



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 06:57 PM UTC

PDB ID : 8UFT / pdb_00008uft
Title : Structure of human endothelial nitric oxide synthase P370N mutant heme domain in complex with 4-methyl-7-(4-methyl-2,3,4,5-tetrahydrobenzo[f][1,4]oxazepin-7-yl)quinolin-2-amine
Authors : Li, H.; Poulos, T.L.
Deposited on : 2023-10-04
Resolution : 1.78 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

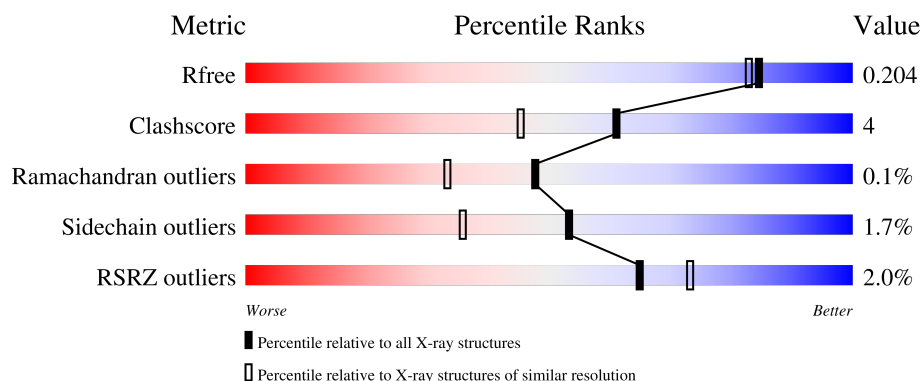
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.78 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	1365 (1.78-1.78)
Clashscore	190562	1395 (1.78-1.78)
Ramachandran outliers	187476	1382 (1.78-1.78)
Sidechain outliers	187428	1382 (1.78-1.78)
RSRZ outliers	180081	1365 (1.78-1.78)

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	440	3% 85% 5% 8%
1	B	440	% 84% 7% 8%
1	C	440	2% 84% 8% 8%
1	D	440	% 85% 6% 8%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	ACT	B	501	-	-	X	-
5	ACT	C	505	-	-	X	-

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 14776 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.

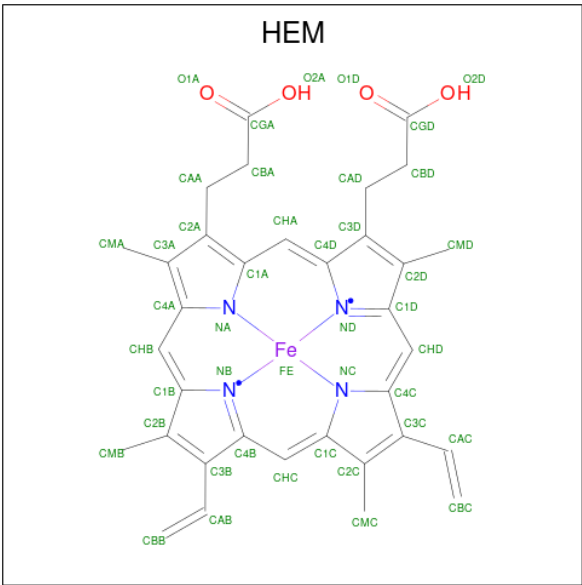
- Molecule 1 is a protein called Nitric oxide synthase 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	403	Total	C	N	O	S	0	2	0
			3223	2052	567	588	16			
1	B	403	Total	C	N	O	S	0	6	0
			3243	2065	569	593	16			
1	C	403	Total	C	N	O	S	0	4	0
			3234	2059	567	592	16			
1	D	403	Total	C	N	O	S	0	5	0
			3238	2061	568	592	17			

There are 8 discrepancies between the modelled and reference sequences:

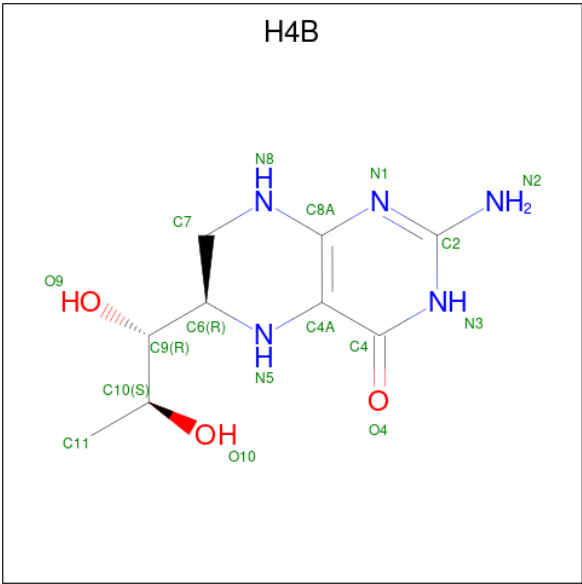
Chain	Residue	Modelled	Actual	Comment	Reference
A	298	GLU	ASP	variant	UNP P29474
A	370	ASN	PRO	engineered mutation	UNP P29474
B	298	GLU	ASP	variant	UNP P29474
B	370	ASN	PRO	engineered mutation	UNP P29474
C	298	GLU	ASP	variant	UNP P29474
C	370	ASN	PRO	engineered mutation	UNP P29474
D	298	GLU	ASP	variant	UNP P29474
D	370	ASN	PRO	engineered mutation	UNP P29474

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



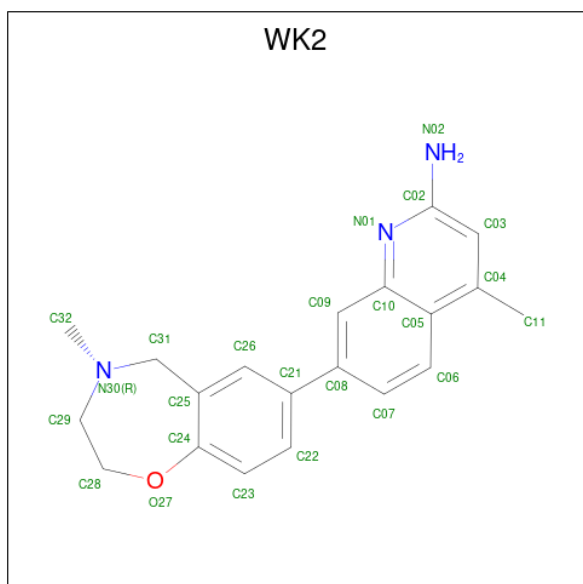
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

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



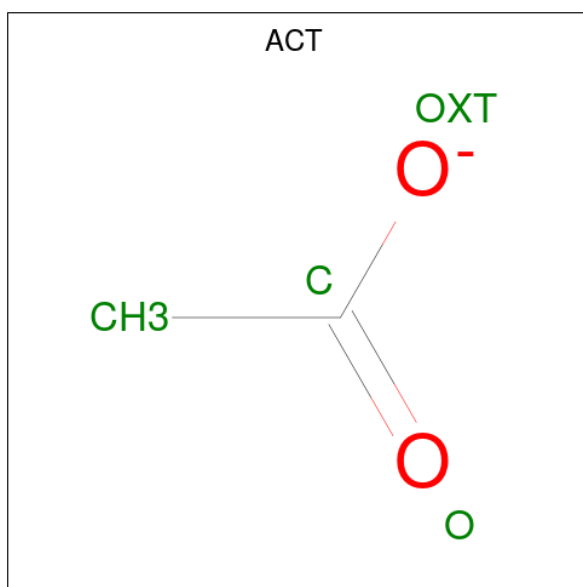
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			17	9	5	3		
3	B	1	Total	C	N	O	0	0
			17	9	5	3		
3	C	1	Total	C	N	O	0	0
			17	9	5	3		
3	D	1	Total	C	N	O	0	0
			17	9	5	3		

- Molecule 4 is (7M)-4-methyl-7-(4-methyl-2,3,4,5-tetrahydro-1,4-benzoxazepin-7-yl)quinolin-2-amine (CCD ID: WK2) (formula: $C_{20}H_{21}N_3O$) (labeled as "Ligand of Interest" by depositor).



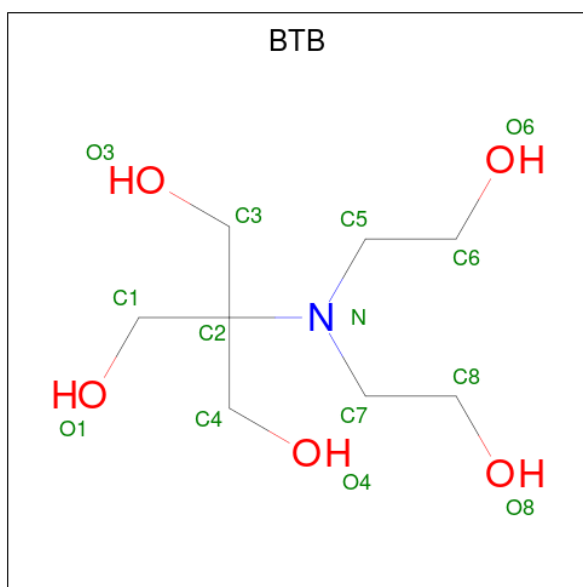
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			24	20	3	1		
4	B	1	Total	C	N	O	0	0
			24	20	3	1		
4	C	1	Total	C	N	O	0	0
			24	20	3	1		
4	D	1	Total	C	N	O	0	0
			24	20	3	1		

- Molecule 5 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



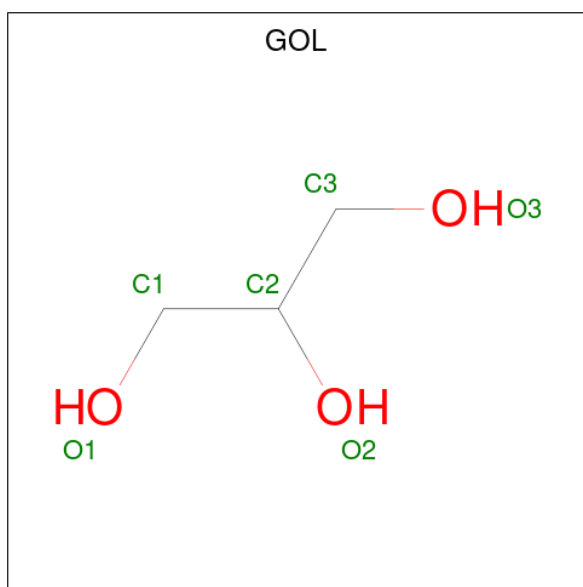
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	C	1	Total	C	O	0	0
			4	2	2		
5	C	1	Total	C	O	0	0
			4	2	2		
5	D	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is 2-[BIS-(2-HYDROXY-ETHYL)-AMINO]-2-HYDROXYMETHYL-PROPAN E-1,3-DIOL (CCD ID: BTB) (formula: C₈H₁₉NO₅).



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

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: $\text{C}_3\text{H}_8\text{O}_3$).



