Full wwPDB X-ray Structure Validation Report (i)

Oct 4, 2023 – 12:07 AM EDT

PDB ID	:	6POB
Title	:	Structure of human neuronal nitric oxide synthase R354A/G357D mutant heme domain in complex with 7-(3-(Aminomethyl)-4-(thiazol-4-ylmethoxy)phenyl)-4-methylquinolin-2-amine
Authors	:	Li, H.; Poulos, T.L.
Deposited on	:	2019-07-03
Resolution	:	1.95 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 (i) 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 (i)) were used in the production of this report:

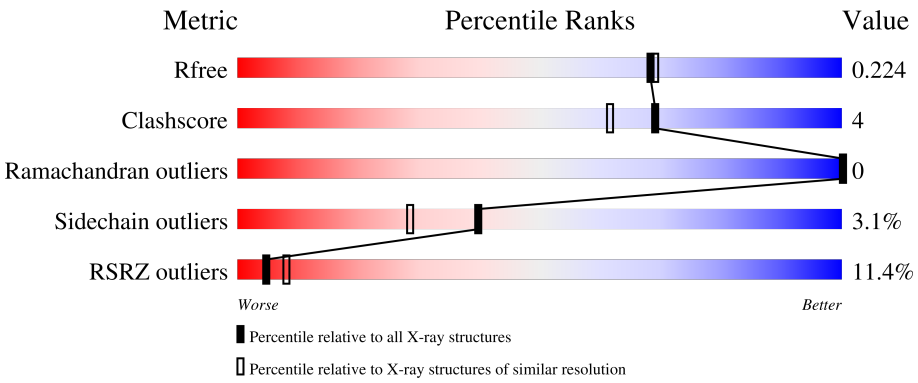
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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}	130704	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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 <=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	7% 90% 8% •
1	B	421	15% 85% 12% ••

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 7465 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.

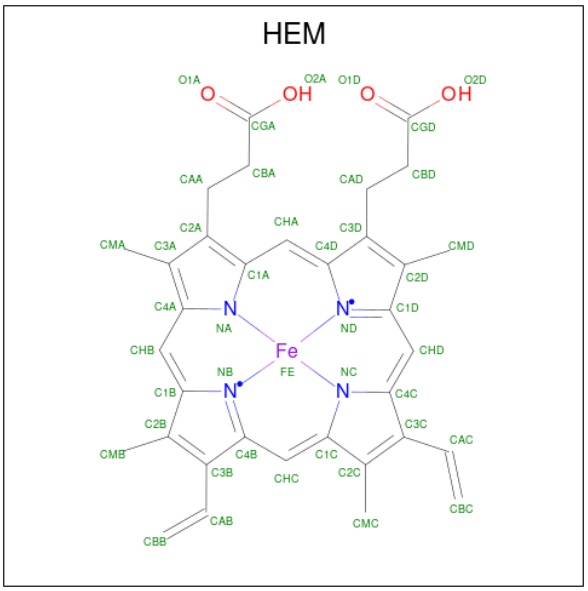
- Molecule 1 is a protein called Nitric oxide synthase, brain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	416	3405	2180	580	622	23	0	4	0
1	B	411	3362	2154	572	615	21	0	3	0

There are 4 discrepancies between the modelled and reference sequences:

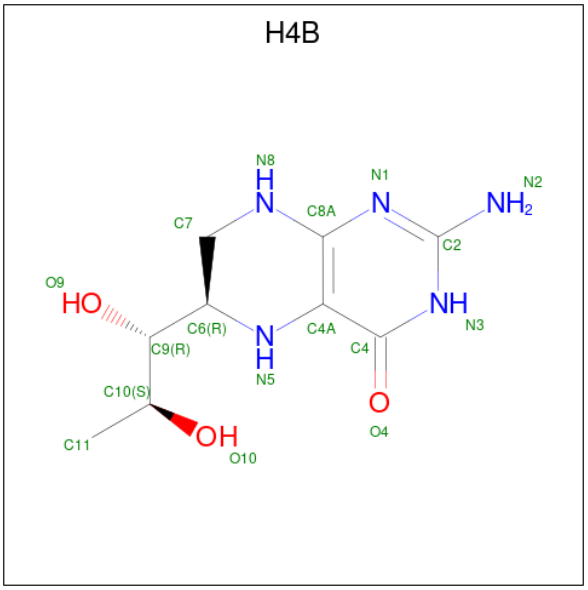
Chain	Residue	Modelled	Actual	Comment	Reference
A	354	ALA	ARG	engineered mutation	UNP P29475
A	357	ASP	GLY	engineered mutation	UNP P29475
B	354	ALA	ARG	engineered mutation	UNP P29475
B	357	ASP	GLY	engineered mutation	UNP P29475

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



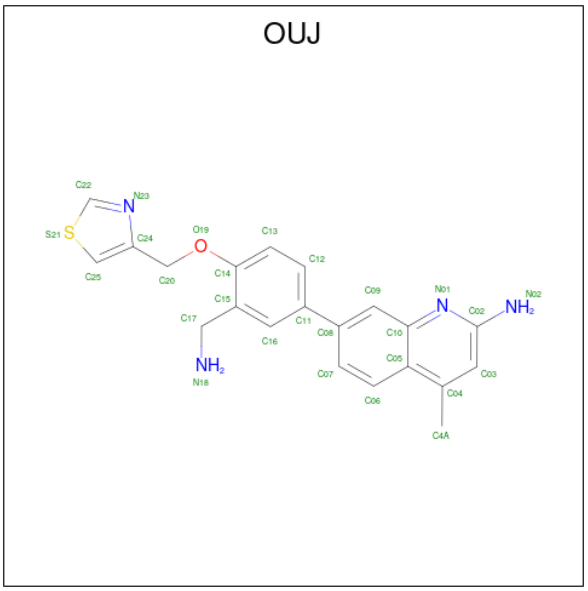
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Fe	N	O		
2	A	1	43	34	1	4	4	0	0
			Total	C	Fe	N	O		
2	B	1	43	34	1	4	4	0	0

