



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:27 PM UTC

PDB ID : 5AD4 / pdb_00005ad4
Title : Structure of rat neuronal nitric oxide synthase heme domain in complex with
7-((3-(2-(Dimethylamino)ethyl)phenoxy)methyl)quinolin-2- amine
Authors : Li, H.; Poulos, T.L.
Deposited on : 2015-08-20
Resolution : 1.98 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

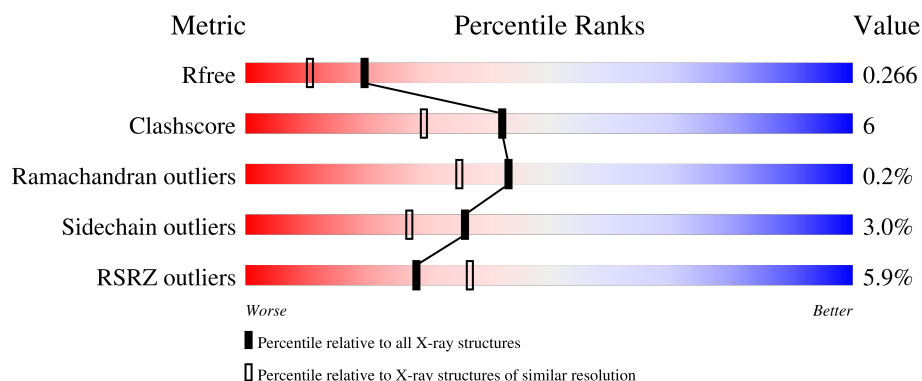
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.98 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	1506 (1.98-1.98)
Clashscore	190562	1534 (1.98-1.98)
Ramachandran outliers	187476	1518 (1.98-1.98)
Sidechain outliers	187428	1518 (1.98-1.98)
RSRZ outliers	180081	1506 (1.98-1.98)

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.

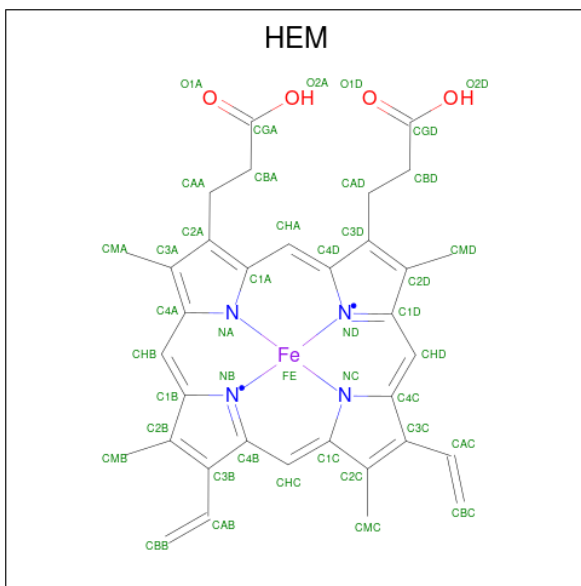
Mol	Chain	Length	Quality of chain
1	A	422	
1	B	422	

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	408	Total 3326	C 2129	N 567	O 608	S 22	0	3	1
1	B	411	Total 3357	C 2148	N 574	O 614	S 21	0	3	0

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



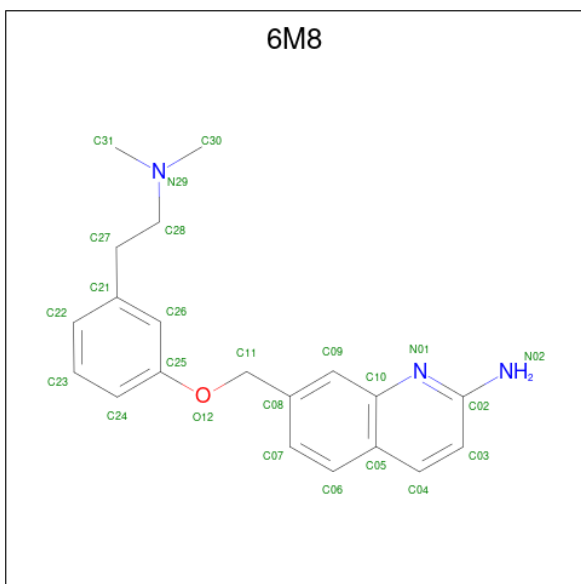
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 3 is 5,6,7,8-TETRAHYDROBIOPTERIN (CCD ID: H4B) (formula: $\text{C}_9\text{H}_{15}\text{N}_5\text{O}_3$).



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-[2-(dimethylamino)ethyl]phenoxy]methyl]quinolin-2-amine (CCD ID: 6M8) (formula: C₂₀H₂₃N₃O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			24	20	3	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	N	O	0	0
			24	20	3	1		

- Molecule 5 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		

- 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		

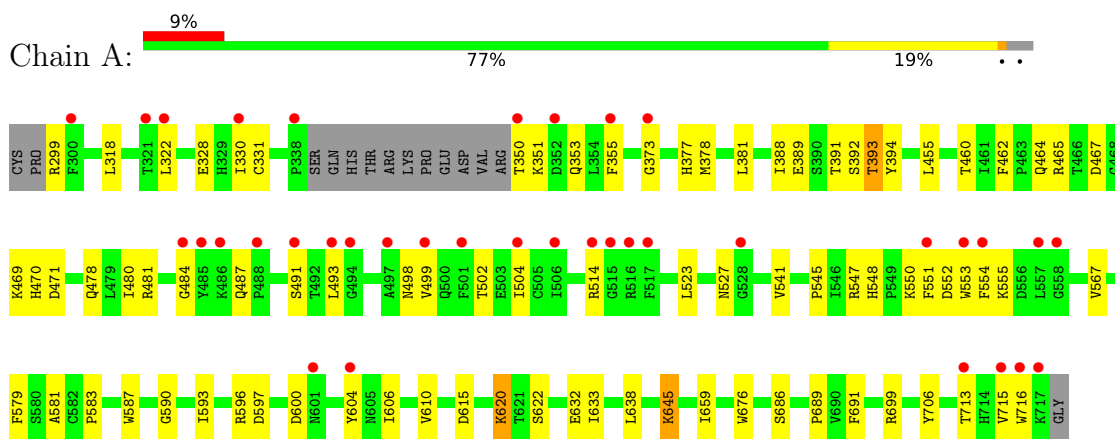
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	129	Total	O	0	0
			129	129		
7	B	159	Total	O	0	0
			159	159		

