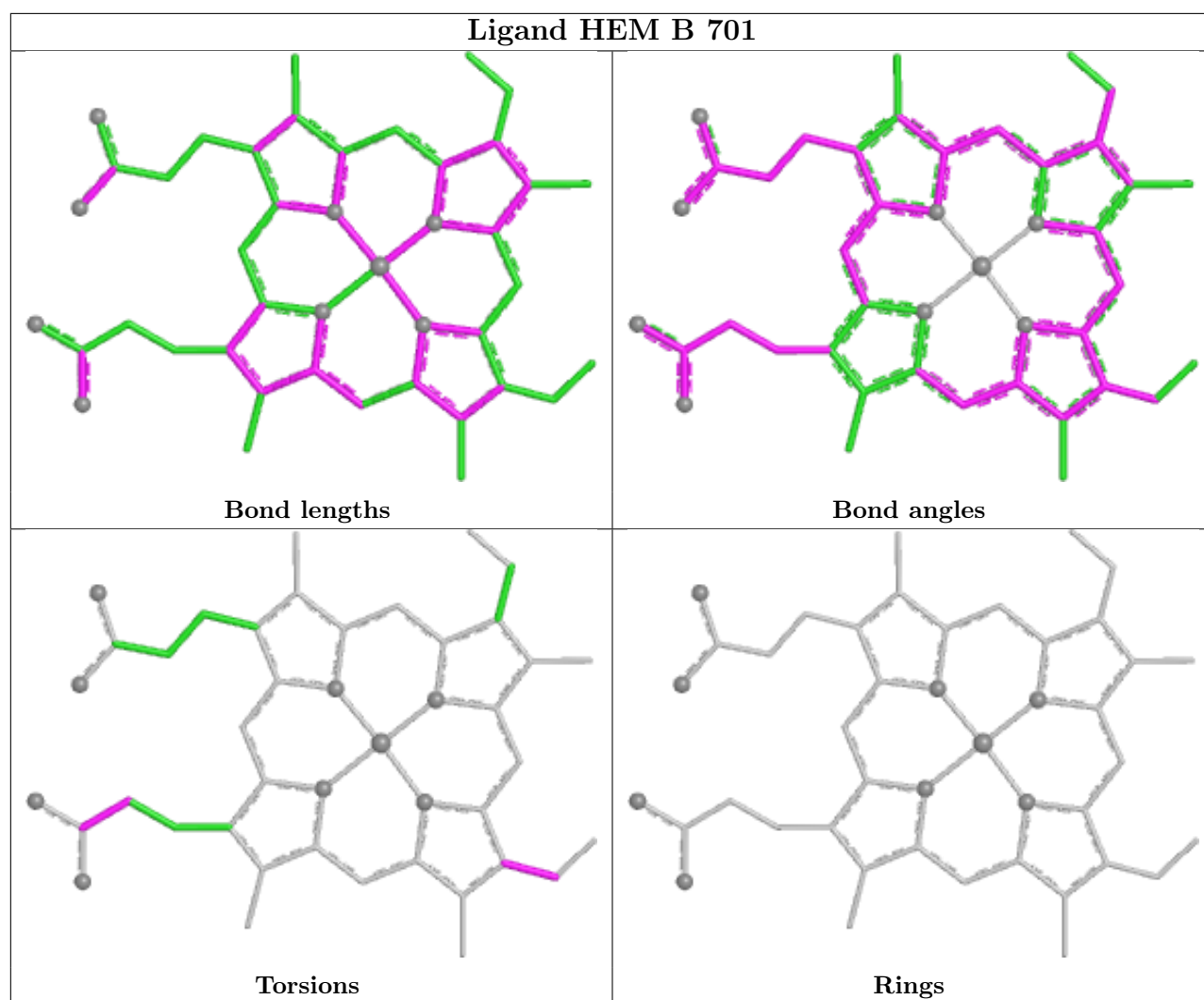


Ligand DXS D 706



Ligand DXS B 706





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%)	-0.01	10 (1%) 67 65	16, 31, 53, 69	0
1	B	552/587 (94%)	0.13	18 (3%) 49 46	15, 33, 57, 70	0
1	C	552/587 (94%)	0.03	9 (1%) 70 68	17, 32, 53, 67	0
1	D	552/587 (94%)	0.15	9 (1%) 70 68	18, 36, 59, 86	0
All	All	2208/2348 (94%)	0.08	46 (2%) 63 61	15, 33, 56, 86	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	122	TYR	4.0
1	D	583	GLN	3.5
1	A	122	TYR	3.5
1	C	122	TYR	3.4
1	C	121	SER	3.3
1	B	33	ALA	3.2
1	A	583	GLN	3.2
1	C	79	LYS	3.1
1	B	282	ASN	3.0
1	C	583	GLN	3.0
1	D	121	SER	3.0
1	B	107	PHE	2.9
1	D	74	PHE	2.9
1	D	107	PHE	2.9
1	B	280	PRO	2.8
1	B	583	GLN	2.7
1	B	122	TYR	2.7
1	A	121	SER	2.6
1	B	101	ASN	2.5
1	B	74	PHE	2.5
1	B	121	SER	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	120	ARG	2.4
1	D	120	ARG	2.4
1	A	33	ALA	2.3
1	A	107	PHE	2.3
1	D	33	ALA	2.3
1	B	77	ARG	2.3
1	B	384	LEU	2.3
1	B	364	GLU	2.3
1	B	281	GLU	2.2
1	D	80	LEU	2.2
1	B	105	ASN	2.2
1	C	364	GLU	2.2
1	A	384	LEU	2.2
1	C	77	ARG	2.2
1	A	409	TYR	2.2
1	D	123	LEU	2.2
1	C	107	PHE	2.2
1	C	33	ALA	2.1
1	B	82	LEU	2.1
1	B	81	LEU	2.1
1	A	74	PHE	2.1
1	B	283	LEU	2.1
1	A	123	LEU	2.0
1	B	79	LYS	2.0
1	C	318	GLN	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

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.

