



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 25, 2020 – 03:27 PM BST

PDB ID : 4FVX  
Title : Structure of rat nNOS heme domain in complex with N(omega)-ethoxy-L-arginine  
Authors : Li, H.; Poulos, T.L.  
Deposited on : 2012-06-29  
Resolution : 2.00 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.

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The following versions of software and data (see [references ⓘ](#)) 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 : 2.13  
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) : 2.13

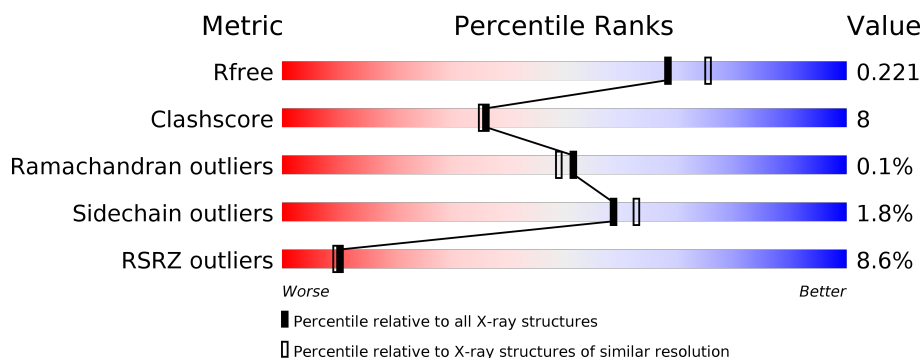
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 on 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	422	<div> <div>12%</div> <div> <div></div> <div>76%</div> <div>19%</div> <div>• •</div> </div> </div>
1	B	422	<div> <div>5%</div> <div> <div></div> <div>85%</div> <div>12%</div> <div>•</div> </div> </div>

## 2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 7211 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	407	Total	C	N	O	S	0	0	0
			3313	2121	566	605	21			
1	B	411	Total	C	N	O	S	0	0	0
			3345	2140	574	610	21			

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



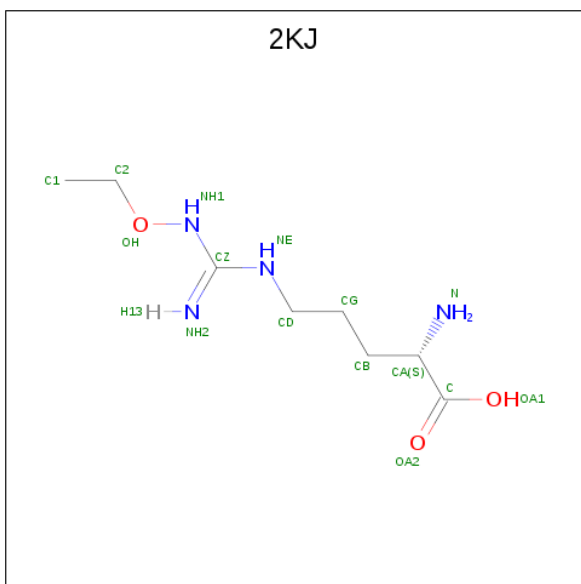
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- 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
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 N 5 -(N-ethoxycarbamimidoyl)-L-ornithine (three-letter code: 2KJ) (formula:  $C_8H_{18}N_4O_3$ ).



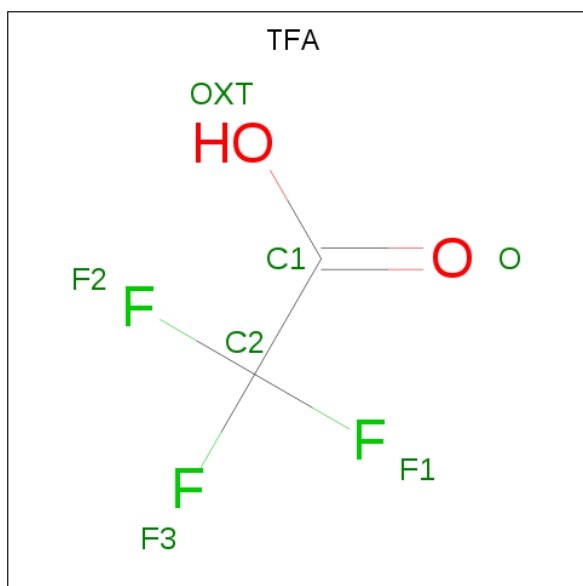
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			15	8	4	3		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	N	O	0	0
			15	8	4	3		

- Molecule 5 is trifluoroacetic acid (three-letter code: TFA) (formula:  $C_2HF_3O_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	F	O	0	0
			7	2	3	2		
5	B	1	Total	C	F	O	0	0
			7	2	3	2		