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



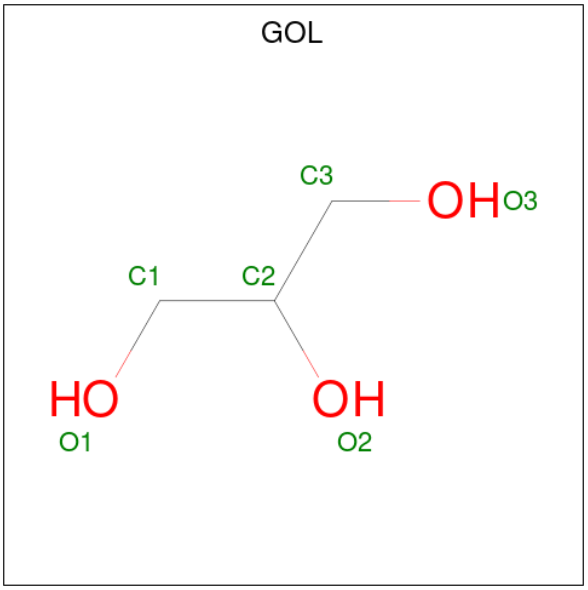
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O			
3	A	1	17	9	5	3		0	0
			Total	C	N	O			
3	B	1	17	9	5	3		0	0

- Molecule 4 is 7-{3-(aminomethyl)-4-[(1,3-thiazol-4-yl)methoxy]phenyl}-4-methylquinolin-2-amine (three-letter code: OUJ) (formula: $C_{21}H_{20}N_4OS$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
4	A	1	27	21	4	1	1	0	0
			Total	C	N	O	S		
4	B	1	27	21	4	1	1	0	0

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

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

- Molecule 6 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 7 is water.

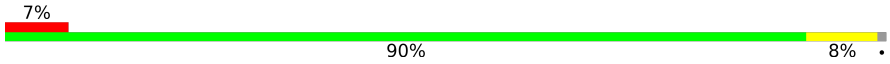
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
7	A	297	297	297	0	0
			Total	O		
7	B	208	208	208	0	0

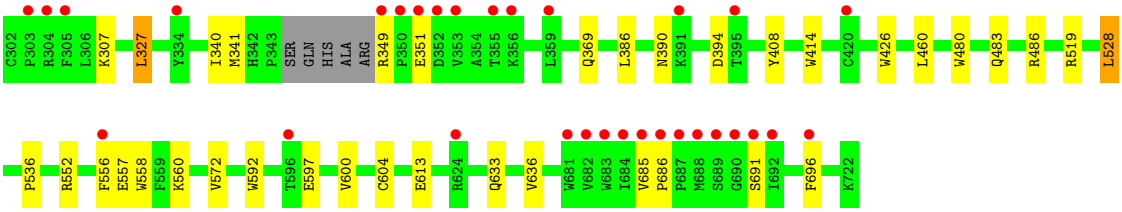
3 Residue-property plots (i)

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

- Molecule 1: Nitric oxide synthase, brain

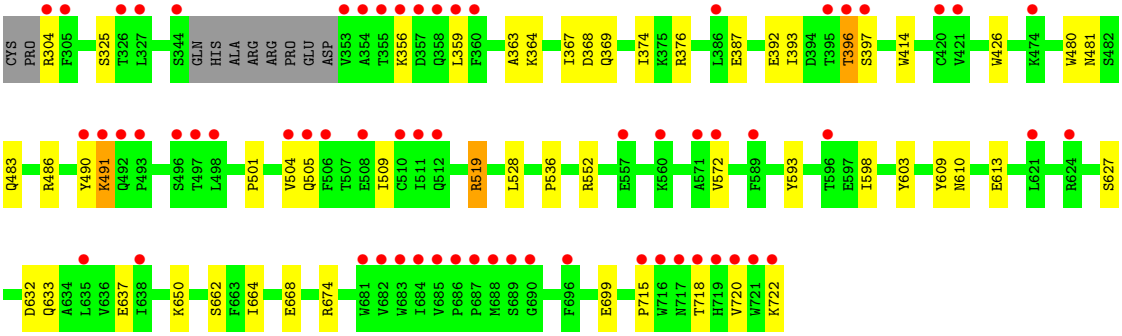
Chain A:





- Molecule 1: Nitric oxide synthase, brain

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	52.28Å 121.92Å 165.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.69 – 1.95 39.69 – 1.95	Depositor EDS
% Data completeness (in resolution range)	99.3 (39.69-1.95) 99.5 (39.69-1.95)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	0.11	Depositor
$< I/\sigma(I) >$ [1]	1.14 (at 1.95Å)	Xtriage
Refinement program	PHENIX (1.11.1-2575_1496: ???)	Depositor
R, R_{free}	0.179 , 0.222 0.181 , 0.224	Depositor DCC
R_{free} test set	3836 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å^2)	39.8	Xtriage
Anisotropy	0.799	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 50.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	7465	wwPDB-VP
Average B, all atoms (Å^2)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.24% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