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

- Molecule 8 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

- Molecule 9 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 10 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	2	Total 2	Ca 2	0	0
10	B	1	Total 1	Ca 1	0	0

- Molecule 11 is GADOLINIUM ATOM (CCD ID: GD) (formula: Gd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	B	1	Total 1	Gd 1	0	0
11	C	2	Total 2	Gd 2	0	0
11	D	1	Total 1	Gd 1	0	0

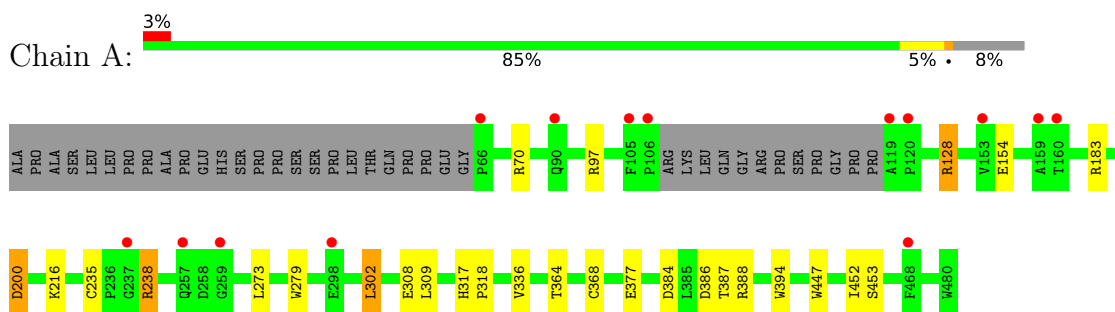
- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	262	Total 262	O 262	0	0
12	B	406	Total 406	O 406	0	0
12	C	284	Total 284	O 284	0	0
12	D	367	Total 367	O 367	0	0

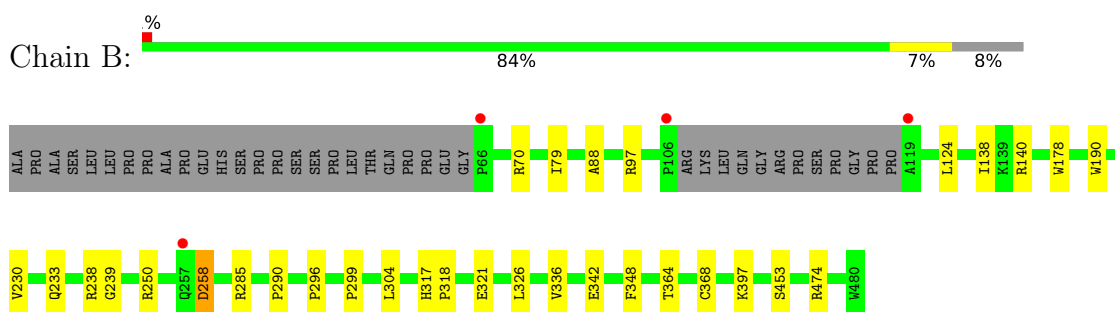
3 Residue-property plots [i](#)

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

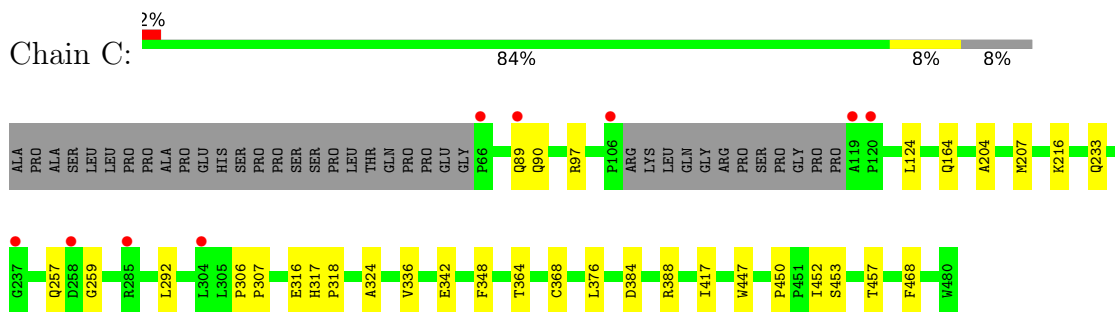
• Molecule 1: Nitric oxide synthase 3



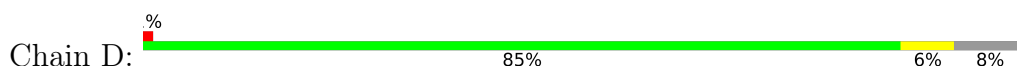
• Molecule 1: Nitric oxide synthase 3

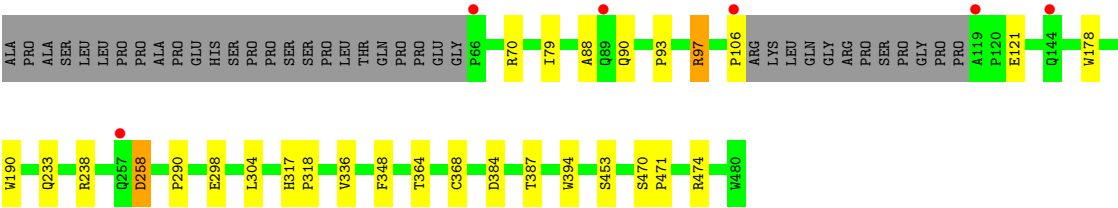


• Molecule 1: Nitric oxide synthase 3



• Molecule 1: Nitric oxide synthase 3





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	60.06Å 154.61Å 108.58Å 90.00° 90.73° 90.00°	Depositor
Resolution (Å)	40.02 – 1.78 40.02 – 1.78	Depositor EDS
% Data completeness (in resolution range)	97.5 (40.02-1.78) 98.0 (40.02-1.78)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 1.78Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.175 , 0.210 0.170 , 0.204	Depositor DCC
R_{free} test set	9300 reflections (4.24%)	wwPDB-VP
Wilson B-factor (Å ²)	25.6	Xtriage
Anisotropy	0.548	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 50.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.066 for h,-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	14776	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.30% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GD, CA, WK2, H4B, HEM, ZN, CL, ACT, GOL, BTB

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	0.25	0/3321	0.44	0/4524
1	B	0.32	0/3353	0.50	0/4567
1	C	0.24	0/3338	0.44	0/4547
1	D	0.31	0/3342	0.49	0/4553
All	All	0.28	0/13354	0.47	0/18191

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3223	0	3128	22	0
1	B	3243	0	3153	26	0
1	C	3234	0	3138	18	0
1	D	3238	0	3144	20	0
2	A	43	0	30	3	0
2	B	43	0	30	3	0
2	C	43	0	30	3	0
2	D	43	0	30	11	0
3	A	17	0	15	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	17	0	15	0	0
3	C	17	0	15	2	0
3	D	17	0	15	4	0
4	A	24	0	0	1	0
4	B	24	0	0	2	0
4	C	24	0	0	1	0
4	D	24	0	0	5	0
5	A	4	0	3	0	0
5	B	8	0	6	8	0
5	C	8	0	6	5	0
5	D	4	0	3	1	0
6	A	14	0	19	2	0
6	B	28	0	36	3	0
6	C	28	0	38	1	0
6	D	28	0	36	8	0
7	A	12	0	16	1	0
7	B	6	0	8	0	0
7	C	18	0	24	0	0
7	D	12	0	16	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	A	1	0	0	0	0
9	C	1	0	0	0	0
10	A	2	0	0	0	0
10	B	1	0	0	0	0
11	B	1	0	0	0	0
11	C	2	0	0	0	0
11	D	1	0	0	0	0
12	A	262	0	0	2	0
12	B	406	0	0	6	0
12	C	284	0	0	2	0
12	D	367	0	0	1	0
All	All	14776	0	12954	110	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 (110) 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)
2:D:501:HEM:O1A	3:D:502:H4B:N2	2.15	0.79
1:C:452:ILE:HB	5:C:505:ACT:H2	1.68	0.75
1:B:238:ARG:NH2	12:B:601:HOH:O	2.20	0.74
1:B:258:ASP:N	1:B:258:ASP:OD1	2.20	0.72
2:D:501:HEM:O2A	4:D:503:WK2:C25	2.39	0.71
1:A:453:SER:H	5:B:501:ACT:H1	1.56	0.70
1:B:453:SER:N	5:B:501:ACT:H2	2.06	0.70
1:A:200:ASP:OD1	1:A:200:ASP:N	2.25	0.69
2:B:502:HEM:HBC2	2:B:502:HEM:HMC2	1.75	0.69
1:B:453:SER:H	5:B:501:ACT:H2	1.57	0.69
1:B:290:PRO:HB3	1:B:304:LEU:HD23	1.75	0.67
1:D:258:ASP:OD1	1:D:258:ASP:N	2.27	0.67
2:A:501:HEM:HMC2	2:A:501:HEM:HBC2	1.76	0.66
1:A:384:ASP:HB3	12:A:764:HOH:O	1.97	0.64
1:A:452:ILE:HB	5:B:501:ACT:H3	1.79	0.64
2:D:501:HEM:HMC2	2:D:501:HEM:HBC2	1.80	0.64
2:C:501:HEM:HBC2	2:C:501:HEM:HMC2	1.81	0.63
1:A:453:SER:H	5:B:501:ACT:CH3	2.10	0.63
1:C:90:GLN:HB3	1:C:468:PHE:CD2	2.36	0.61
2:C:501:HEM:HBB2	2:C:501:HEM:HHC	1.82	0.61
1:D:336:VAL:HG21	4:D:503:WK2:C07	2.32	0.60
1:C:384[A]:ASP:OD1	12:C:601:HOH:O	2.17	0.60
1:B:342[B]:GLU:OE1	1:B:474:ARG:NH1	2.31	0.59
2:D:501:HEM:O1A	3:D:502:H4B:N3	2.35	0.59
2:D:501:HEM:O2A	4:D:503:WK2:C31	2.50	0.59
5:C:505:ACT:H1	1:D:453:SER:H	1.69	0.58
1:D:97:ARG:HH11	1:D:97:ARG:HB3	1.67	0.58
2:B:502:HEM:HBA1	4:B:504:WK2:C09	2.34	0.58
1:C:204:ALA:HA	1:C:207:MET:HE3	1.85	0.58
1:D:70:ARG:HD2	1:D:79:ILE:HD13	1.86	0.57
1:D:93:PRO:HG3	1:D:106:PRO:HB3	1.86	0.57
2:A:501:HEM:HBB2	2:A:501:HEM:HHC	1.87	0.57
1:D:298:GLU:OE2	6:D:506:BTB:H32	2.06	0.56
1:B:336:VAL:HG21	4:B:504:WK2:C07	2.38	0.54
1:A:336:VAL:HG21	4:A:503:WK2:C07	2.38	0.54
1:D:290:PRO:HB3	1:D:304:LEU:HD23	1.89	0.54
1:A:235:CYS:SG	1:A:238:ARG:HD2	2.48	0.53
1:B:453:SER:H	5:B:501:ACT:CH3	2.21	0.53
6:A:505:BTB:O8	1:D:384:ASP:OD2	2.26	0.53
1:C:364:THR:O	1:C:368:CYS:HB2	2.08	0.53
1:A:128:ARG:HH22	1:A:154:GLU:CD	2.18	0.52
5:C:505:ACT:H1	1:D:453:SER:N	2.23	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:298:GLU:OE1	6:D:506:BTB:H41	2.08	0.52
1:A:453:SER:N	5:B:501:ACT:H1	2.24	0.52
2:D:501:HEM:HBB2	2:D:501:HEM:HHC	1.91	0.52
2:D:501:HEM:O2A	4:D:503:WK2:C26	2.59	0.51
2:D:501:HEM:O1A	3:D:502:H4B:C2	2.59	0.51
1:C:97:ARG:HG3	1:D:88:ALA:HB3	1.93	0.50
1:B:138:ILE:HD12	1:B:140:ARG:HD3	1.93	0.50
6:D:506:BTB:O4	6:D:506:BTB:O3	2.22	0.50
1:A:377:GLU:OE2	6:A:505:BTB:O1	2.29	0.50
1:A:70:ARG:NH1	12:A:609:HOH:O	2.44	0.50
1:C:450:PRO:HG2	1:C:457:THR:HG21	1.95	0.49
2:D:501:HEM:CGA	3:D:502:H4B:HN3	2.25	0.49
1:B:238:ARG:HD3	1:B:296:PRO:HB3	1.95	0.49
1:C:336:VAL:HG21	4:C:503:WK2:C07	2.43	0.49
1:A:183:ARG:HD3	1:A:447:TRP:CD2	2.47	0.48
1:D:387:THR:HA	1:D:394:TRP:CD1	2.48	0.48
1:A:447:TRP:CZ3	2:A:501:HEM:HBA2	2.48	0.48
1:A:364:THR:O	1:A:368:CYS:HB2	2.13	0.48
1:C:447:TRP:HA	3:C:502:H4B:N1	2.29	0.47
1:C:388:ARG:HG2	6:C:506:BTB:H81	1.95	0.47
6:D:505:BTB:H82	6:D:505:BTB:H41	1.96	0.47
1:B:285:ARG:NH1	12:B:613:HOH:O	2.47	0.47
1:C:316[B]:GLU:HG2	1:C:324:ALA:HB2	1.97	0.47
1:C:453:SER:H	5:C:505:ACT:CH3	2.28	0.47
5:C:505:ACT:CH3	1:D:453:SER:H	2.27	0.47
1:B:397:LYS:NZ	12:B:611:HOH:O	2.47	0.47
2:B:502:HEM:HHC	2:B:502:HEM:HBB2	1.98	0.46
1:C:368:CYS:SG	1:C:376:LEU:HD13	2.55	0.46
1:B:70:ARG:NH1	12:B:617:HOH:O	2.49	0.46
2:D:501:HEM:HMB2	5:D:504:ACT:H2	1.97	0.45
1:C:233:GLN:HB3	1:C:348:PHE:CE2	2.51	0.45
1:A:386:ASP:OD2	1:A:388:ARG:HG2	2.15	0.45
1:A:387:THR:HA	1:A:394:TRP:CD1	2.52	0.45
1:B:233:GLN:HB3	1:B:348:PHE:CE2	2.51	0.45
2:C:501:HEM:CGA	3:C:502:H4B:HN3	2.30	0.45
1:B:321:GLU:OE1	6:B:506:BTB:O8	2.35	0.45
1:B:299:PRO:HD2	6:B:507:BTB:H82	1.97	0.45
6:D:505:BTB:H11	6:D:505:BTB:H72	1.72	0.45
1:B:238:ARG:HD3	1:B:296:PRO:CB	2.47	0.44
1:B:178:TRP:CE3	1:B:190:TRP:HA	2.53	0.44
6:D:506:BTB:H11	6:D:506:BTB:H51	1.29	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:317:HIS:CG	1:D:318:PRO:HD2	2.52	0.44
1:B:250:ARG:HG3	12:B:841:HOH:O	2.18	0.43
1:D:470:SER:HA	1:D:471:PRO:C	2.43	0.43
6:D:506:BTB:HO3	6:D:506:BTB:HO4	1.60	0.43
1:B:238:ARG:HD2	1:B:239:GLY:O	2.18	0.43
1:B:364:THR:O	1:B:368:CYS:HB2	2.18	0.43
1:D:233:GLN:HB3	1:D:348:PHE:CE2	2.54	0.43
1:A:273:LEU:HD23	1:A:273:LEU:HA	1.88	0.43
1:A:279:TRP:HB2	1:A:302:LEU:HD11	2.01	0.43
1:C:317:HIS:CG	1:C:318:PRO:HD2	2.54	0.42
1:B:97:ARG:HG3	12:B:960:HOH:O	2.18	0.42
1:D:364:THR:O	1:D:368:CYS:HB2	2.20	0.42
1:B:124:LEU:HD23	1:B:124:LEU:HA	1.92	0.42
1:A:384:ASP:OD2	7:A:507:GOL:O1	2.29	0.41
1:B:317:HIS:CG	1:B:318:PRO:HD2	2.55	0.41
1:C:306:PRO:HA	1:C:307:PRO:HD3	1.95	0.41
1:A:452:ILE:CB	5:B:501:ACT:H3	2.50	0.41
1:C:292:LEU:HD23	1:C:292:LEU:HA	1.97	0.41
1:D:178:TRP:CE3	1:D:190:TRP:HA	2.56	0.41
1:D:474:ARG:HD2	12:D:636:HOH:O	2.21	0.41
6:D:506:BTB:H41	6:D:506:BTB:H72	1.47	0.41
1:A:97:ARG:HG3	1:B:88:ALA:HB3	2.03	0.40
2:D:501:HEM:HBA2	4:D:503:WK2:C09	2.51	0.40
6:B:506:BTB:H32	6:B:506:BTB:H51	1.67	0.40
1:A:317:HIS:CG	1:A:318:PRO:HD2	2.56	0.40
1:B:238:ARG:HD3	1:B:296:PRO:CG	2.51	0.40
1:C:376:LEU:HB2	12:C:656:HOH:O	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	401/440 (91%)	391 (98%)	10 (2%)	0	100	100
1	B	405/440 (92%)	403 (100%)	2 (0%)	0	100	100
1	C	403/440 (92%)	391 (97%)	11 (3%)	1 (0%)	43	29
1	D	404/440 (92%)	399 (99%)	5 (1%)	0	100	100
All	All	1613/1760 (92%)	1584 (98%)	28 (2%)	1 (0%)	48	33

