



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:33 PM UTC

PDB ID : 6PO5 / pdb_00006po5
Title : Structure of human neuronal nitric oxide synthase R354A/G357D mutant heme domain in complex with 7-(3-(Aminomethyl)-4-(cyclobutylmethoxy)phenyl)-4-methylquinolin-2-amine
Authors : Li, H.; Poulos, T.L.
Deposited on : 2019-07-03
Resolution : 1.82 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

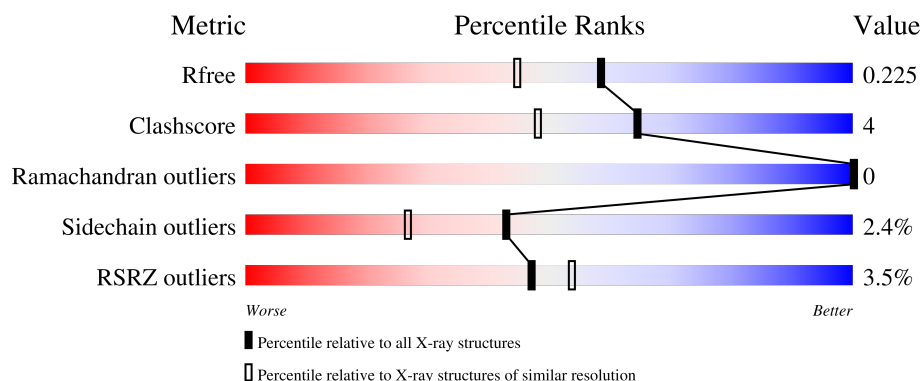
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.82 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	1112 (1.82-1.82)
Clashscore	190562	1148 (1.82-1.82)
Ramachandran outliers	187476	1140 (1.82-1.82)
Sidechain outliers	187428	1140 (1.82-1.82)
RSRZ outliers	180081	1112 (1.82-1.82)

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	421	3% 89% 9% ..
1	B	421	4% 87% 11% .

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 7701 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.

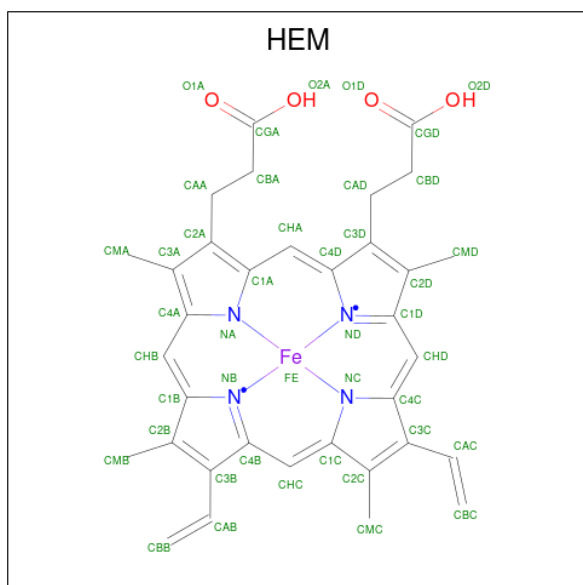
- Molecule 1 is a protein called Nitric oxide synthase, brain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	417	Total	C	N	O	S	0	5	0
			3416	2186	582	625	23			
1	B	411	Total	C	N	O	S	0	3	0
			3362	2154	572	615	21			

There are 4 discrepancies between the modelled and reference sequences:

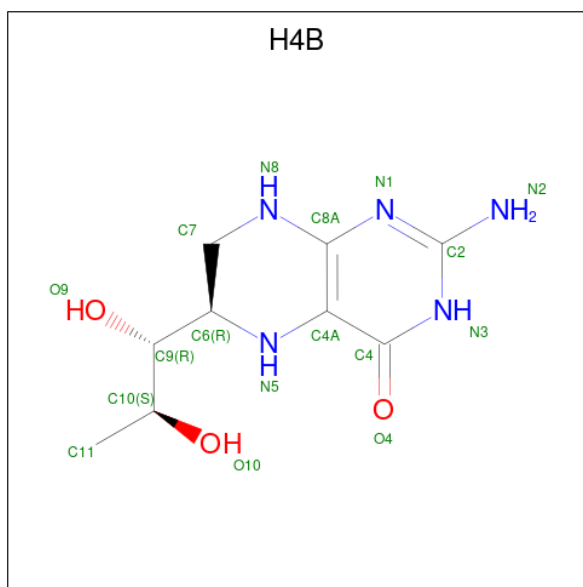
Chain	Residue	Modelled	Actual	Comment	Reference
A	354	ALA	ARG	engineered mutation	UNP P29475
A	357	ASP	GLY	engineered mutation	UNP P29475
B	354	ALA	ARG	engineered mutation	UNP P29475
B	357	ASP	GLY	engineered mutation	UNP P29475

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



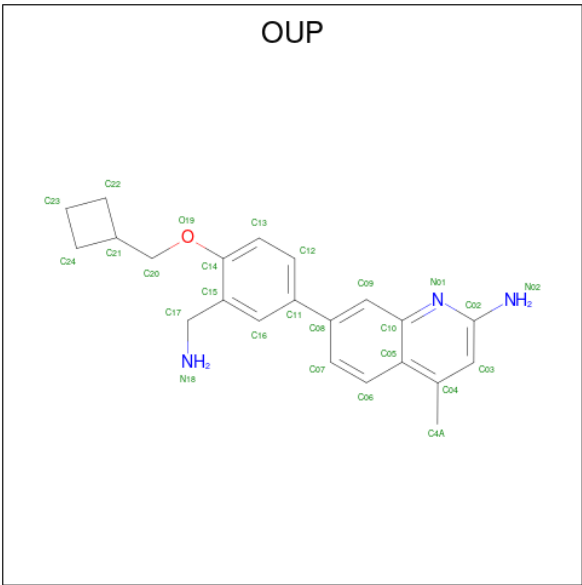
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

