



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:03 PM GMT

PDB ID : 4KCM  
Title : Structure of neuronal nitric oxide synthase heme domain in complex with N-(4-(2-(ethyl(3-(thiophene-2-carboximidamido)benzyl)amino)ethyl)phenyl)thiophene-2-carboximidamide  
Authors : Li, H.; Poulos, T.L.  
Deposited on : 2013-04-24  
Resolution : 2.07 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  
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

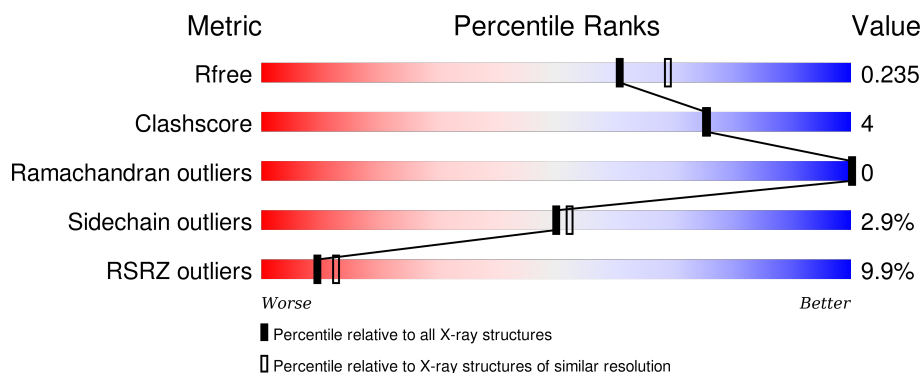
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.07 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))
$R_{free}$	91344	4546 (2.10-2.06)
Clashscore	102246	5101 (2.10-2.06)
Ramachandran outliers	100387	5048 (2.10-2.06)
Sidechain outliers	100360	5049 (2.10-2.06)
RSRZ outliers	91569	4556 (2.10-2.06)

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. 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>85%</div> <div>10%</div> <div>• •</div> </div>
1	B	422	<div> <div>7%</div> <div>87%</div> <div>10%</div> <div>•</div> </div>

## 2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 7081 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.

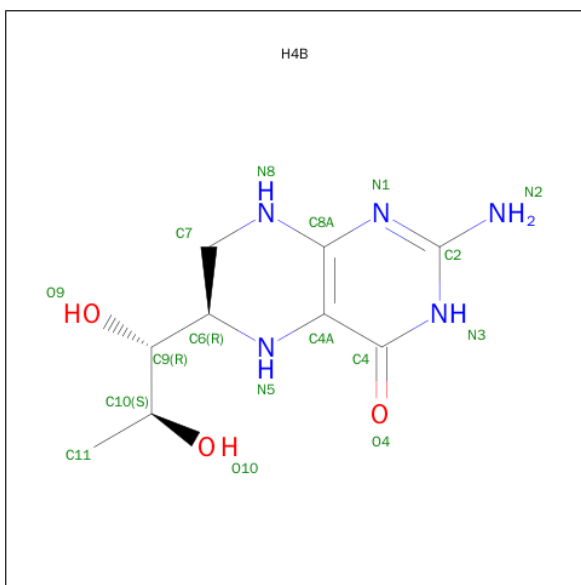
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	407	Total	C	N	O	S	0	2	0
			3322	2127	567	607	21			
1	B	411	Total	C	N	O	S	0	0	0
			3345	2140	574	610	21			