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	259	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	344/373 (92%)	337 (98%)	7 (2%)	48	29
1	B	348/373 (93%)	344 (99%)	4 (1%)	65	51
1	C	346/373 (93%)	339 (98%)	7 (2%)	48	29
1	D	347/373 (93%)	342 (99%)	5 (1%)	59	44
All	All	1385/1492 (93%)	1362 (98%)	23 (2%)	53	36

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	128	ARG
1	A	200	ASP
1	A	216	LYS
1	A	238	ARG
1	A	302	LEU
1	A	308	GLU
1	A	309	LEU
1	B	79	ILE
1	B	230	VAL
1	B	258	ASP

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Mol	Chain	Res	Type
1	B	326	LEU
1	C	89	GLN
1	C	124	LEU
1	C	164	GLN
1	C	216	LYS
1	C	257	GLN
1	C	342	GLU
1	C	417	ILE
1	D	90	GLN
1	D	97	ARG
1	D	121	GLU
1	D	238	ARG
1	D	258	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	126	GLN
1	B	205	GLN
1	B	370	ASN
1	C	126	GLN
1	C	256	GLN
1	D	126	GLN
1	D	205	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 46 ligands modelled in this entry, 13 are monoatomic - leaving 33 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
2	HEM	C	501	1	50,50,50	1.66	9 (18%)	67,82,82	1.24	5 (7%)
5	ACT	C	504	-	3,3,3	0.80	0	3,3,3	0.67	0
6	BTB	B	507	-	13,13,13	0.53	0	7,16,16	1.20	1 (14%)
5	ACT	C	505	-	3,3,3	0.91	0	3,3,3	1.04	0
7	GOL	A	507	-	5,5,5	0.33	0	5,5,5	0.94	0
4	WK2	D	503	-	27,27,27	1.00	1 (3%)	35,39,39	1.32	5 (14%)
3	H4B	A	502	-	17,18,18	0.86	0	14,26,26	1.80	4 (28%)
2	HEM	D	501	1	50,50,50	1.62	9 (18%)	67,82,82	1.25	5 (7%)
6	BTB	D	506	-	13,13,13	0.38	0	7,16,16	0.81	0
3	H4B	C	502	-	17,18,18	0.87	0	14,26,26	1.81	4 (28%)
7	GOL	B	508	-	5,5,5	0.40	0	5,5,5	0.18	0
5	ACT	D	504	-	3,3,3	0.83	0	3,3,3	0.66	0
6	BTB	B	506	11	13,13,13	0.37	0	7,16,16	0.83	0
4	WK2	C	503	-	27,27,27	1.00	1 (3%)	35,39,39	1.22	4 (11%)
3	H4B	D	502	-	17,18,18	0.85	0	14,26,26	1.80	4 (28%)
5	ACT	B	501	-	3,3,3	0.92	0	3,3,3	0.94	0
5	ACT	A	504	-	3,3,3	0.84	0	3,3,3	0.76	0
4	WK2	A	503	-	27,27,27	0.95	1 (3%)	35,39,39	1.44	5 (14%)
7	GOL	C	509	-	5,5,5	0.44	0	5,5,5	0.45	0
7	GOL	D	508	-	5,5,5	0.50	0	5,5,5	0.34	0
4	WK2	B	504	-	27,27,27	0.98	0	35,39,39	1.23	4 (11%)
6	BTB	C	507	11	13,13,13	0.41	0	7,16,16	0.83	0
6	BTB	A	505	-	13,13,13	0.68	0	7,16,16	0.77	0
7	GOL	A	506	-	5,5,5	0.38	0	5,5,5	0.54	0
5	ACT	B	505	-	3,3,3	0.84	0	3,3,3	0.76	0
7	GOL	C	510	-	5,5,5	0.36	0	5,5,5	0.42	0
7	GOL	C	508	-	5,5,5	0.33	0	5,5,5	0.62	0
6	BTB	D	505	11	13,13,13	0.44	0	7,16,16	0.45	0
7	GOL	D	507	-	5,5,5	0.41	0	5,5,5	0.34	0
3	H4B	B	503	-	17,18,18	0.98	1 (5%)	14,26,26	1.66	3 (21%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	HEM	A	501	1	50,50,50	1.66	9 (18%)	67,82,82	1.33	6 (8%)
6	BTB	C	506	11	13,13,13	0.34	0	7,16,16	0.54	0
2	HEM	B	502	1	50,50,50	1.64	10 (20%)	67,82,82	1.38	9 (13%)

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	HEM	C	501	1	-	1/14/54/54	-
6	BTB	B	507	-	-	7/21/21/21	-
7	GOL	A	507	-	-	4/4/4/4	-
4	WK2	D	503	-	-	0/4/14/14	0/3/4/4
3	H4B	A	502	-	-	3/8/17/17	0/2/2/2
2	HEM	D	501	1	-	3/14/54/54	-
6	BTB	D	506	-	-	12/21/21/21	-
3	H4B	C	502	-	-	2/8/17/17	0/2/2/2
7	GOL	B	508	-	-	2/4/4/4	-
6	BTB	B	506	11	-	5/21/21/21	-
4	WK2	C	503	-	-	0/4/14/14	0/3/4/4
3	H4B	D	502	-	-	2/8/17/17	0/2/2/2
4	WK2	A	503	-	-	0/4/14/14	0/3/4/4
7	GOL	C	509	-	-	4/4/4/4	-
7	GOL	D	508	-	-	2/4/4/4	-
4	WK2	B	504	-	-	0/4/14/14	0/3/4/4
6	BTB	C	507	11	-	0/21/21/21	-
6	BTB	A	505	-	-	7/21/21/21	-
7	GOL	A	506	-	-	3/4/4/4	-
7	GOL	C	510	-	-	4/4/4/4	-
7	GOL	C	508	-	-	4/4/4/4	-
6	BTB	D	505	11	-	5/21/21/21	-
7	GOL	D	507	-	-	4/4/4/4	-
3	H4B	B	503	-	-	0/8/17/17	0/2/2/2
2	HEM	A	501	1	-	3/14/54/54	-
6	BTB	C	506	11	-	1/21/21/21	-
2	HEM	B	502	1	-	3/14/54/54	-

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	501	HEM	FE-NC	5.40	2.13	1.95
2	D	501	HEM	FE-NB	5.37	2.11	1.94
2	C	501	HEM	FE-NB	5.15	2.10	1.94
2	B	502	HEM	FE-ND	4.68	2.09	1.94
2	A	501	HEM	FE-NB	4.58	2.09	1.94
2	B	502	HEM	FE-NB	4.48	2.08	1.94
2	B	502	HEM	FE-NC	4.41	2.09	1.95
2	A	501	HEM	FE-NC	4.39	2.09	1.95
2	A	501	HEM	FE-NA	4.17	2.08	1.95
2	D	501	HEM	FE-NC	4.12	2.08	1.95
2	A	501	HEM	FE-ND	3.91	2.06	1.94
2	D	501	HEM	FE-ND	3.83	2.06	1.94
2	C	501	HEM	FE-ND	3.59	2.05	1.94
2	D	501	HEM	CAC-C3C	3.18	1.55	1.47
2	A	501	HEM	CAB-C3B	3.15	1.55	1.47
2	C	501	HEM	CAB-C3B	3.11	1.55	1.47
2	B	502	HEM	FE-NA	3.07	2.05	1.95
2	C	501	HEM	CAC-C3C	3.06	1.55	1.47
2	A	501	HEM	CAC-C3C	3.02	1.55	1.47
2	B	502	HEM	CAB-C3B	2.92	1.55	1.47
2	D	501	HEM	CAB-C3B	2.87	1.55	1.47
2	B	502	HEM	CAC-C3C	2.83	1.54	1.47
2	D	501	HEM	FE-NA	2.75	2.04	1.95
2	D	501	HEM	CMB-C2B	2.68	1.56	1.50
2	C	501	HEM	CMC-C2C	2.45	1.55	1.50
2	B	502	HEM	CMD-C2D	2.45	1.55	1.50
4	C	503	WK2	C05-C10	-2.44	1.38	1.42
2	A	501	HEM	CMC-C2C	2.31	1.55	1.50
3	B	503	H4B	C4-N3	-2.29	1.34	1.38
2	D	501	HEM	CMA-C3A	2.29	1.55	1.50
2	B	502	HEM	CMC-C2C	2.23	1.55	1.50
2	B	502	HEM	C2A-C3A	-2.23	1.33	1.38
2	B	502	HEM	CMB-C2B	2.22	1.55	1.50
2	C	501	HEM	C2A-C3A	-2.19	1.33	1.38
2	C	501	HEM	CMB-C2B	2.17	1.55	1.50
2	C	501	HEM	FE-NA	2.16	2.02	1.95
2	D	501	HEM	CMD-C2D	2.13	1.55	1.50
4	A	503	WK2	C05-C10	-2.09	1.39	1.42
2	A	501	HEM	CMB-C2B	2.06	1.55	1.50
4	D	503	WK2	C09-C10	-2.02	1.38	1.41
2	A	501	HEM	C2A-C3A	-2.00	1.33	1.38

