



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:19 PM UTC

PDB ID : 4CFT / pdb\_00004cft  
Title : Structure of bovine endothelial nitric oxide synthase heme domain in complex with 7-((3-Fluorophenethylamino)ethyl)quinolin-2-amine  
Authors : Li, H.; Poulos, T.L.  
Deposited on : 2013-11-19  
Resolution : 1.79 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](mailto: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)  
Xtrriage (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

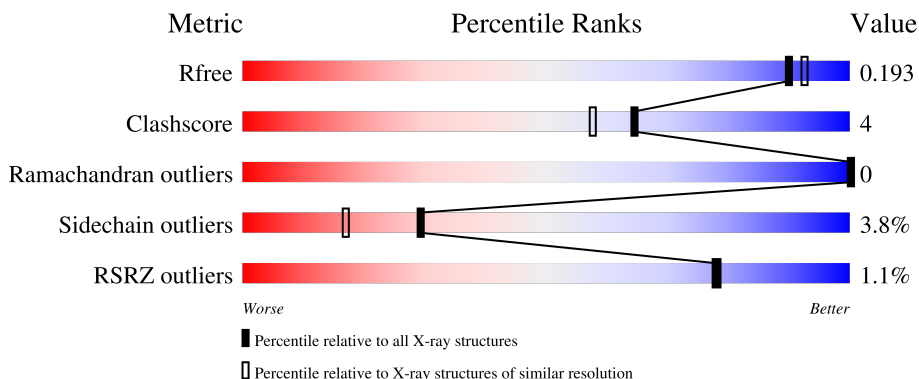
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*


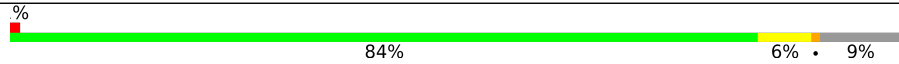
The reported resolution of this entry is 1.79 Å.

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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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  $\geq 3$ , 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	443	 83% 7% • 9%
2	B	443	 84% 6% • 9%

## 2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 6952 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 ENDOTHELIAL NITRIC OXIDE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	As	C	N	O				S
1	A	405	3223	1	2049	568	589	16	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	100	ARG	CYS	conflict	UNP P29473

- Molecule 2 is a protein called ENDOTHELIAL NITRIC OXIDE SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	403	3209	2040	566	587	16	0	0	0

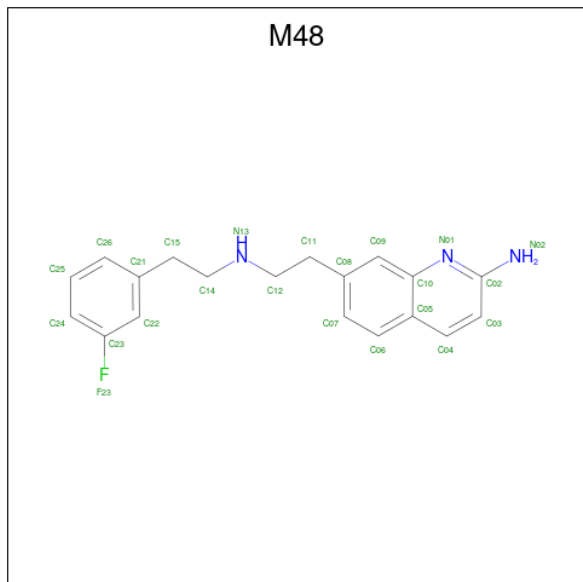
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	100	ARG	CYS	conflict	UNP P29473

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula:  $C_{34}H_{32}FeN_4O_4$ ).

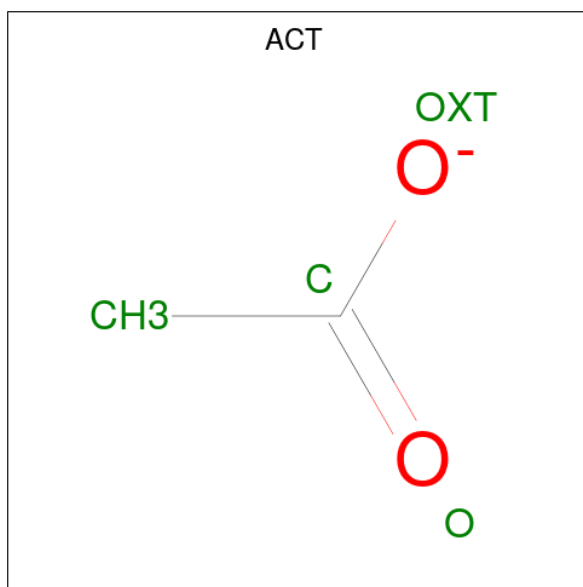


- Molecule 5 is 7-[2-[2-(3-fluorophenyl)ethylamino]ethyl]quinolin-2-amine (CCD ID: M48) (formula:  $C_{19}H_{20}FN_3$ ).



