



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 20, 2020 – 04:36 PM EDT

PDB ID : 6PP4
Title : Structure of human endothelial nitric oxide synthase heme domain in complex with 7-(3-(Aminomethyl)-4-(pyridin-3-ylmethoxy)phenyl)-4-methylquinolin-2-amine
Authors : Chreifi, G.; Li, H.; Poulos, T.L.
Deposited on : 2019-07-05
Resolution : 2.20 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.10.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.10.1

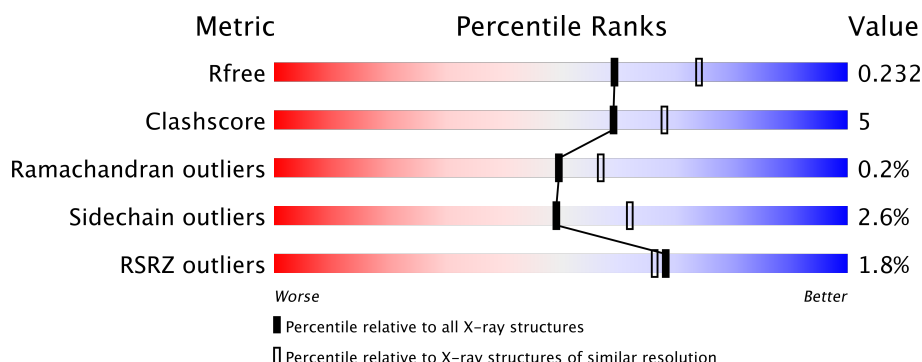
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	111664	4343 (2.20-2.20)
Clashscore	122126	5027 (2.20-2.20)
Ramachandran outliers	120053	4952 (2.20-2.20)
Sidechain outliers	120020	4953 (2.20-2.20)
RSRZ outliers	108989	4245 (2.20-2.20)

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 on 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	440	<div> <div>5%</div> <div> <div></div> <div>79%</div> <div>12%</div> <div>8%</div> </div> </div>
1	B	440	<div> <div>81%</div> <div>10%</div> <div>9%</div> </div>
1	C	440	<div> <div>79%</div> <div>12%</div> <div>9%</div> </div>
1	D	440	<div> <div>%</div> <div>81%</div> <div>10%</div> <div>8%</div> </div>

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 13926 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.

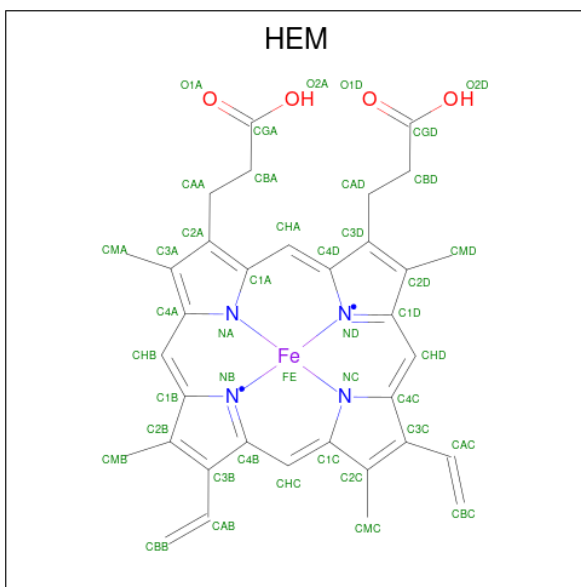
- Molecule 1 is a protein called Nitric oxide synthase, endothelial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	404	Total	C	N	O	S	0	2	0
			3237	2062	570	589	16			
1	B	402	Total	C	N	O	S	0	3	0
			3221	2051	566	587	17			
1	C	401	Total	C	N	O	S	0	2	0
			3209	2044	563	586	16			
1	D	403	Total	C	N	O	S	0	2	0
			3226	2054	569	587	16			

There are 4 discrepancies between the modelled and reference sequences:

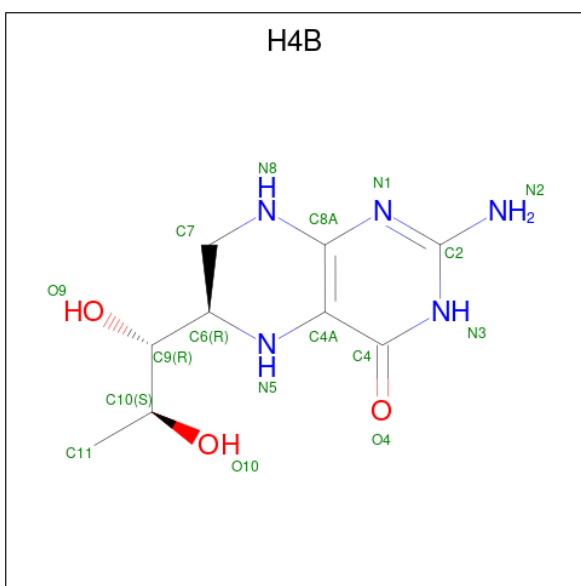
Chain	Residue	Modelled	Actual	Comment	Reference
A	298	GLU	ASP	variant	UNP P29474
B	298	GLU	ASP	variant	UNP P29474
C	298	GLU	ASP	variant	UNP P29474
D	298	GLU	ASP	variant	UNP P29474

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



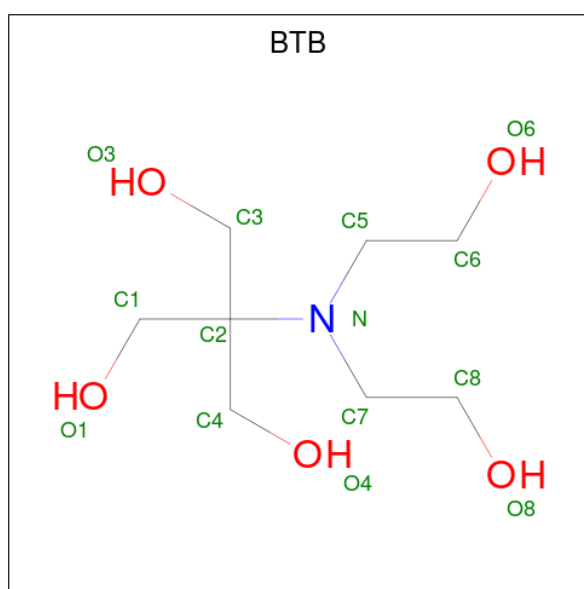
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 3 is 5,6,7,8-TETRAHYDROBIOPTERIN (three-letter code: H4B) (formula: $C_9H_{15}N_5O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			17	9	5	3		
3	B	1	Total	C	N	O	0	0
			17	9	5	3		
3	C	1	Total	C	N	O	0	0
			17	9	5	3		
3	D	1	Total	C	N	O	0	0
			17	9	5	3		

- Molecule 4 is 2-[BIS-(2-HYDROXY-ETHYL)-AMINO]-2-HYDROXYMETHYL-PROPAN E-1,3-DIOL (three-letter code: BTB) (formula: C₈H₁₉NO₅).



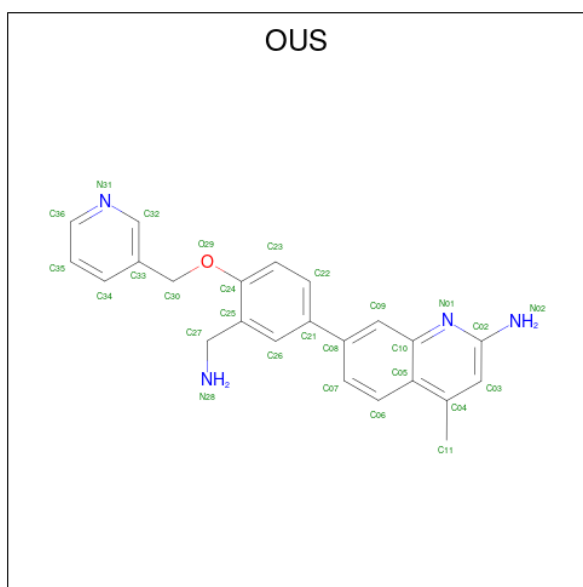
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		
4	C	1	Total	C	N	O	0	0
			14	8	1	5		
4	D	1	Total	C	N	O	0	0
			14	8	1	5		

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

- Molecule 5 is 7-{3-(aminomethyl)-4-[(pyridin-3-yl)methoxy]phenyl}-4-methylquinolin-2-amine (three-letter code: OUS) (formula: C₂₃H₂₂N₄O) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			28	23	4	1		
5	B	1	Total	C	N	O	0	0
			28	23	4	1		
5	C	1	Total	C	N	O	0	0
			28	23	4	1		
5	D	1	Total	C	N	O	0	0
			28	23	4	1		

- Molecule 6 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Zn	0	0
			1	1		
6	C	1	Total	Zn	0	0
			1	1		

- Molecule 7 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 8 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	B	1	Total	Cl	0	0
			1	1		
8	A	1	Total	Cl	0	0
			1	1		
8	D	1	Total	Cl	0	0
			1	1		
8	C	1	Total	Cl	0	0
			1	1		

- Molecule 9 is GADOLINIUM ATOM (three-letter code: GD) (formula: Gd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	B	1	Total	Gd	0	0
			1	1		
9	D	1	Total	Gd	0	0
			1	1		
9	C	1	Total	Gd	0	0
			1	1		

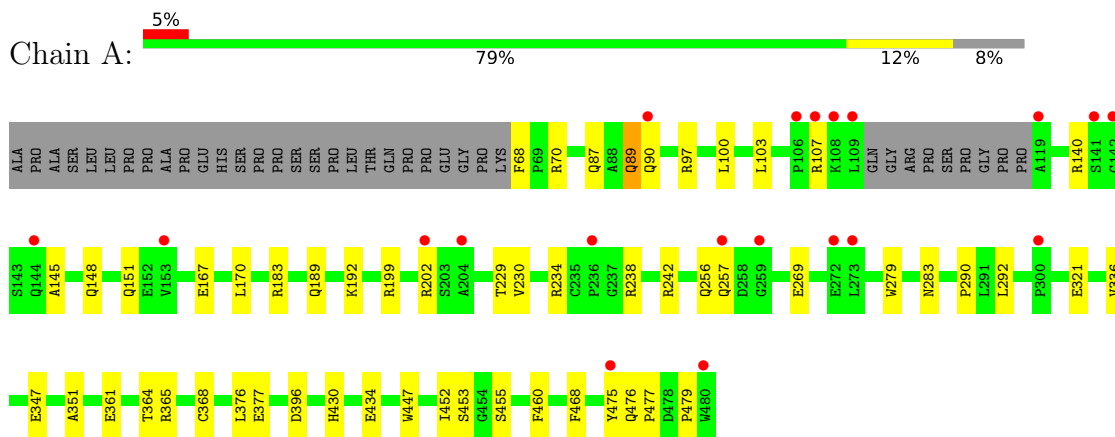
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	91	Total 91	O 91	0	0
10	B	156	Total 156	O 156	0	0
10	C	141	Total 141	O 141	0	0
10	D	160	Total 160	O 160	0	0

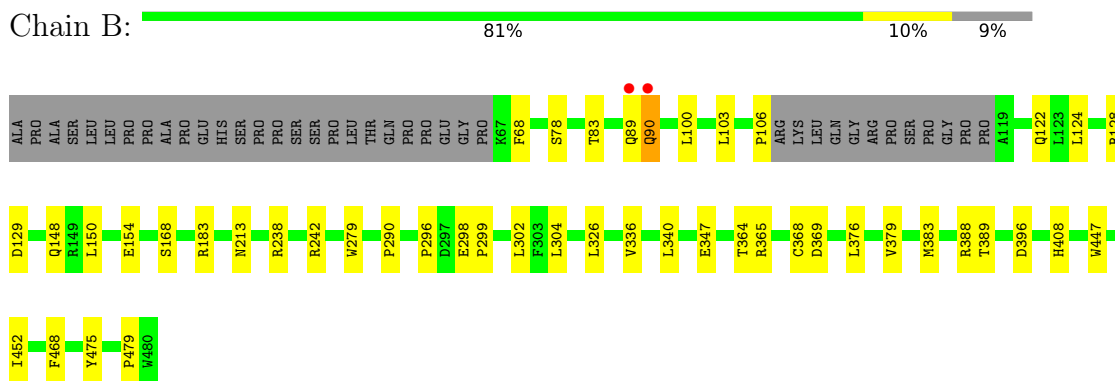
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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.