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	503	WK2	C25-C31-N30	5.02	124.56	115.09
3	D	502	H4B	C2-N1-C8A	4.14	120.67	113.36
3	C	502	H4B	C2-N1-C8A	3.98	120.38	113.36
3	A	502	H4B	C2-N1-C8A	3.83	120.12	113.36
3	B	503	H4B	C2-N1-C8A	3.65	119.80	113.36
4	B	504	WK2	C05-C10-N01	-3.42	119.18	122.80
2	D	501	HEM	C2A-C1A-NA	-3.39	106.39	110.15
2	B	502	HEM	CHC-C1C-NC	3.31	128.06	124.45
2	A	501	HEM	C2A-C1A-NA	-3.27	106.52	110.15
4	D	503	WK2	C28-O27-C24	3.19	121.42	116.13
2	B	502	HEM	C2A-C1A-NA	-3.17	106.64	110.15
3	C	502	H4B	O4-C4-C4A	-3.13	119.72	127.26
4	C	503	WK2	C28-O27-C24	3.02	121.14	116.13
4	A	503	WK2	C05-C10-N01	-3.01	119.62	122.80
3	A	502	H4B	O4-C4-C4A	-2.96	120.13	127.26
4	D	503	WK2	C05-C10-N01	-2.96	119.67	122.80
2	A	501	HEM	C3D-C4D-ND	-2.95	106.93	110.17
2	C	501	HEM	C3B-C4B-NB	-2.93	107.36	109.47
4	A	503	WK2	C31-C25-C24	2.88	124.90	119.82
3	D	502	H4B	O4-C4-C4A	-2.77	120.58	127.26
2	D	501	HEM	CHC-C1C-NC	2.69	127.39	124.45
4	B	504	WK2	C04-C05-C10	2.67	119.63	118.00
3	B	503	H4B	O4-C4-C4A	-2.64	120.90	127.26
2	A	501	HEM	C4D-ND-C1D	2.61	108.30	105.21
2	A	501	HEM	C3B-C2B-C1B	2.58	108.35	106.41
3	C	502	H4B	C4A-C4-N3	2.52	119.06	112.13
2	B	502	HEM	C4D-ND-C1D	2.52	108.19	105.21
2	C	501	HEM	C3B-C2B-C1B	2.52	108.30	106.41
3	D	502	H4B	C4A-C4-N3	2.51	119.03	112.13
2	D	501	HEM	CBA-CAA-C2A	-2.45	105.76	112.53
2	C	501	HEM	C2A-C1A-NA	-2.44	107.45	110.15
2	B	502	HEM	C3B-C2B-C1B	2.43	108.23	106.41
2	B	502	HEM	CHD-C1D-ND	2.42	127.03	124.42
2	A	501	HEM	C3B-C4B-NB	-2.39	107.75	109.47
4	C	503	WK2	C05-C10-N01	-2.38	120.28	122.80
4	A	503	WK2	C04-C05-C10	2.35	119.43	118.00
4	D	503	WK2	C26-C21-C08	-2.35	116.91	120.84
4	D	503	WK2	C32-N30-C31	-2.32	107.25	111.04
3	A	502	H4B	C4A-C4-N3	2.31	118.48	112.13
2	D	501	HEM	CAA-CBA-CGA	-2.31	107.54	113.67
2	C	501	HEM	C1B-NB-C4B	2.30	107.93	105.21
3	B	503	H4B	C4A-C4-N3	2.26	118.34	112.13

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	503	WK2	C03-C04-C05	2.22	120.08	117.84
2	B	502	HEM	C2D-C1D-ND	-2.20	107.36	109.90
4	B	504	WK2	C32-N30-C31	-2.18	107.47	111.04
2	B	502	HEM	C3D-C4D-ND	-2.18	107.78	110.17
3	C	502	H4B	C2-N3-C4	-2.16	121.20	125.11
3	D	502	H4B	C2-N3-C4	-2.15	121.21	125.11
4	C	503	WK2	C25-C31-N30	-2.15	111.03	115.09
6	B	507	BTB	C8-C7-N	-2.14	103.22	111.59
2	C	501	HEM	C4D-ND-C1D	2.11	107.71	105.21
3	A	502	H4B	C2-N3-C4	-2.11	121.28	125.11
2	B	502	HEM	C3A-C4A-NA	-2.10	106.77	110.14
4	B	504	WK2	C03-C04-C05	2.09	119.95	117.84
4	A	503	WK2	C03-C04-C05	2.03	119.89	117.84
2	B	502	HEM	C4A-NA-C1A	2.02	109.12	105.82
4	C	503	WK2	C04-C05-C10	2.02	119.23	118.00
2	D	501	HEM	C4A-NA-C1A	2.00	109.09	105.82
2	A	501	HEM	C1B-NB-C4B	2.00	107.58	105.21

There are no chirality outliers.

All (81) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	H4B	C7-C6-C9-C10
6	A	505	BTB	O1-C1-C2-N
6	B	506	BTB	O1-C1-C2-C3
6	B	506	BTB	O1-C1-C2-C4
6	B	506	BTB	O1-C1-C2-N
6	B	507	BTB	O1-C1-C2-C4
6	B	507	BTB	C1-C2-C4-O4
6	B	507	BTB	N-C2-C4-O4
6	D	505	BTB	O1-C1-C2-C3
6	D	505	BTB	O1-C1-C2-C4
6	D	505	BTB	O1-C1-C2-N
6	D	506	BTB	O1-C1-C2-C3
6	D	506	BTB	O1-C1-C2-C4
6	D	506	BTB	O1-C1-C2-N
6	D	506	BTB	N-C2-C4-O4
6	D	506	BTB	C1-C2-N-C5
6	D	506	BTB	C1-C2-N-C7
6	D	506	BTB	C3-C2-N-C5
6	D	506	BTB	C3-C2-N-C7
6	D	506	BTB	C4-C2-N-C5

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Mol	Chain	Res	Type	Atoms
6	D	506	BTB	C4-C2-N-C7
7	A	506	GOL	O1-C1-C2-O2
7	A	506	GOL	O1-C1-C2-C3
7	A	507	GOL	C1-C2-C3-O3
7	B	508	GOL	C1-C2-C3-O3
7	C	508	GOL	C1-C2-C3-O3
7	C	509	GOL	O1-C1-C2-C3
7	C	509	GOL	C1-C2-C3-O3
7	C	510	GOL	O1-C1-C2-C3
7	D	507	GOL	O1-C1-C2-C3
7	D	507	GOL	C1-C2-C3-O3
7	D	508	GOL	C1-C2-C3-O3
2	B	502	HEM	C2A-CAA-CBA-CGA
6	B	506	BTB	N-C5-C6-O6
6	D	506	BTB	N-C7-C8-O8
7	A	507	GOL	O2-C2-C3-O3
7	B	508	GOL	O2-C2-C3-O3
7	C	509	GOL	O1-C1-C2-O2
7	D	507	GOL	O1-C1-C2-O2
6	D	506	BTB	N-C5-C6-O6
3	A	502	H4B	C7-C6-C9-O9
6	B	506	BTB	N-C7-C8-O8
2	A	501	HEM	C2A-CAA-CBA-CGA
7	A	507	GOL	O1-C1-C2-C3
7	C	508	GOL	O1-C1-C2-C3
7	C	510	GOL	C1-C2-C3-O3
6	B	507	BTB	N-C5-C6-O6
7	A	507	GOL	O1-C1-C2-O2
7	C	510	GOL	O1-C1-C2-O2
7	D	508	GOL	O2-C2-C3-O3
2	C	501	HEM	C4B-C3B-CAB-CBB
7	C	508	GOL	O1-C1-C2-O2
7	C	508	GOL	O2-C2-C3-O3
7	C	509	GOL	O2-C2-C3-O3
7	D	507	GOL	O2-C2-C3-O3
6	D	505	BTB	C1-C2-C4-O4
7	C	510	GOL	O2-C2-C3-O3
6	A	505	BTB	N-C5-C6-O6
6	C	506	BTB	N-C7-C8-O8
6	A	505	BTB	C4-C2-N-C5
6	B	507	BTB	O1-C1-C2-N
6	A	505	BTB	N-C7-C8-O8

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Mol	Chain	Res	Type	Atoms
6	B	507	BTB	O1-C1-C2-C3
2	D	501	HEM	C4B-C3B-CAB-CBB
6	B	507	BTB	N-C7-C8-O8
2	A	501	HEM	CAA-CBA-CGA-O2A
2	A	501	HEM	CAA-CBA-CGA-O1A
7	A	506	GOL	O2-C2-C3-O3
2	B	502	HEM	CAA-CBA-CGA-O2A
2	D	501	HEM	CAA-CBA-CGA-O2A
3	C	502	H4B	C7-C6-C9-C10
3	D	502	H4B	C7-C6-C9-C10
2	B	502	HEM	CAA-CBA-CGA-O1A
2	D	501	HEM	CAA-CBA-CGA-O1A
6	A	505	BTB	O1-C1-C2-C3
3	A	502	H4B	N5-C6-C9-O9
3	C	502	H4B	N5-C6-C9-O9
3	D	502	H4B	N5-C6-C9-O9
6	A	505	BTB	C1-C2-N-C5
6	A	505	BTB	C3-C2-N-C5
6	D	505	BTB	N-C2-C4-O4

There are no ring outliers.

