



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:22 PM GMT

PDB ID : 4CAM
Title : Structure of rat neuronal nitric oxide synthase heme domain in complex with
7-((3-Fluorophenethylamino)methyl)quinolin-2-amine
Authors : Li, H.; Poulos, T.L.
Deposited on : 2013-10-08
Resolution : 1.83 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

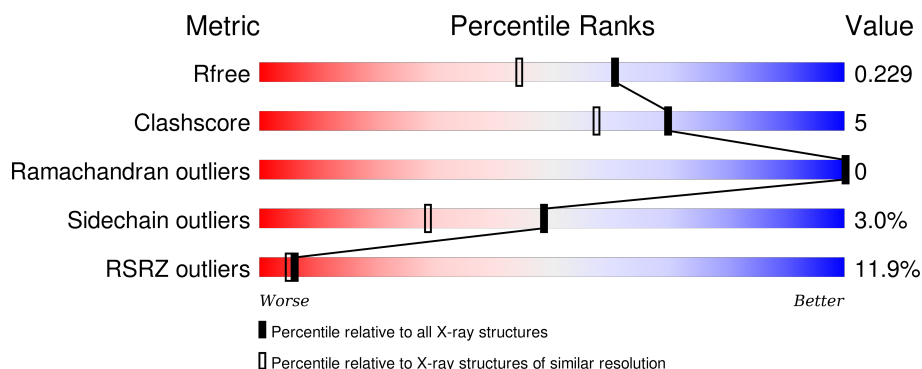
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.83 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	2634 (1.86-1.82)
Clashscore	102246	2862 (1.86-1.82)
Ramachandran outliers	100387	2831 (1.86-1.82)
Sidechain outliers	100360	2832 (1.86-1.82)
RSRZ outliers	91569	2639 (1.86-1.82)

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 ≥ 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>16%</div> <div>83%</div> <div>13%</div> <div>..</div> </div>
1	B	422	<div> <div>7%</div> <div>90%</div> <div>7%</div> <div>..</div> </div>

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 7180 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.

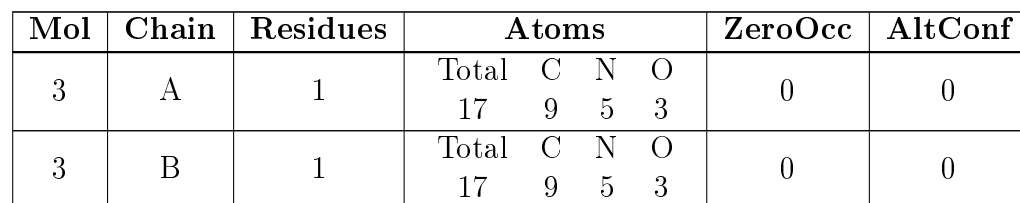
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	407	Total	C	N	O	S	0	4	0
			3331	2133	567	609	22			
1	B	411	Total	C	N	O	S	0	3	0
			3359	2152	574	611	22			

- 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: $C_9H_{15}N_5O_3$).



- M87
-
- Chemical structure of M87, a fluorinated indole derivative. The structure consists of an indole ring system (labeled C02 to C11 and N01) substituted with a 2-amino group (N02) at the 2-position and a 2-(2-fluorophenyl)ethyl group at the 3-position. The ethyl chain is labeled C12 and C13. The phenyl ring is labeled C21 to C26, and the fluorine atom is labeled F23.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	F	N	0	0
			22	18	1	3		



Continued from previous page...