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



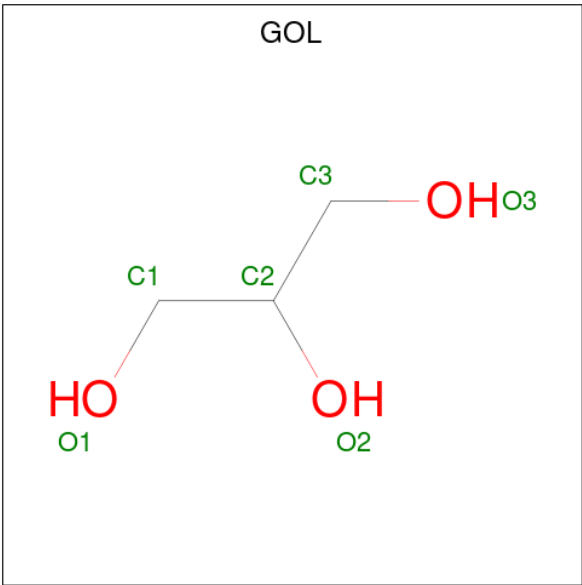
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			17	9	5	3		
3	B	1	Total	C	N	O	0	0
			17	9	5	3		

- Molecule 4 is 7-[3-(aminomethyl)-4-(cyclobutylmethoxy)phenyl]-4-methylquinolin-2-amine (CCD ID: OUP) (formula: $C_{22}H_{25}N_3O$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			26	22	3	1		
4	B	1	Total	C	N	O	0	0
			26	22	3	1		

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

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

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

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

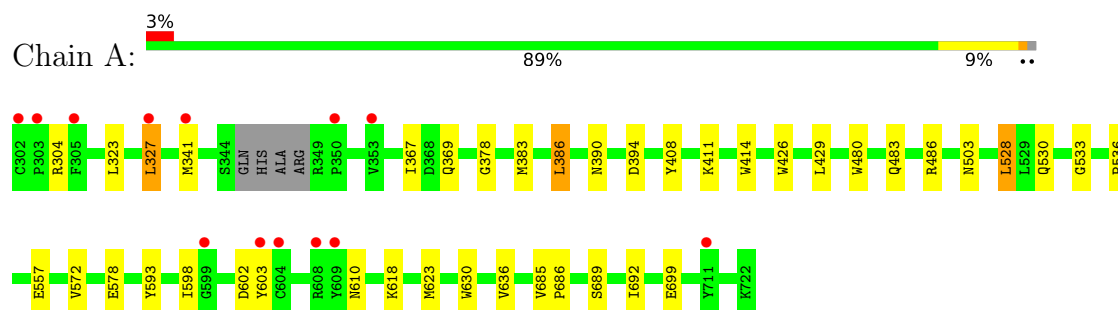
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	407	Total	O	0	0
			407	407		
7	B	325	Total	O	0	0
			325	325		

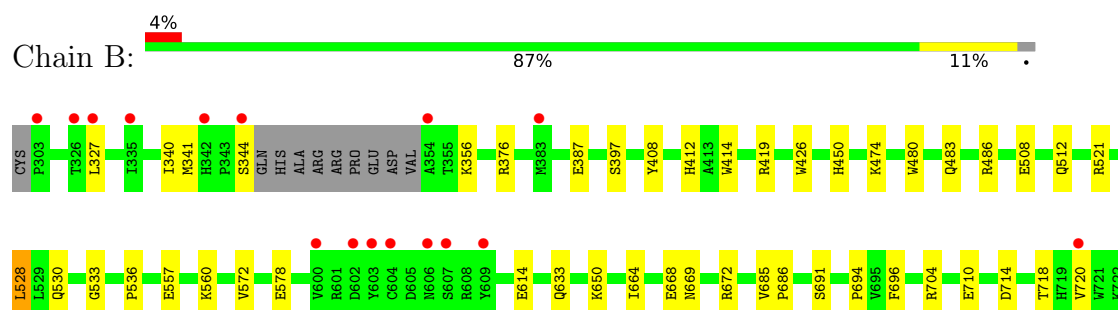
3 Residue-property plots

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, brain



- Molecule 1: Nitric oxide synthase, brain



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	52.24Å 123.73Å 165.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.58 – 1.82 41.58 – 1.82	Depositor EDS
% Data completeness (in resolution range)	98.8 (41.58-1.82) 99.6 (41.58-1.82)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.43 (at 1.82Å)	Xtriage
Refinement program	PHENIX (1.11.1-2575_1496: ???)	Depositor
R, R_{free}	0.189 , 0.228 (Not available) , 0.225	Depositor DCC
R_{free} test set	4758 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	31.9	Xtriage
Anisotropy	0.589	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 50.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7701	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.40% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, H4B, ZN, HEM, OUP

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	0.27	0/3528	0.48	0/4788
1	B	0.25	0/3467	0.44	0/4702
All	All	0.26	0/6995	0.46	0/9490