20 monomers are involved in 53 short contacts:

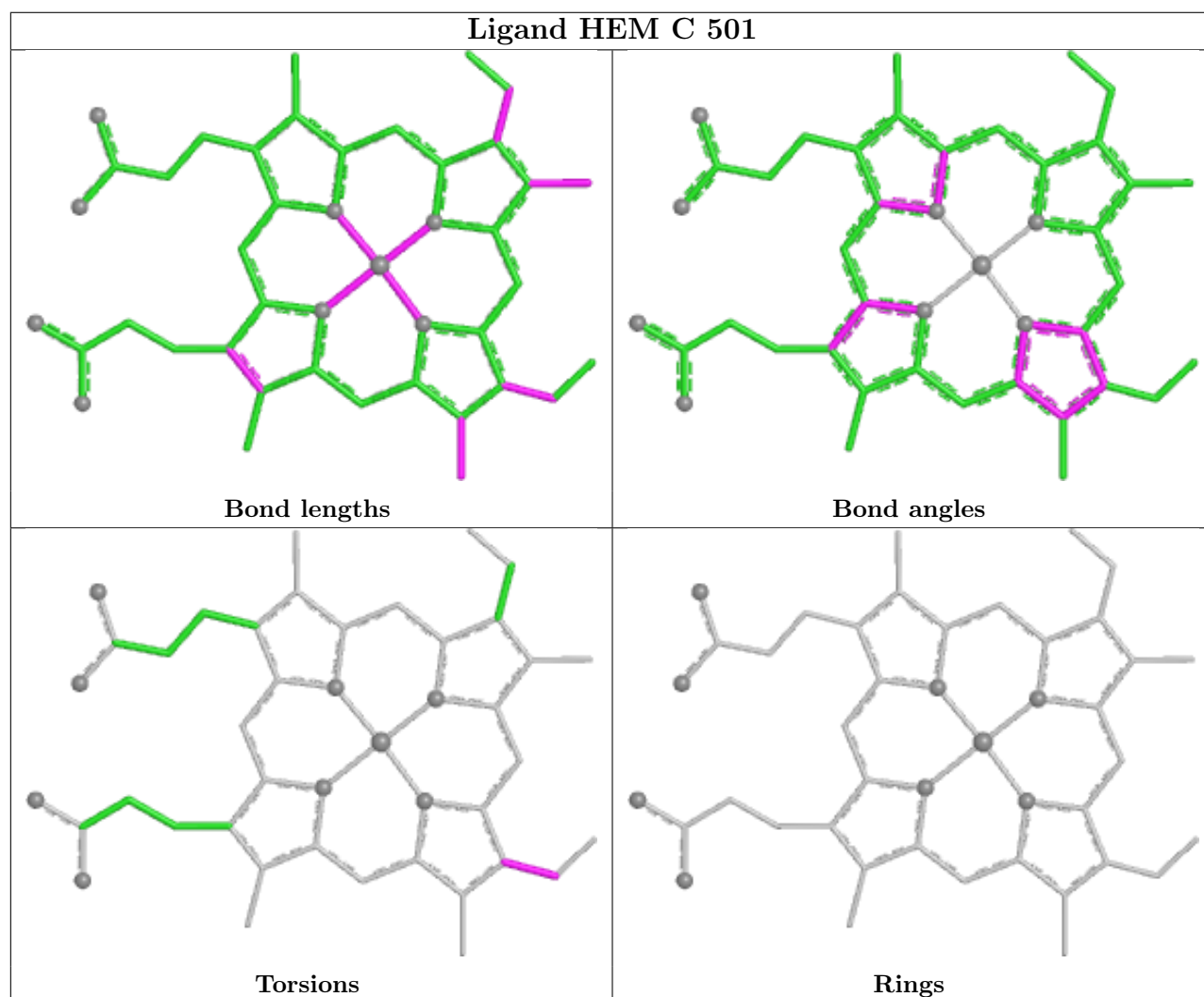
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	501	HEM	3	0
6	B	507	BTB	1	0
5	C	505	ACT	5	0
7	A	507	GOL	1	0
4	D	503	WK2	5	0
2	D	501	HEM	11	0
6	D	506	BTB	6	0
3	C	502	H4B	2	0
5	D	504	ACT	1	0
6	B	506	BTB	2	0
4	C	503	WK2	1	0
3	D	502	H4B	4	0
5	B	501	ACT	8	0
4	A	503	WK2	1	0
4	B	504	WK2	2	0
6	A	505	BTB	2	0
6	D	505	BTB	2	0
2	A	501	HEM	3	0