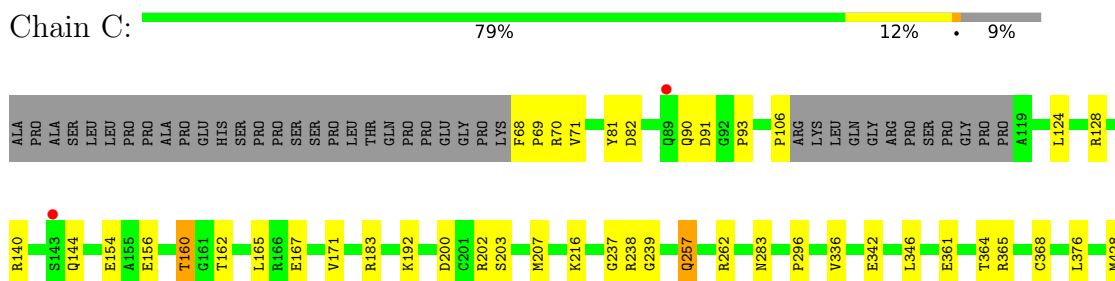
- Molecule 1: Nitric oxide synthase, endothelial



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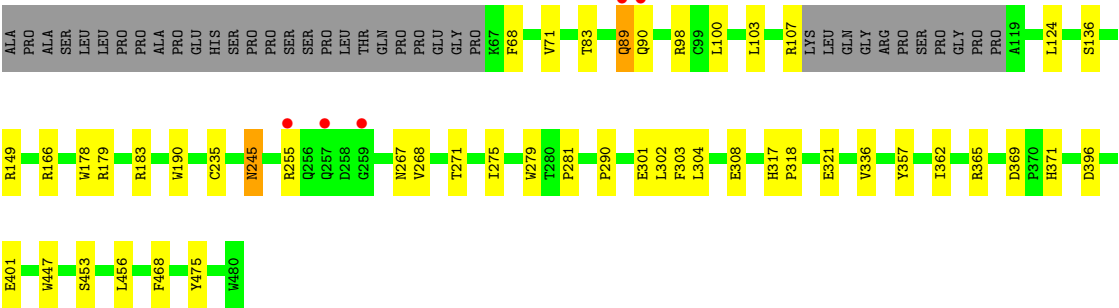
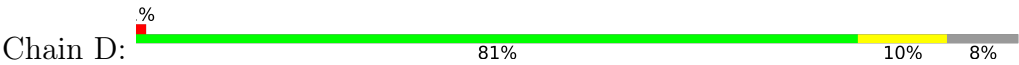


- Molecule 1: Nitric oxide synthase, endothelial





● Molecule 1: Nitric oxide synthase, endothelial



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.17Å 152.48Å 109.14Å 90.00° 90.86° 90.00°	Depositor
Resolution (Å)	43.15 – 2.20 88.74 – 2.20	Depositor EDS
% Data completeness (in resolution range)	95.5 (43.15-2.20) 96.5 (88.74-2.20)	Depositor EDS
R_{merge}	0.19	Depositor
R_{sym}	0.19	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.78 (at 2.20Å)	Xtriage
Refinement program	PHENIX (1.11.1-2575_1496: ???)	Depositor
R, R_{free}	0.187 , 0.238 0.179 , 0.232	Depositor DCC
R_{free} test set	4725 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	33.4	Xtriage
Anisotropy	0.793	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 63.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.097 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13926	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.84% 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: GOL, ZN, H4B, CL, GD, BTB, OUS, 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	0.35	0/3335	0.52	0/4543
1	B	0.41	0/3319	0.55	0/4523
1	C	0.39	0/3307	0.54	0/4507
1	D	0.44	0/3324	0.58	1/4528 (0.0%)
All	All	0.40	0/13285	0.55	1/18101 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	D	179	ARG	NE-CZ-NH2	-5.17	117.71	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	89	GLN	Peptide

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	3237	0	3146	35	0
1	B	3221	0	3126	28	0
1	C	3209	0	3109	29	0
1	D	3226	0	3134	26	0
2	A	43	0	30	4	0
2	B	43	0	30	3	0
2	C	43	0	30	4	0
2	D	43	0	30	2	0
3	A	17	0	15	1	0
3	B	17	0	15	0	0
3	C	17	0	15	1	0
3	D	17	0	15	0	0
4	A	28	0	38	3	0
4	B	42	0	54	6	0
4	C	14	0	19	0	0
4	D	28	0	36	4	0
5	A	28	0	0	2	0
5	B	28	0	0	2	0
5	C	28	0	0	3	0
5	D	28	0	0	1	0
6	A	1	0	0	0	0
6	C	1	0	0	0	0
7	A	6	0	8	1	0
7	C	6	0	8	0	0
8	A	1	0	0	0	0
8	B	1	0	0	0	0
8	C	1	0	0	0	0
8	D	1	0	0	0	0
9	B	1	0	0	0	0
9	C	1	0	0	0	0
9	D	1	0	0	0	0
10	A	91	0	0	4	0
10	B	156	0	0	3	0
10	C	141	0	0	5	0
10	D	160	0	0	1	0
All	All	13926	0	12858	134	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 5.

