



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 01:45 PM UTC

PDB ID : 8FGQ / pdb_00008fgq
Title : Structure of human endothelial nitric oxide synthase heme domain in complex with 4-(5-(2-(dimethylamino)ethyl)-2,3-difluorophenethyl)-6-methylpyrimidin-2-amine
Authors : Li, H.; Poulos, T.L.
Deposited on : 2022-12-12
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 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

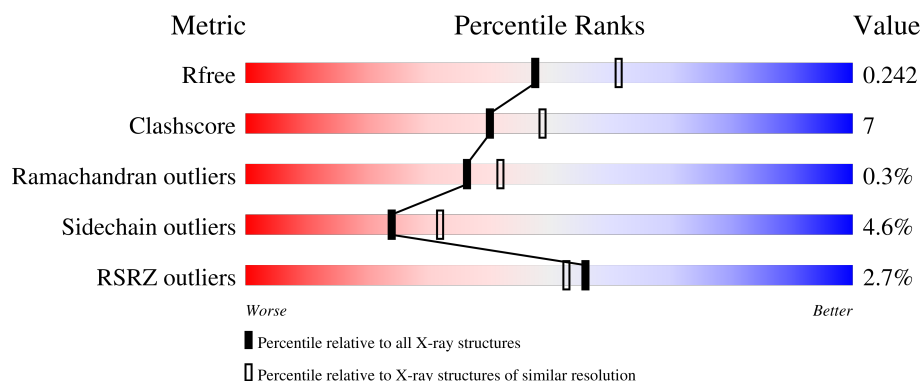
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (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 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	5% 72% 17% • 9%
1	B	440	79% 12% 9%
1	C	440	2% 74% 16% • 9%
1	D	440	• 77% 14% 9%

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 13987 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	400	Total	C	N	O	S	0	2	0
			3204	2041	562	585	16			
1	B	402	Total	C	N	O	S	0	3	0
			3220	2051	566	587	16			
1	C	402	Total	C	N	O	S	0	2	0
			3218	2050	565	587	16			
1	D	401	Total	C	N	O	S	0	2	0
			3206	2042	563	585	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 (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



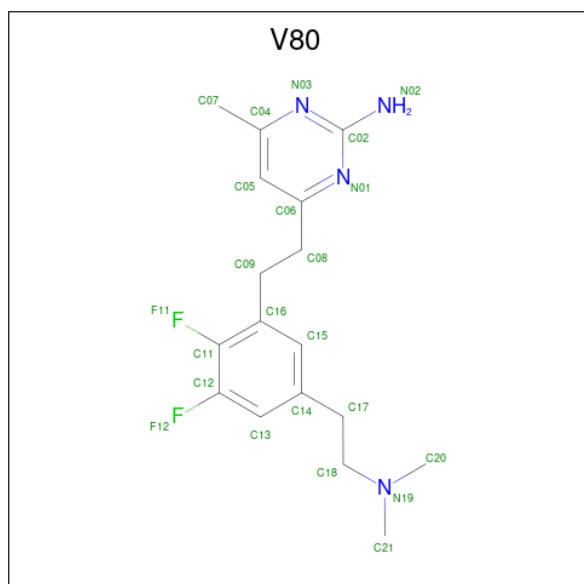
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 3 is 5,6,7,8-TETRAHYDROBIOPTERIN (CCD ID: H4B) (formula: C₉H₁₅N₅O₃).



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 4-(2-{5-[2-(dimethylamino)ethyl]-2,3-difluorophenyl}ethyl)-6-methylpyrimidin-2-amine (CCD ID: V80) (formula: $C_{17}H_{22}F_2N_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	F	N	0	0
			23	17	2	4		
4	B	1	Total	C	F	N	0	0
			23	17	2	4		
4	C	1	Total	C	F	N	0	0
			23	17	2	4		
4	D	1	Total	C	F	N	0	0
			23	17	2	4		

- Molecule 5 is 2-[BIS-(2-HYDROXY-ETHYL)-AMINO]-2-HYDROXYMETHYL-PROPAN E-1,3-DIOL (CCD ID: BTB) (formula: $C_8H_{19}NO_5$).



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

- Molecule 6 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

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

- Molecule 7 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

- Molecule 8 is GADOLINIUM ATOM (CCD ID: GD) (formula: Gd).

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

- Molecule 9 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

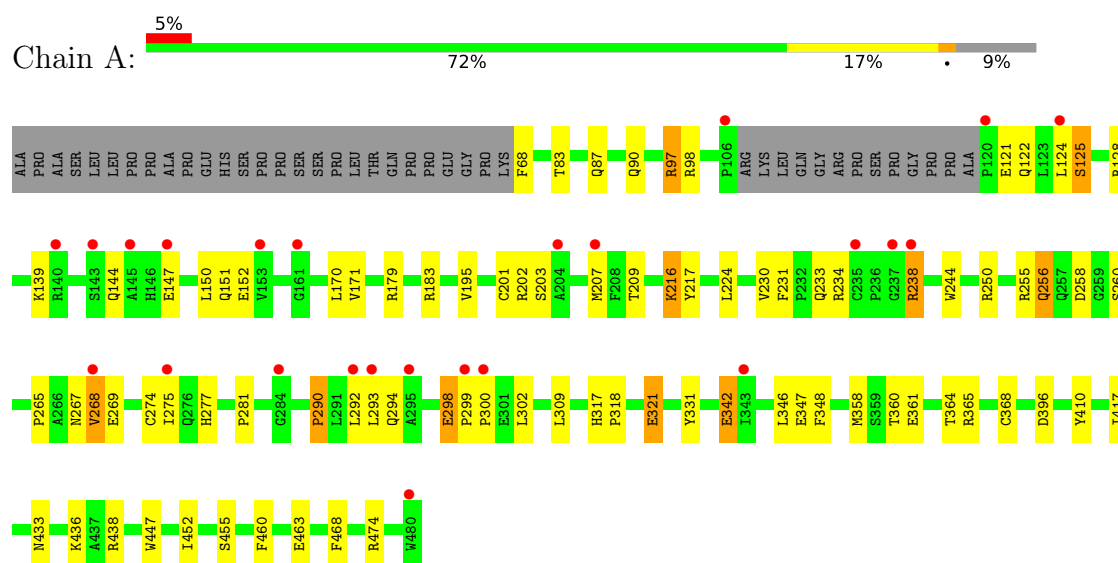
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	112	Total	O	0	0
			112	112		
10	B	182	Total	O	0	0
			182	182		
10	C	147	Total	O	0	0
			147	147		
10	D	196	Total	O	0	0
			196	196		

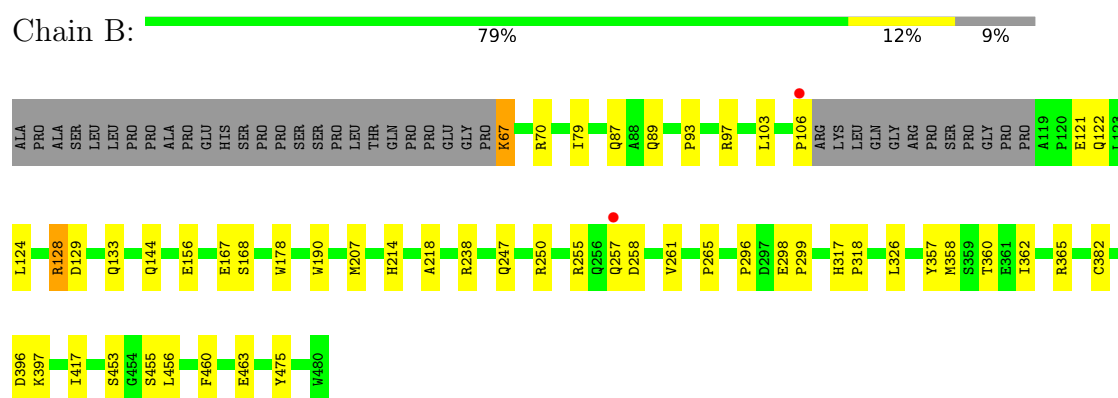
3 Residue-property plots [i](#)

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.

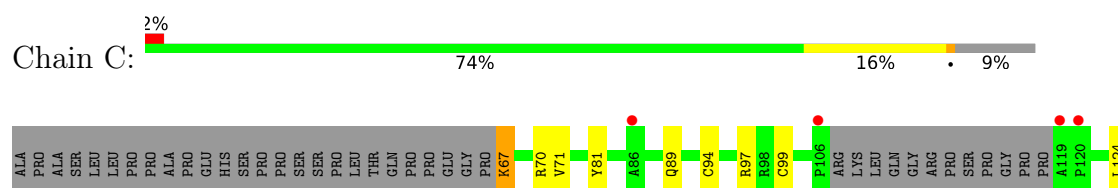
- Molecule 1: Nitric oxide synthase, endothelial

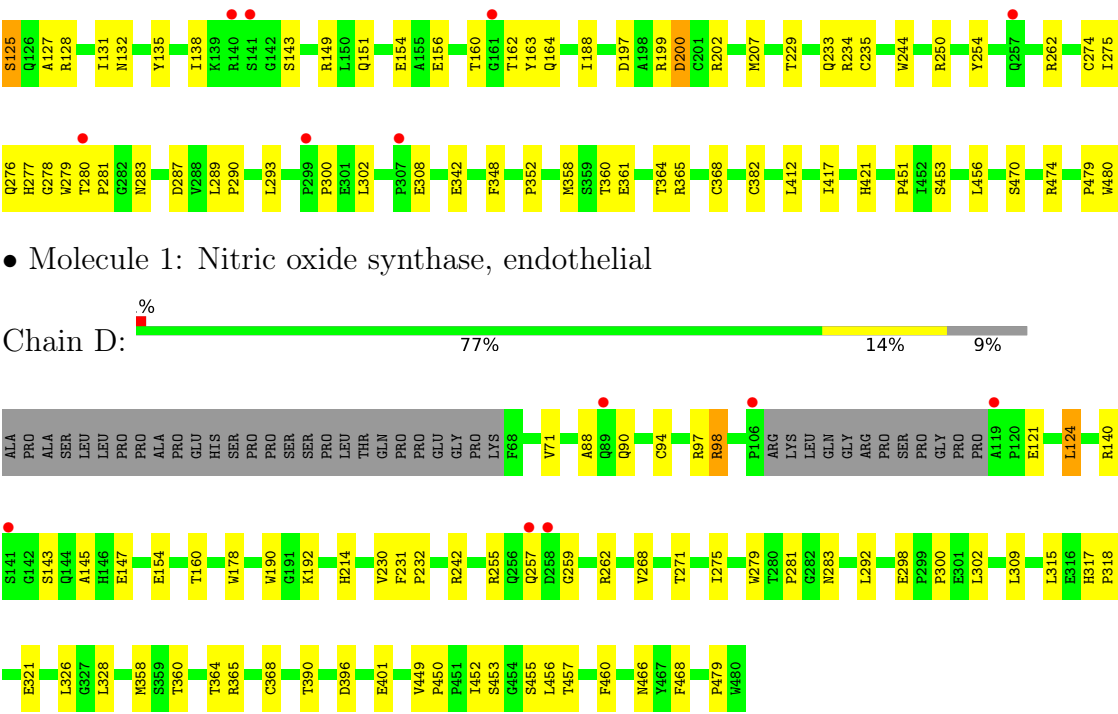


- Molecule 1: Nitric oxide synthase, endothelial



- 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	59.57Å 151.48Å 108.35Å 90.00° 91.08° 90.00°	Depositor
Resolution (Å)	75.78 – 2.20 75.78 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.2 (75.78-2.20) 99.9 (75.78-2.20)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.54 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.193 , 0.246 0.186 , 0.242	Depositor DCC
R_{free} test set	4853 reflections (2.56%)	wwPDB-VP
Wilson B-factor (Å ²)	28.3	Xtriage
Anisotropy	0.765	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 56.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.125 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13987	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% 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: H4B, ZN, GD, GOL, HEM, CL, BTB, V80

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	0.31	0/3302	0.53	0/4499
1	B	0.36	0/3321	0.55	0/4525
1	C	0.31	0/3316	0.51	0/4518
1	D	0.39	0/3304	0.59	0/4503
All	All	0.35	0/13243	0.55	0/18045

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	3204	0	3105	48	1
1	B	3220	0	3127	27	0
1	C	3218	0	3122	48	0
1	D	3206	0	3108	37	0
2	A	43	0	30	6	0
2	B	43	0	30	3	0
2	C	43	0	30	3	0
2	D	43	0	30	3	0
3	A	17	0	15	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	17	0	15	2	0
3	C	17	0	15	1	0
3	D	17	0	15	1	0
4	A	23	0	0	1	0
4	B	23	0	0	0	0
4	C	23	0	0	2	0
4	D	23	0	0	1	0
5	A	42	0	57	2	0
5	B	42	0	55	5	0
5	C	42	0	55	7	0
5	D	28	0	36	6	1
6	A	1	0	0	0	0
6	C	1	0	0	0	0
7	A	1	0	0	0	0
7	B	1	0	0	0	0
7	C	1	0	0	0	0
7	D	1	0	0	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	C	6	0	8	0	0
10	A	112	0	0	1	0
10	B	182	0	0	3	0
10	C	147	0	0	2	0
10	D	196	0	0	1	0
All	All	13987	0	12853	174	1

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 7.

