



# Full wwPDB X-ray Structure Validation Report i

Feb 26, 2014 – 04:34 PM GMT

PDB ID : 3JWZ  
Title : Structure of endothelial nitric oxide synthase heme domain complexed with  
N1-[(3' S,4' R)-4'-((6''-amino-4''-methylpyridin-2''-yl)methyl)pyrrolidin-3'-yl]-  
N2-(3'-fluorophenethyl)ethane-1,2-diaminetetrahydrochloride  
Authors : Delker, S.L.; Li, H.; Poulos, T.L.  
Deposited on : 2009-09-18  
Resolution : 2.40 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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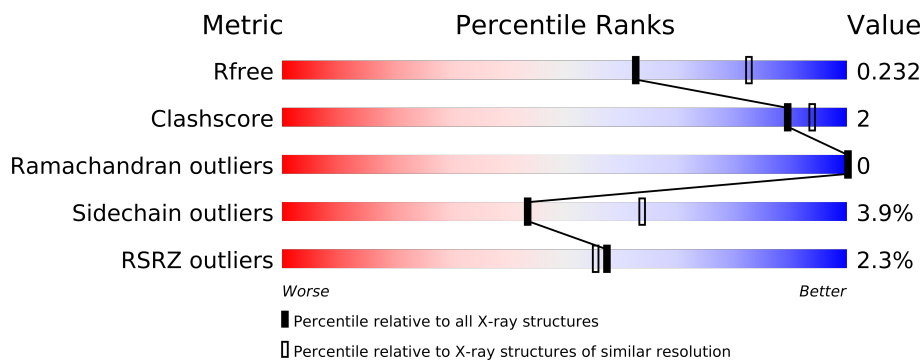
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.40 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))
$R_{free}$	66092	2207 (2.40-2.40)
Clashscore	79885	2789 (2.40-2.40)
Ramachandran outliers	78287	2736 (2.40-2.40)
Sidechain outliers	78261	2737 (2.40-2.40)
RSRZ outliers	66119	2210 (2.40-2.40)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	444	
1	B	444	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
6	J14	B	800	-	X

## 2 Entry composition i

There are 9 unique types of molecules in this entry. The entry contains 6922 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

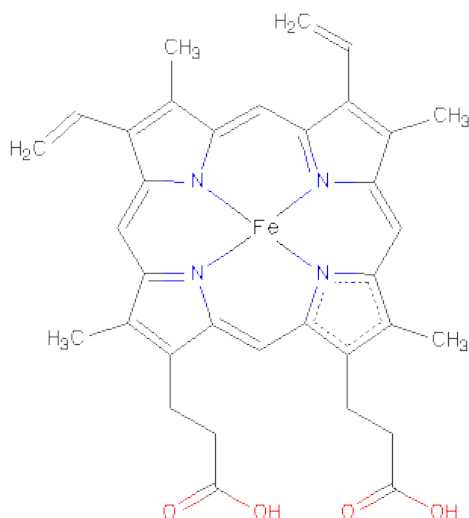
- Molecule 1 is a protein called Nitric oxide synthase, endothelial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	405	Total	C	N	O	S	0	0	0
			3216	2046	565	589	16			
1	B	403	Total	C	N	O	S	0	0	0
			3205	2039	563	587	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	100	ARG	CYS	SEE REMARK 999	UNP P29473
B	100	ARG	CYS	SEE REMARK 999	UNP P29473

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



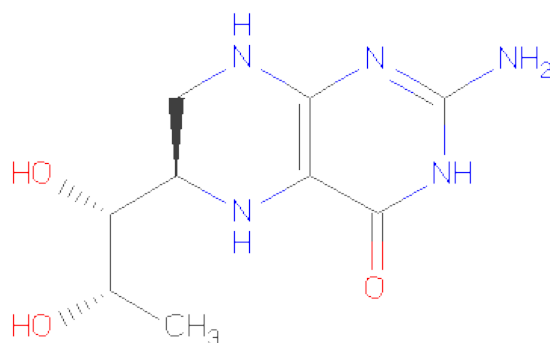
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

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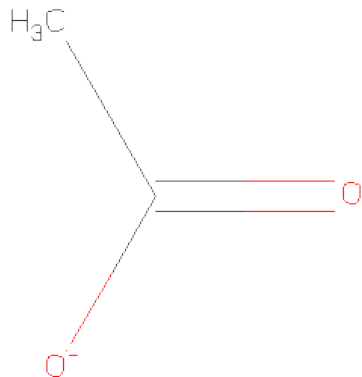
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

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



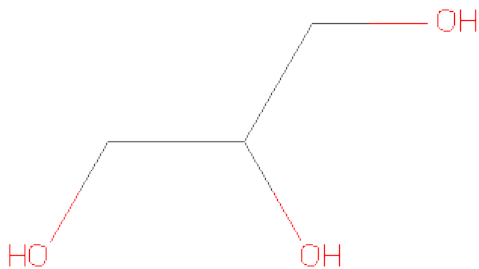
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			17	9	5	3		
3	B	1	Total	C	N	O	0	0
			17	9	5	3		

- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula:  $C_2H_3O_2$ ).