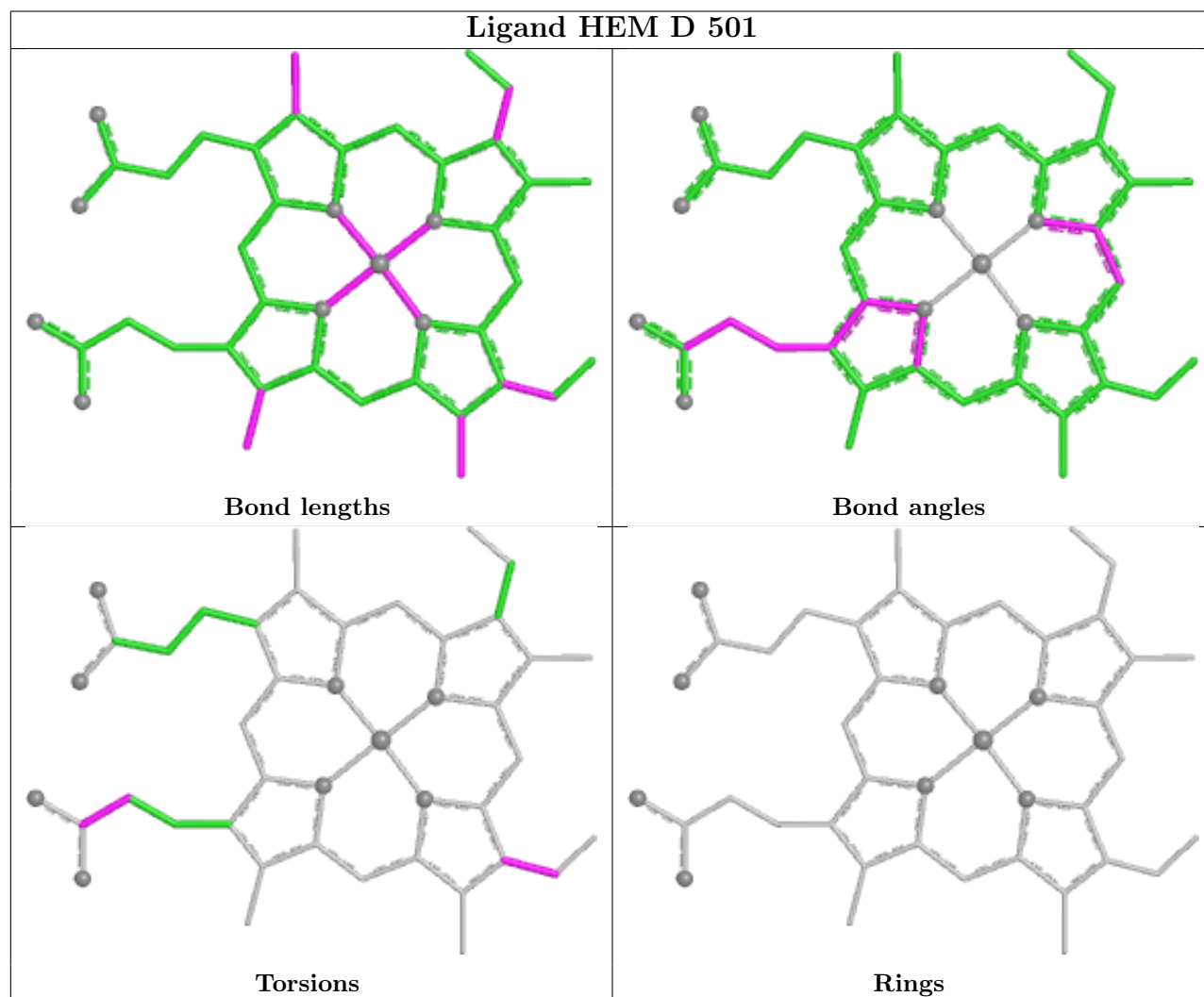
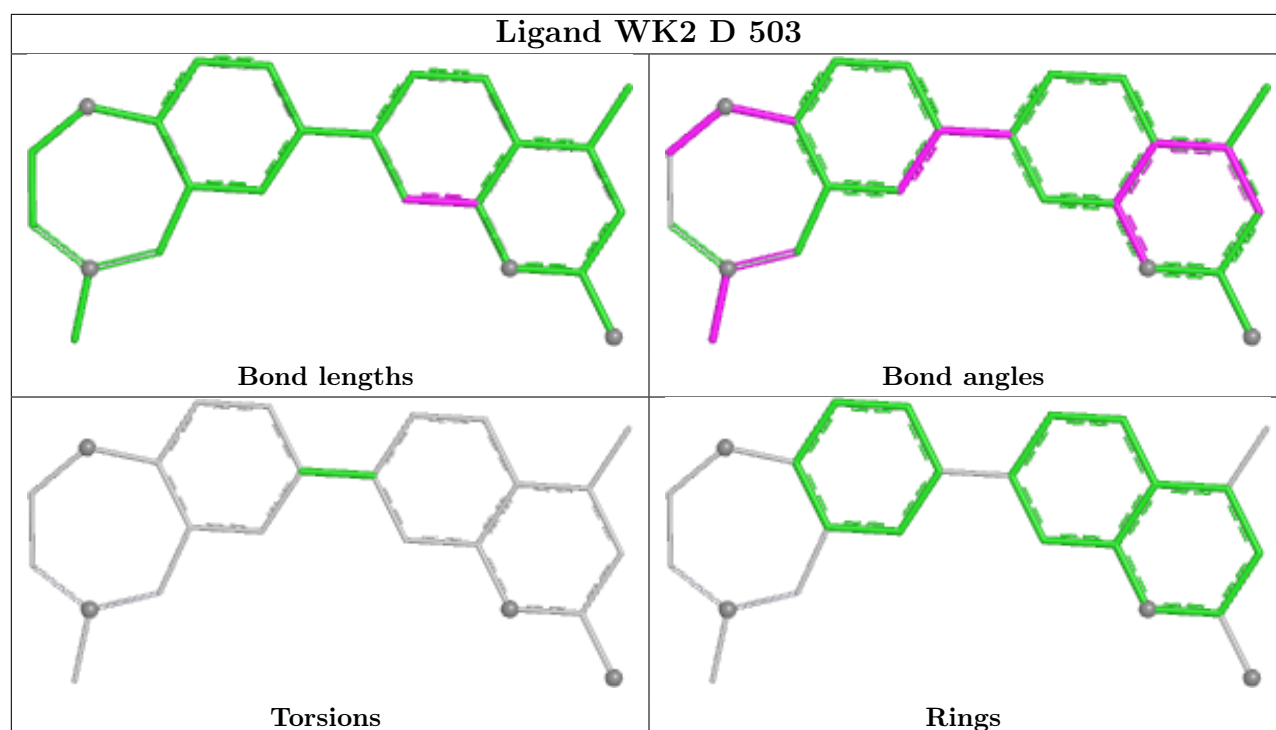
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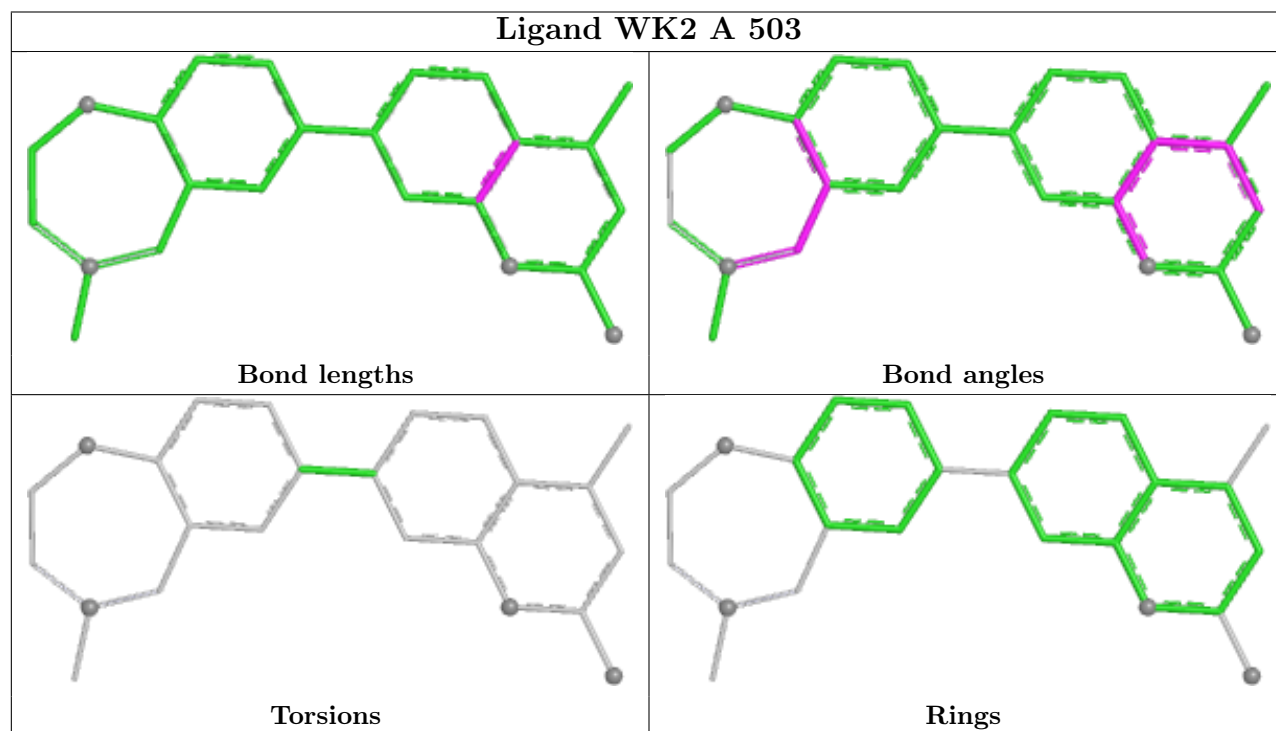
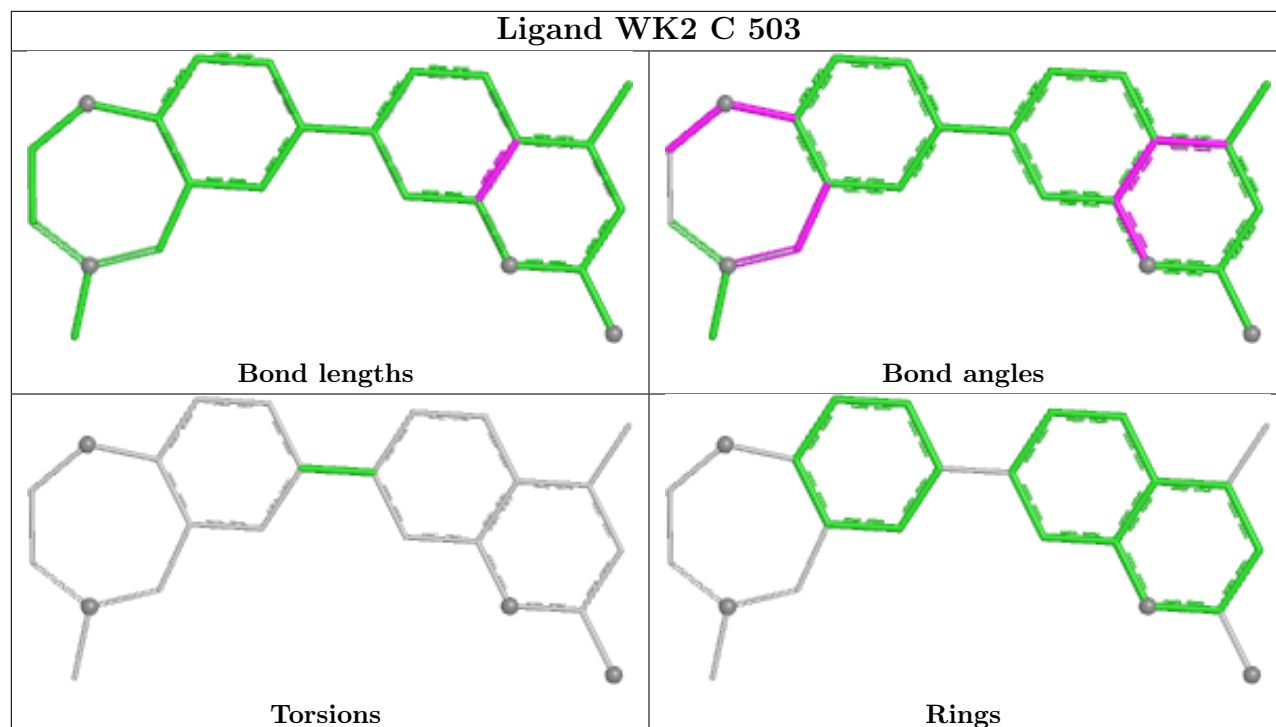
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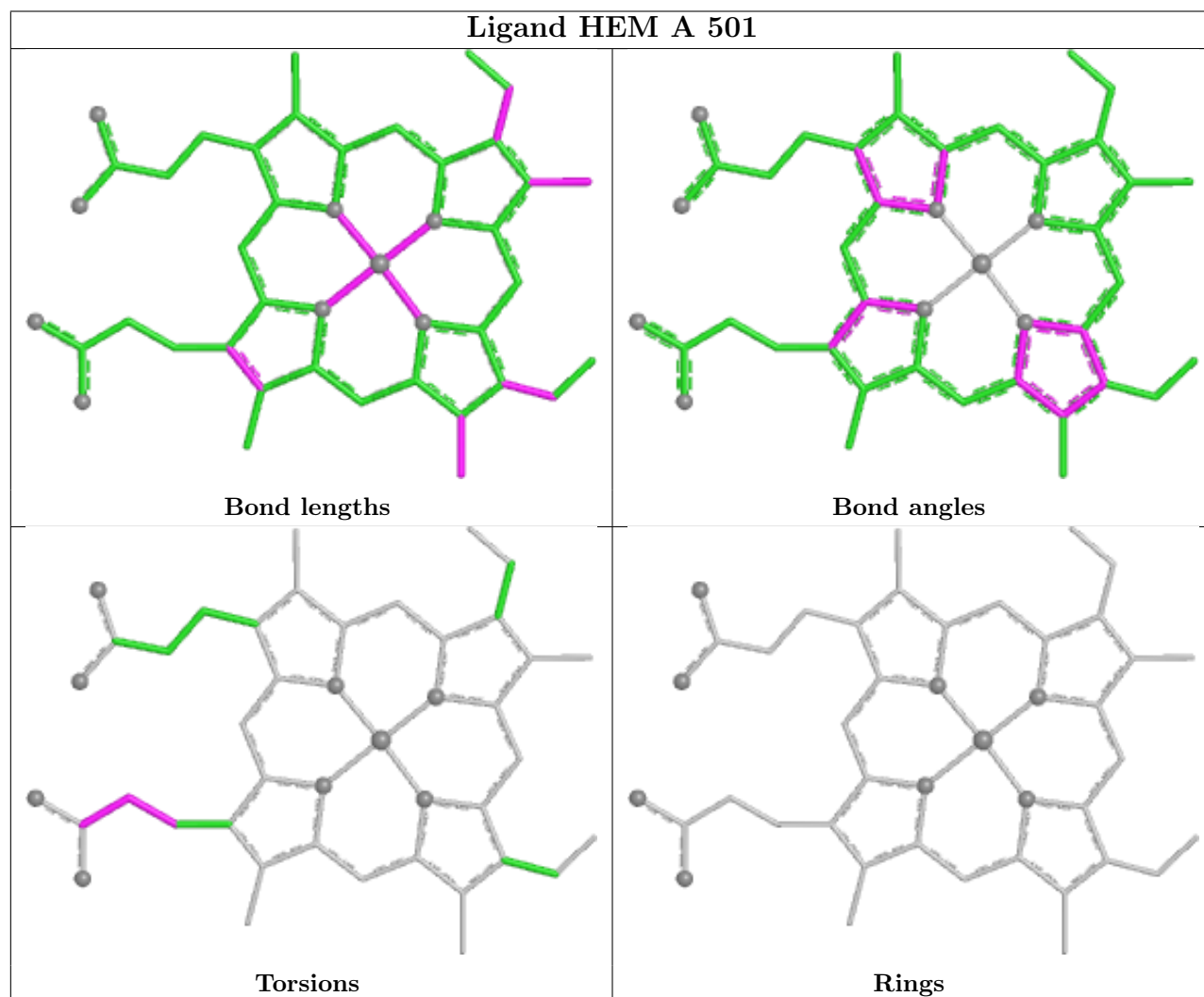
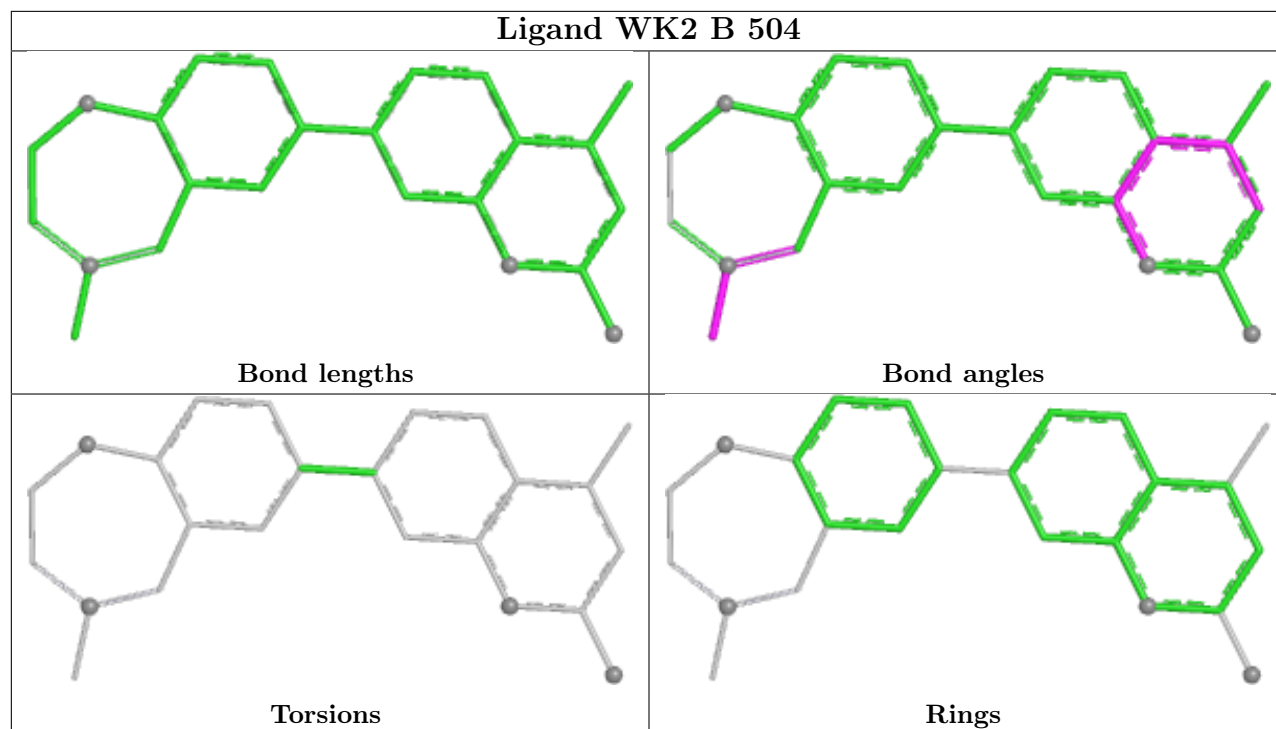
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	506	BTB	1	0
2	B	502	HEM	3	0

