



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:23 AM UTC

PDB ID : 7CVR / pdb_00007cvr
Title : Structure of the CYP102A1 Haem Domain with N-Carboxybenzyl-L-Prolyl-L-Phenylalanine in complex with (S)-1-Tetralylamine
Authors : Stanfield, J.K.; Sugimoto, H.; Shoji, O.
Deposited on : 2020-08-26
Resolution : 1.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

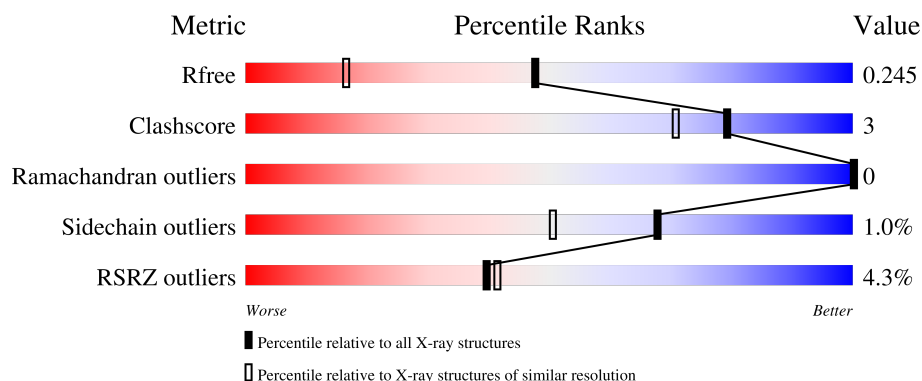
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	4673 (1.60-1.60)
Clashscore	190562	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)
RSRZ outliers	180081	4672 (1.60-1.60)

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	456	5% 95% ..
1	B	456	3% 95% ..

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 8260 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 Bifunctional cytochrome P450/NADPH-P450 reductase.

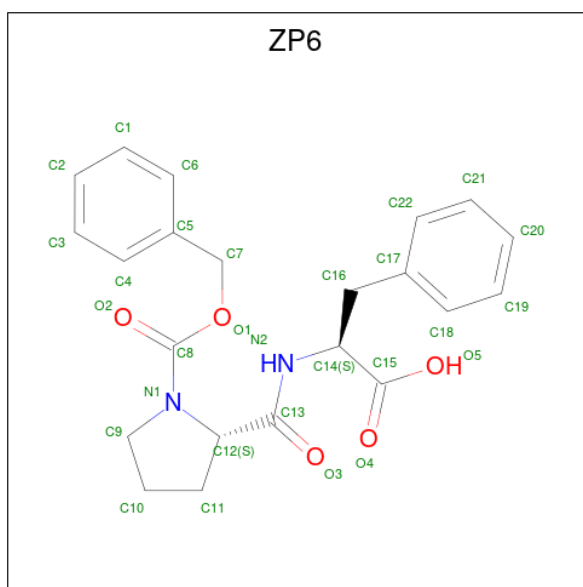
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	453	Total	C	N	O	S	0	6	0
			3702	2363	629	692	18			
1	B	453	Total	C	N	O	S	0	7	0
			3711	2370	631	693	17			

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



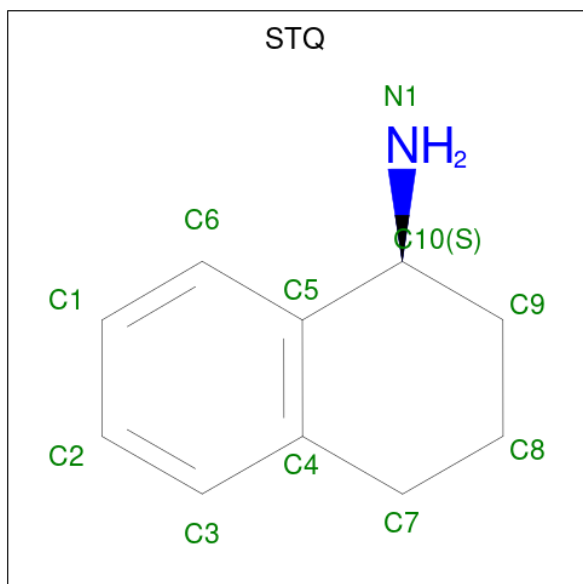
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	1
			86	68	2	8	8		
2	B	1	Total	C	Fe	N	O	0	1
			86	68	2	8	8		

- Molecule 3 is (2S)-3-phenyl-2-[[[(2S)-1-phenylmethoxycarbonylpyrrolidin-2-yl]carbonylamino]propanoic acid (CCD ID: ZP6) (formula: $C_{22}H_{24}N_2O_5$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			29	22	2	5		
3	B	1	Total	C	N	O	0	0
			29	22	2	5		

- Molecule 4 is (1 {S})-1,2,3,4-tetrahydronaphthalen-1-amine (CCD ID: STQ) (formula: C₁₀H₁₃N) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N		0	0
			11	10	1			

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	N	0	0
			11	10	1		

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

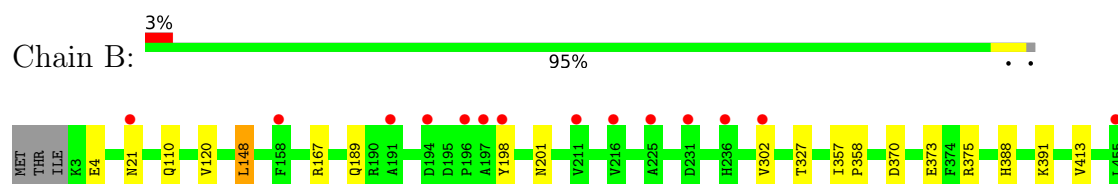


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	266	Total	O	0	0
			266	266		
6	B	299	Total	O	0	0
			299	299		

- Molecule 1: Bifunctional cytochrome P450/NADPH-P450 reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.92Å 127.65Å 147.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.33 – 1.60 48.33 – 1.60	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.33-1.60) 100.0 (48.33-1.60)	Depositor EDS
R_{merge}	0.24	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.204 , 0.238 0.211 , 0.245	Depositor DCC
R_{free} test set	7397 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	21.6	Xtriage
Anisotropy	0.287	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 30.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8260	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 40.05 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.8994e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: ZP6, HEM, STQ, GOL

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	1.03	1/3790 (0.0%)	1.24	3/5121 (0.1%)
1	B	1.02	0/3799	1.21	0/5135
All	All	1.02	1/7589 (0.0%)	1.23	3/10256 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	385	ILE	N-CA	5.06	1.50	1.46

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	357	ILE	N-CA-CB	5.97	114.26	110.50
1	A	250	ASP	CA-CB-CG	5.22	117.82	112.60
1	A	288	GLN	CB-CG-CD	5.20	121.44	112.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3702	0	3670	11	0
1	B	3711	0	3686	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	86	0	60	11	0
2	B	86	0	60	10	0
3	A	29	0	0	0	0
3	B	29	0	0	0	0
4	A	11	0	0	0	0
4	B	11	0	0	0	0
5	A	6	0	8	1	0
5	B	24	0	32	0	0
6	A	266	0	0	0	0
6	B	299	0	0	1	0
All	All	8260	0	7516	41	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 3.

All (41) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:501[A]:HEM:HMC1	2:A:501[A]:HEM:HBC2	1.52	0.91
1:B:327[B]:THR:HG23	2:B:501[B]:HEM:CBC	2.03	0.86
1:B:373[B]:GLU:HG2	1:B:375[B]:ARG:HG3	1.58	0.83
2:A:501[A]:HEM:HBC2	2:A:501[A]:HEM:CMC	2.13	0.77
2:A:501[B]:HEM:HBB2	2:A:501[B]:HEM:HMB2	1.69	0.74
2:A:501[A]:HEM:HBB2	2:A:501[A]:HEM:HMB1	1.73	0.70
1:A:170:PRO:HB3	5:A:504:GOL:H11	1.73	0.69
2:B:501[B]:HEM:HBB2	2:B:501[B]:HEM:HMB2	1.75	0.69
2:B:501[B]:HEM:HMC2	2:B:501[B]:HEM:HBC2	1.76	0.68
2:A:501[B]:HEM:HBB2	2:A:501[B]:HEM:CMB	2.23	0.67
1:A:21:ASN:HA	1:A:189:GLN:HE21	1.61	0.66
1:A:198:TYR:HA	1:A:201:ASN:HD22	1.62	0.65
2:B:501[A]:HEM:HMC1	2:B:501[A]:HEM:HBC2	1.79	0.65
2:A:501[B]:HEM:HMC2	2:A:501[B]:HEM:HBC2	1.79	0.64
1:B:120:VAL:HG11	1:B:302:VAL:HG13	1.79	0.63
1:B:148:LEU:HD21	1:B:413:VAL:HG21	1.80	0.62
1:A:148:LEU:HD21	1:A:413:VAL:HG21	1.83	0.60
1:B:21:ASN:OD1	1:B:189:GLN:NE2	2.35	0.59
2:B:501[B]:HEM:HBB2	2:B:501[B]:HEM:CMB	2.32	0.58
1:A:327[A]:THR:HG23	2:A:501[A]:HEM:HMB3	1.88	0.56
2:B:501[B]:HEM:HBC2	2:B:501[B]:HEM:CMC	2.36	0.56
2:A:501[A]:HEM:HMC1	2:A:501[A]:HEM:CBC	2.31	0.55
2:B:501[A]:HEM:HMB2	2:B:501[A]:HEM:HBB2	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:327[B]:THR:HG23	2:B:501[B]:HEM:HBC1	1.89	0.52
2:A:501[A]:HEM:HBB2	2:A:501[A]:HEM:CMB	2.41	0.49
2:A:501[B]:HEM:HBC2	2:A:501[B]:HEM:CMC	2.40	0.49
1:B:373[B]:GLU:CG	1:B:375[B]:ARG:HG3	2.38	0.49
2:B:501[A]:HEM:HBB2	2:B:501[A]:HEM:CMB	2.44	0.48
1:A:269:THR:HG21	1:A:409[B]:GLU:CG	2.44	0.47
2:B:501[A]:HEM:HBC2	2:B:501[A]:HEM:CMC	2.42	0.47
1:A:269:THR:HG21	1:A:409[B]:GLU:HG3	1.97	0.47
1:A:47:ARG:HG2	1:A:48:VAL:N	2.29	0.47
1:B:198:TYR:HA	1:B:201:ASN:HD22	1.80	0.47
2:A:501[B]:HEM:HMB2	2:A:501[B]:HEM:CBB	2.43	0.46
1:B:388:HIS:HA	1:B:391[B]:LYS:HD3	1.98	0.45
1:B:110:GLN:HG2	6:B:813:HOH:O	2.17	0.44
1:A:81:PHE:HB3	1:A:209:ILE:HG12	2.00	0.43
1:B:370:ASP:OD2	1:B:375[B]:ARG:NH1	2.51	0.43
1:A:269:THR:HG21	1:A:409[B]:GLU:CD	2.44	0.42
1:B:357:ILE:N	1:B:358:PRO:CD	2.83	0.42
1:A:198:TYR:HA	1:A:201:ASN:ND2	2.32	0.41