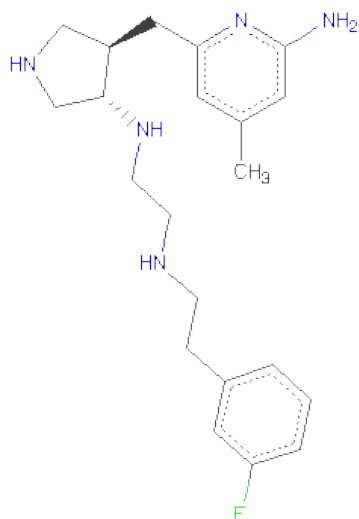
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		
4	B	1	Total	C	O	0	0
			4	2	2		

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



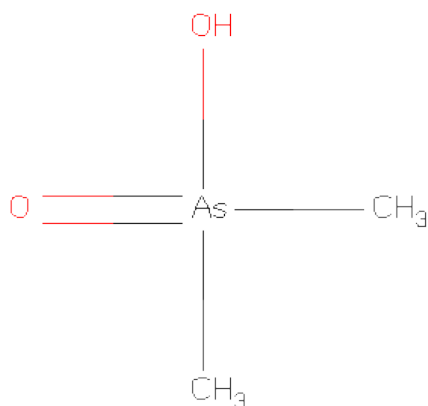
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is N-{(3S,4R)-4-[(6-AMINO-4-METHYLPYRIDIN-2-YL)METHYL]PYRROLIDIN-3-YL}-N'-[2-(3-FLUOROPHENYL)ETHYL]ETHANE-1,2-DIAMINE (three-letter code: J14) (formula: C<sub>21</sub>H<sub>30</sub>FN<sub>5</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	F	N	0	0
			27	21	1	5		
6	B	1	Total	C	F	N	0	0
			27	21	1	5		

- Molecule 7 is CACODYLIC ACID (three-letter code: CAD) (formula: C<sub>2</sub>H<sub>7</sub>AsO<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total 3	As 1	C 2	0	0
7	B	1	Total 3	As 1	C 2	0	0

- Molecule 8 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 9 is water.

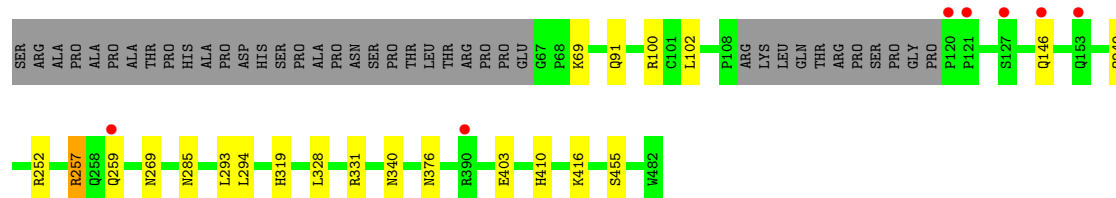
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	164	Total 164	O 164	0	0
9	B	136	Total 136	O 136	0	0

### 3 Residue-property plots

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 ( $RSRZ > 2$ ). 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.

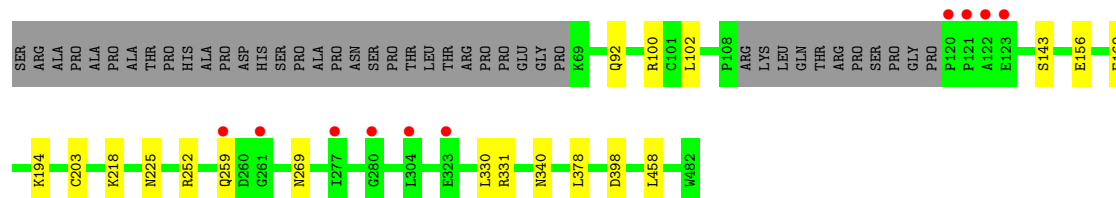
- Molecule 1: Nitric oxide synthase, endothelial

Chain A: 