All (174) 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:B:247:GLN:HB2	1:B:250:ARG:HD3	1.55	0.87
1:B:475:TYR:OH	2:B:501:HEM:O1D	1.93	0.85
1:D:275:ILE:HD12	1:D:281:PRO:HG3	1.66	0.78
2:D:501:HEM:HBB2	2:D:501:HEM:HHC	1.69	0.74
1:B:298:GLU:OE2	5:B:505:BTB:N	2.22	0.73
1:A:321:GLU:H	1:A:321:GLU:CD	1.95	0.72
1:C:207:MET:HE3	1:C:293:LEU:HB3	1.69	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:93:PRO:HG3	1:B:106:PRO:HB3	1.71	0.71
1:A:447:TRP:HE3	2:A:501:HEM:HBA2	1.57	0.70
2:A:501:HEM:HHA	2:A:501:HEM:HBA1	1.74	0.70
2:B:501:HEM:HHC	2:B:501:HEM:HBB2	1.73	0.68
1:A:275:ILE:HD11	1:A:281:PRO:HB3	1.77	0.67
1:C:262:ARG:NH1	1:C:283:ASN:O	2.28	0.67
1:A:433:ASN:HA	1:A:436:LYS:HE3	1.76	0.67
1:C:365:ARG:HH12	3:C:502:H4B:C4	2.09	0.66
1:A:97:ARG:HG2	1:A:98:ARG:HG2	1.78	0.66
2:B:501:HEM:HBC2	2:B:501:HEM:HMC2	1.78	0.65
1:A:238:ARG:NH2	10:A:603:HOH:O	2.30	0.65
1:C:200:ASP:OD1	1:C:200:ASP:N	2.28	0.65
1:C:342:GLU:HG3	1:C:474:ARG:NH2	2.10	0.65
2:D:501:HEM:HMC2	2:D:501:HEM:HBC2	1.78	0.65
1:D:257:GLN:H	1:D:257:GLN:CD	2.06	0.63
1:C:342:GLU:OE2	1:C:470:SER:OG	2.16	0.63
2:A:501:HEM:HBB2	2:A:501:HEM:HHC	1.81	0.63
2:C:501:HEM:HBB2	2:C:501:HEM:HHC	1.82	0.61
1:A:244:TRP:NE1	1:A:294:GLN:OE1	2.32	0.61
2:A:501:HEM:HMC2	2:A:501:HEM:HBC2	1.84	0.60
1:B:70:ARG:HD2	1:B:79:ILE:HD13	1.84	0.60
1:D:298:GLU:OE2	5:D:505:BTB:H41	2.02	0.59
1:D:321:GLU:OE2	5:D:504:BTB:O4	2.19	0.59
1:C:128:ARG:O	1:C:132:ASN:ND2	2.36	0.59
1:A:447:TRP:CE3	2:A:501:HEM:HBA2	2.37	0.58
1:C:160:THR:HG23	1:C:162:THR:H	1.69	0.58
1:B:238:ARG:HG2	1:B:296:PRO:HB3	1.85	0.58
1:B:298:GLU:HG3	1:B:299:PRO:HD2	1.86	0.58
1:A:147:GLU:O	1:A:151:GLN:NE2	2.36	0.58
1:A:147:GLU:HA	1:A:150:LEU:HD12	1.86	0.57
1:A:298:GLU:HG3	1:A:299:PRO:HD2	1.86	0.57
1:D:292:LEU:HD22	1:D:300:PRO:HB2	1.85	0.57
1:A:364:THR:HG21	1:A:452:ILE:HG23	1.86	0.57
1:D:317:HIS:NE2	1:D:401:GLU:OE1	2.36	0.57
1:C:250:ARG:HB2	1:C:289:LEU:HD12	1.87	0.57
1:B:298:GLU:HG2	10:B:715:HOH:O	2.04	0.57
1:D:140:ARG:HH12	1:D:145:ALA:HB3	1.70	0.56
1:A:183:ARG:HD3	1:A:447:TRP:CD2	2.40	0.56
1:A:277:HIS:CD2	1:A:300:PRO:HG2	2.41	0.55
1:D:365:ARG:HH12	3:D:502:H4B:C4	2.18	0.55
1:C:233:GLN:HB3	1:C:348:PHE:CE2	2.42	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:501:HEM:HBC2	2:C:501:HEM:HMC2	1.87	0.54
1:C:67:LYS:HD2	1:D:98:ARG:HH21	1.74	0.53
1:A:292:LEU:HD22	1:A:300:PRO:HB2	1.91	0.53
2:A:501:HEM:HBA1	2:A:501:HEM:CHA	2.36	0.53
5:D:505:BTB:O6	5:D:505:BTB:H32	2.09	0.52
1:C:244:TRP:CZ2	1:C:300:PRO:HG3	2.45	0.51
1:C:453:SER:HB3	1:C:456:LEU:HD12	1.92	0.51
1:A:234:ARG:NH1	1:A:347:GLU:OE1	2.43	0.51
1:B:67:LYS:N	1:B:67:LYS:HE3	2.26	0.51
1:C:364:THR:O	1:C:368:CYS:HB2	2.11	0.51
1:C:197:ASP:OD2	1:C:199:ARG:NH2	2.34	0.51
1:C:149:ARG:NH2	1:C:164:GLN:O	2.39	0.50
1:A:358:MET:HE3	1:A:360:THR:OG1	2.12	0.50
1:A:250:ARG:HA	1:A:250:ARG:HE	1.77	0.50
1:D:358:MET:HE3	1:D:360:THR:OG1	2.12	0.49
1:B:124:LEU:HB3	1:B:128:ARG:HH22	1.76	0.49
1:D:143:SER:O	1:D:147:GLU:HG2	2.13	0.49
1:B:365:ARG:HH12	3:B:502:H4B:C4	2.26	0.49
1:A:125:SER:HA	1:A:128:ARG:HD2	1.93	0.49
1:C:156:GLU:O	1:C:160:THR:HG22	2.12	0.49
1:C:382:CYS:HA	5:C:504:BTB:C1	2.43	0.49
1:C:358:MET:HE3	1:C:360:THR:OG1	2.12	0.49
1:C:382:CYS:HA	5:C:504:BTB:H11	1.93	0.49
1:A:68:PHE:CD1	1:A:83:THR:HG22	2.48	0.48
1:D:90:GLN:HB3	1:D:468:PHE:CD2	2.48	0.48
1:D:242:ARG:NH2	1:D:479:PRO:HD3	2.28	0.48
1:D:257:GLN:C	1:D:259:GLY:H	2.20	0.48
1:A:233:GLN:HB3	1:A:348:PHE:CE2	2.48	0.48
1:B:178:TRP:CE3	1:B:190:TRP:HA	2.49	0.48
1:C:202:ARG:HH11	1:C:202:ARG:HA	1.78	0.48
1:B:453:SER:HB3	1:B:456:LEU:HD12	1.95	0.48
1:D:262:ARG:HD3	1:D:283:ASN:O	2.14	0.48
1:D:317:HIS:CG	1:D:318:PRO:HD2	2.49	0.48
1:A:256:GLN:C	1:A:258:ASP:H	2.22	0.48
1:A:265:PRO:HA	1:A:268:VAL:HG23	1.94	0.48
3:A:502:H4B:H113	1:B:463:GLU:HG2	1.96	0.48
2:D:501:HEM:O2D	4:D:503:V80:F12	2.21	0.48
1:B:382:CYS:SG	5:C:504:BTB:H42	2.54	0.47
1:A:256:GLN:HB2	1:A:260:SER:O	2.14	0.47
1:A:258:ASP:OD1	1:A:258:ASP:N	2.47	0.47
1:C:162:THR:OG1	1:C:163:TYR:N	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:331:TYR:O	1:A:410:TYR:OH	2.25	0.47
1:C:99:CYS:HB3	1:D:466:ASN:HB3	1.97	0.47
1:D:178:TRP:CE3	1:D:190:TRP:HA	2.49	0.47
1:A:171:VAL:HG22	1:A:195:VAL:HB	1.96	0.47
1:A:365:ARG:HH12	3:A:502:H4B:C4	2.26	0.47
1:B:358:MET:HE3	1:B:360:THR:OG1	2.15	0.47
1:A:463:GLU:HG3	3:B:502:H4B:H113	1.96	0.47
1:C:279:TRP:HB2	1:C:302:LEU:HD11	1.97	0.47
1:C:229:THR:O	1:C:352:PRO:HD2	2.16	0.46
1:A:364:THR:O	1:A:368:CYS:HB2	2.15	0.46
1:B:261:VAL:HG11	1:B:265:PRO:HA	1.97	0.46
1:A:216:LYS:HG3	1:A:217:TYR:N	2.29	0.46
1:D:124:LEU:HD11	1:D:154:GLU:HG3	1.97	0.46
1:B:214:HIS:CD2	1:B:214:HIS:C	2.94	0.46
5:B:506:BTB:H11	5:B:506:BTB:H72	1.50	0.46
1:A:207:MET:HG2	1:A:231:PHE:CE1	2.51	0.46
1:C:275:ILE:HD11	1:C:281:PRO:HB3	1.98	0.46
1:C:70:ARG:HB2	1:C:81:TYR:CE2	2.50	0.46
1:D:214:HIS:C	1:D:214:HIS:CD2	2.94	0.46
1:A:207:MET:HE2	1:A:293:LEU:HD13	1.98	0.46
1:D:271:THR:O	1:D:275:ILE:HG12	2.15	0.45
1:D:364:THR:O	1:D:368:CYS:HB2	2.16	0.45
1:A:361:GLU:OE2	4:A:503:V80:N02	2.49	0.45
1:A:179:ARG:NH2	1:A:438:ARG:HG3	2.31	0.45
1:C:421:HIS:HB2	1:D:390:THR:HB	1.99	0.45
1:C:479:PRO:HD2	1:C:480:TRP:CZ3	2.52	0.45
1:D:298:GLU:OE1	5:D:505:BTB:H52	2.16	0.45
2:C:501:HEM:O2D	4:C:503:V80:F12	2.25	0.45
5:C:505:BTB:H41	5:C:505:BTB:H72	1.59	0.45
1:D:298:GLU:CD	5:D:505:BTB:H41	2.41	0.45
1:C:202:ARG:HA	1:C:202:ARG:NH1	2.32	0.44
1:C:254:TYR:OH	1:C:287:ASP:OD2	2.21	0.44
1:C:94:CYS:HB2	10:D:649:HOH:O	2.17	0.44
1:D:279:TRP:HB2	1:D:302:LEU:HD21	1.99	0.44
1:C:127:ALA:O	1:C:131:ILE:HG12	2.17	0.44
1:A:170:LEU:HD11	1:A:230:VAL:HG21	1.99	0.44
1:C:361:GLU:OE2	4:C:503:V80:N02	2.50	0.44
1:C:453:SER:HA	1:D:452:ILE:HG22	1.99	0.44
1:A:267:ASN:O	1:A:269:GLU:N	2.51	0.44
1:A:290:PRO:HB2	1:A:302:LEU:HD22	2.00	0.44
5:D:504:BTB:H32	5:D:504:BTB:H51	1.78	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:450:PRO:HG2	1:D:457:THR:HG21	2.00	0.43
5:A:505:BTB:H72	5:A:505:BTB:H41	1.43	0.43
1:B:298:GLU:OE1	5:B:505:BTB:O3	2.35	0.43
1:C:67:LYS:HD2	1:D:98:ARG:NH2	2.34	0.43
1:D:453:SER:HB3	1:D:456:LEU:HD12	1.98	0.43
1:C:451:PRO:HB2	1:D:455:SER:OG	2.17	0.43
1:A:124:LEU:O	1:A:128:ARG:HG3	2.18	0.43
1:A:342:GLU:HA	1:A:346:LEU:O	2.18	0.43
1:B:455:SER:HA	1:B:460:PHE:CG	2.54	0.43
1:C:135:TYR:HA	1:C:138:ILE:HG12	2.00	0.43
5:C:505:BTB:H32	5:C:505:BTB:O6	2.18	0.43
1:A:90:GLN:HB3	1:A:468:PHE:CD2	2.54	0.43
1:C:276:GLN:O	1:C:278:GLY:N	2.47	0.43
1:A:463:GLU:HG2	1:B:103:LEU:HA	2.01	0.43
1:D:455:SER:HA	1:D:460:PHE:CG	2.54	0.42
1:A:207:MET:HE2	1:A:207:MET:HB3	1.88	0.42
1:C:143:SER:HB2	10:C:607:HOH:O	2.18	0.42
1:C:94:CYS:HB3	1:D:94:CYS:HB3	2.01	0.42
1:C:233:GLN:OE1	10:C:601:HOH:O	2.21	0.42
1:A:274:CYS:SG	1:A:290:PRO:HG2	2.60	0.42
1:B:397:LYS:NZ	10:B:612:HOH:O	2.53	0.42
5:B:506:BTB:H32	5:B:506:BTB:H51	1.77	0.42
5:C:505:BTB:O4	5:C:505:BTB:O3	2.12	0.42
1:A:317:HIS:CG	1:A:318:PRO:HD2	2.55	0.42
1:C:188:ILE:HD12	1:C:188:ILE:HA	1.92	0.42
1:B:129:ASP:O	1:B:133:GLN:HG3	2.20	0.42
1:A:474:ARG:HE	1:A:474:ARG:HB3	1.46	0.42
5:A:505:BTB:H11	5:A:505:BTB:H51	1.62	0.42
1:B:357:TYR:CD2	1:B:362:ILE:HD11	2.54	0.42
1:C:97:ARG:HG2	1:D:88:ALA:HB3	2.01	0.41
1:A:201:CYS:O	1:A:202:ARG:HG3	2.20	0.41
1:B:218:ALA:HB1	10:B:623:HOH:O	2.21	0.41
1:C:279:TRP:CG	1:C:290:PRO:HG3	2.56	0.41
1:B:124:LEU:HD23	1:B:124:LEU:HA	1.89	0.41
1:C:274:CYS:O	1:C:279:TRP:N	2.50	0.41
5:C:506:BTB:H72	5:C:506:BTB:H32	1.77	0.41
1:D:231:PHE:HB3	1:D:232:PRO:CD	2.51	0.41
1:D:315:LEU:HD12	1:D:328:LEU:HB3	2.02	0.41
1:B:317:HIS:CG	1:B:318:PRO:HD2	2.56	0.41
1:A:455:SER:HA	1:A:460:PHE:CG	2.55	0.40
1:C:125:SER:O	1:C:128:ARG:HG3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:449:VAL:HA	1:D:450:PRO:HD3	1.95	0.40
5:B:505:BTB:H32	5:B:505:BTB:H51	1.99	0.40
1:C:276:GLN:C	1:C:278:GLY:H	2.28	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:GLU:OE2	5:D:505:BTB:O4[2_851]	2.15	0.05