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	457/456 (100%)	442 (97%)	15 (3%)	0	100	100
1	B	458/456 (100%)	446 (97%)	12 (3%)	0	100	100
All	All	915/912 (100%)	888 (97%)	27 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	403/400 (101%)	396 (98%)	7 (2%)	53	30
1	B	404/400 (101%)	401 (99%)	3 (1%)	76	63
All	All	807/800 (101%)	797 (99%)	10 (1%)	68	43

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	148	LEU
1	A	167	ARG
1	A	168	ASP
1	A	190[A]	ARG
1	A	190[B]	ARG
1	A	327[A]	THR
1	A	327[B]	THR
1	B	4	GLU
1	B	148	LEU
1	B	167	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	GLN
1	A	169	GLN
1	A	189	GLN
1	A	201	ASN
1	A	204	GLN
1	A	206	GLN
1	A	359	GLN
1	A	387	GLN
1	B	73	GLN
1	B	189	GLN
1	B	201	ASN
1	B	204	GLN

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Mol	Chain	Res	Type
1	B	206	GLN
1	B	288	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

13 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ZP6	B	502	-	31,31,31	1.28	1 (3%)	41,41,41	1.52	5 (12%)
2	HEM	B	501[A]	4,1	50,50,50	1.53	8 (16%)	67,82,82	1.61	12 (17%)
5	GOL	B	505	-	5,5,5	0.09	0	5,5,5	0.36	0
2	HEM	B	501[B]	4,1	50,50,50	1.56	8 (16%)	67,82,82	1.53	8 (11%)
2	HEM	A	501[A]	4,1	50,50,50	1.58	11 (22%)	67,82,82	1.61	12 (17%)
2	HEM	A	501[B]	4,1	50,50,50	1.59	10 (20%)	67,82,82	1.57	10 (14%)
5	GOL	B	504	-	5,5,5	0.07	0	5,5,5	0.34	0
5	GOL	A	504	-	5,5,5	0.15	0	5,5,5	0.41	0
5	GOL	B	507	-	5,5,5	0.06	0	5,5,5	0.22	0
4	STQ	B	503	2	12,12,12	1.05	1 (8%)	15,16,16	1.17	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ZP6	A	502	-	31,31,31	1.37	1 (3%)	41,41,41	1.48	6 (14%)
5	GOL	B	506	-	5,5,5	0.16	0	5,5,5	0.35	0
4	STQ	A	503	2	12,12,12	0.78	0	15,16,16	0.83	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	ZP6	B	502	-	-	0/25/35/35	0/3/3/3
2	HEM	B	501[A]	4,1	-	3/14/54/54	-
5	GOL	B	505	-	-	4/4/4/4	-
2	HEM	B	501[B]	4,1	-	2/14/54/54	-
2	HEM	A	501[A]	4,1	-	2/14/54/54	-
2	HEM	A	501[B]	4,1	-	2/14/54/54	-
5	GOL	B	504	-	-	4/4/4/4	-
5	GOL	A	504	-	-	0/4/4/4	-
5	GOL	B	507	-	-	3/4/4/4	-
4	STQ	B	503	2	-	-	0/2/2/2
3	ZP6	A	502	-	-	0/25/35/35	0/3/3/3
5	GOL	B	506	-	-	1/4/4/4	-
4	STQ	A	503	2	-	-	0/2/2/2

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502	ZP6	O1-C8	6.22	1.46	1.34
3	B	502	ZP6	O1-C8	5.47	1.45	1.34
2	A	501[B]	HEM	FE-NB	4.77	2.09	1.94
2	B	501[A]	HEM	FE-NB	4.75	2.09	1.94
2	A	501[A]	HEM	FE-NB	4.45	2.08	1.94
2	B	501[B]	HEM	FE-NB	4.39	2.08	1.94
2	A	501[A]	HEM	FE-NC	3.95	2.08	1.95
2	A	501[A]	HEM	C1B-NB	-3.85	1.33	1.40
2	B	501[B]	HEM	FE-NC	3.71	2.07	1.95
2	B	501[A]	HEM	C1B-NB	-3.64	1.34	1.40
2	A	501[B]	HEM	FE-NC	3.40	2.06	1.95
2	B	501[A]	HEM	FE-NC	3.39	2.06	1.95
2	A	501[B]	HEM	C4B-NB	-3.29	1.32	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	503	STQ	C5-C10	-3.26	1.46	1.51
2	B	501[B]	HEM	C1B-NB	-3.08	1.34	1.40
2	B	501[B]	HEM	C4D-C3D	3.05	1.50	1.45
2	A	501[A]	HEM	C4D-C3D	3.04	1.50	1.45
2	A	501[B]	HEM	C1B-NB	-3.00	1.35	1.40
2	A	501[B]	HEM	C3C-C4C	-2.99	1.40	1.46
2	B	501[B]	HEM	FE-NA	2.77	2.04	1.95
2	A	501[B]	HEM	C1C-C2C	-2.68	1.40	1.45
2	A	501[B]	HEM	C4D-C3D	2.68	1.49	1.45
2	B	501[B]	HEM	C4D-ND	-2.63	1.35	1.40
2	B	501[A]	HEM	FE-NA	2.59	2.03	1.95
2	A	501[A]	HEM	C1D-ND	-2.45	1.34	1.38
2	A	501[A]	HEM	C1C-NC	-2.45	1.35	1.39
2	A	501[A]	HEM	C1D-C2D	2.45	1.49	1.44
2	A	501[A]	HEM	C4D-ND	-2.43	1.36	1.40
2	A	501[A]	HEM	FE-NA	2.41	2.03	1.95
2	B	501[A]	HEM	C4D-C3D	2.38	1.49	1.45
2	B	501[A]	HEM	C4B-NB	-2.36	1.34	1.38
2	B	501[A]	HEM	C1D-ND	-2.33	1.34	1.38
2	B	501[B]	HEM	C4B-NB	-2.31	1.34	1.38
2	A	501[B]	HEM	FE-NA	2.26	2.02	1.95
2	B	501[B]	HEM	C3B-C4B	2.22	1.49	1.44
2	B	501[A]	HEM	C3C-C4C	-2.18	1.42	1.46
2	A	501[A]	HEM	O1A-CGA	2.16	1.29	1.22
2	A	501[B]	HEM	C1D-ND	-2.13	1.34	1.38
2	A	501[B]	HEM	O1D-CGD	2.13	1.29	1.22
2	A	501[A]	HEM	C4B-NB	-2.09	1.34	1.38