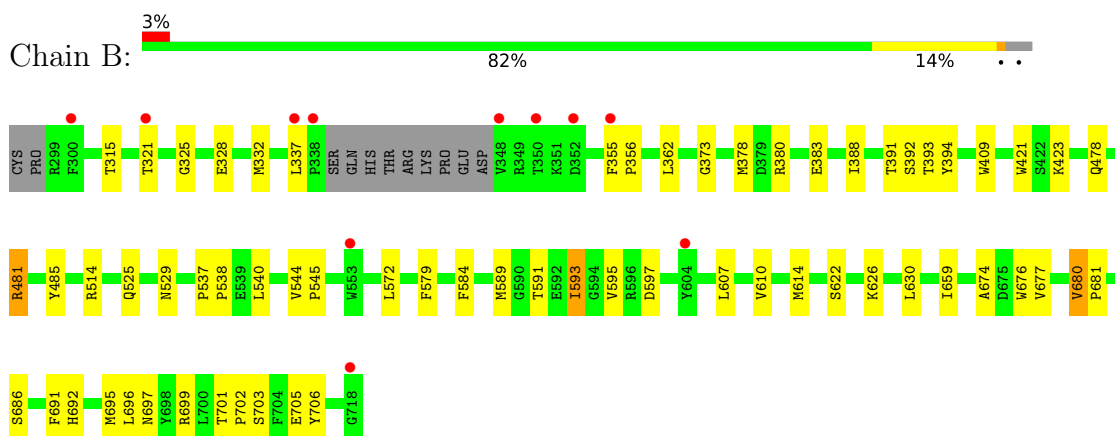
3 Residue-property plots

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	51.71Å 111.76Å 164.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.12 – 1.98 39.12 – 1.98	Depositor EDS
% Data completeness (in resolution range)	72.2 (39.12-1.98) 72.2 (39.12-1.98)	Depositor EDS
R_{merge}	0.24	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 1.98Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.202 , 0.266 0.202 , 0.266	Depositor DCC
R_{free} test set	2376 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	24.8	Xtriage
Anisotropy	0.053	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 52.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7148	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.61% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ACT, 6M8, H4B, ZN, HEM

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	0.47	1/3428 (0.0%)	0.88	0/4651
1	B	0.50	0/3459	0.86	2/4689 (0.0%)
All	All	0.48	1/6887 (0.0%)	0.87	2/9340 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	716	TRP	C-N	-6.98	1.23	1.33

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	680	VAL	CA-C-N	5.67	123.74	119.66
1	B	680	VAL	C-N-CA	5.67	123.74	119.66