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



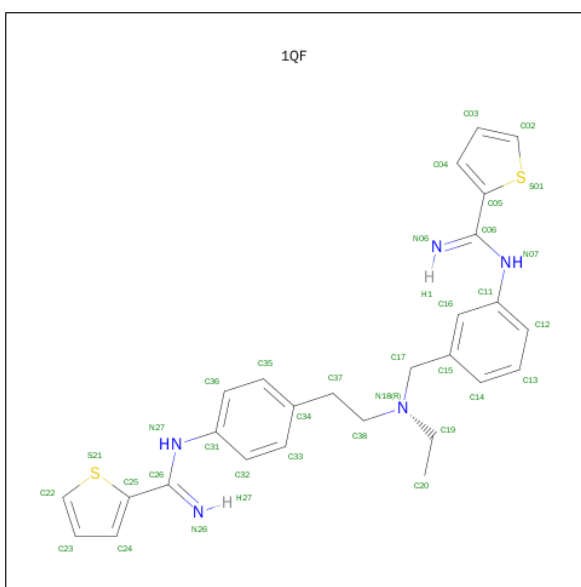
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- 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 17	C 9	N 5	O 3	0	0
3	B	1	Total 17	C 9	N 5	O 3	0	0

- Molecule 4 is N-(4-{2-[ETHYL(3-{[(E)-IMINO(THIOPHEN-2-YL)METHYL]AMINO}BENZYL)AMINO]ETHYL}PHENYL)THIOPHENE-2-CARBOXIMIDAMIDE (three-letter code: 1QF) (formula: C<sub>27</sub>H<sub>29</sub>N<sub>5</sub>S<sub>2</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	S	0	0
			34	27	5	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	N	S	0	0
			34	27	5	2		

- Molecule 5 is BROMIDE ION (three-letter code: BR) (formula: Br).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Br	0	0
			1	1		
5	A	1	Total	Br	0	0
			1	1		

- 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		

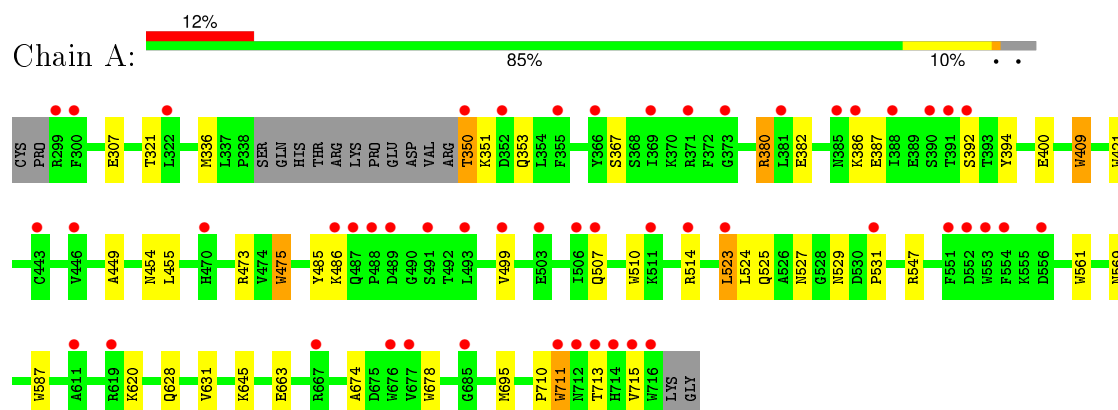
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	92	Total	O	0	0
			92	92		
7	B	131	Total	O	0	0
			131	131		