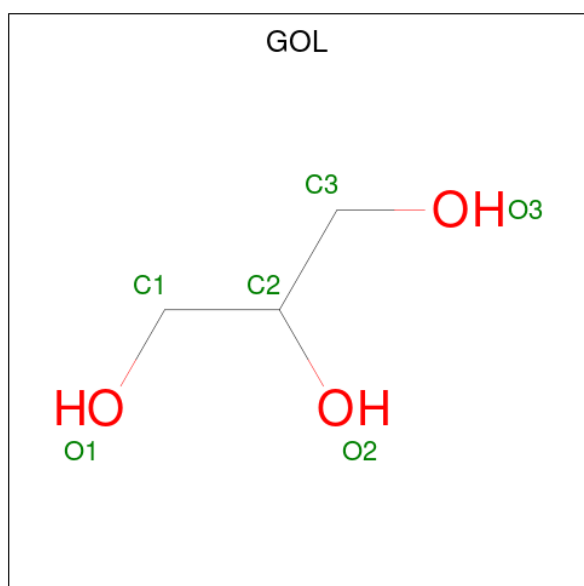
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	F	N		
5	A	1	23	19	1	3	0	0
5	B	1	23	19	1	3	0	0

- Molecule 6 is ACETATE ION (CCD ID: ACT) (formula:  $C_2H_3O_2$ ).



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

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

- Molecule 8 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total Zn 1 1	0	0

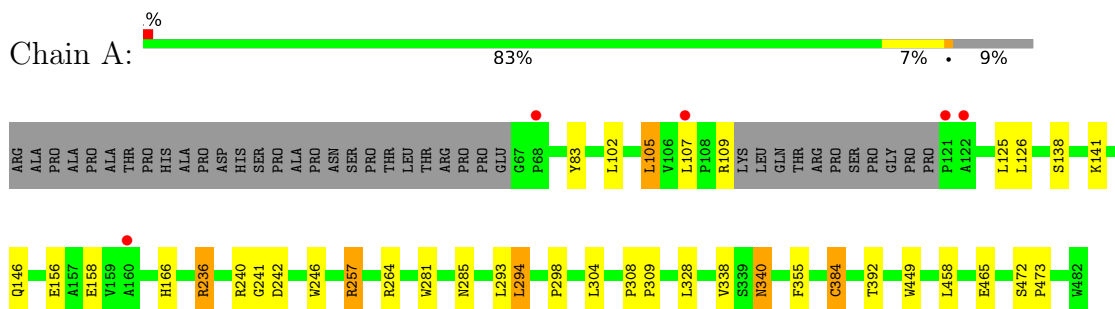
- Molecule 9 is water.

<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
9	A	177	Total 177	O 177	0	0
9	B	148	Total 148	O 148	0	0

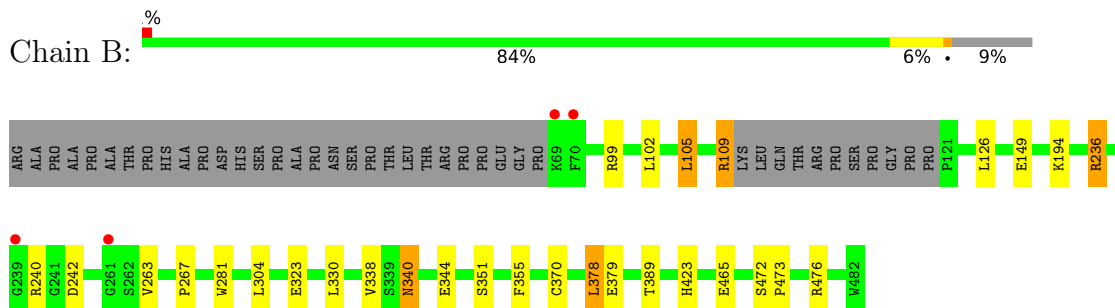
### 3 Residue-property plots [i](#)

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: ENDOTHELIAL NITRIC OXIDE SYNTHASE



- Molecule 2: ENDOTHELIAL NITRIC OXIDE SYNTHASE



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	57.81Å 106.51Å 156.95Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	88.13 – 1.79 88.13 – 1.79	Depositor EDS
% Data completeness (in resolution range)	98.7 (88.13-1.79) 98.8 (88.13-1.79)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.34 (at 1.79Å)	Xtrriage
Refinement program	REFMAC 5.8.0049	Depositor
R, $R_{free}$	0.165 , 0.192 0.165 , 0.193	Depositor DCC
$R_{free}$ test set	4578 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	40.9	Xtrriage
Anisotropy	0.179	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 36.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.98	EDS
Total number of atoms	6952	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	46.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.56% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: HEM, ZN, ACT, M48, GOL, CAS, H4B

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	0.67	0/3303	0.79	0/4497
2	B	0.64	0/3298	0.81	0/4491
All	All	0.65	0/6601	0.80	0/8988