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

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

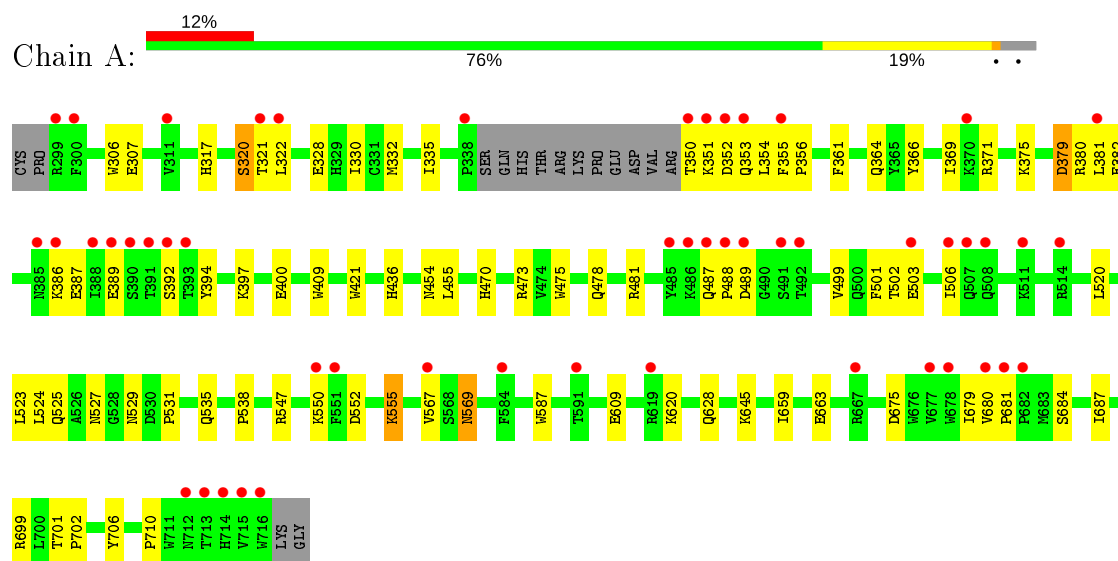
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	161	Total	O	0	0
			161	161		
7	B	227	Total	O	0	0
			227	227		

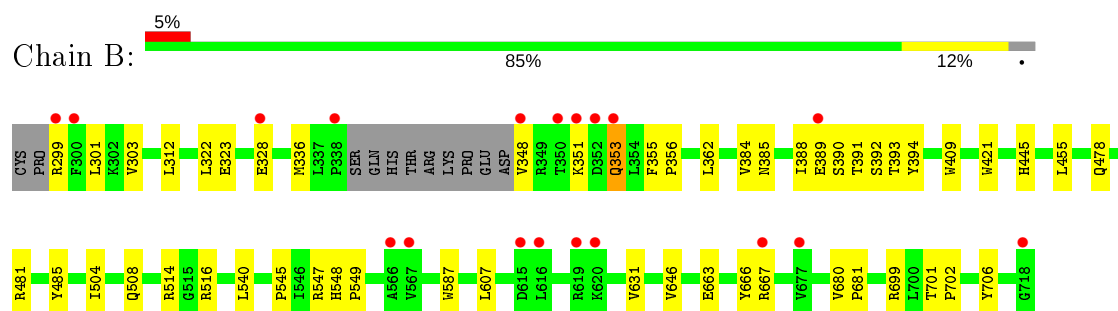
### 3 Residue-property plots [i](#)

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



- 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, $\alpha$ , $\beta$ , $\gamma$	52.00Å 111.28Å 165.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.80 – 2.00 37.80 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.5 (37.80-2.00) 99.5 (37.80-2.00)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.52 (at 2.00Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.181 , 0.221 0.180 , 0.221	Depositor DCC
$R_{free}$ test set	3215 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	32.2	Xtriage
Anisotropy	0.742	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 51.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	7211	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.32% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, ZN, TFA, H4B, 2KJ

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	0.60	0/3406	0.67	2/4621 (0.0%)
1	B	0.66	0/3438	0.67	0/4661
All	All	0.63	0/6844	0.67	2/9282 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	379	ASP	CB-CG-OD2	6.66	124.30	118.30
1	A	381	LEU	CA-CB-CG	5.32	127.54	115.30

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	3313	0	3221	68	0
1	B	3345	0	3259	43	0
2	A	43	0	30	3	0
2	B	43	0	30	3	0
3	A	17	0	15	0	0
3	B	17	0	15	0	0
4	A	15	0	16	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	15	0	16	0	0
5	A	7	0	0	0	0
5	B	7	0	0	0	0
6	A	1	0	0	0	0
7	A	161	0	0	2	0
7	B	227	0	0	9	0
All	All	7211	0	6602	107	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 8.