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	F	N	0	0
			22	18	1	3		

- Molecule 5 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



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

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

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

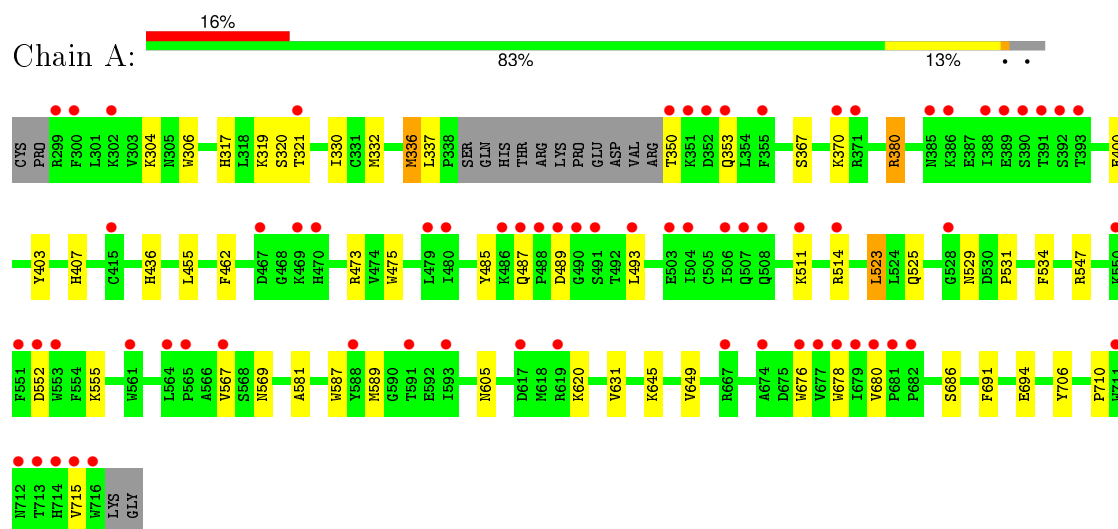
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	141	Total	O	0	0
			141	141		
7	B	176	Total	O	0	0
			176	176		

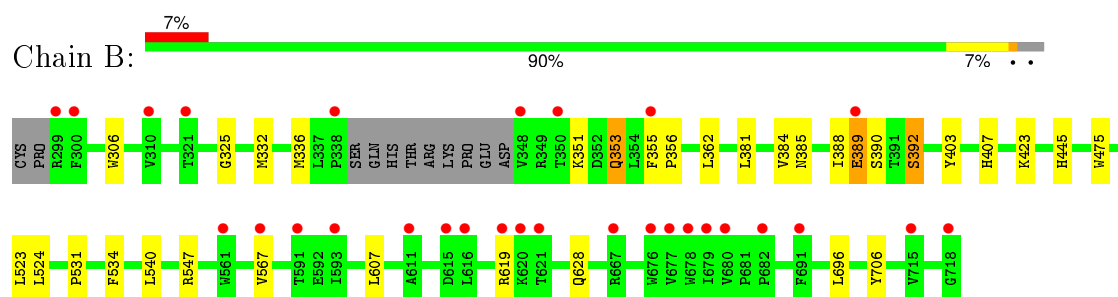
3 Residue-property plots

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, BRAIN



• Molecule 1: NITRIC OXIDE SYNTHASE, BRAIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	51.93Å 111.13Å 164.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.11 – 1.83 39.08 – 1.83	Depositor EDS
% Data completeness (in resolution range)	99.3 (39.11-1.83) 99.4 (39.08-1.83)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 1.83Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.194 , 0.230 0.193 , 0.229	Depositor DCC
R_{free} test set	4196 reflections (5.22%)	DCC
Wilson B-factor (Å ²)	32.2	Xtriage
Anisotropy	0.615	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	2 of 84598 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7180	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.39% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 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: M87, ZN, H4B, HEM, ACT

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	0.52	0/3436	0.65	0/4663
1	B	0.60	0/3462	0.70	0/4694
All	All	0.56	0/6898	0.68	0/9357

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

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	3331	0	3245	36	0
1	B	3359	0	3278	26	0
2	A	43	0	30	2	0
2	B	43	0	30	4	0
3	A	17	0	15	0	0
3	B	17	0	15	0	0
4	A	22	0	18	3	0
4	B	22	0	18	2	0
5	A	4	0	3	0	0
5	B	4	0	3	0	0
6	A	1	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	A	141	0	0	4	0
7	B	176	0	0	0	0
All	All	7180	0	6655	61	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 5.