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 [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	3223	0	3127	36	0
2	B	3209	0	3118	16	0
3	A	43	0	30	0	0
3	B	43	0	30	2	0
4	A	17	0	15	1	0
4	B	17	0	15	0	0
5	A	23	0	20	3	0
5	B	23	0	20	4	0
6	A	8	0	6	0	0
6	B	8	0	6	0	0
7	A	6	0	8	0	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	B	6	0	8	0	0
8	A	1	0	0	0	0
9	A	177	0	0	1	0
9	B	148	0	0	2	0
All	All	6952	0	6403	51	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 (51) 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:384:CAS:SG	1:A:384:CAS:AS	2.51	1.27
1:A:236:ARG:HH11	1:A:236:ARG:HG3	1.09	1.16
1:A:257:ARG:HG3	1:A:257:ARG:HH11	1.27	0.96
1:A:236:ARG:HG3	1:A:236:ARG:NH1	1.82	0.84
1:A:240:ARG:HD3	1:A:298:PRO:HB3	1.70	0.74
1:A:257:ARG:HG3	1:A:257:ARG:NH1	2.01	0.73
1:A:236:ARG:HH11	1:A:236:ARG:CG	1.95	0.71
2:B:236:ARG:HD2	2:B:242:ASP:OD1	1.91	0.70
1:A:109:ARG:HH11	1:A:109:ARG:HA	1.57	0.69
9:A:2001:HOH:O	2:B:109:ARG:HG3	1.94	0.68
1:A:240:ARG:HD3	1:A:298:PRO:CB	2.28	0.63
1:A:294:LEU:HD13	1:A:304:LEU:HD13	1.79	0.62
1:A:236:ARG:NH1	1:A:236:ARG:CG	2.58	0.59
1:A:281:TRP:HB2	1:A:304:LEU:HD21	1.85	0.58
1:A:246:TRP:HB2	1:A:294:LEU:HB3	1.84	0.58
1:A:338:VAL:HG21	5:A:801:M48:C07	2.34	0.57
2:B:338:VAL:HG21	5:B:801:M48:C07	2.36	0.56
2:B:281:TRP:HB2	2:B:304:LEU:HD21	1.87	0.55
1:A:107:LEU:HD21	5:A:801:M48:H25	1.88	0.55
1:A:109:ARG:HA	1:A:109:ARG:NH1	2.20	0.54
1:A:294:LEU:HD13	1:A:304:LEU:CD1	2.38	0.52
1:A:355:PHE:HE1	5:A:801:M48:H06	1.74	0.52
2:B:378:LEU:HB2	9:B:2118:HOH:O	2.09	0.52
2:B:355:PHE:HE1	5:B:801:M48:H06	1.74	0.52
1:A:105:LEU:HD22	2:B:465:GLU:HB3	1.91	0.51
1:A:240:ARG:HD2	1:A:241:GLY:O	2.10	0.51
1:A:384:CAS:AS	1:A:384:CAS:CB	3.19	0.51
1:A:257:ARG:NH1	1:A:257:ARG:CG	2.74	0.50
1:A:138:SER:O	1:A:141:LYS:HE3	2.10	0.50

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B:801:M48:H151	9:B:2147:HOH:O	2.12	0.50
1:A:83:TYR:HE2	2:B:109:ARG:HH21	1.60	0.49
1:A:126:LEU:HD11	1:A:156:GLU:HA	1.94	0.49
1:A:384:CAS:SG	1:A:384:CAS:CE2	3.01	0.48
1:A:236:ARG:HG2	1:A:242:ASP:OD1	2.13	0.48
1:A:340:ASN:H	1:A:340:ASN:HD22	1.62	0.48
1:A:472:SER:HA	1:A:473:PRO:C	2.38	0.48
1:A:240:ARG:HD3	1:A:298:PRO:CG	2.44	0.47
2:B:344:GLU:OE1	2:B:476:ARG:NH2	2.49	0.46
3:B:500:HEM:HBA2	5:B:801:M48:C09	2.45	0.46
1:A:392:THR:HB	2:B:423:HIS:HB2	1.97	0.45
2:B:263:VAL:HG11	2:B:267:PRO:HA	1.99	0.45
1:A:465:GLU:HB3	2:B:105:LEU:HD22	1.99	0.44
2:B:370:CYS:SG	2:B:378:LEU:HD13	2.56	0.44
2:B:236:ARG:HD3	2:B:351:SER:HB3	1.99	0.44
2:B:472:SER:HA	2:B:473:PRO:C	2.43	0.44
1:A:264:ARG:HD3	1:A:285:ASN:O	2.18	0.43
1:A:308:PRO:HA	1:A:309:PRO:HD3	1.93	0.43
3:B:500:HEM:HBC2	3:B:500:HEM:CMC	2.48	0.42
2:B:340:ASN:HD22	2:B:340:ASN:H	1.67	0.42
1:A:158:GLU:OE2	1:A:166:HIS:HD2	2.02	0.42
1:A:449:TRP:HA	4:A:600:H4B:N1	2.35	0.41

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	400/443 (90%)	393 (98%)	7 (2%)	0	100	100
2	B	399/443 (90%)	389 (98%)	10 (2%)	0	100	100
All	All	799/886 (90%)	782 (98%)	17 (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	343/375 (92%)	332 (97%)	11 (3%)	34	22
2	B	343/376 (91%)	328 (96%)	15 (4%)	25	13
All	All	686/751 (91%)	660 (96%)	26 (4%)	29	17

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	102	LEU
1	A	105	LEU
1	A	125	LEU
1	A	146	GLN
1	A	236	ARG
1	A	257	ARG
1	A	293	LEU
1	A	294	LEU
1	A	328	LEU
1	A	340	ASN
1	A	458	LEU
2	B	99	ARG
2	B	102	LEU
2	B	105	LEU
2	B	109	ARG
2	B	126	LEU
2	B	149	GLU
2	B	194	LYS
2	B	236	ARG
2	B	240	ARG
2	B	323	GLU
2	B	330	LEU
2	B	340	ASN
2	B	378	LEU
2	B	379	GLU

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
2	B	389	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	GLN
1	A	92	GLN
1	A	128	GLN
1	A	166	HIS
1	A	191	GLN
1	A	207	GLN
1	A	340	ASN
1	A	376	ASN
1	A	413	GLN
1	A	468	ASN
2	B	178	GLN
2	B	191	GLN
2	B	222	ASN
2	B	225	ASN
2	B	340	ASN
2	B	376	ASN
2	B	478	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](#)

1 non-standard protein/DNA/RNA residue is 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
1	CAS	A	384	1	5,8,9	1.26	1 (20%)	1,9,11	0.69	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
1	CAS	A	384	1	-	0/0/7/9	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	384	CAS	AS-CE1	2.36	2.02	1.96

