



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:19 PM UTC

PDB ID : 8UFU / pdb_00008ufu
Title : Structure of human endothelial nitric oxide synthase heme domain in complex with 7-(9-amino-6,7,8,9-tetrahydro-5H-benzo[7]annulen-2-yl)-4-methylquinolin-2-amine
Authors : Li, H.; Poulos, T.L.
Deposited on : 2023-10-04
Resolution : 2.05 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

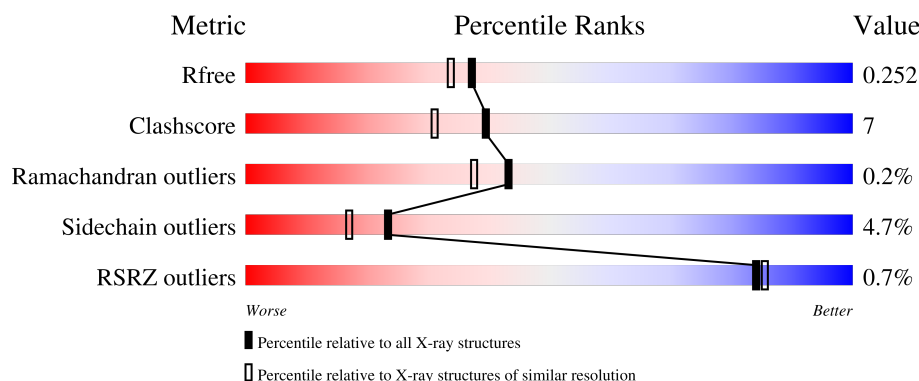
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	% 74% 15% • 9%
1	B	440	77% 13% • 9%
1	C	440	% 73% 16% • 9%
1	D	440	78% 12% • 9%

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 13720 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.

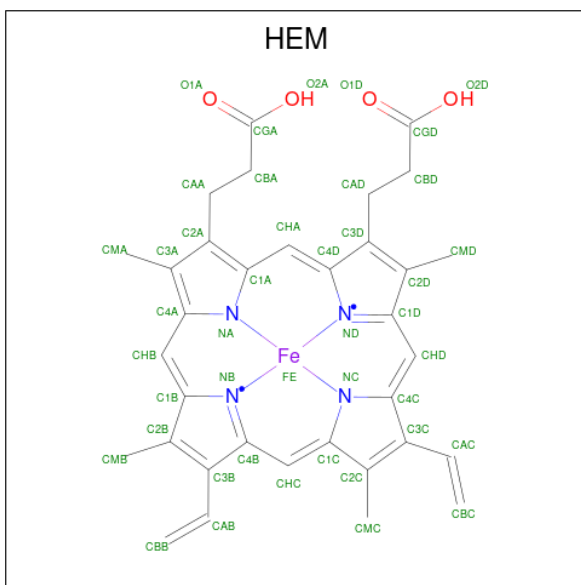
- Molecule 1 is a protein called Nitric oxide synthase, endothelial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	401	Total	C	N	O	S	0	2	0
			3215	2047	566	586	16			
1	B	401	Total	C	N	O	S	0	3	0
			3211	2045	564	586	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:

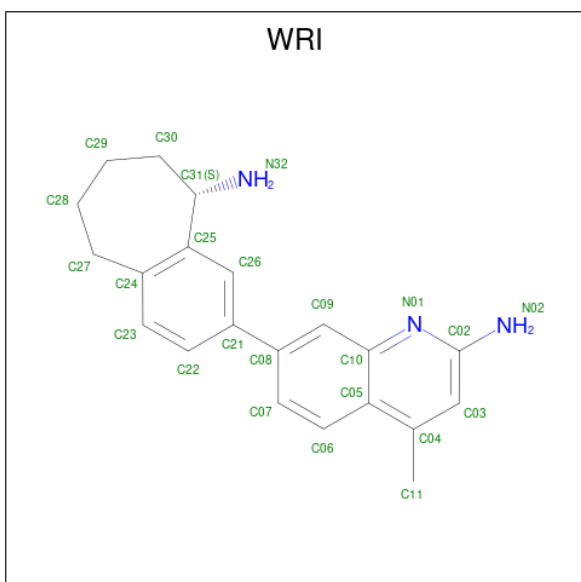
Chain	Residue	Modelled	Actual	Comment	Reference
A	298	GLU	ASP	variant	UNP P29474
B	298	GLU	ASP	variant	UNP P29474
C	298	GLU	ASP	variant	UNP P29474
D	298	GLU	ASP	variant	UNP P29474

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



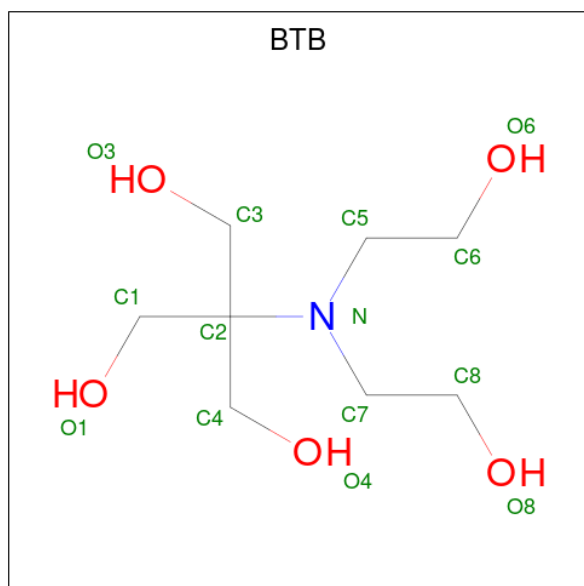
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 3 is (7M)-7-[(9S)-9-amino-6,7,8,9-tetrahydro-5H-benzo[7]annulen-2-yl]-4-methylquinolin-2-amine (CCD ID: WRI) (formula: C₂₁H₂₃N₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			24	21	3		
3	A	1	Total	C	N	0	0
			24	21	3		
3	B	1	Total	C	N	0	0
			24	21	3		
3	B	1	Total	C	N	0	0
			24	21	3		
3	C	1	Total	C	N	0	0
			24	21	3		
3	C	1	Total	C	N	0	0
			24	21	3		
3	C	1	Total	C	N	0	0
			24	21	3		
3	D	1	Total	C	N	0	0
			24	21	3		

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



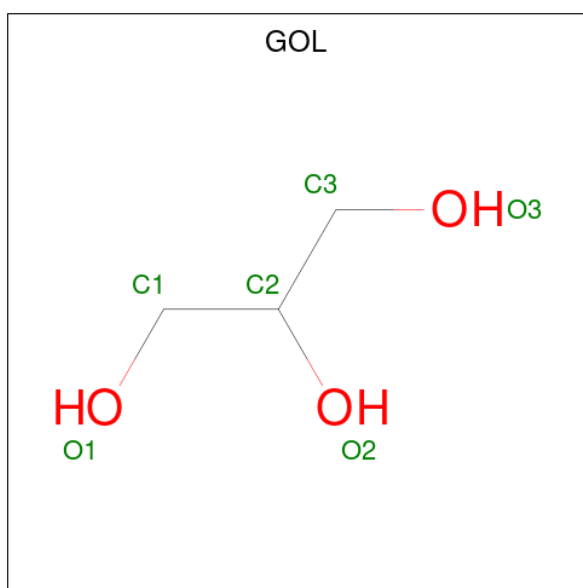
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		

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

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 6 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

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

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

- Molecule 7 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	2	Total 2	Zn 2	0	0
7	B	1	Total 1	Zn 1	0	0
7	C	3	Total 3	Zn 3	0	0

- Molecule 8 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	1	Total 1	Ca 1	0	0
8	C	1	Total 1	Ca 1	0	0

- Molecule 9 is GADOLINIUM ATOM (CCD ID: GD) (formula: Gd).

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

- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	50	Total 50	O 50	0	0
10	B	118	Total 118	O 118	0	0
10	C	75	Total 75	O 75	0	0

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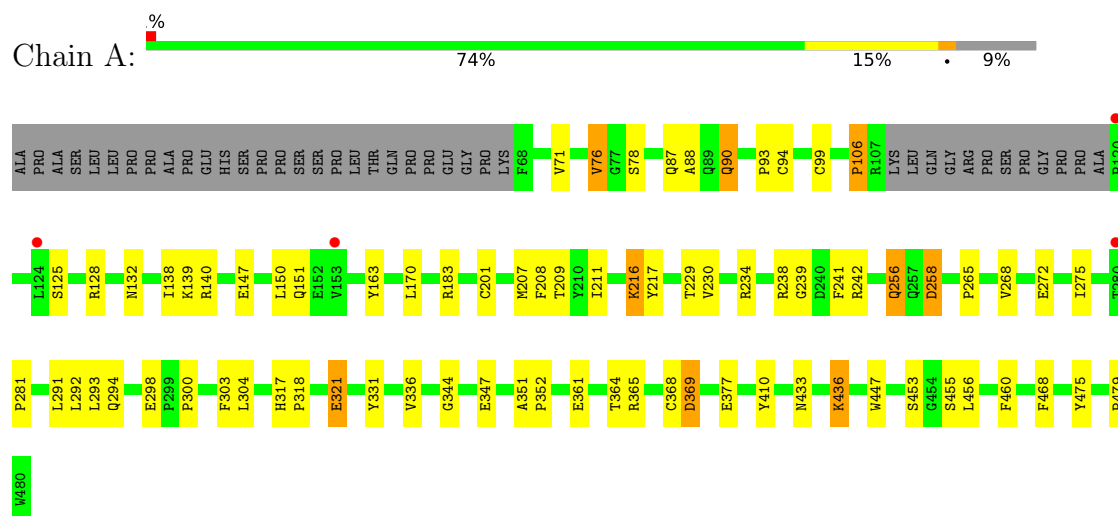
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	D	119	Total	O	0	0
			119	119		

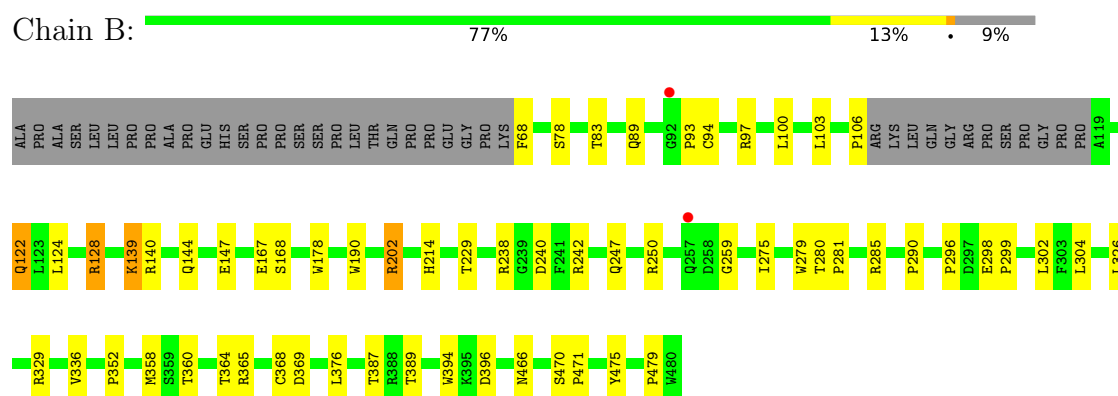
3 Residue-property plots [i](#)

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.

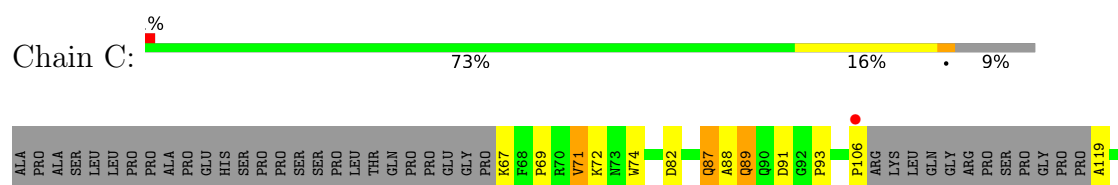
- Molecule 1: Nitric oxide synthase, endothelial

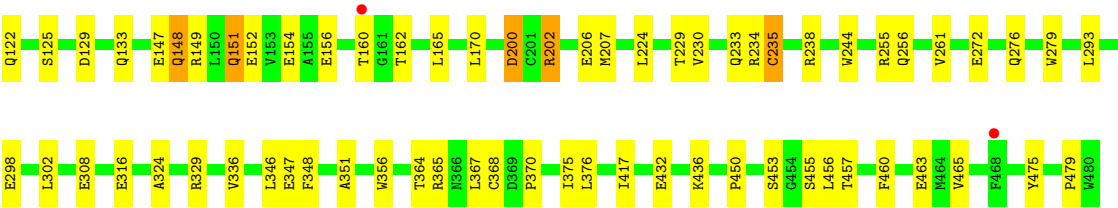


- Molecule 1: Nitric oxide synthase, endothelial

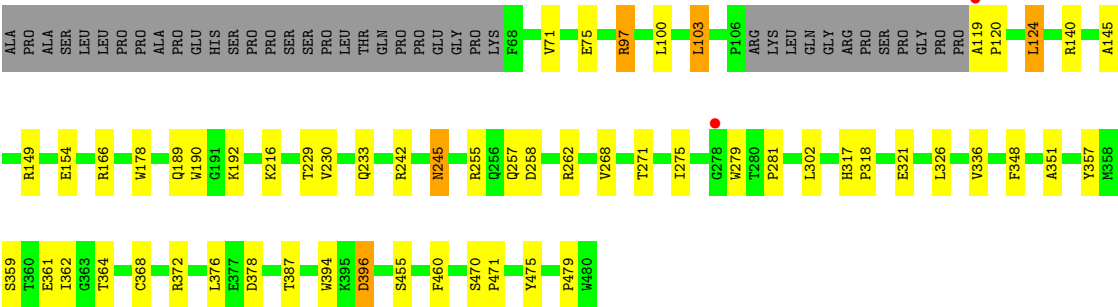
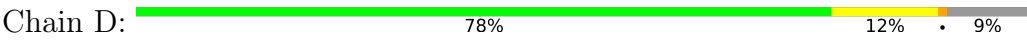


- Molecule 1: Nitric oxide synthase, endothelial





● 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	59.67Å 152.83Å 109.16Å 90.00° 90.82° 90.00°	Depositor
Resolution (Å)	38.74 – 2.05 38.74 – 2.05	Depositor EDS
% Data completeness (in resolution range)	94.2 (38.74-2.05) 94.9 (38.74-2.05)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.40 (at 2.05Å)	Xtriage
Refinement program	PHENIX (1.11.1_2575: ???)	Depositor
R, R_{free}	0.205 , 0.256 0.200 , 0.252	Depositor DCC
R_{free} test set	5867 reflections (3.09%)	wwPDB-VP
Wilson B-factor (Å ²)	42.0	Xtriage
Anisotropy	0.551	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 55.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.108 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13720	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.81% 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: ZN, CL, GOL, WRI, GD, HEM, CA, 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	0.30	0/3313	0.49	0/4513
1	B	0.34	0/3312	0.52	0/4514
1	C	0.32	0/3316	0.50	0/4518
1	D	0.40	1/3304 (0.0%)	0.55	0/4503
All	All	0.34	1/13245 (0.0%)	0.52	0/18048