All (107) 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:B:706:TYR:OH	2:B:801:HEM:O1D	1.92	0.86
1:B:303:VAL:HG12	7:B:1049:HOH:O	1.82	0.79
1:A:350:THR:HG22	1:A:352:ASP:H	1.49	0.76
1:A:335:ILE:HD11	7:B:1049:HOH:O	1.92	0.70
1:B:646:VAL:HG12	7:B:1033:HOH:O	1.93	0.67
1:A:350:THR:HB	1:A:353:GLN:HG3	1.76	0.67
1:A:355:PHE:N	1:A:356:PRO:HD2	2.10	0.66
1:A:350:THR:HB	1:A:353:GLN:CG	2.27	0.65
1:A:436:HIS:ND1	7:A:980:HOH:O	2.30	0.65
1:A:478:GLN:HB2	1:A:481:ARG:HG3	1.80	0.64
1:A:350:THR:HB	1:A:353:GLN:CD	2.22	0.60
1:A:307:GLU:HG3	7:B:939:HOH:O	2.01	0.60
1:A:355:PHE:N	1:A:356:PRO:CD	2.65	0.59
1:A:322:LEU:HD13	1:A:699:ARG:NH2	2.17	0.59
1:A:380:ARG:HD3	1:A:400:GLU:OE1	2.03	0.58
1:A:675:ASP:O	1:A:679:ILE:HG12	2.03	0.58
1:A:523:LEU:HD22	1:A:531:PRO:HB2	1.84	0.58
1:A:569:ASN:HD22	1:A:569:ASN:H	1.50	0.58
1:B:504:ILE:HD11	7:B:1052:HOH:O	2.03	0.58
1:A:353:GLN:O	1:A:356:PRO:HG2	2.04	0.57
1:A:473:ARG:NH2	1:A:710:PRO:HD3	2.19	0.57
1:B:328:GLU:OE1	1:B:328:GLU:N	2.37	0.57
1:A:455:LEU:HD12	1:A:587:TRP:HB3	1.87	0.57
1:A:306:TRP:CD2	1:B:336:MET:HE2	2.40	0.56
1:B:355:PHE:N	1:B:356:PRO:HD2	2.19	0.56
1:A:628:GLN:HG2	1:B:631:VAL:HG11	1.86	0.56
1:B:355:PHE:CE1	1:B:385:ASN:HB2	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:317:HIS:O	1:A:320:SER:HB3	2.06	0.55
1:B:646:VAL:CG1	7:B:1033:HOH:O	2.53	0.55
1:A:361:PHE:O	1:A:364:GLN:HG2	2.06	0.54
1:A:609:GLU:HG3	7:A:974:HOH:O	2.07	0.54
1:A:350:THR:HG22	1:A:352:ASP:N	2.20	0.54
1:A:306:TRP:CG	1:B:336:MET:HE1	2.43	0.54
1:A:354:LEU:C	1:A:356:PRO:HD2	2.28	0.54
1:A:567:VAL:HG23	4:A:803:2KJ:H19	1.90	0.53
1:A:371:ARG:HG3	1:A:371:ARG:HH21	1.71	0.53
1:A:555:LYS:HB3	1:A:555:LYS:NZ	2.23	0.53
1:B:299:ARG:HB3	1:B:299:ARG:NH1	2.23	0.53
1:A:306:TRP:CE2	1:B:336:MET:HE2	2.44	0.52
1:A:535:GLN:HE21	1:A:538:PRO:HD3	1.75	0.52
1:B:548:HIS:CG	1:B:549:PRO:HD2	2.46	0.51
1:A:332:MET:HE2	1:B:301:LEU:HD13	1.92	0.51
1:B:391:THR:O	1:B:392:SER:HB2	2.11	0.51
1:B:323:GLU:O	1:B:699:ARG:HD3	2.11	0.50
1:A:525:GLN:HG3	1:A:529:ASN:O	2.11	0.50
2:A:801:HEM:HMC1	2:A:801:HEM:HBC2	1.93	0.50
1:A:487:GLN:HB3	1:A:488:PRO:HD2	1.93	0.50
1:B:701:THR:HA	1:B:702:PRO:C	2.32	0.49
2:A:801:HEM:CMC	2:A:801:HEM:HBC2	2.43	0.49
2:B:801:HEM:HBC2	2:B:801:HEM:CMC	2.42	0.49
1:B:445:HIS:CD2	1:B:445:HIS:C	2.87	0.48
1:B:516:ARG:HD2	7:B:1111:HOH:O	2.11	0.48
1:A:455:LEU:HD12	1:A:587:TRP:CB	2.44	0.47
1:A:380:ARG:NH1	1:A:397:LYS:HG2	2.29	0.47
1:A:366:TYR:HA	1:A:369:ILE:HG12	1.96	0.46
1:A:659:ILE:O	1:A:663:GLU:HG3	2.15	0.46
1:A:524:LEU:O	1:A:531:PRO:HA	2.15	0.46
1:A:328:GLU:H	1:A:328:GLU:CD	2.18	0.46
1:A:470:HIS:HB3	1:A:527:ASN:ND2	2.31	0.46
1:A:680:VAL:HA	1:A:681:PRO:HD3	1.83	0.46
1:A:502:THR:O	1:A:506:ILE:HG13	2.15	0.45
1:A:382:GLU:O	1:A:386:LYS:HG3	2.17	0.45
1:B:389:GLU:HG3	1:B:390:SER:N	2.30	0.45
1:A:353:GLN:C	1:A:356:PRO:HD2	2.38	0.45
1:A:306:TRP:CD2	1:B:336:MET:CE	2.99	0.45
1:B:348:VAL:HG13	1:B:348:VAL:O	2.18	0.44
1:A:684:SER:HB3	1:A:687:ILE:HG12	2.00	0.44
1:B:388:ILE:O	1:B:392:SER:N	2.45	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:801:HEM:HBC2	2:B:801:HEM:HMC1	2.00	0.44
1:A:569:ASN:ND2	1:A:569:ASN:H	2.15	0.43
1:B:478:GLN:HB2	1:B:481:ARG:HG3	2.00	0.43
1:A:375:LYS:NZ	1:A:379:ASP:OD1	2.49	0.43
1:A:501:PHE:HD2	1:A:520:LEU:HD13	1.83	0.43
1:B:362:LEU:HD11	1:B:384:VAL:HG21	2.00	0.43
1:B:504:ILE:O	1:B:508:GLN:HG2	2.18	0.43
1:A:475:TRP:HB2	1:A:523:LEU:HB3	2.00	0.43
1:A:569:ASN:N	1:A:569:ASN:HD22	2.09	0.43
1:A:330:ILE:O	1:A:330:ILE:HG23	2.17	0.43
1:A:409:TRP:CE3	1:A:421:TRP:HA	2.53	0.43
1:B:353:GLN:HB3	1:B:353:GLN:HE21	1.55	0.43
1:A:487:GLN:HB3	1:A:488:PRO:CD	2.49	0.43
1:B:322:LEU:HD13	1:B:699:ARG:NH2	2.34	0.43
1:B:485:TYR:CZ	1:B:514:ARG:HA	2.54	0.43
1:B:699:ARG:NE	7:B:1026:HOH:O	2.52	0.43
1:B:351:LYS:HE3	1:B:392:SER:OG	2.18	0.42
1:A:306:TRP:CD1	1:B:336:MET:HE1	2.54	0.42
1:A:321:THR:HG23	1:A:322:LEU:N	2.35	0.42
1:A:353:GLN:O	1:A:356:PRO:HD2	2.20	0.42
1:A:706:TYR:OH	2:A:801:HEM:O1D	2.13	0.41
1:A:321:THR:HG23	1:A:322:LEU:HG	2.02	0.41
1:B:393:THR:OG1	1:B:394:TYR:N	2.52	0.41
1:A:351:LYS:HE3	1:A:392:SER:HB3	2.01	0.41
1:A:550:LYS:HB2	1:A:550:LYS:HE3	1.50	0.41
1:B:409:TRP:CE3	1:B:421:TRP:HA	2.55	0.41
1:A:701:THR:HA	1:A:702:PRO:C	2.41	0.41
1:A:322:LEU:HB2	1:A:699:ARG:HB2	2.01	0.41
1:B:322:LEU:HB2	1:B:699:ARG:HB2	2.03	0.41
1:A:306:TRP:CG	1:B:336:MET:CE	3.04	0.41
1:B:680:VAL:HA	1:B:681:PRO:HD3	1.85	0.41
1:A:684:SER:HB3	1:A:687:ILE:CG1	2.51	0.41
1:B:516:ARG:CD	7:B:1111:HOH:O	2.66	0.41
1:B:455:LEU:HD23	1:B:587:TRP:HB3	2.03	0.41
1:B:312:LEU:HB3	1:B:666:TYR:CD2	2.56	0.40
1:B:663:GLU:O	1:B:667:ARG:HD2	2.22	0.40
1:A:499:VAL:O	1:A:503:GLU:HG3	2.22	0.40
1:A:387:GLU:OE1	1:A:394:TYR:HA	2.22	0.40
1:B:607:LEU:HA	1:B:607:LEU:HD23	1.97	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	403/422 (96%)	390 (97%)	12 (3%)	1 (0%)	47	44
1	B	407/422 (96%)	399 (98%)	8 (2%)	0	100	100
All	All	810/844 (96%)	789 (97%)	20 (2%)	1 (0%)	51	49