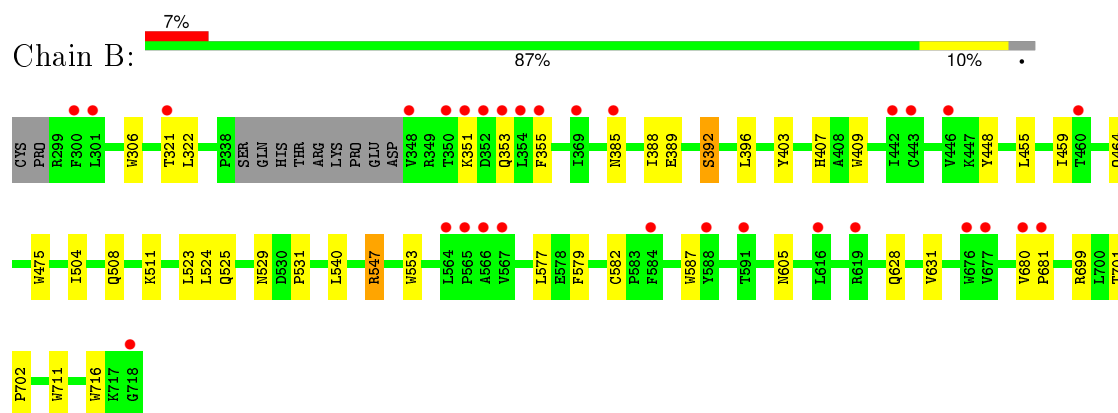
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ( $\text{RSRZ} > 2$ ). 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, $\alpha$ , $\beta$ , $\gamma$	51.74Å 110.79Å 165.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	92.19 – 2.07 46.01 – 2.07	Depositor EDS
% Data completeness (in resolution range)	99.0 (92.19-2.07) 99.1 (46.01-2.07)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.96 (at 2.07Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, $R_{free}$	0.191 , 0.236 0.191 , 0.235	Depositor DCC
$R_{free}$ test set	2863 reflections (5.19%)	DCC
Wilson B-factor (Å <sup>2</sup> )	37.7	Xtriage
Anisotropy	0.749	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 46.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 58005 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	7081	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.81% 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, 0.375 respectively for untwinned datasets, and 0.333, 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, 1QF, H4B, BR

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	0.70	5/3421 (0.1%)	0.70	0/4643
1	B	0.74	4/3438 (0.1%)	0.71	0/4661
All	All	0.72	9/6859 (0.1%)	0.70	0/9304

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	561	TRP	CD2-CE2	5.67	1.48	1.41
1	A	409	TRP	CD2-CE2	5.61	1.48	1.41
1	B	711	TRP	CD2-CE2	5.53	1.48	1.41
1	A	510	TRP	CD2-CE2	5.49	1.48	1.41
1	A	475	TRP	CD2-CE2	5.40	1.47	1.41
1	B	553	TRP	CD2-CE2	5.38	1.47	1.41
1	B	306	TRP	CD2-CE2	5.15	1.47	1.41
1	A	711	TRP	CD2-CE2	5.07	1.47	1.41
1	B	716	TRP	CD2-CE2	5.03	1.47	1.41