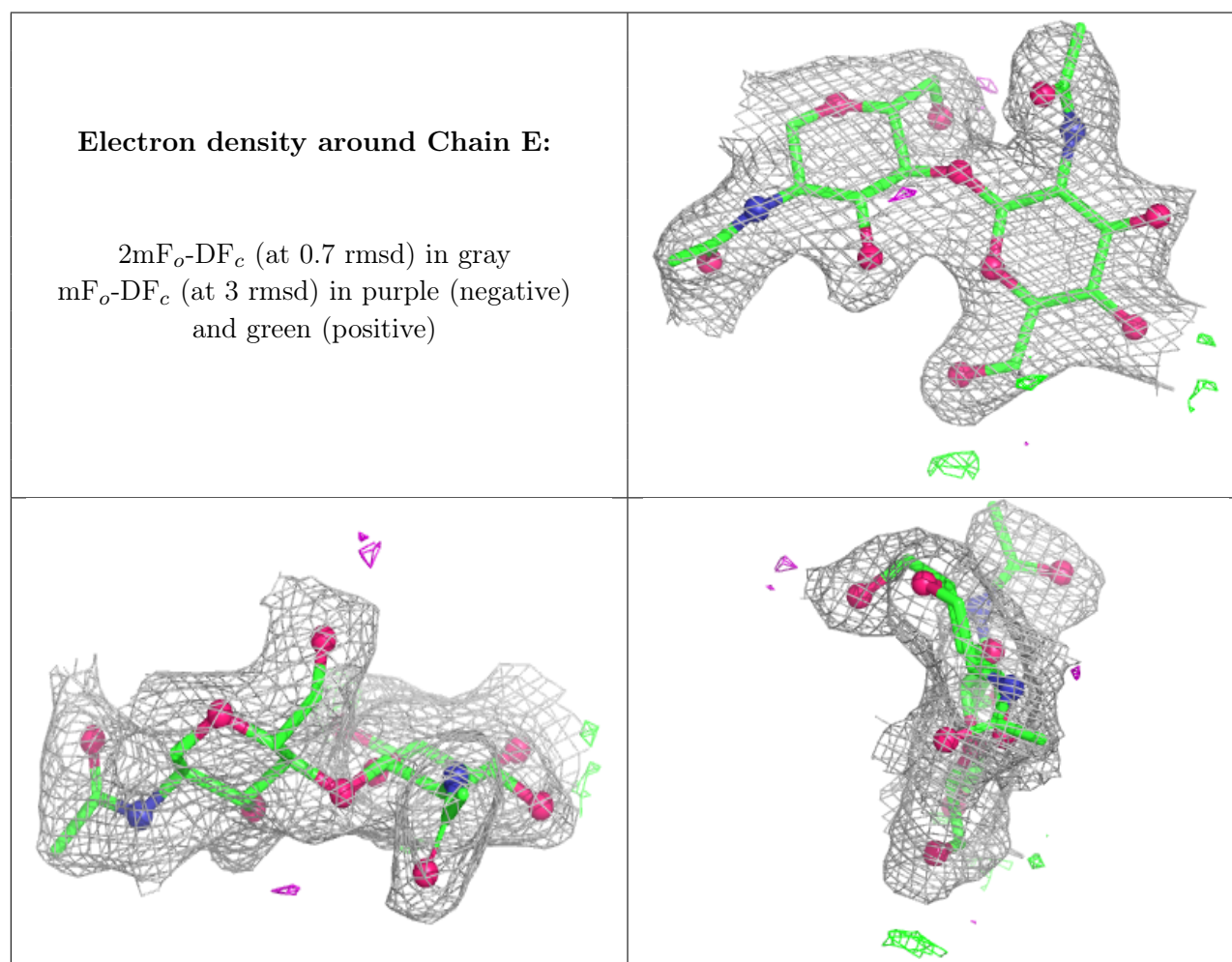
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	H	2	14/15	0.82	0.14	49,56,59,60	0
2	NAG	F	2	14/15	0.83	0.13	43,50,57,58	0
2	NAG	G	2	14/15	0.84	0.12	45,50,53,57	0