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	398/440 (90%)	371 (93%)	23 (6%)	4 (1%)	12	11
1	B	401/440 (91%)	393 (98%)	8 (2%)	0	100	100
1	C	400/440 (91%)	382 (96%)	17 (4%)	1 (0%)	36	42
1	D	399/440 (91%)	389 (98%)	10 (2%)	0	100	100
All	All	1598/1760 (91%)	1535 (96%)	58 (4%)	5 (0%)	36	42

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	203	SER
1	A	268	VAL
1	A	144	GLN
1	A	290	PRO
1	C	277	HIS

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	342/373 (92%)	324 (95%)	18 (5%)	20	26
1	B	344/373 (92%)	325 (94%)	19 (6%)	19	24
1	C	343/373 (92%)	329 (96%)	14 (4%)	27	37
1	D	342/373 (92%)	329 (96%)	13 (4%)	29	40
All	All	1371/1492 (92%)	1307 (95%)	64 (5%)	24	31

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	87	GLN
1	A	97	ARG
1	A	121	GLU
1	A	122	GLN
1	A	125	SER
1	A	139	LYS
1	A	209	THR
1	A	216	LYS
1	A	224	LEU
1	A	238	ARG
1	A	255	ARG
1	A	256	GLN
1	A	298	GLU
1	A	309	LEU
1	A	321	GLU
1	A	342	GLU
1	A	396	ASP
1	A	417	ILE
1	B	67	LYS
1	B	87	GLN
1	B	89	GLN
1	B	97	ARG
1	B	121	GLU
1	B	122	GLN

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Mol	Chain	Res	Type
1	B	128	ARG
1	B	144	GLN
1	B	156	GLU
1	B	167	GLU
1	B	168[A]	SER
1	B	168[B]	SER
1	B	207	MET
1	B	255	ARG
1	B	257	GLN
1	B	258	ASP
1	B	326	LEU
1	B	396	ASP
1	B	417	ILE
1	C	67	LYS
1	C	71	VAL
1	C	89	GLN
1	C	124	LEU
1	C	125	SER
1	C	151	GLN
1	C	154	GLU
1	C	200	ASP
1	C	234	ARG
1	C	235	CYS
1	C	280	THR
1	C	308	GLU
1	C	412	LEU
1	C	417	ILE
1	D	71	VAL
1	D	97	ARG
1	D	98	ARG
1	D	121	GLU
1	D	124	LEU
1	D	160	THR
1	D	192	LYS
1	D	230	VAL
1	D	255	ARG
1	D	268	VAL
1	D	309	LEU
1	D	326	LEU
1	D	396	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	133	GLN
1	A	220	ASN
1	A	257	GLN
1	A	435	GLN
1	B	148	GLN
1	D	126	GLN
1	D	466	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 34 ligands modelled in this entry, 10 are monoatomic - leaving 24 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
5	BTB	A	505	-	13,13,13	0.57	0	7,16,16	0.94	1 (14%)
3	H4B	A	502	-	17,18,18	0.91	0	14,26,26	2.20	5 (35%)
2	HEM	C	501	1	50,50,50	1.74	9 (18%)	67,82,82	1.30	6 (8%)
2	HEM	B	501	1	50,50,50	1.72	10 (20%)	67,82,82	1.35	8 (11%)
4	V80	A	503	-	24,24,24	0.42	0	33,33,33	1.77	6 (18%)
5	BTB	C	505	-	13,13,13	0.72	0	7,16,16	1.62	2 (28%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	BTB	C	506	-	13,13,13	0.42	0	7,16,16	0.27	0
2	HEM	A	501	1	50,50,50	1.78	10 (20%)	67,82,82	1.25	8 (11%)
5	BTB	A	504	8	13,13,13	0.50	0	7,16,16	0.70	0
3	H4B	D	502	-	17,18,18	0.83	0	14,26,26	1.80	4 (28%)
5	BTB	A	506	-	13,13,13	0.39	0	7,16,16	0.41	0
4	V80	D	503	-	24,24,24	0.59	0	33,33,33	1.57	7 (21%)
4	V80	C	503	-	24,24,24	0.43	0	33,33,33	1.92	9 (27%)
5	BTB	B	506	-	13,13,13	0.45	0	7,16,16	0.46	0
5	BTB	B	505	-	13,13,13	0.42	0	7,16,16	0.87	0
5	BTB	D	505	-	13,13,13	0.57	0	7,16,16	0.89	0
4	V80	B	503	-	24,24,24	0.69	0	33,33,33	1.71	8 (24%)
2	HEM	D	501	1	50,50,50	1.67	9 (18%)	67,82,82	1.28	6 (8%)
3	H4B	C	502	-	17,18,18	0.87	0	14,26,26	1.94	5 (35%)
9	GOL	C	508	-	5,5,5	0.42	0	5,5,5	0.25	0
5	BTB	C	504	8	13,13,13	0.61	0	7,16,16	1.33	1 (14%)
3	H4B	B	502	-	17,18,18	0.95	0	14,26,26	2.22	5 (35%)
5	BTB	B	504	8	13,13,13	0.48	0	7,16,16	0.34	0
5	BTB	D	504	8	13,13,13	0.42	0	7,16,16	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
5	BTB	A	505	-	-	11/21/21/21	-
3	H4B	A	502	-	-	3/8/17/17	0/2/2/2
2	HEM	C	501	1	-	7/14/54/54	-
2	HEM	B	501	1	-	6/14/54/54	-
4	V80	A	503	-	-	4/10/10/10	0/2/2/2
5	BTB	C	505	-	-	10/21/21/21	-
5	BTB	C	506	-	-	8/21/21/21	-
2	HEM	A	501	1	-	5/14/54/54	-
5	BTB	A	504	8	-	4/21/21/21	-
3	H4B	D	502	-	-	3/8/17/17	0/2/2/2
5	BTB	A	506	-	-	9/21/21/21	-
4	V80	D	503	-	-	3/10/10/10	0/2/2/2
4	V80	C	503	-	-	2/10/10/10	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	BTB	B	506	-	-	10/21/21/21	-
5	BTB	B	505	-	-	8/21/21/21	-
5	BTB	D	505	-	-	10/21/21/21	-
4	V80	B	503	-	-	4/10/10/10	0/2/2/2
2	HEM	D	501	1	-	6/14/54/54	-
3	H4B	C	502	-	-	3/8/17/17	0/2/2/2
9	GOL	C	508	-	-	4/4/4/4	-
5	BTB	C	504	8	-	6/21/21/21	-
3	H4B	B	502	-	-	3/8/17/17	0/2/2/2
5	BTB	B	504	8	-	3/21/21/21	-
5	BTB	D	504	8	-	4/21/21/21	-

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	HEM	FE-NA	6.66	2.17	1.95
2	A	501	HEM	FE-NB	6.43	2.14	1.94
2	C	501	HEM	FE-NB	5.48	2.11	1.94
2	C	501	HEM	FE-NA	5.32	2.12	1.95
2	D	501	HEM	FE-NB	4.94	2.10	1.94
2	A	501	HEM	FE-NA	4.94	2.11	1.95
2	D	501	HEM	FE-NC	4.34	2.09	1.95
2	C	501	HEM	FE-NC	4.05	2.08	1.95
2	B	501	HEM	FE-NB	4.01	2.07	1.94
2	D	501	HEM	FE-ND	3.96	2.07	1.94
2	A	501	HEM	FE-ND	3.72	2.06	1.94
2	D	501	HEM	FE-NA	3.69	2.07	1.95
2	C	501	HEM	FE-ND	3.47	2.05	1.94
2	B	501	HEM	FE-ND	3.43	2.05	1.94
2	A	501	HEM	FE-NC	3.33	2.06	1.95
2	B	501	HEM	CAC-C3C	3.28	1.56	1.47
2	A	501	HEM	CAB-C3B	3.06	1.55	1.47
2	C	501	HEM	CAB-C3B	3.03	1.55	1.47
2	D	501	HEM	CAB-C3B	3.01	1.55	1.47
2	D	501	HEM	CAC-C3C	2.97	1.55	1.47
2	A	501	HEM	CAC-C3C	2.90	1.55	1.47
2	C	501	HEM	CAC-C3C	2.88	1.55	1.47
2	B	501	HEM	CAB-C3B	2.79	1.54	1.47
2	D	501	HEM	CMA-C3A	2.35	1.55	1.50
2	D	501	HEM	CMB-C2B	2.27	1.55	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	HEM	CMB-C2B	2.26	1.55	1.50
2	C	501	HEM	CMC-C2C	2.25	1.55	1.50
2	A	501	HEM	CMC-C2C	2.24	1.55	1.50
2	C	501	HEM	CMB-C2B	2.23	1.55	1.50
2	D	501	HEM	CMC-C2C	2.21	1.55	1.50
2	B	501	HEM	CMD-C2D	2.10	1.55	1.50
2	A	501	HEM	CMD-C2D	2.06	1.55	1.50
2	A	501	HEM	C2A-C3A	-2.06	1.33	1.38
2	B	501	HEM	CMA-C3A	2.05	1.55	1.50
2	A	501	HEM	CMB-C2B	2.02	1.54	1.50
2	C	501	HEM	CMA-C3A	2.02	1.54	1.50
2	B	501	HEM	CMC-C2C	2.02	1.54	1.50
2	B	501	HEM	FE-NC	2.01	2.01	1.95