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	502	ZP6	O1-C8-N1	5.62	117.52	111.02
2	A	501[A]	HEM	CHC-C4B-NB	5.60	130.45	124.42
2	A	501[B]	HEM	CHC-C4B-NB	5.09	129.90	124.42
2	B	501[B]	HEM	C1B-NB-C4B	5.04	111.17	105.21
2	B	501[B]	HEM	CHC-C4B-NB	5.00	129.80	124.42
3	A	502	ZP6	O1-C8-N1	5.00	116.80	111.02
2	B	501[A]	HEM	CHC-C4B-NB	4.79	129.57	124.42
3	B	502	ZP6	O1-C8-O2	-4.37	117.16	124.76
2	A	501[B]	HEM	C1B-NB-C4B	4.36	110.37	105.21
2	B	501[A]	HEM	CHD-C4C-NC	4.28	129.12	124.45
2	B	501[A]	HEM	C1B-NB-C4B	4.21	110.19	105.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501[A]	HEM	C1B-NB-C4B	4.07	110.02	105.21
2	B	501[B]	HEM	C3B-C4B-NB	-3.89	106.68	109.47
3	A	502	ZP6	O1-C8-O2	-3.77	118.21	124.76
2	B	501[B]	HEM	CHB-C1B-NB	3.69	128.94	124.37
2	A	501[B]	HEM	CHD-C4C-NC	3.59	128.36	124.45
2	A	501[A]	HEM	CHD-C4C-NC	3.40	128.16	124.45
2	A	501[A]	HEM	CHB-C1B-NB	3.20	128.32	124.37
2	A	501[B]	HEM	CHD-C1D-ND	3.12	127.78	124.42
3	A	502	ZP6	C7-O1-C8	3.09	123.28	115.51
2	B	501[B]	HEM	CHD-C1D-ND	3.06	127.71	124.42
2	A	501[B]	HEM	CHB-C1B-NB	3.00	128.08	124.37
2	A	501[B]	HEM	CHD-C1D-C2D	-2.88	120.48	125.03
2	B	501[A]	HEM	CHD-C1D-C2D	-2.87	120.49	125.03
2	B	501[B]	HEM	CHD-C1D-C2D	-2.84	120.55	125.03
2	A	501[A]	HEM	C3B-C4B-NB	-2.82	107.44	109.47
2	B	501[A]	HEM	O2A-CGA-CBA	2.76	122.73	114.00
2	B	501[A]	HEM	C4C-NC-C1C	2.65	110.14	105.82
2	A	501[B]	HEM	C3B-C4B-NB	-2.62	107.59	109.47
2	A	501[A]	HEM	CHD-C1D-C2D	-2.55	121.00	125.03
2	B	501[A]	HEM	C2D-C1D-ND	2.55	112.84	109.90
2	A	501[A]	HEM	CHD-C1D-ND	2.45	127.06	124.42
2	A	501[B]	HEM	CHA-C4D-ND	2.44	127.39	124.37
2	A	501[A]	HEM	CAD-C3D-C4D	2.40	128.88	124.70
2	B	501[B]	HEM	CHA-C4D-ND	2.39	127.32	124.37
3	B	502	ZP6	C7-O1-C8	2.38	121.51	115.51
2	B	501[A]	HEM	CHA-C4D-ND	2.36	127.28	124.37
4	B	503	STQ	C6-C5-C4	2.34	121.66	118.78
2	A	501[A]	HEM	CHA-C4D-ND	2.28	127.18	124.37
3	A	502	ZP6	C11-C12-C13	-2.25	106.43	111.21
2	B	501[A]	HEM	CHA-C4D-C3D	-2.22	121.13	125.23
2	A	501[A]	HEM	CHA-C4D-C3D	-2.21	121.16	125.23
2	A	501[B]	HEM	O2D-CGD-CBD	2.20	120.97	114.00
2	A	501[A]	HEM	O2A-CGA-CBA	2.19	120.91	114.00
2	A	501[B]	HEM	CHA-C4D-C3D	-2.12	121.31	125.23
2	B	501[B]	HEM	C4C-NC-C1C	2.12	109.28	105.82
3	B	502	ZP6	C4-C5-C6	2.11	121.37	118.23
2	A	501[A]	HEM	O2D-CGD-CBD	2.11	120.67	114.00
2	B	501[A]	HEM	O2D-CGD-CBD	2.11	120.66	114.00
3	A	502	ZP6	C16-C14-N2	-2.09	106.51	110.83
2	B	501[A]	HEM	CHD-C1D-ND	2.08	126.66	124.42
2	B	501[A]	HEM	CAD-C3D-C4D	2.07	128.31	124.70
3	B	502	ZP6	C14-N2-C13	2.04	126.03	121.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	ZP6	C10-C9-N1	2.01	106.70	103.24

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	504	GOL	O1-C1-C2-C3
5	B	504	GOL	C1-C2-C3-O3
5	B	505	GOL	O1-C1-C2-C3
5	B	507	GOL	O1-C1-C2-C3
5	B	504	GOL	O2-C2-C3-O3
5	B	505	GOL	O1-C1-C2-O2
5	B	505	GOL	O2-C2-C3-O3
5	B	505	GOL	C1-C2-C3-O3
5	B	507	GOL	C1-C2-C3-O3
5	B	504	GOL	O1-C1-C2-O2
5	B	506	GOL	O2-C2-C3-O3
5	B	507	GOL	O1-C1-C2-O2
2	A	501[B]	HEM	CAA-CBA-CGA-O1A
2	B	501[B]	HEM	CAA-CBA-CGA-O1A
2	B	501[B]	HEM	CAA-CBA-CGA-O2A
2	A	501[A]	HEM	CAD-CBD-CGD-O2D
2	A	501[B]	HEM	CAA-CBA-CGA-O2A
2	B	501[A]	HEM	CAD-CBD-CGD-O2D
2	A	501[A]	HEM	CAD-CBD-CGD-O1D
2	B	501[A]	HEM	CAD-CBD-CGD-O1D
2	B	501[A]	HEM	CAA-CBA-CGA-O2A

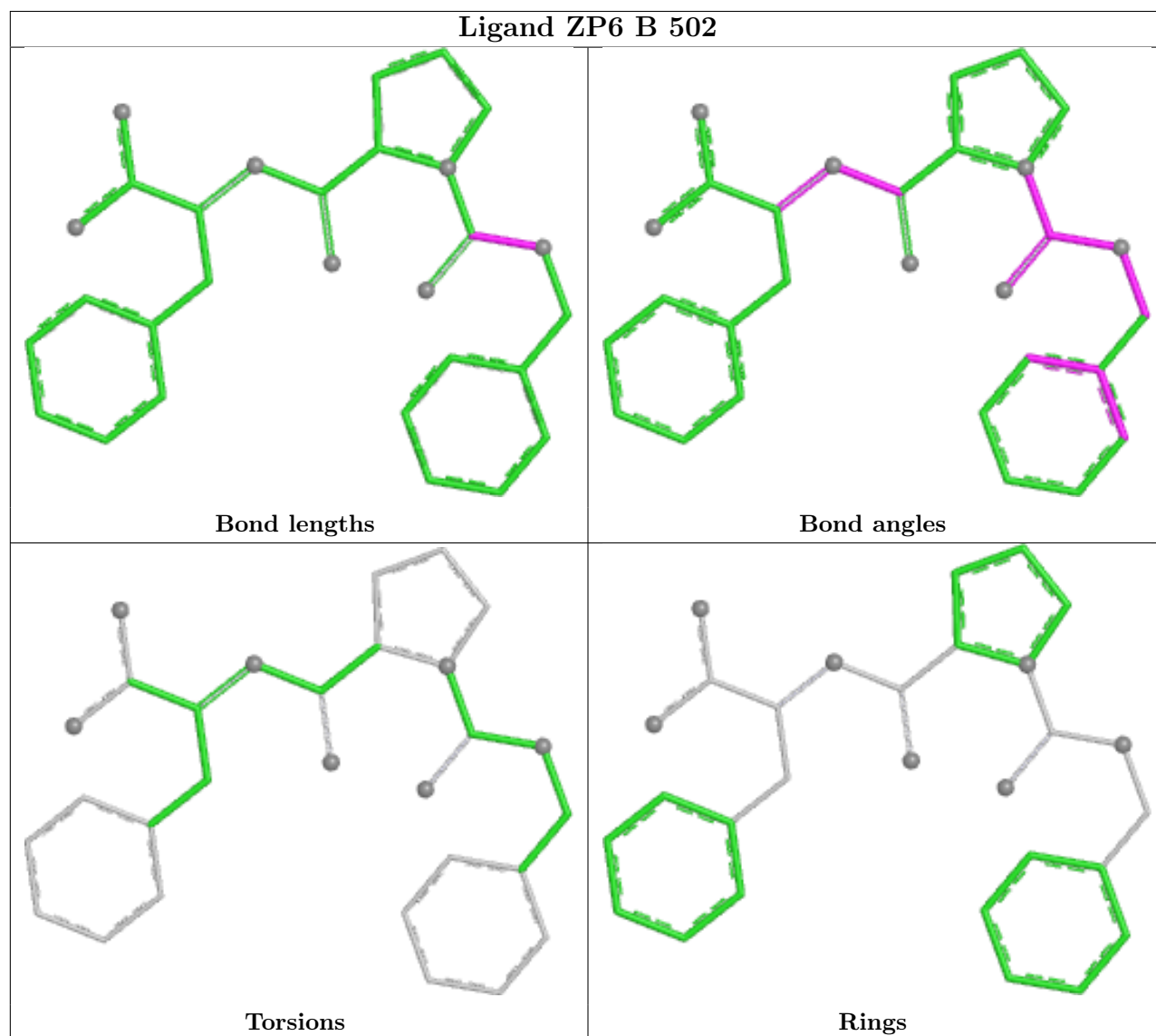
There are no ring outliers.