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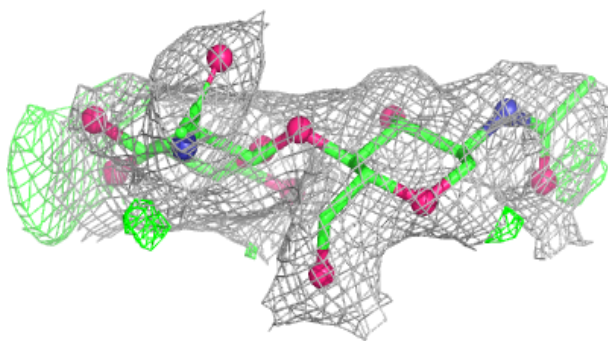
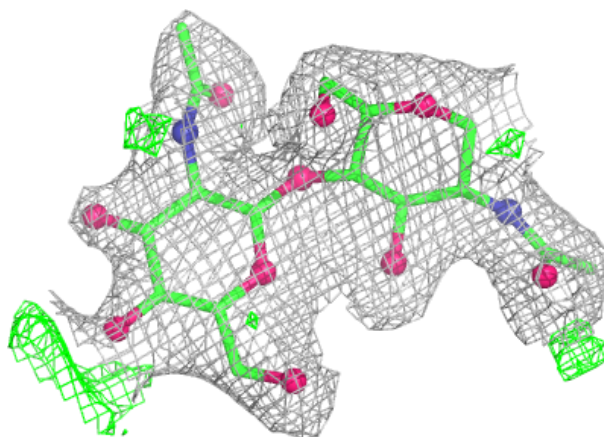
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	E	2	14/15	0.87	0.10	39,43,50,50	0
2	NAG	F	1	14/15	0.93	0.09	22,31,35,42	0
2	NAG	G	1	14/15	0.94	0.08	22,30,34,41	0
2	NAG	H	1	14/15	0.95	0.07	26,35,40,44	0
2	NAG	E	1	14/15	0.96	0.07	25,30,35,36	0

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.



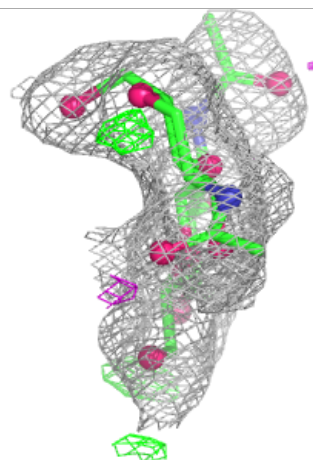
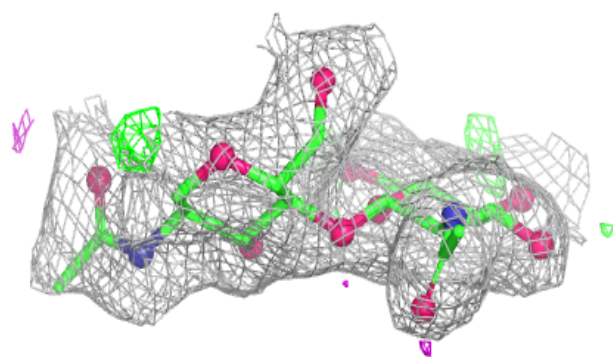
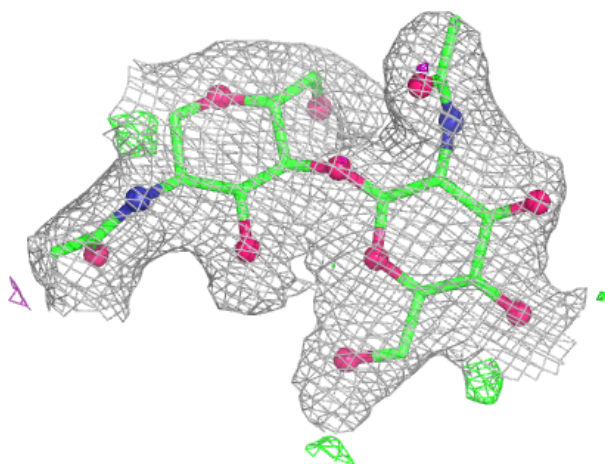
Electron density around Chain F:

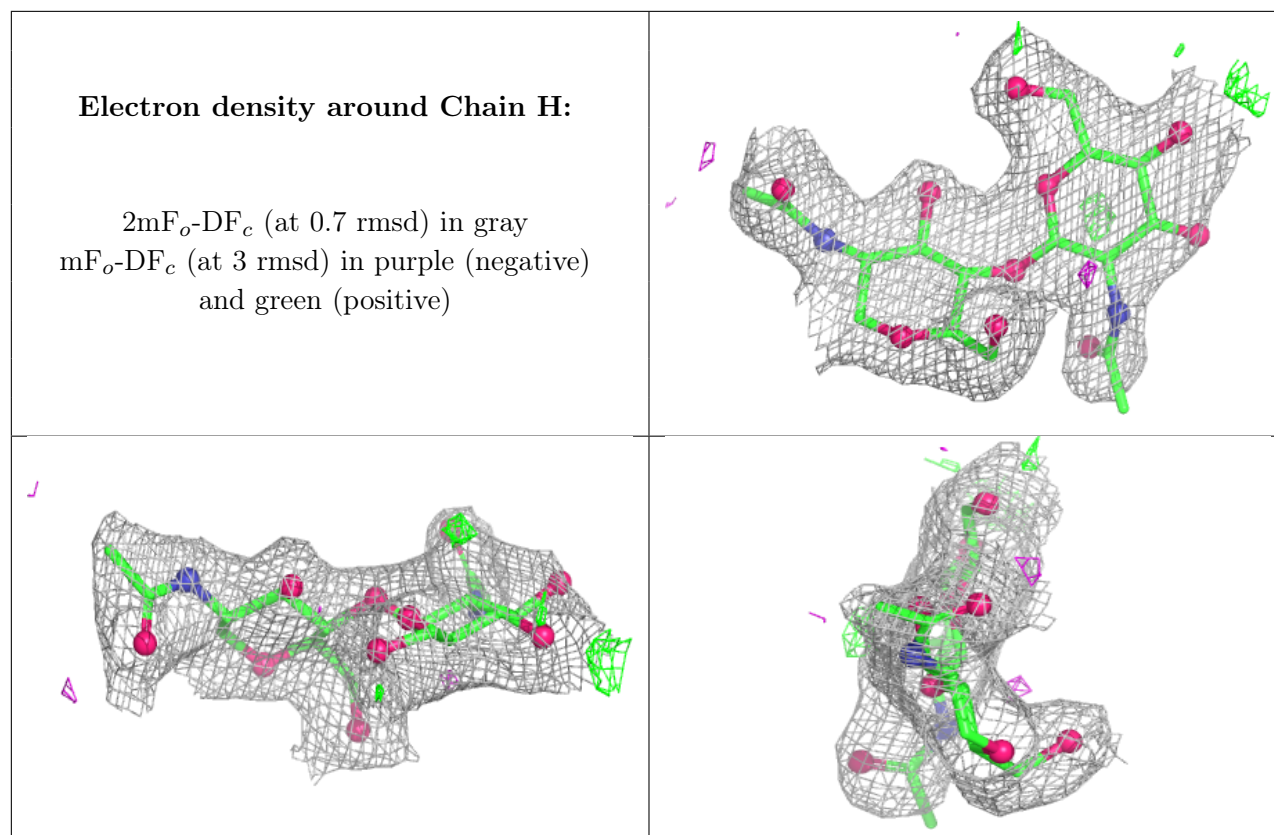
$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



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)





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
6	BOG	D	707	20/20	0.72	0.18	55,68,74,75	0
4	NAG	D	702	14/15	0.75	0.17	47,58,62,62	0
4	NAG	C	702	14/15	0.77	0.15	46,55,59,60	0
6	BOG	C	707	20/20	0.78	0.18	55,65,69,70	0
4	NAG	D	705	14/15	0.78	0.14	45,54,61,64	0
6	BOG	A	707	20/20	0.80	0.14	52,58,64,64	0
4	NAG	B	702	14/15	0.80	0.13	48,54,62,64	0
4	NAG	A	705	14/15	0.80	0.15	43,48,63,63	0
4	NAG	A	702	14/15	0.82	0.14	45,50,60,60	0
4	NAG	B	705	14/15	0.86	0.11	38,48,51,55	0
5	DXS	D	706	44/44	0.88	0.13	25,40,63,67	0
6	BOG	D	708	20/20	0.89	0.12	41,44,46,46	0
5	DXS	B	706	44/44	0.90	0.14	28,39,64,66	0
5	DXS	A	706	44/44	0.90	0.13	21,38,60,63	0

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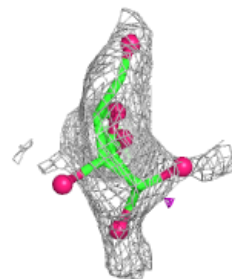
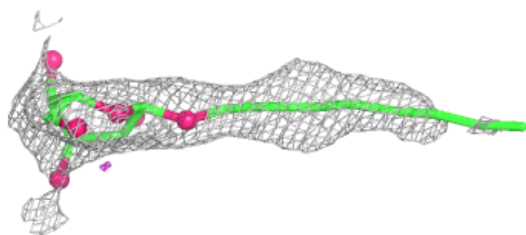
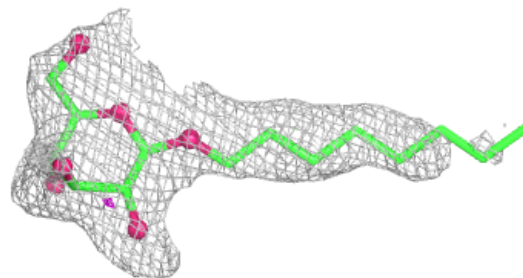
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	NAG	C	705	14/15	0.92	0.08	32,40,45,51	0
5	DXS	C	706	44/44	0.92	0.12	24,39,61,63	0
6	BOG	A	708	20/20	0.92	0.11	23,28,35,35	20
3	HEM	D	701	43/43	0.96	0.09	22,31,48,59	0
3	HEM	A	701	43/43	0.97	0.08	25,28,44,57	0
3	HEM	B	701	43/43	0.97	0.09	20,28,43,50	0
3	HEM	C	701	43/43	0.97	0.09	19,27,48,57	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.