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

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	3416	0	3331	25	0
1	B	3362	0	3276	24	0
2	A	43	0	30	3	0
2	B	43	0	30	4	0
3	A	17	0	15	0	0
3	B	17	0	15	0	0
4	A	26	0	0	2	0
4	B	26	0	0	2	0
5	A	12	0	16	2	0
5	B	6	0	8	0	0
6	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	A	407	0	0	3	0
7	B	325	0	0	3	0
All	All	7701	0	6721	53	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (53) 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:B:802:HEM:HHC	2:B:802:HEM:HBB2	1.70	0.74
1:A:390[A]:ASN:ND2	7:A:901:HOH:O	2.18	0.74
1:A:528:LEU:HD22	1:A:536:PRO:HB2	1.70	0.72
1:B:327:LEU:HB2	1:B:704:ARG:HG2	1.74	0.69
1:B:528:LEU:HD22	1:B:536:PRO:HB2	1.75	0.69
1:A:429:LEU:O	5:A:805:GOL:H12	1.93	0.69
1:B:704:ARG:NH1	1:B:710:GLU:OE2	2.28	0.67
1:B:714:ASP:OD1	7:B:901:HOH:O	2.11	0.66
2:A:801:HEM:HMC2	2:A:801:HEM:HBC2	1.82	0.62
2:B:802:HEM:HMC2	2:B:802:HEM:HBC2	1.82	0.59
1:B:664:ILE:O	1:B:668[B]:GLU:HG2	2.02	0.59
1:B:483:GLN:HB2	1:B:486:ARG:HG3	1.85	0.57
1:B:356:LYS:HG3	1:B:397:SER:HA	1.87	0.56
1:B:572:VAL:HG21	4:B:804:OUP:C07	2.36	0.56
1:A:411:LYS:NZ	5:A:805:GOL:H11	2.22	0.55
1:B:512:GLN:NE2	7:B:905:HOH:O	2.40	0.54
1:B:530:GLN:HE21	1:B:533:GLY:HA2	1.73	0.53
1:A:572:VAL:HG21	4:A:803:OUP:C07	2.39	0.53
1:A:480:TRP:HB2	1:A:528:LEU:HB3	1.92	0.52
1:B:614:GLU:HG2	7:B:1196:HOH:O	2.10	0.51
1:B:419:ARG:HB2	2:B:802:HEM:HBD2	1.94	0.49
1:A:483:GLN:HB2	1:A:486:ARG:HG2	1.93	0.49
1:A:369:GLN:NE2	7:A:907:HOH:O	2.42	0.49
2:A:801:HEM:HBA2	4:A:803:OUP:C12	2.43	0.49
1:B:414:TRP:CE3	1:B:426:TRP:HA	2.47	0.49
1:B:691:SER:HA	1:B:696:PHE:CG	2.48	0.48
1:A:689:SER:HB3	1:A:692:ILE:HD11	1.95	0.48
1:A:598:ILE:HA	1:A:602:ASP:HB2	1.96	0.47
1:A:623:MET:HA	1:A:630:TRP:CD1	2.49	0.47
1:A:327:LEU:H	1:A:327:LEU:HD23	1.79	0.47
1:A:378:GLY:HA2	1:A:383:MET:HE3	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:341:MET:HE3	1:B:341:MET:HB2	1.75	0.47
1:A:699:GLU:HB3	1:B:340:ILE:HD13	1.98	0.46
1:B:480:TRP:HB2	1:B:528:LEU:HB3	1.97	0.46
1:A:304:ARG:HG3	1:A:323:LEU:HD11	1.97	0.46
1:A:414:TRP:CE3	1:A:426:TRP:HA	2.50	0.46
1:B:557:GLU:HG3	1:B:560:LYS:NZ	2.31	0.45
1:A:603:TYR:O	1:A:610:ASN:N	2.49	0.45
2:A:801:HEM:HMD1	2:A:801:HEM:HBD1	1.98	0.44
2:B:802:HEM:HBA2	4:B:804:OUP:C12	2.49	0.43
1:A:341:MET:HE1	7:A:1047:HOH:O	2.19	0.43
1:A:636:VAL:HG11	1:B:633:GLN:HG3	2.01	0.42
1:B:408:TYR:CE1	1:B:412:HIS:CE1	3.08	0.42
1:A:486:ARG:HE	1:A:503:ASN:HD21	1.68	0.42
1:B:664:ILE:HG13	1:B:694:PRO:HB2	2.00	0.42
1:A:593:TYR:CD1	1:A:598:ILE:HD11	2.55	0.42
1:A:685:VAL:HA	1:A:686:PRO:HD3	1.93	0.41
1:B:450:HIS:CD2	1:B:450:HIS:C	2.98	0.41
1:A:530:GLN:HE21	1:A:533:GLY:HA2	1.85	0.41
1:B:685:VAL:HA	1:B:686:PRO:HD3	1.91	0.41
1:A:369:GLN:HG2	1:A:408:TYR:OH	2.21	0.41
1:B:669:ASN:OD1	1:B:672:ARG:NH2	2.53	0.41
1:A:367:ILE:HG21	1:A:386:LEU:HD13	2.03	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	418/421 (99%)	411 (98%)	7 (2%)	0	100	100
1	B	410/421 (97%)	401 (98%)	9 (2%)	0	100	100
All	All	828/842 (98%)	812 (98%)	16 (2%)	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	378/376 (100%)	371 (98%)	7 (2%)	50	37
1	B	370/376 (98%)	359 (97%)	11 (3%)	36	17
All	All	748/752 (100%)	730 (98%)	18 (2%)	43	27

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	327	LEU
1	A	386	LEU
1	A	394	ASP
1	A	528	LEU
1	A	557	GLU
1	A	578	GLU
1	A	618	LYS
1	B	344	SER
1	B	376	ARG
1	B	387	GLU
1	B	474	LYS
1	B	508	GLU
1	B	521	ARG
1	B	528	LEU
1	B	578	GLU
1	B	650	LYS
1	B	718	THR
1	B	720	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	412	HIS

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Mol	Chain	Res	Type
1	B	412	HIS
1	B	513	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 10 ligands modelled in this entry, 1 is monoatomic - leaving 9 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	H4B	A	802	-	17,18,18	0.83	1 (5%)	14,26,26	1.66	3 (21%)
2	HEM	B	802	1	50,50,50	1.61	11 (22%)	67,82,82	1.36	6 (8%)
2	HEM	A	801	1	50,50,50	1.45	8 (16%)	67,82,82	1.35	8 (11%)
5	GOL	A	805	-	5,5,5	0.44	0	5,5,5	0.58	0
4	OUP	B	804	-	29,29,29	0.98	2 (6%)	40,41,41	1.45	5 (12%)
4	OUP	A	803	-	29,29,29	0.95	1 (3%)	40,41,41	1.42	4 (10%)
5	GOL	A	804	-	5,5,5	0.37	0	5,5,5	0.29	0
5	GOL	B	805	-	5,5,5	0.37	0	5,5,5	0.25	0
3	H4B	B	803	-	17,18,18	0.87	0	14,26,26	1.81	5 (35%)

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	H4B	A	802	-	-	3/8/17/17	0/2/2/2
2	HEM	B	802	1	-	5/14/54/54	-
2	HEM	A	801	1	-	4/14/54/54	-
5	GOL	A	805	-	-	4/4/4/4	-
4	OUP	B	804	-	-	6/11/17/17	0/4/4/4
4	OUP	A	803	-	-	3/11/17/17	0/4/4/4
5	GOL	A	804	-	-	2/4/4/4	-
5	GOL	B	805	-	-	2/4/4/4	-
3	H4B	B	803	-	-	3/8/17/17	0/2/2/2