- Molecule 1: Nitric oxide synthase, endothelial

Chain B: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	58.01Å 106.49Å 156.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.65 – 2.40 46.65 – 2.40	Depositor EDS
% Data completeness (in resolution range)	97.3 (46.65-2.40) 97.3 (46.65-2.40)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.04 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.5.0089	Depositor
R, $R_{free}$	0.172 , 0.228 0.183 , 0.232	Depositor DCC
$R_{free}$ test set	1879 reflections (5.23%)	DCC
Wilson B-factor (Å <sup>2</sup> )	36.2	Xtriage
Anisotropy	0.340	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 16.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 37777 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6922	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.06% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, ZN, H4B, ACT, HEM, J14, CAD

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	0.57	0/3307	0.64	0/4506
1	B	0.59	1/3295 (0.0%)	0.63	0/4489
All	All	0.58	1/6602 (0.0%)	0.64	0/8995

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	203	CYS	CB-SG	-5.69	1.72	1.81

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3216	0	0	10	0
1	B	3205	0	0	3	1
2	A	43	0	30	2	0
2	B	43	0	30	1	0
3	A	17	0	15	0	0
3	B	17	0	15	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	4	0	3	0	0
4	B	4	0	3	0	0
5	A	6	0	8	0	0
5	B	6	0	8	0	0
6	A	27	0	30	3	0
6	B	27	0	30	0	0
7	A	3	0	0	0	0
7	B	3	0	0	0	0
8	B	1	0	0	0	0
9	A	164	0	0	3	1
9	B	136	0	0	1	0
All	All	6922	0	172	15	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 2.

All (15) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:410:HIS:ND1	9:A:1146:HOH:O	2.18	0.77
2:A:500:HEM:HBB2	2:A:500:HEM:HHC	1.88	0.55
1:B:252:ARG:NH1	1:B:269:ASN:OD1	2.41	0.54
1:A:249:GLN:NE2	6:A:800:J14:N1'	2.57	0.53
1:A:252:ARG:NH1	1:A:269:ASN:OD1	2.43	0.51
1:B:252:ARG:NH2	9:B:1108:HOH:O	2.44	0.49
1:A:249:GLN:NE2	6:A:800:J14:H2'A	2.31	0.46
1:A:455:SER:OG	1:B:398:ASP:OD2	2.35	0.44
2:A:500:HEM:CBB	2:A:500:HEM:HHC	2.49	0.43
2:B:500:HEM:HHC	2:B:500:HEM:HBB2	2.00	0.43
1:A:319:HIS:NE2	1:A:403:GLU:OE1	2.52	0.42
1:A:376:ASN:ND2	9:A:1005:HOH:O	2.52	0.42
1:A:257:ARG:CG	1:A:257:ARG:NH1	2.83	0.42
1:A:249:GLN:NE2	6:A:800:J14:C2'	2.83	0.41
1:A:285:ASN:CB	9:A:1012:HOH:O	2.69	0.41

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	Distance(Å)	Clash(Å)
1:B:169:GLU:OE1	9:A:1149:HOH:O[2_555]	2.18	0.02

## 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	401/444 (90%)	394 (98%)	7 (2%)	0	100	100
1	B	399/444 (90%)	388 (97%)	11 (3%)	0	100	100
All	All	800/888 (90%)	782 (98%)	18 (2%)	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	344/377 (91%)	331 (96%)	13 (4%)	44	65
1	B	343/377 (91%)	329 (96%)	14 (4%)	41	61
All	All	687/754 (91%)	660 (96%)	27 (4%)	43	64

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	69	LYS
1	A	91	GLN
1	A	100	ARG
1	A	102	LEU
1	A	146	GLN
1	A	257	ARG
1	A	259	GLN
1	A	293	LEU
1	A	294	LEU
1	A	328	LEU
1	A	331	ARG