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	489	ASP

### 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	363/377 (96%)	354 (98%)	9 (2%)	47	49
1	B	366/377 (97%)	362 (99%)	4 (1%)	73	78
All	All	729/754 (97%)	716 (98%)	13 (2%)	59	63

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	320	SER
1	A	389	GLU
1	A	454	ASN
1	A	547	ARG
1	A	552	ASP

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Mol	Chain	Res	Type
1	A	555	LYS
1	A	569	ASN
1	A	620	LYS
1	A	645	LYS
1	B	353	GLN
1	B	540	LEU
1	B	545	PRO
1	B	547	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	353	GLN
1	A	425	GLN
1	A	454	ASN
1	A	507	GLN
1	A	527	ASN
1	A	535	GLN
1	A	569	ASN
1	A	605	ASN
1	A	628	GLN
1	A	642	GLN
1	A	697	ASN
1	B	364	GLN
1	B	385	ASN
1	B	425	GLN
1	B	454	ASN
1	B	507	GLN
1	B	535	GLN
1	B	601	ASN
1	B	605	ASN
1	B	642	GLN
1	B	697	ASN

### 5.3.3 RNA ⓘ

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 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	$\# Z  > 2$
5	TFA	B	804	-	3,6,6	0.10	0	3,9,9	1.39	0
5	TFA	A	804	-	3,6,6	0.13	0	3,9,9	1.13	0
4	2KJ	A	803	-	8,14,14	0.63	0	7,16,16	1.08	1 (14%)
2	HEM	A	801	1	27,50,50	2.13	5 (18%)	17,82,82	2.61	7 (41%)
3	H4B	B	802	-	16,18,18	1.42	3 (18%)	11,26,26	2.30	6 (54%)
2	HEM	B	801	1	27,50,50	2.26	8 (29%)	17,82,82	2.29	5 (29%)
4	2KJ	B	803	-	8,14,14	0.67	0	7,16,16	1.31	1 (14%)
3	H4B	A	802	-	16,18,18	0.82	0	11,26,26	2.66	6 (54%)