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	802	HEM	FE-NA	5.25	2.12	1.95
2	A	801	HEM	FE-NC	4.70	2.10	1.95
2	B	802	HEM	FE-NC	4.29	2.09	1.95
2	A	801	HEM	FE-NA	3.23	2.05	1.95
2	B	802	HEM	CAC-C3C	3.09	1.55	1.47
2	B	802	HEM	CAB-C3B	3.04	1.55	1.47
2	A	801	HEM	CAB-C3B	3.04	1.55	1.47
2	B	802	HEM	FE-NB	3.03	2.04	1.94
2	A	801	HEM	CAC-C3C	2.77	1.54	1.47
2	A	801	HEM	CMC-C2C	2.59	1.56	1.50
2	B	802	HEM	CMC-C2C	2.45	1.55	1.50
2	B	802	HEM	CMD-C2D	2.45	1.55	1.50
4	B	804	OUP	C05-C10	-2.41	1.38	1.42
2	B	802	HEM	FE-ND	2.38	2.02	1.94
2	A	801	HEM	C3B-C2B	-2.19	1.32	1.37
2	B	802	HEM	CMB-C2B	2.19	1.55	1.50
2	A	801	HEM	C2A-C3A	-2.13	1.33	1.38
4	A	803	OUP	C05-C10	-2.11	1.39	1.42
2	B	802	HEM	C2A-C3A	-2.10	1.33	1.38
2	A	801	HEM	CMB-C2B	2.10	1.55	1.50
3	A	802	H4B	C4-N3	-2.08	1.34	1.38
2	B	802	HEM	C3B-C2B	-2.05	1.33	1.37
4	B	804	OUP	C02-N01	2.00	1.36	1.33

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	802	HEM	CBA-CAA-C2A	-4.68	99.59	112.53
3	B	803	H4B	C2-N1-C8A	4.07	120.54	113.36
4	A	803	OUP	O19-C14-C15	4.06	121.14	115.89
4	B	804	OUP	C04-C05-C10	3.97	120.42	118.00
4	A	803	OUP	C04-C05-C10	3.72	120.27	118.00
4	B	804	OUP	C05-C10-N01	-3.67	118.91	122.80
2	A	801	HEM	CBA-CAA-C2A	-3.64	102.47	112.53
3	A	802	H4B	C2-N1-C8A	3.64	119.77	113.36
4	A	803	OUP	C05-C10-N01	-3.44	119.15	122.80
2	B	802	HEM	C3B-C2B-C1B	3.12	108.75	106.41
4	B	804	OUP	O19-C14-C15	2.98	119.74	115.89
2	A	801	HEM	CAD-C3D-C4D	-2.91	119.62	124.70
2	A	801	HEM	CAD-C3D-C2D	2.80	133.12	127.87
2	A	801	HEM	CHA-C4D-ND	2.64	127.63	124.37
2	B	802	HEM	C2A-C1A-NA	-2.61	107.26	110.15
2	A	801	HEM	CBD-CAD-C3D	2.53	119.53	112.53
3	B	803	H4B	O4-C4-C4A	-2.49	121.26	127.26
3	A	802	H4B	C4-C4A-N5	2.41	122.85	116.27
2	B	802	HEM	CHA-C4D-ND	2.34	127.26	124.37
3	B	803	H4B	C4A-C4-N3	2.34	118.55	112.13
3	A	802	H4B	C4A-C4-N3	2.31	118.47	112.13
2	A	801	HEM	C2A-C1A-NA	-2.30	107.60	110.15
3	B	803	H4B	N2-C2-N3	2.25	121.51	116.76
2	B	802	HEM	C4D-ND-C1D	2.24	107.86	105.21
4	B	804	OUP	C07-C08-C11	-2.21	117.44	121.25
2	A	801	HEM	C3B-C2B-C1B	2.19	108.06	106.41
2	B	802	HEM	O1D-CGD-CBD	-2.18	116.16	123.09
4	A	803	OUP	O19-C14-C13	-2.13	119.30	123.95
2	A	801	HEM	CAB-C3B-C2B	-2.04	121.78	128.43
3	B	803	H4B	N3-C2-N1	-2.04	119.59	123.32
4	B	804	OUP	C06-C05-C04	-2.01	119.72	123.61

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	801	HEM	C2D-C3D-CAD-CBD
2	B	802	HEM	C2D-C3D-CAD-CBD
2	B	802	HEM	C4D-C3D-CAD-CBD
4	A	803	OUP	O19-C20-C21-C22
4	B	804	OUP	O19-C20-C21-C22

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Mol	Chain	Res	Type	Atoms
4	B	804	OUP	O19-C20-C21-C24
5	A	804	GOL	O1-C1-C2-C3
5	A	805	GOL	O1-C1-C2-C3
5	B	805	GOL	O1-C1-C2-C3
2	A	801	HEM	C4D-C3D-CAD-CBD
5	A	805	GOL	C1-C2-C3-O3
5	A	805	GOL	O1-C1-C2-O2
5	A	805	GOL	O2-C2-C3-O3
5	B	805	GOL	O1-C1-C2-O2
4	B	804	OUP	C13-C14-O19-C20
3	B	803	H4B	C7-C6-C9-O9
5	A	804	GOL	O1-C1-C2-O2
4	B	804	OUP	C15-C14-O19-C20
4	B	804	OUP	C21-C20-O19-C14
3	A	802	H4B	C7-C6-C9-C10
3	B	803	H4B	C7-C6-C9-C10
3	A	802	H4B	C7-C6-C9-O9
2	B	802	HEM	C4B-C3B-CAB-CBB
4	A	803	OUP	C13-C14-O19-C20
4	A	803	OUP	C15-C14-O19-C20
2	A	801	HEM	C3D-CAD-CBD-CGD
4	B	804	OUP	C14-C15-C17-N18
3	A	802	H4B	N5-C6-C9-O9
3	B	803	H4B	N5-C6-C9-O9
2	B	802	HEM	CAD-CBD-CGD-O2D
2	B	802	HEM	CAD-CBD-CGD-O1D
2	A	801	HEM	CAD-CBD-CGD-O2D