5 monomers are involved in 22 short contacts:

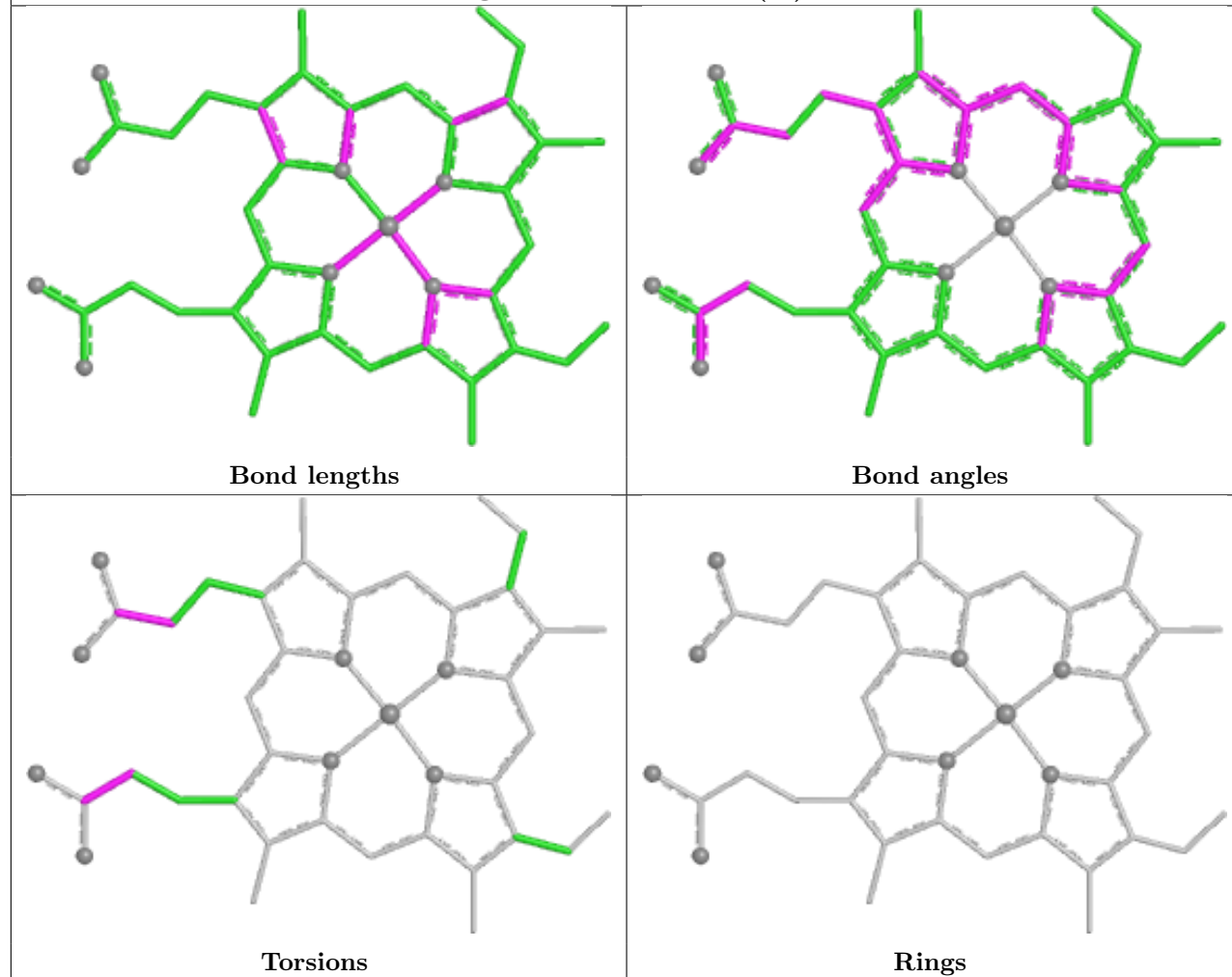
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501[A]	HEM	4	0
2	B	501[B]	HEM	6	0
2	A	501[A]	HEM	6	0
2	A	501[B]	HEM	5	0
5	A	504	GOL	1	0

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

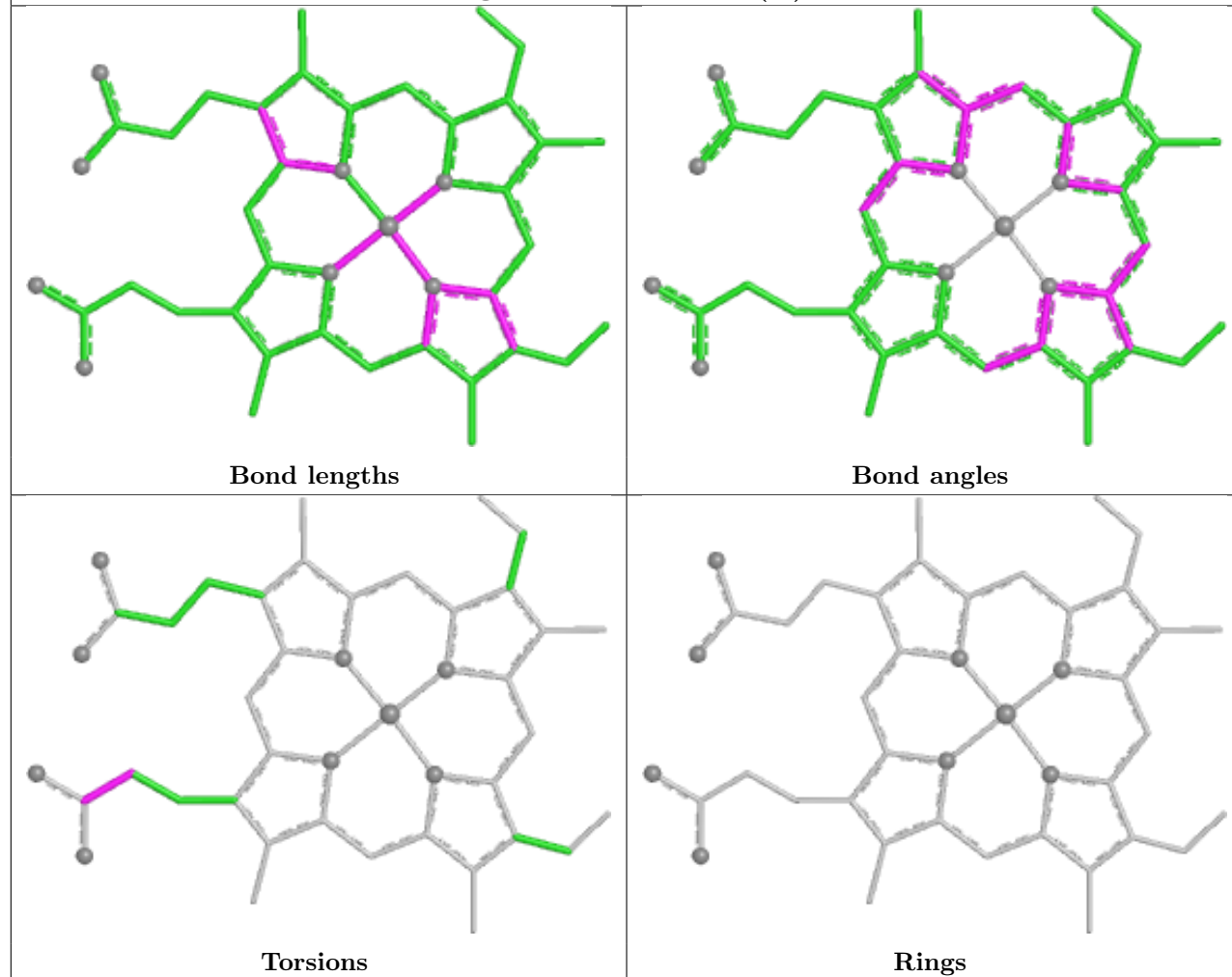
also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



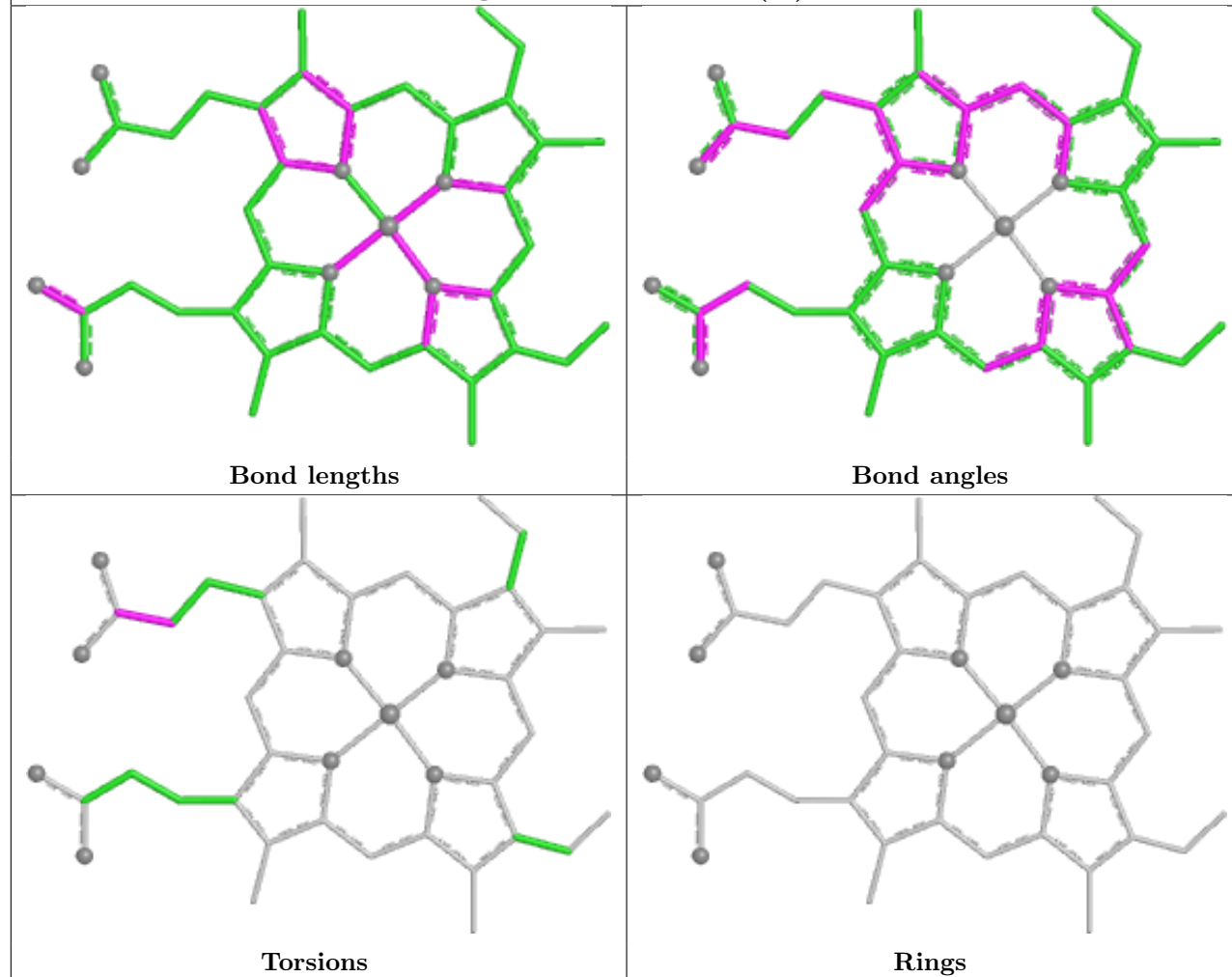
Ligand HEM B 501 (A)