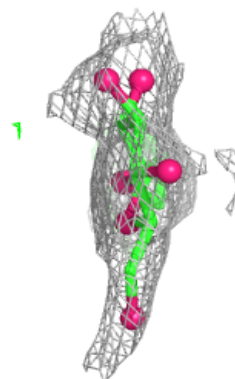
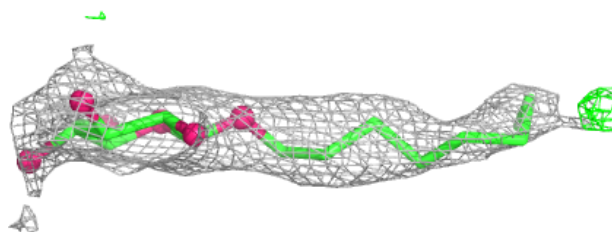
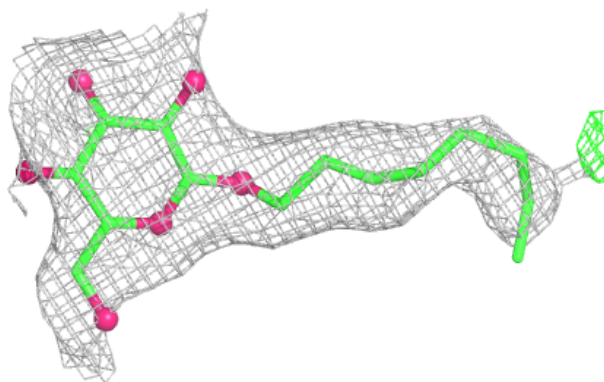
Electron density around BOG D 707:

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

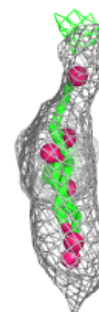
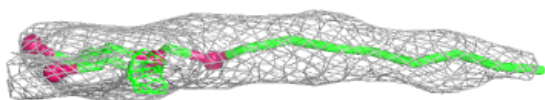
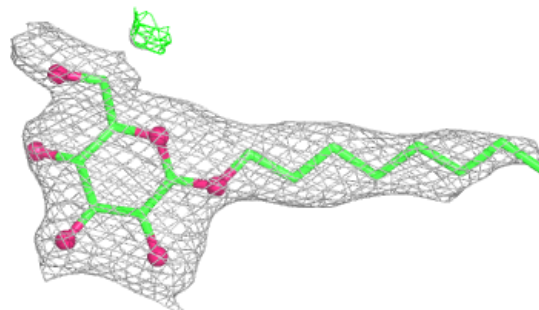


Electron density around BOG C 707:

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

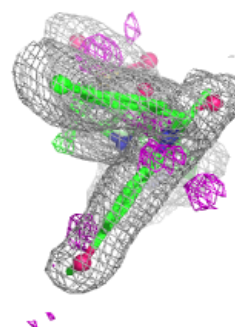
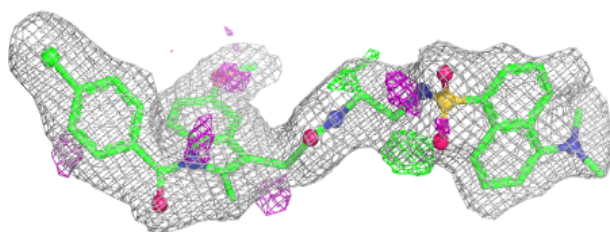
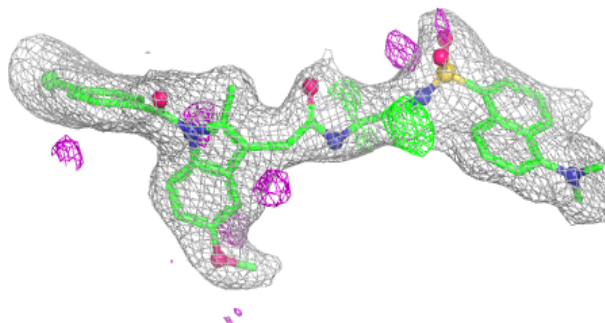
**Electron density around BOG A 707:**

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

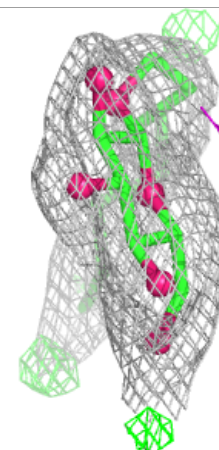
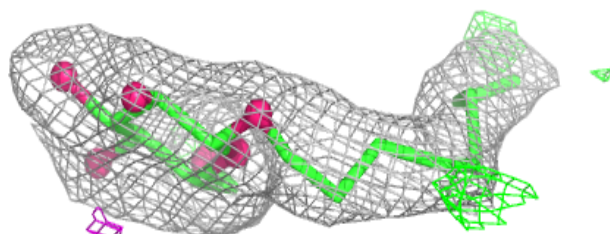
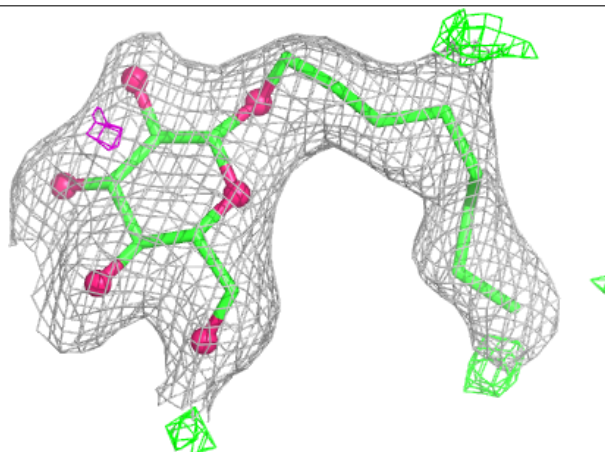