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	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	372	ARG	CA-C	10.18	1.59	1.53

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	240	ASP	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	3215	0	3118	45	0
1	B	3211	0	3114	35	0
1	C	3218	0	3122	45	0
1	D	3206	0	3108	32	0
2	A	43	0	30	4	0
2	B	43	0	30	3	0
2	C	43	0	30	3	0
2	D	43	0	30	2	0
3	A	48	0	0	3	0
3	B	48	0	0	1	0
3	C	72	0	0	1	0
3	D	24	0	0	2	0
4	A	28	0	38	6	0
4	B	28	0	36	5	0
4	C	28	0	38	7	0
4	D	28	0	36	5	0
5	A	6	0	8	0	0
5	C	6	0	8	0	0
5	D	6	0	8	1	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
6	C	1	0	0	1	0
6	D	1	0	0	0	0
7	A	2	0	0	0	0
7	B	1	0	0	0	0
7	C	3	0	0	0	0
8	A	1	0	0	0	0
8	C	1	0	0	0	0
9	B	1	0	0	0	0
9	D	1	0	0	0	0
10	A	50	0	0	1	0
10	B	118	0	0	1	0
10	C	75	0	0	2	0
10	D	119	0	0	2	0
All	All	13720	0	12754	174	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 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:D:321:GLU:OE2	4:D:503:BTB:O4	2.04	0.76
1:C:234:ARG:NH1	1:C:347:GLU:OE1	2.20	0.75
2:D:501:HEM:HBB2	2:D:501:HEM:HHC	1.69	0.74
1:D:475:TYR:OH	2:D:501:HEM:O2D	2.06	0.73
1:B:124:LEU:HB3	1:B:128:ARG:HH12	1.52	0.73
1:D:100:LEU:HB3	1:D:103:LEU:HD22	1.73	0.70
1:A:433:ASN:HA	1:A:436:LYS:HE3	1.74	0.68
1:C:91:ASP:OD1	10:C:601:HOH:O	2.10	0.68
1:D:154:GLU:OE2	10:D:601:HOH:O	2.11	0.68
1:B:475:TYR:OH	2:B:501:HEM:O2D	2.11	0.67
1:C:475:TYR:OH	2:C:501:HEM:O2D	2.13	0.67
1:C:453:SER:HB3	1:C:456:LEU:HD12	1.77	0.66
1:C:88:ALA:HB3	1:D:97:ARG:HD2	1.77	0.66
1:D:255:ARG:HD3	1:D:268:VAL:HG11	1.77	0.66
2:A:501:HEM:HBB2	2:A:501:HEM:HHC	1.77	0.66
1:B:247:GLN:HB2	1:B:250:ARG:HD3	1.78	0.65
2:B:501:HEM:HHC	2:B:501:HEM:HBB2	1.77	0.65
1:C:200:ASP:OD1	1:C:200:ASP:N	2.29	0.65
1:C:450:PRO:HG2	1:C:457:THR:HG21	1.79	0.64
1:C:93:PRO:HG3	1:C:106:PRO:HB3	1.81	0.63
1:D:279:TRP:HB2	1:D:302:LEU:HD21	1.81	0.62
2:B:501:HEM:HBC2	2:B:501:HEM:HMC2	1.81	0.62
1:C:365:ARG:NH1	2:C:501:HEM:O2A	2.33	0.62
1:C:119:ALA:HB1	1:C:122:GLN:HB2	1.82	0.61
1:B:100:LEU:HB3	1:B:103:LEU:HD22	1.83	0.61
1:B:298:GLU:OE2	4:B:505:BTB:N	2.34	0.61
2:A:501:HEM:HMC2	2:A:501:HEM:HBC2	1.81	0.60
1:A:128:ARG:O	1:A:132:ASN:ND2	2.34	0.60
1:A:275:ILE:HD11	1:A:281:PRO:HB3	1.84	0.60
1:A:234:ARG:NH1	1:A:347:GLU:OE1	2.34	0.59
1:A:475:TYR:OH	2:A:501:HEM:O2D	2.18	0.59
1:D:140:ARG:HH12	1:D:145:ALA:HB3	1.67	0.59
1:A:321:GLU:H	1:A:321:GLU:CD	2.09	0.59
4:C:504:BTB:HO3	4:C:504:BTB:HO8	1.51	0.59
1:C:147:GLU:O	1:C:151:GLN:HG2	2.04	0.58
1:A:90:GLN:HB3	1:A:468:PHE:CD2	2.38	0.58
1:C:370:PRO:HG2	1:D:75:GLU:HG2	1.86	0.58
1:D:124:LEU:HD11	1:D:154:GLU:HG3	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:285:ARG:NH2	10:B:601:HOH:O	2.23	0.57
1:B:93:PRO:HG3	1:B:106:PRO:HB3	1.85	0.56
1:A:377:GLU:OE1	4:A:504:BTB:H11	2.05	0.56
4:B:505:BTB:H81	4:B:505:BTB:O6	2.06	0.56
1:C:160:THR:HG23	1:C:162:THR:H	1.69	0.56
1:A:292:LEU:HD22	1:A:300:PRO:HB2	1.87	0.56
1:A:256:GLN:C	1:A:258:ASP:H	2.13	0.55
1:A:453:SER:HB3	1:A:456:LEU:HD12	1.89	0.55
1:A:238:ARG:HH21	1:A:241:PHE:HD1	1.56	0.53
1:B:279:TRP:CD1	1:B:290:PRO:HG3	2.43	0.53
1:B:279:TRP:HB2	1:B:302:LEU:HD21	1.89	0.53
1:C:156:GLU:O	1:C:160:THR:HG22	2.08	0.53
1:B:229:THR:O	1:B:352:PRO:HD2	2.08	0.53
1:B:336:VAL:HG21	3:B:502:WRI:C07	2.39	0.52
1:A:331:TYR:O	1:A:410:TYR:OH	2.28	0.52
1:B:238:ARG:HG2	1:B:296:PRO:HB3	1.92	0.52
1:A:88:ALA:O	1:B:97:ARG:NH1	2.43	0.52
1:B:298:GLU:HG3	1:B:299:PRO:HD2	1.92	0.51
1:A:242:ARG:NH2	1:A:479:PRO:HD3	2.25	0.51
1:B:242:ARG:NH2	1:B:479:PRO:HD3	2.25	0.51
1:A:93:PRO:HG3	1:A:106:PRO:HB3	1.92	0.51
1:A:258:ASP:OD1	1:A:258:ASP:N	2.44	0.50
1:A:361:GLU:OE2	3:A:502:WRI:N02	2.44	0.50
1:C:87:GLN:O	1:C:89:GLN:NE2	2.44	0.50
1:D:387:THR:HA	1:D:394:TRP:CD1	2.46	0.50
1:B:368:CYS:SG	1:B:376:LEU:HD13	2.52	0.50
1:D:455:SER:HA	1:D:460:PHE:CG	2.46	0.50
1:C:356:TRP:CZ2	2:C:501:HEM:HBB1	2.47	0.49
1:A:147:GLU:HA	1:A:150:LEU:HD12	1.95	0.49
1:A:291:LEU:HD23	1:A:293:LEU:HD21	1.92	0.49
1:B:364:THR:O	1:B:368:CYS:HB2	2.12	0.49
1:C:336:VAL:HG21	3:C:502:WRI:C07	2.42	0.49
1:A:365:ARG:O	1:A:369:ASP:HB2	2.12	0.49
4:C:504:BTB:O8	4:C:504:BTB:O3	2.18	0.49
1:C:207:MET:HE3	1:C:293:LEU:HB3	1.94	0.49
1:D:178:TRP:CE3	1:D:190:TRP:HA	2.48	0.48
1:D:245:ASN:N	1:D:245:ASN:OD1	2.46	0.48
1:C:453:SER:OG	1:D:396:ASP:OD1	2.29	0.48
1:B:298:GLU:CD	4:B:505:BTB:H31	2.39	0.48
1:C:279:TRP:HB2	1:C:302:LEU:HD11	1.96	0.48
1:C:364:THR:O	1:C:368:CYS:HB2	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:316[B]:GLU:HG2	1:C:324:ALA:HB2	1.94	0.48
1:A:76:VAL:HG12	1:A:78:SER:H	1.78	0.47
1:D:271:THR:O	1:D:275:ILE:HG12	2.14	0.47
1:C:71:VAL:HG21	1:C:465:VAL:HG23	1.96	0.47
1:A:170:LEU:HD11	1:A:230:VAL:HG11	1.96	0.47
1:C:133:GLN:NE2	10:C:605:HOH:O	2.40	0.47
4:D:504:BTB:H31	4:D:504:BTB:H71	1.36	0.47
4:A:505:BTB:H41	4:A:505:BTB:H72	1.60	0.47
1:C:202:ARG:HG2	1:C:206:GLU:OE1	2.15	0.47
1:A:183:ARG:HG2	1:A:447:TRP:CG	2.50	0.47
1:A:128:ARG:HE	1:A:128:ARG:HB3	1.53	0.46
1:C:298:GLU:OE1	4:C:505:BTB:O1	2.31	0.46
1:A:336:VAL:HG21	3:A:502:WRI:C07	2.46	0.46
1:B:122:GLN:CD	1:B:122:GLN:H	2.23	0.46
1:B:202:ARG:HB2	1:B:202:ARG:NH1	2.31	0.46
1:D:189:GLN:HB3	10:D:666:HOH:O	2.14	0.46
4:C:504:BTB:H12	4:C:504:BTB:H51	1.82	0.46
1:D:368:CYS:SG	1:D:376:LEU:HD13	2.55	0.46
1:C:233:GLN:HB3	1:C:348:PHE:CE2	2.51	0.46
4:C:505:BTB:H42	4:C:505:BTB:H51	1.48	0.46
1:B:275:ILE:HG12	1:B:281:PRO:HG3	1.97	0.45
1:A:455:SER:HA	1:A:460:PHE:CG	2.51	0.45
2:A:501:HEM:HBD1	3:A:502:WRI:C22	2.47	0.45
1:C:148:GLN:C	1:C:148:GLN:HE21	2.25	0.45
1:C:235:CYS:SG	1:C:238:ARG:HD3	2.56	0.45
1:A:364:THR:O	1:A:368:CYS:HB2	2.16	0.45
4:A:504:BTB:O3	4:A:504:BTB:O1	2.32	0.45
4:D:504:BTB:H52	4:D:504:BTB:H82	1.43	0.45
1:B:214:HIS:CD2	1:B:214:HIS:C	2.94	0.45
1:D:336:VAL:HG21	3:D:502:WRI:C07	2.47	0.45
1:C:368:CYS:SG	1:C:376:LEU:HD13	2.57	0.45
1:C:229:THR:O	1:C:351:ALA:HA	2.17	0.45
4:D:504:BTB:H51	4:D:504:BTB:H42	1.38	0.44
1:C:455:SER:HA	1:C:460:PHE:CG	2.52	0.44
1:A:147:GLU:O	1:A:151:GLN:HG2	2.17	0.44
1:A:216:LYS:HG3	1:A:217:TYR:N	2.31	0.44
1:A:265:PRO:HA	1:A:268:VAL:HG23	1.99	0.44
4:B:505:BTB:O3	4:B:505:BTB:H51	2.17	0.44
1:A:138:ILE:HG13	1:A:140:ARG:HB2	1.99	0.44
1:B:358:MET:HE3	1:B:360:THR:OG1	2.18	0.44
1:C:272:GLU:O	1:C:276:GLN:HG3	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:208:PHE:CE1	1:A:303:PHE:HB3	2.53	0.43
4:A:504:BTB:H12	10:A:647:HOH:O	2.18	0.43
1:A:163:TYR:OH	1:A:344:GLY:O	2.33	0.43
4:A:505:BTB:H11	4:A:505:BTB:H51	1.75	0.43
1:C:432:GLU:O	1:C:436:LYS:HE3	2.17	0.43
4:C:505:BTB:H31	4:C:505:BTB:H72	1.39	0.43
1:B:202:ARG:HB2	1:B:202:ARG:HH11	1.84	0.43
1:A:217:TYR:C	1:A:217:TYR:CD2	2.96	0.43
1:D:233:GLN:HB3	1:D:348:PHE:CE2	2.53	0.43
1:D:242:ARG:NH2	1:D:479:PRO:HD3	2.33	0.43
1:D:470:SER:HA	1:D:471:PRO:C	2.43	0.43
1:A:317:HIS:CG	1:A:318:PRO:HD2	2.53	0.43
1:D:361:GLU:OE2	3:D:502:WRI:N02	2.52	0.43
1:A:99:CYS:HB3	1:B:466:ASN:HB3	2.01	0.43
1:B:139:LYS:O	1:B:140:ARG:HD2	2.19	0.43
1:C:255:ARG:HG3	1:C:261:VAL:HG22	2.00	0.42
1:B:279:TRP:CG	1:B:290:PRO:HG3	2.54	0.42
4:C:504:BTB:H72	4:C:504:BTB:H41	1.57	0.42
1:C:463:GLU:HB3	1:D:103:LEU:HD12	2.02	0.42
1:A:211:ILE:HG21	1:A:291:LEU:HD21	2.01	0.42
1:A:139:LYS:HB2	1:A:139:LYS:HE3	1.82	0.42
1:A:207:MET:HB3	1:A:207:MET:HE2	1.81	0.42
4:A:504:BTB:H72	4:A:504:BTB:H41	1.61	0.42
1:B:280:THR:HA	1:B:281:PRO:HD3	1.89	0.42
1:C:244:TRP:CD1	1:C:479:PRO:HG2	2.54	0.42
1:A:234:ARG:HA	1:A:238:ARG:HH12	1.85	0.42
1:C:235:CYS:SG	1:C:238:ARG:NH1	2.92	0.42
1:A:242:ARG:HB2	1:A:294:GLN:NE2	2.34	0.41
1:C:367:LEU:O	1:C:375:ILE:HG12	2.20	0.41
1:D:364:THR:O	1:D:368:CYS:HB2	2.20	0.41
1:B:68:PHE:CD1	1:B:83:THR:HG22	2.55	0.41
1:D:357:TYR:CD2	1:D:362:ILE:HD11	2.55	0.41
1:A:229:THR:O	1:A:351:ALA:HA	2.21	0.41
1:C:69:PRO:HG2	1:C:82:ASP:HB3	2.02	0.41
1:D:149:ARG:HD3	1:D:166:ARG:CZ	2.49	0.41
1:A:94:CYS:HB3	1:B:94:CYS:HB3	2.03	0.41
1:A:229:THR:O	1:A:352:PRO:HD2	2.20	0.41
1:B:470:SER:HA	1:B:471:PRO:C	2.46	0.41
1:D:229:THR:O	1:D:351:ALA:HA	2.20	0.41
1:B:365:ARG:O	1:B:369:ASP:HB2	2.20	0.41
1:B:124:LEU:HD23	1:B:124:LEU:HA	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:89:GLN:H	1:C:89:GLN:HG2	1.60	0.41
4:D:503:BTB:H12	4:D:503:BTB:H72	1.74	0.41
1:C:170:LEU:HD11	1:C:230:VAL:HG11	2.03	0.41
1:D:119:ALA:N	1:D:120:PRO:HD3	2.36	0.41
4:B:505:BTB:H42	4:B:505:BTB:H72	1.82	0.40
1:D:275:ILE:HD12	1:D:281:PRO:HG3	2.03	0.40
1:C:165:LEU:HG	1:C:346:LEU:HD12	2.04	0.40
1:D:317:HIS:CG	1:D:318:PRO:HD2	2.56	0.40
1:B:387:THR:HA	1:B:394:TRP:CD1	2.57	0.40
1:D:378:ASP:OD2	5:D:505:GOL:O2	2.39	0.40
1:B:178:TRP:CE3	1:B:190:TRP:HA	2.57	0.40
1:C:72:LYS:HE2	1:C:74:TRP:CE3	2.57	0.40
1:C:365:ARG:NH2	6:C:507:CL:CL	2.92	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	399/440 (91%)	381 (96%)	16 (4%)	2 (0%)	24	16
1	B	400/440 (91%)	388 (97%)	11 (3%)	1 (0%)	36	30
1	C	400/440 (91%)	392 (98%)	8 (2%)	0	100	100
1	D	399/440 (91%)	389 (98%)	10 (2%)	0	100	100
All	All	1598/1760 (91%)	1550 (97%)	45 (3%)	3 (0%)	43	37