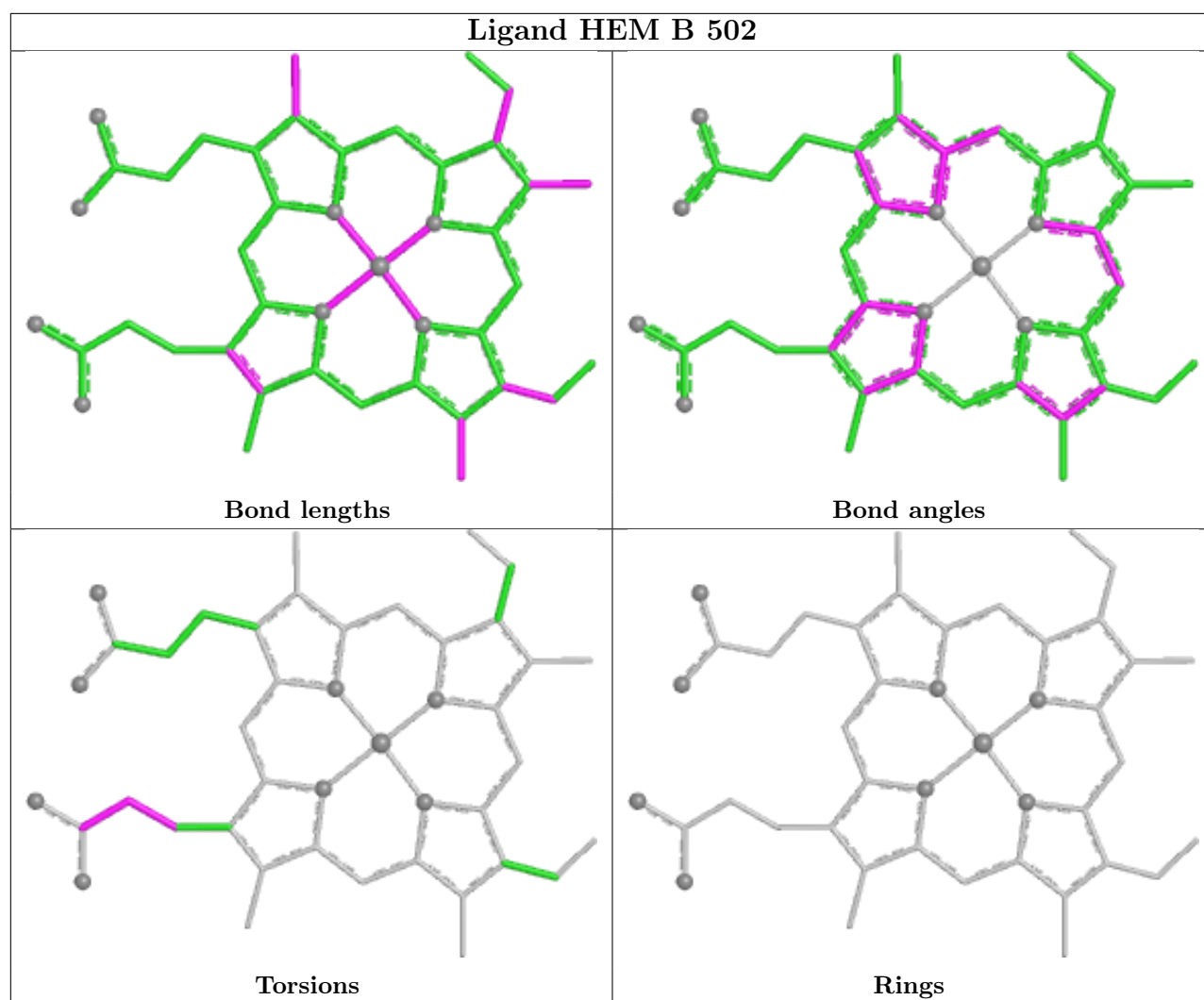
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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











5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	403/440 (91%)	0.20	14 (3%) 47 53	18, 39, 74, 94	2 (0%)
1	B	403/440 (91%)	-0.27	4 (0%) 79 85	15, 26, 51, 72	6 (1%)
1	C	403/440 (91%)	0.14	9 (2%) 62 70	18, 39, 71, 95	4 (0%)
1	D	403/440 (91%)	-0.26	6 (1%) 72 78	16, 27, 54, 81	5 (1%)
All	All	1612/1760 (91%)	-0.05	33 (2%) 65 73	15, 32, 67, 95	17 (1%)

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	66	PRO	5.0
1	C	66	PRO	4.8
1	B	66	PRO	4.6
1	B	119	ALA	4.6
1	B	106	PRO	4.3
1	D	119	ALA	4.1
1	A	66	PRO	4.1
1	A	106	PRO	3.9
1	C	119	ALA	3.7
1	A	119	ALA	3.3
1	C	120	PRO	3.0
1	C	89	GLN	3.0
1	A	159	ALA	2.9
1	D	106	PRO	2.7
1	A	120	PRO	2.6
1	C	304	LEU	2.6
1	B	257[A]	GLN	2.5
1	D	89	GLN	2.4
1	C	237	GLY	2.4
1	C	106	PRO	2.4
1	A	160	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	153	VAL	2.3
1	A	105	PHE	2.3
1	A	298	GLU	2.3
1	D	257	GLN	2.2
1	A	90	GLN	2.1
1	A	257	GLN	2.1
1	C	258	ASP	2.1
1	C	285	ARG	2.1
1	A	468	PHE	2.1
1	A	259	GLY	2.1
1	A	237	GLY	2.0
1	D	144	GLN	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 oligosaccharides 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
7	GOL	A	506	6/6	0.53	0.14	71,77,80,83	0
7	GOL	C	508	6/6	0.65	0.19	53,68,74,76	0
6	BTB	D	506	14/14	0.68	0.19	49,75,88,94	0
6	BTB	B	507	14/14	0.71	0.21	57,66,78,86	0
7	GOL	B	508	6/6	0.76	0.15	48,56,58,65	0
7	GOL	A	507	6/6	0.77	0.17	46,55,57,60	0
6	BTB	C	506	14/14	0.79	0.18	53,77,90,90	0
6	BTB	A	505	14/14	0.82	0.21	21,57,69,78	0
3	H4B	A	502	17/17	0.83	0.16	48,64,74,75	0
7	GOL	D	507	6/6	0.83	0.12	59,63,65,71	0
7	GOL	C	509	6/6	0.84	0.17	47,61,73,76	0

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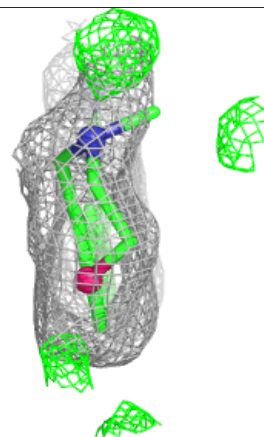
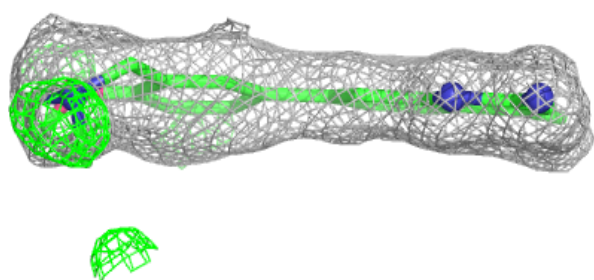
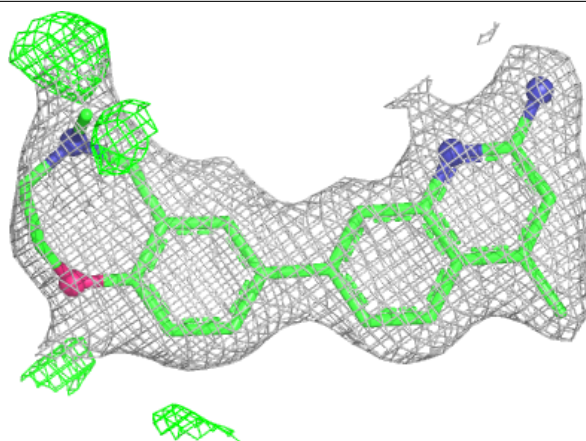
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	BTB	C	507	14/14	0.85	0.13	58,66,81,82	0
5	ACT	C	505	4/4	0.85	0.18	28,31,44,46	0
3	H4B	C	502	17/17	0.86	0.16	39,68,75,75	0
6	BTB	B	506	14/14	0.86	0.17	42,59,68,80	0
3	H4B	B	503	17/17	0.86	0.14	41,54,67,67	0
7	GOL	D	508	6/6	0.86	0.12	38,49,55,56	0
6	BTB	D	505	14/14	0.87	0.17	37,51,75,84	0
7	GOL	C	510	6/6	0.87	0.17	67,69,72,80	0
3	H4B	D	502	17/17	0.88	0.14	32,57,67,69	0
4	WK2	A	503	24/24	0.90	0.12	28,46,67,74	0
4	WK2	C	503	24/24	0.91	0.13	27,45,75,76	0
4	WK2	B	504	24/24	0.94	0.09	19,31,56,58	0
4	WK2	D	503	24/24	0.94	0.10	21,33,67,69	0
5	ACT	B	501	4/4	0.95	0.11	16,17,36,40	0
8	CL	C	511	1/1	0.95	0.10	55,55,55,55	0
5	ACT	C	504	4/4	0.96	0.10	39,39,40,41	0
5	ACT	A	504	4/4	0.96	0.07	34,38,40,41	0
11	GD	C	513	1/1	0.96	0.11	67,67,67,67	0
8	CL	D	509	1/1	0.97	0.07	37,37,37,37	0
5	ACT	D	504	4/4	0.97	0.06	29,33,38,43	0
2	HEM	D	501	43/43	0.98	0.07	16,20,54,75	0
2	HEM	A	501	43/43	0.98	0.09	23,32,68,99	0
8	CL	A	508	1/1	0.98	0.05	45,45,45,45	0
8	CL	B	509	1/1	0.98	0.05	36,36,36,36	0
2	HEM	B	502	43/43	0.98	0.07	15,21,54,73	0
5	ACT	B	505	4/4	0.98	0.06	28,32,37,39	0
10	CA	A	510	1/1	0.98	0.10	25,25,25,25	0
11	GD	B	510	1/1	0.98	0.06	29,29,29,29	0
11	GD	C	512	1/1	0.98	0.12	64,64,64,64	0
2	HEM	C	501	43/43	0.98	0.08	25,34,58,76	0
11	GD	D	510	1/1	0.98	0.06	29,29,29,29	0
10	CA	A	511	1/1	0.99	0.03	20,20,20,20	0
10	CA	B	511	1/1	0.99	0.03	22,22,22,22	0
9	ZN	C	514	1/1	0.99	0.04	22,22,22,22	0
9	ZN	A	509	1/1	1.00	0.05	21,21,21,21	0

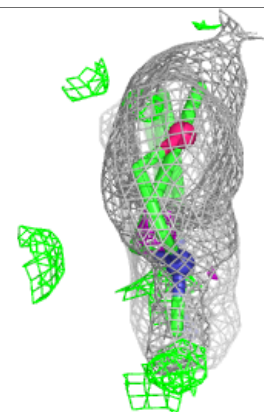
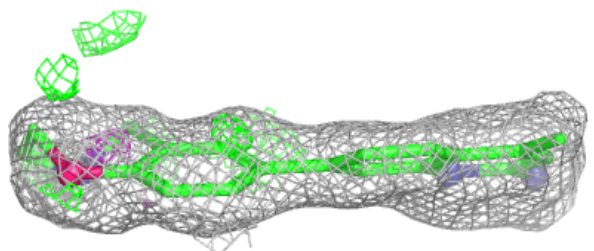
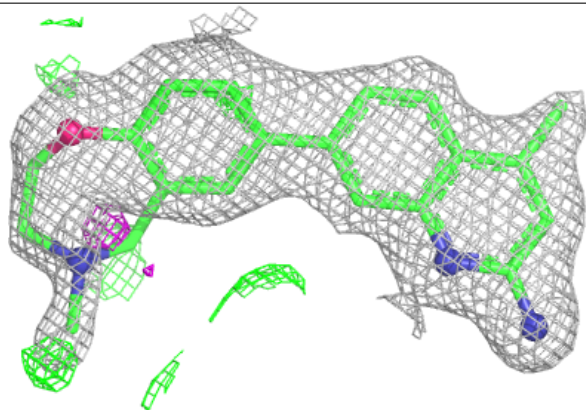
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around WK2 A 503:

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

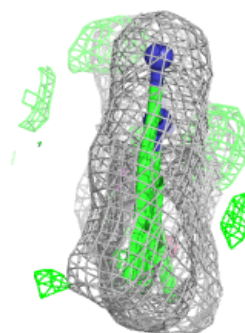
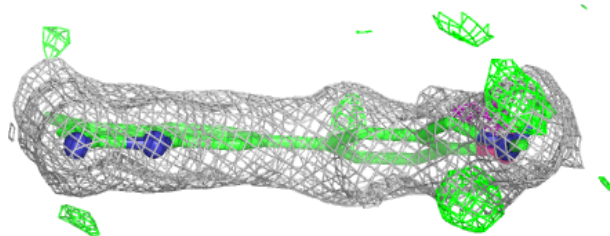
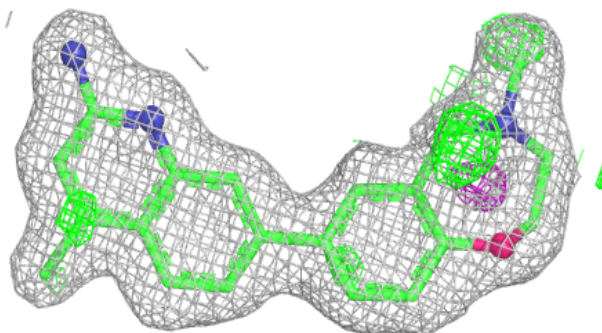
**Electron density around WK2 C 503:**