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
5	TFA	B	804	-	-	0/0/6/6	-
5	TFA	A	804	-	-	0/0/6/6	-
4	2KJ	A	803	-	-	3/9/15/15	-
2	HEM	A	801	1	-	0/6/54/54	-
3	H4B	B	802	-	-	0/8/17/17	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	B	801	1	-	0/6/54/54	-
4	2KJ	B	803	-	-	2/9/15/15	-
3	H4B	A	802	-	-	0/8/17/17	0/2/2/2

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	HEM	C3B-C2B	-5.12	1.33	1.40
2	B	801	HEM	C3D-C2D	4.95	1.52	1.37
2	A	801	HEM	C3D-C2D	4.78	1.51	1.37
2	B	801	HEM	C3C-C2C	-4.65	1.33	1.40
2	B	801	HEM	C3B-C2B	-4.10	1.34	1.40
2	A	801	HEM	C3C-CAC	3.78	1.55	1.47
2	A	801	HEM	C3B-CAB	3.70	1.55	1.47
2	B	801	HEM	C3C-CAC	3.64	1.55	1.47
3	B	802	H4B	C7-C6	3.27	1.55	1.52
2	B	801	HEM	CAD-C3D	3.25	1.57	1.52
2	B	801	HEM	C3B-CAB	2.69	1.53	1.47
2	B	801	HEM	CMD-C2D	2.69	1.57	1.51
3	B	802	H4B	C7-N8	2.62	1.49	1.44
2	B	801	HEM	CMB-C2B	2.61	1.57	1.51
3	B	802	H4B	C4-N3	2.55	1.37	1.33
2	A	801	HEM	C3C-C2C	-2.46	1.37	1.40

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	HEM	CBA-CAA-C2A	-6.14	101.15	112.49
2	A	801	HEM	CBA-CAA-C2A	-6.07	101.29	112.49
2	A	801	HEM	CBD-CAD-C3D	-5.91	101.59	112.48
3	A	802	H4B	C4-C4A-C8A	4.61	118.66	114.57
2	B	801	HEM	CBD-CAD-C3D	-4.48	104.23	112.48
3	B	802	H4B	C4-C4A-C8A	4.26	118.36	114.57
3	A	802	H4B	C4-C4A-N5	3.57	122.12	119.12
3	A	802	H4B	C2-N1-C8A	3.38	122.12	114.54
3	A	802	H4B	N3-C2-N1	-3.11	120.55	125.42
3	B	802	H4B	C4-C4A-N5	3.03	121.66	119.12
2	B	801	HEM	C1D-C2D-C3D	-3.02	104.89	107.00
3	A	802	H4B	N2-C2-N3	3.01	121.93	117.25
2	A	801	HEM	C4C-C3C-C2C	2.96	108.96	106.90
3	B	802	H4B	N3-C2-N1	-2.90	120.87	125.42
3	B	802	H4B	C2-N1-C8A	2.90	121.04	114.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	HEM	CMA-C3A-C4A	-2.83	124.11	128.46
2	A	801	HEM	CMC-C2C-C3C	2.67	129.67	124.68
2	A	801	HEM	C1D-C2D-C3D	-2.66	105.15	107.00
3	A	802	H4B	C4-N3-C2	2.61	120.08	115.93
2	A	801	HEM	C4A-C3A-C2A	2.37	108.65	107.00
2	B	801	HEM	CMA-C3A-C4A	-2.35	124.85	128.46
4	A	803	2KJ	CD-NE-CZ	-2.30	119.29	123.50
3	B	802	H4B	C4-N3-C2	2.26	119.51	115.93
4	B	803	2KJ	CD-NE-CZ	-2.25	119.39	123.50
3	B	802	H4B	N2-C2-N3	2.15	120.60	117.25
2	B	801	HEM	CAD-CBD-CGD	2.07	116.14	112.67

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	803	2KJ	CZ-NH1-OH-C2
4	B	803	2KJ	CZ-NH1-OH-C2
4	A	803	2KJ	NE-CD-CG-CB
4	B	803	2KJ	NE-CD-CG-CB
4	A	803	2KJ	C1-C2-OH-NH1