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	106	PRO
1	A	239	GLY

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Mol	Chain	Res	Type
1	B	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	343/373 (92%)	327 (95%)	16 (5%)	23	17
1	B	343/373 (92%)	327 (95%)	16 (5%)	23	17
1	C	343/373 (92%)	324 (94%)	19 (6%)	19	12
1	D	342/373 (92%)	328 (96%)	14 (4%)	27	21
All	All	1371/1492 (92%)	1306 (95%)	65 (5%)	23	17

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	VAL
1	A	76	VAL
1	A	87	GLN
1	A	90	GLN
1	A	125	SER
1	A	201	CYS
1	A	209	THR
1	A	216	LYS
1	A	256	GLN
1	A	258	ASP
1	A	272	GLU
1	A	298	GLU
1	A	304	LEU
1	A	321	GLU
1	A	369	ASP
1	A	436	LYS
1	B	78	SER
1	B	89	GLN
1	B	122	GLN
1	B	128	ARG

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Mol	Chain	Res	Type
1	B	139	LYS
1	B	144	GLN
1	B	147	GLU
1	B	167	GLU
1	B	168[A]	SER
1	B	168[B]	SER
1	B	202	ARG
1	B	304	LEU
1	B	326	LEU
1	B	329	ARG
1	B	389	THR
1	B	396	ASP
1	C	67	LYS
1	C	71	VAL
1	C	87	GLN
1	C	89	GLN
1	C	125	SER
1	C	129	ASP
1	C	148	GLN
1	C	149	ARG
1	C	151	GLN
1	C	152	GLU
1	C	154	GLU
1	C	200	ASP
1	C	202	ARG
1	C	224	LEU
1	C	235	CYS
1	C	256	GLN
1	C	308	GLU
1	C	329	ARG
1	C	417	ILE
1	D	71	VAL
1	D	97	ARG
1	D	103	LEU
1	D	124	LEU
1	D	192	LYS
1	D	216	LYS
1	D	230	VAL
1	D	245	ASN
1	D	257	GLN
1	D	258	ASP
1	D	262	ARG

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Mol	Chain	Res	Type
1	D	326	LEU
1	D	359	SER
1	D	396	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	133	GLN
1	A	256	GLN
1	B	148	GLN
1	B	247	GLN
1	C	148	GLN
1	C	257	GLN
1	C	433	ASN
1	D	126	GLN

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 37 ligands modelled in this entry, 14 are monoatomic - leaving 23 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
3	WRI	C	510	-	27,27,27	0.95	1 (3%)	36,39,39	1.47	7 (19%)
5	GOL	A	506	-	5,5,5	0.34	0	5,5,5	0.42	0
4	BTB	A	505	-	13,13,13	0.37	0	7,16,16	0.51	0
4	BTB	C	504	-	13,13,13	0.72	0	7,16,16	1.76	2 (28%)
2	HEM	C	501	1	50,50,50	1.71	10 (20%)	67,82,82	1.56	14 (20%)
3	WRI	B	502	-	27,27,27	1.06	3 (11%)	36,39,39	2.19	13 (36%)
4	BTB	D	503	9	13,13,13	0.43	0	7,16,16	1.09	0
3	WRI	D	502	-	27,27,27	0.85	0	36,39,39	1.46	4 (11%)
4	BTB	B	505	-	13,13,13	0.56	0	7,16,16	1.00	0
3	WRI	C	502	-	27,27,27	0.92	2 (7%)	36,39,39	1.48	7 (19%)
3	WRI	A	502	-	27,27,27	0.93	1 (3%)	36,39,39	1.72	7 (19%)
4	BTB	C	505	-	13,13,13	0.41	0	7,16,16	0.70	0
3	WRI	A	503	-	27,27,27	0.94	0	36,39,39	1.29	5 (13%)
5	GOL	D	505	-	5,5,5	0.37	0	5,5,5	0.29	0
5	GOL	C	506	-	5,5,5	0.43	0	5,5,5	0.43	0
2	HEM	B	501	1	50,50,50	1.78	8 (16%)	67,82,82	1.52	11 (16%)
4	BTB	D	504	-	13,13,13	0.53	0	7,16,16	1.06	1 (14%)
2	HEM	A	501	1	50,50,50	1.63	7 (14%)	67,82,82	1.44	11 (16%)
2	HEM	D	501	1	50,50,50	1.76	8 (16%)	67,82,82	1.43	15 (22%)
4	BTB	B	504	9	13,13,13	0.59	0	7,16,16	0.67	0
3	WRI	B	503	-	27,27,27	0.91	1 (3%)	36,39,39	1.18	5 (13%)
3	WRI	C	503	-	27,27,27	0.96	1 (3%)	36,39,39	1.18	3 (8%)
4	BTB	A	504	-	13,13,13	0.53	0	7,16,16	0.86	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
3	WRI	C	510	-	-	4/4/15/15	0/4/4/4
5	GOL	A	506	-	-	4/4/4/4	-
4	BTB	A	505	-	-	3/21/21/21	-
4	BTB	C	504	-	-	10/21/21/21	-
2	HEM	C	501	1	-	2/14/54/54	-
3	WRI	B	502	-	-	0/4/15/15	0/4/4/4
4	BTB	D	503	9	-	1/21/21/21	-
3	WRI	D	502	-	-	0/4/15/15	0/4/4/4

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	BTB	B	505	-	-	3/21/21/21	-
3	WRI	C	502	-	-	0/4/15/15	0/4/4/4
3	WRI	A	502	-	-	0/4/15/15	0/4/4/4
4	BTB	C	505	-	-	13/21/21/21	-
3	WRI	A	503	-	-	4/4/15/15	1/4/4/4
5	GOL	D	505	-	-	2/4/4/4	-
5	GOL	C	506	-	-	4/4/4/4	-
2	HEM	B	501	1	-	2/14/54/54	-
4	BTB	D	504	-	-	13/21/21/21	-
2	HEM	A	501	1	-	1/14/54/54	-
2	HEM	D	501	1	-	1/14/54/54	-
4	BTB	B	504	9	-	3/21/21/21	-
3	WRI	B	503	-	-	4/4/15/15	0/4/4/4
3	WRI	C	503	-	-	4/4/15/15	0/4/4/4
4	BTB	A	504	-	-	4/21/21/21	-

All (42) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	501	HEM	FE-NB	5.97	2.13	1.94
2	B	501	HEM	FE-NB	5.81	2.12	1.94
2	B	501	HEM	FE-NC	5.71	2.14	1.95
2	C	501	HEM	FE-NC	5.41	2.13	1.95
2	A	501	HEM	FE-NC	4.88	2.11	1.95
2	D	501	HEM	FE-NC	4.61	2.10	1.95
2	D	501	HEM	FE-NA	4.57	2.10	1.95
2	C	501	HEM	FE-NA	4.51	2.10	1.95
2	A	501	HEM	FE-NB	4.41	2.08	1.94
2	C	501	HEM	FE-ND	4.36	2.08	1.94
2	A	501	HEM	FE-NA	4.34	2.09	1.95
2	B	501	HEM	FE-NA	3.98	2.08	1.95
2	C	501	HEM	FE-NB	3.86	2.06	1.94
2	B	501	HEM	FE-ND	3.56	2.05	1.94
2	B	501	HEM	CAB-C3B	3.12	1.55	1.47
2	A	501	HEM	CAB-C3B	3.12	1.55	1.47
2	C	501	HEM	CAB-C3B	3.06	1.55	1.47
2	D	501	HEM	CAB-C3B	3.00	1.55	1.47
2	B	501	HEM	CAC-C3C	2.99	1.55	1.47
2	A	501	HEM	CAC-C3C	2.98	1.55	1.47
2	D	501	HEM	CAC-C3C	2.89	1.55	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	HEM	FE-ND	2.88	2.03	1.94
2	C	501	HEM	CAC-C3C	2.72	1.54	1.47
2	D	501	HEM	FE-ND	2.65	2.03	1.94
3	B	502	WRI	C05-C10	-2.64	1.38	1.42
3	C	510	WRI	C02-N01	2.61	1.36	1.33
2	D	501	HEM	CMA-C3A	2.51	1.55	1.50
3	B	502	WRI	C02-N01	2.44	1.36	1.33
3	C	502	WRI	C02-N01	2.39	1.36	1.33
2	B	501	HEM	CMA-C3A	2.20	1.55	1.50
3	B	502	WRI	C04-C05	-2.14	1.38	1.42
3	A	502	WRI	C05-C10	-2.13	1.39	1.42
2	C	501	HEM	CMB-C2B	2.11	1.55	1.50
2	D	501	HEM	CMB-C2B	2.11	1.55	1.50
3	C	502	WRI	C05-C10	-2.10	1.39	1.42
3	C	503	WRI	C03-C04	2.09	1.40	1.37
3	B	503	WRI	C02-N01	2.08	1.36	1.33
2	C	501	HEM	CMC-C2C	2.07	1.55	1.50
2	C	501	HEM	CMD-C2D	2.06	1.55	1.50
2	C	501	HEM	CMA-C3A	2.05	1.55	1.50
2	A	501	HEM	CMC-C2C	2.05	1.55	1.50
2	B	501	HEM	CMB-C2B	2.03	1.54	1.50

All (105) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	502	WRI	C22-C21-C08	-4.94	112.73	121.25
3	B	502	WRI	N02-C02-N01	4.88	122.27	118.24
3	B	502	WRI	C26-C21-C08	4.28	128.01	120.84
2	C	501	HEM	C3D-C4D-ND	-4.26	105.50	110.17
3	A	502	WRI	C04-C05-C10	4.26	120.59	118.00
3	A	502	WRI	C05-C10-N01	-4.16	118.40	122.80
3	B	502	WRI	C29-C28-C27	-3.67	107.06	114.27
2	C	501	HEM	C2A-C1A-NA	-3.65	106.11	110.15
2	A	501	HEM	C3B-C2B-C1B	3.62	109.13	106.41
2	B	501	HEM	C3B-C2B-C1B	3.56	109.08	106.41
3	D	502	WRI	C05-C10-N01	-3.56	119.03	122.80
2	D	501	HEM	C3D-C4D-ND	-3.54	106.29	110.17
2	A	501	HEM	C3B-C4B-NB	-3.53	106.93	109.47
3	C	510	WRI	C22-C21-C08	-3.51	115.19	121.25
2	B	501	HEM	C3B-C4B-NB	-3.46	106.99	109.47
3	A	502	WRI	C07-C08-C21	-3.34	115.48	121.25
3	B	502	WRI	C05-C10-N01	-3.33	119.27	122.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	502	WRI	C04-C05-C10	3.30	120.01	118.00
2	C	501	HEM	C4D-ND-C1D	3.29	109.11	105.21
3	C	503	WRI	C22-C21-C08	-3.29	115.57	121.25
3	C	502	WRI	C05-C10-N01	-3.28	119.33	122.80
2	B	501	HEM	C1B-NB-C4B	3.24	109.04	105.21
4	C	504	BTB	O3-C3-C2	-3.20	103.87	111.40
2	B	501	HEM	C4D-ND-C1D	3.19	108.98	105.21
2	A	501	HEM	C1B-NB-C4B	3.18	108.98	105.21
3	D	502	WRI	C29-C30-C31	-3.18	108.15	114.25
2	D	501	HEM	C3B-C2B-C1B	3.15	108.78	106.41
3	B	502	WRI	C07-C08-C21	-3.15	115.82	121.25
3	A	503	WRI	C22-C21-C08	-3.11	115.89	121.25
3	C	510	WRI	N02-C02-N01	3.10	120.81	118.24
3	D	502	WRI	C04-C05-C10	3.10	119.89	118.00
2	D	501	HEM	C4D-ND-C1D	3.08	108.85	105.21
2	C	501	HEM	CHC-C4B-NB	3.07	127.73	124.42
3	C	502	WRI	C28-C27-C24	-3.04	107.54	114.03
2	C	501	HEM	CAD-C3D-C2D	-3.00	122.25	127.87
3	C	503	WRI	C26-C21-C08	2.97	125.82	120.84
2	C	501	HEM	C1B-NB-C4B	2.94	108.68	105.21
3	C	502	WRI	N02-C02-N01	2.90	120.64	118.24
3	A	502	WRI	N02-C02-N01	2.88	120.62	118.24
2	A	501	HEM	C3D-C4D-ND	-2.88	107.01	110.17
2	B	501	HEM	C3D-C4D-ND	-2.86	107.03	110.17
3	C	502	WRI	C04-C05-C10	2.86	119.74	118.00
3	C	510	WRI	C24-C25-C31	-2.81	120.81	124.73
2	C	501	HEM	CAD-CBD-CGD	-2.76	106.34	113.67
2	B	501	HEM	C2D-C1D-ND	-2.69	106.80	109.90
2	A	501	HEM	C4D-ND-C1D	2.69	108.39	105.21
3	B	502	WRI	C24-C25-C31	-2.69	120.98	124.73
3	B	502	WRI	C03-C04-C05	2.68	120.55	117.84
3	A	502	WRI	C24-C25-C31	-2.67	121.00	124.73
2	D	501	HEM	CAD-CBD-CGD	-2.64	106.67	113.67
2	B	501	HEM	CHD-C1D-ND	2.62	127.24	124.42
4	C	504	BTB	O4-C4-C2	-2.59	105.31	111.40
3	C	510	WRI	C26-C21-C08	2.58	125.17	120.84
3	B	502	WRI	C06-C05-C10	2.55	121.19	118.36
2	D	501	HEM	C2A-C1A-NA	-2.54	107.33	110.15
2	C	501	HEM	C4A-NA-C1A	2.52	109.94	105.82
2	A	501	HEM	CHD-C1D-ND	2.51	127.13	124.42
3	B	502	WRI	C06-C05-C04	-2.49	118.79	123.61
3	C	510	WRI	C05-C10-N01	-2.49	120.17	122.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	HEM	CHB-C1B-NB	2.46	127.41	124.37
2	D	501	HEM	CHC-C1C-NC	2.46	127.13	124.45
2	B	501	HEM	C2B-C1B-NB	-2.46	107.01	109.84
3	D	502	WRI	C22-C21-C08	-2.46	117.02	121.25
3	B	503	WRI	C04-C05-C10	2.42	119.48	118.00
3	B	503	WRI	C05-C10-N01	-2.42	120.23	122.80
2	A	501	HEM	C2B-C1B-NB	-2.42	107.06	109.84
2	C	501	HEM	CHD-C1D-ND	2.39	127.00	124.42
2	C	501	HEM	C3B-C4B-NB	-2.39	107.75	109.47
2	A	501	HEM	CAD-CBD-CGD	-2.34	107.46	113.67
2	D	501	HEM	CHB-C1B-NB	2.34	127.26	124.37
3	C	502	WRI	C29-C28-C27	-2.32	109.72	114.27
3	A	503	WRI	C29-C28-C27	-2.31	109.72	114.27
2	D	501	HEM	CHD-C1D-ND	2.29	126.89	124.42
4	D	504	BTB	O4-C4-C2	-2.29	106.02	111.40
3	B	503	WRI	C29-C28-C27	-2.25	109.86	114.27
3	A	503	WRI	C09-C08-C21	-2.23	116.05	121.05
3	A	502	WRI	C06-C05-C04	-2.22	119.30	123.61
3	C	502	WRI	C22-C21-C08	-2.22	117.43	121.25
2	A	501	HEM	CAA-CBA-CGA	-2.20	107.83	113.67
2	B	501	HEM	CAD-CBD-CGD	-2.19	107.86	113.67
2	D	501	HEM	C4A-NA-C1A	2.15	109.33	105.82
2	C	501	HEM	CHA-C4D-ND	2.13	127.01	124.37
2	B	501	HEM	C2A-C1A-NA	-2.13	107.79	110.15
2	D	501	HEM	C1B-NB-C4B	2.13	107.72	105.21
3	C	510	WRI	C03-C04-C05	2.12	119.98	117.84
2	D	501	HEM	C4C-NC-C1C	2.10	109.25	105.82
3	A	502	WRI	C22-C21-C08	-2.09	117.65	121.25
3	C	502	WRI	C28-C29-C30	-2.09	109.58	115.77
2	A	501	HEM	C2A-C1A-NA	-2.09	107.84	110.15
2	D	501	HEM	CHA-C4D-ND	2.08	126.95	124.37
3	A	503	WRI	C26-C21-C08	2.08	124.33	120.84
3	B	502	WRI	C09-C10-N01	2.07	121.81	118.78
2	C	501	HEM	C4D-C3D-C2D	2.07	109.90	106.89
2	D	501	HEM	C2D-C1D-ND	-2.06	107.53	109.90
3	C	510	WRI	C28-C27-C24	-2.05	109.64	114.03
2	D	501	HEM	CBD-CAD-C3D	-2.04	106.88	112.53
2	C	501	HEM	C3A-C4A-NA	-2.04	106.87	110.14
3	B	502	WRI	C03-C02-N02	-2.03	117.43	121.81
3	A	503	WRI	C28-C29-C30	-2.03	109.75	115.77
3	B	503	WRI	C22-C21-C08	-2.02	117.77	121.25
3	B	503	WRI	C28-C29-C30	-2.02	109.79	115.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	503	WRI	C29-C28-C27	-2.01	110.32	114.27
2	B	501	HEM	C4C-NC-C1C	2.01	109.10	105.82
2	C	501	HEM	C2D-C1D-ND	-2.01	107.58	109.90
2	D	501	HEM	CAD-C3D-C2D	-2.01	124.11	127.87