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	384	CAS	3	0

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 13 ligands modelled in this entry, 1 is monoatomic - leaving 12 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
7	GOL	B	880	-	5,5,5	0.21	0	5,5,5	0.82	0
6	ACT	B	860	-	3,3,3	0.71	0	3,3,3	0.90	0
4	H4B	B	600	-	17,18,18	0.77	0	14,26,26	1.95	5 (35%)
4	H4B	A	600	-	17,18,18	1.06	1 (5%)	14,26,26	1.95	6 (42%)
7	GOL	A	880	-	5,5,5	0.22	0	5,5,5	0.64	0
3	HEM	B	500	2	50,50,50	1.67	8 (16%)	67,82,82	1.60	13 (19%)
3	HEM	A	500	1	50,50,50	1.52	7 (14%)	67,82,82	1.61	16 (23%)
6	ACT	B	861	-	3,3,3	0.87	0	3,3,3	0.51	0
5	M48	B	801	-	25,25,25	0.97	0	30,33,33	1.03	1 (3%)
6	ACT	A	861	-	3,3,3	0.82	0	3,3,3	0.75	0
5	M48	A	801	-	25,25,25	1.01	1 (4%)	30,33,33	1.28	4 (13%)
6	ACT	A	860	-	3,3,3	0.71	0	3,3,3	0.98	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
7	GOL	B	880	-	-	0/4/4/4	-
4	H4B	B	600	-	-	0/8/17/17	0/2/2/2
4	H4B	A	600	-	-	0/8/17/17	0/2/2/2
7	GOL	A	880	-	-	0/4/4/4	-
3	HEM	B	500	2	-	1/14/54/54	-
3	HEM	A	500	1	-	0/14/54/54	-
5	M48	B	801	-	-	3/8/8/8	0/3/3/3
5	M48	A	801	-	-	5/8/8/8	0/3/3/3

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	500	HEM	FE-NB	4.96	2.10	1.94
3	B	500	HEM	C1B-NB	-4.42	1.32	1.40
3	B	500	HEM	FE-NC	4.36	2.09	1.95
3	A	500	HEM	FE-NC	4.30	2.09	1.95
3	A	500	HEM	C1B-NB	-3.99	1.33	1.40
3	A	500	HEM	FE-NB	3.94	2.07	1.94
3	B	500	HEM	C4D-ND	-3.23	1.34	1.40
3	A	500	HEM	C4C-NC	-3.06	1.33	1.39
3	A	500	HEM	FE-NA	2.90	2.04	1.95

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	600	H4B	O4-C4	2.75	1.28	1.23
3	B	500	HEM	C4C-NC	-2.71	1.34	1.39
3	B	500	HEM	FE-NA	2.49	2.03	1.95
3	B	500	HEM	C3B-C4B	2.40	1.49	1.44
5	A	801	M48	C02-N01	2.27	1.36	1.33
3	B	500	HEM	C4D-C3D	2.27	1.48	1.45
3	A	500	HEM	C1A-NA	-2.17	1.35	1.39
3	A	500	HEM	C4D-ND	-2.02	1.36	1.40

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	500	HEM	CHC-C4B-NB	4.43	129.19	124.42
4	A	600	H4B	C2-N1-C8A	4.42	121.17	113.36
4	B	600	H4B	C2-N1-C8A	4.11	120.61	113.36
3	A	500	HEM	CHD-C1D-ND	3.90	128.62	124.42
3	B	500	HEM	CHD-C1D-ND	3.64	128.34	124.42
3	B	500	HEM	CBA-CAA-C2A	-3.61	102.56	112.53
3	A	500	HEM	CHD-C1D-C2D	-3.56	119.42	125.03
3	A	500	HEM	CHC-C4B-NB	3.35	128.03	124.42
3	B	500	HEM	C1B-NB-C4B	3.32	109.14	105.21
3	A	500	HEM	C3B-C4B-NB	-3.27	107.12	109.47
3	B	500	HEM	CHD-C1D-C2D	-3.23	119.93	125.03
3	B	500	HEM	C3B-C4B-NB	-3.10	107.24	109.47
4	B	600	H4B	O4-C4-C4A	-3.06	119.88	127.26
4	B	600	H4B	C4A-C4-N3	2.98	120.32	112.13
3	A	500	HEM	CHC-C1C-NC	2.97	127.69	124.45
3	B	500	HEM	C1A-CHA-C4D	-2.93	119.35	126.25
3	B	500	HEM	CHC-C1C-NC	2.83	127.54	124.45
3	A	500	HEM	CHA-C4D-C3D	-2.74	120.17	125.23
3	A	500	HEM	C4C-CHD-C1D	-2.68	120.32	126.02
3	A	500	HEM	C1A-CHA-C4D	-2.65	120.01	126.25
5	B	801	M48	C24-C23-C22	-2.62	119.77	123.23
3	A	500	HEM	O2D-CGD-CBD	2.60	122.22	114.00
3	B	500	HEM	O2D-CGD-CBD	2.58	122.14	114.00
3	A	500	HEM	C1B-NB-C4B	2.53	108.20	105.21
3	A	500	HEM	C3D-C4D-ND	2.52	112.94	110.17
5	A	801	M48	C05-C10-N01	-2.51	118.80	122.25
4	A	600	H4B	N2-C2-N3	2.49	122.03	116.76
5	A	801	M48	C24-C23-C22	-2.47	119.97	123.23
4	A	600	H4B	O4-C4-C4A	-2.40	121.46	127.26
3	B	500	HEM	C4C-NC-C1C	2.40	109.74	105.82