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	3326	0	3237	43	0
1	B	3357	0	3275	36	0
2	A	43	0	30	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	43	0	30	5	0
3	A	17	0	15	0	0
3	B	17	0	15	0	0
4	A	24	0	23	3	0
4	B	24	0	23	5	0
5	A	4	0	3	0	0
5	B	4	0	3	0	0
6	B	1	0	0	0	0
7	A	129	0	0	2	1
7	B	159	0	0	2	1
All	All	7148	0	6654	85	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (85) 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:596:ARG:NH2	1:A:600:ASP:OD2	2.19	0.76
1:B:584:PHE:HE1	4:B:800:6M8:H06	1.50	0.76
2:B:750:HEM:HMC2	2:B:750:HEM:HBC2	1.70	0.74
2:B:750:HEM:HHC	2:B:750:HEM:HBB2	1.69	0.73
1:B:706:TYR:OH	2:B:750:HEM:O1D	2.07	0.70
2:A:750:HEM:HBB2	2:A:750:HEM:HHC	1.73	0.69
1:B:607:LEU:HD13	1:B:626:LYS:HG2	1.75	0.68
2:B:750:HEM:HBD1	4:B:800:6M8:H111	1.76	0.67
1:B:593:ILE:HA	1:B:597:ASP:HB2	1.77	0.66
1:A:478:GLN:HB2	1:A:481:ARG:HG3	1.78	0.66
1:A:393:THR:OG1	1:A:394:TYR:N	2.30	0.63
1:B:478:GLN:HB2	1:B:481:ARG:HG3	1.80	0.62
1:A:487:GLN:HE21	1:A:493:LEU:HB2	1.65	0.62
1:A:351:LYS:NZ	1:A:389:GLU:O	2.35	0.60
2:B:750:HEM:O2A	4:B:800:6M8:H24	2.01	0.59
7:A:2100:HOH:O	1:B:692:HIS:ND1	2.30	0.58
1:A:706:TYR:OH	2:A:750:HEM:O1D	2.18	0.57
1:A:455:LEU:HD12	1:A:587:TRP:HB3	1.87	0.57
1:A:567:VAL:HG21	4:A:800:6M8:C07	2.33	0.57
1:B:373:GLY:HA2	1:B:378:MET:HE2	1.87	0.57
1:A:355:PHE:HD1	1:A:388:ILE:HD12	1.70	0.56
1:A:593:ILE:HA	1:A:597:ASP:HB2	1.88	0.55
1:A:373:GLY:HA2	1:A:378:MET:HE3	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:330:ILE:HD11	1:B:696:LEU:HB3	1.89	0.55
1:B:699:ARG:NH2	1:B:705:GLU:OE1	2.42	0.52
1:A:465:ARG:NH2	1:A:471:ASP:OD2	2.35	0.52
1:B:572:LEU:HB3	1:B:579:PHE:HB2	1.92	0.52
1:B:393:THR:OG1	1:B:394:TYR:N	2.37	0.51
1:B:485:TYR:CE1	1:B:514:ARG:HA	2.47	0.49
1:B:610:VAL:O	1:B:614:MET:HG3	2.11	0.49
1:A:686:SER:HA	1:A:691:PHE:CG	2.47	0.49
2:A:750:HEM:HMC2	2:A:750:HEM:HBC2	1.93	0.49
1:A:484:GLY:O	1:A:499:VAL:HA	2.14	0.48
1:A:350:THR:N	1:A:353:GLN:OE1	2.48	0.47
1:B:674:ALA:HB3	1:B:695:MET:HB3	1.96	0.47
1:A:377:HIS:CD2	1:A:381:LEU:HD13	2.50	0.46
1:B:595:VAL:HG13	1:B:630:LEU:HD11	1.98	0.46
1:B:686:SER:HA	1:B:691:PHE:CG	2.51	0.46
1:B:391:THR:HG22	7:B:2027:HOH:O	2.15	0.46
1:A:462:PHE:HB2	1:A:581:ALA:HB3	1.97	0.46
1:A:659:ILE:HG13	1:A:689:PRO:HB2	1.97	0.46
1:A:470:HIS:HB3	1:A:527:ASN:OD1	2.16	0.46
1:A:590:GLY:HA3	1:A:638:LEU:HD21	1.98	0.45
1:A:596:ARG:O	1:A:600:ASP:HB2	2.16	0.45
1:B:485:TYR:CZ	1:B:514:ARG:HA	2.52	0.45
4:B:800:6M8:H282	4:B:800:6M8:H22	1.60	0.45
1:A:610:VAL:HG21	1:A:633:ILE:HD11	1.98	0.45
1:A:391:THR:O	1:A:392:SER:OG	2.29	0.45
1:A:331:CYS:HB3	1:B:697:ASN:HB3	1.99	0.45
1:A:551:PHE:HB3	1:A:553:TRP:NE1	2.32	0.44
1:A:467:ASP:OD1	1:A:470:HIS:ND1	2.51	0.44
1:A:676:TRP:CE3	1:B:677:VAL:HG22	2.53	0.44
1:A:322:LEU:HB2	1:A:699:ARG:HB2	1.99	0.44
1:A:498:ASN:O	1:A:502:THR:OG1	2.31	0.44
1:B:701:THR:HA	1:B:702:PRO:C	2.43	0.44
1:B:589:MET:HE3	1:B:591:THR:OG1	2.17	0.44
1:A:554:PHE:HB3	7:A:2086:HOH:O	2.17	0.44
1:A:464:GLN:HB3	1:A:579:PHE:CE1	2.53	0.43
1:A:455:LEU:HD12	1:A:587:TRP:CB	2.48	0.43
1:A:548:HIS:NE2	1:A:632:GLU:OE1	2.51	0.43
1:A:620:LYS:HB2	1:A:620:LYS:NZ	2.34	0.43
2:A:750:HEM:HBD1	4:A:800:6M8:H111	2.00	0.42
1:A:606:ILE:O	1:A:610:VAL:HG23	2.19	0.42
1:A:480:ILE:HD13	1:A:541:VAL:HG13	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:604:TYR:O	1:A:606:ILE:HG23	2.19	0.42
2:A:750:HEM:O2A	4:A:800:6M8:H24	2.19	0.42
1:B:380:ARG:NH1	1:B:383:GLU:OE1	2.51	0.42
1:B:409:TRP:CE3	1:B:421:TRP:HA	2.55	0.42
1:B:362:LEU:HA	1:B:362:LEU:HD23	1.83	0.41
1:B:537:PRO:HA	1:B:538:PRO:HD3	1.90	0.41
1:B:659:ILE:HD13	1:B:659:ILE:HA	1.88	0.41
1:A:467:ASP:C	1:A:469:LYS:H	2.28	0.41
1:B:388:ILE:O	1:B:392:SER:N	2.48	0.41
1:A:299:ARG:HG3	1:A:318:LEU:HD21	2.03	0.41
1:A:545:PRO:HG2	1:A:547:ARG:NH1	2.36	0.41
1:B:676:TRP:CE2	1:B:680:VAL:HG21	2.56	0.41
1:B:680:VAL:HA	1:B:681:PRO:HD3	1.86	0.41
1:A:645:LYS:HE3	1:A:645:LYS:HB2	1.89	0.41
1:B:584:PHE:CE1	4:B:800:6M8:H06	2.41	0.41
1:B:703:SER:HB2	7:B:2035:HOH:O	2.20	0.41
1:A:460:THR:O	1:A:583:PRO:HD2	2.21	0.40
1:B:325:GLY:O	1:B:332:MET:HE2	2.21	0.40
1:B:544:VAL:HA	1:B:545:PRO:HD3	1.93	0.40
1:B:355:PHE:N	1:B:356:PRO:HD2	2.36	0.40
1:B:525:GLN:HG3	1:B:529:ASN:O	2.21	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:2055:HOH:O	7:B:2056:HOH:O[2_554]	2.15	0.05

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	407/422 (96%)	387 (95%)	18 (4%)	2 (0%)	24	14
1	B	410/422 (97%)	394 (96%)	16 (4%)	0	100	100
All	All	817/844 (97%)	781 (96%)	34 (4%)	2 (0%)	43	35