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts ⓘ

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



Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3322	0	3234	23	0
1	B	3345	0	3259	23	0
2	A	43	0	30	4	0
2	B	43	0	30	5	0
3	A	17	0	15	0	0
3	B	17	0	15	0	0
4	A	34	0	27	0	0
4	B	34	0	27	4	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	1	0	0	0	0
7	A	92	0	0	1	0
7	B	131	0	0	3	0
All	All	7081	0	6637	52	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 (52) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:801:HEM:HBC2	2:A:801:HEM:HMC2	1.61	0.83
1:A:350:THR:N	1:A:353:GLN:HE21	1.82	0.77
1:B:351:LYS:HE2	1:B:389:GLU:HA	1.74	0.69
1:A:351:LYS:HE2	1:A:392:SER:HB3	1.75	0.68
1:A:336:MET:CE	1:A:678:TRP:HZ2	2.12	0.62
1:B:388:ILE:O	1:B:392:SER:N	2.31	0.62
1:B:605:ASN:ND2	7:B:1029:HOH:O	2.27	0.61
1:A:523:LEU:HD22	1:A:531:PRO:HB2	1.85	0.59
1:B:587:TRP:H	2:B:801:HEM:HAB	1.68	0.58
2:A:801:HEM:HHC	2:A:801:HEM:HBB2	1.87	0.57
1:A:387:GLU:OE1	1:A:394:TYR:HA	2.05	0.57
2:B:801:HEM:O1D	4:B:803:1QF:H9	2.07	0.54
1:A:350:THR:N	1:A:353:GLN:NE2	2.54	0.54
1:A:473:ARG:NH2	1:A:710:PRO:HD3	2.24	0.53
1:A:380:ARG:HD3	1:A:400:GLU:OE1	2.08	0.53
1:B:322:LEU:HD13	1:B:699:ARG:HH21	1.73	0.53
1:A:631:VAL:HG11	1:B:628:GLN:HG2	1.93	0.51
1:B:475:TRP:HB2	1:B:523:LEU:HB3	1.92	0.51
1:B:525:GLN:HG3	1:B:529:ASN:O	2.12	0.50
1:A:499:VAL:N	7:A:988:HOH:O	2.44	0.49
1:A:307:GLU:HG3	7:B:913:HOH:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:620:LYS:HB2	1:A:620:LYS:NZ	2.28	0.49
2:B:801:HEM:C4B	4:B:803:1QF:S21	3.08	0.47
1:A:628:GLN:HG3	1:B:631:VAL:HG11	1.97	0.46
1:B:351:LYS:HE3	1:B:392:SER:OG	2.15	0.46
1:B:459:ILE:HD11	1:B:582:CYS:HB2	1.97	0.45
1:A:674:ALA:HB3	1:A:695:MET:HB3	1.99	0.45
2:A:801:HEM:CMC	2:A:801:HEM:HBC2	2.37	0.45
1:A:382:GLU:OE2	1:A:386:LYS:HE3	2.17	0.45
1:B:464:GLN:HB3	1:B:579:PHE:CE2	2.53	0.44
1:B:701:THR:HA	1:B:702:PRO:C	2.37	0.44
1:B:403:TYR:CE1	1:B:407:HIS:CE1	3.05	0.44
1:B:524:LEU:O	1:B:531:PRO:HA	2.17	0.43
1:A:525:GLN:HG3	1:A:529:ASN:O	2.19	0.43
1:B:355:PHE:CE1	1:B:385:ASN:HB2	2.53	0.43
1:B:547:ARG:HD3	1:B:547:ARG:H	1.83	0.43
2:A:801:HEM:HHC	2:A:801:HEM:CBB	2.49	0.43
1:A:449:ALA:O	1:A:455:LEU:HA	2.19	0.43
1:A:485:TYR:CZ	1:A:514:ARG:HA	2.54	0.43
1:A:336:MET:CE	1:A:678:TRP:CZ2	2.99	0.42
1:B:455:LEU:HD23	1:B:587:TRP:HB3	2.00	0.42
1:B:409:TRP:HB2	2:B:801:HEM:CBC	2.49	0.42
2:B:801:HEM:CHC	4:B:803:1QF:S21	3.08	0.42
1:B:396:LEU:HG	1:B:577:LEU:HD12	2.02	0.42
1:A:409:TRP:CE3	1:A:421:TRP:HA	2.54	0.42
1:B:504:ILE:O	1:B:508:GLN:HG2	2.20	0.41
1:A:455:LEU:HD12	1:A:587:TRP:HB3	2.03	0.41
1:B:448:TYR:C	1:B:448:TYR:CD2	2.93	0.41
4:B:803:1QF:H14	7:B:1020:HOH:O	2.21	0.41
1:B:680:VAL:HA	1:B:681:PRO:HD3	1.95	0.41
1:A:524:LEU:O	1:A:531:PRO:HA	2.22	0.40
1:A:475:TRP:CZ3	1:A:711:TRP:HB3	2.57	0.40

There are no symmetry-related clashes.