*Continued on next page...*

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	500	HEM	CHA-C4D-C3D	-2.33	120.93	125.23
3	B	500	HEM	C4B-C3B-C2B	-2.31	105.16	107.28
4	A	600	H4B	C4A-C4-N3	2.27	118.36	112.13
4	B	600	H4B	C2-N3-C4	-2.24	121.04	125.11
3	A	500	HEM	O2A-CGA-O1A	-2.19	117.70	123.33
3	A	500	HEM	O2A-CGA-CBA	2.15	120.78	114.00
4	A	600	H4B	C4-C4A-N5	2.10	121.99	116.27
4	B	600	H4B	C4-C4A-N5	2.09	121.96	116.27
3	A	500	HEM	C4C-NC-C1C	2.08	109.21	105.82
3	A	500	HEM	C4A-CHB-C1B	-2.07	121.38	126.25
3	B	500	HEM	CHA-C4D-ND	2.06	126.92	124.37
5	A	801	M48	C12-N13-C14	2.06	123.15	113.40
5	A	801	M48	C09-C10-N01	2.04	121.76	118.78
3	A	500	HEM	CBA-CAA-C2A	-2.02	106.96	112.53
4	A	600	H4B	N3-C2-N1	-2.01	119.64	123.32

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	801	M48	N13-C14-C15-C21
5	A	801	M48	C11-C12-N13-C14
5	B	801	M48	N13-C14-C15-C21
5	A	801	M48	C15-C14-N13-C12
5	B	801	M48	C11-C12-N13-C14
5	B	801	M48	C15-C14-N13-C12
5	A	801	M48	C07-C08-C11-C12
5	A	801	M48	C09-C08-C11-C12
3	B	500	HEM	CAD-CBD-CGD-O2D

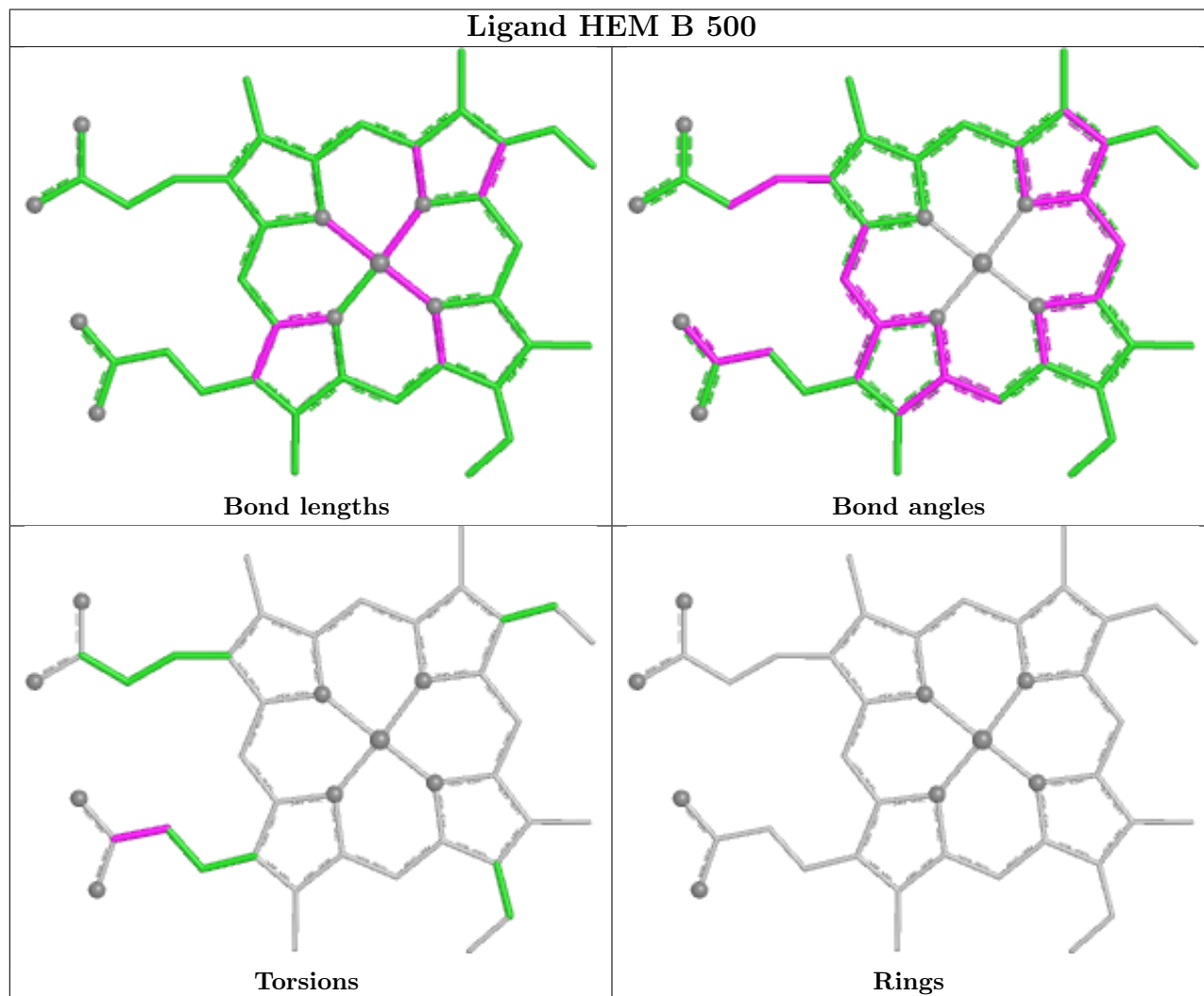
There are no ring outliers.