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

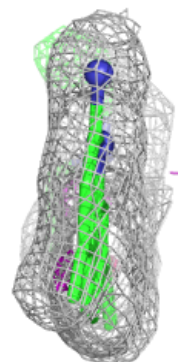
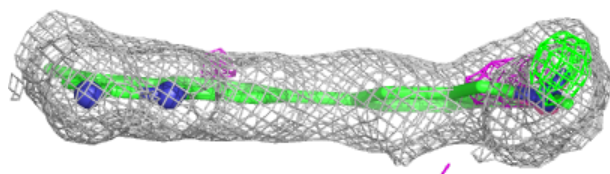
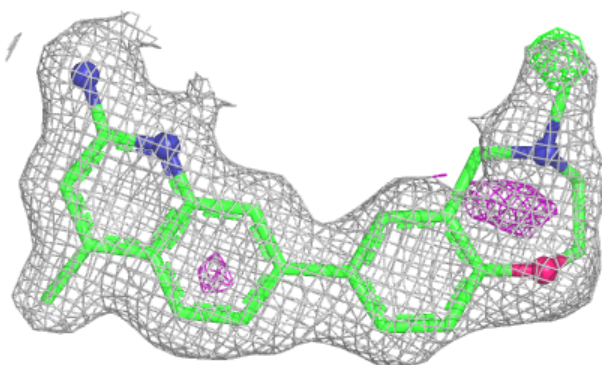


Electron density around WK2 B 504:

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

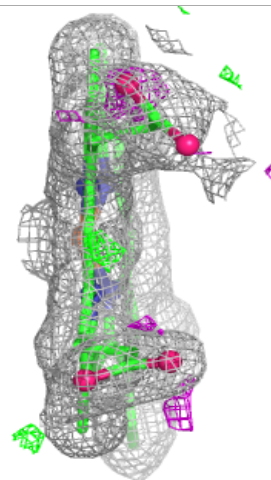
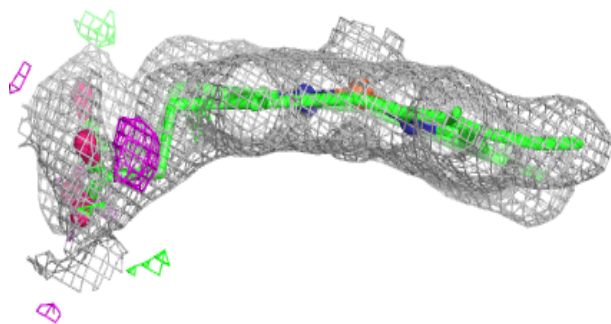
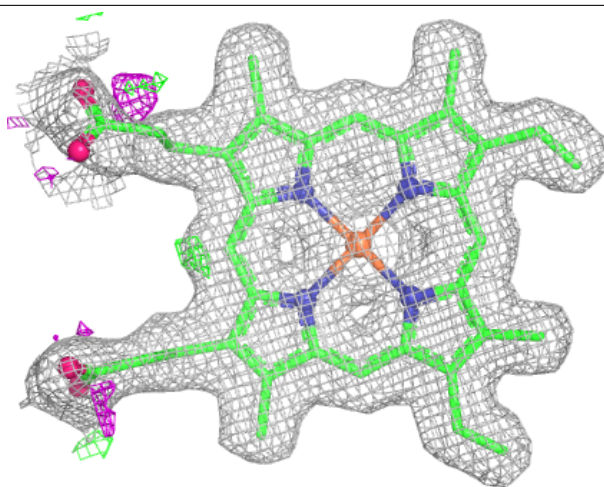
**Electron density around WK2 D 503:**

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



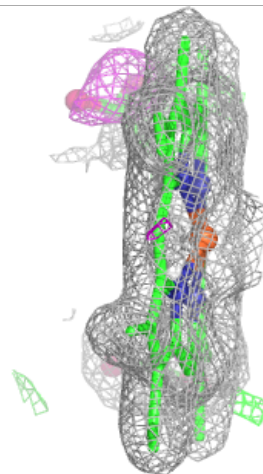
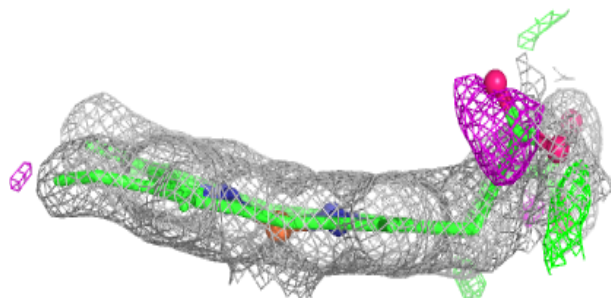
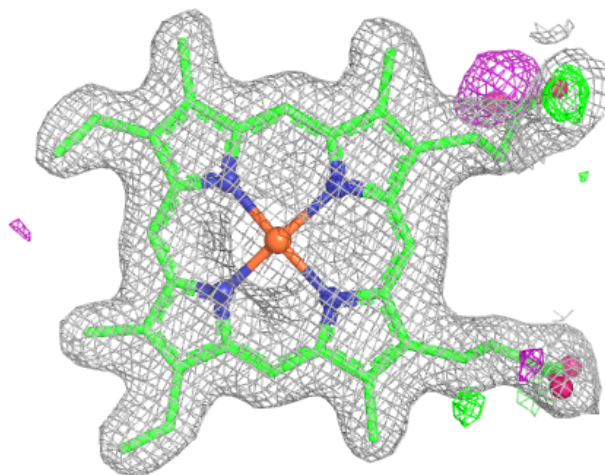
Electron density around HEM D 501:

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



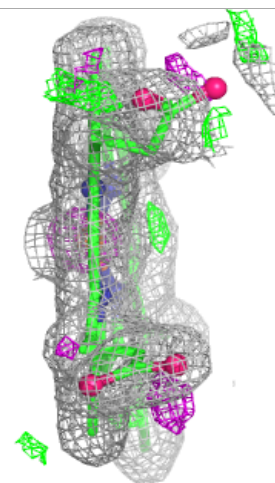
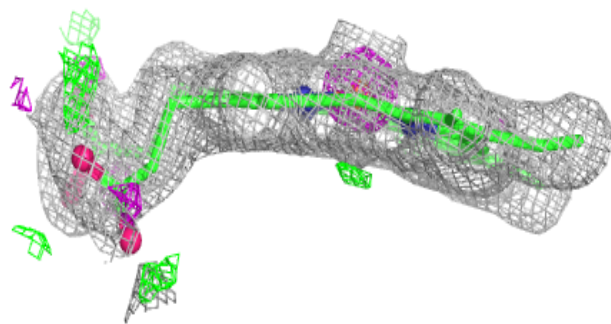
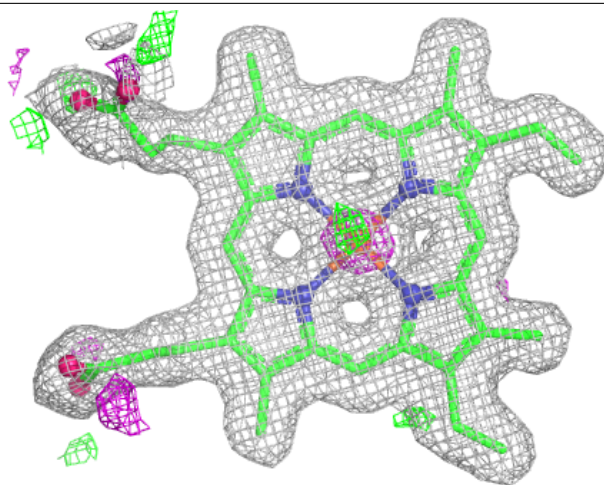
Electron density around HEM A 501:

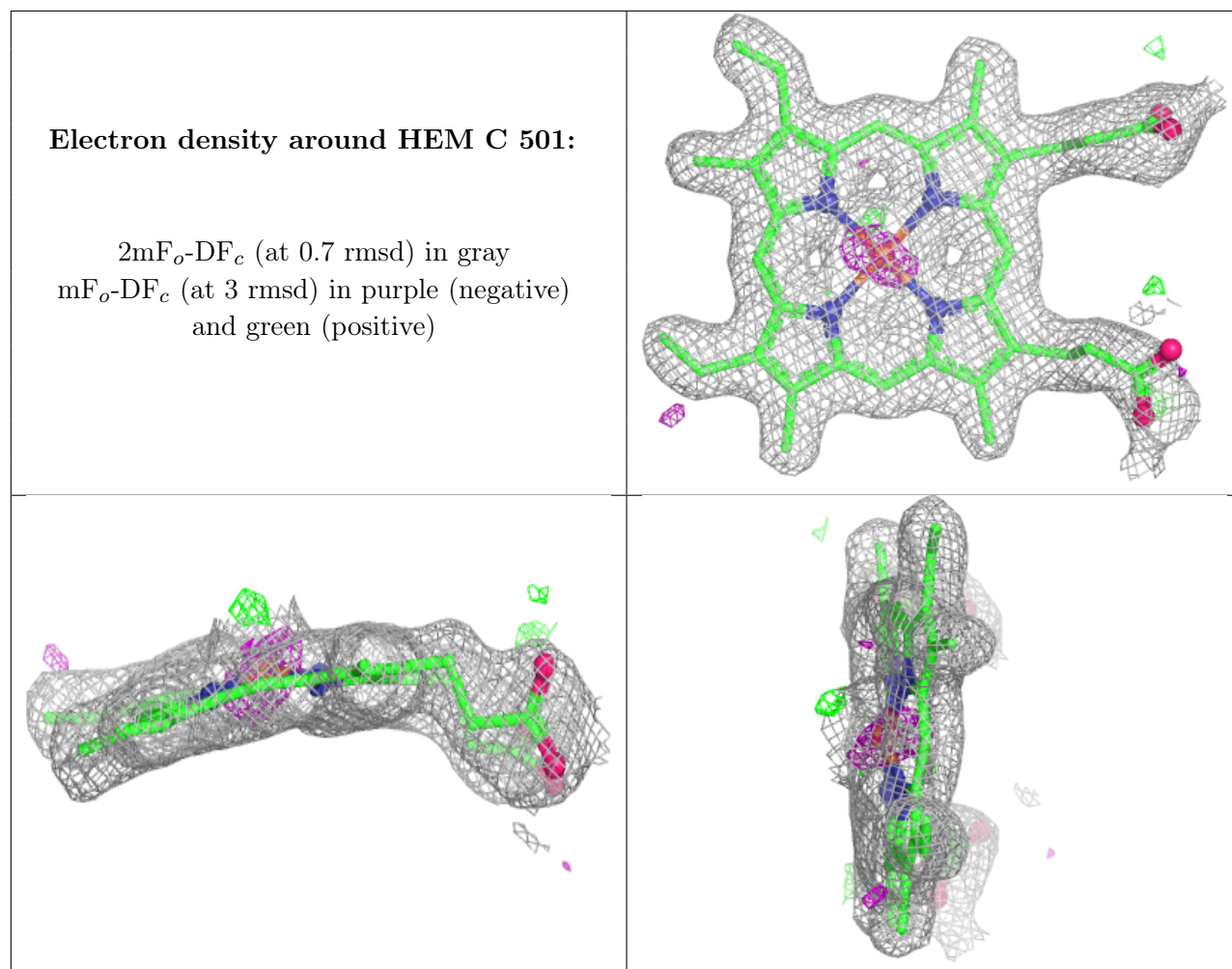
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM B 502:

$2mF_o-DF_c$ (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.