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	514	ARG
1	A	491	SER

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	366/377 (97%)	353 (96%)	13 (4%)	31	20
1	B	369/377 (98%)	360 (98%)	9 (2%)	43	35
All	All	735/754 (98%)	713 (97%)	22 (3%)	36	27

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	328	GLU
1	A	393	THR
1	A	504	ILE
1	A	523	LEU
1	A	550	LYS
1	A	552	ASP
1	A	555	LYS
1	A	615	ASP
1	A	620	LYS
1	A	622	SER
1	A	645	LYS
1	A	713	THR
1	A	715	VAL
1	B	315	THR

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Mol	Chain	Res	Type
1	B	321	THR
1	B	328	GLU
1	B	337	LEU
1	B	423	LYS
1	B	481	ARG
1	B	540	LEU
1	B	593	ILE
1	B	622	SER

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	407	HIS
1	A	507	GLN
1	B	364	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 1 is monoatomic - leaving 8 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
2	HEM	B	750	1	50,50,50	1.90	12 (24%)	67,82,82	1.30	8 (11%)
5	ACT	A	860	-	3,3,3	0.81	0	3,3,3	0.78	0
5	ACT	B	860	-	3,3,3	0.79	0	3,3,3	0.88	0
3	H4B	B	760	-	17,18,18	0.71	0	14,26,26	1.98	6 (42%)
2	HEM	A	750	1	50,50,50	2.17	10 (20%)	67,82,82	1.53	10 (14%)
4	6M8	A	800	-	26,26,26	0.92	1 (3%)	34,35,35	0.87	1 (2%)
4	6M8	B	800	-	26,26,26	0.77	1 (3%)	34,35,35	0.99	1 (2%)
3	H4B	A	760	-	17,18,18	0.63	0	14,26,26	2.13	6 (42%)

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
2	HEM	B	750	1	-	4/14/54/54	-
3	H4B	B	760	-	-	2/8/17/17	0/2/2/2
2	HEM	A	750	1	-	3/14/54/54	-
4	6M8	A	800	-	-	4/10/10/10	0/3/3/3
4	6M8	B	800	-	-	2/10/10/10	0/3/3/3
3	H4B	A	760	-	-	0/8/17/17	0/2/2/2

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	750	HEM	C3D-C2D	7.95	1.53	1.36
2	B	750	HEM	C3D-C2D	7.40	1.52	1.36
2	A	750	HEM	FE-NB	5.78	2.12	1.94
2	A	750	HEM	FE-ND	5.70	2.12	1.94
2	A	750	HEM	FE-NA	5.64	2.13	1.95
2	B	750	HEM	FE-ND	4.80	2.09	1.94
2	A	750	HEM	FE-NC	3.85	2.07	1.95
2	B	750	HEM	FE-NB	3.82	2.06	1.94
2	B	750	HEM	FE-NC	3.73	2.07	1.95
2	B	750	HEM	FE-NA	3.73	2.07	1.95
2	A	750	HEM	CAB-C3B	3.05	1.55	1.47
2	B	750	HEM	CAC-C3C	2.86	1.55	1.47
2	A	750	HEM	CAC-C3C	2.85	1.55	1.47
2	B	750	HEM	CAB-C3B	2.84	1.55	1.47
4	A	800	6M8	C02-N01	2.62	1.36	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	750	HEM	CMC-C2C	2.55	1.56	1.50
2	A	750	HEM	CMA-C3A	2.46	1.55	1.50
2	A	750	HEM	CMC-C2C	2.45	1.55	1.50
2	B	750	HEM	CMD-C2D	2.38	1.55	1.50
2	B	750	HEM	CMA-C3A	2.20	1.55	1.50
4	B	800	6M8	C02-N01	2.16	1.36	1.33
2	A	750	HEM	C2A-C3A	-2.02	1.33	1.38
2	B	750	HEM	CMB-C2B	2.02	1.54	1.50
2	B	750	HEM	C2A-C3A	-2.02	1.33	1.38

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	750	HEM	C4D-ND-C1D	6.28	112.65	105.21
3	A	760	H4B	C2-N1-C8A	4.57	121.42	113.36
2	B	750	HEM	C4D-ND-C1D	4.51	110.55	105.21
3	B	760	H4B	C2-N1-C8A	4.17	120.73	113.36
3	A	760	H4B	O4-C4-C4A	-3.16	119.63	127.26
2	A	750	HEM	C3B-C2B-C1B	3.14	108.77	106.41
2	B	750	HEM	CHA-C4D-ND	3.11	128.22	124.37
3	A	760	H4B	C4A-C4-N3	3.03	120.46	112.13
3	A	760	H4B	C2-N3-C4	-2.89	119.87	125.11
3	B	760	H4B	O4-C4-C4A	-2.85	120.39	127.26
3	B	760	H4B	C4A-C4-N3	2.80	119.83	112.13
3	B	760	H4B	C2-N3-C4	-2.76	120.11	125.11
2	A	750	HEM	CBD-CAD-C3D	-2.72	105.01	112.53
2	A	750	HEM	CHA-C4D-ND	2.67	127.66	124.37
2	A	750	HEM	CHD-C1D-ND	2.59	127.21	124.42
2	B	750	HEM	CAA-CBA-CGA	-2.57	106.85	113.67
2	A	750	HEM	C1B-NB-C4B	2.36	108.00	105.21
3	A	760	H4B	N2-C2-N3	2.35	121.73	116.76
3	A	760	H4B	C4-C4A-N5	2.32	122.60	116.27
2	A	750	HEM	C3D-C4D-ND	-2.31	107.64	110.17
4	A	800	6M8	C05-C10-N01	-2.30	119.09	122.25
2	B	750	HEM	C2A-C1A-NA	-2.28	107.62	110.15
2	B	750	HEM	C1D-C2D-C3D	-2.26	104.60	106.98
2	A	750	HEM	CBA-CAA-C2A	-2.26	106.30	112.53
2	A	750	HEM	C3B-C4B-NB	-2.24	107.86	109.47
2	A	750	HEM	O1D-CGD-CBD	-2.23	116.02	123.09
3	B	760	H4B	C4-C4A-N5	2.23	122.34	116.27
2	B	750	HEM	O1D-CGD-CBD	-2.20	116.11	123.09
2	B	750	HEM	CBA-CAA-C2A	-2.17	106.55	112.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	800	6M8	N02-C02-N01	2.16	120.02	118.24
2	B	750	HEM	C3B-C2B-C1B	2.15	108.02	106.41
3	B	760	H4B	N2-C2-N3	2.09	121.18	116.76