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Mol	Chain	Res	Type
1	A	340	ASN
1	A	416	LYS
1	B	92	GLN
1	B	100	ARG
1	B	102	LEU
1	B	143	SER
1	B	156	GLU
1	B	194	LYS
1	B	218	LYS
1	B	225	ASN
1	B	259	GLN
1	B	330	LEU
1	B	331	ARG
1	B	340	ASN
1	B	378	LEU
1	B	458	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA chains 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 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	$\# Z  > 2$
2	HEM	A	500	1	49,50,50	2.36	16 (32%)	46,82,82	2.13	12 (26%)
3	H4B	A	600	-	18,18,18	0.78	0	24,26,26	1.64	4 (16%)
6	J14	A	800	-	29,29,29	0.70	0	38,38,38	1.38	6 (15%)
4	ACT	A	860	-	1,3,3	0.95	0	0,3,3	0.00	-
5	GOL	A	880	-	5,5,5	0.41	0	5,5,5	0.36	0
7	CAD	A	950	-	2,2,4	0.10	0	0,1,6	0.00	-
2	HEM	B	500	1	49,50,50	2.33	14 (28%)	46,82,82	1.92	6 (13%)
3	H4B	B	600	-	18,18,18	0.98	1 (5%)	24,26,26	1.66	5 (20%)
6	J14	B	800	-	29,29,29	0.67	0	38,38,38	1.48	6 (15%)
4	ACT	B	860	-	1,3,3	1.02	0	0,3,3	0.00	-
5	GOL	B	880	-	5,5,5	0.43	0	5,5,5	0.33	0
7	CAD	B	950	-	2,2,4	0.10	0	0,1,6	0.00	-

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	500	1	-	0/14/114/114	0/0/8/8
3	H4B	A	600	-	-	0/8/17/17	0/0/2/2
6	J14	A	800	-	-	0/13/23/23	0/3/3/3
4	ACT	A	860	-	-	0/0/0/0	0/0/0/0
5	GOL	A	880	-	-	0/4/4/4	0/0/0/0
7	CAD	A	950	-	-	0/0/0/0	0/0/0/0
2	HEM	B	500	1	-	0/14/114/114	0/0/8/8
3	H4B	B	600	-	-	0/8/17/17	0/0/2/2
6	J14	B	800	-	-	0/13/23/23	0/3/3/3
4	ACT	B	860	-	-	0/0/0/0	0/0/0/0
5	GOL	B	880	-	-	0/4/4/4	0/0/0/0
7	CAD	B	950	-	-	0/0/0/0	0/0/0/0

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	HEM	C3B-C2B	-6.11	1.33	1.43
2	B	500	HEM	C3B-C2B	-5.92	1.33	1.43
2	B	500	HEM	C3D-C2D	5.59	1.53	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500	HEM	C3C-C2C	-5.36	1.34	1.43
2	A	500	HEM	C2B-C1B	5.19	1.45	1.44
2	B	500	HEM	FE-NA	5.15	2.14	1.92
2	B	500	HEM	C3B-CAB	4.98	1.56	1.40
2	B	500	HEM	C3C-CAC	4.81	1.55	1.40
2	A	500	HEM	C3C-CAC	4.76	1.55	1.40
2	A	500	HEM	C3B-CAB	4.74	1.55	1.40
2	A	500	HEM	C3C-C2C	-4.68	1.35	1.43
2	A	500	HEM	FE-NA	4.56	2.12	1.92
2	A	500	HEM	C3D-C2D	4.37	1.51	1.43
2	B	500	HEM	C4A-C3A	3.96	1.45	1.40
2	B	500	HEM	C3D-C4D	3.88	1.45	1.44
2	A	500	HEM	C4A-C3A	3.69	1.44	1.40
2	A	500	HEM	C2D-C1D	3.30	1.45	1.44
2	A	500	HEM	FE-NB	3.23	2.09	1.97
2	B	500	HEM	FE-NC	3.00	2.09	1.97
2	A	500	HEM	FE-NC	2.89	2.08	1.97
3	B	600	H4B	C2-N2	2.59	1.36	1.32
2	A	500	HEM	CMC-C2C	2.58	1.55	1.47
2	A	500	HEM	FE-ND	2.56	2.07	1.97
2	B	500	HEM	CMD-C2D	2.47	1.55	1.47
2	B	500	HEM	CHB-C1B	2.44	1.39	1.35
2	B	500	HEM	CMB-C2B	2.41	1.54	1.47
2	A	500	HEM	CMD-C2D	2.33	1.54	1.47
2	B	500	HEM	CHA-C4D	2.32	1.39	1.35
2	A	500	HEM	CHB-C1B	2.22	1.39	1.35
2	A	500	HEM	CMB-C2B	2.18	1.54	1.47
2	B	500	HEM	CMC-C2C	2.13	1.54	1.47