All (81) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	503	V80	C02-N01-C06	5.33	121.58	117.06
4	C	503	V80	C02-N01-C06	4.92	121.24	117.06
4	B	503	V80	C02-N01-C06	4.58	120.95	117.06
3	A	502	H4B	C11-C10-C9	-4.48	106.63	112.11
3	B	502	H4B	C2-N1-C8A	4.46	121.23	113.36
4	C	503	V80	C02-N03-C04	4.37	120.73	117.11
4	D	503	V80	C02-N01-C06	4.18	120.61	117.06
3	C	502	H4B	C2-N1-C8A	4.17	120.72	113.36
3	D	502	H4B	C2-N1-C8A	4.11	120.61	113.36
3	A	502	H4B	C2-N1-C8A	4.05	120.50	113.36
4	A	503	V80	C05-C06-N01	-3.88	118.30	122.73
4	A	503	V80	C02-N03-C04	3.74	120.21	117.11
3	B	502	H4B	O4-C4-C4A	-3.72	118.28	127.26
4	D	503	V80	C02-N03-C04	3.60	120.09	117.11
4	C	503	V80	N03-C02-N01	-3.39	120.29	125.48
2	B	501	HEM	C3D-C4D-ND	-3.36	106.48	110.17
3	B	502	H4B	C11-C10-C9	-3.33	108.04	112.11
2	B	501	HEM	CBD-CAD-C3D	-3.24	103.58	112.53
3	A	502	H4B	O4-C4-C4A	-3.21	119.52	127.26
4	B	503	V80	N03-C02-N01	-3.18	120.61	125.48
4	A	503	V80	N03-C02-N01	-3.14	120.67	125.48
4	C	503	V80	C05-C06-N01	-3.06	119.24	122.73
4	B	503	V80	C02-N03-C04	3.05	119.64	117.11
4	B	503	V80	N02-C02-N03	3.03	121.77	117.22
2	C	501	HEM	C3B-C4B-NB	-2.99	107.32	109.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	502	H4B	O4-C4-C4A	-2.94	120.16	127.26
4	C	503	V80	F11-C11-C16	2.88	121.90	118.00
5	C	504	BTB	O4-C4-C2	2.88	118.17	111.40
2	B	501	HEM	C4D-ND-C1D	2.88	108.61	105.21
4	D	503	V80	N03-C02-N01	-2.85	121.11	125.48
2	D	501	HEM	C3B-C4B-NB	-2.85	107.42	109.47
4	A	503	V80	C08-C06-N01	2.77	120.26	116.06
2	D	501	HEM	C1B-NB-C4B	2.73	108.44	105.21
2	A	501	HEM	CAA-C2A-C1A	2.71	130.22	124.94
2	A	501	HEM	C3B-C4B-NB	-2.70	107.53	109.47
4	C	503	V80	C18-C17-C14	-2.69	104.56	112.15
4	D	503	V80	C05-C06-N01	-2.68	119.68	122.73
3	B	502	H4B	C4A-C4-N3	2.67	119.45	112.13
2	B	501	HEM	CBA-CAA-C2A	-2.66	105.17	112.53
2	C	501	HEM	C2A-C1A-NA	-2.65	107.22	110.15
4	A	503	V80	F11-C11-C16	2.64	121.57	118.00
2	D	501	HEM	CBD-CAD-C3D	-2.63	105.27	112.53
2	D	501	HEM	C3D-C4D-ND	-2.62	107.29	110.17
5	C	505	BTB	O3-C3-C2	-2.62	105.25	111.40
3	D	502	H4B	O4-C4-C4A	-2.60	121.00	127.26
4	D	503	V80	N02-C02-N03	2.56	121.07	117.22
2	C	501	HEM	C1B-NB-C4B	2.54	108.21	105.21
3	B	502	H4B	C2-N3-C4	-2.54	120.51	125.11
2	A	501	HEM	C3B-C2B-C1B	2.53	108.31	106.41
5	C	505	BTB	O4-C4-C2	-2.52	105.47	111.40
2	A	501	HEM	C1B-NB-C4B	2.50	108.17	105.21
2	D	501	HEM	C4D-ND-C1D	2.50	108.17	105.21
2	B	501	HEM	C3B-C2B-C1B	2.49	108.28	106.41
2	A	501	HEM	C4D-ND-C1D	2.47	108.14	105.21
3	C	502	H4B	C4A-C4-N3	2.47	118.92	112.13
3	A	502	H4B	C4A-C4-N3	2.44	118.83	112.13
4	D	503	V80	F11-C11-C16	2.41	121.26	118.00
4	B	503	V80	C09-C08-C06	-2.41	107.66	113.01
2	C	501	HEM	C3D-C4D-ND	-2.37	107.57	110.17
4	B	503	V80	C09-C16-C11	2.36	122.63	120.73
3	D	502	H4B	C4A-C4-N3	2.34	118.56	112.13
3	A	502	H4B	C2-N3-C4	-2.34	120.88	125.11
2	C	501	HEM	C4D-ND-C1D	2.33	107.97	105.21
4	B	503	V80	C05-C06-N01	-2.31	120.10	122.73
2	A	501	HEM	C2A-C1A-NA	-2.31	107.59	110.15
3	C	502	H4B	C11-C10-C9	-2.29	109.32	112.11
2	B	501	HEM	CAD-CBD-CGD	-2.27	107.65	113.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	HEM	C3B-C2B-C1B	2.25	108.10	106.41
3	C	502	H4B	C2-N3-C4	-2.23	121.06	125.11
2	A	501	HEM	CAA-C2A-C3A	-2.21	122.11	127.07
4	C	503	V80	C08-C06-N01	2.18	119.37	116.06
4	B	503	V80	C07-C04-N03	2.16	119.83	116.56
2	B	501	HEM	CAA-CBA-CGA	2.16	119.40	113.67
2	A	501	HEM	C3D-C4D-ND	-2.16	107.80	110.17
4	C	503	V80	C09-C16-C11	2.14	122.45	120.73
2	D	501	HEM	C2A-C1A-NA	-2.13	107.78	110.15
2	B	501	HEM	CHB-C1B-NB	2.10	126.97	124.37
3	D	502	H4B	N3-C2-N1	-2.05	119.57	123.32
4	C	503	V80	C17-C18-N19	-2.04	107.18	113.87
4	D	503	V80	C07-C04-N03	2.03	119.62	116.56
5	A	505	BTB	O3-C3-C2	-2.01	106.67	111.40

There are no chirality outliers.

All (136) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	HEM	C1A-C2A-CAA-CBA
2	C	501	HEM	C1A-C2A-CAA-CBA
3	C	502	H4B	C7-C6-C9-O9
3	C	502	H4B	C7-C6-C9-C10
3	D	502	H4B	C7-C6-C9-O9
3	D	502	H4B	C7-C6-C9-C10
5	A	504	BTB	C1-C2-C4-O4
5	A	504	BTB	C3-C2-C4-O4
5	A	504	BTB	N-C2-C4-O4
5	A	505	BTB	O1-C1-C2-C4
5	A	505	BTB	C1-C2-N-C5
5	A	505	BTB	C1-C2-N-C7
5	A	505	BTB	C3-C2-N-C5
5	A	505	BTB	C3-C2-N-C7
5	A	505	BTB	C4-C2-N-C5
5	A	505	BTB	C4-C2-N-C7
5	A	506	BTB	O1-C1-C2-C3
5	A	506	BTB	C1-C2-C3-O3
5	A	506	BTB	C4-C2-C3-O3
5	A	506	BTB	N-C2-C3-O3
5	A	506	BTB	C1-C2-C4-O4
5	A	506	BTB	C3-C2-C4-O4
5	A	506	BTB	N-C2-C4-O4

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Mol	Chain	Res	Type	Atoms
5	B	504	BTB	O1-C1-C2-C3
5	B	504	BTB	O1-C1-C2-N
5	B	505	BTB	O1-C1-C2-N
5	B	505	BTB	C1-C2-C4-O4
5	B	505	BTB	C3-C2-C4-O4
5	B	505	BTB	N-C2-C4-O4
5	B	506	BTB	C1-C2-C4-O4
5	B	506	BTB	C3-C2-C4-O4
5	B	506	BTB	N-C2-C4-O4
5	B	506	BTB	C1-C2-N-C5
5	B	506	BTB	C1-C2-N-C7
5	B	506	BTB	C3-C2-N-C5
5	B	506	BTB	C3-C2-N-C7
5	B	506	BTB	C4-C2-N-C5
5	B	506	BTB	C4-C2-N-C7
5	C	504	BTB	O1-C1-C2-C3
5	C	504	BTB	O1-C1-C2-C4
5	C	504	BTB	O1-C1-C2-N
5	C	504	BTB	N-C2-C3-O3
5	C	505	BTB	O1-C1-C2-C3
5	C	505	BTB	O1-C1-C2-C4
5	C	505	BTB	O1-C1-C2-N
5	C	506	BTB	O1-C1-C2-C3
5	C	506	BTB	O1-C1-C2-C4
5	C	506	BTB	O1-C1-C2-N
5	C	506	BTB	N-C2-C3-O3
5	C	506	BTB	N-C2-C4-O4
5	D	504	BTB	O1-C1-C2-C3
5	D	504	BTB	O1-C1-C2-C4
5	D	504	BTB	O1-C1-C2-N
5	D	505	BTB	O1-C1-C2-C3
5	D	505	BTB	O1-C1-C2-C4
5	D	505	BTB	O1-C1-C2-N
5	D	505	BTB	N-C2-C3-O3
5	D	505	BTB	C6-C5-N-C7
9	C	508	GOL	O1-C1-C2-C3
5	B	505	BTB	N-C5-C6-O6
2	A	501	HEM	C3A-C2A-CAA-CBA
4	B	503	V80	C08-C09-C16-C15
4	D	503	V80	C17-C18-N19-C21
2	B	501	HEM	C3D-CAD-CBD-CGD
2	D	501	HEM	C2A-CAA-CBA-CGA

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Mol	Chain	Res	Type	Atoms
4	C	503	V80	C08-C09-C16-C15
2	C	501	HEM	C2A-CAA-CBA-CGA
9	C	508	GOL	C1-C2-C3-O3
5	B	505	BTB	N-C7-C8-O8
5	B	506	BTB	N-C7-C8-O8
4	B	503	V80	C17-C18-N19-C21
5	B	505	BTB	C4-C2-C3-O3
5	C	504	BTB	C1-C2-C3-O3
5	C	506	BTB	C3-C2-C4-O4
2	C	501	HEM	C3A-C2A-CAA-CBA
2	A	501	HEM	C2A-CAA-CBA-CGA
2	C	501	HEM	C3D-CAD-CBD-CGD
5	A	505	BTB	N-C7-C8-O8
5	C	504	BTB	N-C7-C8-O8
4	C	503	V80	C14-C17-C18-N19
4	D	503	V80	C14-C17-C18-N19
4	A	503	V80	C08-C09-C16-C15
3	B	502	H4B	C7-C6-C9-O9
2	A	501	HEM	C3D-CAD-CBD-CGD
9	C	508	GOL	O1-C1-C2-O2
9	C	508	GOL	O2-C2-C3-O3
5	A	505	BTB	N-C5-C6-O6
5	B	505	BTB	O1-C1-C2-C4
4	A	503	V80	C14-C17-C18-N19
3	A	502	H4B	C7-C6-C9-O9
4	A	503	V80	C15-C14-C17-C18
2	D	501	HEM	C3D-CAD-CBD-CGD
4	A	503	V80	C13-C14-C17-C18
5	A	506	BTB	N-C7-C8-O8
2	A	501	HEM	C4B-C3B-CAB-CBB
2	B	501	HEM	C4B-C3B-CAB-CBB
2	C	501	HEM	C4B-C3B-CAB-CBB
2	D	501	HEM	C4B-C3B-CAB-CBB
5	A	505	BTB	O1-C1-C2-C3
5	D	504	BTB	C3-C2-C4-O4
5	D	505	BTB	C1-C2-C3-O3
3	D	502	H4B	N5-C6-C9-O9
5	A	505	BTB	O1-C1-C2-N
5	A	506	BTB	O1-C1-C2-N
5	C	505	BTB	C1-C2-N-C5
5	C	505	BTB	C1-C2-N-C7
5	C	505	BTB	C3-C2-N-C5

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Mol	Chain	Res	Type	Atoms
5	C	505	BTB	C4-C2-N-C7
5	D	505	BTB	C3-C2-N-C5
4	B	503	V80	C08-C09-C16-C11
5	C	506	BTB	N-C7-C8-O8
3	A	502	H4B	C7-C6-C9-C10
3	B	502	H4B	C7-C6-C9-C10
5	B	504	BTB	C3-C2-C4-O4
5	C	506	BTB	C1-C2-C3-O3
4	B	503	V80	C14-C17-C18-N19
2	B	501	HEM	CAD-CBD-CGD-O2D
2	D	501	HEM	C1A-C2A-CAA-CBA
2	D	501	HEM	CAA-CBA-CGA-O1A
4	D	503	V80	C08-C09-C16-C15
2	D	501	HEM	CAA-CBA-CGA-O2A
5	C	505	BTB	C6-C5-N-C7
2	B	501	HEM	CAD-CBD-CGD-O1D
2	B	501	HEM	CAA-CBA-CGA-O1A
2	B	501	HEM	CAA-CBA-CGA-O2A
3	A	502	H4B	N5-C6-C9-O9
3	B	502	H4B	N5-C6-C9-O9
3	C	502	H4B	N5-C6-C9-O9
2	C	501	HEM	CAA-CBA-CGA-O2A
5	A	504	BTB	C3-C2-N-C5
5	C	505	BTB	C3-C2-N-C7
5	C	505	BTB	C4-C2-N-C5
5	D	505	BTB	N-C2-C4-O4
5	D	505	BTB	C1-C2-N-C5
5	D	505	BTB	C4-C2-N-C5
2	C	501	HEM	CAA-CBA-CGA-O1A

There are no ring outliers.