There are no chirality outliers.

All (82) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	504	BTB	O1-C1-C2-C3
4	A	504	BTB	O1-C1-C2-C4
4	A	504	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
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	505	BTB	C1-C2-C4-O4
4	B	505	BTB	N-C2-C4-O4
4	C	504	BTB	C1-C2-C3-O3
4	C	504	BTB	C4-C2-C3-O3
4	C	504	BTB	N-C2-C3-O3
4	C	504	BTB	N-C2-C4-O4
4	C	504	BTB	C1-C2-N-C5
4	C	504	BTB	C1-C2-N-C7
4	C	504	BTB	C3-C2-N-C5
4	C	504	BTB	C3-C2-N-C7
4	C	504	BTB	C4-C2-N-C5
4	C	504	BTB	C4-C2-N-C7
4	C	505	BTB	O1-C1-C2-C3
4	C	505	BTB	O1-C1-C2-C4
4	C	505	BTB	O1-C1-C2-N
4	C	505	BTB	C1-C2-C4-O4
4	C	505	BTB	C3-C2-C4-O4
4	C	505	BTB	N-C2-C4-O4
4	D	504	BTB	O1-C1-C2-C3
4	D	504	BTB	C1-C2-C3-O3
4	D	504	BTB	C4-C2-C3-O3
4	D	504	BTB	N-C2-C3-O3
4	D	504	BTB	C1-C2-N-C5
4	D	504	BTB	C1-C2-N-C7

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Mol	Chain	Res	Type	Atoms
4	D	504	BTB	C3-C2-N-C5
4	D	504	BTB	C3-C2-N-C7
4	D	504	BTB	C4-C2-N-C5
4	D	504	BTB	C4-C2-N-C7
4	D	504	BTB	C8-C7-N-C5
5	A	506	GOL	O1-C1-C2-C3
5	A	506	GOL	C1-C2-C3-O3
5	C	506	GOL	C1-C2-C3-O3
5	D	505	GOL	O1-C1-C2-C3
3	B	503	WRI	C07-C08-C21-C22
3	C	510	WRI	C07-C08-C21-C22
3	C	510	WRI	C09-C08-C21-C26
3	C	510	WRI	C07-C08-C21-C26
3	C	510	WRI	C09-C08-C21-C22
3	B	503	WRI	C07-C08-C21-C26
3	B	503	WRI	C09-C08-C21-C26
3	B	503	WRI	C09-C08-C21-C22
3	A	503	WRI	C07-C08-C21-C26
3	A	503	WRI	C07-C08-C21-C22
3	A	503	WRI	C09-C08-C21-C26
3	A	503	WRI	C09-C08-C21-C22
5	C	506	GOL	O1-C1-C2-C3
5	A	506	GOL	O1-C1-C2-O2
5	A	506	GOL	O2-C2-C3-O3
5	C	506	GOL	O2-C2-C3-O3
2	B	501	HEM	C4B-C3B-CAB-CBB
5	C	506	GOL	O1-C1-C2-O2
5	D	505	GOL	O1-C1-C2-O2
4	D	504	BTB	C3-C2-C4-O4
3	C	503	WRI	C07-C08-C21-C26
3	C	503	WRI	C07-C08-C21-C22
4	A	504	BTB	N-C7-C8-O8
3	C	503	WRI	C09-C08-C21-C22
4	D	503	BTB	N-C7-C8-O8
2	A	501	HEM	C4B-C3B-CAB-CBB
2	D	501	HEM	C4B-C3B-CAB-CBB
2	B	501	HEM	C2A-CAA-CBA-CGA
4	C	505	BTB	C1-C2-N-C5
4	C	505	BTB	C3-C2-N-C7
4	C	505	BTB	C4-C2-N-C5
4	C	505	BTB	C4-C2-N-C7
3	C	503	WRI	C09-C08-C21-C26

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Mol	Chain	Res	Type	Atoms
4	B	505	BTB	C3-C2-C4-O4
2	C	501	HEM	C2B-C3B-CAB-CBB
4	C	505	BTB	N-C7-C8-O8
2	C	501	HEM	C4B-C3B-CAB-CBB
4	C	505	BTB	C1-C2-N-C7
4	C	505	BTB	C3-C2-N-C5
4	D	504	BTB	O1-C1-C2-N

All (1) ring outliers are listed below:

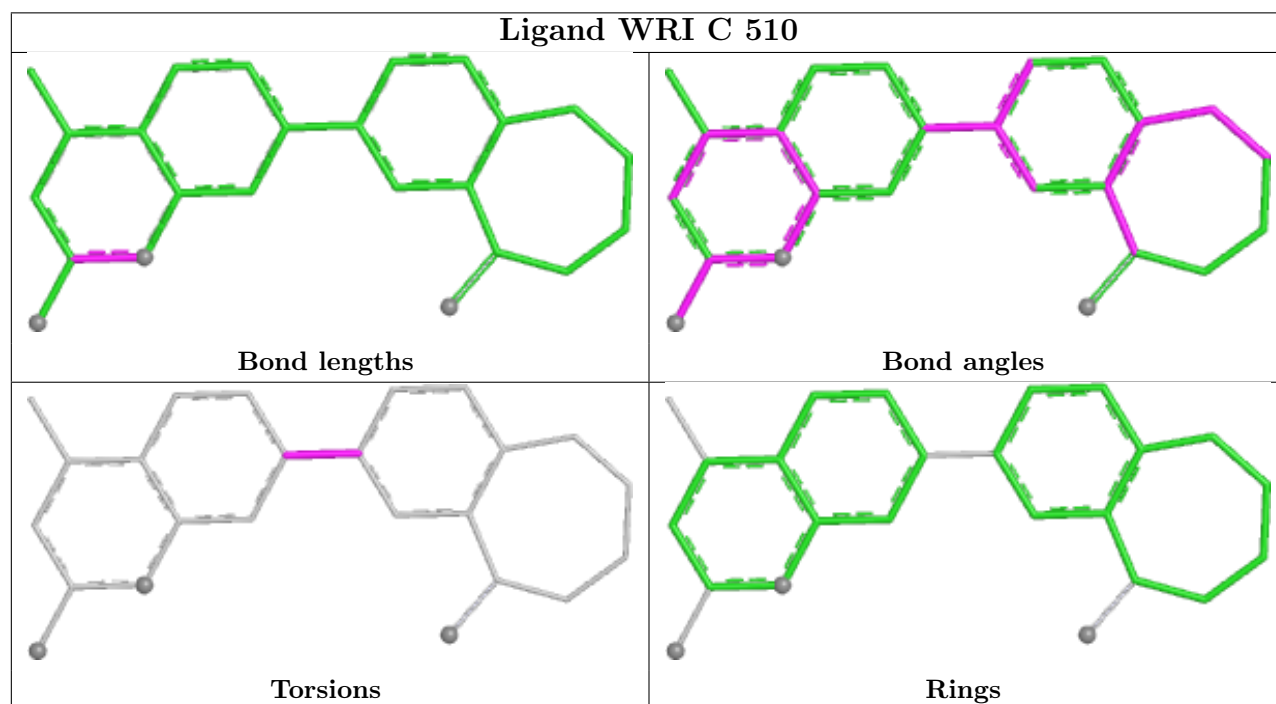
Mol	Chain	Res	Type	Atoms
3	A	503	WRI	C24-C25-C27-C28-C29-C30-C31

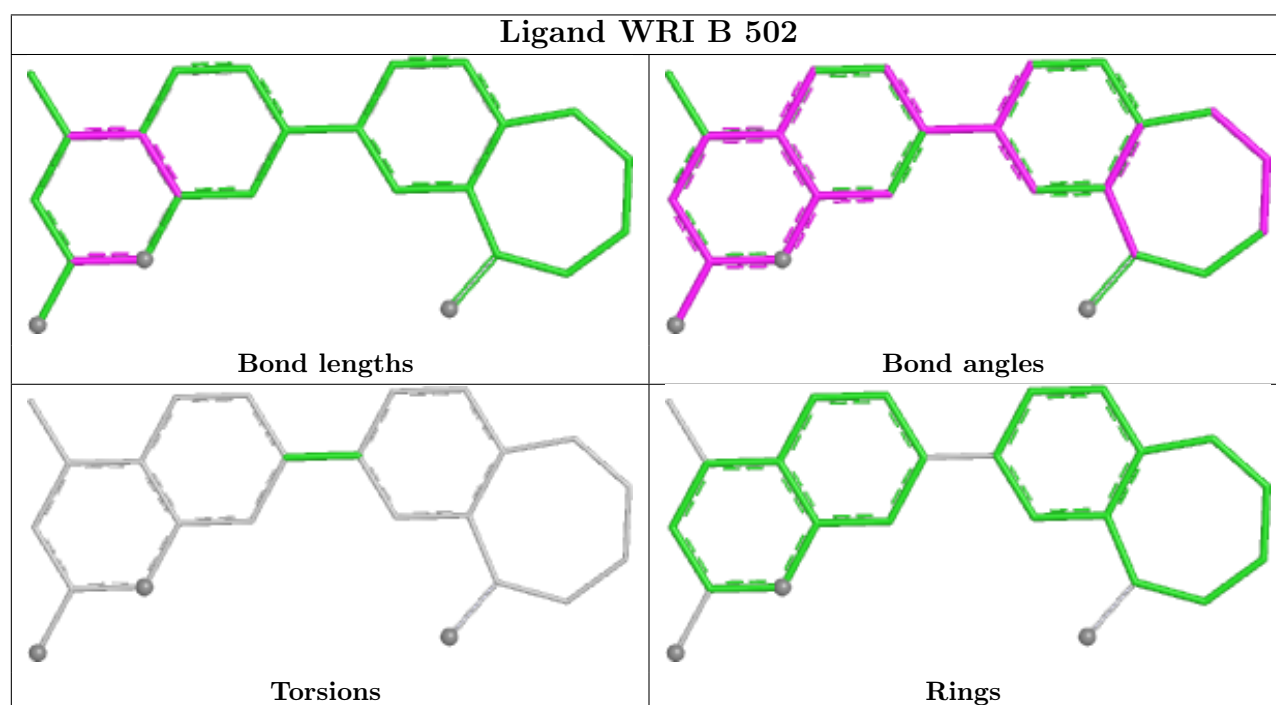
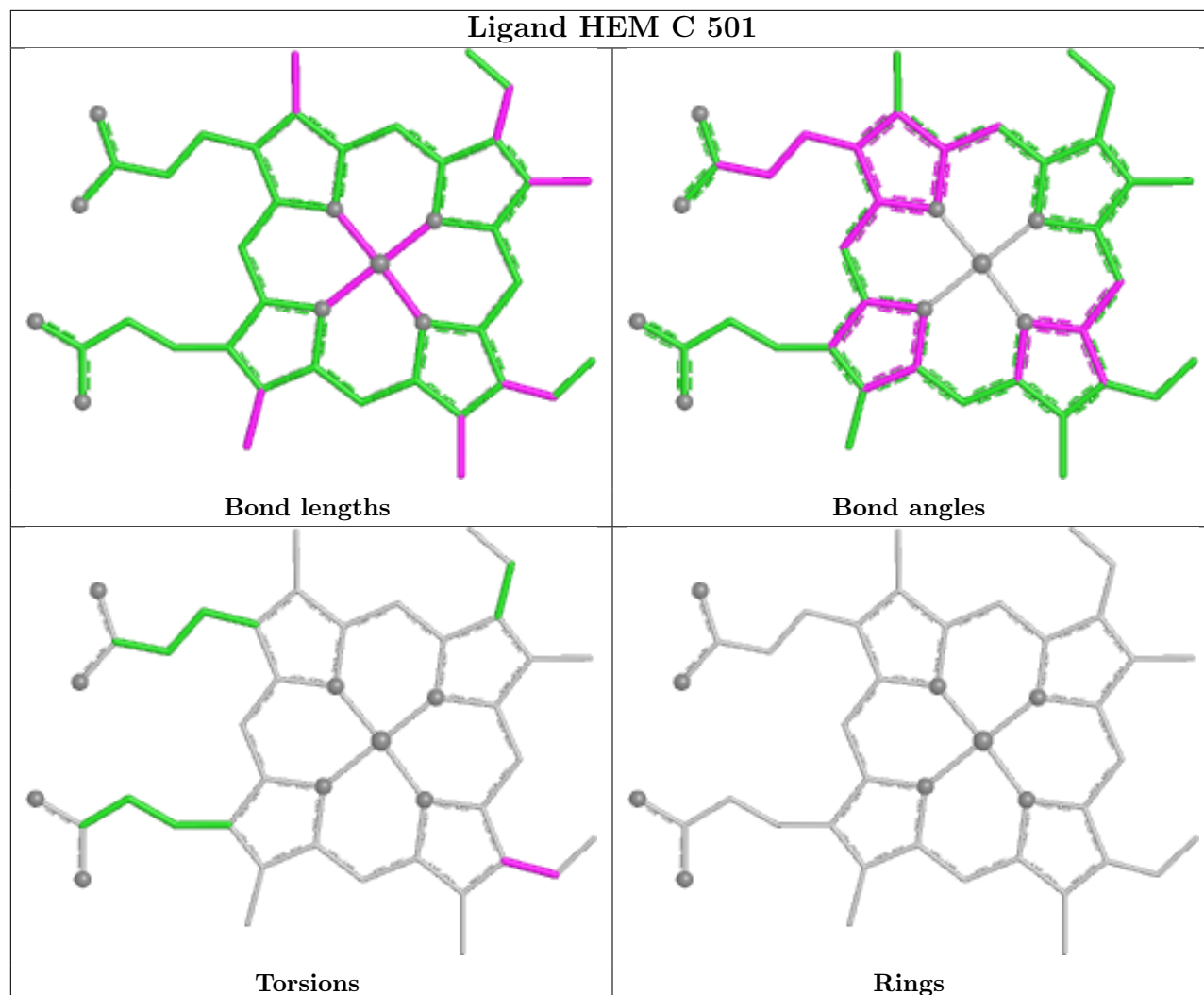
16 monomers are involved in 42 short contacts:

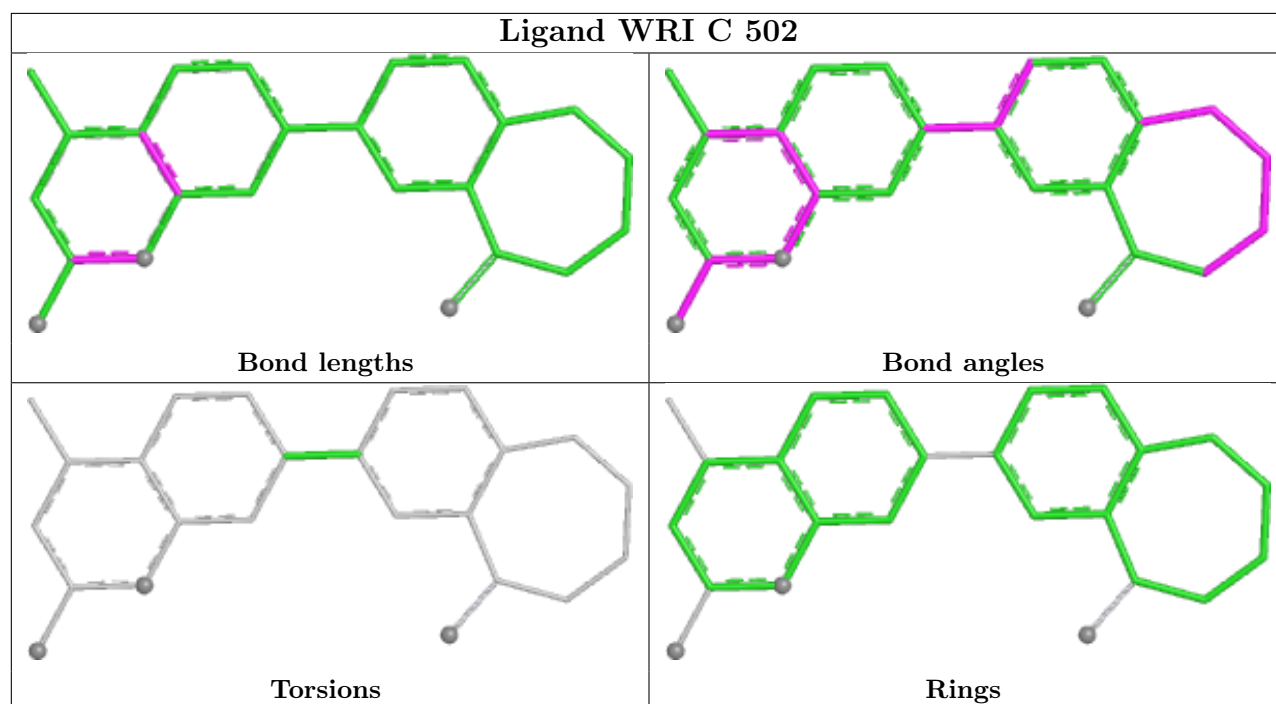
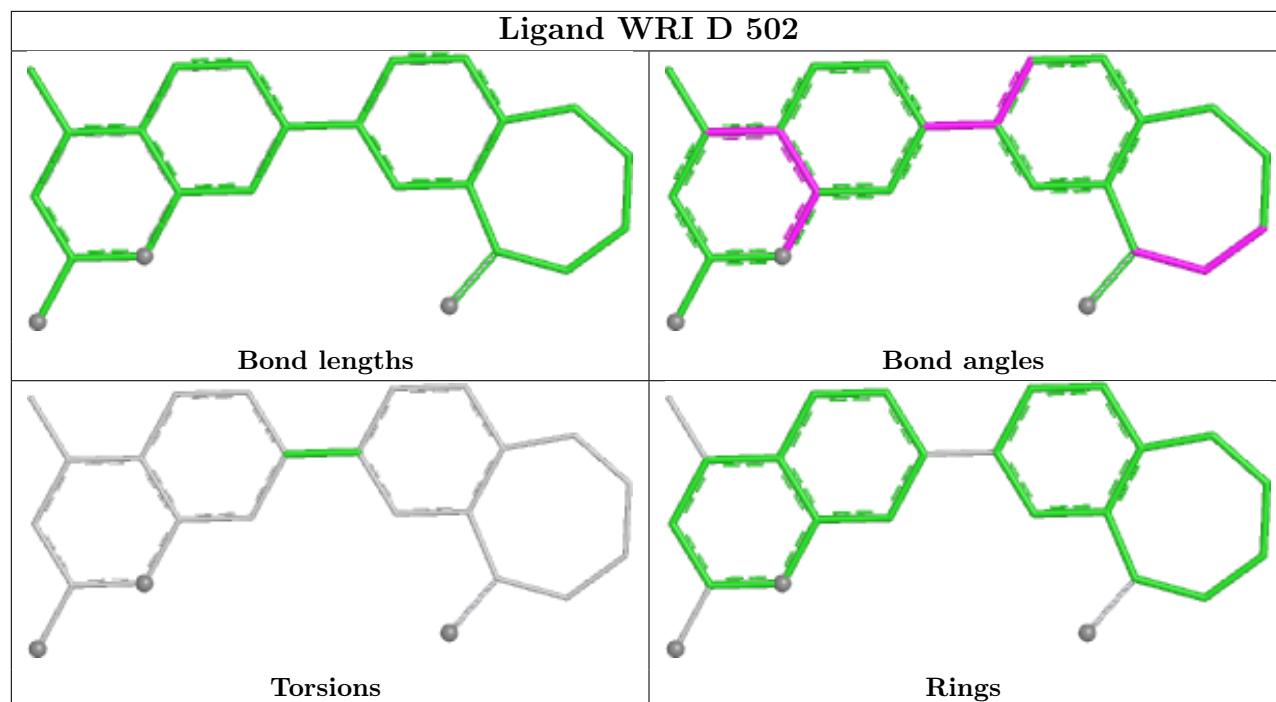
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	505	BTB	2	0
4	C	504	BTB	4	0
2	C	501	HEM	3	0
3	B	502	WRI	1	0
4	D	503	BTB	2	0
3	D	502	WRI	2	0
4	B	505	BTB	5	0
3	C	502	WRI	1	0
3	A	502	WRI	3	0
4	C	505	BTB	3	0
5	D	505	GOL	1	0
2	B	501	HEM	3	0
4	D	504	BTB	3	0
2	A	501	HEM	4	0
2	D	501	HEM	2	0
4	A	504	BTB	4	0

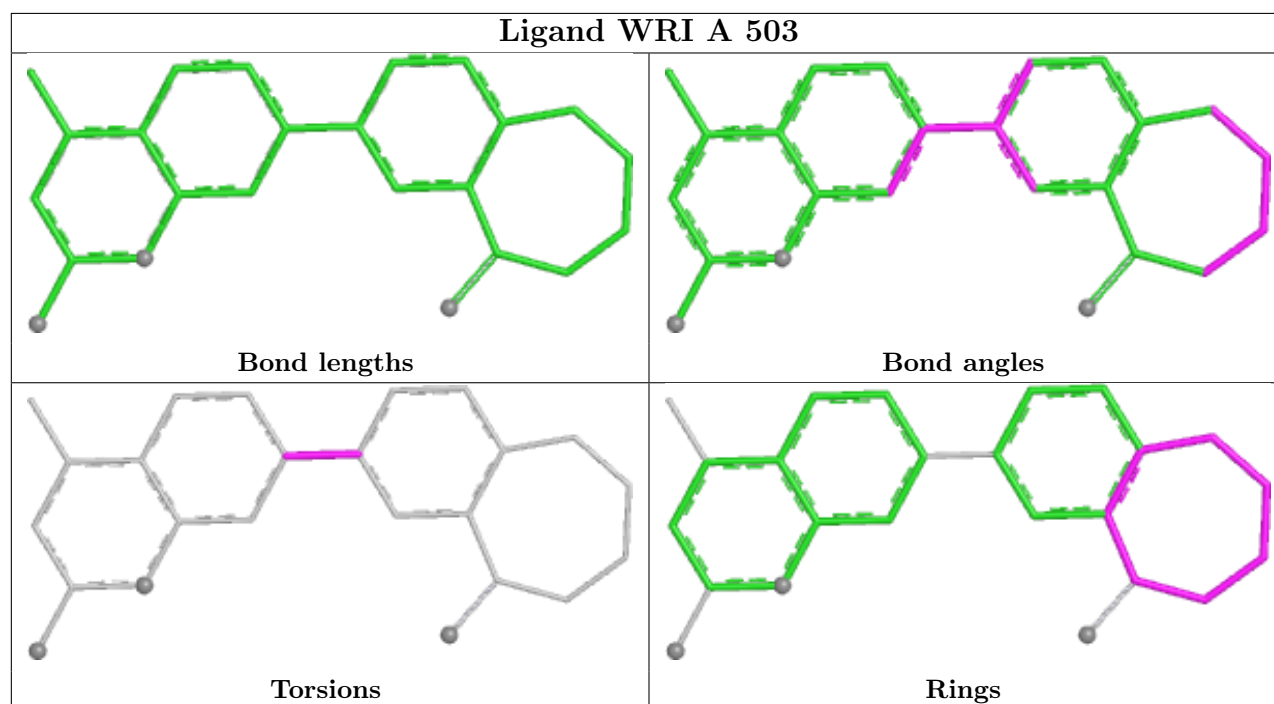
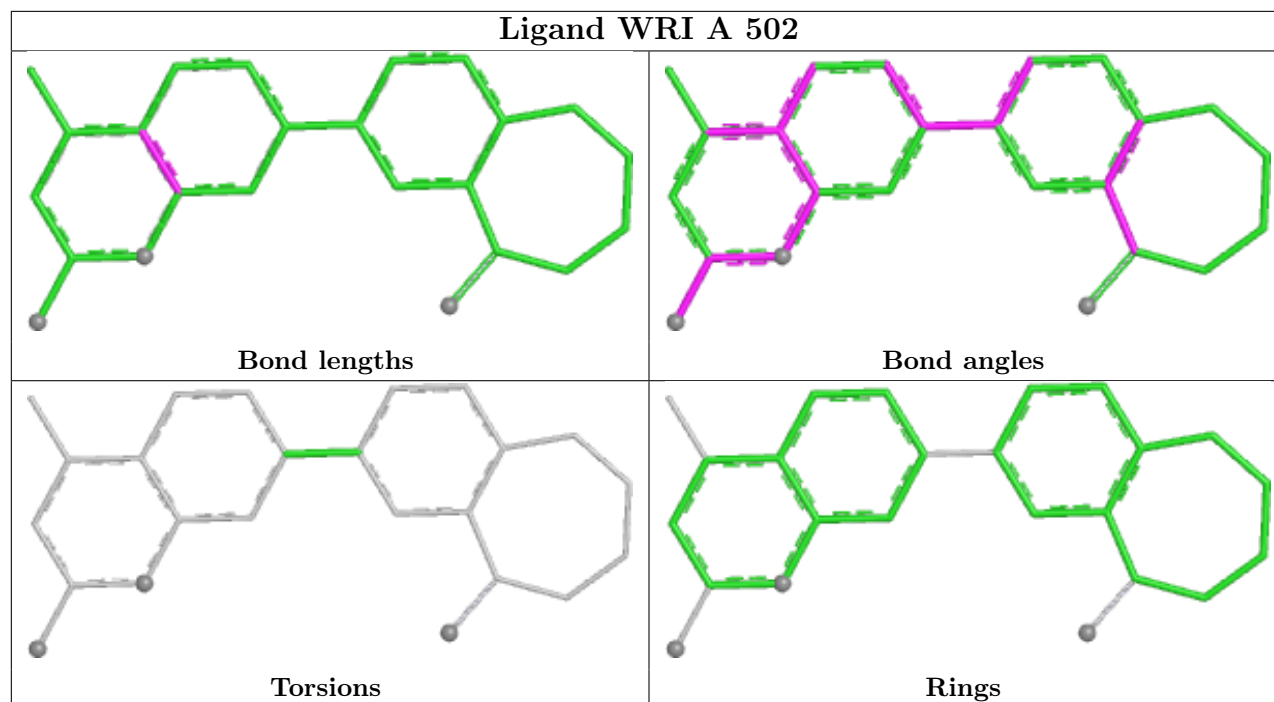
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