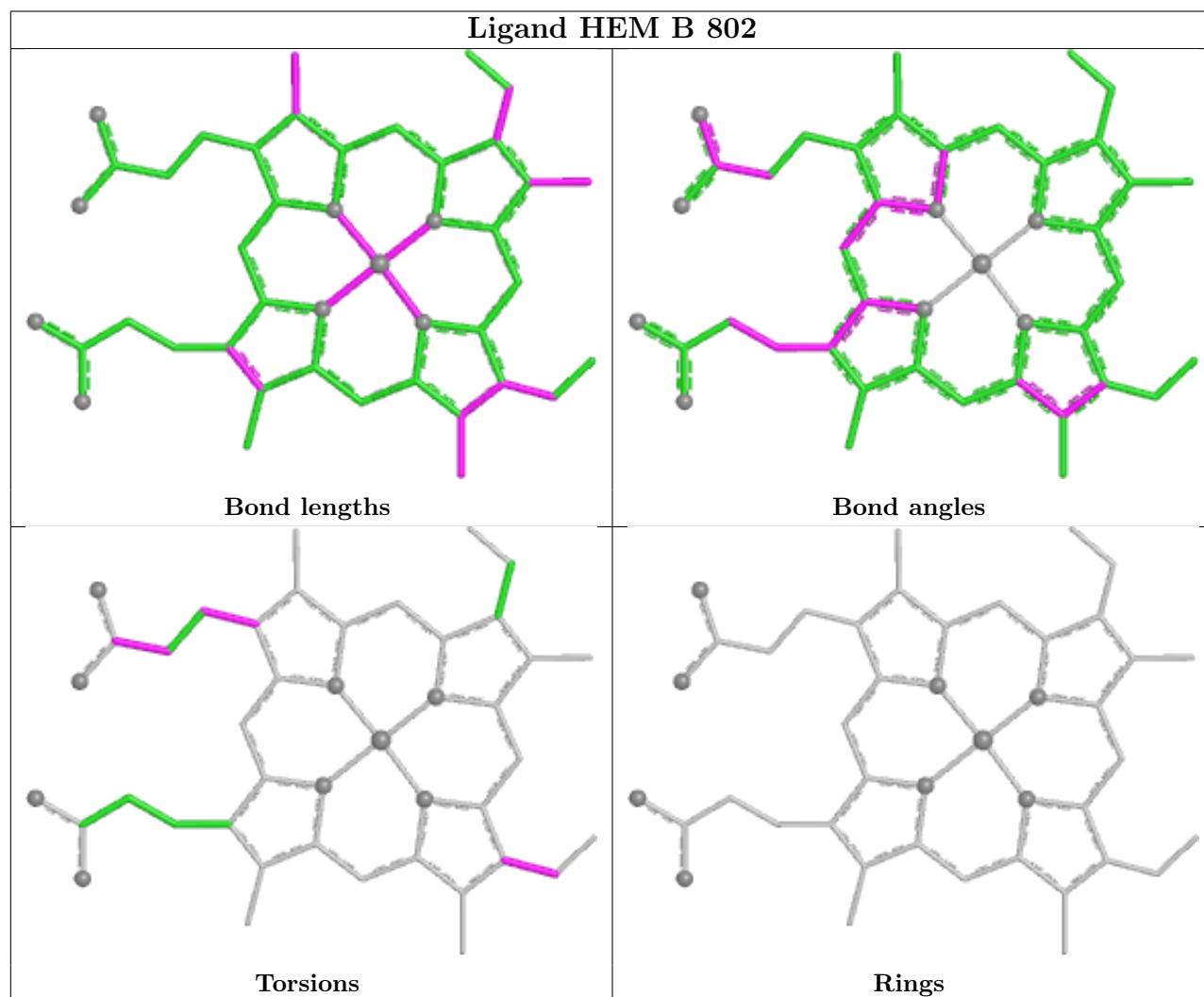
There are no ring outliers.

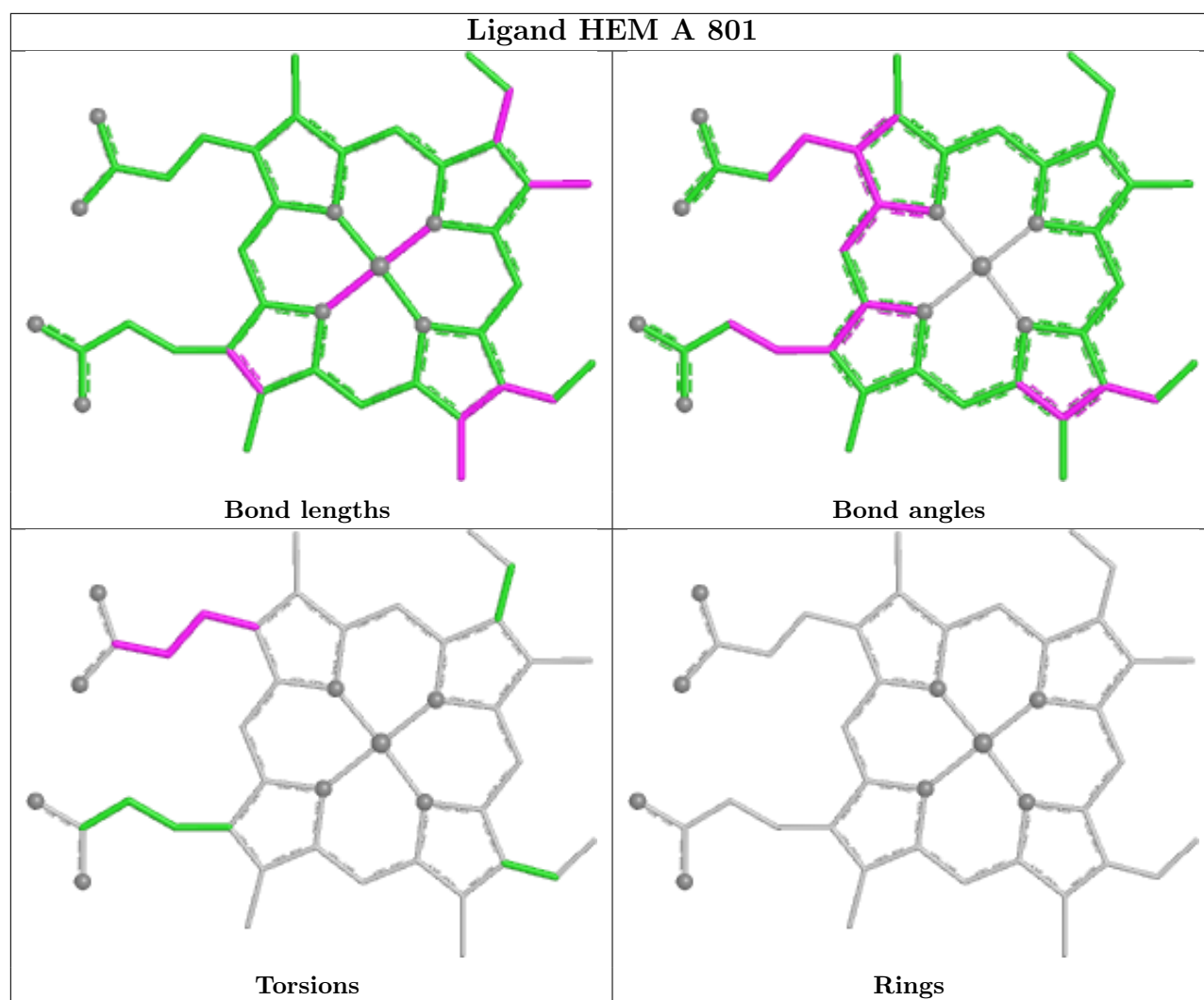
5 monomers are involved in 11 short contacts:

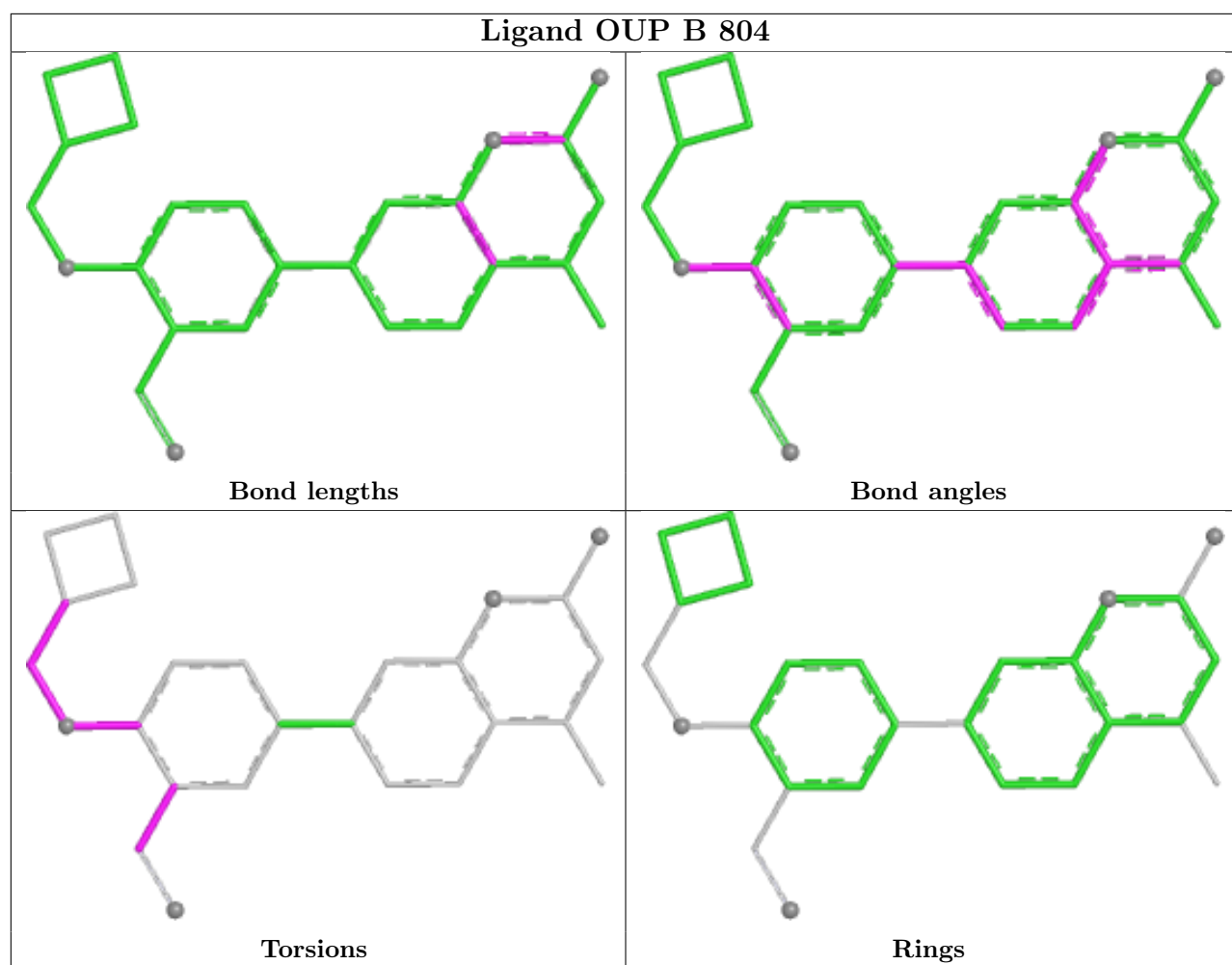
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	802	HEM	4	0
2	A	801	HEM	3	0
5	A	805	GOL	2	0
4	B	804	OUP	2	0
4	A	803	OUP	2	0