All (134) 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:C:183:ARG:HB2	2:C:502:HEM:HBD2	1.56	0.86
1:D:275:ILE:HD12	1:D:281:PRO:HG3	1.65	0.78
1:A:242:ARG:NH2	1:A:477:PRO:O	2.18	0.77
1:B:365:ARG:NH2	1:B:369:ASP:OD2	2.19	0.74
1:C:90:GLN:NE2	10:C:603:HOH:O	2.23	0.72
1:D:279:TRP:HB2	1:D:302:LEU:HD21	1.72	0.71
1:B:408:HIS:ND1	10:B:602:HOH:O	2.24	0.69
1:A:242:ARG:HD3	1:A:479:PRO:HB3	1.76	0.68
1:D:475:TYR:OH	2:D:501:HEM:O2D	2.11	0.67
1:D:365:ARG:NH2	1:D:369:ASP:OD2	2.28	0.67
1:B:475:TYR:OH	2:B:501:HEM:O2D	2.12	0.66
4:B:504:BTB:O4	4:B:504:BTB:O3	2.12	0.66
1:C:144:GLN:NE2	10:C:604:HOH:O	2.30	0.64
1:B:100:LEU:HB3	1:B:103:LEU:HD22	1.79	0.63
2:D:501:HEM:HHC	2:D:501:HEM:HBB2	1.79	0.63
1:A:70:ARG:NH1	10:A:607:HOH:O	2.30	0.63
1:A:183:ARG:HD3	1:A:447:TRP:CD2	2.36	0.61
1:B:279:TRP:HB2	1:B:302:LEU:HD21	1.82	0.60
1:D:90:GLN:HB3	1:D:468:PHE:CD1	2.37	0.60
1:D:336:VAL:HG21	5:D:503:OUS:C07	2.32	0.60
2:C:502:HEM:HBB2	2:C:502:HEM:HHC	1.84	0.59
1:C:336:VAL:HG21	5:C:504:OUS:C07	2.32	0.59
2:A:501:HEM:HBC2	2:A:501:HEM:HMC2	1.84	0.59
1:A:183:ARG:HB2	2:A:501:HEM:HBD2	1.84	0.58
2:B:501:HEM:HBB2	2:B:501:HEM:HHC	1.83	0.58
1:A:364:THR:O	1:A:368:CYS:HB2	2.03	0.58
1:A:377:GLU:HB2	10:A:636:HOH:O	2.03	0.58
1:C:262:ARG:NH1	1:C:283:ASN:O	2.35	0.57
1:A:238:ARG:NH1	10:A:606:HOH:O	2.30	0.56
2:C:502:HEM:O1A	10:C:601:HOH:O	2.18	0.56
2:A:501:HEM:HBB2	2:A:501:HEM:HHC	1.86	0.56
1:D:267:ASN:O	1:D:271:THR:OG1	2.18	0.56
1:C:70:ARG:HB2	1:C:81:TYR:CE2	2.40	0.56
1:D:290:PRO:HB3	1:D:304:LEU:HD23	1.88	0.56
1:C:365:ARG:HH12	3:C:503:H4B:C4	2.19	0.55
1:A:145:ALA:HA	1:A:148:GLN:HB3	1.88	0.55
1:B:106:PRO:O	10:B:601:HOH:O	2.18	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:336:VAL:HG21	5:B:503:OUS:C07	2.37	0.54
1:D:100:LEU:HB3	1:D:103:LEU:HD22	1.89	0.54
10:C:713:HOH:O	1:D:371:HIS:HB3	2.07	0.54
1:A:183:ARG:NH1	1:A:475:TYR:OH	2.42	0.53
1:A:242:ARG:NH1	1:A:476:GLN:OE1	2.40	0.53
1:C:453:SER:HB3	1:C:456:LEU:HD12	1.91	0.53
1:A:242:ARG:HH22	1:A:476:GLN:HG3	1.75	0.52
1:B:150:LEU:O	1:B:154:GLU:HG2	2.10	0.52
1:A:199:ARG:O	1:A:202:ARG:NH1	2.28	0.51
1:B:90:GLN:HB2	1:B:468:PHE:CD1	2.46	0.51
1:D:271:THR:O	1:D:275:ILE:HG12	2.10	0.51
1:D:317:HIS:CG	1:D:318:PRO:HD2	2.46	0.51
1:B:238:ARG:HG2	1:B:296:PRO:HB3	1.93	0.51
2:B:501:HEM:HBD1	5:B:503:OUS:C22	2.40	0.50
1:B:89:GLN:NE2	1:B:129:ASP:OD2	2.44	0.49
1:A:321:GLU:CD	1:A:321:GLU:H	2.15	0.49
1:C:428:MET:HG3	1:C:458:PRO:HB2	1.94	0.49
1:A:336:VAL:HG21	5:A:504:OUS:C07	2.43	0.49
1:B:340:LEU:HD21	1:B:347:GLU:HG2	1.94	0.48
1:C:156:GLU:O	1:C:160:THR:HG22	2.14	0.48
1:A:361:GLU:OE2	5:A:504:OUS:N02	2.47	0.48
1:A:68:PHE:N	10:A:617:HOH:O	2.47	0.48
1:A:170:LEU:HD11	1:A:230:VAL:HG11	1.96	0.48
1:A:368:CYS:SG	1:A:376:LEU:HD13	2.54	0.48
1:B:298:GLU:HG3	1:B:299:PRO:HD2	1.97	0.47
1:C:450:PRO:HG2	1:C:457:THR:HG21	1.97	0.47
1:C:342:GLU:OE2	1:C:470:SER:OG	2.33	0.47
1:D:317:HIS:NE2	1:D:401:GLU:OE1	2.44	0.47
4:B:508:BTB:H52	4:B:508:BTB:H82	1.41	0.47
1:C:165:LEU:HG	1:C:346:LEU:HD12	1.98	0.46
1:A:229:THR:O	1:A:351:ALA:HA	2.15	0.46
4:A:503:BTB:H82	1:C:257:GLN:HE22	1.81	0.46
1:B:68:PHE:CD2	1:B:83:THR:HG22	2.49	0.46
4:D:505:BTB:H52	4:D:505:BTB:H81	1.44	0.46
1:C:368:CYS:SG	1:C:376:LEU:HD13	2.56	0.46
1:D:357:TYR:CD2	1:D:362:ILE:HD11	2.51	0.46
1:D:149:ARG:HD3	1:D:166:ARG:CZ	2.46	0.45
1:D:453:SER:HB3	1:D:456:LEU:HD12	1.99	0.45
1:B:242:ARG:NH2	1:B:479:PRO:HD3	2.32	0.45
1:B:290:PRO:HB3	1:B:304:LEU:HD23	1.98	0.45
1:D:255:ARG:HB2	1:D:268:VAL:CG1	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:505:BTB:H72	4:A:505:BTB:H12	1.70	0.45
4:B:508:BTB:O8	4:B:508:BTB:H41	2.18	0.44
1:C:455:SER:HA	1:C:460:PHE:CG	2.53	0.44
1:B:298:GLU:OE2	4:B:505:BTB:N	2.51	0.44
1:C:183:ARG:HD3	1:C:447:TRP:CD2	2.53	0.44
4:D:505:BTB:H11	4:D:505:BTB:H71	1.59	0.44
4:D:505:BTB:H51	4:D:505:BTB:H32	1.48	0.44
1:A:234:ARG:NH1	1:A:347:GLU:OE1	2.51	0.44
1:A:455:SER:HA	1:A:460:PHE:CG	2.52	0.44
1:D:371:HIS:CE1	10:D:609:HOH:O	2.71	0.44
1:B:379:VAL:O	1:B:383:MET:HG3	2.18	0.43
1:D:245:ASN:N	1:D:245:ASN:OD1	2.51	0.43
1:B:298:GLU:CD	4:B:505:BTB:H31	2.38	0.43
4:A:503:BTB:H82	4:A:503:BTB:O4	2.18	0.43
1:C:364:THR:O	1:C:368:CYS:HB2	2.19	0.43
1:D:178:TRP:CE3	1:D:190:TRP:HA	2.54	0.43
1:A:151:GLN:OE1	1:A:151:GLN:N	2.52	0.43
1:A:292:LEU:HA	1:A:292:LEU:HD23	1.86	0.43
1:A:87:GLN:O	1:A:89:GLN:NE2	2.52	0.42
1:B:213:ASN:HB2	10:B:717:HOH:O	2.19	0.42
1:C:445:TRP:CE2	1:C:449:VAL:HG21	2.54	0.42
1:A:279:TRP:CG	1:A:290:PRO:HG3	2.55	0.42
1:C:167:GLU:O	1:C:171:VAL:HG23	2.19	0.42
1:C:239:GLY:HA3	1:C:296:PRO:HB3	2.01	0.42
1:B:124:LEU:HD23	1:B:124:LEU:HA	1.79	0.42
1:B:68:PHE:CG	1:B:83:THR:HG22	2.55	0.42
1:D:255:ARG:HB2	1:D:268:VAL:HG11	2.01	0.42
1:B:388:ARG:HE	1:B:388:ARG:HB2	1.68	0.42
1:C:361:GLU:OE2	5:C:504:OUS:N02	2.51	0.42
1:C:69:PRO:HD2	1:C:82:ASP:O	2.20	0.42
1:D:321:GLU:OE2	4:D:504:BTB:O4	2.37	0.42
1:D:301:GLU:HB3	1:D:303:PHE:HE1	1.84	0.42
1:D:183:ARG:HD3	1:D:447:TRP:CD2	2.55	0.42
1:A:365:ARG:HH12	3:A:502:H4B:C4	2.32	0.42
1:C:200:ASP:OD1	1:C:200:ASP:N	2.41	0.42
1:A:364:THR:HG21	1:A:452:ILE:HG23	2.02	0.41
1:C:160:THR:HG23	1:C:162:THR:H	1.84	0.41
1:A:167:GLU:OE1	7:A:507:GOL:H12	2.20	0.41
1:B:364:THR:O	1:B:368:CYS:HB2	2.21	0.41
1:A:90:GLN:HG3	1:A:468:PHE:CE1	2.55	0.41
1:B:368:CYS:SG	1:B:376:LEU:HD13	2.59	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:183:ARG:HB2	2:A:501:HEM:CBD	2.48	0.41
1:C:124:LEU:O	1:C:128:ARG:HG3	2.21	0.41
1:B:183:ARG:HD3	1:B:447:TRP:CD2	2.56	0.41
1:C:91:ASP:OD1	10:C:602:HOH:O	2.22	0.41
1:A:453:SER:HA	1:B:452:ILE:HG22	2.02	0.41
1:D:68:PHE:CD2	1:D:83:THR:HG22	2.55	0.40
1:A:189:GLN:OE1	1:A:192:LYS:HE2	2.21	0.40
2:C:502:HEM:O2A	5:C:504:OUS:N28	2.54	0.40
1:A:100:LEU:HB3	1:A:103:LEU:HD22	2.04	0.40
1:B:124:LEU:HD13	1:B:128:ARG:HH21	1.85	0.40
4:B:508:BTB:H12	4:B:508:BTB:H51	1.77	0.40
1:A:430:HIS:CE1	1:A:434:GLU:HG3	2.56	0.40
1:C:93:PRO:HG3	1:C:106:PRO:HB3	2.03	0.40
1:C:68:PHE:HA	1:C:69:PRO:HD3	1.95	0.40
1:D:308:GLU:OE1	1:D:308:GLU:N	2.35	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	402/440 (91%)	377 (94%)	23 (6%)	2 (0%)	31	33
1	B	401/440 (91%)	393 (98%)	8 (2%)	0	100	100
1	C	399/440 (91%)	388 (97%)	10 (2%)	1 (0%)	43	48
1	D	401/440 (91%)	390 (97%)	11 (3%)	0	100	100
All	All	1603/1760 (91%)	1548 (97%)	52 (3%)	3 (0%)	49	57

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	107	ARG
1	A	283	ASN
1	C	237	GLY

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	345/373 (92%)	338 (98%)	7 (2%)	58	72
1	B	344/373 (92%)	335 (97%)	9 (3%)	49	62
1	C	342/373 (92%)	330 (96%)	12 (4%)	39	49
1	D	344/373 (92%)	335 (97%)	9 (3%)	49	62
All	All	1375/1492 (92%)	1338 (97%)	37 (3%)	49	60

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	89	GLN
1	A	97	ARG
1	A	140	ARG
1	A	256	GLN
1	A	257	GLN
1	A	269	GLU
1	A	396	ASP
1	B	78	SER
1	B	90	GLN
1	B	122	GLN
1	B	148	GLN
1	B	168[A]	SER
1	B	168[B]	SER
1	B	326	LEU
1	B	389	THR
1	B	396	ASP
1	C	71	VAL
1	C	140	ARG
1	C	154	GLU

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Mol	Chain	Res	Type
1	C	160	THR
1	C	192	LYS
1	C	202	ARG
1	C	203	SER
1	C	207	MET
1	C	216	LYS
1	C	238	ARG
1	C	257	GLN
1	C	470	SER
1	D	71	VAL
1	D	89	GLN
1	D	98	ARG
1	D	107	ARG
1	D	124	LEU
1	D	136	SER
1	D	235	CYS
1	D	245	ASN
1	D	396	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 31 ligands modelled in this entry, 9 are monoatomic - leaving 22 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	BTB	C	505	-	13,13,13	0.33	0	7,16,16	0.36	0
3	H4B	A	502	-	16,18,18	0.95	0	11,26,26	2.75	6 (54%)
4	BTB	A	505	-	13,13,13	0.54	0	7,16,16	0.83	0
3	H4B	B	502	-	16,18,18	0.87	0	11,26,26	2.80	7 (63%)
3	H4B	D	502	-	16,18,18	0.74	0	11,26,26	2.73	6 (54%)
7	GOL	A	507	-	5,5,5	0.33	0	5,5,5	0.31	0
4	BTB	B	505	-	13,13,13	0.44	0	7,16,16	0.63	0
5	OUS	B	503	-	31,31,31	0.84	0	42,43,43	1.50	7 (16%)
7	GOL	C	507	-	5,5,5	0.49	0	5,5,5	0.39	0
4	BTB	D	504	9	13,13,13	0.48	0	7,16,16	0.50	0
4	BTB	B	504	9	13,13,13	0.48	0	7,16,16	0.25	0
3	H4B	C	503	-	16,18,18	0.95	0	11,26,26	2.74	6 (54%)
4	BTB	B	508	-	13,13,13	1.00	2 (15%)	7,16,16	1.49	1 (14%)
2	HEM	C	502	1	27,50,50	1.94	6 (22%)	17,82,82	2.04	4 (23%)
5	OUS	C	504	-	31,31,31	0.84	0	42,43,43	1.29	4 (9%)
2	HEM	D	501	1	27,50,50	1.78	4 (14%)	17,82,82	2.39	6 (35%)
4	BTB	D	505	-	13,13,13	0.44	0	7,16,16	0.88	0
5	OUS	D	503	-	31,31,31	0.91	0	42,43,43	1.19	4 (9%)
2	HEM	A	501	1	27,50,50	1.86	4 (14%)	17,82,82	1.65	4 (23%)
5	OUS	A	504	-	31,31,31	0.87	2 (6%)	42,43,43	1.62	8 (19%)
4	BTB	A	503	9	13,13,13	0.62	0	7,16,16	0.97	0
2	HEM	B	501	1	27,50,50	1.83	4 (14%)	17,82,82	1.80	5 (29%)

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	BTB	C	505	-	-	3/21/21/21	-
3	H4B	A	502	-	-	4/8/17/17	0/2/2/2
4	BTB	A	505	-	-	7/21/21/21	-
3	H4B	B	502	-	-	4/8/17/17	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	H4B	D	502	-	-	4/8/17/17	0/2/2/2
7	GOL	A	507	-	-	1/4/4/4	-
4	BTB	B	505	-	-	7/21/21/21	-
5	OUS	B	503	-	-	2/11/11/11	0/4/4/4
7	GOL	C	507	-	-	4/4/4/4	-
4	BTB	D	504	9	-	3/21/21/21	-
4	BTB	B	504	9	-	6/21/21/21	-
3	H4B	C	503	-	-	3/8/17/17	0/2/2/2
4	BTB	B	508	-	-	7/21/21/21	-
2	HEM	C	502	1	-	3/6/54/54	-
5	OUS	C	504	-	-	2/11/11/11	0/4/4/4
2	HEM	D	501	1	-	0/6/54/54	-
4	BTB	D	505	-	-	14/21/21/21	-
5	OUS	D	503	-	-	0/11/11/11	0/4/4/4
2	HEM	A	501	1	-	0/6/54/54	-
5	OUS	A	504	-	-	2/11/11/11	0/4/4/4
4	BTB	A	503	9	-	6/21/21/21	-
2	HEM	B	501	1	-	0/6/54/54	-