All (61) 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:336:MET:CE	1:A:678:TRP:HZ2	1.85	0.89
1:A:567:VAL:HG21	4:A:800:M87:C07	2.11	0.81
1:A:336:MET:CE	1:A:678:TRP:CZ2	2.68	0.77
1:B:567:VAL:HG21	4:B:800:M87:C07	2.14	0.76
2:B:750:HEM:HHB	2:B:750:HEM:HHB2	1.68	0.74
1:A:523:LEU:HD22	1:A:531:PRO:HB2	1.70	0.72
1:B:355[A]:PHE:CE1	1:B:385:ASN:HB2	2.29	0.68
4:A:800:M87:H26	7:A:2069:HOH:O	1.95	0.67
1:A:321:THR:HG22	7:A:2006:HOH:O	1.98	0.62
1:A:336:MET:HE3	1:A:678:TRP:HZ2	1.62	0.62
1:A:380:ARG:HD3	1:A:400:GLU:OE2	2.01	0.60
1:A:332:MET:HE2	1:B:696:LEU:HD22	1.84	0.59
1:B:355[B]:PHE:HB2	1:B:356:PRO:CD	2.33	0.58
1:A:332:MET:HE2	1:B:696:LEU:CD2	2.33	0.57
1:B:706:TYR:OH	2:B:750:HEM:O2D	2.20	0.57
1:A:336:MET:HE1	1:A:678:TRP:CZ2	2.39	0.56
1:A:706:TYR:OH	2:A:750:HEM:O2D	2.18	0.56
1:A:620:LYS:HB2	1:A:620:LYS:NZ	2.20	0.55
1:B:351:LYS:HE2	1:B:389:GLU:HA	1.89	0.55
1:B:403:TYR:CE1	1:B:407:HIS:CE1	2.94	0.55
1:A:321:THR:HG21	7:A:2007:HOH:O	2.06	0.55
1:A:631:VAL:HG11	1:B:628:GLN:HG3	1.88	0.55
1:A:473:ARG:NH2	1:A:710:PRO:HD3	2.23	0.53
1:B:362:LEU:HD12	1:B:381:LEU:HD23	1.91	0.53
2:B:750:HEM:HHB	2:B:750:HEM:CBB	2.38	0.52
1:A:485:TYR:HB3	1:A:514:ARG:NH2	2.25	0.52
1:B:475:TRP:HB2	1:B:523:LEU:HB3	1.92	0.51
1:A:605:ASN:ND2	7:A:2111:HOH:O	2.46	0.48
1:A:336:MET:HG3	1:B:306:TRP:NE1	2.29	0.47
1:A:525:GLN:HG3	1:A:529:ASN:O	2.15	0.47
1:A:475:TRP:HB2	1:A:523:LEU:HB3	1.95	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:336:MET:HE3	1:A:678:TRP:CZ2	2.45	0.46
1:B:388:ILE:O	1:B:392:SER:N	2.45	0.46
1:A:462:PHE:HB2	1:A:581:ALA:HB3	1.97	0.45
1:B:325:GLY:O	1:B:332:MET:HG3	2.16	0.45
1:A:306:TRP:CD2	1:B:336:MET:HE3	2.52	0.45
1:A:304:LYS:O	1:A:694:GLU:HG3	2.17	0.45
1:B:355[B]:PHE:HB2	1:B:356:PRO:HD3	1.99	0.44
1:B:524:LEU:HD12	1:B:534:PHE:CD2	2.52	0.44
1:A:436:HIS:CD2	1:A:534:PHE:HE2	2.36	0.44
1:B:353:GLN:HB3	1:B:353:GLN:HE21	1.64	0.43
1:A:330:ILE:HD11	1:B:696:LEU:HB3	2.01	0.43
1:A:686:SER:HA	1:A:691:PHE:CG	2.53	0.43
2:A:750:HEM:C4B	4:A:800:M87:H03	2.54	0.42
1:A:589:MET:HA	1:A:649:VAL:O	2.19	0.42
1:A:403:TYR:CE1	1:A:407:HIS:CE1	3.07	0.42
2:B:750:HEM:CHC	2:B:750:HEM:HBB2	2.40	0.42
1:A:676:TRP:CE2	1:A:680:VAL:HG21	2.55	0.41
1:A:455:LEU:HD12	1:A:587:TRP:HB3	2.01	0.41
1:A:487:GLN:HE21	1:A:493:LEU:HB2	1.85	0.41
1:B:619:ARG:HB2	1:B:619:ARG:HE	1.62	0.41
1:B:445:HIS:C	1:B:445:HIS:CD2	2.93	0.41
1:B:362:LEU:HD11	1:B:384:VAL:HG21	2.01	0.41
1:B:351:LYS:HE3	1:B:392:SER:OG	2.21	0.41
1:B:567:VAL:HG21	4:B:800:M87:C06	2.49	0.41
1:A:350:THR:N	1:A:353:GLN:HE21	2.19	0.41
1:B:607:LEU:HA	1:B:607:LEU:HD23	1.94	0.41
1:A:317:HIS:O	1:A:320:SER:HB3	2.20	0.41
1:B:524:LEU:O	1:B:531:PRO:HA	2.20	0.41
1:A:676:TRP:CZ2	1:A:680:VAL:HG21	2.56	0.41
1:A:475:TRP:CZ2	1:A:531:PRO:HG3	2.56	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	407/422 (96%)	396 (97%)	11 (3%)	0	100	100
1	B	410/422 (97%)	403 (98%)	7 (2%)	0	100	100
All	All	817/844 (97%)	799 (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	367/377 (97%)	352 (96%)	15 (4%)	37	16
1	B	369/377 (98%)	362 (98%)	7 (2%)	65	50
All	All	736/754 (98%)	714 (97%)	22 (3%)	48	29