19 monomers are involved in 44 short contacts:

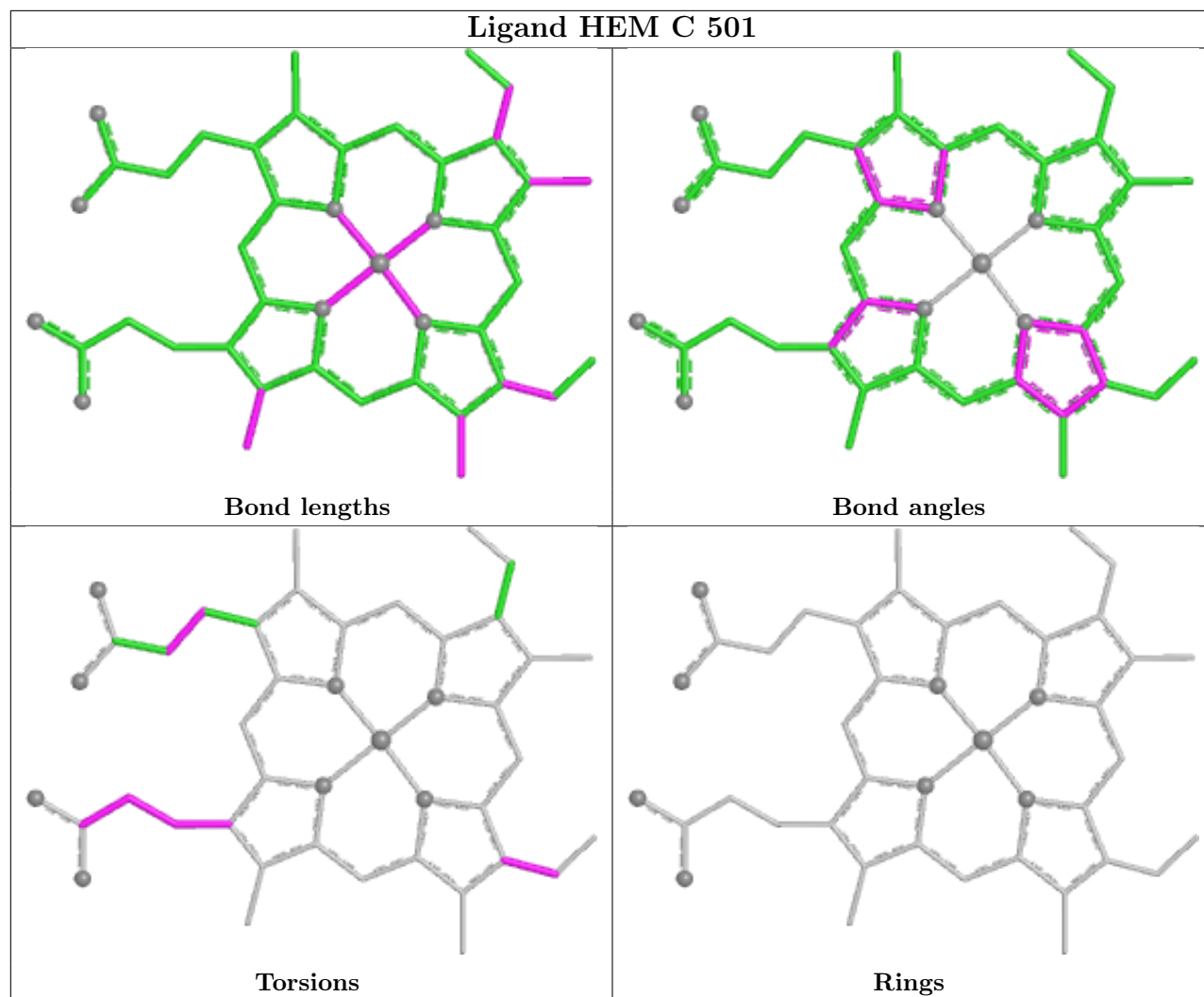
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	505	BTB	2	0
3	A	502	H4B	2	0
2	C	501	HEM	3	0
2	B	501	HEM	3	0
4	A	503	V80	1	0
5	C	505	BTB	3	0
5	C	506	BTB	1	0
2	A	501	HEM	6	0

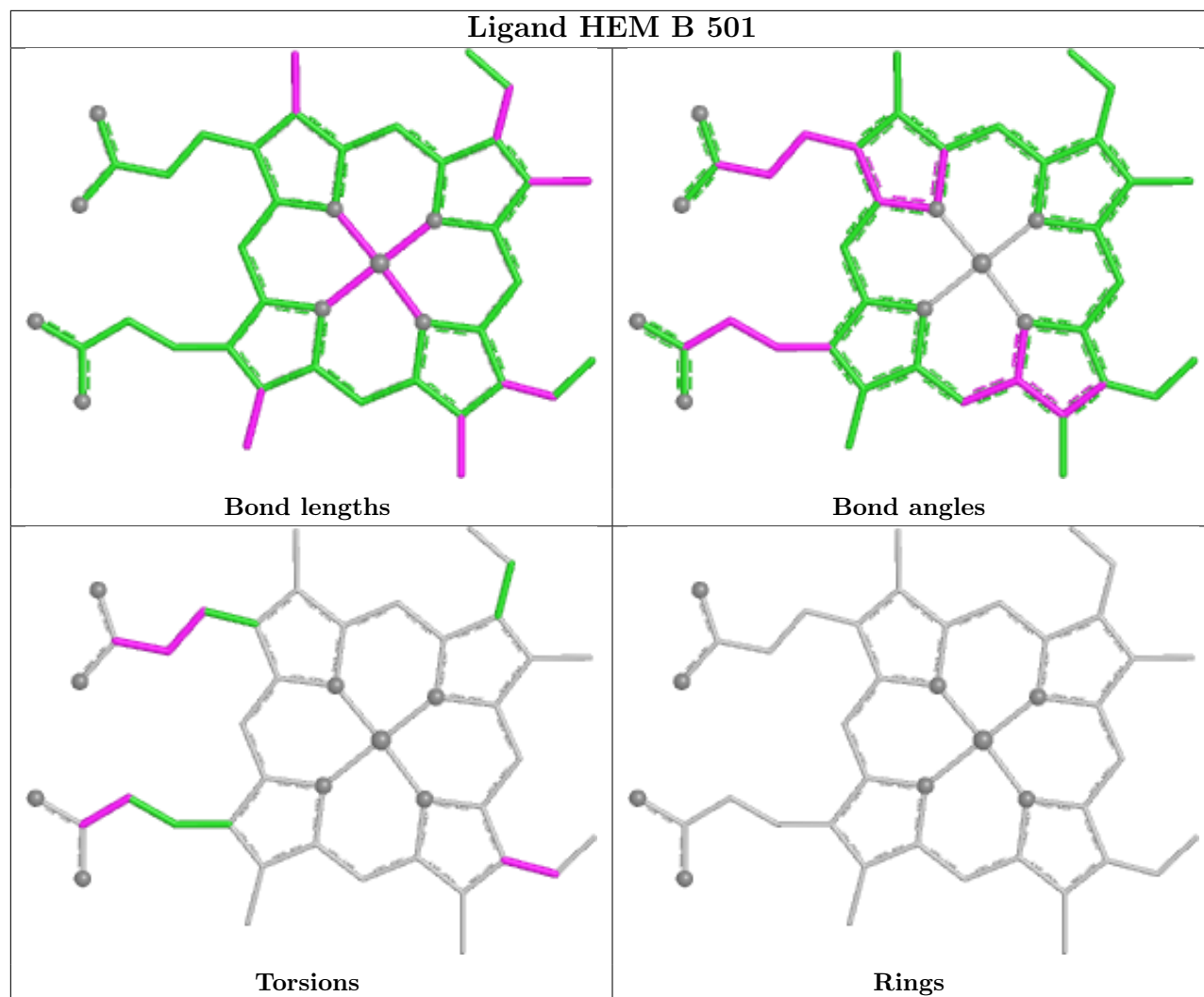
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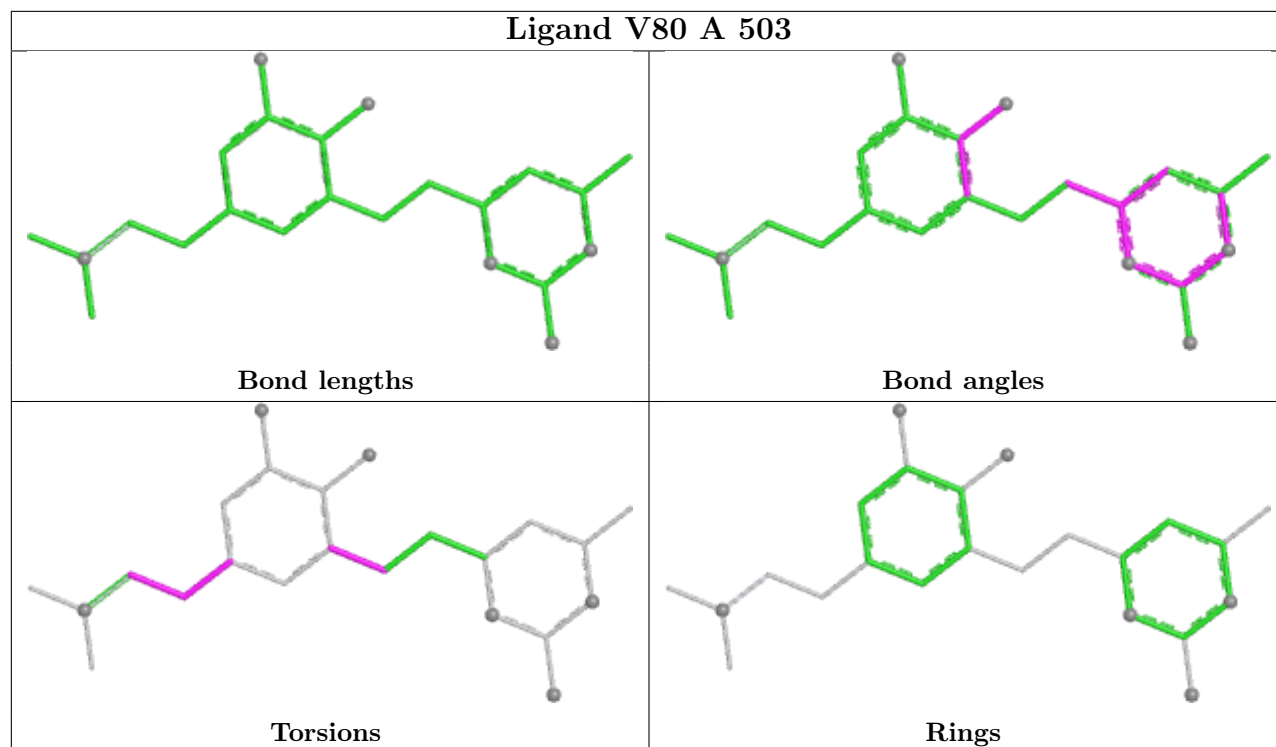
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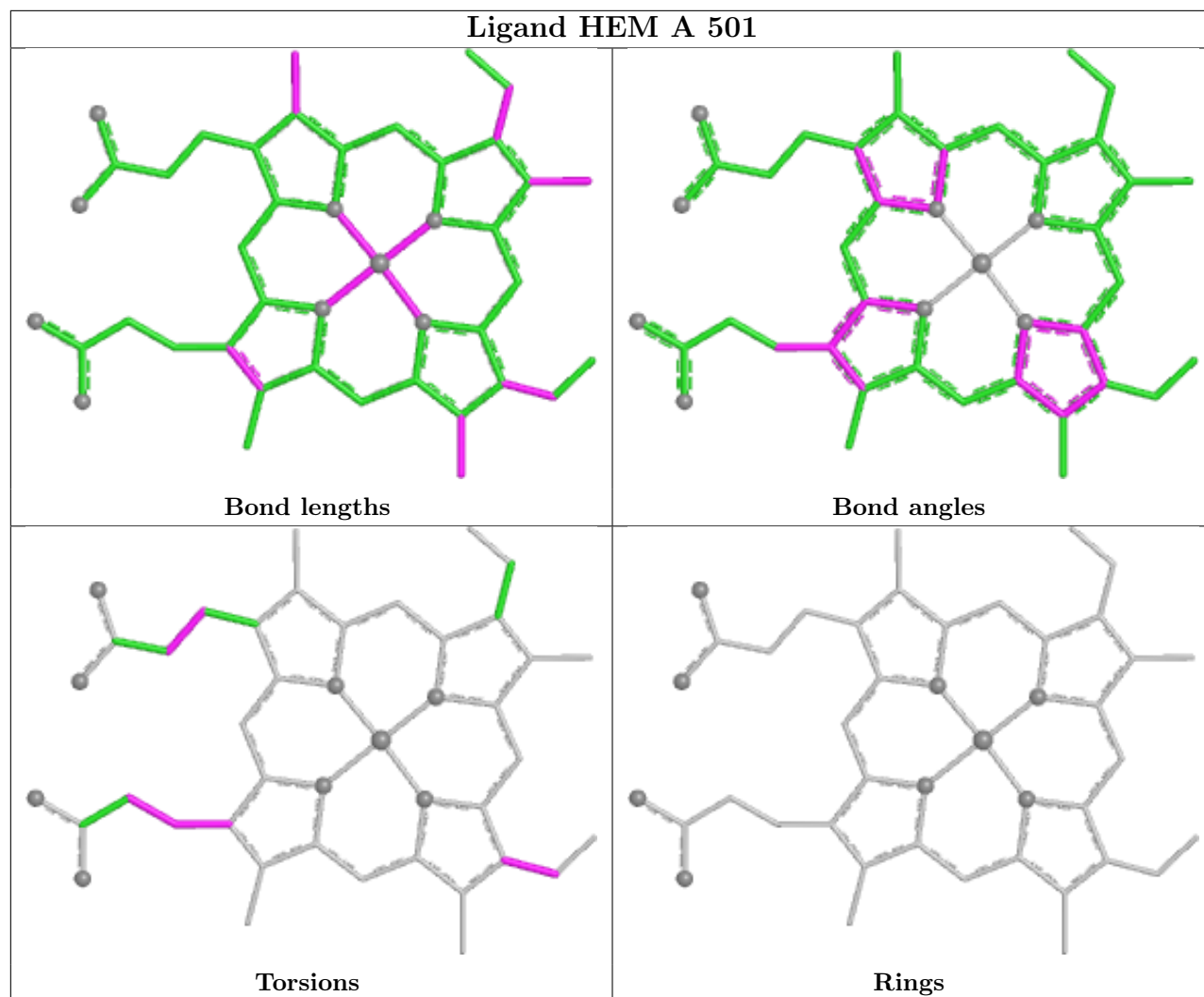
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	502	H4B	1	0
4	D	503	V80	1	0
4	C	503	V80	2	0
5	B	506	BTB	2	0
5	B	505	BTB	3	0
5	D	505	BTB	4	1
2	D	501	HEM	3	0
3	C	502	H4B	1	0
5	C	504	BTB	3	0
3	B	502	H4B	2	0
5	D	504	BTB	2	0