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	HEM	C3B-C4B-NB	-7.56	108.59	114.00
2	B	500	HEM	C3B-C4B-NB	-6.93	109.04	114.00
2	A	500	HEM	C4D-ND-C1D	6.15	111.46	105.16
2	B	500	HEM	CBA-CAA-C2A	-5.74	102.59	112.69
2	B	500	HEM	C4D-ND-C1D	4.63	109.90	105.16
3	A	600	H4B	C4-C4A-C8A	4.56	118.79	114.56
3	B	600	H4B	C4-C4A-C8A	4.27	118.52	114.56
6	B	800	J14	C1-N1-C3'	3.75	119.47	113.21
2	A	500	HEM	CBA-CAA-C2A	-3.71	106.15	112.69
2	A	500	HEM	C2D-C1D-ND	-3.62	108.66	112.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	800	J14	C71-C21-N11	3.47	124.19	116.71
2	A	500	HEM	CBD-CAD-C3D	-3.46	106.82	114.37
6	A	800	J14	C1-N1-C3'	3.14	118.44	113.21
6	A	800	J14	C71-C21-N11	3.03	123.23	116.71
3	B	600	H4B	N8-C8A-N1	2.99	120.20	115.82
6	B	800	J14	C5'-N1'-C2'	2.88	113.20	105.92
3	B	600	H4B	O9-C9-C10	-2.82	104.28	108.86
2	B	500	HEM	CBD-CAD-C3D	-2.80	108.27	114.37
6	B	800	J14	C14-C13-C12	-2.76	119.52	123.32
2	B	500	HEM	CHA-C4D-ND	2.74	128.07	124.31
6	A	800	J14	C5'-N1'-C2'	2.71	112.76	105.92
2	A	500	HEM	CAD-C3D-C4D	2.65	129.28	124.53
2	A	500	HEM	CMA-C3A-C4A	-2.60	124.63	128.62
3	A	600	H4B	N2-C2-N3	2.55	120.67	117.86
2	A	500	HEM	C4A-CHB-C1B	-2.54	124.12	127.47
2	A	500	HEM	C4C-NC-C1C	2.54	108.17	105.53
6	A	800	J14	C2-C1-N1	2.49	115.03	110.97
6	B	800	J14	C61-N11-C21	2.45	119.95	118.23
3	A	600	H4B	C7-C6-C9	-2.30	110.11	113.66
6	A	800	J14	C4'-C5'-N1'	-2.30	101.12	105.75
3	B	600	H4B	C4A-C4-N3	2.30	119.95	114.06
6	B	800	J14	C31-C21-N11	-2.25	120.49	122.99
2	A	500	HEM	O1D-CGD-CBD	-2.23	115.37	123.03
6	A	800	J14	C14-C13-C12	-2.22	120.26	123.32
3	A	600	H4B	C4-C4A-N5	2.15	122.22	119.10
2	A	500	HEM	CHC-C4B-NB	2.14	126.36	124.58
2	A	500	HEM	C3A-C4A-NA	-2.09	107.83	109.41
3	B	600	H4B	C2-N1-C8A	2.08	120.56	117.61
2	B	500	HEM	CHC-C4B-NB	2.08	126.31	124.58

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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	405/444 (91%)	-0.05	7 (1%) 67 65	22, 35, 60, 82	0
1	B	403/444 (90%)	-0.04	10 (2%) 54 52	22, 38, 63, 86	0
All	All	808/888 (90%)	-0.05	17 (2%) 57 58	22, 36, 61, 86	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	120	PRO	7.8
1	B	259	GLN	4.7
1	A	120	PRO	4.1
1	B	121	PRO	4.0
1	A	259	GLN	3.0
1	B	123	GLU	2.9
1	A	146	GLN	2.9
1	B	261	GLY	2.5
1	B	323	GLU	2.4
1	B	304	LEU	2.2
1	B	277	ILE	2.2
1	A	153	GLN	2.2
1	A	390	ARG	2.1
1	A	121	PRO	2.1
1	B	280	GLY	2.1
1	A	127	SER	2.0
1	B	122	ALA	2.0

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates 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. 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
6	J14	B	800	27/27	0.24	2.54	36,48,71,72	0
6	J14	A	800	27/27	0.22	1.80	29,41,65,66	0
5	GOL	B	880	6/6	0.23	1.75	40,45,47,50	0
2	HEM	A	500	43/43	0.20	1.09	21,26,35,47	0
3	H4B	B	600	17/17	0.22	1.05	21,23,25,28	0
2	HEM	B	500	43/43	0.17	0.74	26,29,39,48	0
5	GOL	A	880	6/6	0.19	0.66	46,47,48,48	0
3	H4B	A	600	17/17	0.20	0.25	22,24,27,29	0
8	ZN	B	900	1/1	0.15	0.01	28,28,28,28	0
7	CAD	A	950	3/5	0.10	-0.43	50,50,52,53	0
7	CAD	B	950	3/5	0.11	-1.06	53,53,54,57	0
4	ACT	A	860	4/4	0.11	-1.17	38,39,39,39	0
4	ACT	B	860	4/4	0.09	-1.62	40,41,42,42	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.