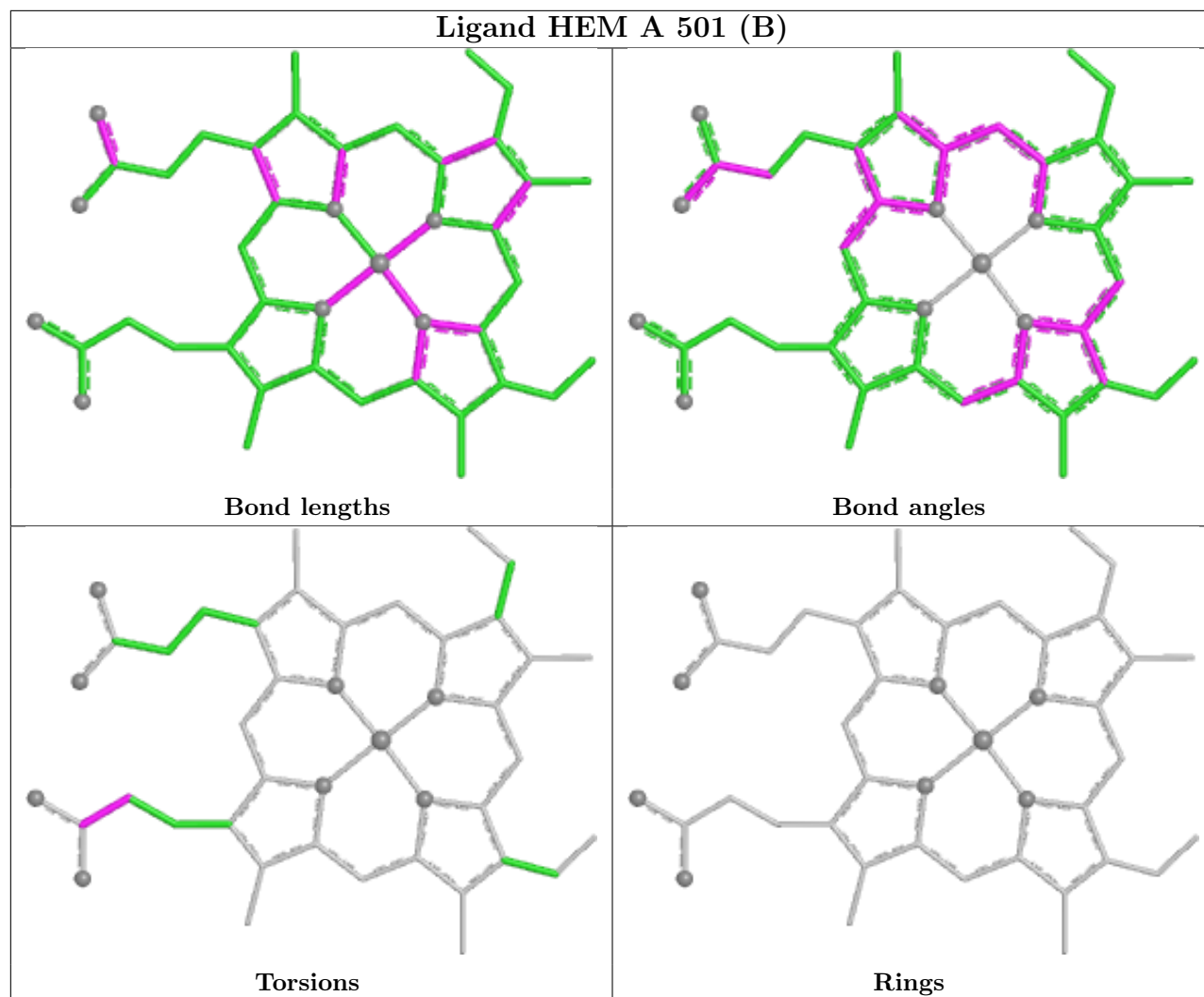
Ligand HEM B 501 (B)



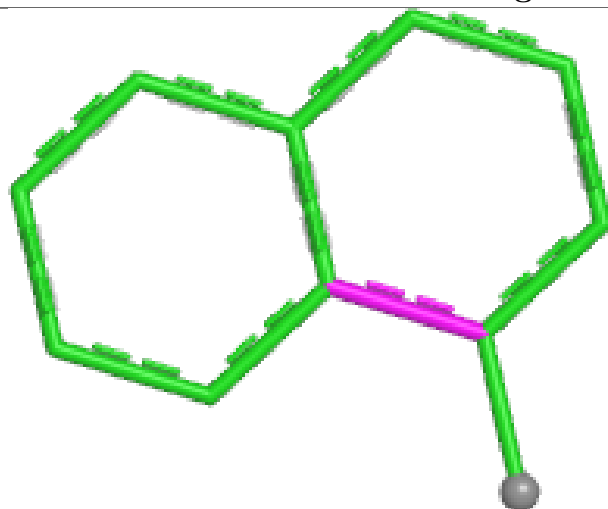
Ligand HEM A 501 (A)



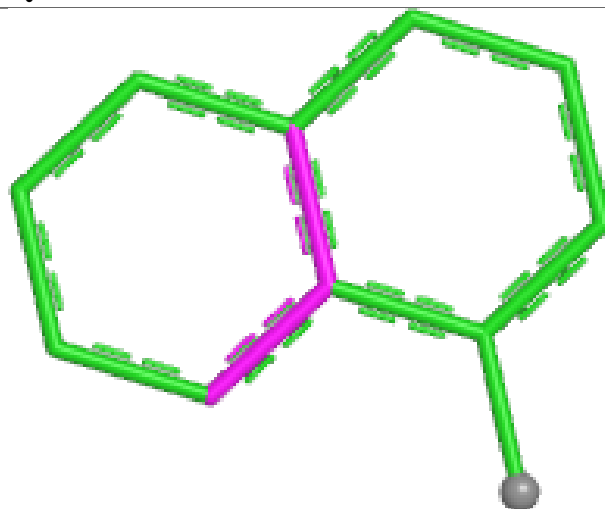
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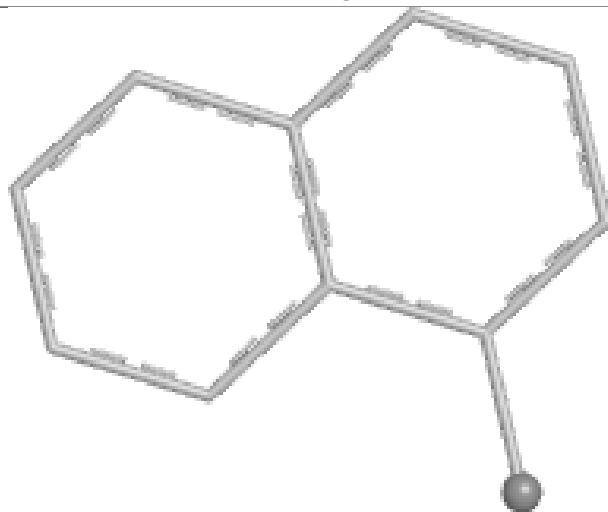
Ligand STQ B 503