## 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	405/422 (96%)	395 (98%)	10 (2%)	0	100	100
1	B	407/422 (96%)	396 (97%)	11 (3%)	0	100	100
All	All	812/844 (96%)	791 (97%)	21 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	365/377 (97%)	350 (96%)	15 (4%)	37	36
1	B	366/377 (97%)	360 (98%)	6 (2%)	70	75
All	All	731/754 (97%)	710 (97%)	21 (3%)	50	52

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	321	THR
1	A	350	THR
1	A	367	SER
1	A	380	ARG
1	A	454	ASN
1	A	486	LYS
1	A	507	GLN
1	A	523	LEU
1	A	527	ASN
1	A	547	ARG
1	A	569	ASN
1	A	645	LYS
1	A	663	GLU
1	A	713	THR
1	A	715	VAL

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Mol	Chain	Res	Type
1	B	321	THR
1	B	353	GLN
1	B	392	SER
1	B	511	LYS
1	B	540	LEU
1	B	547	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	454	ASN
1	A	507	GLN
1	A	527	ASN
1	A	569	ASN
1	A	605	ASN
1	A	697	ASN
1	B	364	GLN
1	B	385	ASN
1	B	454	ASN
1	B	535	GLN
1	B	601	ASN
1	B	605	ASN
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 ⓘ

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 9 ligands modelled in this entry, 3 are monoatomic - leaving 6 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	A	801	1	30,50,50	2.97	14 (46%)	24,82,82	2.77	8 (33%)
3	H4B	A	802	-	13,18,18	1.05	0	11,26,26	2.65	6 (54%)
4	1QF	A	803	-	35,37,37	1.32	5 (14%)	35,49,49	1.53	2 (5%)
2	HEM	B	801	1	30,50,50	2.79	13 (43%)	24,82,82	2.66	8 (33%)
3	H4B	B	802	-	13,18,18	1.05	1 (7%)	11,26,26	2.21	5 (45%)
4	1QF	B	803	-	35,37,37	1.63	7 (20%)	35,49,49	1.80	5 (14%)

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	A	801	1	-	0/10/54/54	0/0/8/8
3	H4B	A	802	-	-	0/8/17/17	0/2/2/2
4	1QF	A	803	-	-	0/19/27/27	0/4/4/4
2	HEM	B	801	1	-	0/10/54/54	0/0/8/8
3	H4B	B	802	-	-	0/8/17/17	0/2/2/2
4	1QF	B	803	-	-	0/19/27/27	0/4/4/4

All (40) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	HEM	C3D-C4D	-4.55	1.45	1.51
2	A	801	HEM	C3B-C4B	-3.71	1.48	1.51
4	B	803	1QF	C26-N27	-3.27	1.32	1.39
4	A	803	1QF	C11-N07	-3.18	1.35	1.41
4	B	803	1QF	C11-N07	-3.15	1.35	1.41
4	B	803	1QF	C06-N07	-2.74	1.33	1.39
4	A	803	1QF	C26-N27	-2.71	1.33	1.39
2	A	801	HEM	C2D-C3D	-2.64	1.46	1.54
4	A	803	1QF	C06-N07	-2.62	1.33	1.39
4	B	803	1QF	C31-N27	-2.38	1.37	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	801	HEM	C3B-C4B	-2.37	1.49	1.51
2	B	801	HEM	C3D-C4D	-2.34	1.48	1.51
2	B	801	HEM	C2D-C3D	-2.29	1.47	1.54
2	B	801	HEM	C2C-C1C	-2.09	1.48	1.52
2	A	801	HEM	C3C-CAC	-2.04	1.47	1.51
3	B	802	H4B	C4-N3	2.09	1.37	1.33
2	B	801	HEM	FE-NB	2.09	2.08	1.97
4	B	803	1QF	C05-S01	2.13	1.74	1.72
2	B	801	HEM	C1A-CHA	2.34	1.46	1.39
2	A	801	HEM	FE-ND	2.43	2.10	1.97
2	A	801	HEM	C4A-CHB	2.47	1.46	1.39
2	B	801	HEM	FE-ND	2.50	2.10	1.97
2	A	801	HEM	FE-NB	2.69	2.11	1.97
4	A	803	1QF	C25-C26	2.77	1.49	1.45
2	A	801	HEM	CHD-C1D	2.89	1.46	1.38
2	A	801	HEM	C2A-C3A	2.98	1.46	1.37
2	A	801	HEM	CHC-C4B	3.04	1.47	1.38
2	B	801	HEM	CHC-C4B	3.20	1.47	1.38
2	B	801	HEM	C2A-C3A	3.31	1.47	1.37
4	A	803	1QF	C05-C06	3.59	1.50	1.45
4	B	803	1QF	C05-C06	3.78	1.50	1.45
2	B	801	HEM	CHC-C1C	3.78	1.45	1.36
2	B	801	HEM	CHD-C4C	4.25	1.46	1.36
2	A	801	HEM	CHD-C4C	4.47	1.46	1.36
4	B	803	1QF	C25-C26	4.53	1.51	1.45
2	A	801	HEM	CHC-C1C	4.90	1.47	1.36
2	A	801	HEM	C4C-NC	7.21	1.44	1.36
2	B	801	HEM	C4C-NC	7.68	1.45	1.36
2	A	801	HEM	C1C-NC	8.17	1.46	1.36
2	B	801	HEM	C1C-NC	8.28	1.46	1.36