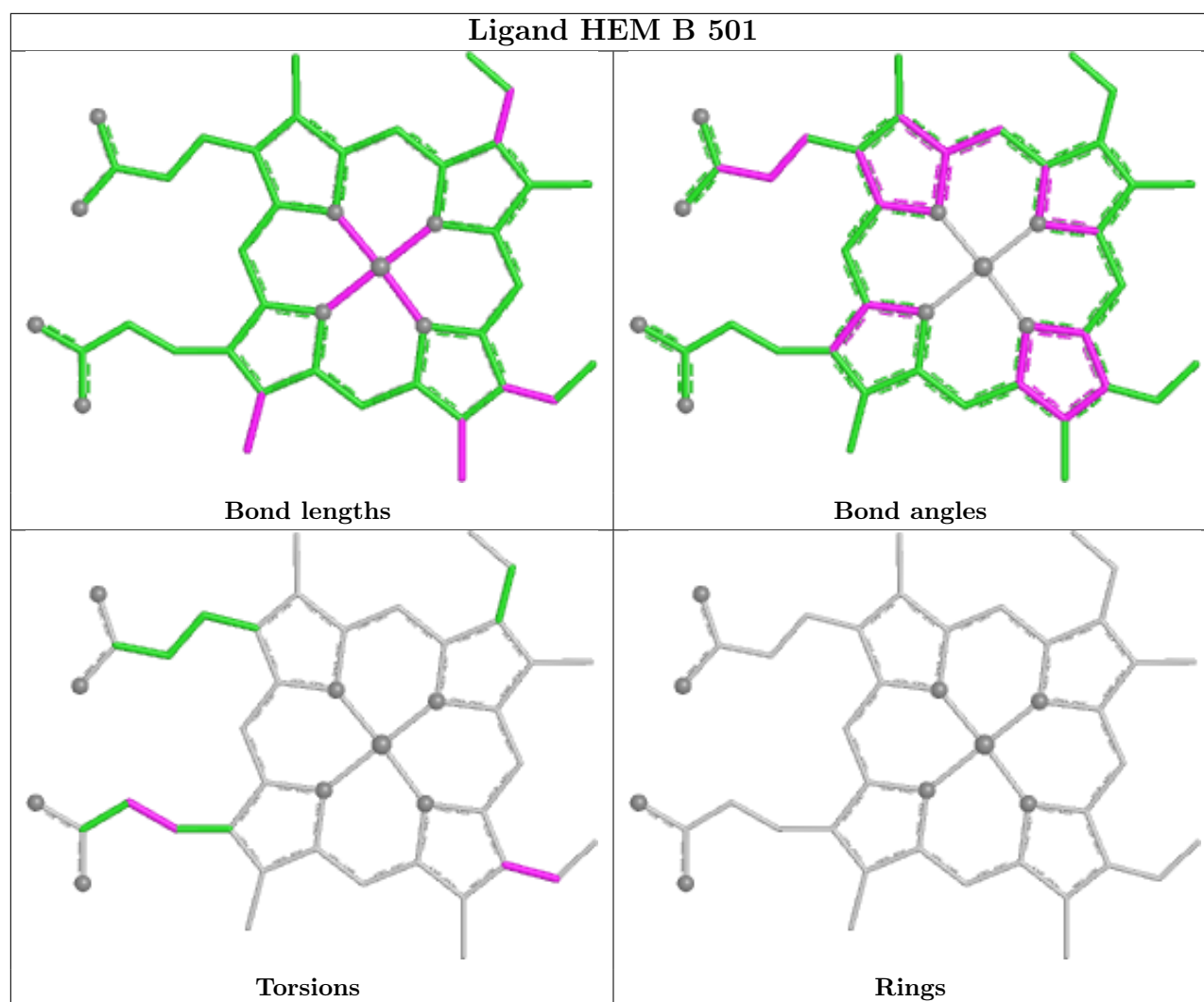
The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

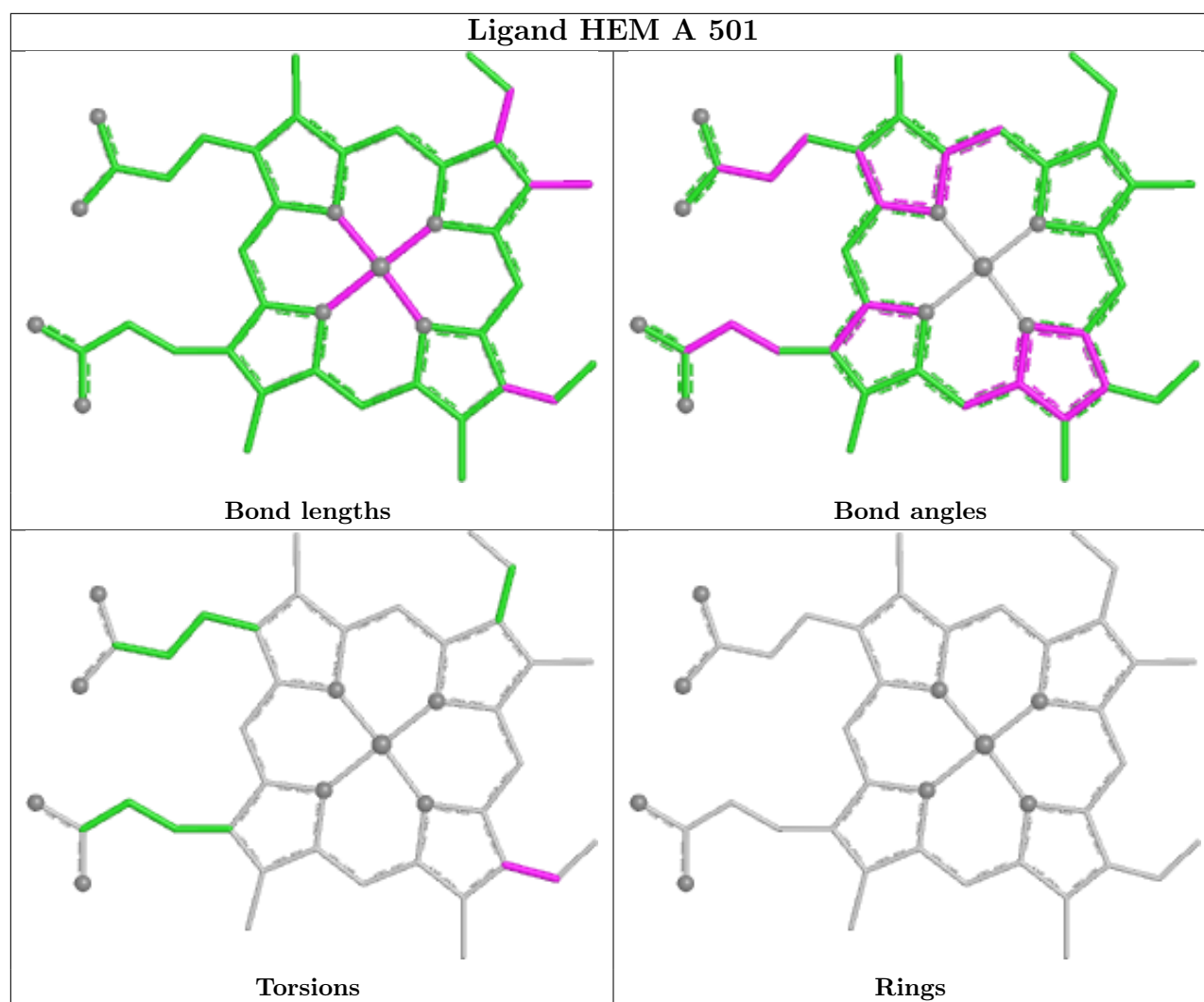


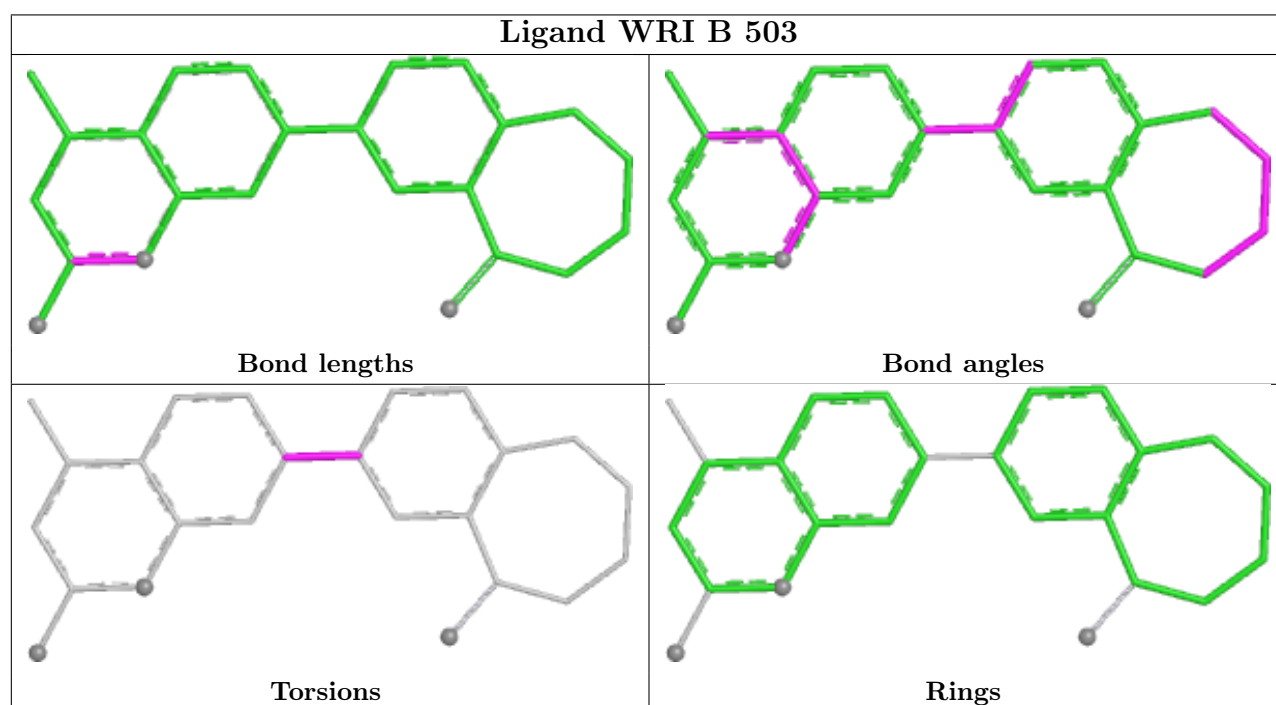
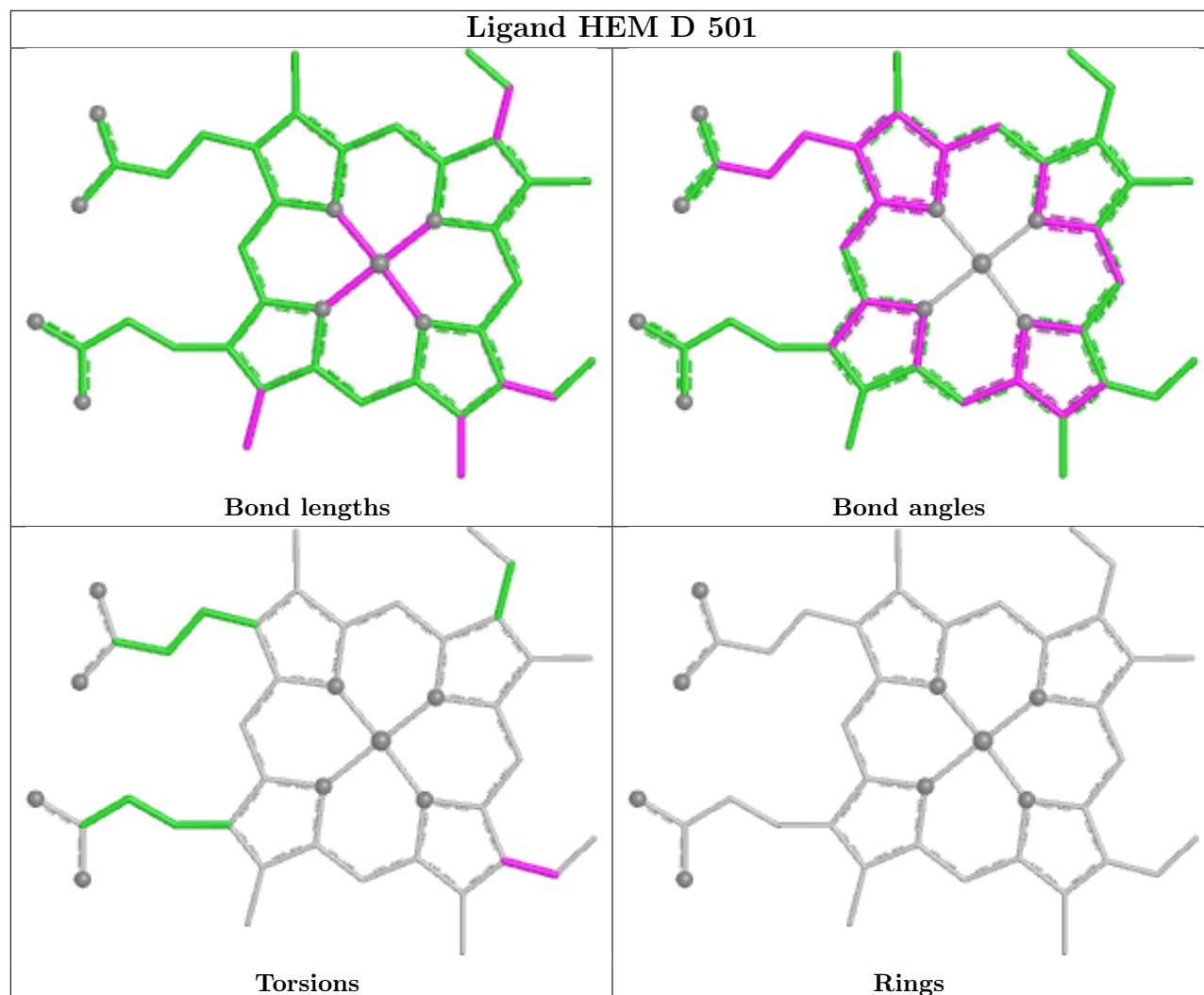