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

Mol	Chain	Res	Type
1	A	319	LYS
1	A	336	MET
1	A	337	LEU
1	A	367	SER
1	A	370	LYS
1	A	380	ARG
1	A	489	ASP
1	A	511	LYS
1	A	523	LEU
1	A	547	ARG
1	A	552	ASP
1	A	555	LYS
1	A	569	ASN
1	A	645	LYS
1	A	715	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	353	GLN
1	B	389	GLU
1	B	390	SER
1	B	392	SER
1	B	423	LYS
1	B	540	LEU
1	B	547	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	425	GLN
1	A	454	ASN
1	A	569	ASN
1	A	605	ASN
1	A	697	ASN
1	B	364	GLN
1	B	454	ASN
1	B	535	GLN
1	B	601	ASN
1	B	605	ASN
1	B	697	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 1 is monoatomic - leaving 8 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	HEM	A	750	1	30,50,50	2.08	7 (23%)	24,82,82	2.33	9 (37%)
3	H4B	A	760	-	13,18,18	1.36	1 (7%)	11,26,26	2.79	6 (54%)
4	M87	A	800	-	24,24,24	1.14	3 (12%)	31,32,32	1.29	4 (12%)
5	ACT	A	860	-	1,3,3	1.48	0	0,3,3	0.00	-
2	HEM	B	750	1	30,50,50	2.13	7 (23%)	24,82,82	2.57	13 (54%)
3	H4B	B	760	-	13,18,18	1.05	0	11,26,26	2.64	5 (45%)
4	M87	B	800	-	24,24,24	1.14	4 (16%)	31,32,32	1.93	10 (32%)
5	ACT	B	860	-	1,3,3	0.99	0	0,3,3	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	750	1	-	0/10/54/54	0/0/8/8
3	H4B	A	760	-	-	0/8/17/17	0/2/2/2
4	M87	A	800	-	-	0/7/7/7	0/3/3/3
5	ACT	A	860	-	-	0/0/0/0	0/0/0/0
2	HEM	B	750	1	-	0/10/54/54	0/0/8/8
3	H4B	B	760	-	-	0/8/17/17	0/2/2/2
4	M87	B	800	-	-	0/7/7/7	0/3/3/3
5	ACT	B	860	-	-	0/0/0/0	0/0/0/0

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	750	HEM	C2D-C3D	-6.42	1.35	1.54
2	B	750	HEM	C2C-C1C	-6.37	1.40	1.52
2	B	750	HEM	C2D-C3D	-6.34	1.35	1.54
2	A	750	HEM	C2C-C1C	-5.56	1.42	1.52
2	B	750	HEM	C2B-C1B	-2.84	1.42	1.51
2	A	750	HEM	C3D-C4D	-2.75	1.48	1.51

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	750	HEM	C3D-C4D	-2.56	1.48	1.51
2	A	750	HEM	C3B-C4B	-2.48	1.49	1.51
2	A	750	HEM	C2B-C1B	-2.46	1.43	1.51
2	B	750	HEM	C3B-C4B	-2.31	1.49	1.51
2	B	750	HEM	FE-NB	2.07	2.08	1.97
4	B	800	M87	C02-N01	2.10	1.35	1.33
4	B	800	M87	C04-C03	2.11	1.41	1.36
4	B	800	M87	C24-C23	2.16	1.41	1.37
4	A	800	M87	C04-C03	2.21	1.41	1.36
4	A	800	M87	C02-N01	2.42	1.36	1.33
2	A	750	HEM	FE-NB	2.44	2.10	1.97
3	A	760	H4B	C2-N2	2.52	1.39	1.34
4	B	800	M87	C06-C07	2.66	1.42	1.36
4	A	800	M87	C06-C07	2.88	1.42	1.36
2	B	750	HEM	FE-NC	3.60	2.10	1.95
2	A	750	HEM	FE-NC	3.68	2.10	1.95