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	750	HEM	C2A-CAA-CBA-CGA
2	B	750	HEM	C2A-CAA-CBA-CGA
4	B	800	6M8	C22-C21-C27-C28
4	B	800	6M8	C26-C21-C27-C28
2	A	750	HEM	C4B-C3B-CAB-CBB
2	B	750	HEM	C4B-C3B-CAB-CBB
4	A	800	6M8	C24-C25-O12-C11
4	A	800	6M8	C26-C25-O12-C11
3	B	760	H4B	C11-C10-C9-O9
3	B	760	H4B	O10-C10-C9-O9
4	A	800	6M8	C22-C21-C27-C28
2	B	750	HEM	CAD-CBD-CGD-O1D
4	A	800	6M8	C26-C21-C27-C28
2	B	750	HEM	CAD-CBD-CGD-O2D
2	A	750	HEM	CAD-CBD-CGD-O2D

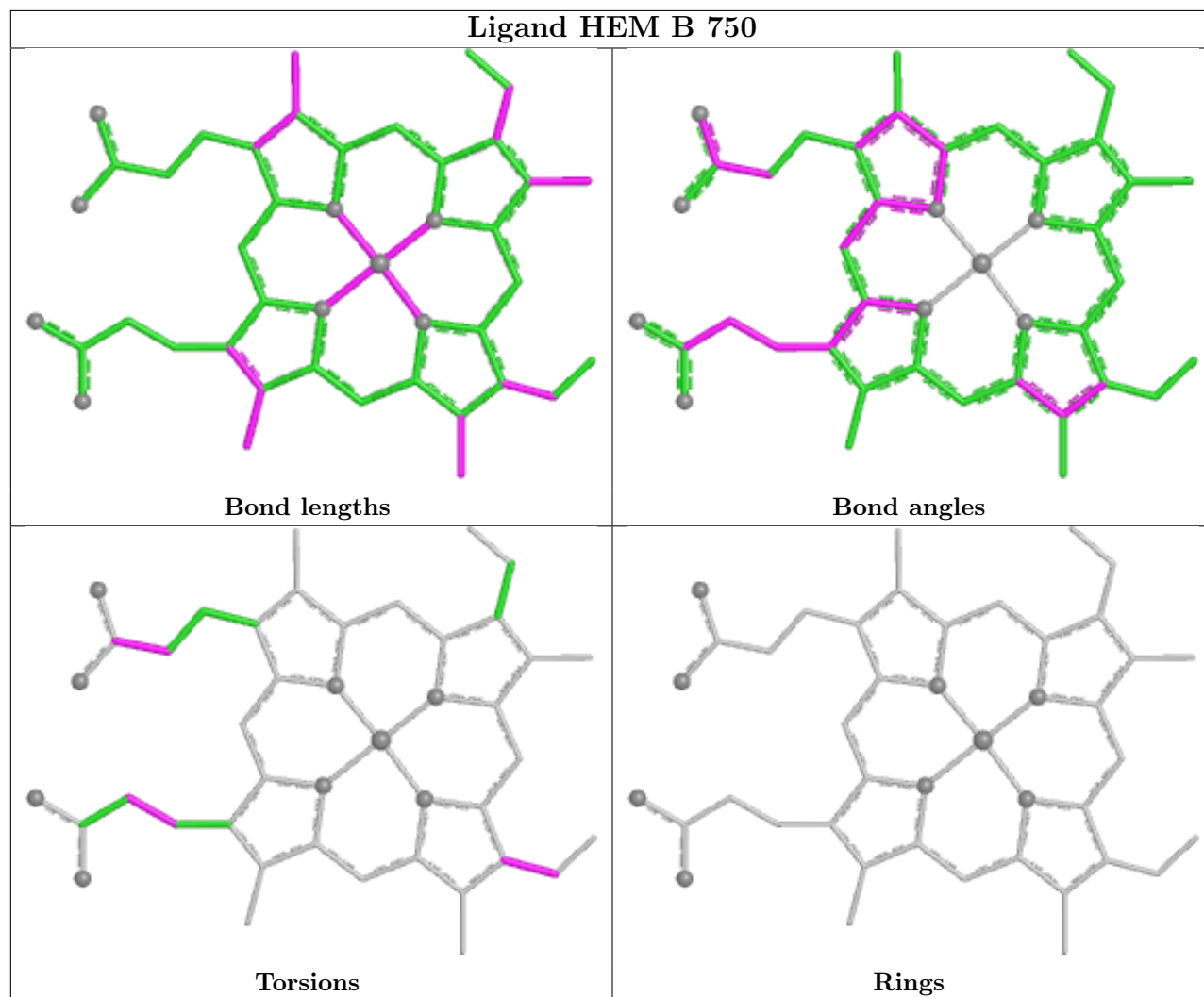
There are no ring outliers.