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	HEM	C3B-C2B	-4.69	1.33	1.40
2	A	501	HEM	C3B-C2B	-4.46	1.34	1.40
2	C	502	HEM	C3B-CAB	4.40	1.56	1.47
2	C	502	HEM	C3B-C2B	-4.27	1.34	1.40
2	D	501	HEM	C3C-C2C	-4.16	1.34	1.40
2	D	501	HEM	C3C-CAC	3.94	1.55	1.47
2	D	501	HEM	C3B-C2B	-3.89	1.35	1.40
2	B	501	HEM	C3B-CAB	3.82	1.55	1.47
2	C	502	HEM	C3C-CAC	3.78	1.55	1.47
2	A	501	HEM	C3C-CAC	3.72	1.55	1.47
2	A	501	HEM	C3B-CAB	3.60	1.55	1.47
2	A	501	HEM	C3C-C2C	-3.60	1.35	1.40
2	B	501	HEM	C3C-C2C	-3.51	1.35	1.40
2	B	501	HEM	C3C-CAC	3.48	1.54	1.47
2	D	501	HEM	C3B-CAB	3.41	1.54	1.47
2	C	502	HEM	C3C-C2C	-3.16	1.36	1.40
5	A	504	OUS	C05-C10	-2.28	1.38	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	508	BTB	C4-C2	-2.26	1.50	1.53
2	C	502	HEM	CMD-C2D	2.20	1.56	1.51
2	C	502	HEM	CAA-C2A	2.18	1.55	1.52
5	A	504	OUS	C10-N01	-2.15	1.34	1.37
4	B	508	BTB	C1-C2	-2.11	1.50	1.53

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	502	HEM	CBD-CAD-C3D	-6.22	101.02	112.48
2	D	501	HEM	CBA-CAA-C2A	-5.43	102.47	112.49
3	A	502	H4B	C4-C4A-C8A	5.42	119.38	114.57
3	B	502	H4B	C4-C4A-C8A	4.83	118.86	114.57
2	B	501	HEM	CAD-CBD-CGD	-4.57	105.01	112.67
5	A	504	OUS	C04-C05-C10	4.56	120.48	118.01
3	C	503	H4B	C4-C4A-C8A	4.53	118.60	114.57
3	D	502	H4B	C4-C4A-C8A	4.28	118.38	114.57
2	D	501	HEM	CAA-CBA-CGA	-4.20	105.62	112.67
3	B	502	H4B	C4-C4A-N5	4.13	122.59	119.12
3	D	502	H4B	C4-C4A-N5	4.10	122.56	119.12
3	C	503	H4B	N3-C2-N1	-3.86	119.37	125.42
2	D	501	HEM	CMA-C3A-C4A	-3.71	122.77	128.46
3	A	502	H4B	N3-C2-N1	-3.68	119.64	125.42
5	B	503	OUS	O29-C24-C25	3.60	120.48	115.78
3	D	502	H4B	C4-N3-C2	3.56	121.59	115.93
5	B	503	OUS	C04-C05-C10	3.54	119.93	118.01
3	C	503	H4B	C4-C4A-N5	3.47	122.03	119.12
3	D	502	H4B	N3-C2-N1	-3.43	120.04	125.42
3	A	502	H4B	C4-N3-C2	3.40	121.33	115.93
3	C	503	H4B	C4-N3-C2	3.40	121.33	115.93
5	C	504	OUS	O29-C24-C25	3.34	120.15	115.78
3	B	502	H4B	N3-C2-N1	-3.34	120.18	125.42
5	A	504	OUS	O29-C24-C25	3.27	120.05	115.78
2	A	501	HEM	CBD-CAD-C3D	-3.06	106.84	112.48
5	D	503	OUS	O29-C24-C25	3.04	119.76	115.78
3	B	502	H4B	N2-C2-N3	3.03	121.96	117.25
3	B	502	H4B	C2-N1-C8A	3.02	121.30	114.54
5	A	504	OUS	C05-C10-N01	-2.97	119.66	122.81
4	B	508	BTB	O4-C4-C2	-2.96	103.33	111.44
5	B	503	OUS	C05-C10-N01	-2.95	119.68	122.81
3	A	502	H4B	C2-N1-C8A	2.95	121.14	114.54
3	C	503	H4B	C2-N1-C8A	2.94	121.12	114.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	504	OUS	C08-C09-C10	-2.92	118.89	121.44
2	D	501	HEM	CAD-CBD-CGD	-2.86	107.88	112.67
2	B	501	HEM	CAA-CBA-CGA	-2.82	107.94	112.67
3	B	502	H4B	C4-N3-C2	2.82	120.41	115.93
5	A	504	OUS	C36-N31-C32	2.78	121.66	116.85
3	A	502	H4B	C4-C4A-N5	2.71	121.39	119.12
3	C	503	H4B	N2-C2-N3	2.68	121.41	117.25
3	D	502	H4B	C2-N1-C8A	2.67	120.51	114.54
2	C	502	HEM	CMA-C3A-C4A	-2.63	124.42	128.46
2	A	501	HEM	CMD-C2D-C1D	-2.61	124.45	128.46
5	C	504	OUS	C26-C21-C08	-2.48	116.77	120.86
2	D	501	HEM	CMA-C3A-C2A	2.44	129.54	124.94
5	B	503	OUS	C07-C08-C21	-2.40	117.20	121.36
2	C	502	HEM	CAA-CBA-CGA	-2.34	108.75	112.67
2	A	501	HEM	CMC-C2C-C3C	2.34	129.05	124.68
2	A	501	HEM	CBA-CAA-C2A	-2.33	108.19	112.49
3	D	502	H4B	C4A-N5-C6	-2.33	114.83	121.16
2	C	502	HEM	CMC-C2C-C3C	2.25	128.89	124.68
5	A	504	OUS	O29-C24-C23	-2.22	119.17	123.97
5	C	504	OUS	C36-N31-C32	2.21	120.66	116.85
5	D	503	OUS	C36-N31-C32	2.20	120.66	116.85
5	D	503	OUS	C05-C10-N01	-2.19	120.49	122.81
2	D	501	HEM	CBD-CAD-C3D	-2.18	108.47	112.48
5	A	504	OUS	C06-C05-C04	-2.17	119.51	123.66
5	B	503	OUS	C36-N31-C32	2.17	120.61	116.85
5	B	503	OUS	C06-C05-C04	-2.17	119.52	123.66
2	B	501	HEM	CBD-CAD-C3D	-2.15	108.52	112.48
3	B	502	H4B	C4A-N5-C6	-2.13	115.35	121.16
3	A	502	H4B	N2-C2-N3	2.10	120.51	117.25
2	B	501	HEM	CMC-C2C-C3C	2.08	128.57	124.68
5	C	504	OUS	C05-C10-N01	-2.05	120.63	122.81
2	B	501	HEM	CMA-C3A-C4A	-2.04	125.33	128.46
5	A	504	OUS	C33-C32-N31	-2.03	119.75	123.72
5	B	503	OUS	C30-O29-C24	2.03	121.75	117.76
5	D	503	OUS	C08-C09-C10	-2.02	119.67	121.44

There are no chirality outliers.

All (82) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	505	BTB	C1-C2-C3-O3
4	C	505	BTB	C4-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
4	C	505	BTB	N-C2-C3-O3
3	A	502	H4B	N5-C6-C9-O9
3	A	502	H4B	C7-C6-C9-O9
3	A	502	H4B	C7-C6-C9-C10
4	A	505	BTB	O1-C1-C2-C3
4	A	505	BTB	O1-C1-C2-C4
4	A	505	BTB	O1-C1-C2-N
4	A	505	BTB	C1-C2-C4-O4
4	A	505	BTB	C3-C2-C4-O4
4	A	505	BTB	N-C2-C4-O4
3	B	502	H4B	N5-C6-C9-O9
3	B	502	H4B	C7-C6-C9-O9
3	B	502	H4B	C7-C6-C9-C10
3	D	502	H4B	N5-C6-C9-O9
3	D	502	H4B	N5-C6-C9-C10
3	D	502	H4B	C7-C6-C9-O9
3	D	502	H4B	C7-C6-C9-C10
4	B	505	BTB	C1-C2-C3-O3
4	B	505	BTB	C4-C2-C3-O3
4	B	505	BTB	N-C2-C3-O3
4	B	505	BTB	C1-C2-C4-O4
4	B	505	BTB	C3-C2-C4-O4
7	C	507	GOL	O1-C1-C2-C3
7	C	507	GOL	C1-C2-C3-O3
4	D	504	BTB	O1-C1-C2-C3
4	D	504	BTB	O1-C1-C2-C4
4	D	504	BTB	O1-C1-C2-N
4	B	504	BTB	O1-C1-C2-C3
4	B	504	BTB	O1-C1-C2-C4
4	B	504	BTB	O1-C1-C2-N
4	B	504	BTB	C1-C2-C3-O3
4	B	504	BTB	N-C2-C3-O3
3	C	503	H4B	C7-C6-C9-O9
3	C	503	H4B	C7-C6-C9-C10
4	B	508	BTB	C1-C2-C3-O3
4	B	508	BTB	N-C2-C3-O3
4	B	508	BTB	C8-C7-N-C5
2	C	502	HEM	C1A-C2A-CAA-CBA
2	C	502	HEM	C3A-C2A-CAA-CBA
2	C	502	HEM	C2A-CAA-CBA-CGA
4	D	505	BTB	O1-C1-C2-C3
4	D	505	BTB	O1-C1-C2-C4

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Mol	Chain	Res	Type	Atoms
4	D	505	BTB	O1-C1-C2-N
4	D	505	BTB	C1-C2-C4-O4
4	D	505	BTB	C3-C2-C4-O4
4	D	505	BTB	N-C2-C4-O4
4	D	505	BTB	C1-C2-N-C7
4	D	505	BTB	C3-C2-N-C5
4	D	505	BTB	C3-C2-N-C7
4	D	505	BTB	C4-C2-N-C7
4	D	505	BTB	C8-C7-N-C5
4	A	503	BTB	O1-C1-C2-C3
4	A	503	BTB	O1-C1-C2-C4
4	A	503	BTB	O1-C1-C2-N
4	A	503	BTB	C4-C2-C3-O3
4	A	503	BTB	N-C2-C3-O3
4	B	505	BTB	N-C7-C8-O8
7	A	507	GOL	C1-C2-C3-O3
7	C	507	GOL	O1-C1-C2-O2
4	A	505	BTB	N-C5-C6-O6
7	C	507	GOL	O2-C2-C3-O3
4	D	505	BTB	N-C7-C8-O8
4	B	508	BTB	N-C7-C8-O8
3	B	502	H4B	N5-C6-C9-C10
5	C	504	OUS	C23-C24-O29-C30
5	C	504	OUS	C25-C24-O29-C30
5	A	504	OUS	C23-C24-O29-C30
4	B	504	BTB	C4-C2-C3-O3
4	A	503	BTB	C1-C2-C3-O3
4	B	505	BTB	N-C2-C4-O4
4	B	508	BTB	N-C2-C4-O4
4	D	505	BTB	C1-C2-N-C5
4	D	505	BTB	C4-C2-N-C5
3	A	502	H4B	N5-C6-C9-C10
5	A	504	OUS	C25-C24-O29-C30
5	B	503	OUS	C23-C24-O29-C30
5	B	503	OUS	C26-C25-C27-N28
3	C	503	H4B	N5-C6-C9-O9
4	B	508	BTB	C4-C2-C3-O3
4	B	508	BTB	C3-C2-C4-O4