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	800	M87	C03-C02-N01	-4.49	117.30	122.20
2	B	750	HEM	C1D-CHD-C4C	-3.67	119.68	125.82
3	A	760	H4B	N3-C2-N1	-3.42	119.92	125.53
4	B	800	M87	C11-C08-C09	-3.38	116.61	121.98
2	A	750	HEM	C1D-CHD-C4C	-3.36	120.20	125.82
4	B	800	M87	C14-C21-C22	-3.32	114.99	120.56
4	A	800	M87	C11-C08-C09	-3.26	116.80	121.98
3	B	760	H4B	N3-C2-N1	-3.19	120.30	125.53
2	B	750	HEM	CBA-CAA-C2A	-2.95	107.24	112.53
2	B	750	HEM	C3B-C4B-NB	-2.74	106.39	111.63
4	A	800	M87	C03-C02-N01	-2.71	119.23	122.20
2	A	750	HEM	CBA-CAA-C2A	-2.57	107.93	112.53
2	B	750	HEM	CMA-C3A-C4A	-2.53	124.17	128.36
4	B	800	M87	C24-C23-C22	-2.23	120.47	123.35
4	A	800	M87	C24-C23-C22	-2.20	120.52	123.35
2	B	750	HEM	C3C-CAC-CBC	-2.19	121.09	124.46
2	A	750	HEM	C3B-C4B-NB	-2.19	107.44	111.63
2	B	750	HEM	C3B-CAB-CBB	-2.11	121.22	124.46
2	B	750	HEM	CAA-C2A-C1A	-2.05	124.78	127.01
3	A	760	H4B	N2-C2-N1	2.07	120.63	117.20
4	B	800	M87	F23-C23-C24	2.11	122.05	118.52
2	A	750	HEM	C2D-C3D-C4D	2.13	105.11	101.50

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	800	M87	C13-C14-C21	2.17	117.34	112.83
3	B	760	H4B	C4A-C8A-N8	2.36	121.21	118.43
2	B	750	HEM	C2D-C3D-C4D	2.48	105.71	101.50
4	B	800	M87	C08-C11-N12	2.63	119.41	112.88
4	B	800	M87	C04-C03-C02	2.66	122.16	119.55
3	A	760	H4B	C2-N1-C8A	2.70	120.61	114.54
2	A	750	HEM	CMD-C2D-C3D	2.73	126.45	114.35
2	B	750	HEM	CMD-C2D-C3D	2.97	127.48	114.35
3	A	760	H4B	C4A-C8A-N8	2.98	121.95	118.43
3	B	760	H4B	C2-N1-C8A	3.01	121.30	114.54
4	A	800	M87	C02-N01-C10	3.36	121.79	117.78
3	B	760	H4B	C4-N3-C2	3.55	120.87	115.94
4	B	800	M87	N02-C02-N01	3.66	120.41	118.03
2	A	750	HEM	CMB-C2B-C3B	3.69	125.74	116.53
4	B	800	M87	C02-N01-C10	3.83	122.35	117.78
2	B	750	HEM	CAD-C3D-C4D	4.43	128.08	112.47
3	A	760	H4B	C4-N3-C2	4.44	122.10	115.94
2	A	750	HEM	CAD-C3D-C2D	4.44	125.99	113.22
2	B	750	HEM	CMC-C2C-C3C	4.44	127.62	116.53
2	B	750	HEM	CMB-C2B-C3B	4.47	127.68	116.53
2	B	750	HEM	CAD-C3D-C2D	4.52	126.20	113.22
2	A	750	HEM	CAD-C3D-C4D	4.66	128.90	112.47
2	A	750	HEM	CMC-C2C-C3C	5.04	129.11	116.53
3	A	760	H4B	C4-C4A-C8A	5.53	119.57	114.56
3	B	760	H4B	C4-C4A-C8A	5.79	119.81	114.56

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	750	HEM	2	0
4	A	800	M87	3	0
2	B	750	HEM	4	0
4	B	800	M87	2	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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	407/422 (96%)	0.79	68 (16%) 2 2	26, 49, 86, 117	0
1	B	411/422 (97%)	0.26	29 (7%) 19 17	24, 38, 66, 89	0
All	All	818/844 (96%)	0.53	97 (11%) 6 5	24, 43, 81, 117	0