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

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	0.36	0/3514	0.53	1/4769 (0.0%)
1	B	0.34	0/3466	0.49	0/4701
All	All	0.35	0/6980	0.51	1/9470 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	327	LEU	CA-CB-CG	6.58	130.44	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3405	0	3320	22	0
1	B	3362	0	3277	26	0
2	A	43	0	30	2	0
2	B	43	0	30	2	0
3	A	17	0	15	1	0
3	B	17	0	15	0	0
4	A	27	0	0	3	0
4	B	27	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	12	0	16	0	0
5	B	6	0	8	1	0
6	B	1	0	0	0	0
7	A	297	0	0	3	0
7	B	208	0	0	1	0
All	All	7465	0	6711	50	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 (50) 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:A:349:ARG:HD3	1:A:351:GLU:HB3	1.62	0.81
2:B:802:HEM:HHC	2:B:802:HEM:HBB2	1.63	0.80
1:B:491:LYS:HD3	1:B:491:LYS:H	1.52	0.71
1:A:528:LEU:HD22	1:A:536:PRO:HB2	1.71	0.71
2:B:802:HEM:HMC2	2:B:802:HEM:HBC2	1.73	0.69
1:A:557:GLU:OE2	1:A:560:LYS:NZ	2.28	0.66
1:B:662[B]:SER:OG	7:B:901:HOH:O	2.14	0.66
1:B:501:PRO:HA	1:B:504:VAL:HG23	1.78	0.66
1:B:392:GLU:O	1:B:396:THR:OG1	2.15	0.65
2:A:801:HEM:HBB2	2:A:801:HEM:HHC	1.79	0.65
2:A:801:HEM:HMC2	2:A:801:HEM:HBC2	1.78	0.64
1:A:483:GLN:HB2	1:A:486:ARG:HG3	1.79	0.64
1:B:664:ILE:O	1:B:668[B]:GLU:HG2	2.00	0.61
1:B:483:GLN:HB2	1:B:486:ARG:HG3	1.83	0.61
1:B:356:LYS:HG2	1:B:397:SER:HA	1.86	0.58
1:B:364:LYS:NZ	1:B:368:ASP:OD2	2.35	0.57
1:B:572:VAL:HG21	4:B:804:OUJ:C07	2.35	0.57
1:A:633:GLN:NE2	1:B:637:GLU:OE2	2.37	0.56
1:A:572:VAL:HG21	4:A:803:OUJ:C07	2.36	0.56
1:B:528:LEU:HD22	1:B:536:PRO:HB2	1.89	0.55
1:A:390:ASN:OD1	7:A:901:HOH:O	2.19	0.53
1:A:460:LEU:HD12	1:A:592:TRP:HB3	1.96	0.48
1:A:480:TRP:HB2	1:A:528:LEU:HB3	1.96	0.48
1:B:490:TYR:CE1	1:B:519:ARG:HA	2.49	0.48
1:B:668[B]:GLU:HG2	1:B:668[B]:GLU:H	1.58	0.46
1:B:603:TYR:HA	1:B:609:TYR:HB2	1.98	0.46
1:B:374:ILE:O	1:B:376:ARG:HG3	2.15	0.46
1:B:505:GLN:O	1:B:509:ILE:HG13	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:414:TRP:CE3	1:A:426:TRP:HA	2.51	0.45
1:A:341:MET:HE2	3:A:802:H4B:H9	1.99	0.44
1:B:359:LEU:HD23	1:B:393:ILE:HG23	1.99	0.44
1:B:480:TRP:CE2	1:B:715:PRO:HB2	2.52	0.44
4:A:803:OUJ:N18	7:A:903:HOH:O	2.36	0.44
1:B:593:TYR:CD1	1:B:598:ILE:HD11	2.53	0.44
1:A:340:ILE:HD13	1:B:699:GLU:HB3	1.99	0.43
1:A:369:GLN:HG2	1:A:408:TYR:OH	2.17	0.43
1:A:636:VAL:HG11	1:B:633:GLN:HG3	2.00	0.43
1:B:603:TYR:O	1:B:610:ASN:N	2.51	0.43
1:B:674:ARG:HE	5:B:805:GOL:H31	1.84	0.43
1:A:341:MET:HB2	1:A:341:MET:HE3	1.90	0.42
1:A:556:PHE:HB3	1:A:558:TRP:CE2	2.54	0.42
1:A:460:LEU:HD12	1:A:592:TRP:CB	2.50	0.41
1:A:691:SER:HA	1:A:696:PHE:CG	2.55	0.41
1:B:356:LYS:HE2	1:B:397:SER:OG	2.19	0.41
1:A:519:ARG:NH1	7:A:909:HOH:O	2.46	0.41
1:A:685:VAL:HA	1:A:686:PRO:HD3	1.93	0.41
1:B:363:ALA:O	1:B:367:ILE:HG12	2.21	0.41
1:A:600:VAL:O	1:A:604:CYS:HB2	2.21	0.40
1:A:597:GLU:OE2	4:A:803:OUJ:N02	2.55	0.40
1:B:414:TRP:CE3	1:B:426:TRP:HA	2.57	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	416/421 (99%)	410 (99%)	6 (1%)	0	100	100
1	B	410/421 (97%)	395 (96%)	15 (4%)	0	100	100
All	All	826/842 (98%)	805 (98%)	21 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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	376/376 (100%)	369 (98%)	7 (2%)	57	50
1	B	370/376 (98%)	354 (96%)	16 (4%)	29	16
All	All	746/752 (99%)	723 (97%)	23 (3%)	40	28

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

Mol	Chain	Res	Type
1	A	307	LYS
1	A	327	LEU
1	A	386	LEU
1	A	394	ASP
1	A	528	LEU
1	A	552	ARG
1	A	613	GLU
1	B	304	ARG
1	B	325	SER
1	B	369	GLN
1	B	387	GLU
1	B	396	THR
1	B	481	ASN
1	B	491	LYS
1	B	519	ARG
1	B	552	ARG
1	B	613	GLU
1	B	627	SER
1	B	632	ASP
1	B	650	LYS
1	B	718	THR
1	B	720	VAL
1	B	722	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are

no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

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	-	16,18,18	0.79	0	11,26,26	2.55	5 (45%)
4	OUJ	A	803	-	26,30,30	0.99	1 (3%)	36,42,42	1.90	9 (25%)
3	H4B	B	803	-	16,18,18	0.88	0	11,26,26	2.54	6 (54%)
5	GOL	A	805	-	5,5,5	0.38	0	5,5,5	0.23	0
5	GOL	B	805	-	5,5,5	0.38	0	5,5,5	0.23	0
2	HEM	A	801	1	41,50,50	1.51	5 (12%)	45,82,82	1.51	8 (17%)
2	HEM	B	802	1	41,50,50	1.56	7 (17%)	45,82,82	1.71	8 (17%)
5	GOL	A	804	-	5,5,5	0.39	0	5,5,5	0.40	0
4	OUJ	B	804	-	26,30,30	0.99	3 (11%)	36,42,42	1.93	10 (27%)

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	-	-	0/8/17/17	0/2/2/2
4	OUJ	A	803	-	-	0/9/11/11	0/4/4/4
3	H4B	B	803	-	-	0/8/17/17	0/2/2/2
5	GOL	A	805	-	-	2/4/4/4	-
5	GOL	B	805	-	-	2/4/4/4	-
2	HEM	A	801	1	-	1/12/54/54	-
2	HEM	B	802	1	-	2/12/54/54	-
5	GOL	A	804	-	-	4/4/4/4	-
4	OUJ	B	804	-	-	0/9/11/11	0/4/4/4

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	802	HEM	C3C-CAC	3.98	1.56	1.47
2	B	802	HEM	C3C-C2C	-3.93	1.34	1.40
2	A	801	HEM	C3C-C2C	-3.90	1.35	1.40
2	A	801	HEM	C3C-CAC	3.66	1.55	1.47
2	A	801	HEM	CAB-C3B	3.06	1.55	1.47
2	B	802	HEM	CAB-C3B	2.86	1.55	1.47
2	B	802	HEM	FE-NB	2.45	2.09	1.96
2	B	802	HEM	CMC-C2C	2.24	1.56	1.51
2	A	801	HEM	C3B-C2B	-2.21	1.32	1.37
2	B	802	HEM	CMB-C2B	2.19	1.55	1.50
4	B	804	OUJ	C02-N01	2.15	1.36	1.33
4	A	803	OUJ	C05-C10	-2.15	1.39	1.42
4	B	804	OUJ	C05-C10	-2.12	1.39	1.42
2	B	802	HEM	C3B-C2B	-2.08	1.33	1.37
2	A	801	HEM	CMB-C2B	2.06	1.55	1.50
4	B	804	OUJ	C25-S21	2.05	1.73	1.70