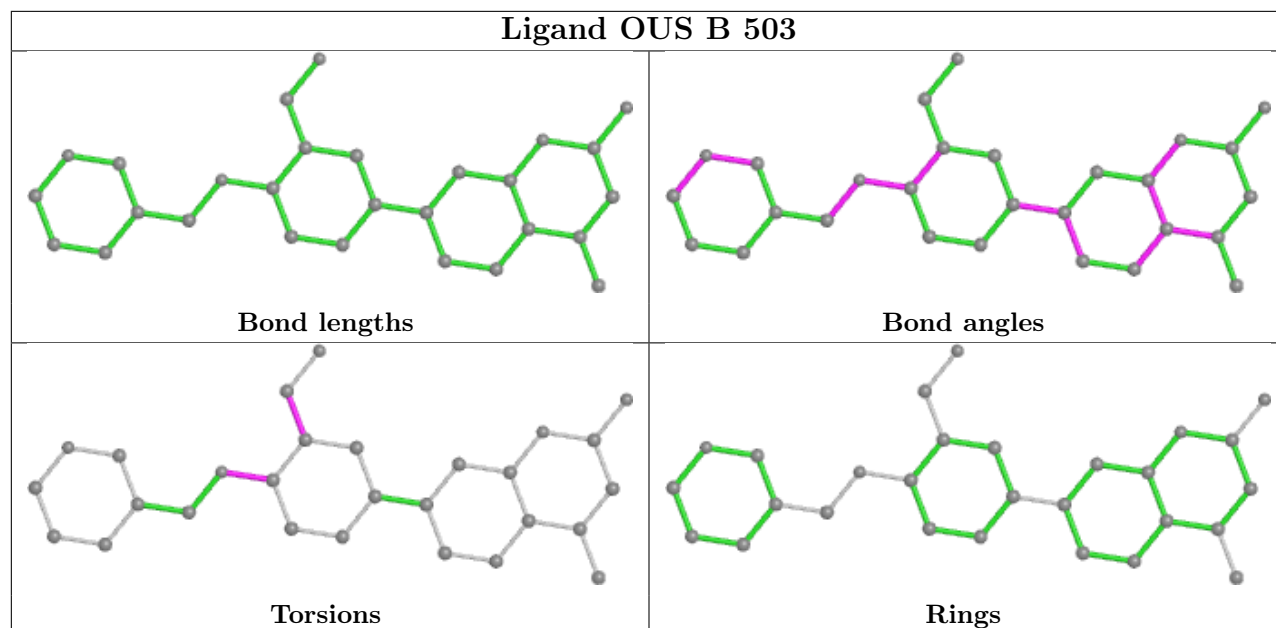
There are no ring outliers.

18 monomers are involved in 35 short contacts:

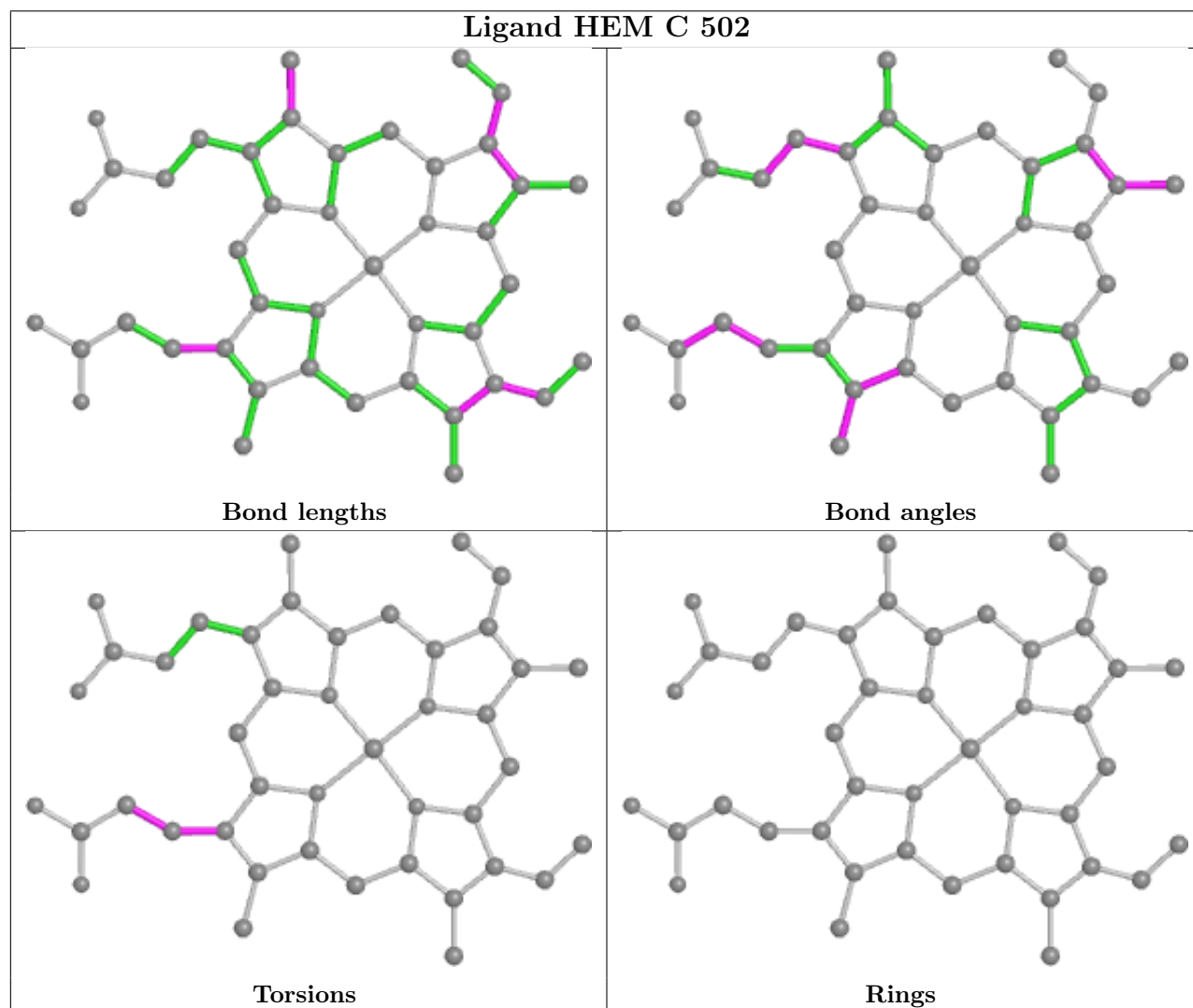
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502	H4B	1	0
4	A	505	BTB	1	0
7	A	507	GOL	1	0
4	B	505	BTB	2	0
5	B	503	OUS	2	0
4	D	504	BTB	1	0
4	B	504	BTB	1	0
3	C	503	H4B	1	0
4	B	508	BTB	3	0
2	C	502	HEM	4	0
5	C	504	OUS	3	0
2	D	501	HEM	2	0
4	D	505	BTB	3	0
5	D	503	OUS	1	0
2	A	501	HEM	4	0
5	A	504	OUS	2	0
4	A	503	BTB	2	0
2	B	501	HEM	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