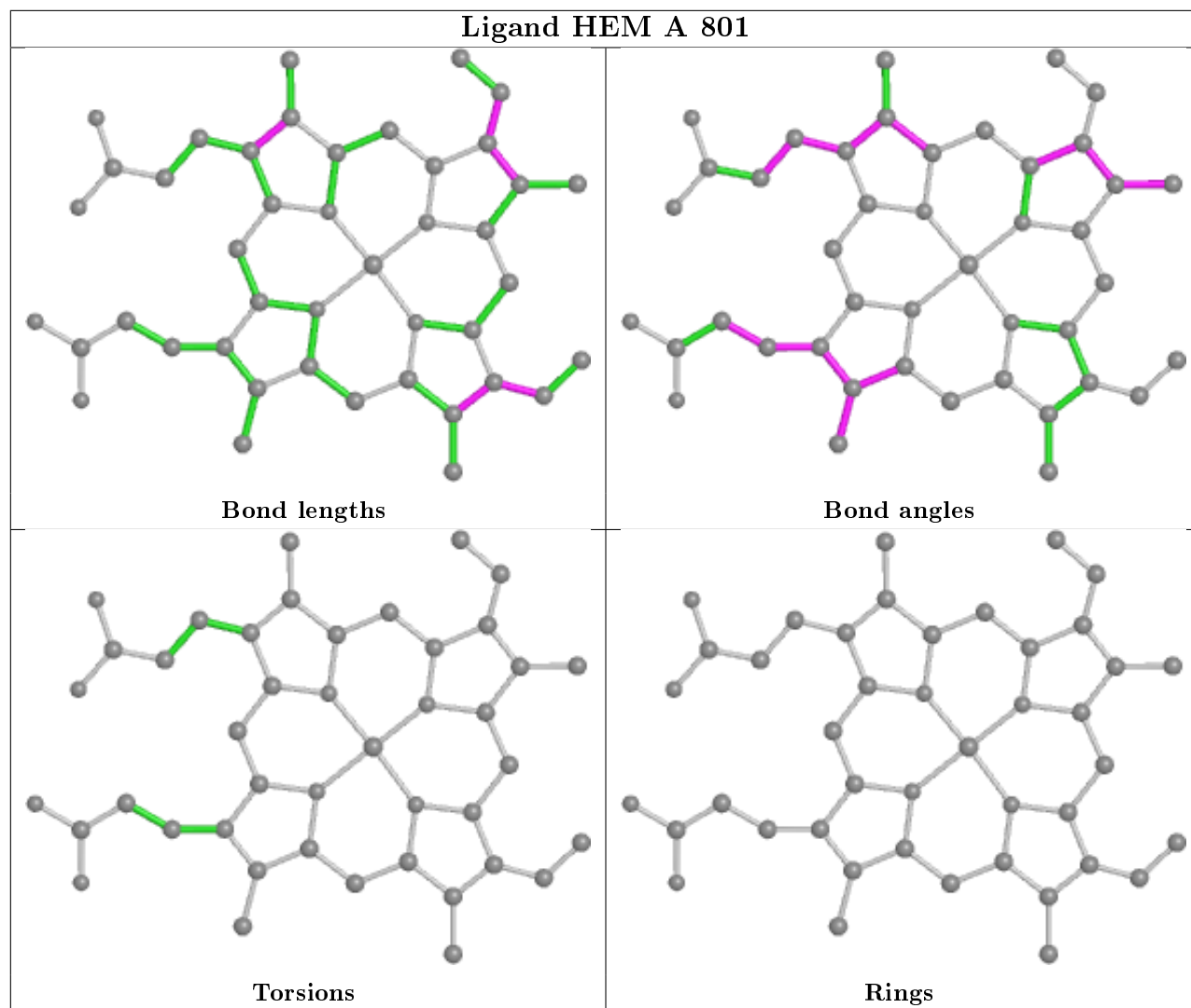
There are no ring outliers.

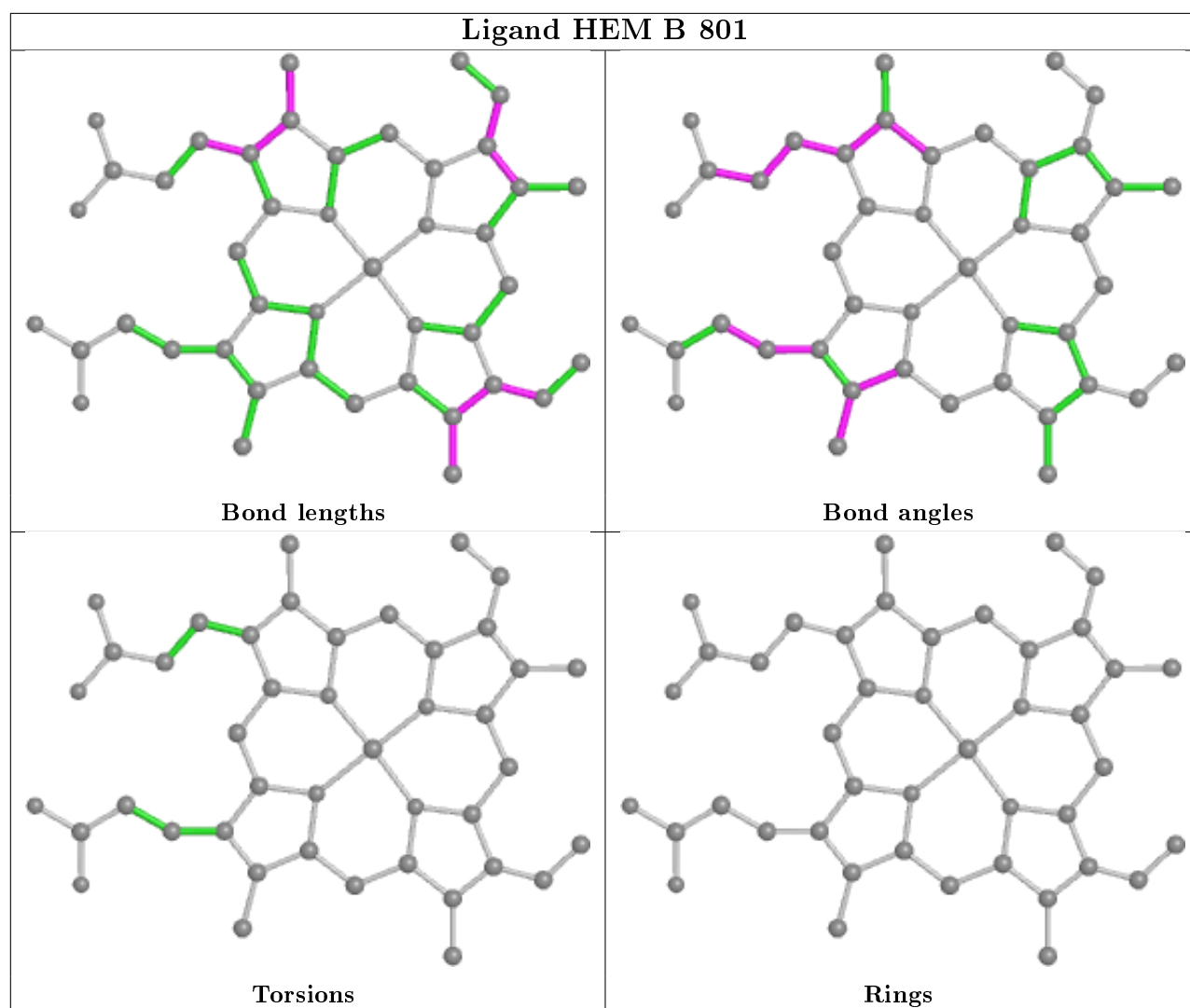
3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	803	2KJ	1	0
2	A	801	HEM	3	0
2	B	801	HEM	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 ⓘ