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.

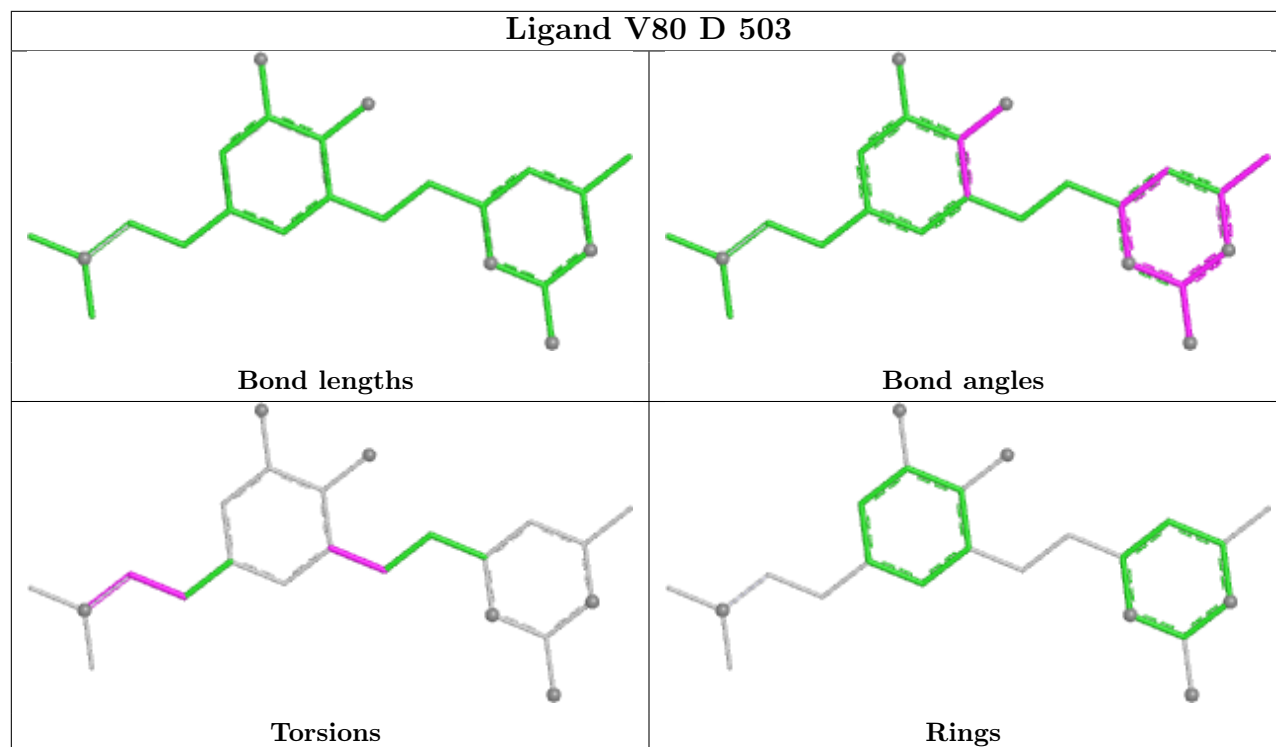




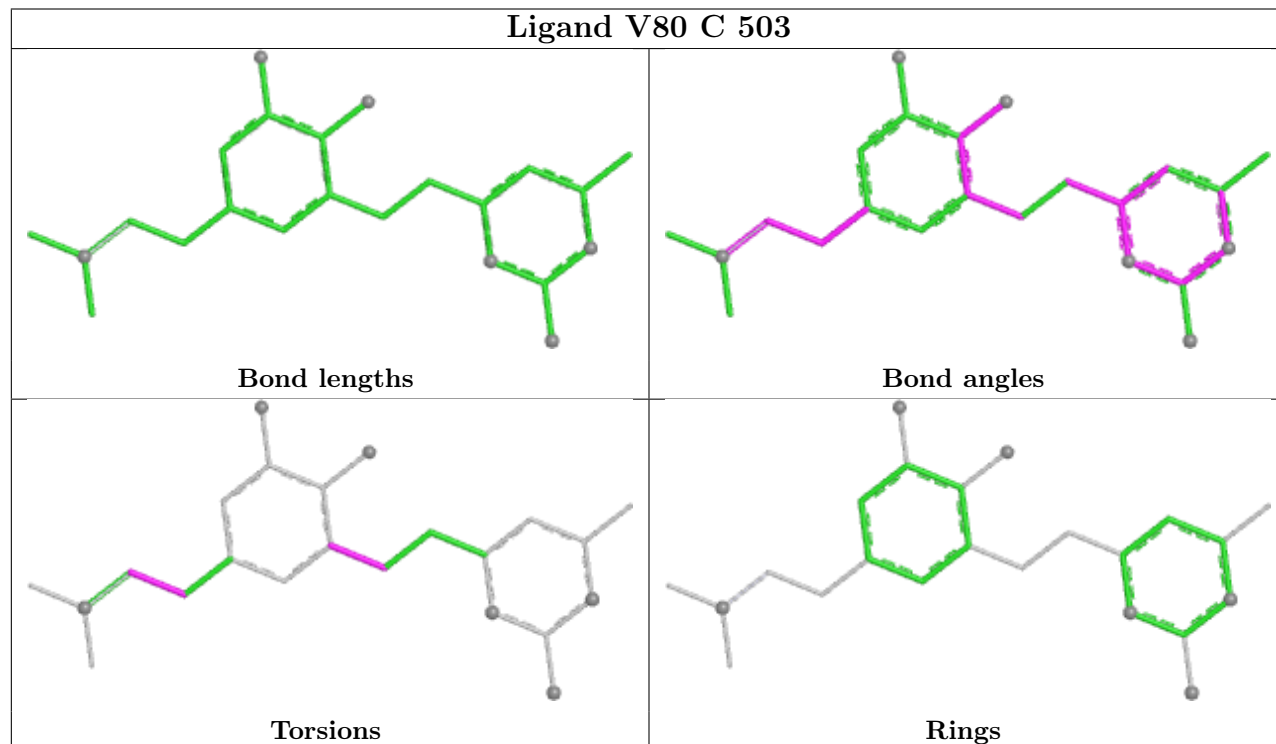


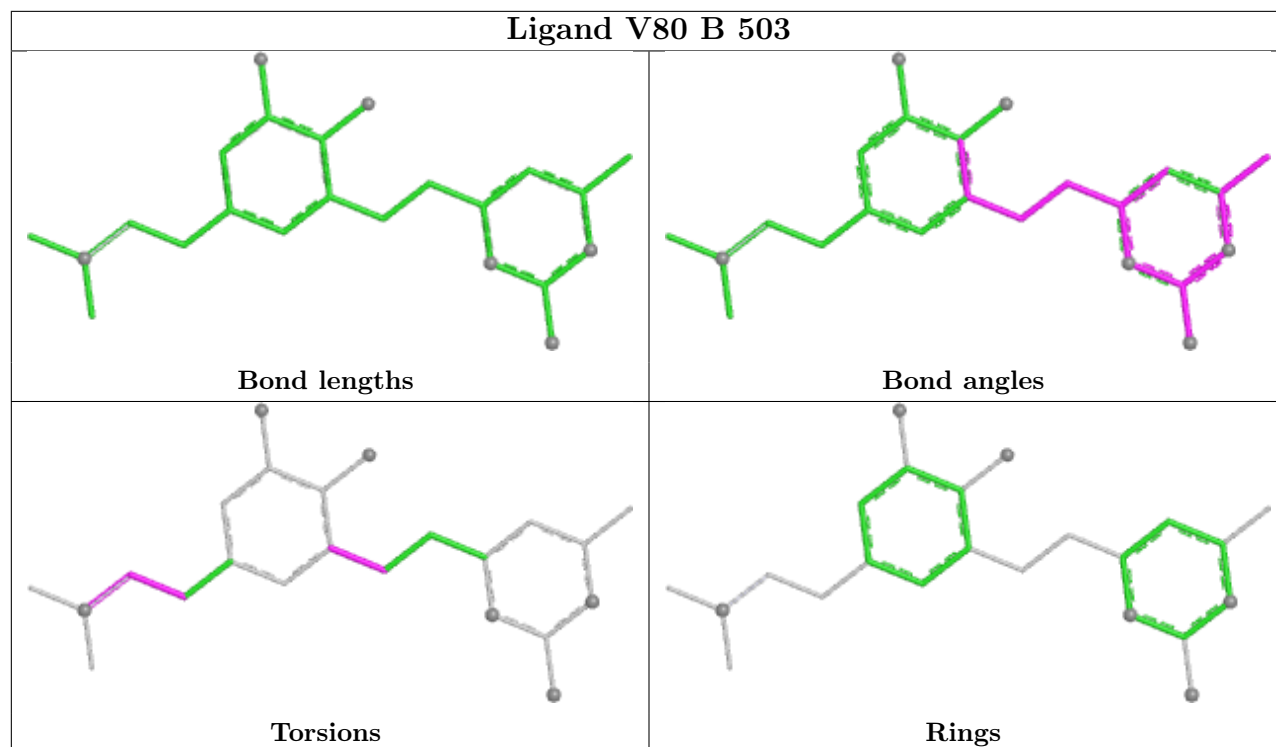


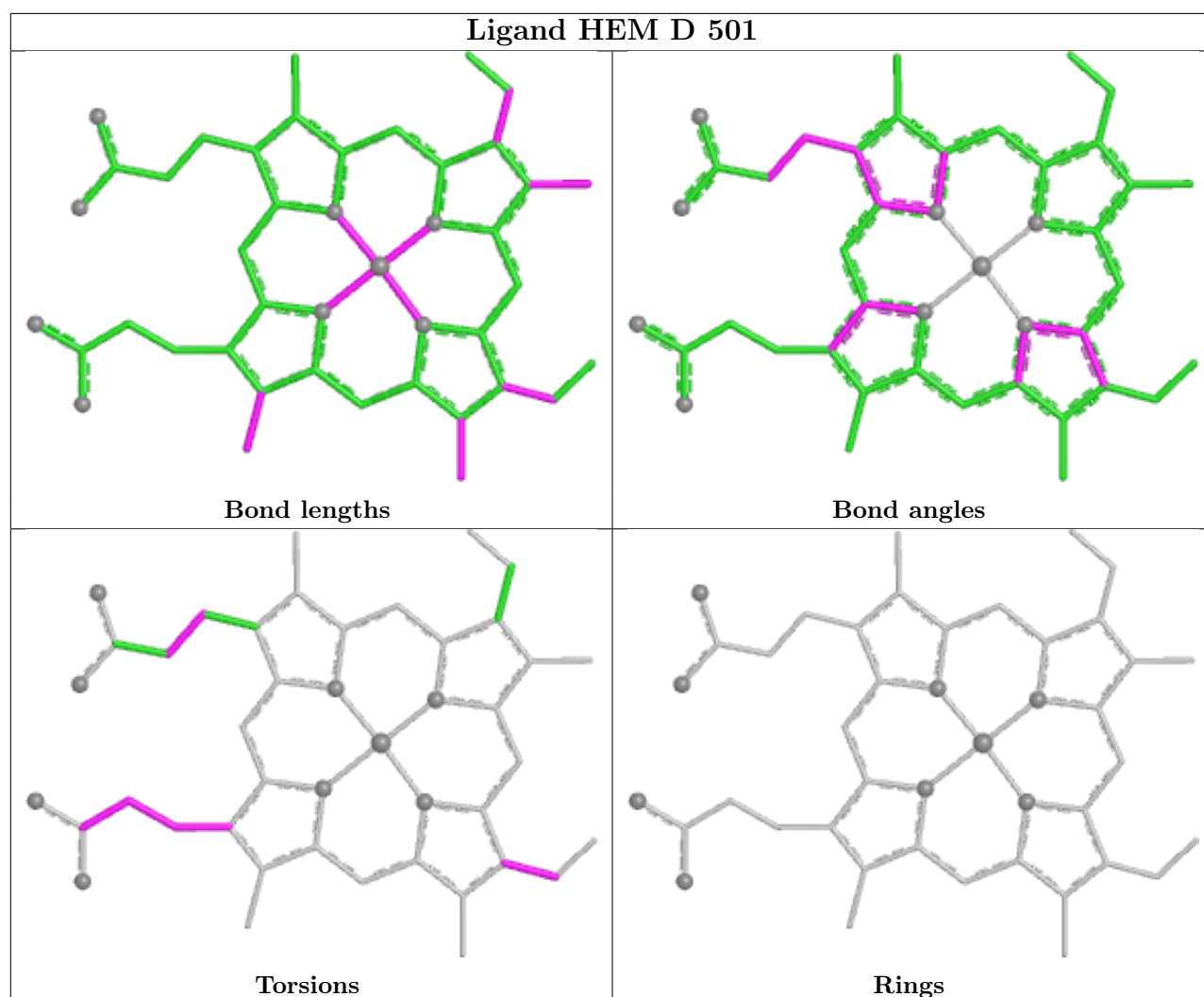
Ligand V80 D 503



Ligand V80 C 503







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	400/440 (90%)	0.52	24 (6%)	27 24	20, 52, 108, 136	2 (0%)
1	B	402/440 (91%)	-0.21	2 (0%)	87 85	13, 32, 66, 104	3 (0%)
1	C	402/440 (91%)	0.20	11 (2%)	56 53	20, 43, 89, 114	2 (0%)
1	D	401/440 (91%)	-0.24	6 (1%)	72 69	16, 32, 58, 121	2 (0%)
All	All	1605/1760 (91%)	0.07	43 (2%)	56 53	13, 38, 88, 136	9 (0%)

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	106	PRO	4.2
1	A	204	ALA	4.0
1	A	120	PRO	3.9
1	C	119	ALA	3.6
1	C	106	PRO	3.3
1	A	153	VAL	3.1
1	B	106	PRO	3.0
1	C	120	PRO	2.8
1	D	89	GLN	2.7
1	A	295	ALA	2.7
1	A	237	GLY	2.7
1	C	299	PRO	2.7
1	D	106	PRO	2.7
1	A	292	LEU	2.7
1	C	161	GLY	2.6
1	A	343	ILE	2.6
1	D	119	ALA	2.5
1	A	300	PRO	2.5
1	A	143	SER	2.5
1	C	140	ARG	2.5
1	D	257	GLN	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	275	ILE	2.4
1	A	299	PRO	2.4
1	B	257	GLN	2.3
1	C	141	SER	2.3
1	D	141[A]	SER	2.3
1	A	124	LEU	2.2
1	A	284	GLY	2.2
1	A	238	ARG	2.2
1	A	293	LEU	2.2
1	D	258	ASP	2.2
1	A	161	GLY	2.2
1	A	140	ARG	2.2
1	A	268	VAL	2.2
1	A	145	ALA	2.2
1	A	480	TRP	2.1
1	C	86	ALA	2.1
1	A	147	GLU	2.1
1	C	280	THR	2.0
1	A	207	MET	2.0
1	A	235	CYS	2.0
1	C	257	GLN	2.0
1	C	307	PRO	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
5	BTB	C	506	14/14	0.58	0.18	87,106,111,113	0

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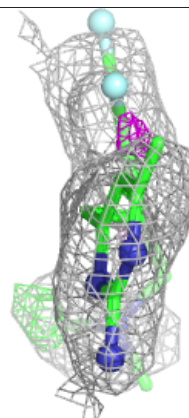
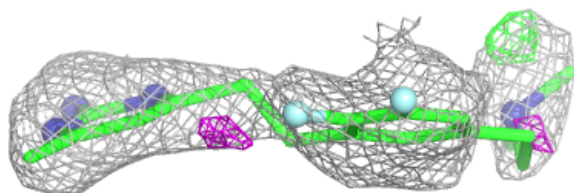
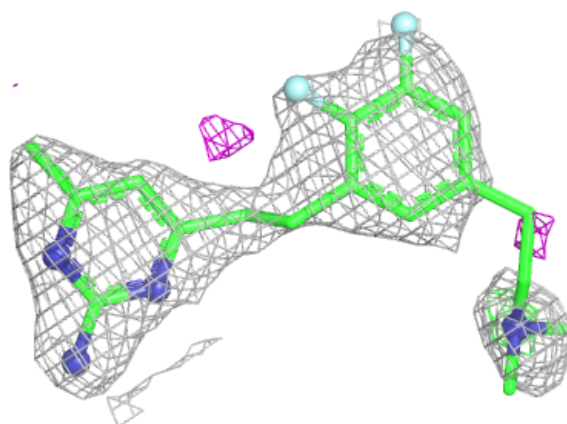
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	BTB	B	506	14/14	0.64	0.16	90,95,100,101	0