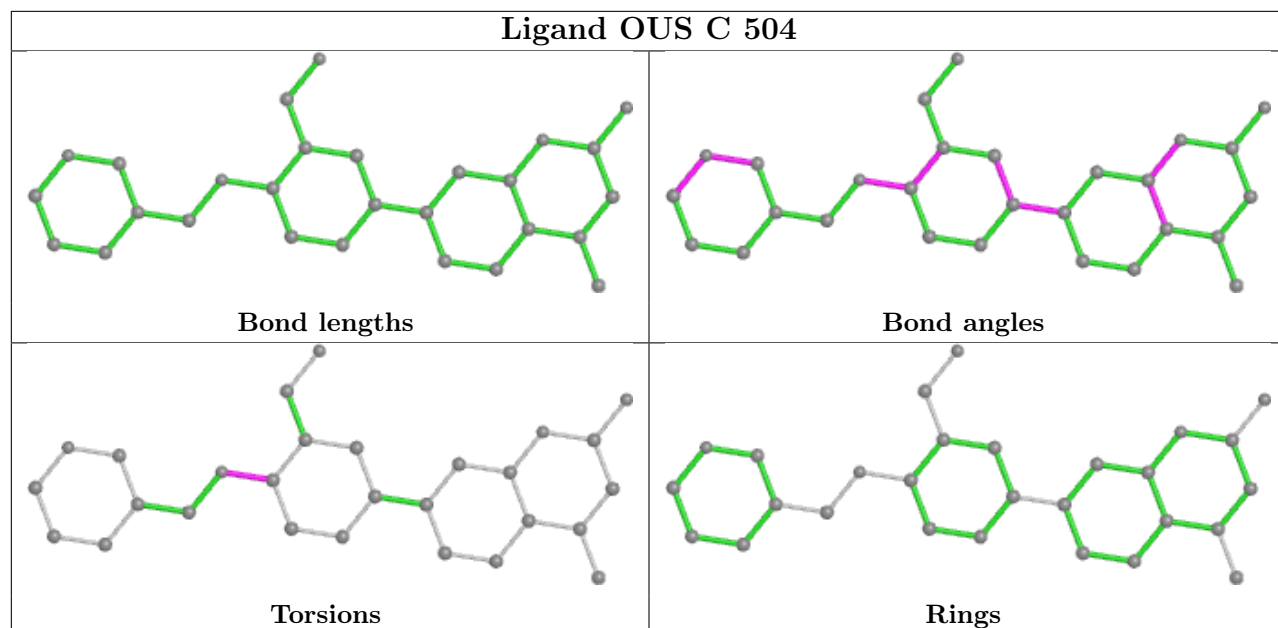
Ligand OUS B 503



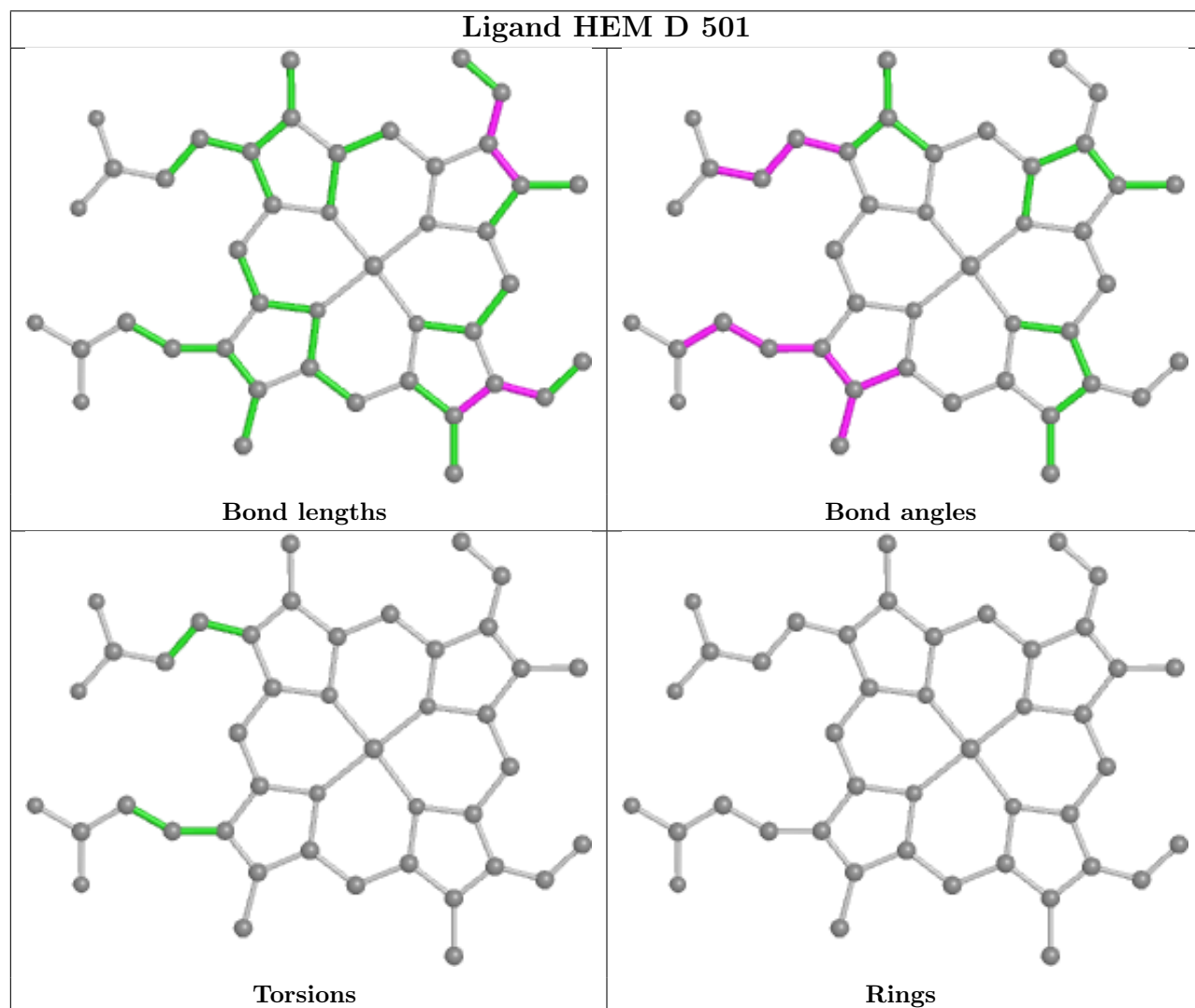
Ligand HEM C 502



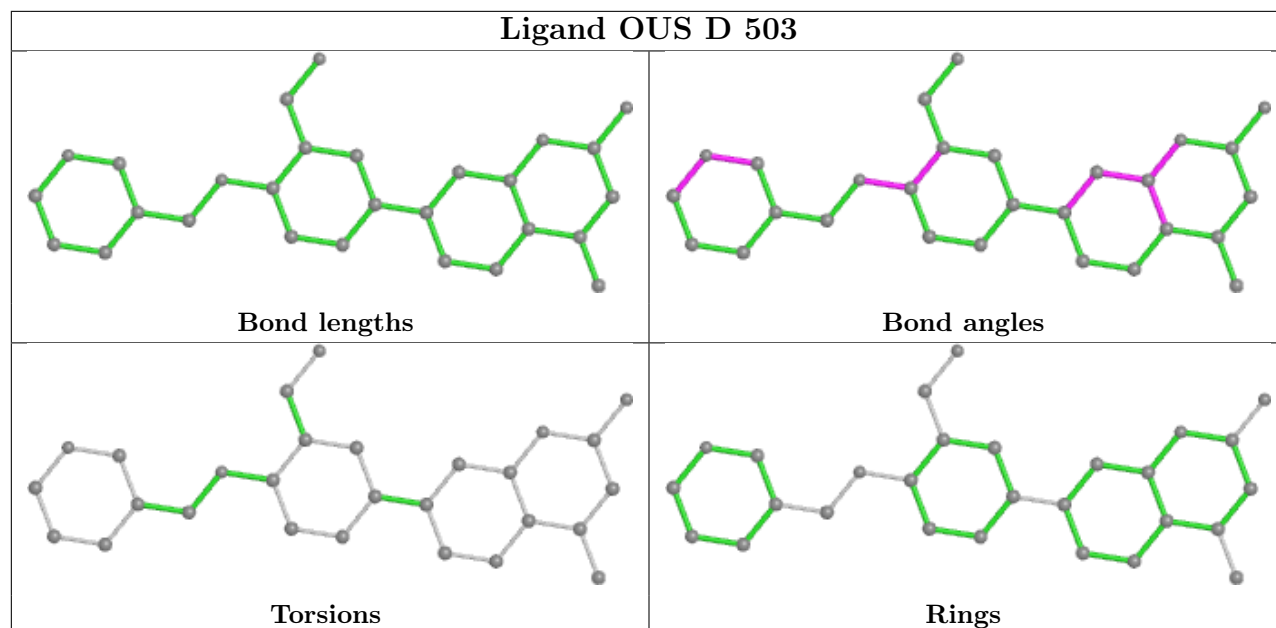
Ligand OUS C 504



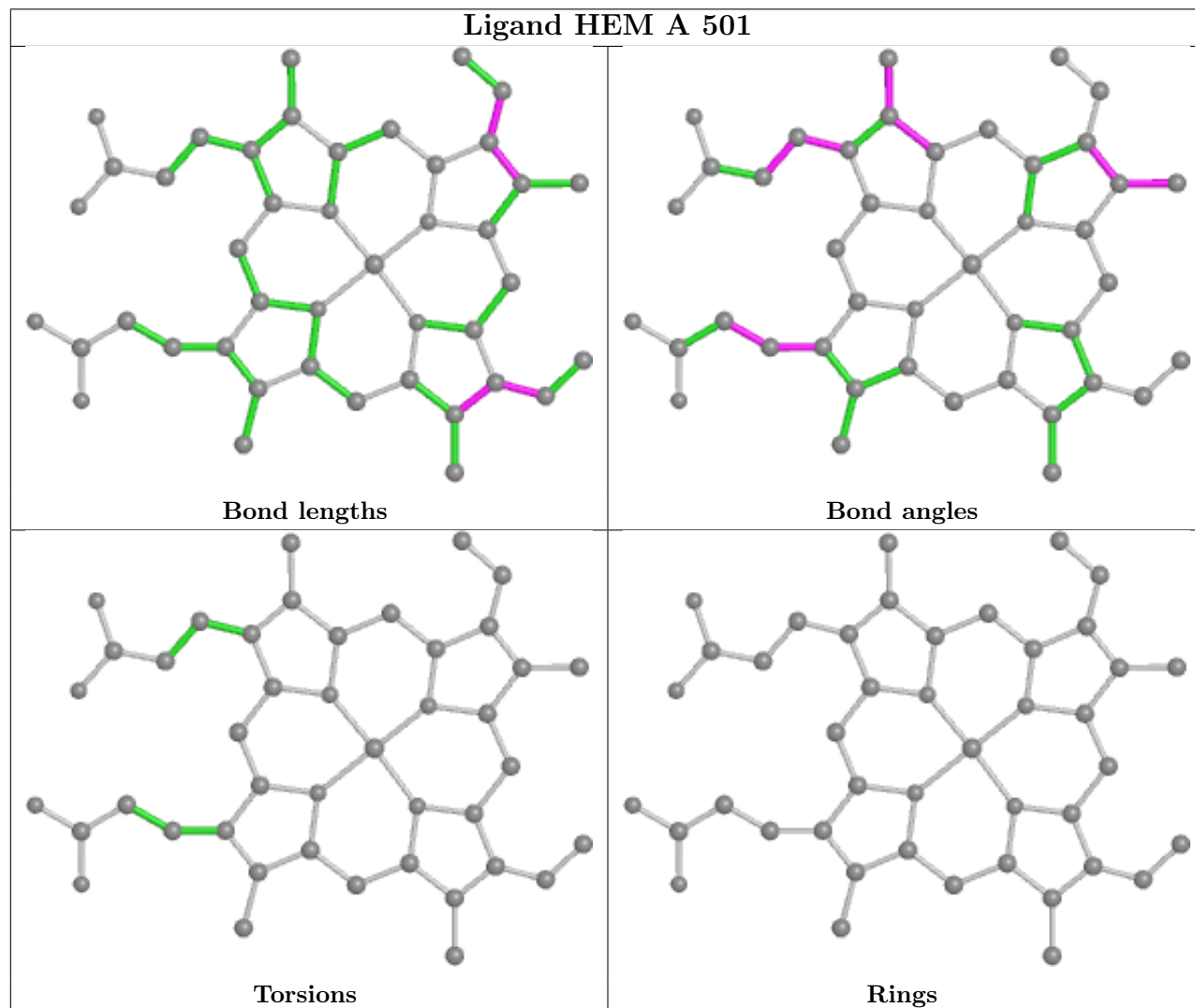
Ligand HEM D 501



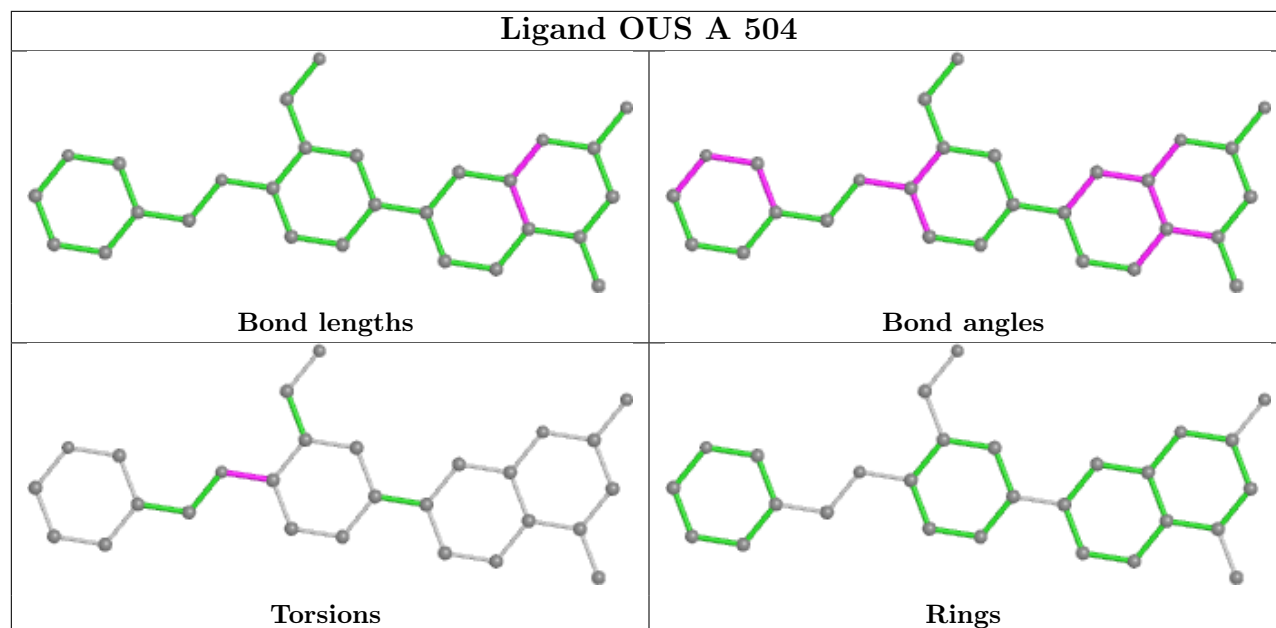
Ligand OUS D 503



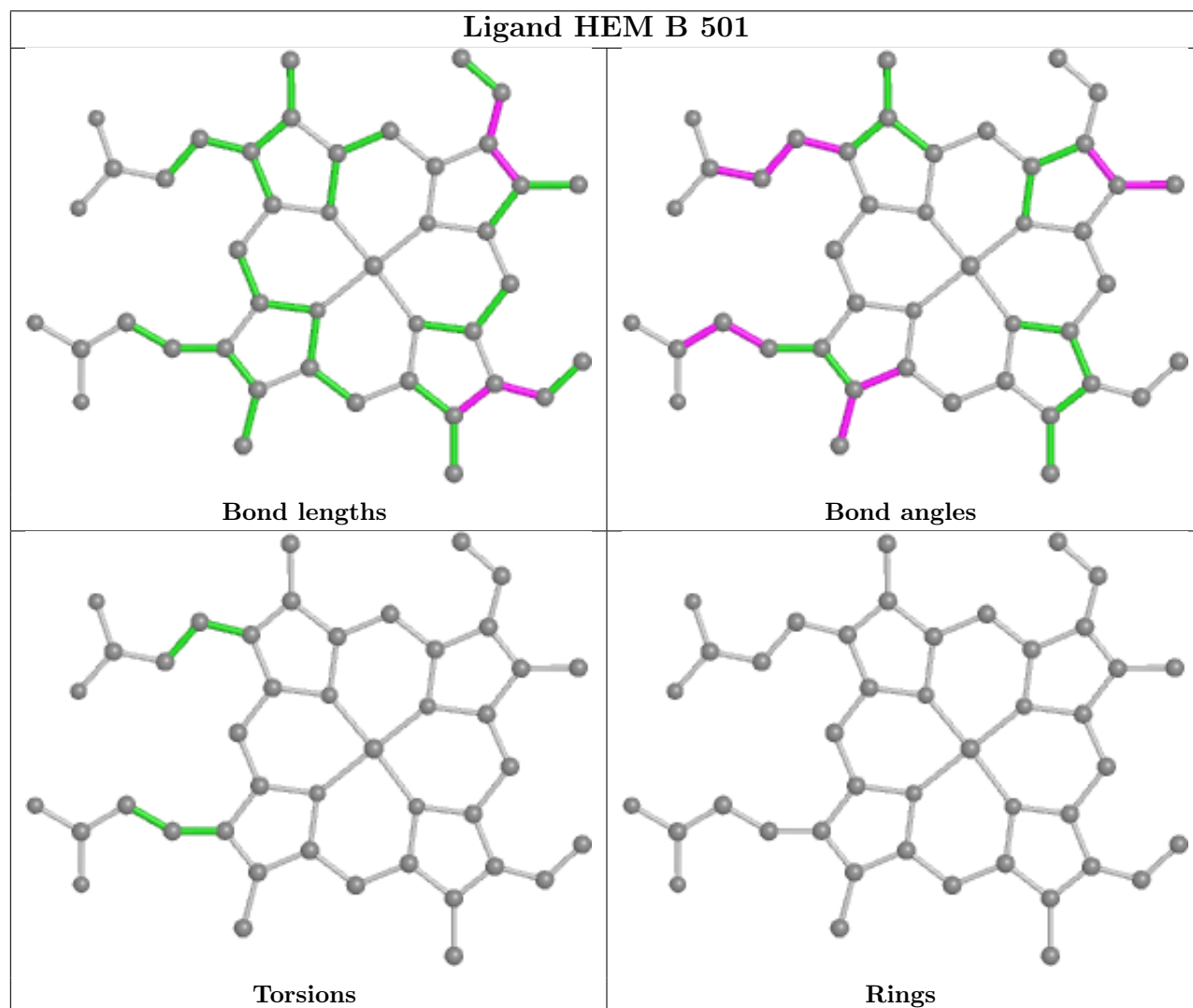
Ligand HEM A 501



Ligand OUS A 504



Ligand HEM B 501



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(Å²)	Q<0.9
1	A	404/440 (91%)	0.12	20 (4%)	29	27	28, 59, 120, 151	0
1	B	402/440 (91%)	-0.21	2 (0%)	90	90	24, 41, 78, 126	0
1	C	401/440 (91%)	-0.23	2 (0%)	90	90	25, 47, 85, 118	0
1	D	403/440 (91%)	-0.23	5 (1%)	79	77	21, 41, 79, 143	0
All	All	1610/1760 (91%)	-0.14	29 (1%)	68	66	21, 47, 100, 151	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	89	GLN	4.9
1	A	109	LEU	4.5
1	A	106	PRO	4.5
1	A	107	ARG	4.5
1	A	108	LYS	4.3
1	D	90	GLN	3.5
1	D	257	GLN	3.4
1	B	89	GLN	3.3
1	A	236	PRO	3.3
1	A	142	GLY	3.1
1	A	144	GLN	3.0
1	C	89	GLN	3.0
1	A	204	ALA	2.9
1	A	480	TRP	2.9
1	A	90	GLN	2.8
1	D	259	GLY	2.8
1	A	141	SER	2.8
1	B	90	GLN	2.7
1	A	259	GLY	2.6
1	A	475	TYR	2.5
1	A	272	GLU	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	273	LEU	2.4
1	A	119	ALA	2.3
1	A	300	PRO	2.2
1	C	143	SER	2.1
1	A	202	ARG	2.1
1	A	153	VAL	2.1
1	A	257	GLN	2.0
1	D	255	ARG	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](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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.

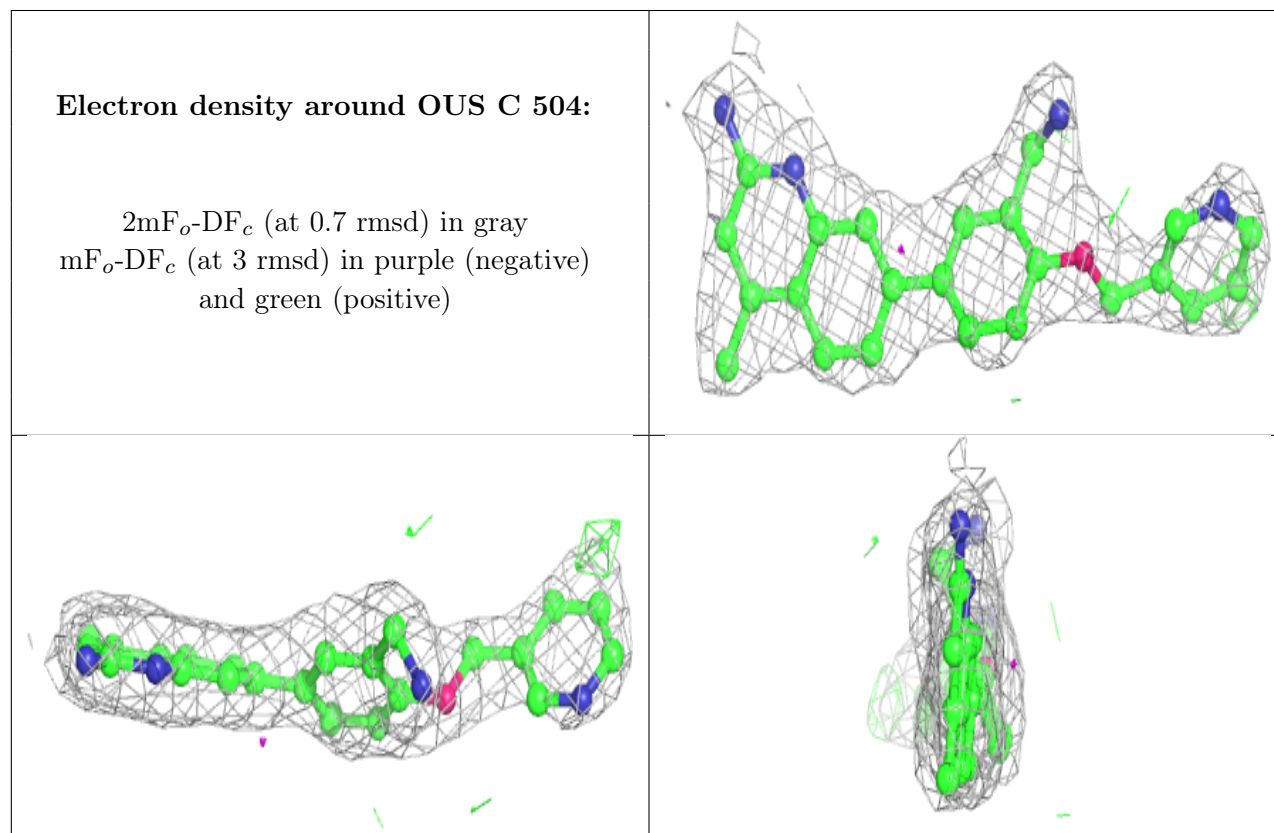
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	BTB	C	505	14/14	0.56	0.20	77,111,119,119	0