4 monomers are involved in 14 short contacts:

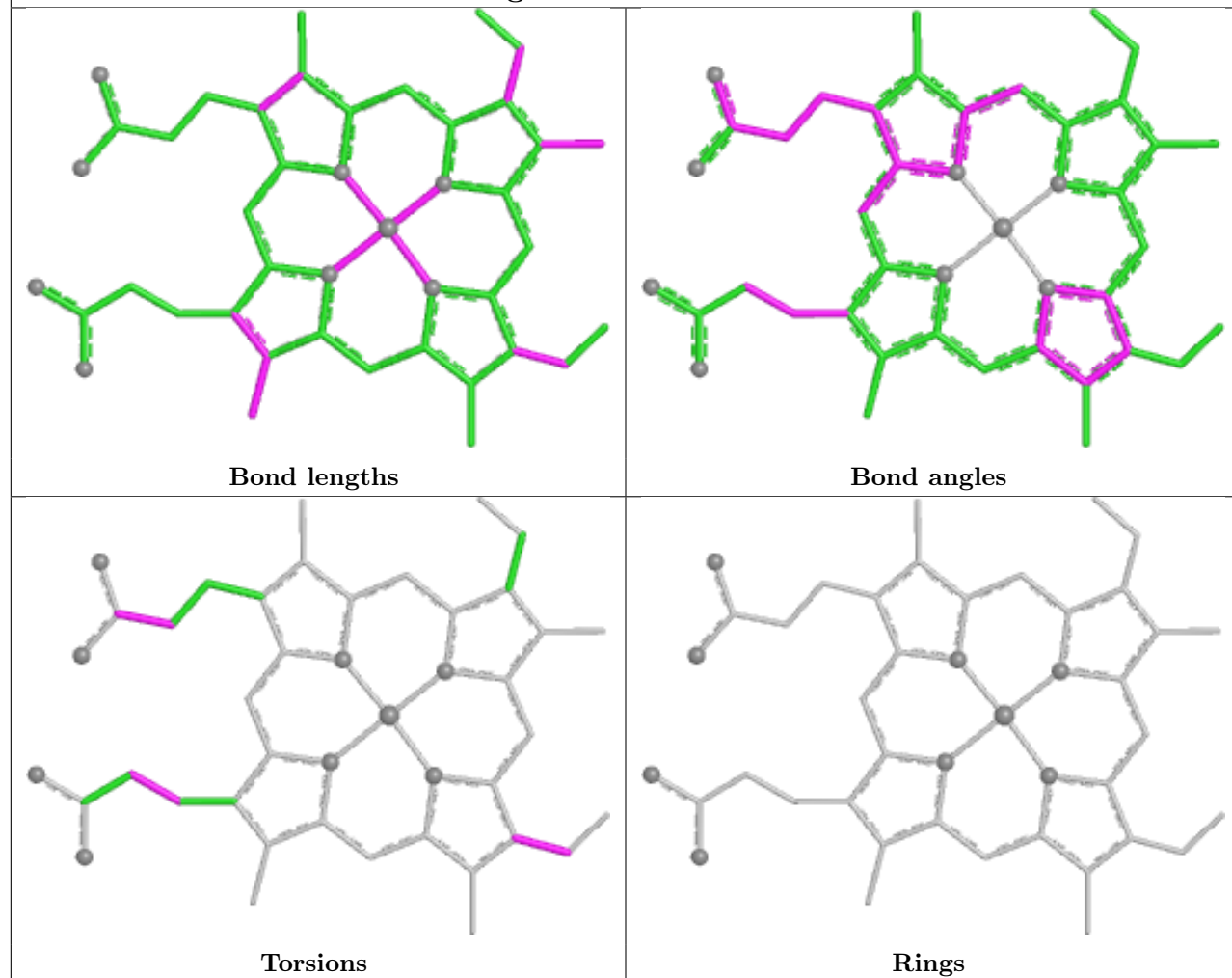
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	750	HEM	5	0
2	A	750	HEM	5	0
4	A	800	6M8	3	0
4	B	800	6M8	5	0

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

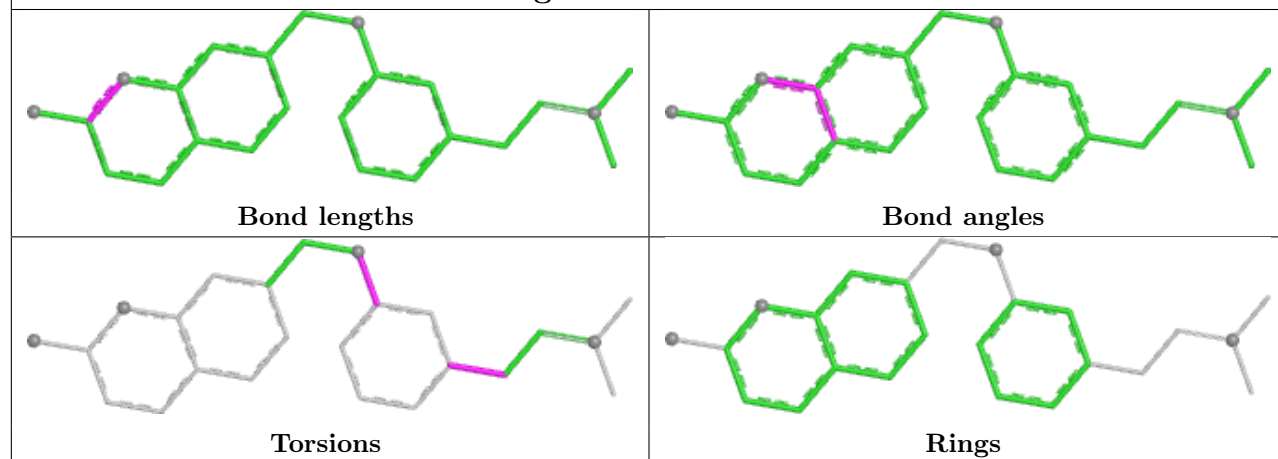
any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