5	BTB	D	505	14/14	0.73	0.15	64,74,80,81	0
3	H4B	A	502	17/17	0.74	0.20	58,86,102,102	0
5	BTB	C	505	14/14	0.74	0.18	53,60,72,72	0
3	H4B	B	502	17/17	0.74	0.20	54,75,93,100	0
5	BTB	B	505	14/14	0.74	0.14	53,58,69,76	0
5	BTB	A	506	14/14	0.76	0.13	72,84,90,91	0
9	GOL	C	508	6/6	0.77	0.14	73,77,78,78	0
3	H4B	D	502	17/17	0.81	0.18	47,74,90,91	0
5	BTB	A	505	14/14	0.81	0.14	62,76,80,82	0
3	H4B	C	502	17/17	0.81	0.17	56,71,90,96	0
4	V80	A	503	23/23	0.86	0.17	38,82,95,102	0
4	V80	C	503	23/23	0.86	0.17	48,83,95,96	0
5	BTB	A	504	14/14	0.87	0.14	52,78,88,90	0
4	V80	B	503	23/23	0.91	0.14	15,49,77,83	0
5	BTB	D	504	14/14	0.91	0.14	34,57,76,79	0
5	BTB	B	504	14/14	0.92	0.12	24,43,62,64	0
4	V80	D	503	23/23	0.92	0.15	18,55,80,85	0
5	BTB	C	504	14/14	0.93	0.10	25,57,65,68	0
2	HEM	A	501	43/43	0.95	0.12	38,51,92,100	0
7	CL	A	508	1/1	0.96	0.09	47,47,47,47	0
7	CL	D	506	1/1	0.97	0.04	22,22,22,22	0
2	HEM	C	501	43/43	0.97	0.09	30,42,85,93	0
2	HEM	B	501	43/43	0.98	0.08	11,24,73,81	0
8	GD	A	509	1/1	0.98	0.05	81,81,81,81	0
8	GD	C	510	1/1	0.98	0.05	74,74,74,74	0
8	GD	D	507	1/1	0.98	0.04	40,40,40,40	0
2	HEM	D	501	43/43	0.98	0.07	16,24,69,81	0
8	GD	B	508	1/1	0.99	0.03	38,38,38,38	0
7	CL	C	509	1/1	0.99	0.04	31,31,31,31	0
6	ZN	A	507	1/1	0.99	0.02	35,35,35,35	0
7	CL	B	507	1/1	0.99	0.04	21,21,21,21	0
6	ZN	C	507	1/1	1.00	0.02	30,30,30,30	0

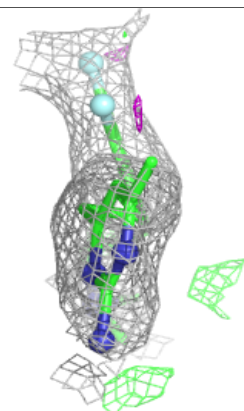
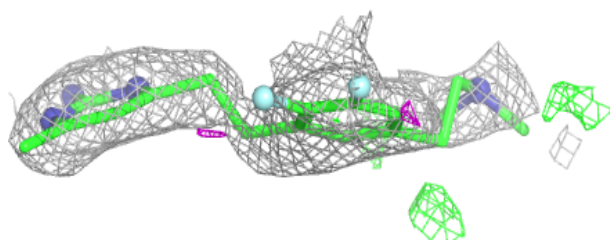
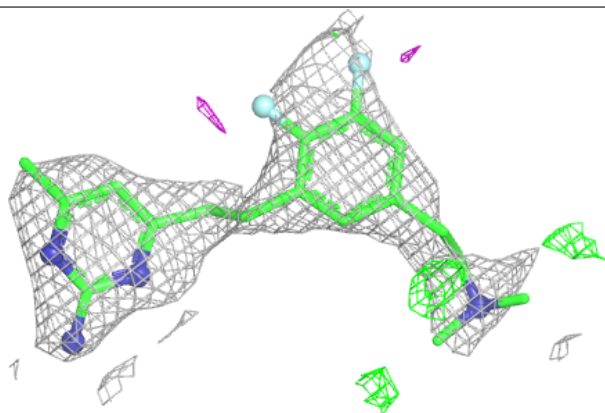
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around V80 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)

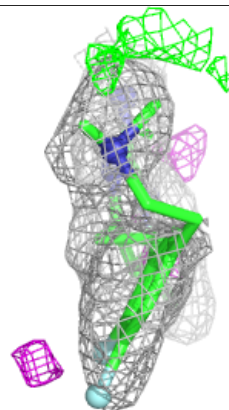
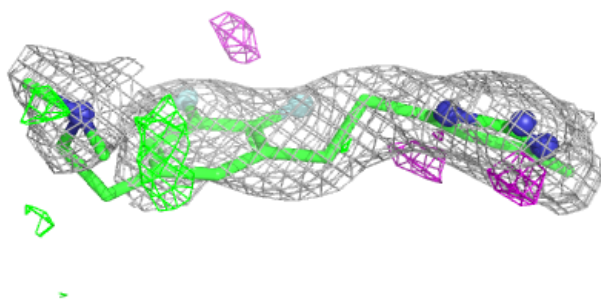
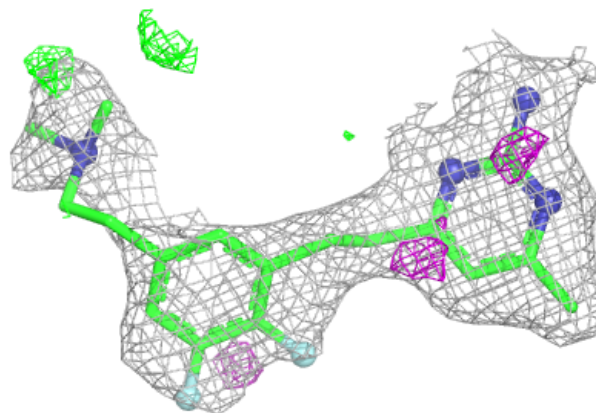
**Electron density around V80 C 503:**