3	H4B	B	502	17/17	0.79	0.34	57,71,102,102	0
4	BTB	D	504	14/14	0.79	0.21	47,69,81,82	0
4	BTB	A	503	14/14	0.79	0.22	66,73,87,94	0
3	H4B	C	503	17/17	0.82	0.34	65,80,90,95	0
7	GOL	A	507	6/6	0.82	0.23	70,79,81,91	0
4	BTB	A	505	14/14	0.85	0.19	75,90,94,95	0
4	BTB	B	504	14/14	0.88	0.16	47,63,84,86	0
3	H4B	A	502	17/17	0.88	0.31	69,78,98,102	0
4	BTB	D	505	14/14	0.88	0.19	60,71,86,86	0
3	H4B	D	502	17/17	0.88	0.27	47,64,89,91	0
7	GOL	C	507	6/6	0.91	0.17	52,58,67,67	0
4	BTB	B	505	14/14	0.92	0.18	23,61,81,87	0
5	OUS	C	504	28/28	0.92	0.14	32,57,81,83	0
4	BTB	B	508	14/14	0.93	0.17	36,64,82,84	0

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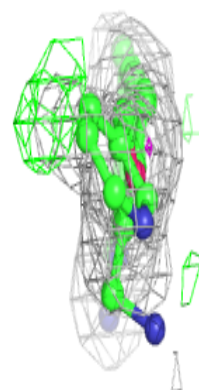
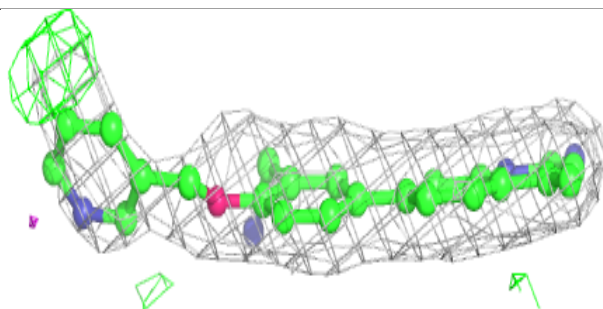
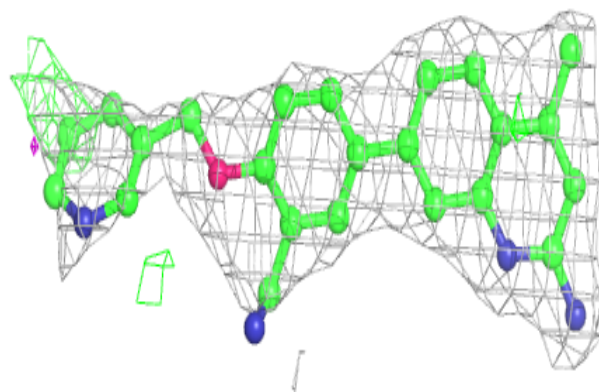
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	OUS	D	503	28/28	0.95	0.15	21,49,88,90	0
5	OUS	A	504	28/28	0.95	0.17	43,63,96,98	0
5	OUS	B	503	28/28	0.95	0.16	26,59,96,97	0
8	CL	A	508	1/1	0.96	0.07	58,58,58,58	0
2	HEM	A	501	43/43	0.97	0.14	27,50,95,103	0
2	HEM	C	502	43/43	0.97	0.13	26,38,85,95	0
8	CL	B	506	1/1	0.97	0.14	43,43,43,43	0
8	CL	D	507	1/1	0.98	0.07	44,44,44,44	0
9	GD	D	506	1/1	0.98	0.18	48,48,48,48	0
2	HEM	D	501	43/43	0.98	0.13	15,30,73,83	0
2	HEM	B	501	43/43	0.98	0.12	18,32,67,88	0
8	CL	C	508	1/1	0.99	0.11	55,55,55,55	0
9	GD	B	507	1/1	0.99	0.17	48,48,48,48	0
6	ZN	A	506	1/1	0.99	0.10	42,42,42,42	0
9	GD	C	501	1/1	0.99	0.07	78,78,78,78	0
6	ZN	C	506	1/1	1.00	0.11	32,32,32,32	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.