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	801	HEM	CBA-CAA-C2A	-5.60	103.06	112.62
3	A	802	H4B	C8A-C4A-C4	5.36	119.33	114.57
3	B	803	H4B	C8A-C4A-C4	5.27	119.25	114.57
4	A	803	OUJ	C20-C24-C25	-4.98	122.33	129.99
2	B	802	HEM	C4B-CHC-C1C	4.93	129.07	122.56
4	A	803	OUJ	C04-C05-C10	4.90	120.67	118.01
4	B	804	OUJ	C04-C05-C10	4.54	120.47	118.01
2	B	802	HEM	CBA-CAA-C2A	-4.53	104.89	112.62
4	B	804	OUJ	C05-C10-N01	-4.43	118.11	122.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	803	OUJ	C05-C10-N01	-4.26	118.29	122.81
4	B	804	OUJ	C20-C24-C25	-4.24	123.46	129.99
4	A	803	OUJ	O19-C14-C15	3.73	120.66	115.78
4	B	804	OUJ	O19-C14-C15	3.43	120.27	115.78
2	B	802	HEM	CMA-C3A-C4A	-3.43	123.19	128.46
3	A	802	H4B	C2-N3-C4	3.36	121.26	115.93
3	B	803	H4B	C2-N3-C4	3.35	121.26	115.93
3	A	802	H4B	N1-C2-N3	-3.07	120.60	125.42
3	B	803	H4B	N1-C2-N3	-2.91	120.86	125.42
4	B	804	OUJ	N02-C02-N01	2.87	120.63	118.26
2	A	801	HEM	C3B-C2B-C1B	2.71	108.50	106.49
2	B	802	HEM	C4D-ND-C1D	2.70	107.86	105.07
4	B	804	OUJ	C07-C08-C11	-2.69	116.70	121.36
2	A	801	HEM	CMC-C2C-C3C	2.62	129.58	124.68
3	B	803	H4B	C4-C4A-N5	2.58	121.28	119.12
2	B	802	HEM	CHD-C1D-ND	2.57	127.22	124.43
3	A	802	H4B	C2-N1-C8A	2.56	120.28	114.54
2	A	801	HEM	CMA-C3A-C4A	-2.50	124.63	128.46
2	B	802	HEM	C3D-C4D-ND	-2.48	107.41	110.17
4	B	804	OUJ	C03-C04-C05	2.39	120.14	117.78
4	A	803	OUJ	C06-C05-C04	-2.36	119.16	123.66
4	B	804	OUJ	C09-C10-N01	2.32	122.25	118.72
3	B	803	H4B	C2-N1-C8A	2.30	119.69	114.54
4	B	804	OUJ	C20-O19-C14	2.25	122.20	117.76
2	B	802	HEM	C3B-C2B-C1B	2.17	108.10	106.49
3	B	803	H4B	N2-C2-N3	2.15	120.60	117.25
4	B	804	OUJ	C06-C05-C04	-2.15	119.55	123.66
4	A	803	OUJ	C20-O19-C14	2.12	121.93	117.76
2	A	801	HEM	C4D-ND-C1D	2.10	107.24	105.07
2	A	801	HEM	C3D-C4D-ND	-2.10	107.83	110.17
4	A	803	OUJ	C25-S21-C22	2.06	96.57	92.37
2	A	801	HEM	CMA-C3A-C2A	2.04	128.79	124.94
4	A	803	OUJ	C08-C09-C10	-2.03	119.66	121.44
2	B	802	HEM	CMA-C3A-C2A	2.03	128.76	124.94
2	A	801	HEM	CAD-CBD-CGD	-2.02	109.25	113.60
4	A	803	OUJ	C07-C08-C11	-2.01	117.87	121.36
3	A	802	H4B	C4-C4A-N5	2.00	120.80	119.12

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	804	GOL	O1-C1-C2-C3
5	A	804	GOL	C1-C2-C3-O3
5	B	805	GOL	O1-C1-C2-C3
5	A	804	GOL	O1-C1-C2-O2
5	A	805	GOL	O1-C1-C2-C3
5	B	805	GOL	O1-C1-C2-O2
5	A	804	GOL	O2-C2-C3-O3
5	A	805	GOL	O1-C1-C2-O2
2	A	801	HEM	C4B-C3B-CAB-CBB
2	B	802	HEM	C4B-C3B-CAB-CBB
2	B	802	HEM	C2A-CAA-CBA-CGA

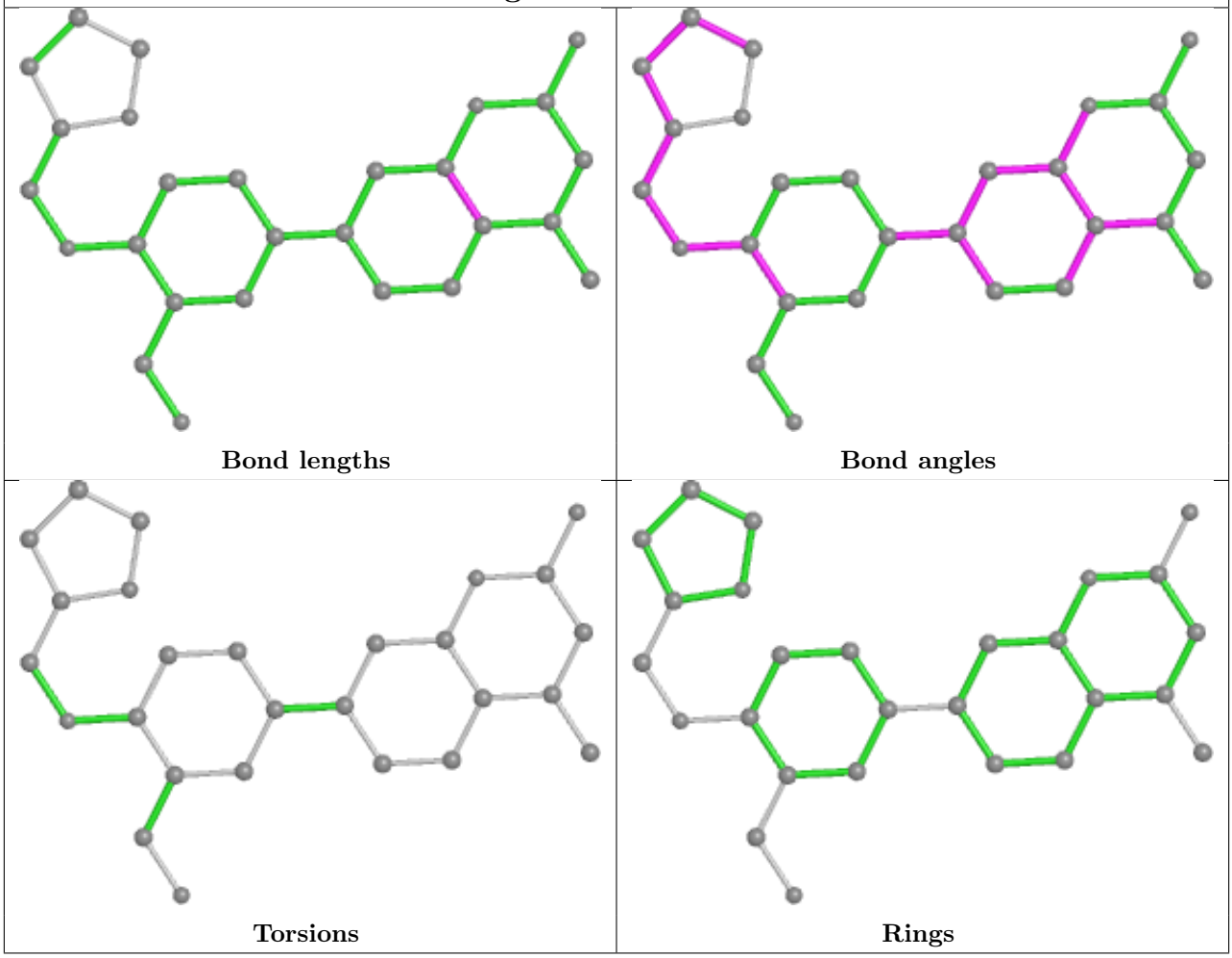
There are no ring outliers.