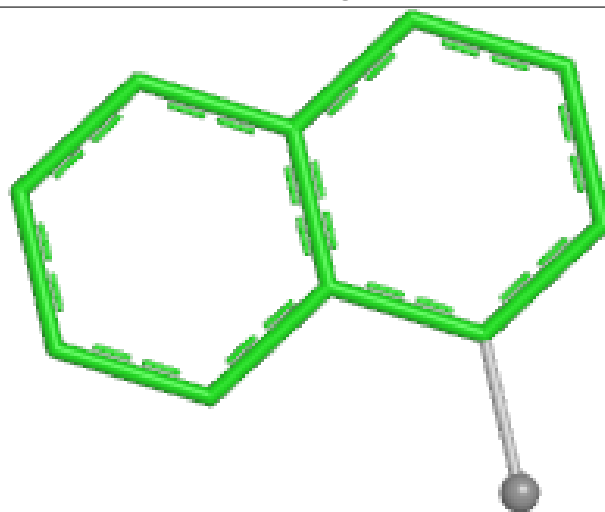
Bond lengths



Bond angles

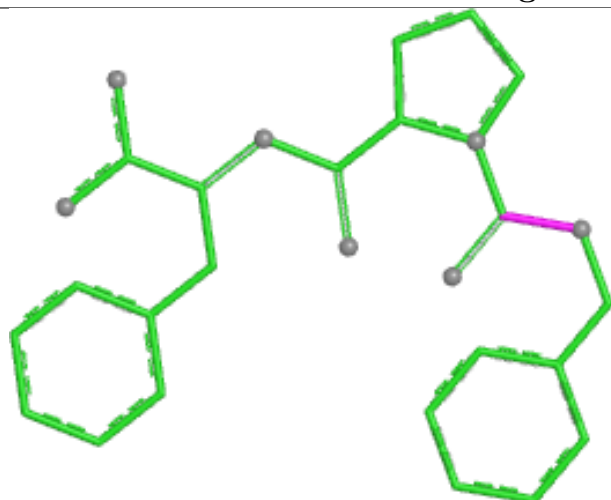


Torsions

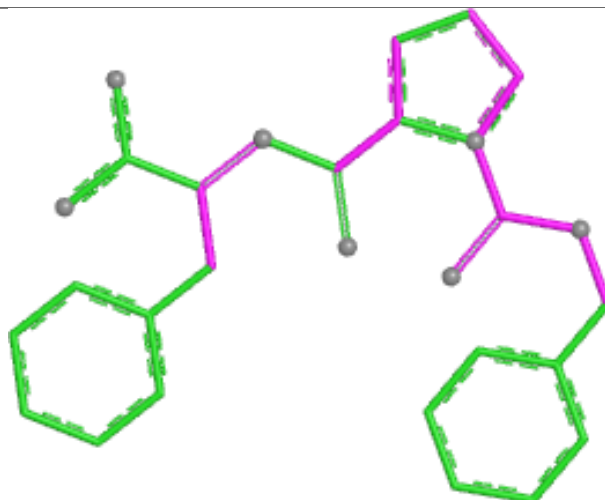


Rings

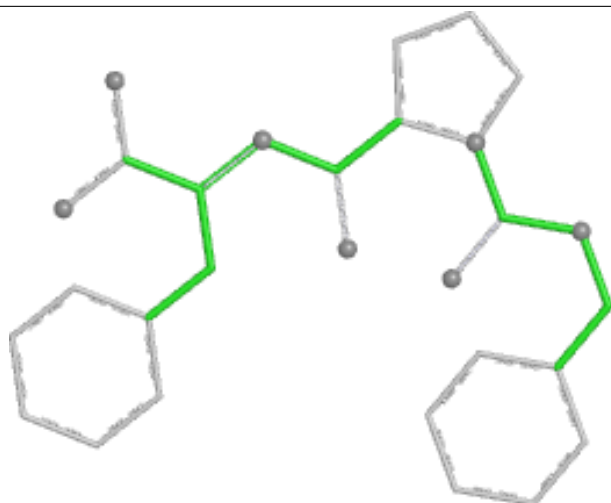
Ligand ZP6 A 502



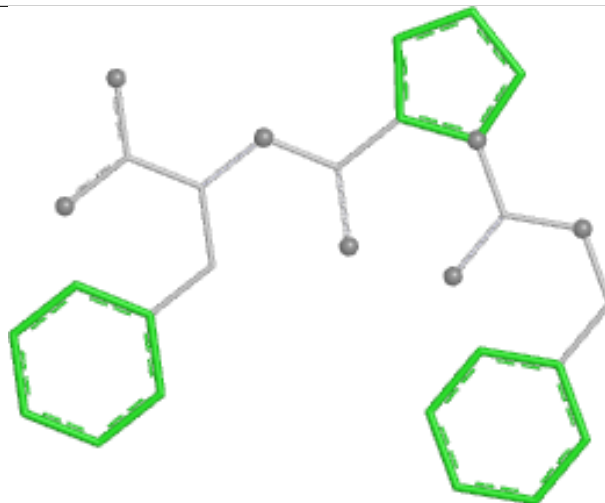
Bond lengths



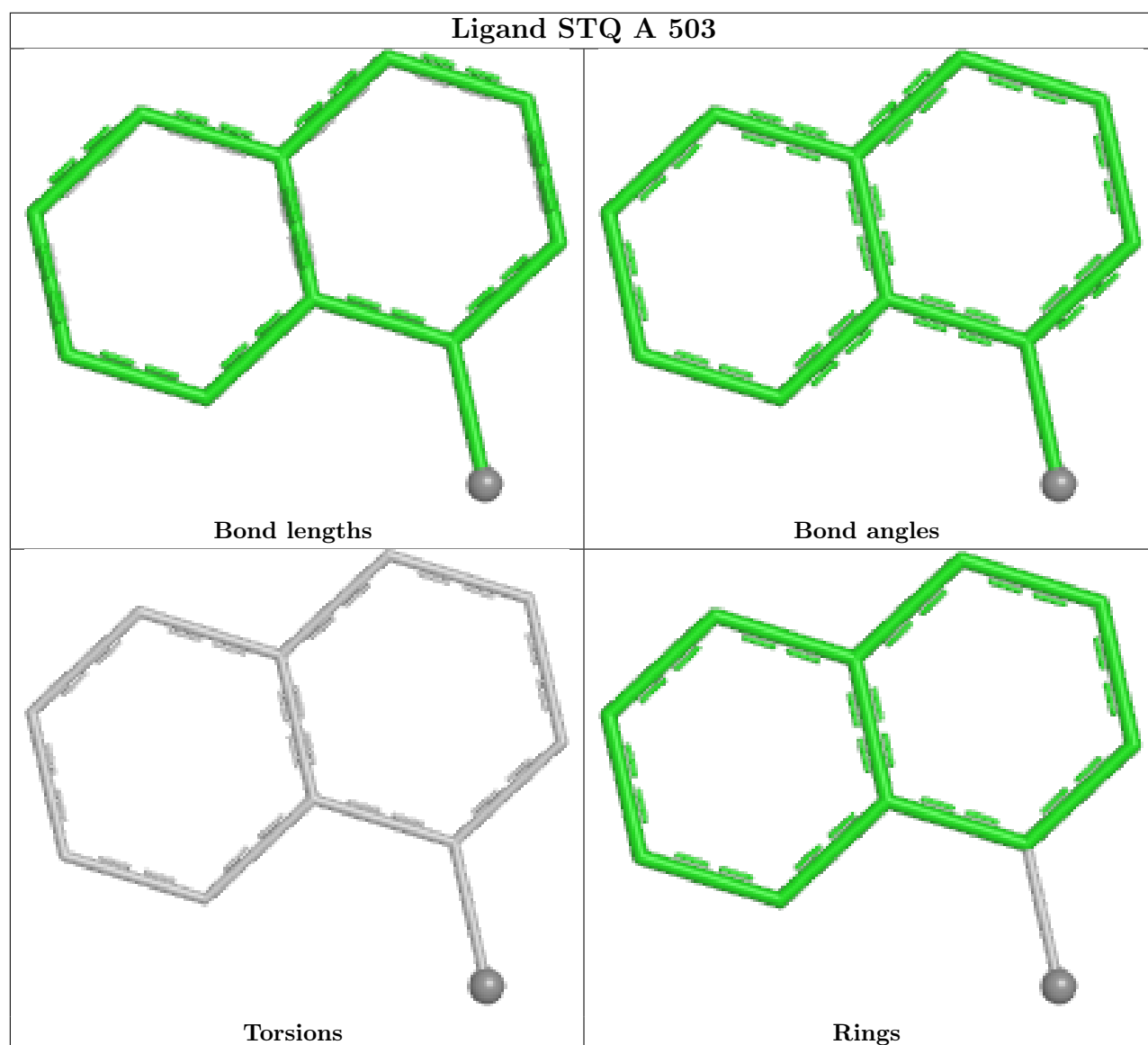
Bond angles



Torsions



Rings



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	453/456 (99%)	0.45	25 (5%)	30 31	11, 26, 50, 75	6 (1%)
1	B	453/456 (99%)	0.29	14 (3%)	51 54	8, 24, 45, 68	7 (1%)
All	All	906/912 (99%)	0.37	39 (4%)	40 41	8, 25, 48, 75	13 (1%)