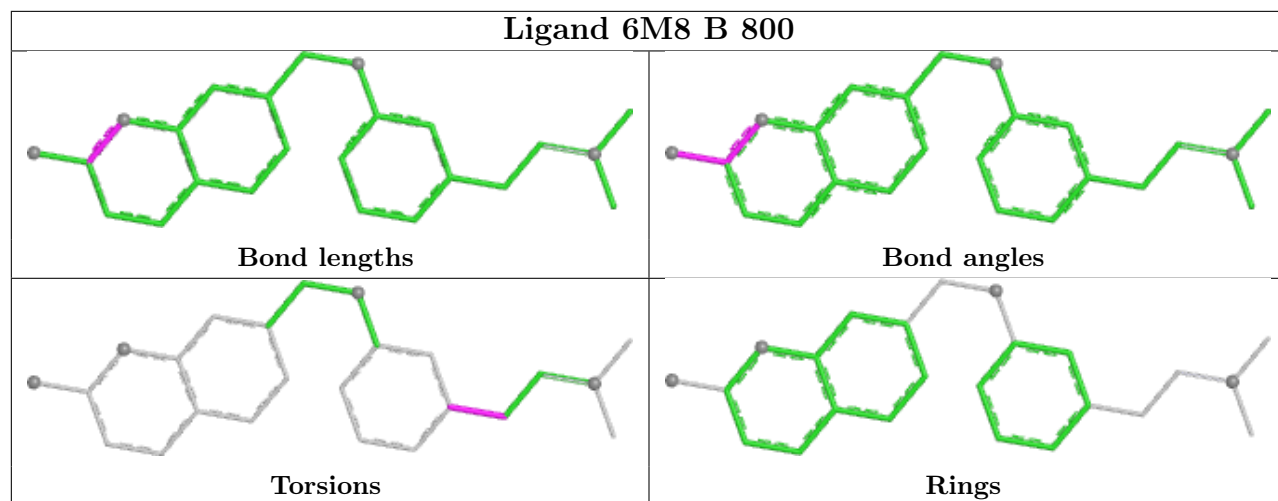


Ligand HEM A 750



Ligand 6M8 A 800





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	408/422 (96%)	0.44	37 (9%) 15 20	12, 35, 76, 104	3 (0%)
1	B	411/422 (97%)	-0.06	11 (2%) 56 66	9, 27, 58, 83	3 (0%)
All	All	819/844 (97%)	0.19	48 (5%) 28 37	9, 31, 72, 104	6 (0%)

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	338	PRO	4.4
1	B	348	VAL	4.3
1	A	551	PHE	4.2
1	B	338	PRO	4.1
1	A	715	VAL	3.8
1	A	493	LEU	3.7
1	A	717	LYS	3.5
1	A	321	THR	3.5
1	A	554	PHE	3.5
1	A	300	PHE	3.4
1	A	716	TRP	3.3
1	A	485	TYR	3.3
1	B	352	ASP	3.3
1	A	604	TYR	3.2
1	A	322	LEU	3.2
1	A	355	PHE	3.2
1	B	300	PHE	3.1
1	A	515	GLY	3.1
1	B	718	GLY	3.0
1	A	488	PRO	2.9
1	A	373	GLY	2.8
1	A	350	THR	2.8
1	A	553	TRP	2.8
1	A	484	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	321	THR	2.6
1	A	558	GLY	2.6
1	A	499	VAL	2.5
1	A	514	ARG	2.5
1	A	486	LYS	2.5
1	A	494	GLY	2.4
1	A	517	PHE	2.4
1	A	504	ILE	2.4
1	A	557	LEU	2.4
1	A	713	THR	2.4
1	A	516	ARG	2.3
1	A	352	ASP	2.3
1	B	604	TYR	2.3
1	A	501	PHE	2.2
1	A	506	ILE	2.2
1	A	601	ASN	2.2
1	B	337	LEU	2.2
1	B	355	PHE	2.1
1	A	528	GLY	2.1
1	B	350	THR	2.1
1	B	553	TRP	2.1
1	A	497	ALA	2.1
1	A	491	SER	2.0
1	A	330	ILE	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](#)

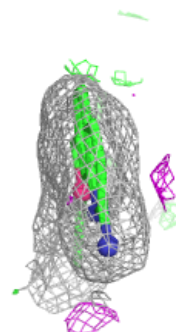
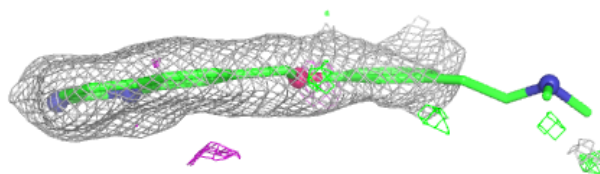
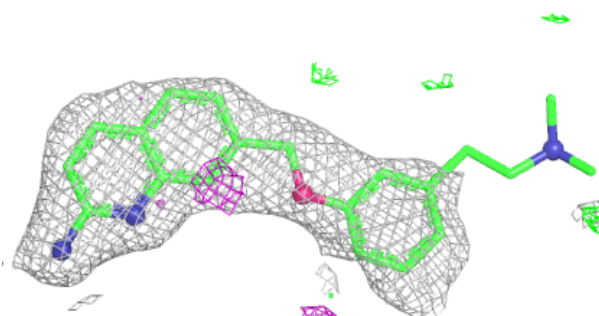
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	H4B	A	760	17/17	0.88	0.14	36,55,79,80	0
3	H4B	B	760	17/17	0.88	0.14	26,54,67,70	0
4	6M8	B	800	24/24	0.93	0.14	9,52,99,102	0
5	ACT	A	860	4/4	0.93	0.07	26,36,39,39	0
5	ACT	B	860	4/4	0.93	0.10	37,43,47,56	0
4	6M8	A	800	24/24	0.94	0.12	11,23,67,78	0
2	HEM	A	750	43/43	0.98	0.06	6,18,42,71	0
2	HEM	B	750	43/43	0.98	0.06	4,16,39,74	0
6	ZN	B	900	1/1	1.00	0.01	24,24,24,24	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.