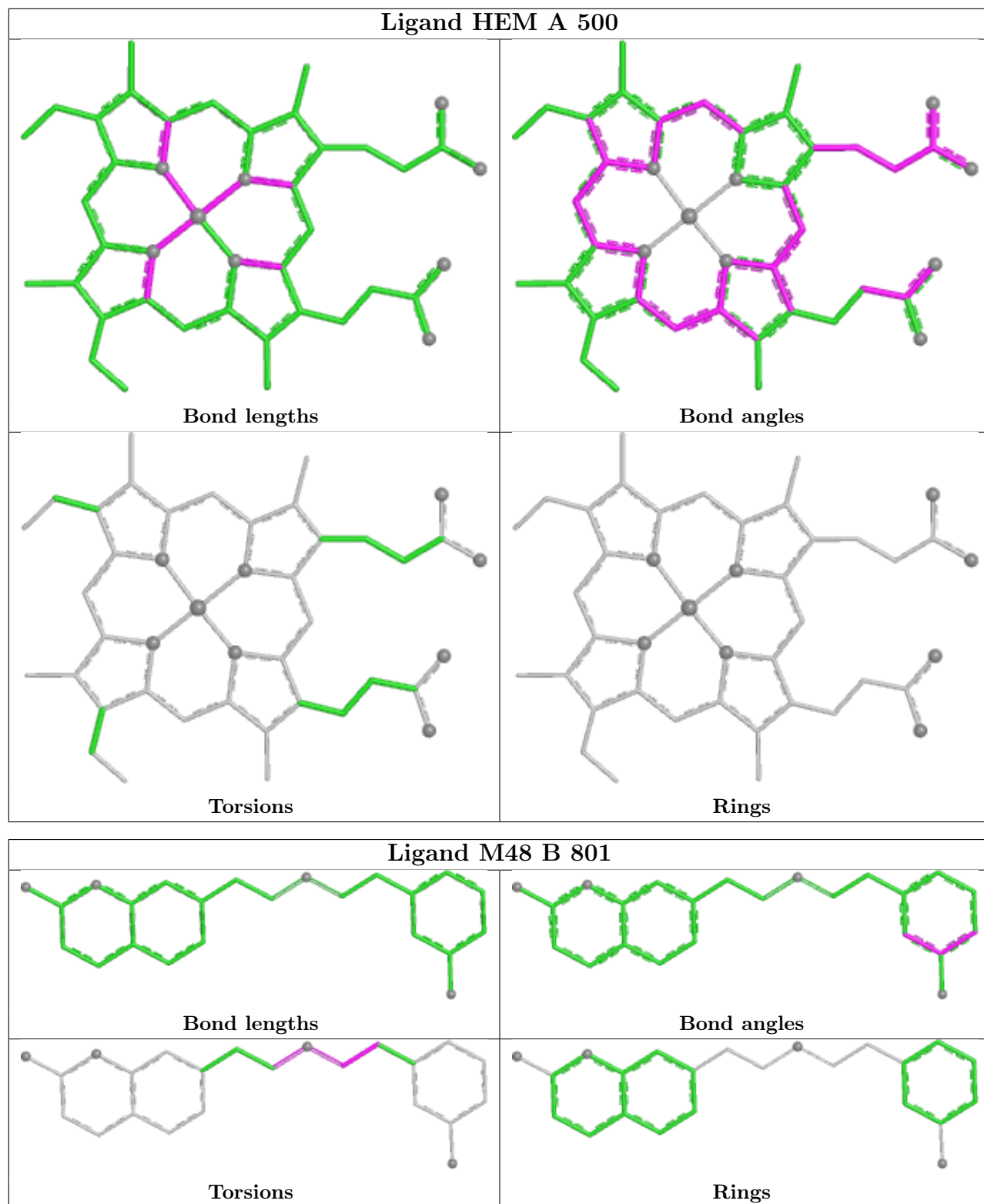
4 monomers are involved in 9 short contacts:

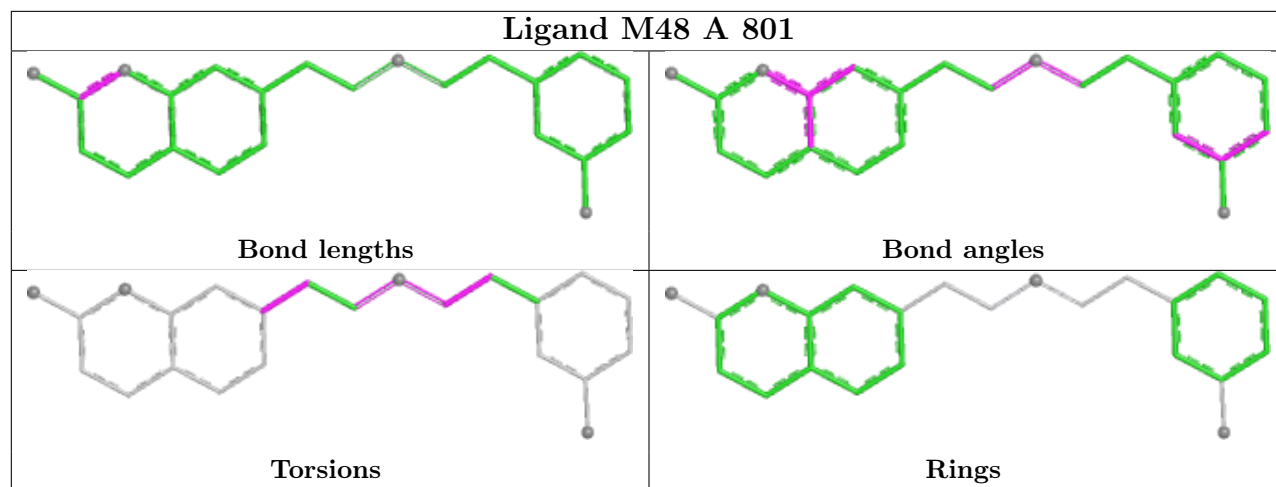
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	600	H4B	1	0
3	B	500	HEM	2	0
5	B	801	M48	4	0
5	A	801	M48	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