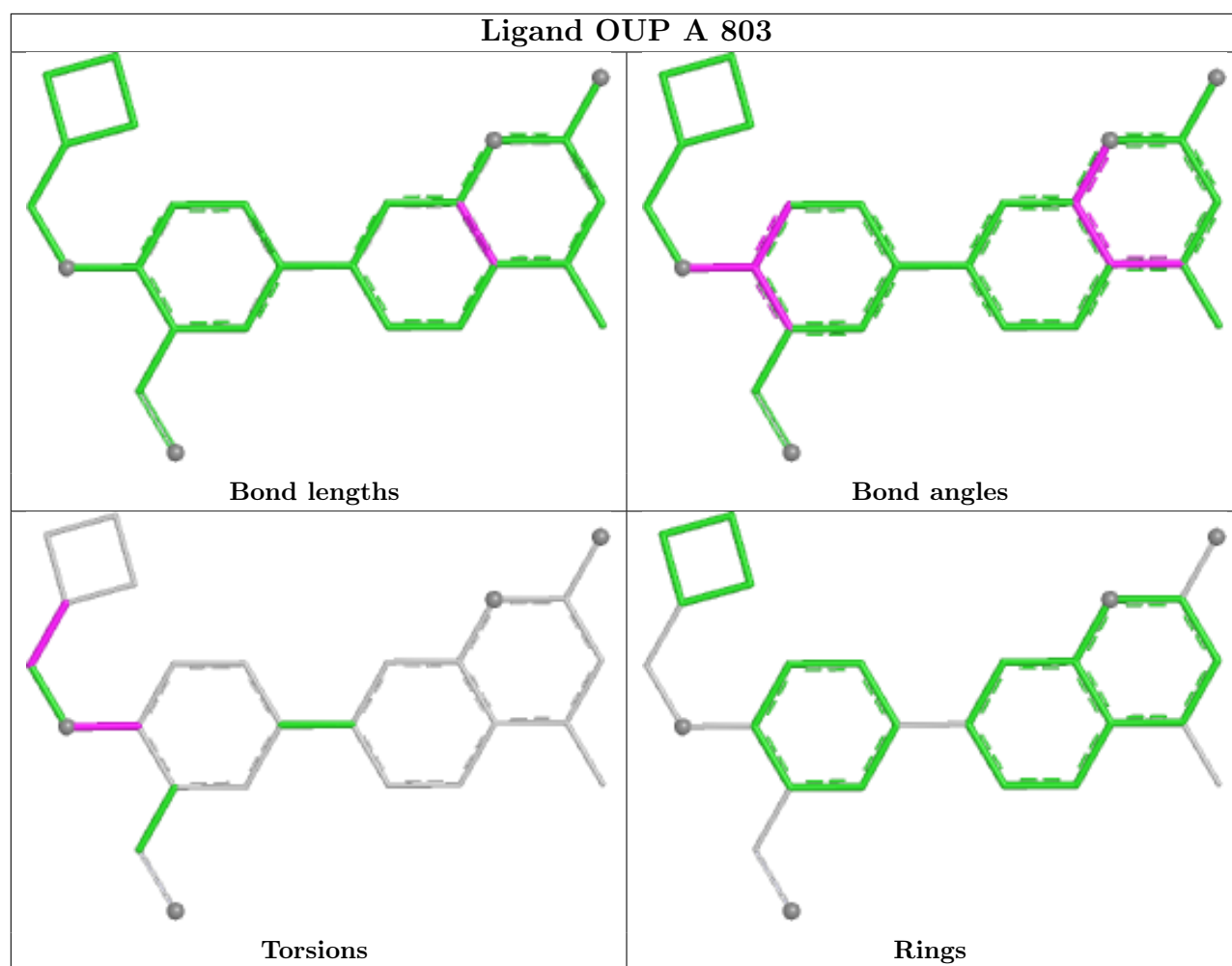
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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









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	417/421 (99%)	-0.04	13 (3%)	51 58	21, 41, 78, 129	5 (1%)
1	B	411/421 (97%)	0.14	16 (3%)	43 49	21, 47, 88, 121	3 (0%)
All	All	828/842 (98%)	0.05	29 (3%)	47 54	21, 44, 86, 129	8 (0%)

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	600	VAL	5.0
1	A	305	PHE	3.9
1	A	604	CYS	3.6
1	B	604	CYS	3.5
1	B	327	LEU	3.3
1	B	609	TYR	3.2
1	A	599	GLY	3.2
1	A	350	PRO	3.1
1	A	603	TYR	3.0
1	A	327	LEU	3.0
1	A	303	PRO	2.8
1	B	606	ASN	2.7
1	B	602	ASP	2.6
1	B	383	MET	2.5
1	B	720	VAL	2.5
1	A	302	CYS	2.5
1	A	341	MET	2.5
1	B	603	TYR	2.4
1	A	609	TYR	2.4
1	B	335	ILE	2.3
1	A	608	ARG	2.2
1	A	711	TYR	2.2
1	B	303	PRO	2.2
1	B	326	THR	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	354	ALA	2.1
1	B	344	SER	2.1
1	B	607	SER	2.0
1	B	342	HIS	2.0
1	A	353	VAL	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](#)

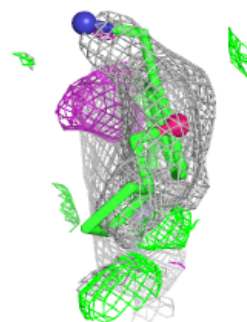
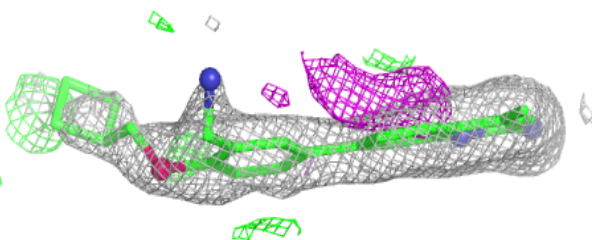
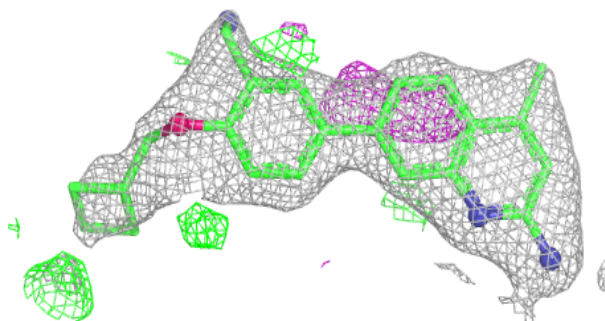
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	B	805	6/6	0.77	0.18	74,78,85,90	0
5	GOL	A	804	6/6	0.83	0.16	65,71,73,76	0
5	GOL	A	805	6/6	0.85	0.14	44,58,63,82	0
4	OUP	B	804	26/26	0.91	0.15	30,57,109,111	0
4	OUP	A	803	26/26	0.91	0.13	26,63,94,95	0
3	H4B	B	803	17/17	0.93	0.10	35,49,68,73	0
3	H4B	A	802	17/17	0.94	0.09	33,44,58,66	0
2	HEM	A	801	43/43	0.97	0.09	25,32,56,66	0
2	HEM	B	802	43/43	0.98	0.06	27,34,54,73	0
6	ZN	B	801	1/1	1.00	0.01	35,35,35,35	0