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	191	ALA	4.8
1	A	191	ALA	4.7
1	A	197	ALA	3.8
1	A	198	TYR	3.5
1	A	158	PHE	3.4
1	A	196	PRO	3.4
1	A	225	ALA	3.3
1	B	158	PHE	3.3
1	A	455	LEU	3.1
1	A	211	VAL	3.0
1	A	136	ASP	2.9
1	A	195	ASP	2.9
1	A	188	LEU	2.8
1	B	455	LEU	2.8
1	B	231	ASP	2.8
1	A	190[A]	ARG	2.7
1	A	46	GLY	2.6
1	A	193	PRO	2.4
1	A	215	LEU	2.3
1	B	225	ALA	2.3
1	B	197	ALA	2.3
1	A	45	PRO	2.2
1	B	196	PRO	2.2
1	B	198	TYR	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	189	GLN	2.2
1	B	216	VAL	2.2
1	A	221	ALA	2.2
1	B	194	ASP	2.2
1	A	4	GLU	2.2
1	B	21	ASN	2.2
1	A	3	LYS	2.2
1	B	211	VAL	2.2
1	A	194	ASP	2.2
1	A	220	ILE	2.2
1	A	180	ALA	2.1
1	A	302	VAL	2.0
1	B	302	VAL	2.0
1	A	20	LEU	2.0
1	B	236	HIS	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	GOL	A	504	6/6	0.80	0.16	41,48,51,56	0
5	GOL	B	504	6/6	0.81	0.20	55,56,61,61	0
5	GOL	B	505	6/6	0.82	0.16	49,54,54,56	0
5	GOL	B	507	6/6	0.85	0.14	38,45,51,54	0
3	ZP6	A	502	29/29	0.86	0.14	29,38,46,49	0
5	GOL	B	506	6/6	0.89	0.12	29,40,42,47	0
3	ZP6	B	502	29/29	0.89	0.10	25,30,35,41	0
4	STQ	A	503	11/11	0.93	0.09	16,19,20,20	0

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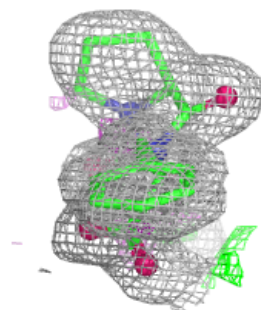
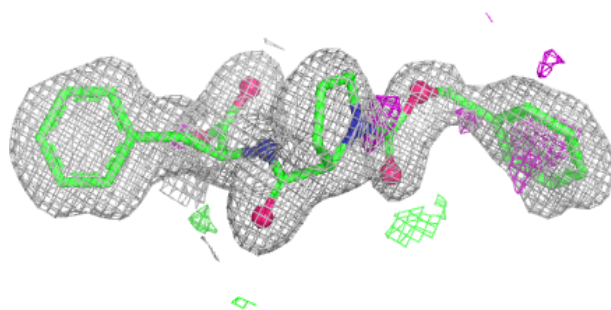
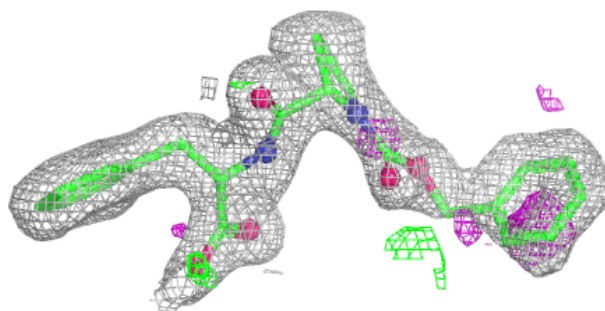
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	STQ	B	503	11/11	0.96	0.06	15,17,18,19	0
2	HEM	A	501[B]	43/43	0.98	0.06	17,18,19,19	43
2	HEM	B	501[A]	43/43	0.98	0.06	15,16,17,20	43
2	HEM	B	501[B]	43/43	0.98	0.06	15,16,17,17	43
2	HEM	A	501[A]	43/43	0.98	0.06	16,17,18,19	43

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.

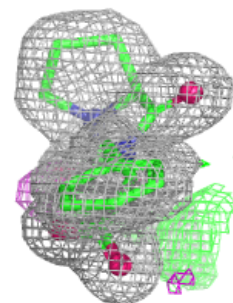
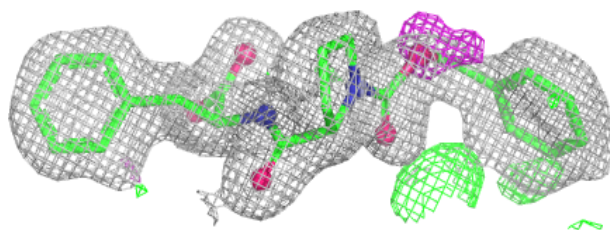
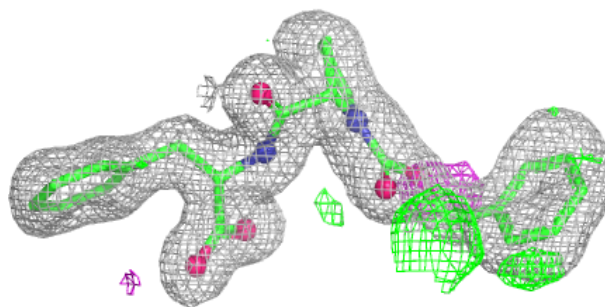
Electron density around ZP6 A 502:

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



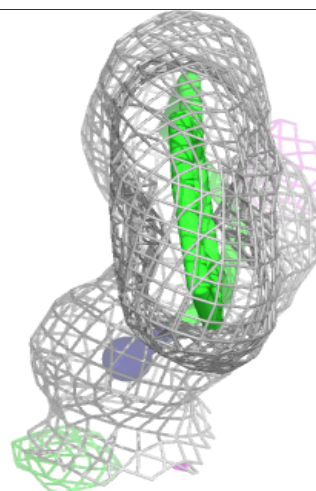
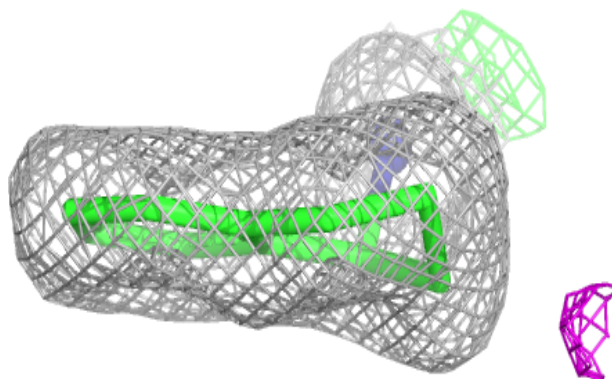
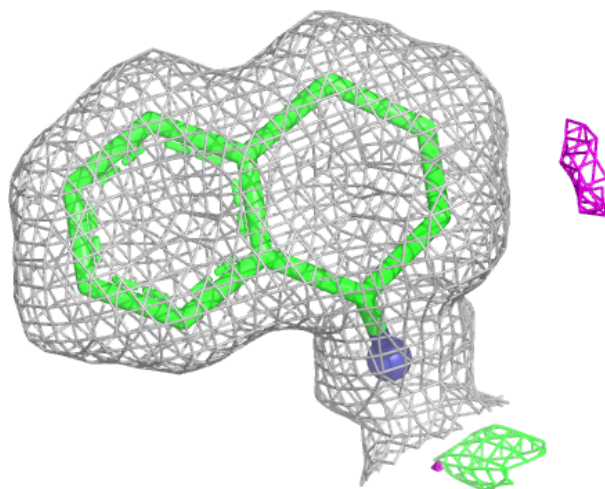
Electron density around ZP6 B 502:

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



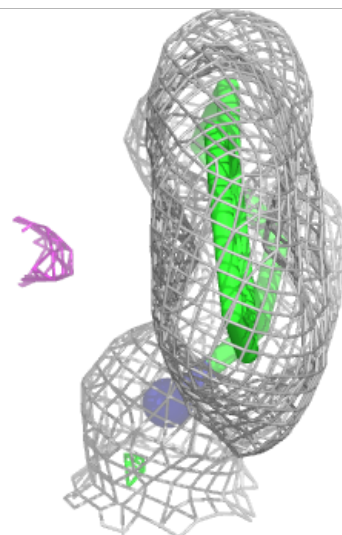
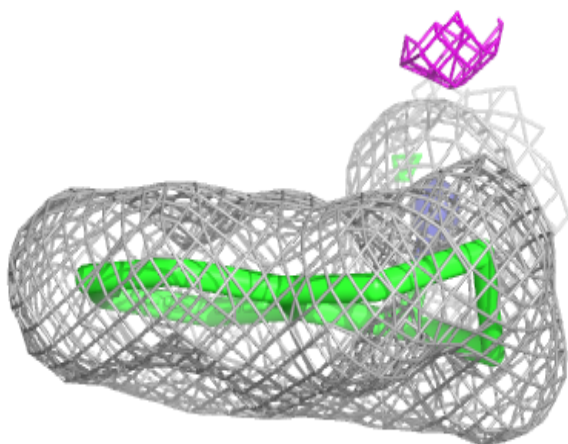
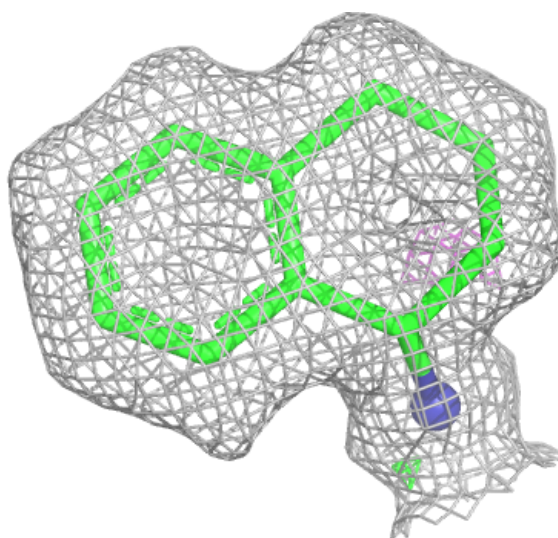
Electron density around STQ A 503:

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



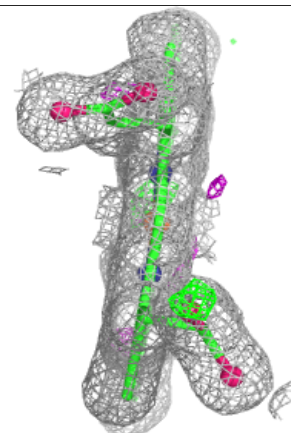
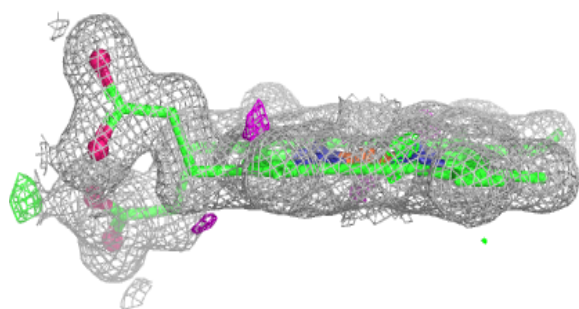
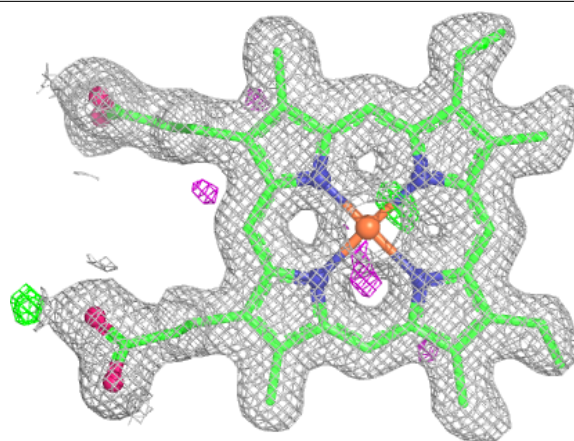
Electron density around STQ B 503:

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and green (positive)



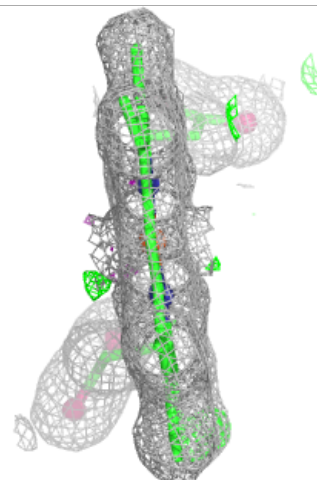
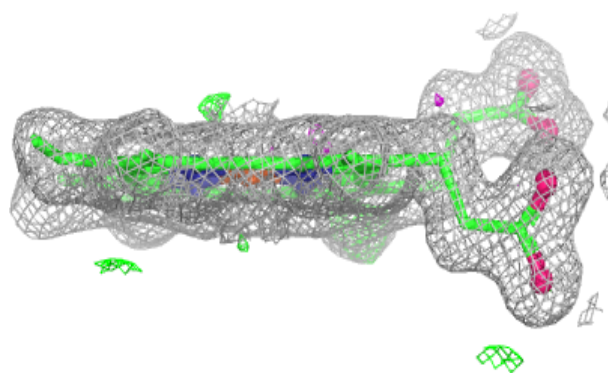
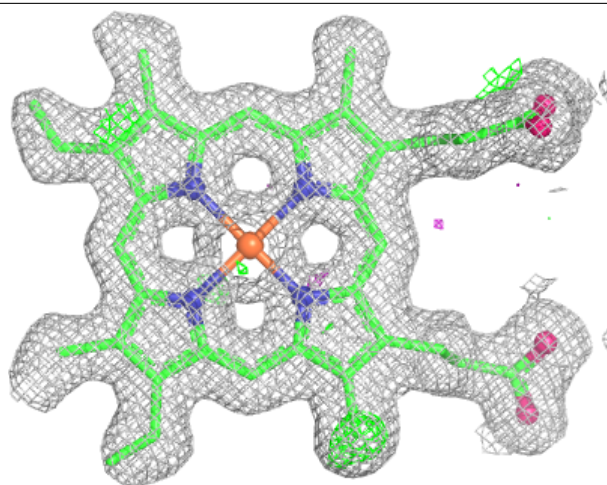
Electron density around HEM A 501 (B):

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



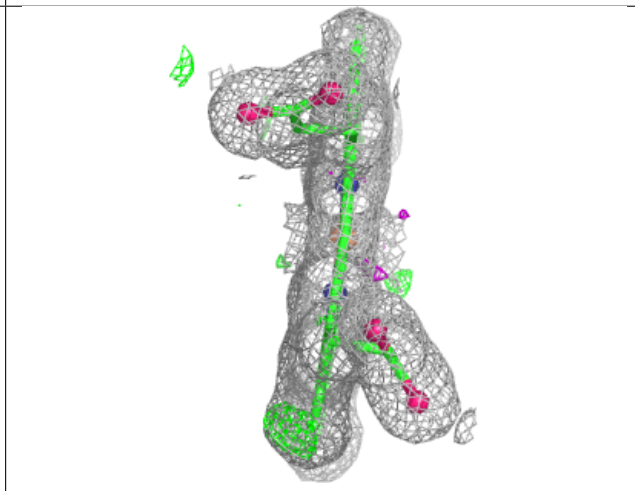
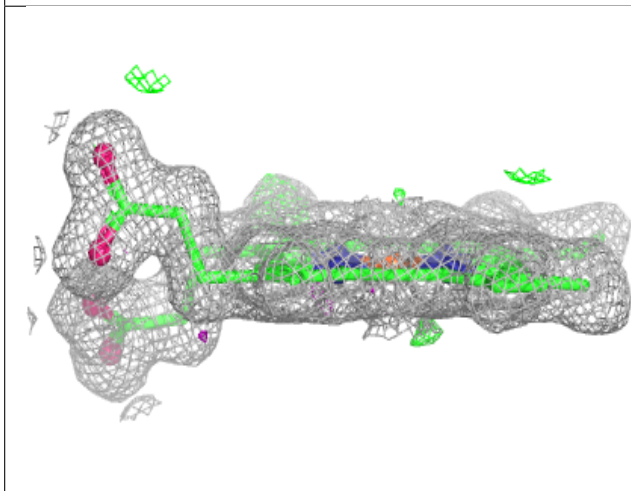
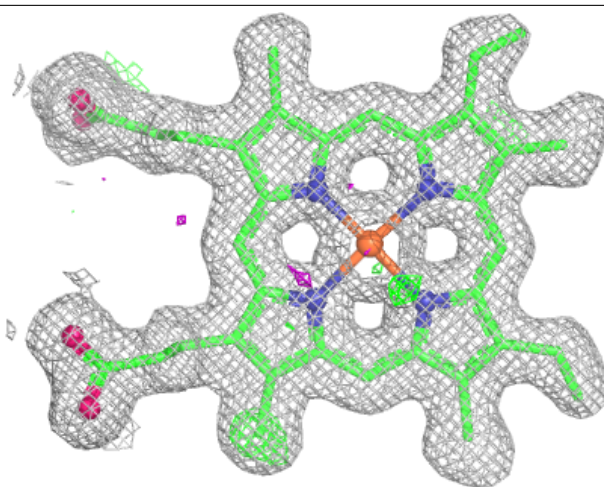
Electron density around HEM B 501 (A):

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



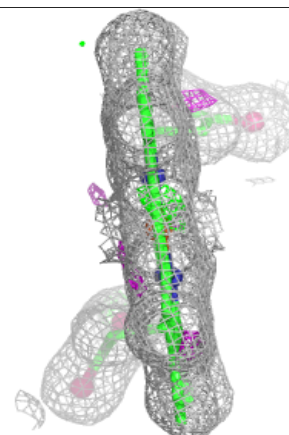
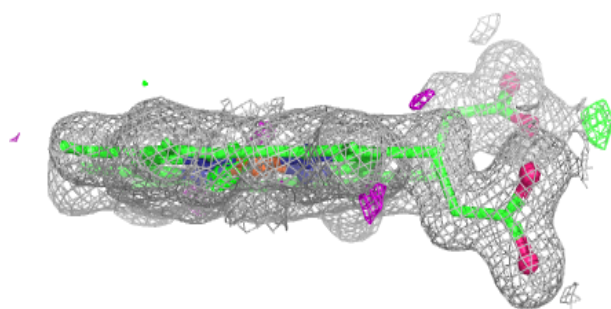
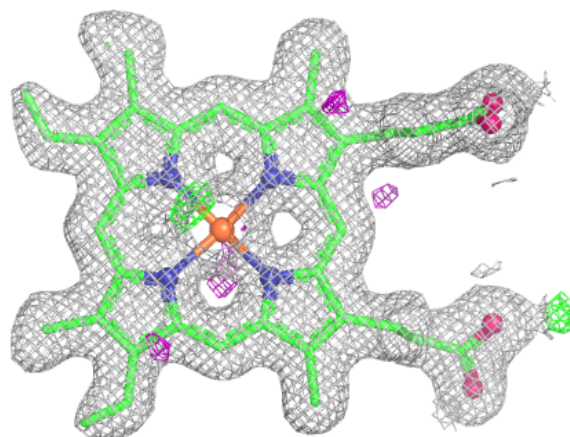
Electron density around HEM B 501 (B):

$2mF_o-DF_c$ (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 (A):

$2mF_o-DF_c$ (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.