addition, ligands with molecular weight  $> 250$  and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	404/443 (91%)	0.06	5 (1%) 76 77	31, 42, 68, 91	0
2	B	403/443 (90%)	0.17	4 (0%) 79 80	31, 44, 69, 99	0
All	All	807/886 (91%)	0.11	9 (1%) 78 78	31, 43, 69, 99	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	121	PRO	4.7
2	B	70	PHE	3.9
1	A	107	LEU	3.7
1	A	160	ALA	3.4
1	A	122	ALA	3.3
2	B	261	GLY	3.3
2	B	239	GLY	3.2
2	B	69	LYS	3.2
1	A	68	PRO	2.5

### 6.2 Non-standard residues in protein, DNA, RNA chains [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	CAS	A	384	9/10	0.88	0.17	40,41,81,87	0

### 6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 6.4 Ligands

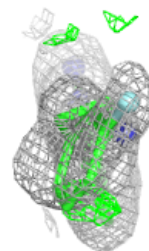
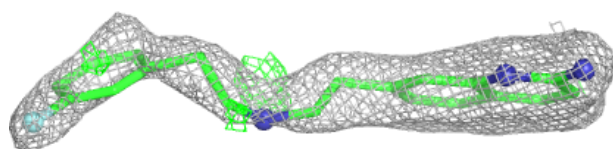
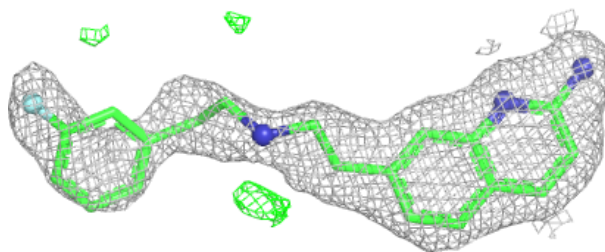
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
6	ACT	B	861	4/4	0.82	0.17	48,52,55,55	0
7	GOL	A	880	6/6	0.85	0.16	54,62,66,69	0
7	GOL	B	880	6/6	0.85	0.15	52,59,62,64	0
5	M48	B	801	23/23	0.92	0.12	39,47,89,92	0
5	M48	A	801	23/23	0.92	0.13	34,47,94,95	0
6	ACT	A	860	4/4	0.95	0.09	46,50,52,56	0
6	ACT	A	861	4/4	0.95	0.09	44,46,46,46	0
6	ACT	B	860	4/4	0.95	0.10	45,52,53,57	0
4	H4B	A	600	17/17	0.96	0.06	33,38,40,42	0
4	H4B	B	600	17/17	0.97	0.06	35,38,41,41	0
3	HEM	A	500	43/43	0.98	0.07	30,34,42,44	0
3	HEM	B	500	43/43	0.98	0.07	33,36,43,48	0
8	ZN	A	1483	1/1	0.99	0.08	40,40,40,40	0

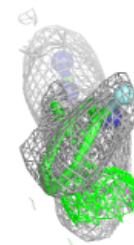
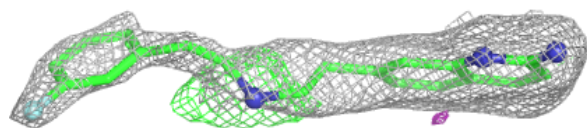
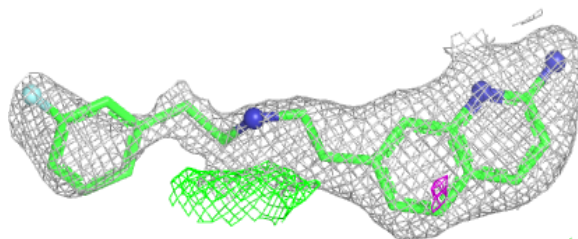
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