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



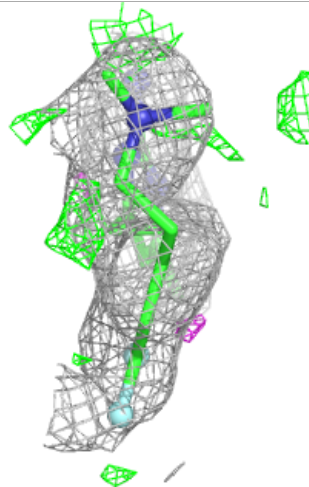
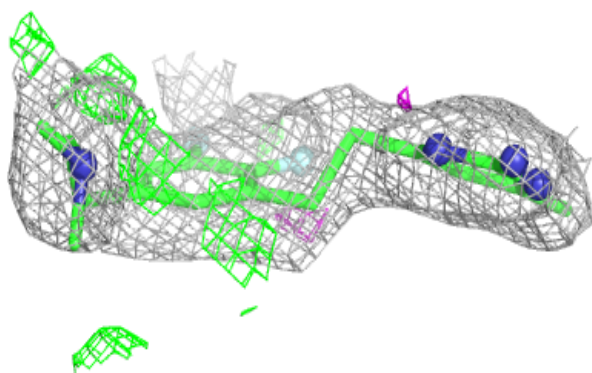
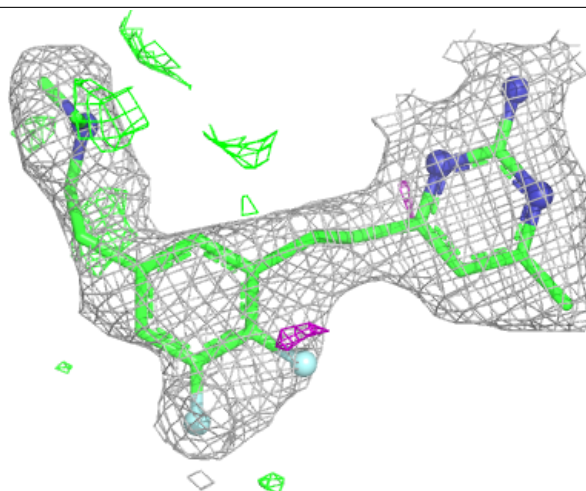
Electron density around V80 B 503:

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



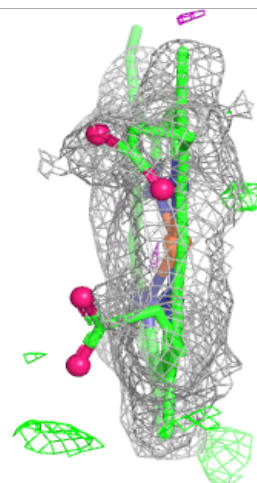
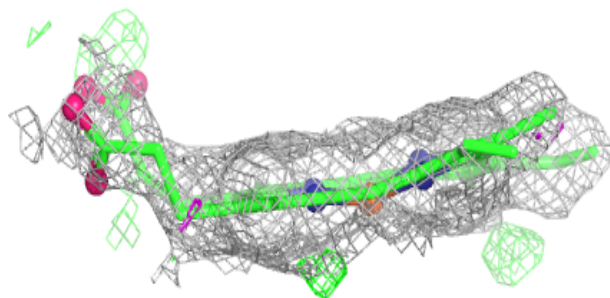
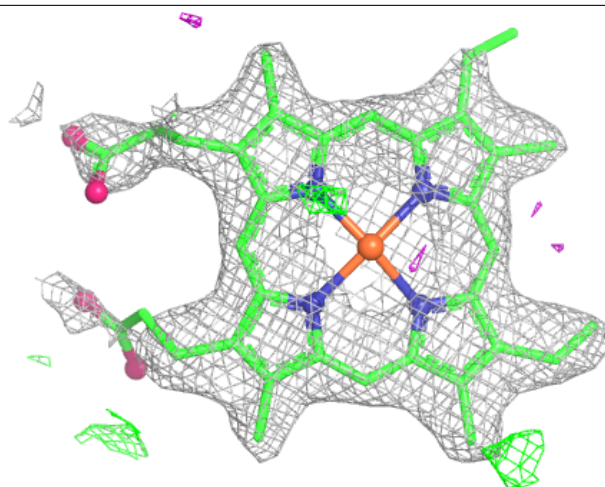
Electron density around V80 D 503:

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



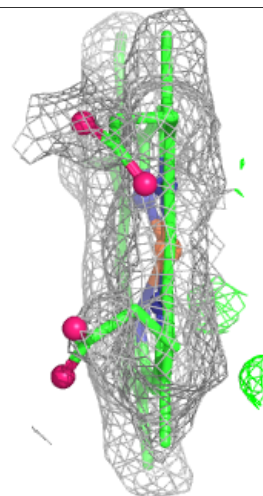
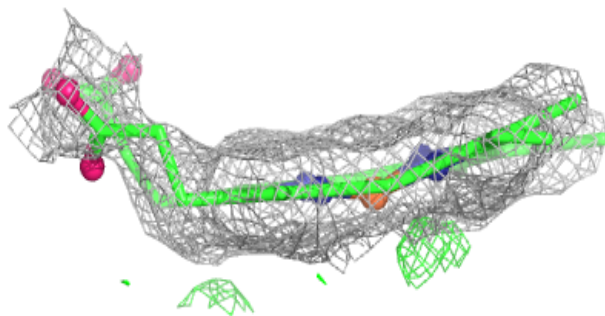
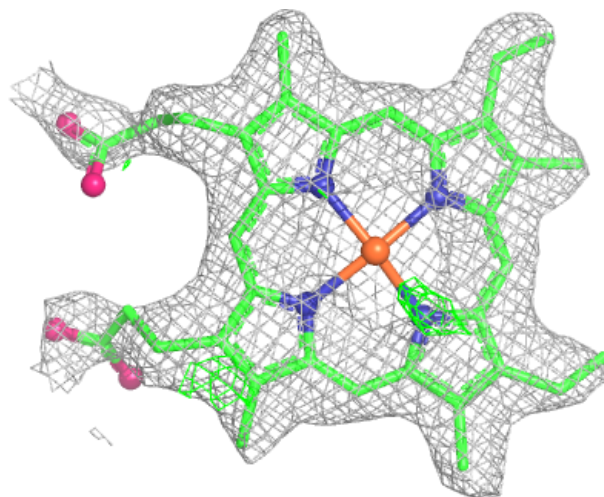
Electron density around HEM A 501:

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



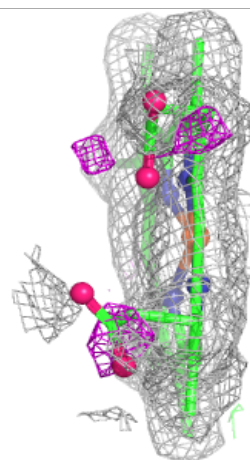
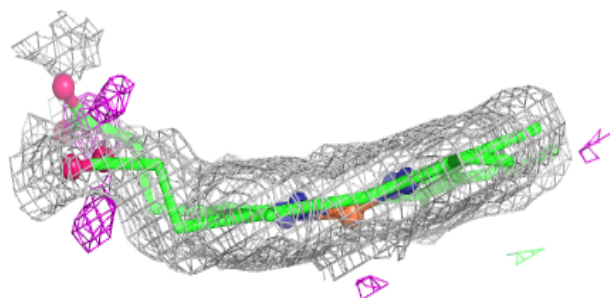
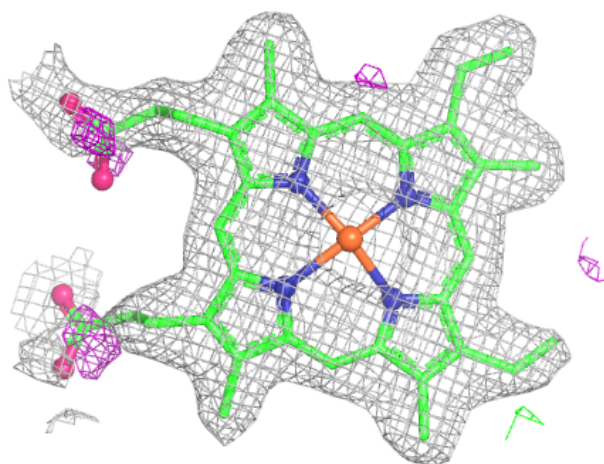
Electron density around HEM C 501:

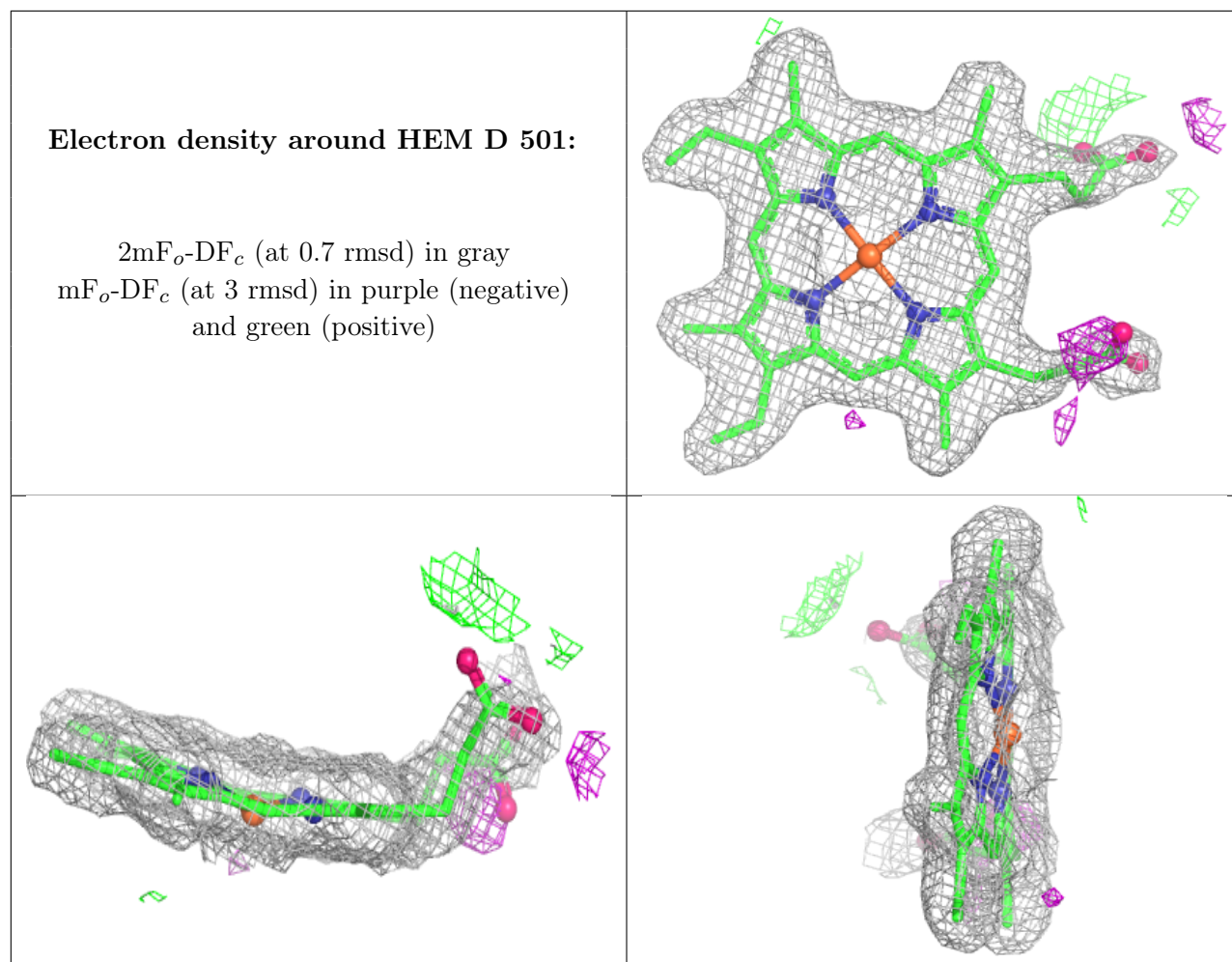
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM B 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.