6 monomers are involved in 10 short contacts:

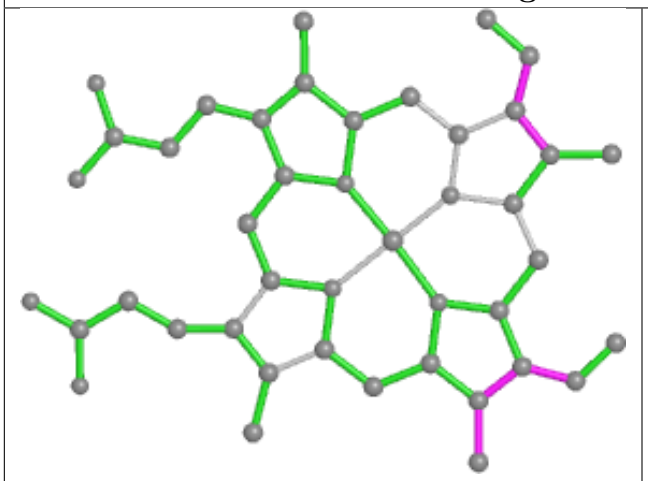
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	802	H4B	1	0
4	A	803	OUJ	3	0
5	B	805	GOL	1	0
2	A	801	HEM	2	0
2	B	802	HEM	2	0
4	B	804	OUJ	1	0

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 then 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.