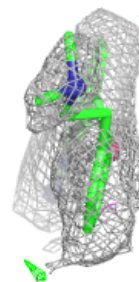
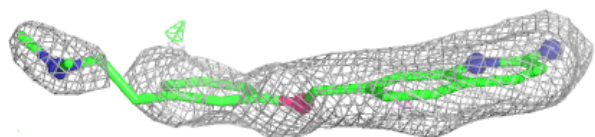
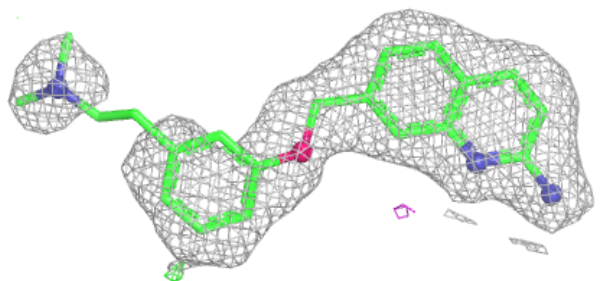
Electron density around 6M8 B 800:

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



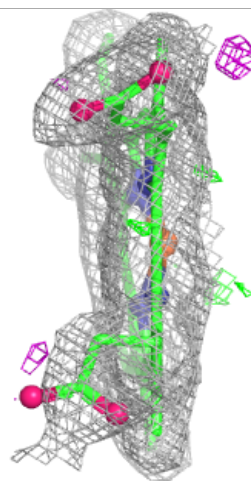
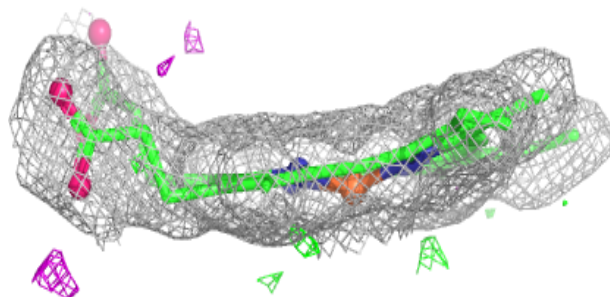
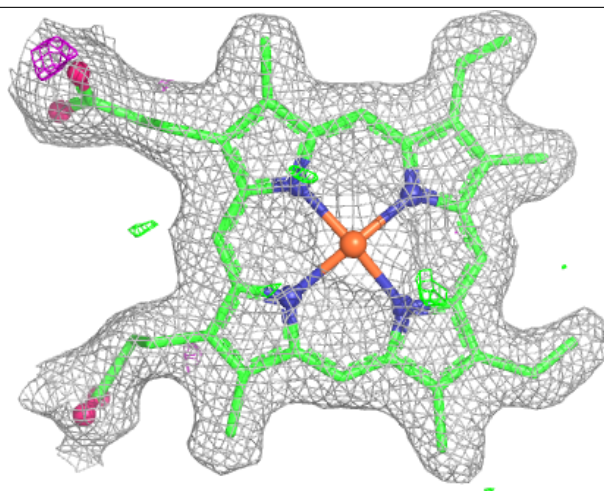
Electron density around 6M8 A 800:

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



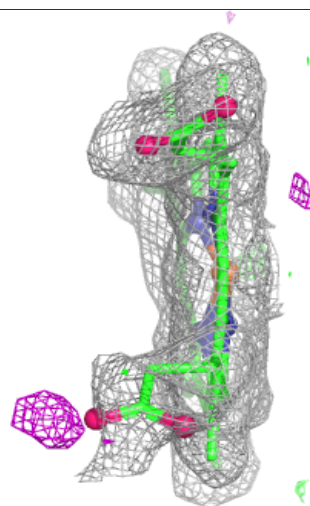
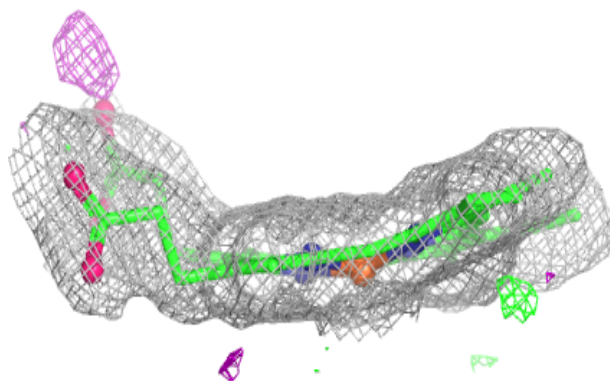
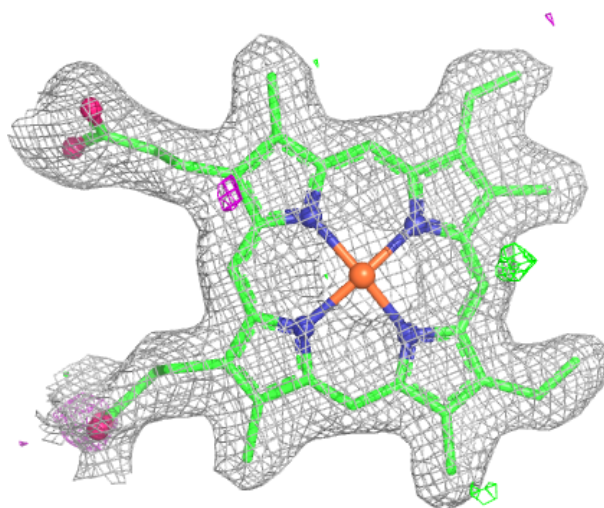
Electron density around HEM A 750:

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



Electron density around HEM B 750:

$2mF_o-DF_c$ (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.