**Electron density around M48 B 801:**

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

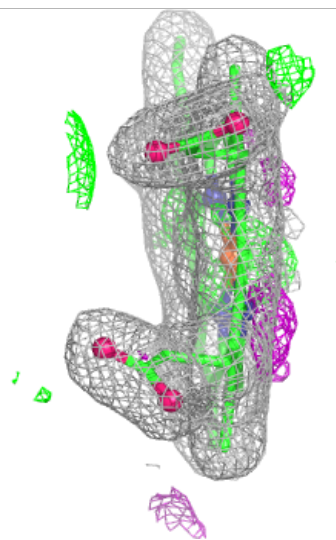
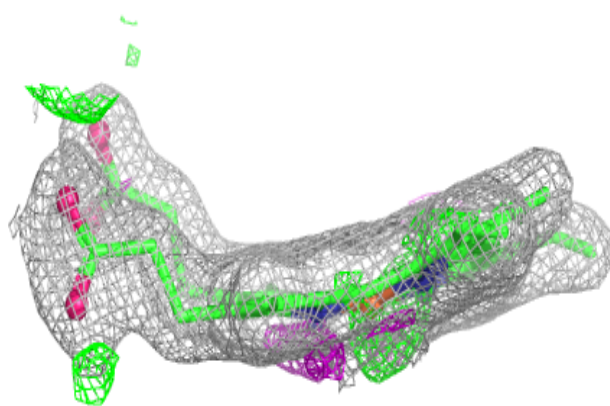
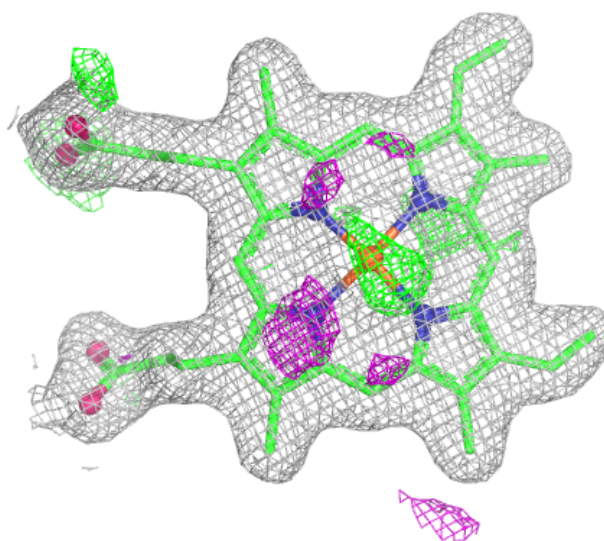
**Electron density around M48 A 801:**

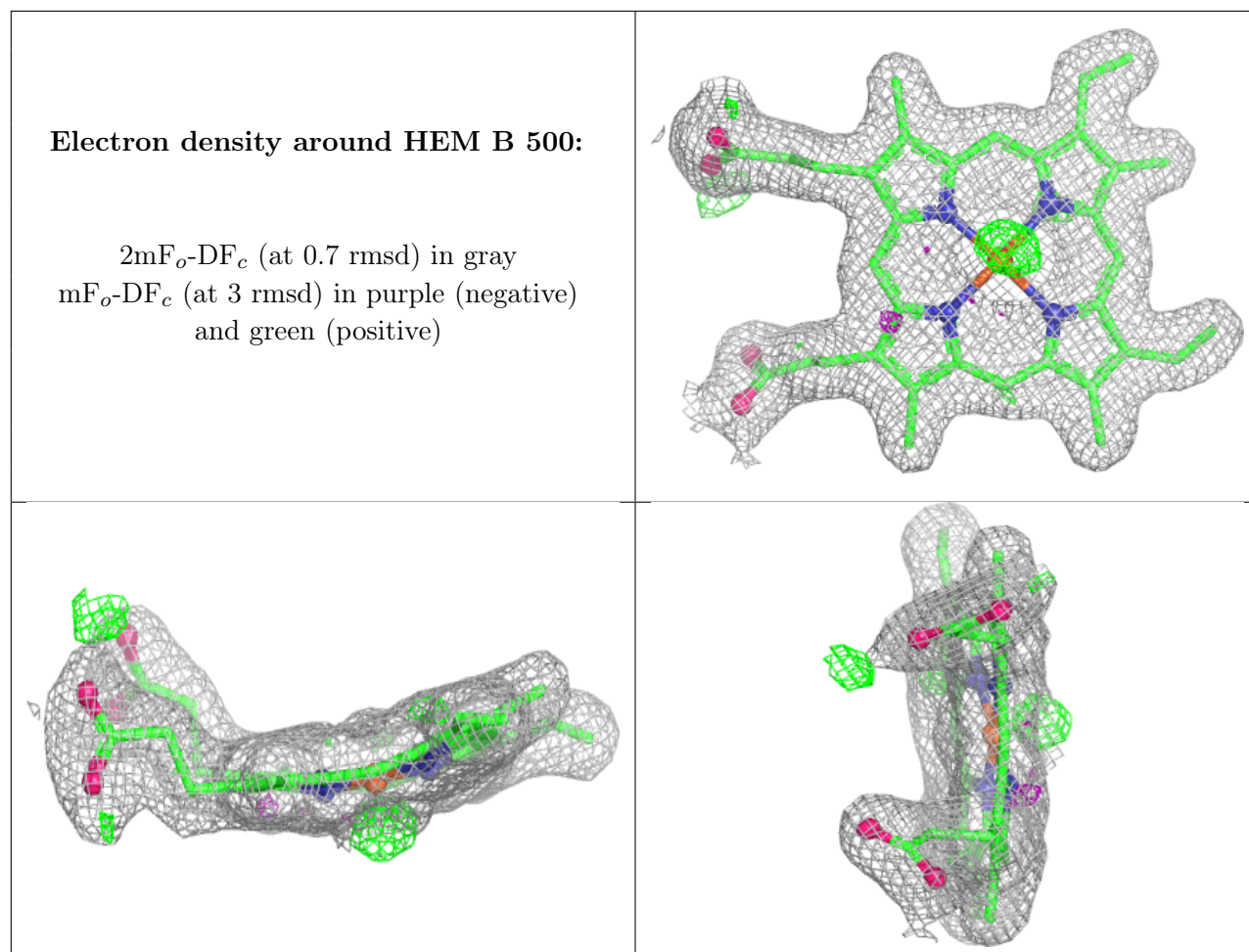
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around HEM A 500:**

$2mF_o-DF_c$  (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.