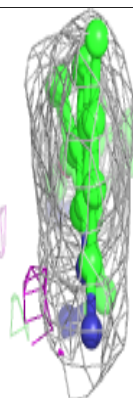
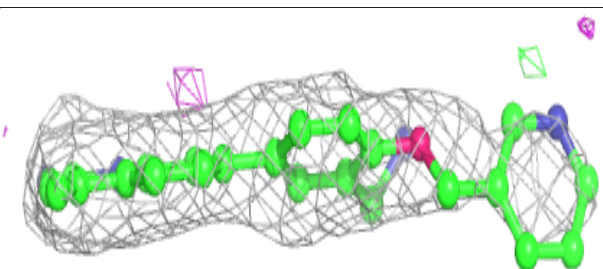
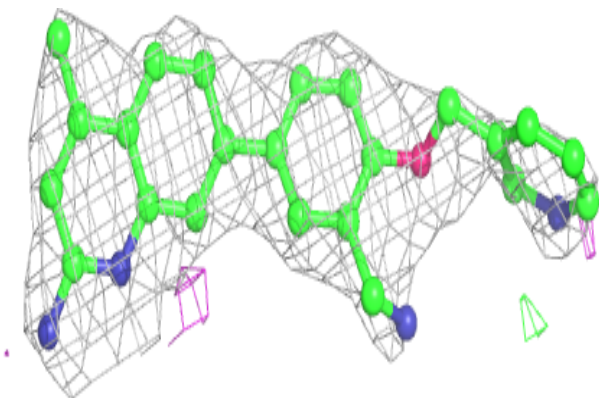


Electron density around OUS D 503:

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

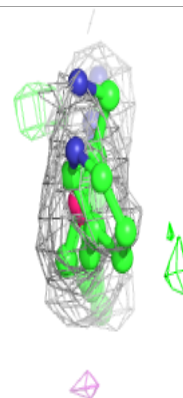
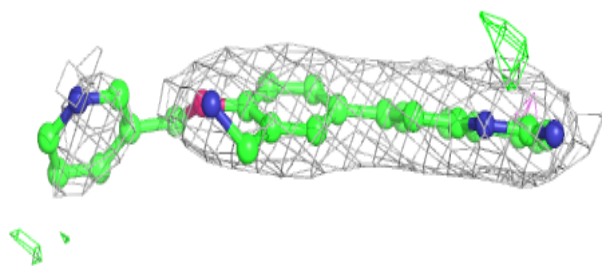
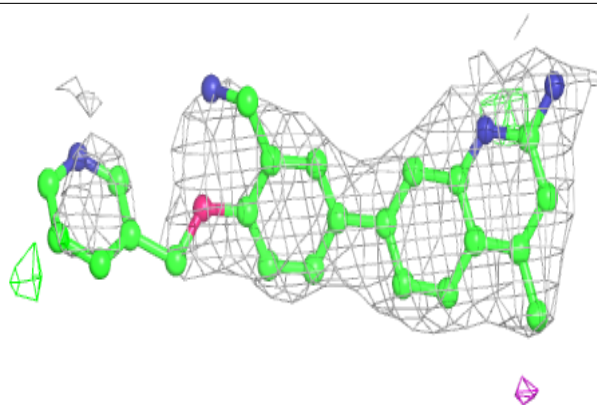
**Electron density around OUS A 504:**

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



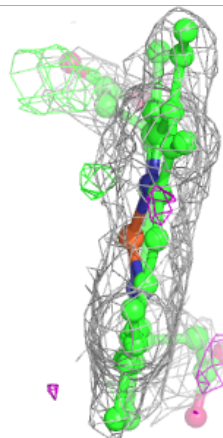
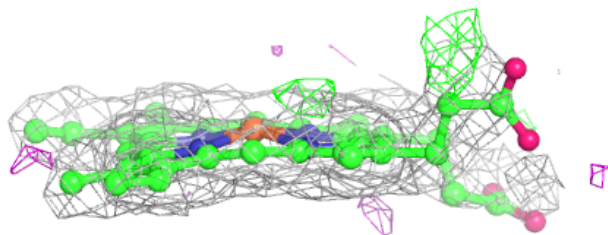
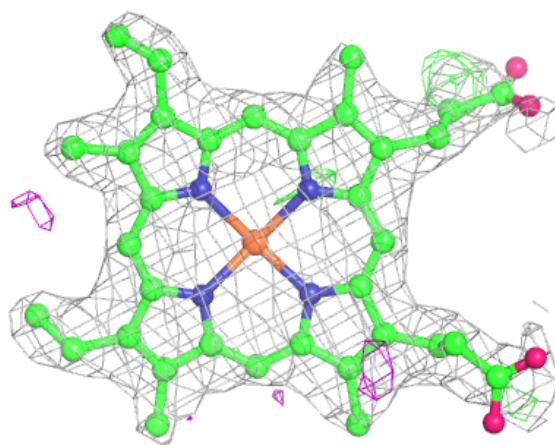
Electron density around OUS B 503:

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



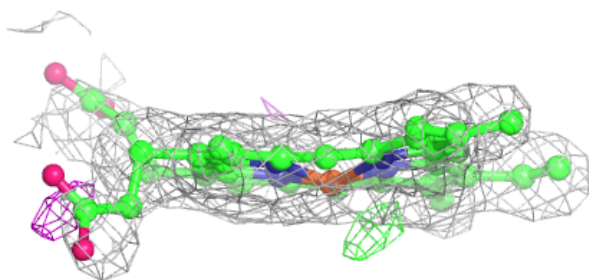
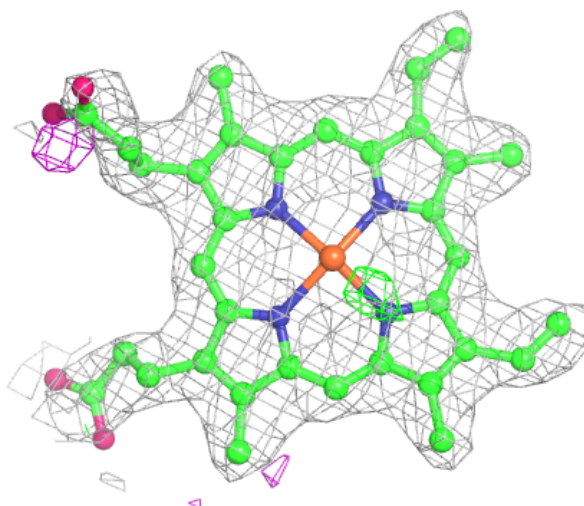
Electron density around HEM A 501:

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



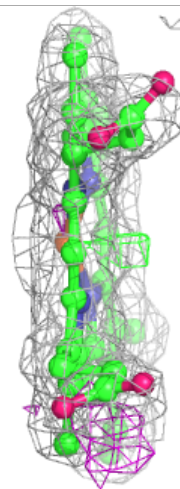
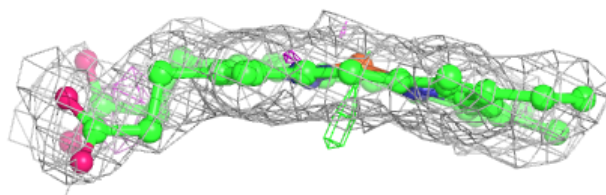
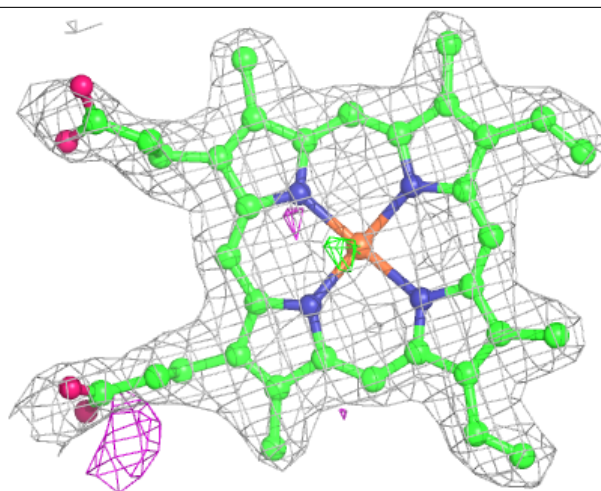
Electron density around HEM C 502:

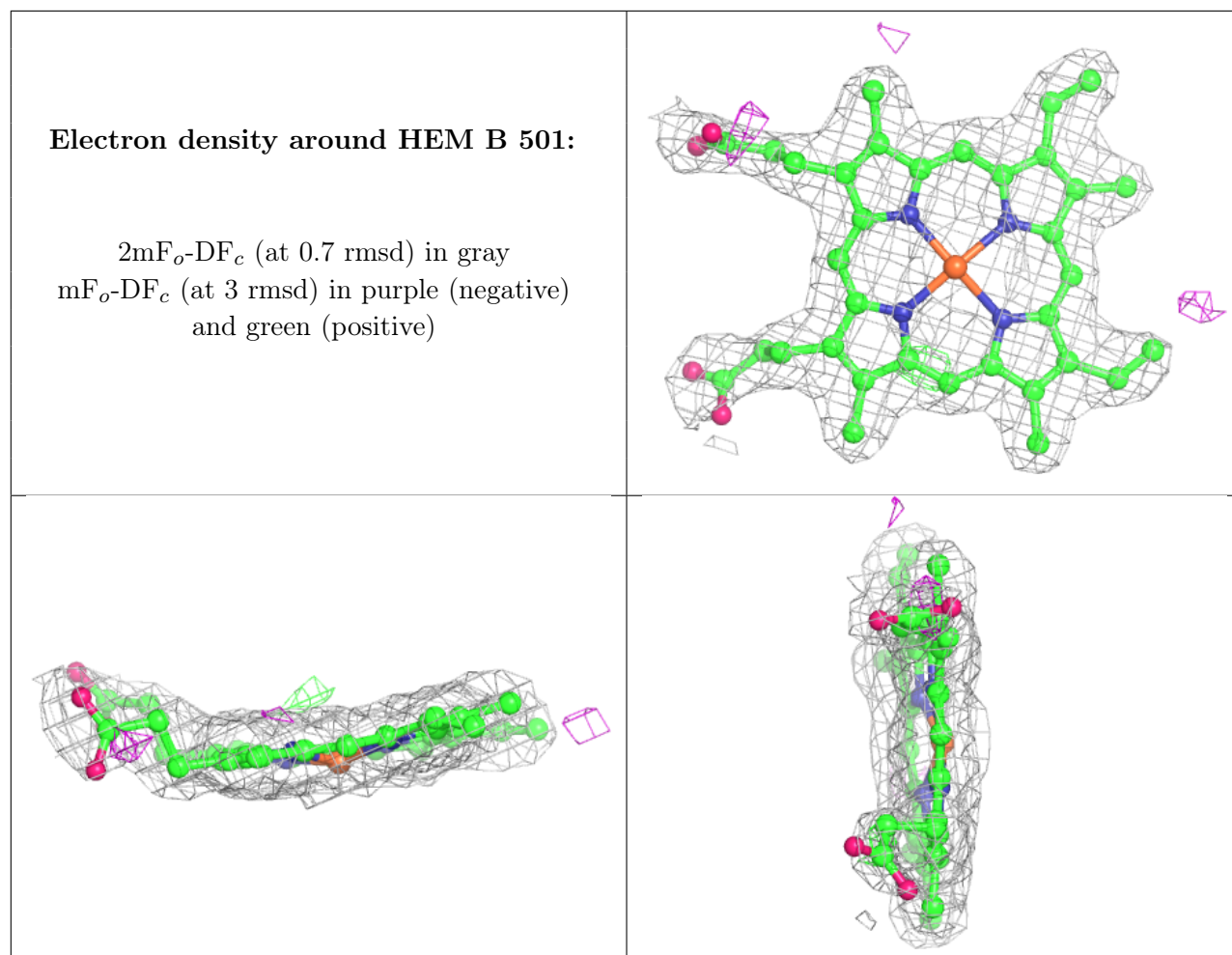
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM D 501:

$2mF_o-DF_c$ (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.