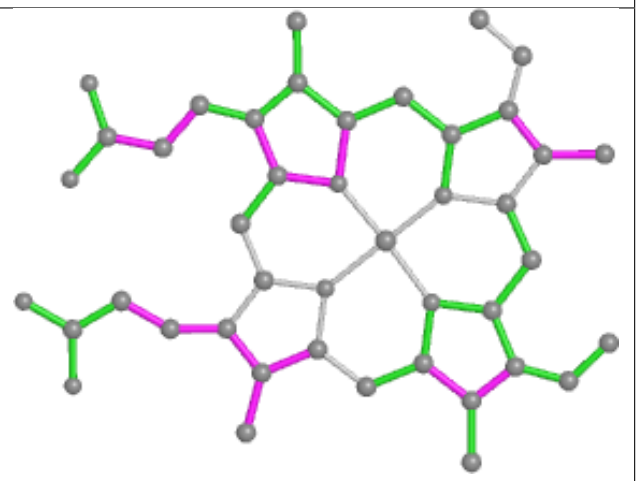
Ligand OUJ A 803



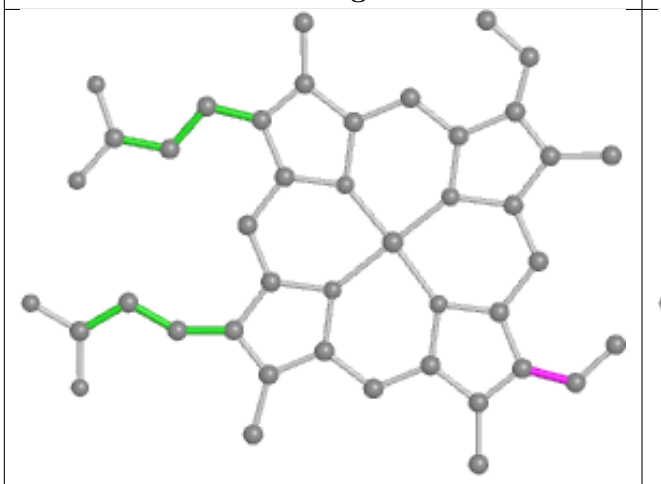
Ligand HEM A 801



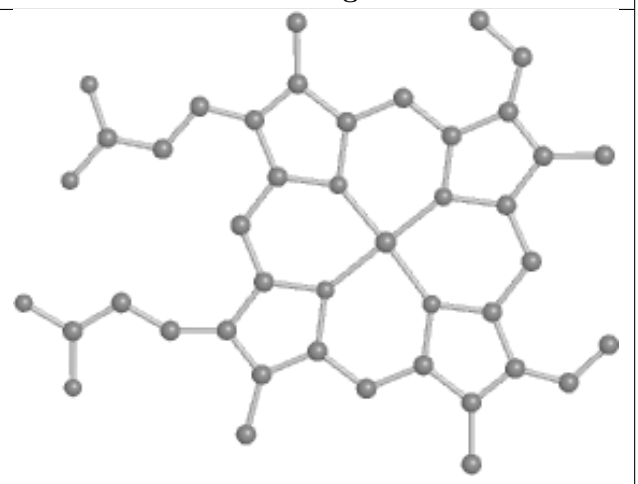
Bond lengths



Bond angles

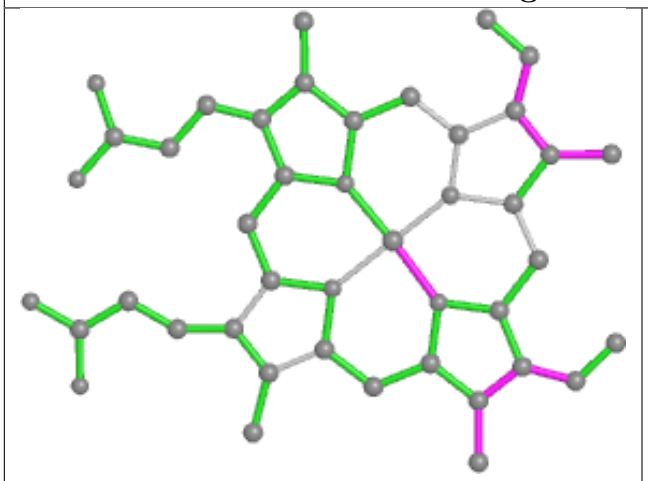


Torsions

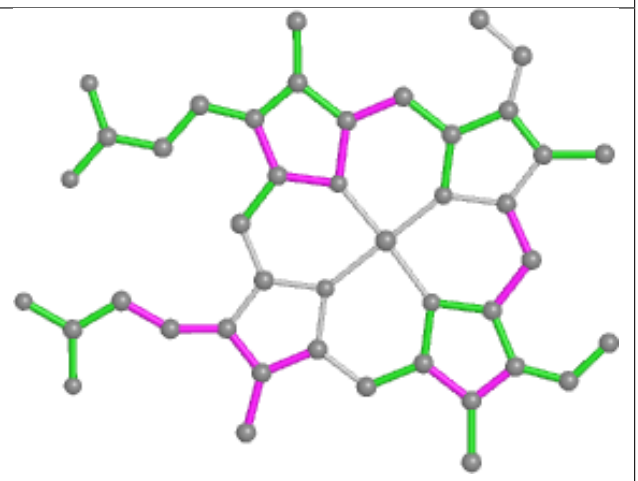


Rings

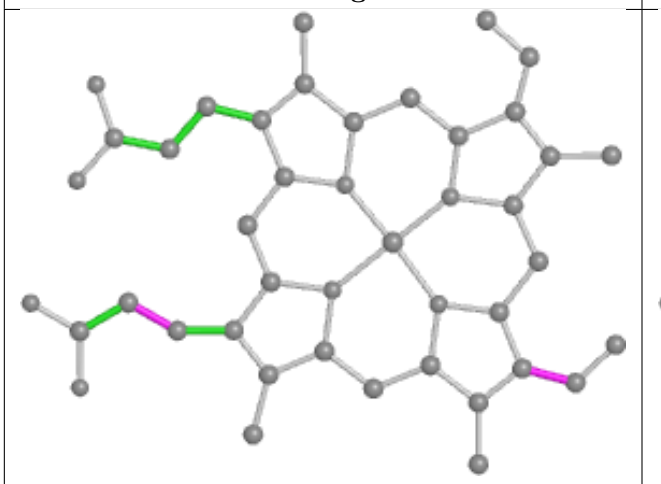
Ligand HEM B 802



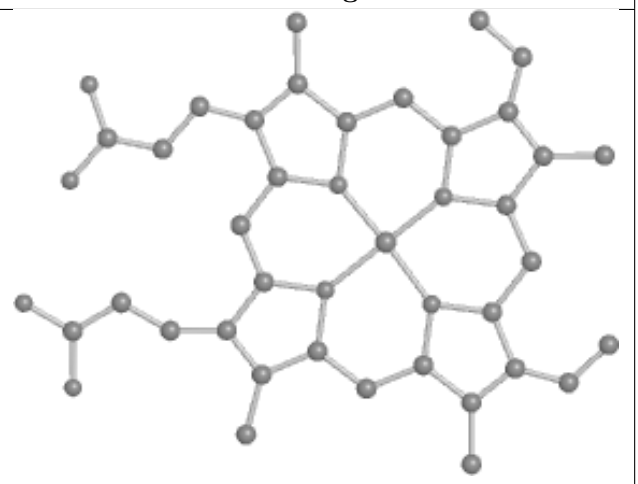
Bond lengths



Bond angles

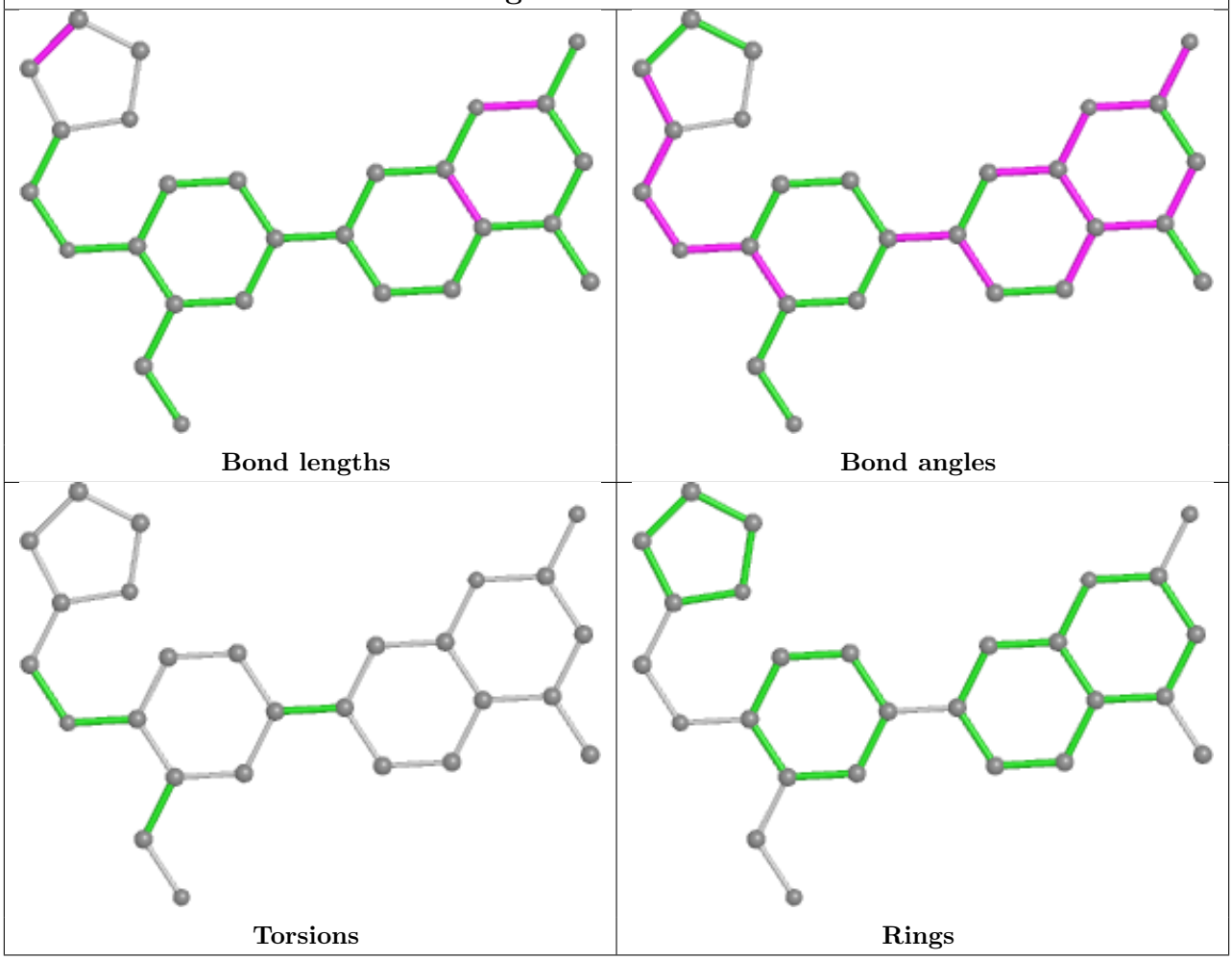


Torsions



Rings

Ligand OUJ B 804



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	416/421 (98%)	0.39	31 (7%)	14	22	33, 49, 81, 135	0
1	B	411/421 (97%)	0.83	63 (15%)	2	3	35, 60, 110, 140	0
All	All	827/842 (98%)	0.61	94 (11%)	5	8	33, 54, 103, 140	0

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	720	VAL	8.4
1	B	721	TRP	8.3
1	A	351	GLU	6.2
1	A	350	PRO	6.1
1	B	493	PRO	6.0
1	B	353	VAL	6.0
1	B	716	TRP	5.9
1	B	511	ILE	5.5
1	B	510	CYS	5.3
1	B	360	PHE	4.9
1	B	357	ASP	4.4
1	B	326	THR	4.4
1	B	683	TRP	4.4
1	A	685	VAL	4.4
1	B	504	VAL	4.3
1	B	508	GLU	4.3
1	B	719	HIS	4.2
1	B	686	PRO	4.2
1	B	718	THR	4.2
1	A	681	TRP	4.2
1	B	304	ARG	4.2
1	B	682	VAL	4.2
1	B	717	ASN	4.0
1	A	305	PHE	4.0

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Mol	Chain	Res	Type	RSRZ
1	A	696	PHE	4.0
1	B	722	LYS	3.8
1	A	682	VAL	3.8
1	B	359	LEU	3.6
1	A	353	VAL	3.6
1	B	397	SER	3.5
1	B	344	SER	3.4
1	A	684	ILE	3.4
1	B	685	VAL	3.4
1	A	349	ARG	3.3
1	B	396	THR	3.3
1	B	498	LEU	3.3
1	A	624	ARG	3.3
1	B	356	LYS	3.3
1	B	684	ILE	3.3
1	A	687	PRO	3.3
1	B	589	PHE	3.2
1	B	624	ARG	3.2
1	B	557	GLU	3.2