### 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, 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	407/422 (96%)	0.50	51 (12%) 3 3	26, 50, 91, 116	0
1	B	411/422 (97%)	0.03	19 (4%) 32 31	25, 39, 66, 85	0
All	All	818/844 (96%)	0.26	70 (8%) 10 9	25, 44, 85, 116	0

All (70) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	300	PHE	8.0
1	A	716	TRP	6.5
1	A	352	ASP	5.6
1	A	488	PRO	5.5
1	A	355	PHE	5.1
1	B	350	THR	4.9
1	B	348	VAL	4.8
1	B	718	GLY	4.7
1	B	619	ARG	4.4
1	A	300	PHE	4.3
1	A	351	LYS	4.2
1	A	388	ILE	4.1
1	A	299	ARG	3.9
1	B	352	ASP	3.7
1	A	507	GLN	3.6
1	A	491	SER	3.6
1	A	715	VAL	3.5
1	A	350	THR	3.4
1	A	514	ARG	3.4
1	A	322	LEU	3.3
1	A	503	GLU	3.3
1	A	392	SER	3.2
1	A	567	VAL	3.2
1	A	389	GLU	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	616	LEU	3.1
1	B	351	LYS	2.9
1	A	390	SER	2.9
1	A	486	LYS	2.9
1	A	712	ASN	2.8
1	A	393	THR	2.8
1	A	713	THR	2.8
1	A	489	ASP	2.8
1	B	338	PRO	2.7
1	A	619	ARG	2.7
1	A	682	PRO	2.7
1	A	678	TRP	2.7
1	A	511	LYS	2.7
1	A	311	VAL	2.6
1	A	391	THR	2.5
1	A	714	HIS	2.5
1	A	667	ARG	2.5
1	B	299	ARG	2.4
1	A	370	LYS	2.3
1	A	385	ASN	2.3
1	B	353	GLN	2.3
1	B	389	GLU	2.3
1	A	584	PHE	2.3
1	A	321	THR	2.3
1	A	386	LYS	2.3
1	A	506	ILE	2.3
1	A	381	LEU	2.2
1	A	508	GLN	2.2
1	A	680	VAL	2.2
1	B	567	VAL	2.2
1	A	338	PRO	2.2
1	B	620	LYS	2.2
1	A	485	TYR	2.2
1	A	551	PHE	2.2
1	B	566	ALA	2.2
1	A	487	GLN	2.1
1	B	615	ASP	2.1
1	A	677	VAL	2.1
1	A	681	PRO	2.1
1	A	353	GLN	2.1
1	A	550	LYS	2.0
1	B	328	GLU	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	677	VAL	2.0
1	A	492	THR	2.0
1	A	591	THR	2.0
1	B	667	ARG	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](#)

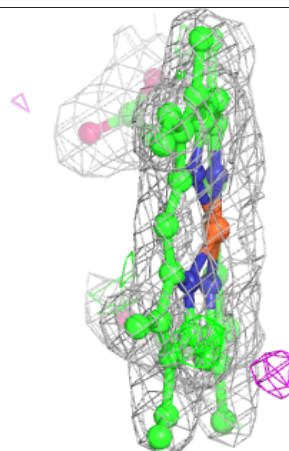
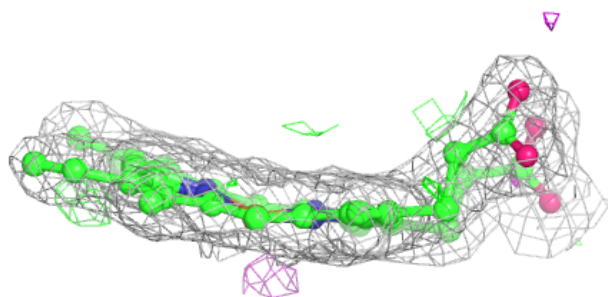
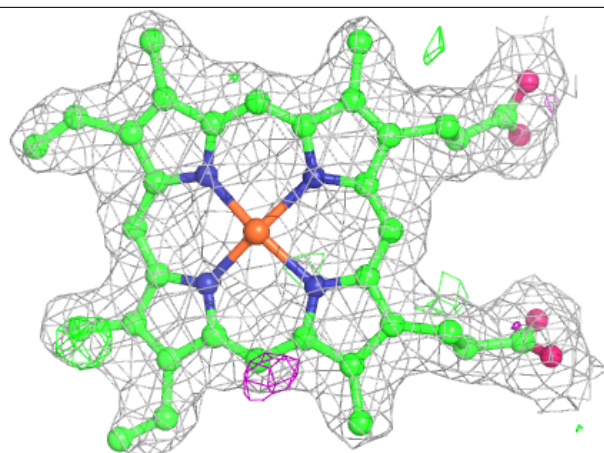
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