Electron density around DXS D 706:

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

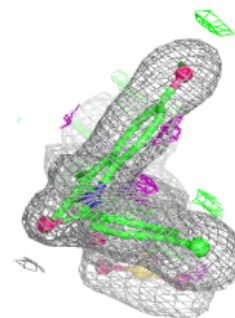
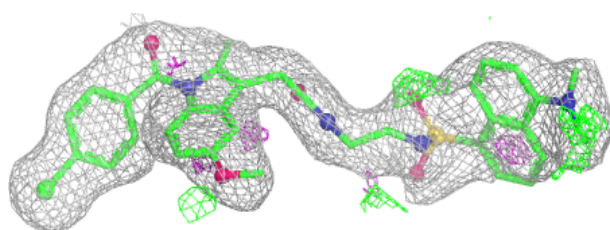
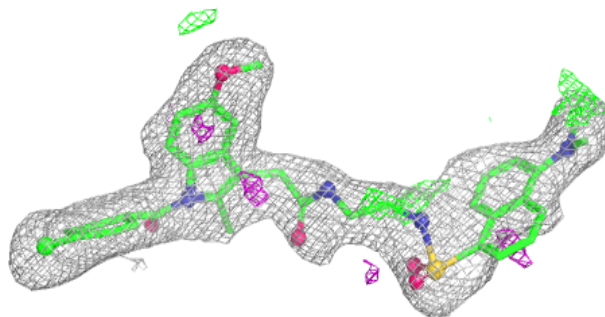
**Electron density around BOG D 708:**

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

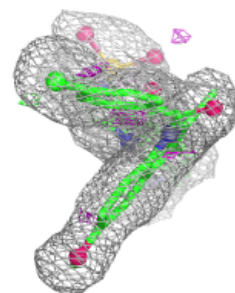
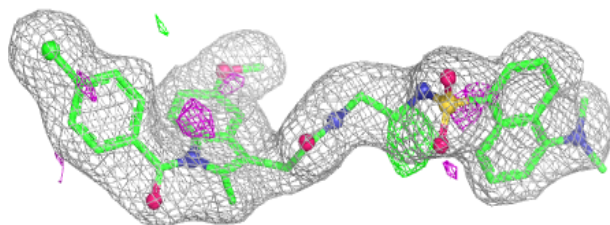
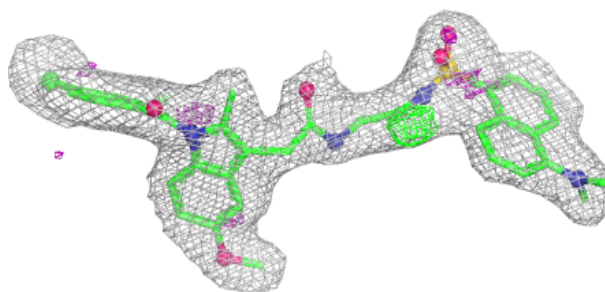


Electron density around DXS B 706:

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

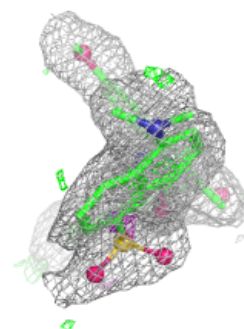
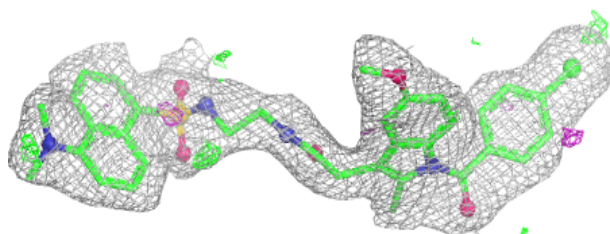
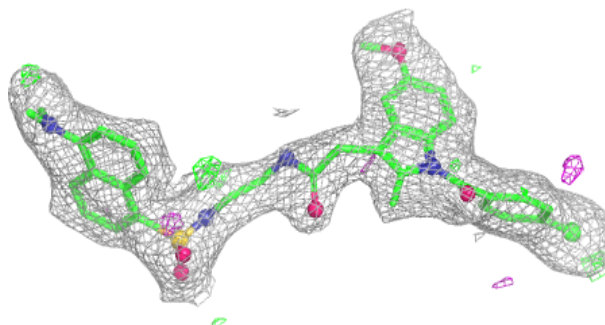
**Electron density around DXS A 706:**