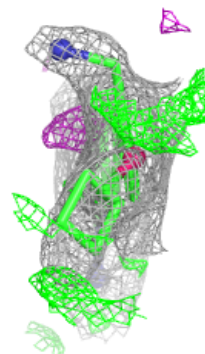
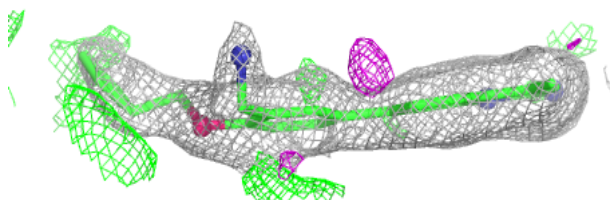
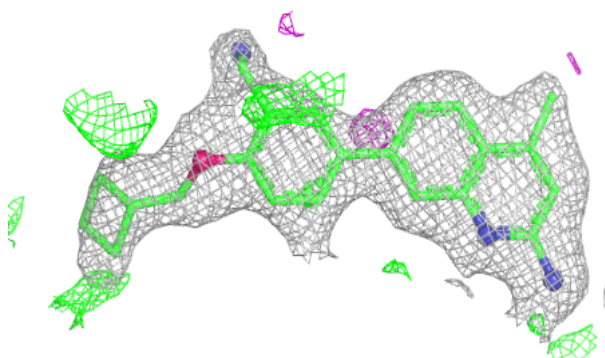
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around OUP B 804:

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

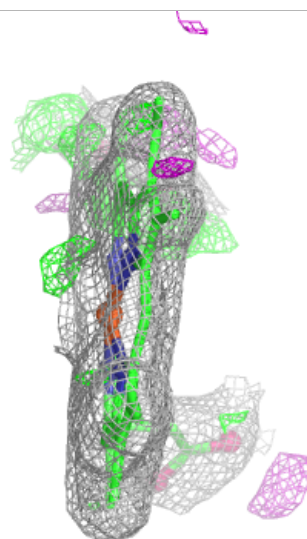
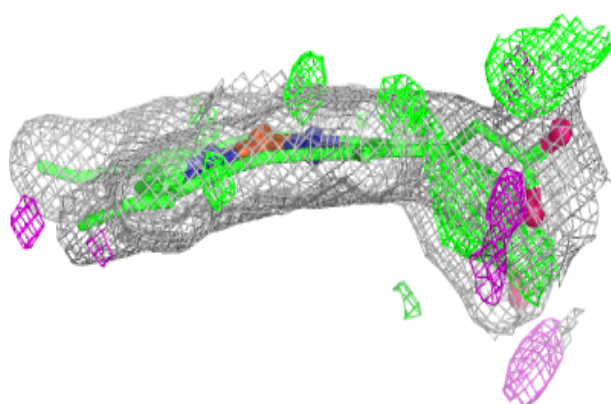
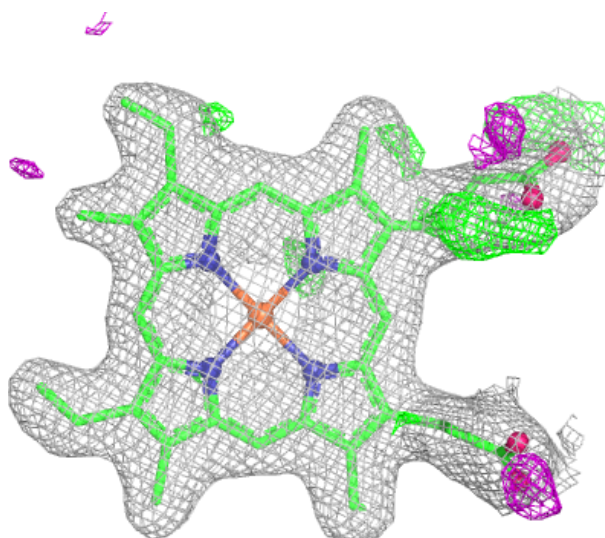
**Electron density around OUP A 803:**

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



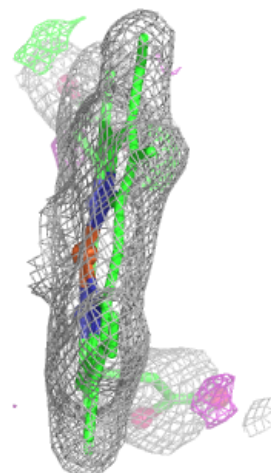
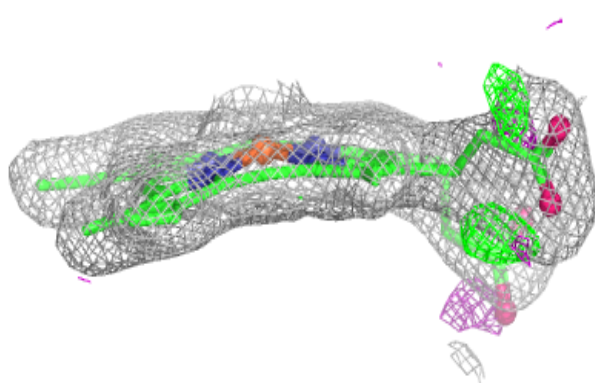
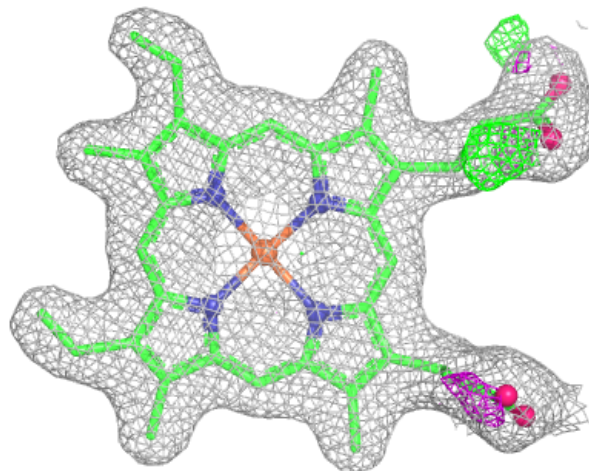
Electron density around HEM A 801:

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



Electron density around HEM B 802:

$2mF_o-DF_c$ (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.