1	B	635	LEU	3.1
1	B	305	PHE	3.1
1	B	690	GLY	3.1
1	B	496	SER	3.1
1	A	686	PRO	3.1
1	A	690	GLY	3.0
1	B	474	LYS	2.9
1	B	358	GLN	2.8
1	A	352	ASP	2.8
1	B	687	PRO	2.8
1	B	688	MET	2.8
1	B	492	GLN	2.8
1	B	689	SER	2.8
1	B	420	CYS	2.8
1	A	304	ARG	2.7
1	B	560	LYS	2.7
1	B	490	TYR	2.7
1	B	497	THR	2.7
1	B	596	THR	2.7
1	B	715	PRO	2.7
1	B	681	TRP	2.6
1	A	355	THR	2.6
1	B	572	VAL	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	354	ALA	2.6
1	B	491	LYS	2.5
1	A	683	TRP	2.5
1	B	505	GLN	2.5
1	A	334	TYR	2.5
1	A	303	PRO	2.4
1	A	356	LYS	2.4
1	A	420	CYS	2.4
1	A	689	SER	2.3
1	B	327	LEU	2.3
1	A	395	THR	2.3
1	B	386	LEU	2.3
1	B	355	THR	2.2
1	B	395	THR	2.2
1	A	391	LYS	2.2
1	B	621	LEU	2.1
1	B	512	GLN	2.1
1	B	421	VAL	2.1
1	B	506	PHE	2.1
1	A	596	THR	2.1
1	A	692	ILE	2.1
1	A	359	LEU	2.1
1	B	696	PHE	2.1
1	A	691	SER	2.1
1	B	638	ILE	2.0
1	A	688	MET	2.0
1	A	556	PHE	2.0
1	B	571	ALA	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 monosaccharides 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,

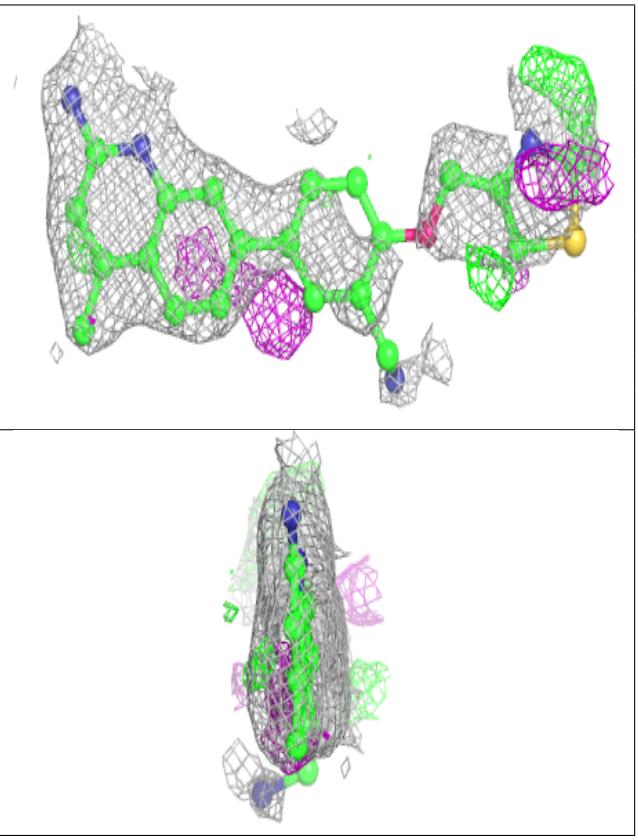
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	GOL	B	805	6/6	0.80	0.17	81,84,85,86	0
5	GOL	A	804	6/6	0.87	0.16	64,68,73,74	0
4	OUJ	A	803	27/27	0.87	0.25	36,85,126,129	0
4	OUJ	B	804	27/27	0.89	0.32	35,82,151,152	0
5	GOL	A	805	6/6	0.93	0.09	55,68,72,72	0
3	H4B	B	803	17/17	0.94	0.20	34,48,68,69	0
3	H4B	A	802	17/17	0.97	0.13	38,45,51,51	0
2	HEM	B	802	43/43	0.98	0.20	31,39,56,73	0
2	HEM	A	801	43/43	0.98	0.18	28,37,54,60	0
6	ZN	B	801	1/1	1.00	0.12	43,43,43,43	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.