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

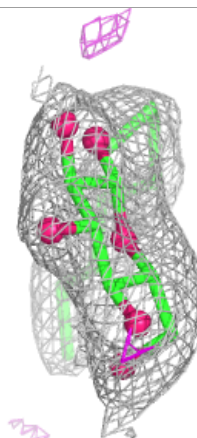
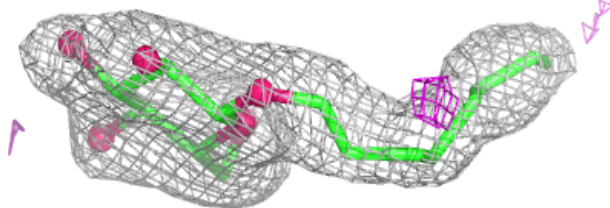
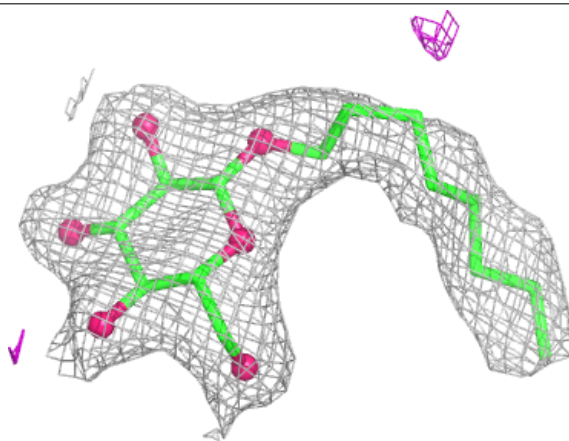


Electron density around DXS C 706:

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

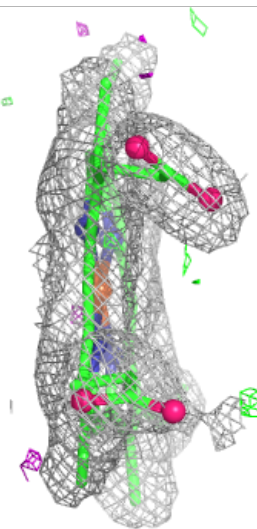
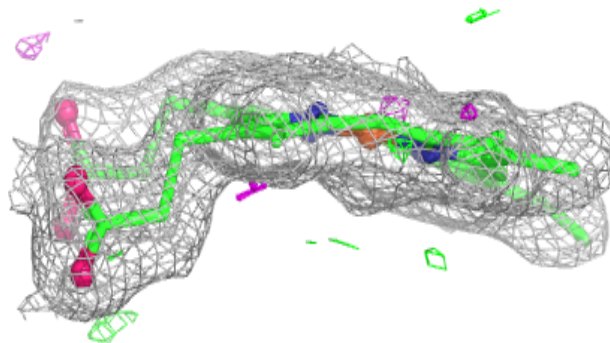
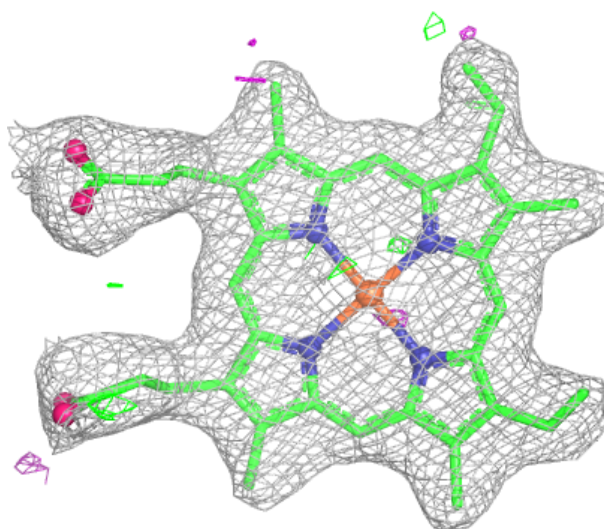
**Electron density around BOG A 708:**

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



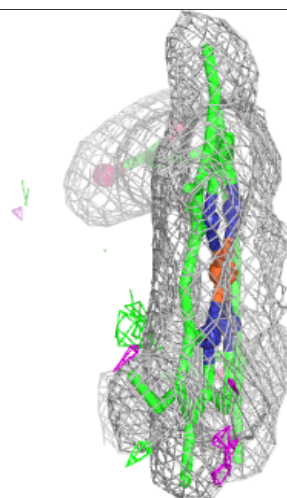
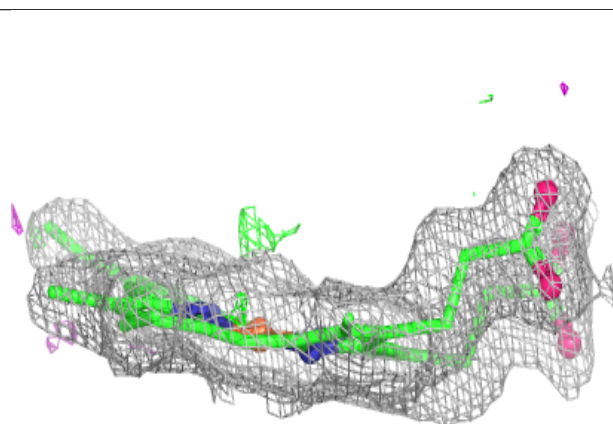
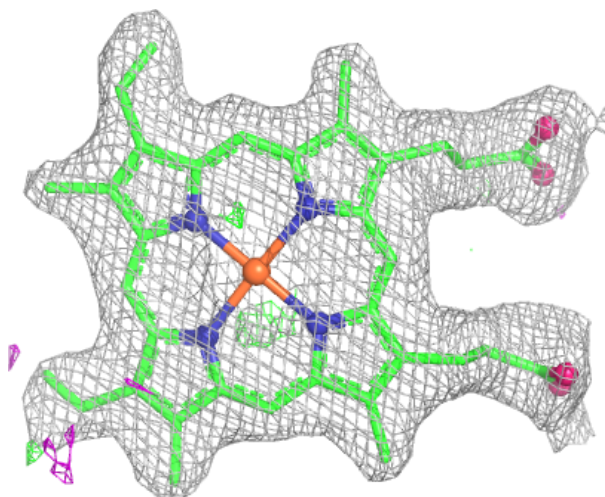
Electron density around HEM D 701:

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



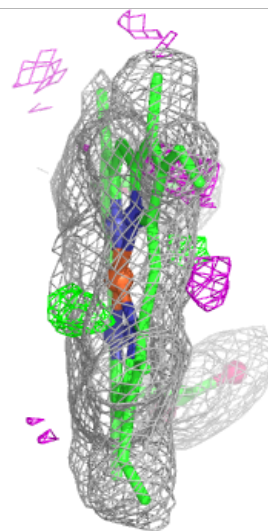
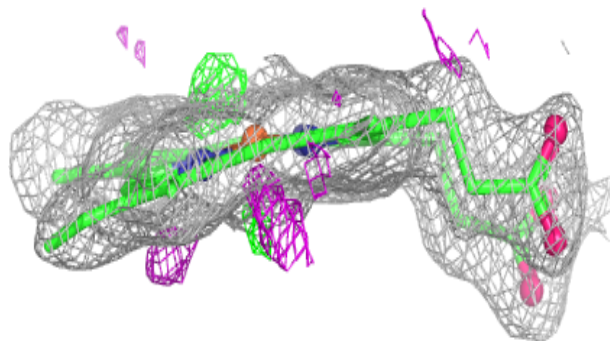
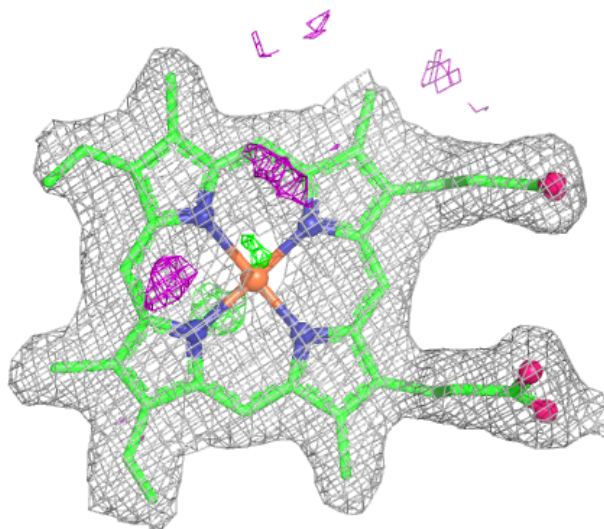
Electron density around HEM A 701:

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



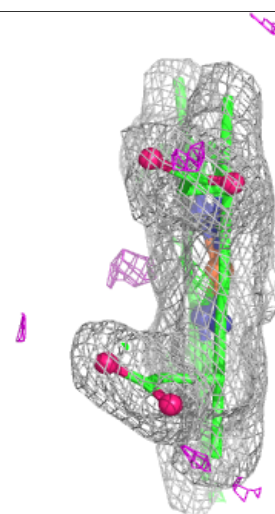
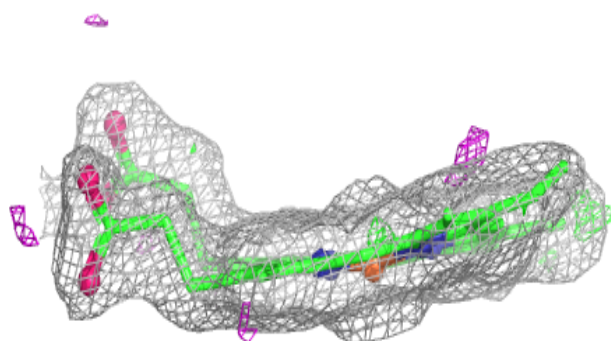
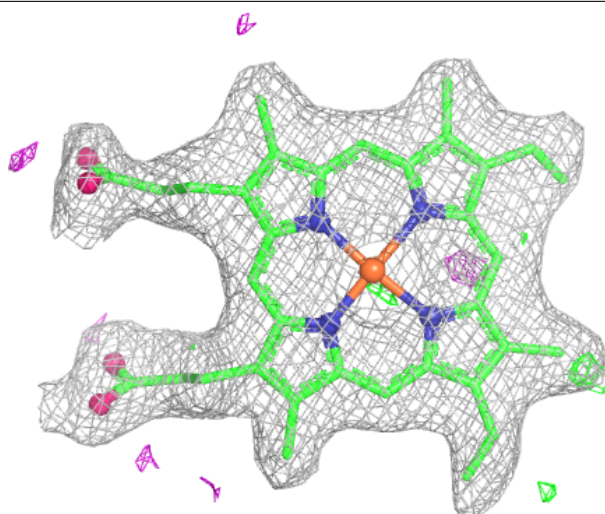
Electron density around HEM B 701:

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



Electron density around HEM 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.