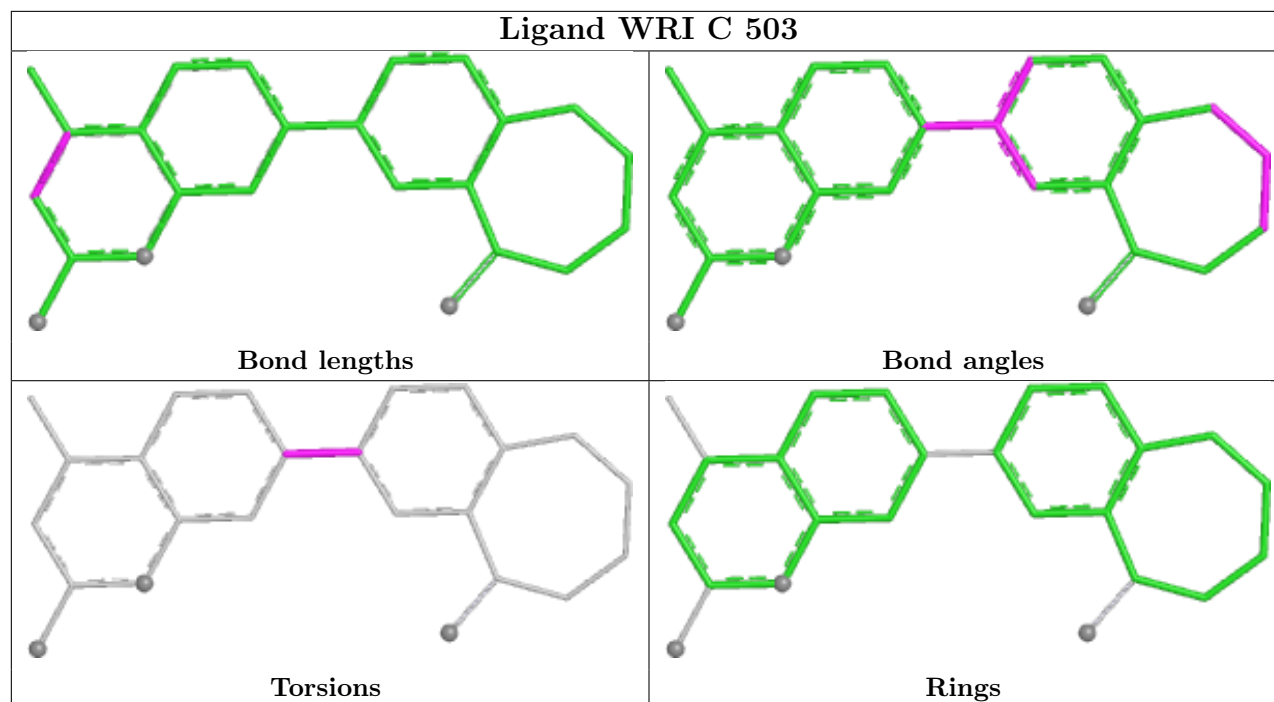












5.7 Other polymers [i](#)

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	401/440 (91%)	0.27	4 (0%) 79 81	30, 71, 120, 154	2 (0%)
1	B	401/440 (91%)	-0.07	2 (0%) 87 89	26, 50, 90, 135	3 (0%)
1	C	402/440 (91%)	0.03	3 (0%) 84 85	33, 61, 101, 130	2 (0%)
1	D	401/440 (91%)	-0.12	2 (0%) 87 89	29, 50, 84, 141	2 (0%)
All	All	1605/1760 (91%)	0.03	11 (0%) 84 85	26, 57, 106, 154	9 (0%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	153	VAL	3.7
1	A	120	PRO	3.3
1	D	119	ALA	2.9
1	B	92	GLY	2.5
1	A	280	THR	2.4
1	C	106	PRO	2.3
1	C	468	PHE	2.2
1	B	257	GLN	2.1
1	D	278	GLY	2.1
1	A	124	LEU	2.1
1	C	160	THR	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](#)

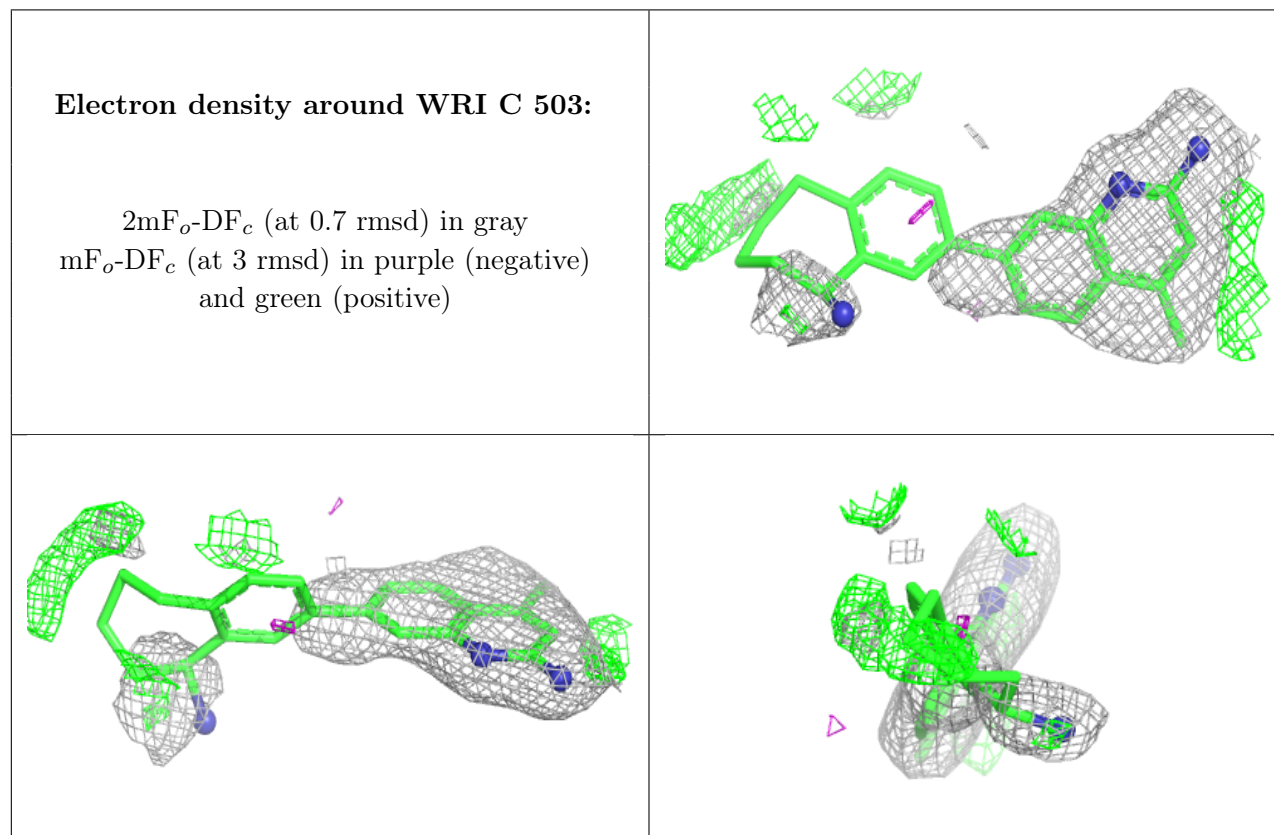
There are no oligosaccharides in this entry.

6.4 Ligands

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.

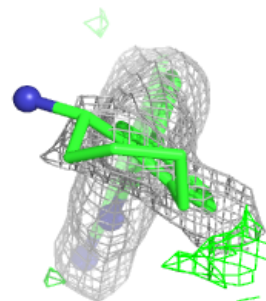
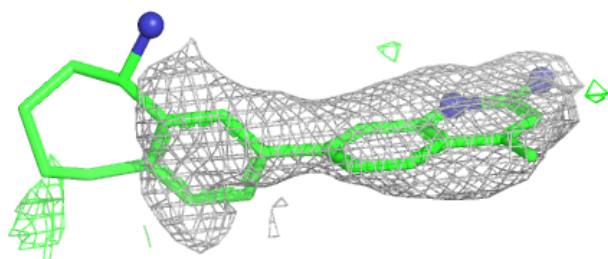
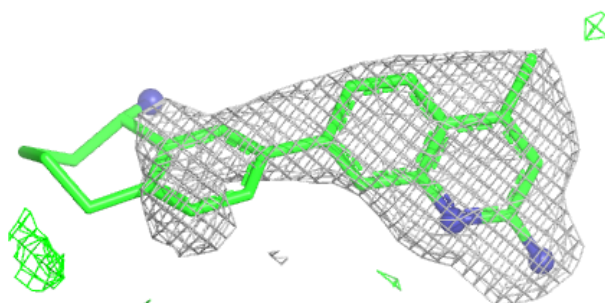
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	BTB	D	504	14/14	0.72	0.14	76,83,91,93	0
4	BTB	C	504	14/14	0.74	0.15	56,79,88,90	0
4	BTB	B	505	14/14	0.75	0.13	55,69,79,80	0
5	GOL	D	505	6/6	0.82	0.09	80,81,90,92	0
4	BTB	A	505	14/14	0.83	0.11	103,119,123,125	0
4	BTB	C	505	14/14	0.83	0.13	81,111,118,119	0
5	GOL	A	506	6/6	0.84	0.09	64,77,80,84	0
3	WRI	C	503	24/24	0.85	0.19	45,94,129,131	0
3	WRI	A	503	24/24	0.85	0.18	53,96,136,139	0
3	WRI	B	503	24/24	0.86	0.17	40,85,128,133	0
4	BTB	A	504	14/14	0.87	0.11	81,91,100,100	0
3	WRI	A	502	24/24	0.88	0.15	64,77,99,107	0
3	WRI	C	510	24/24	0.89	0.15	38,78,126,126	0
3	WRI	C	502	24/24	0.89	0.12	53,67,85,87	0
4	BTB	D	503	14/14	0.89	0.12	42,72,84,87	0
3	WRI	D	502	24/24	0.90	0.13	36,55,93,99	0
5	GOL	C	506	6/6	0.91	0.10	58,65,83,88	0
3	WRI	B	502	24/24	0.91	0.13	41,53,97,104	0
4	BTB	B	504	14/14	0.92	0.10	46,64,79,82	0
6	CL	A	507	1/1	0.94	0.12	70,70,70,70	0
6	CL	D	506	1/1	0.96	0.10	53,53,53,53	0
7	ZN	C	511	1/1	0.96	0.10	68,68,68,68	0
6	CL	B	506	1/1	0.97	0.06	55,55,55,55	0
2	HEM	A	501	43/43	0.97	0.09	44,65,86,89	0
7	ZN	A	509	1/1	0.97	0.10	65,65,65,65	0
7	ZN	B	508	1/1	0.97	0.07	64,64,64,64	0
7	ZN	C	509	1/1	0.97	0.10	61,61,61,61	0
2	HEM	C	501	43/43	0.97	0.08	36,49,86,98	0
6	CL	C	507	1/1	0.98	0.05	60,60,60,60	0
2	HEM	D	501	43/43	0.98	0.07	26,39,77,90	0
2	HEM	B	501	43/43	0.98	0.07	24,41,70,74	0
8	CA	A	510	1/1	0.98	0.12	48,48,48,48	0
7	ZN	A	508	1/1	0.99	0.03	52,52,52,52	0
8	CA	C	512	1/1	0.99	0.05	56,56,56,56	0
9	GD	B	507	1/1	0.99	0.03	55,55,55,55	0
9	GD	D	507	1/1	0.99	0.03	54,54,54,54	0
7	ZN	C	508	1/1	1.00	0.02	45,45,45,45	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.

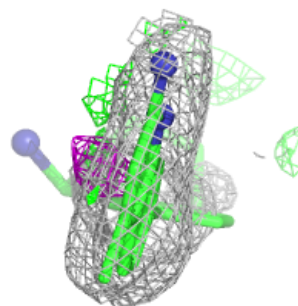
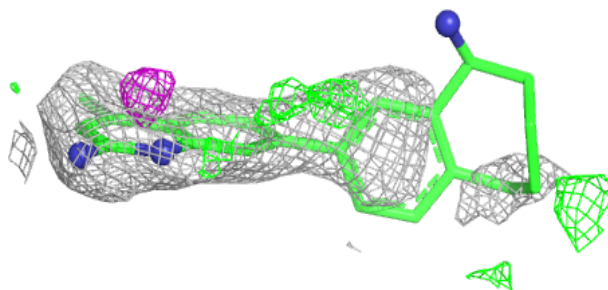
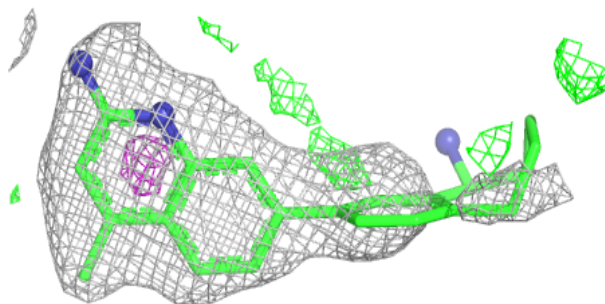


Electron density around WRI A 503:

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

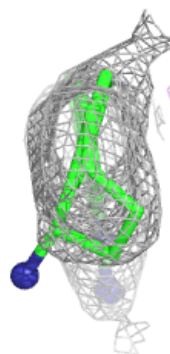
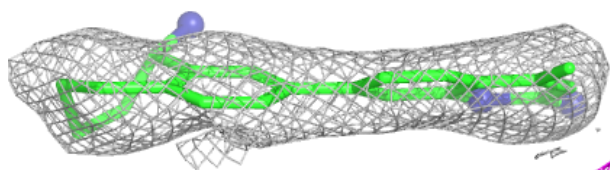
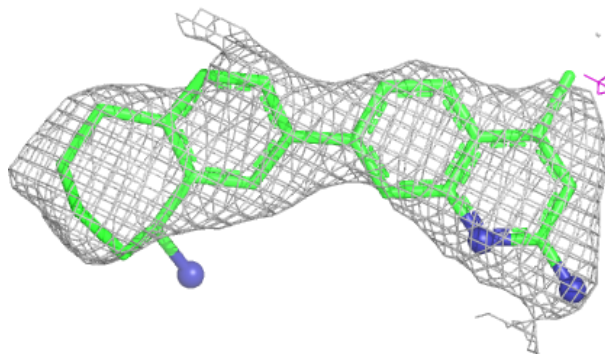
**Electron density around WRI B 503:**

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

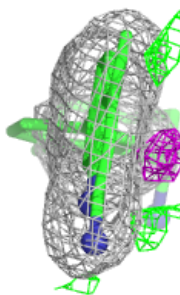
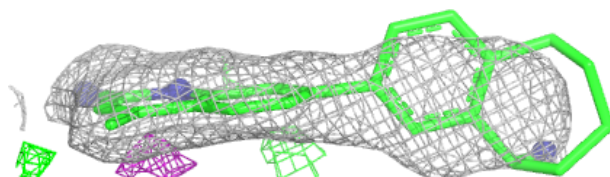
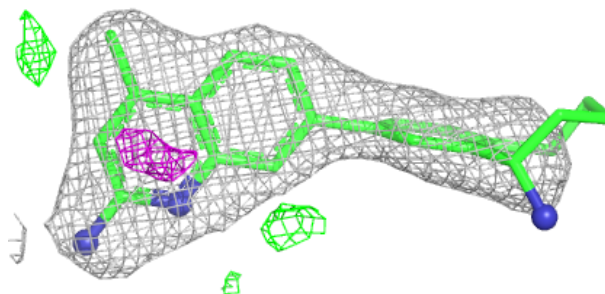


Electron density around WRI A 502:

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

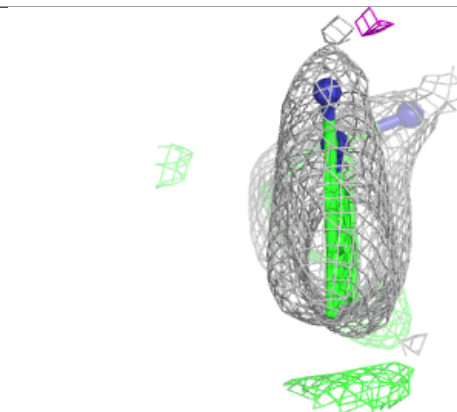
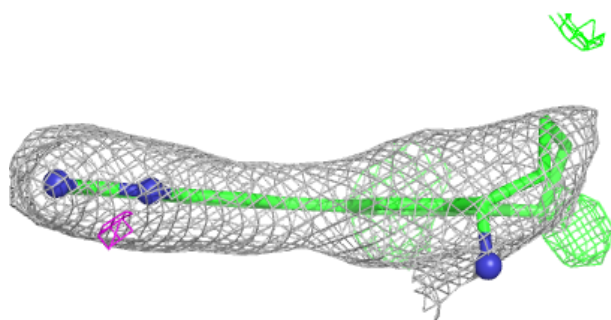
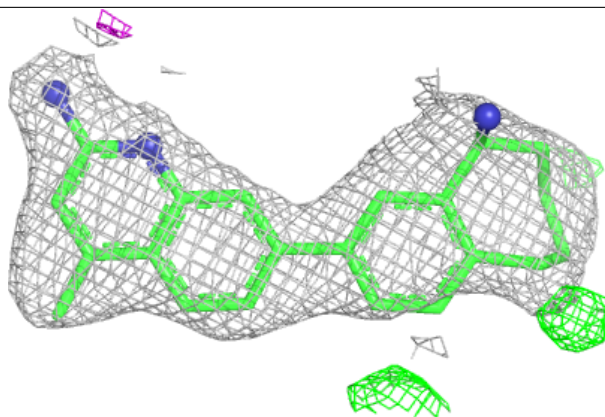
**Electron density around WRI C 510:**

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

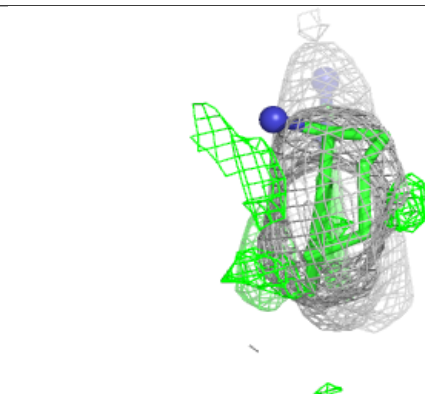
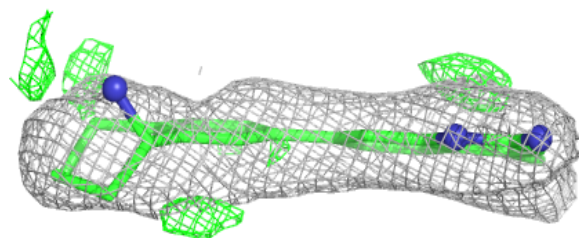
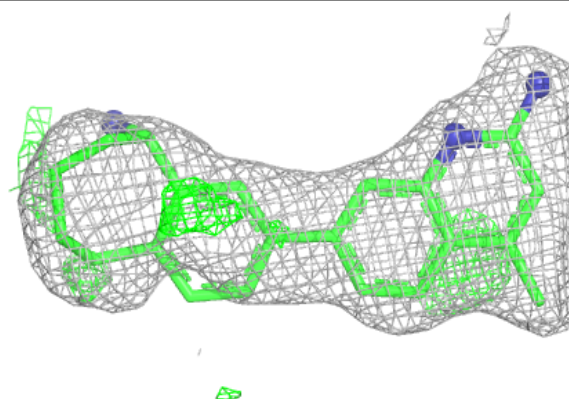


Electron density around WRI C 502:

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

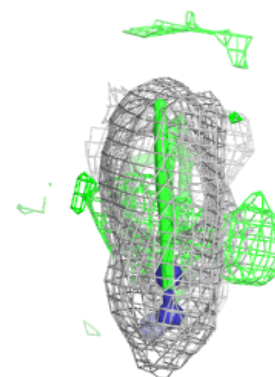
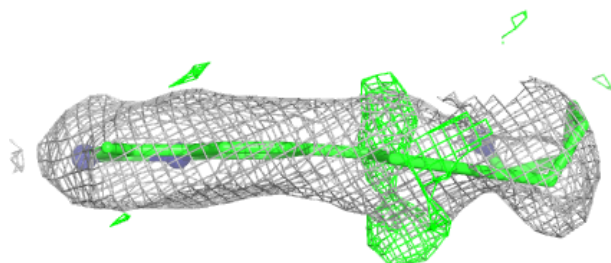
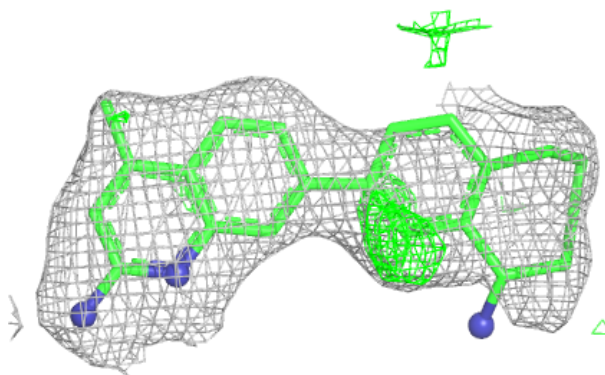
**Electron density around WRI D 502:**

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

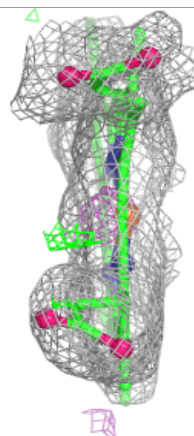
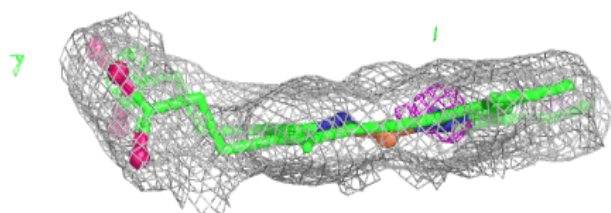
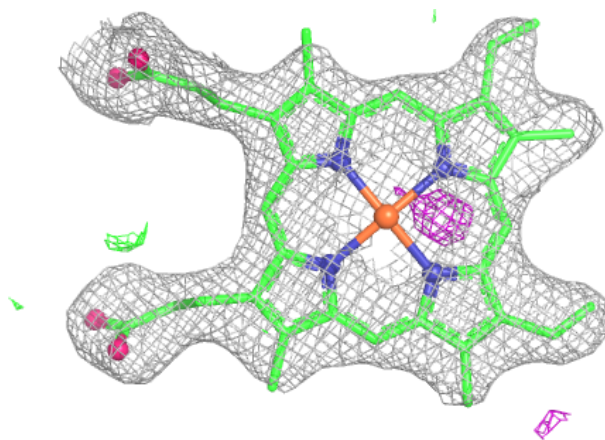


Electron density around WRI B 502:

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

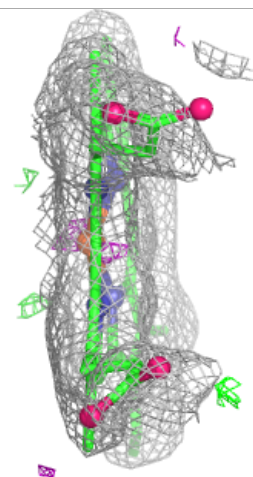
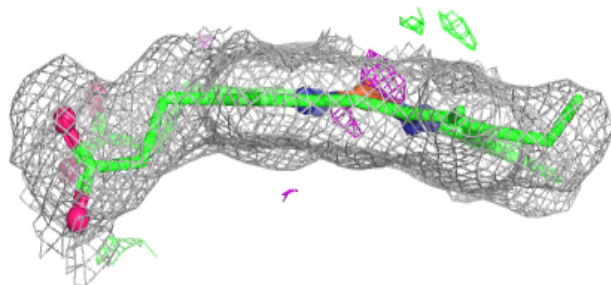
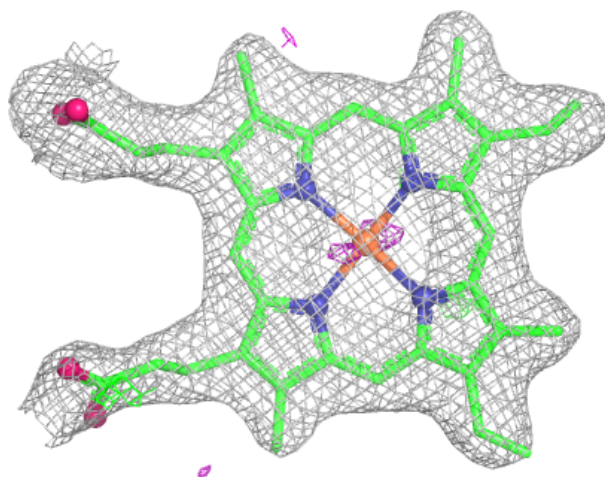
**Electron density around HEM A 501:**

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



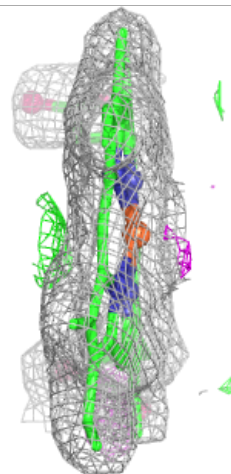
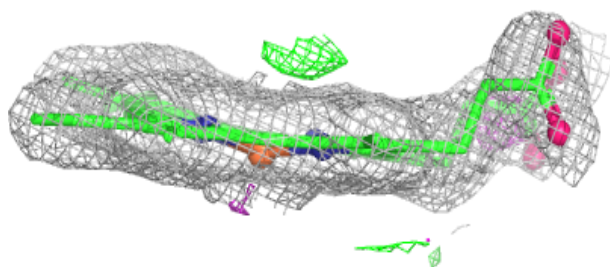
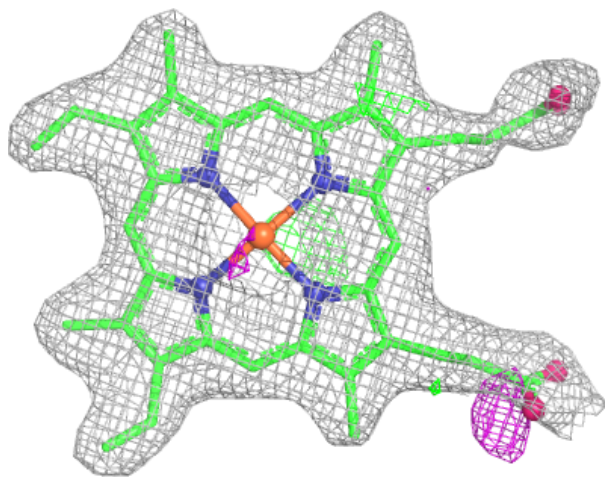
Electron density around HEM C 501:

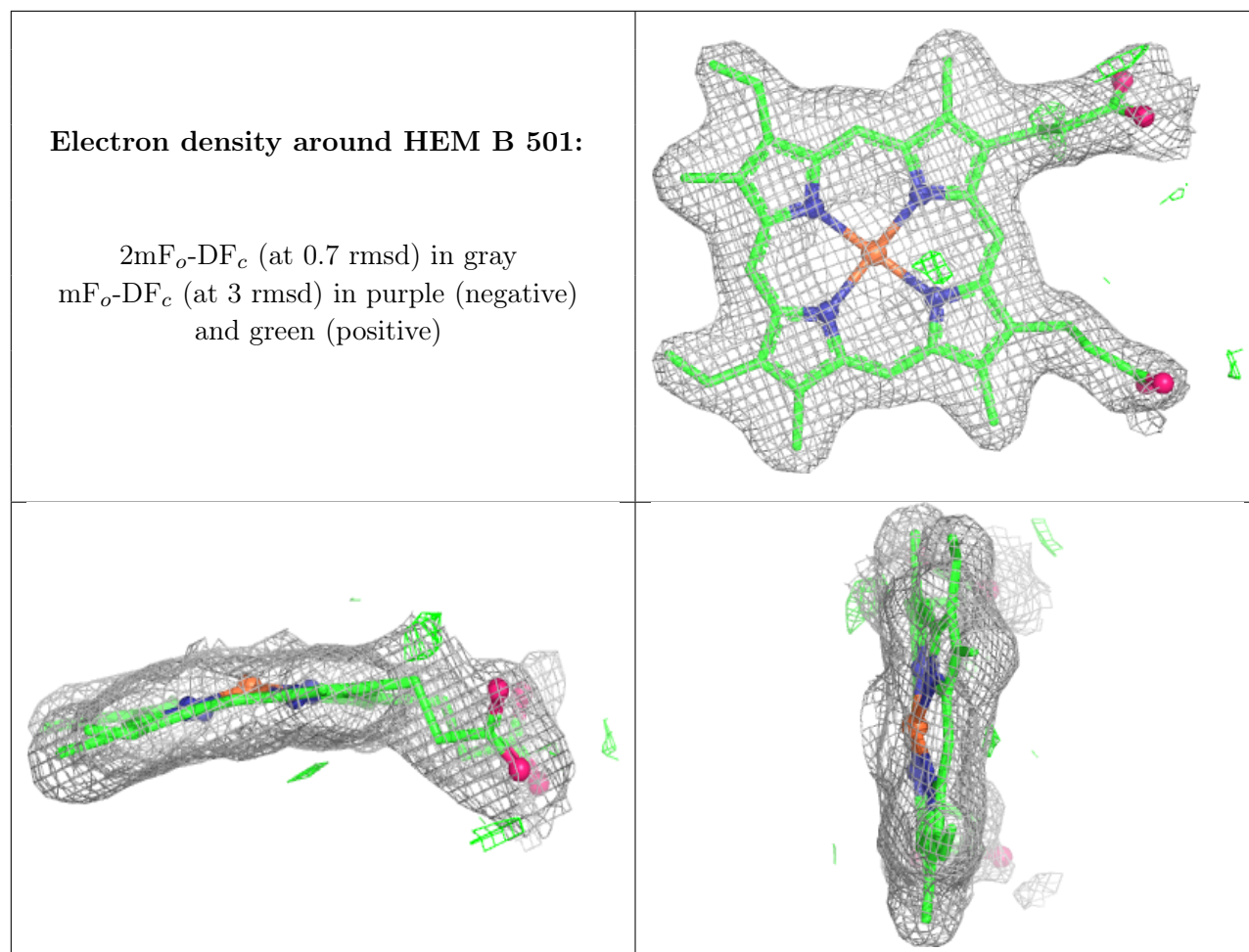
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM 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.