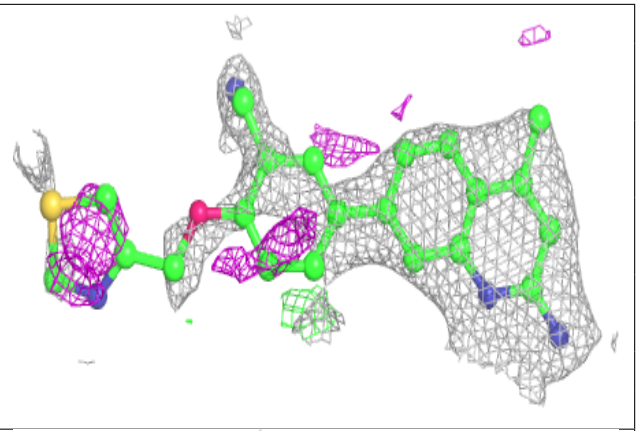
Electron density around OUJ A 803:

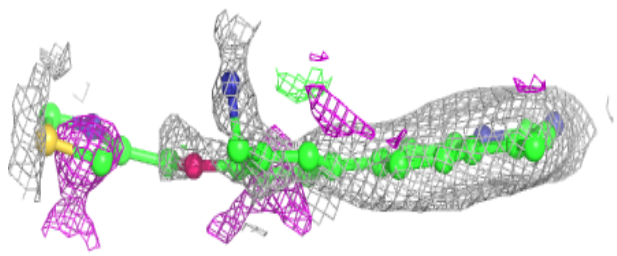
2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative) and green (positive)

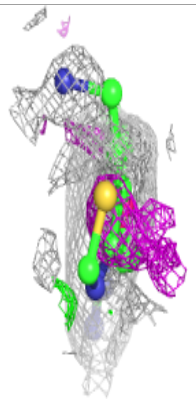


Electron density around OUJ B 804:

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

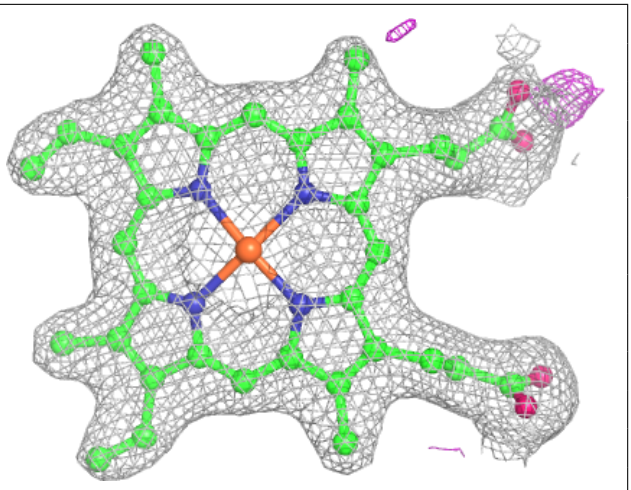


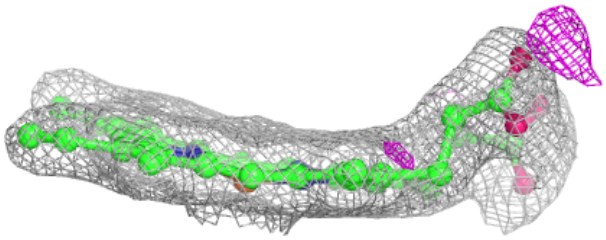


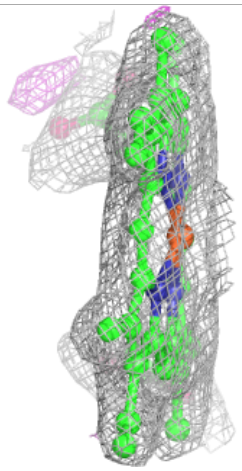


Electron density around HEM B 802:

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







Electron density around HEM A 801:

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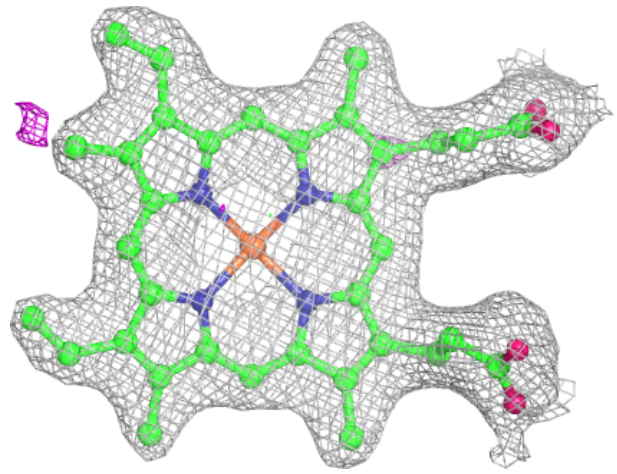
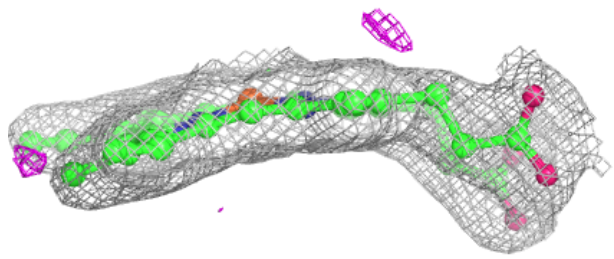
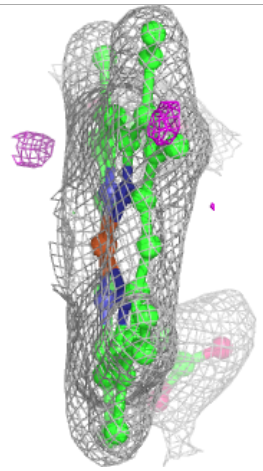
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
and green (positive)







6.5 Other polymers (i)

There are no such residues in this entry.