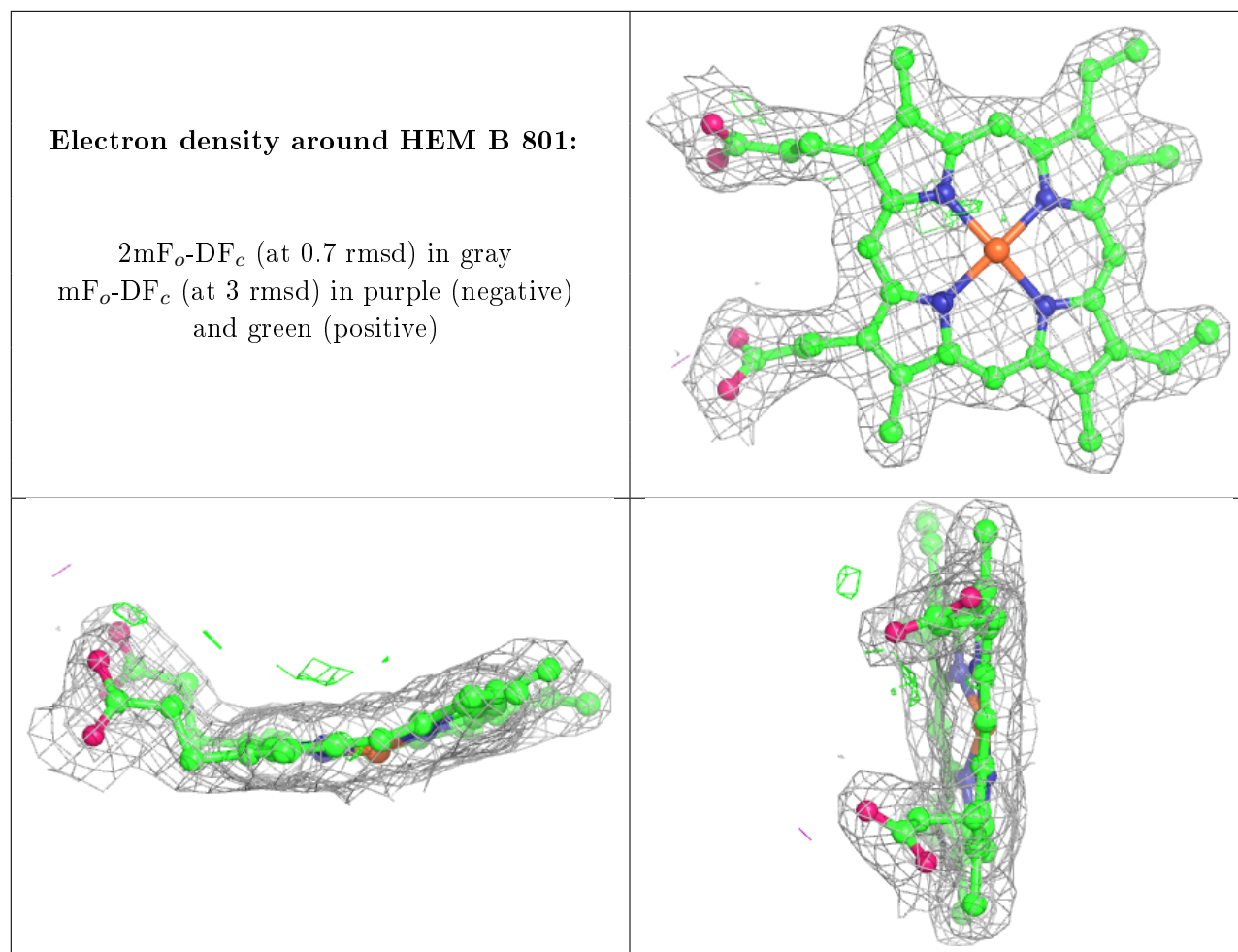
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	TFA	B	804	7/7	0.92	0.20	49,51,54,55	7
4	2KJ	A	803	15/15	0.95	0.26	32,35,44,46	0
4	2KJ	B	803	15/15	0.95	0.18	30,34,44,47	0
3	H4B	A	802	17/17	0.96	0.17	26,30,35,36	0
5	TFA	A	804	7/7	0.97	0.17	60,60,61,61	7
3	H4B	B	802	17/17	0.97	0.14	24,28,32,34	0
2	HEM	A	801	43/43	0.98	0.19	26,31,39,46	0
2	HEM	B	801	43/43	0.98	0.15	23,28,36,44	0
6	ZN	A	805	1/1	0.99	0.07	37,37,37,37	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 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)





## 6.5 Other polymers [i](#)

There are no such residues in this entry.