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	803	1QF	C23-C22-S21	-6.05	106.53	113.23
4	B	803	1QF	C03-C02-S01	-6.02	106.57	113.23
4	A	803	1QF	C23-C22-S21	-5.61	107.03	113.23
2	A	801	HEM	CBA-CAA-C2A	-5.54	102.60	112.53
4	A	803	1QF	C03-C02-S01	-5.52	107.12	113.23
3	A	802	H4B	N3-C2-N1	-3.82	119.27	125.53
2	B	801	HEM	C3C-CAC-CBC	-3.50	119.08	124.46
2	B	801	HEM	CBA-CAA-C2A	-3.45	106.35	112.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	HEM	C3B-CAB-CBB	-2.85	120.09	124.46
2	A	801	HEM	C3C-CAC-CBC	-2.71	120.30	124.46
3	B	802	H4B	N3-C2-N1	-2.71	121.10	125.53
4	B	803	1QF	C32-C33-C34	-2.30	117.90	121.04
4	B	803	1QF	C17-N18-C38	-2.18	106.72	111.28
4	B	803	1QF	C05-C06-N07	2.05	120.92	115.07
3	A	802	H4B	N2-C2-N3	2.13	120.73	117.20
2	A	801	HEM	CAA-C2A-C1A	2.31	129.52	127.01
3	B	802	H4B	C2-N1-C8A	2.36	119.84	114.54
3	B	802	H4B	N2-C2-N1	2.87	121.95	117.20
3	B	802	H4B	C4-N3-C2	3.21	120.40	115.94
2	A	801	HEM	CMD-C2D-C3D	3.22	128.60	114.35
2	B	801	HEM	CMD-C2D-C3D	3.34	129.13	114.35
3	A	802	H4B	C4-C4A-C8A	3.56	117.79	114.56
3	A	802	H4B	C2-N1-C8A	3.61	122.65	114.54
3	A	802	H4B	C4-N3-C2	3.74	121.12	115.94
3	A	802	H4B	C4A-C8A-N8	3.84	122.95	118.43
2	A	801	HEM	CAD-C3D-C4D	3.92	126.31	112.47
3	B	802	H4B	C4-C4A-C8A	4.06	118.24	114.56
2	B	801	HEM	CMC-C2C-C3C	4.07	126.69	116.53
2	B	801	HEM	CAD-C3D-C4D	4.77	129.29	112.47
2	A	801	HEM	CMC-C2C-C3C	5.21	129.53	116.53
2	B	801	HEM	CAD-C3D-C2D	5.34	128.58	113.22
2	A	801	HEM	CMB-C2B-C3B	5.47	130.18	116.53
2	A	801	HEM	CAD-C3D-C2D	5.79	129.85	113.22
2	B	801	HEM	CMB-C2B-C3B	6.24	132.11	116.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	HEM	4	0
2	B	801	HEM	5	0
4	B	803	1QF	4	0

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

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%)	0.71	51 (12%) 5 6	30, 54, 93, 125	0
1	B	411/422 (97%)	0.46	30 (7%) 18 24	31, 46, 74, 96	0
All	All	818/844 (96%)	0.58	81 (9%) 9 13	30, 49, 88, 125	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	300	PHE	8.5
1	A	355	PHE	5.8
1	A	715	VAL	5.4
1	A	716	TRP	4.9
1	A	713	THR	4.9
1	B	348	VAL	4.8
1	B	350	THR	4.6
1	A	506	ILE	4.4
1	B	352	ASP	4.3
1	A	386	LYS	4.2
1	B	718	GLY	3.9
1	A	488	PRO	3.6
1	A	511	LYS	3.6
1	A	299	ARG	3.6
1	B	355	PHE	3.5
1	A	507	GLN	3.4
1	A	712	ASN	3.3
1	A	373	GLY	3.2
1	B	351	LYS	3.1
1	B	567	VAL	3.1
1	B	353	GLN	3.1
1	B	301	LEU	3.1
1	A	371	ARG	3.0
1	A	486	LYS	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	714	HIS	3.0
1	A	388	ILE	2.9
1	A	381	LEU	2.9
1	B	566	ALA	2.8
1	A	366	TYR	2.7
1	A	491	SER	2.7
1	B	619	ARG	2.7
1	B	677	VAL	2.6
1	A	390	SER	2.6
1	A	392	SER	2.6
1	A	487	GLN	2.6
1	A	493	LEU	2.6
1	A	514	ARG	2.6
1	B	616	LEU	2.6
1	A	503	GLU	2.5
1	B	321	THR	2.5
1	A	553	TRP	2.5
1	A	677	VAL	2.5
1	A	322	LEU	2.5
1	B	354	LEU	2.5
1	A	470	HIS	2.5
1	A	385	ASN	2.5
1	A	556	ASP	2.5
1	B	680	VAL	2.4
1	B	565	PRO	2.4
1	A	531	PRO	2.4
1	B	676	TRP	2.4
1	A	554	PHE	2.4
1	A	391	THR	2.4
1	A	551	PHE	2.4
1	B	591	THR	2.4
1	A	446	VAL	2.3
1	A	369	ILE	2.3
1	A	523	LEU	2.3
1	A	552	ASP	2.3
1	B	584	PHE	2.3
1	A	300	PHE	2.2
1	A	685	GLY	2.2
1	B	443	CYS	2.2
1	A	711	TRP	2.2
1	B	460	THR	2.2
1	A	443	CYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	676	TRP	2.2
1	B	446	VAL	2.2
1	A	489	ASP	2.1
1	A	619	ARG	2.1
1	A	352	ASP	2.1
1	B	681	PRO	2.1
1	A	499	VAL	2.1
1	B	369	ILE	2.1
1	B	442	ILE	2.1
1	B	385	ASN	2.1
1	A	667	ARG	2.0
1	A	611	ALA	2.0
1	B	588	TYR	2.0
1	B	564	LEU	2.0
1	A	350	THR	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	HEM	B	801	43/43	0.98	0.21	1.36	31,35,39,45	0
3	H4B	B	802	17/17	0.98	0.21	1.25	32,36,40,43	0
4	1QF	A	803	34/34	0.92	0.18	1.06	31,55,104,109	0
2	HEM	A	801	43/43	0.98	0.18	1.00	29,35,39,43	0
3	H4B	A	802	17/17	0.97	0.18	0.64	32,35,39,40	0
5	BR	A	804	1/1	0.99	0.14	0.41	43,43,43,43	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
5	BR	B	804	1/1	0.98	0.14	0.08	48,48,48,48	0
4	1QF	B	803	34/34	0.93	0.17	-0.16	30,55,81,88	0
6	ZN	A	805	1/1	0.99	0.12	-0.47	43,43,43,43	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.