All (97) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	348	VAL	8.4
1	A	715	VAL	7.7
1	A	488	PRO	7.5
1	B	300	PHE	7.4
1	A	716	TRP	6.7
1	A	355	PHE	5.8
1	A	713	THR	5.4
1	B	619	ARG	5.0
1	B	718	GLY	4.9
1	A	507	GLN	4.8
1	A	352	ASP	4.6
1	A	489	ASP	4.6
1	A	391	THR	4.5
1	A	392	SER	4.4
1	A	486	LYS	4.3
1	A	491	SER	4.2
1	B	350	THR	4.1
1	A	712	ASN	4.1
1	A	487	GLN	4.1
1	B	616	LEU	4.1
1	A	490	GLY	4.1
1	A	388	ILE	3.9
1	A	390	SER	3.9
1	A	677	VAL	3.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	300	PHE	3.4
1	A	551	PHE	3.4
1	A	386	LYS	3.4
1	B	677	VAL	3.4
1	A	503	GLU	3.3
1	B	620	LYS	3.3
1	B	321	THR	3.3
1	A	680	VAL	3.2
1	A	350	THR	3.1
1	A	511	LYS	3.1
1	A	506	ILE	3.1
1	A	676	TRP	3.1
1	A	469	LYS	3.1
1	A	619	ARG	3.1
1	A	528	GLY	3.1
1	A	679	ILE	3.0
1	A	479	LEU	3.0
1	B	611	ALA	2.9
1	A	714	HIS	2.9
1	B	680	VAL	2.9
1	B	355[A]	PHE	2.9
1	B	591	THR	2.9
1	B	615	ASP	2.9
1	A	711	TRP	2.9
1	A	351	LYS	2.8
1	A	593	ILE	2.8
1	A	389	GLU	2.8
1	A	385	ASN	2.8
1	B	389	GLU	2.8
1	A	667	ARG	2.7
1	A	508	GLN	2.7
1	A	678	TRP	2.7
1	B	667	ARG	2.7
1	A	480	ILE	2.7
1	A	588	TYR	2.7
1	A	299	ARG	2.7
1	A	467	ASP	2.6
1	A	353	GLN	2.6
1	B	676	TRP	2.6
1	B	678	TRP	2.6
1	A	567	VAL	2.6
1	A	682	PRO	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	299	ARG	2.5
1	A	321	THR	2.5
1	A	470	HIS	2.4
1	A	550	LYS	2.4
1	A	552	ASP	2.4
1	A	514	ARG	2.4
1	A	561	TRP	2.3
1	B	715	VAL	2.3
1	A	371	ARG	2.3
1	A	565	PRO	2.3
1	B	621	THR	2.3
1	A	393	THR	2.3
1	B	567	VAL	2.2
1	B	310	VAL	2.2
1	A	591	THR	2.2
1	B	682	PRO	2.2
1	A	553	TRP	2.2
1	A	302	LYS	2.1
1	B	691	PHE	2.1
1	B	593	ILE	2.1
1	A	681	PRO	2.1
1	B	338	PRO	2.1
1	A	493	LEU	2.1
1	A	370	LYS	2.1
1	B	561	TRP	2.1
1	A	415	CYS	2.0
1	A	674	ALA	2.0
1	B	679	ILE	2.0
1	A	564	LEU	2.0
1	A	617	ASP	2.0
1	A	504	ILE	2.0

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

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

6.3 Carbohydrates [i](#)

There are no 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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	M87	A	800	22/22	0.92	0.23	1.21	29,34,74,83	0
2	HEM	A	750	43/43	0.97	0.18	0.74	28,30,37,45	0
5	ACT	A	860	4/4	0.96	0.13	0.69	60,61,64,66	0
2	HEM	B	750	43/43	0.97	0.16	0.60	23,27,36,38	0
4	M87	B	800	22/22	0.93	0.18	0.59	28,35,60,64	0
3	H4B	B	760	17/17	0.96	0.17	0.43	28,31,35,37	0
3	H4B	A	760	17/17	0.96	0.16	0.21	27,30,35,36	0
6	ZN	A	1717	1/1	1.00	0.09	-0.52	36,36,36,36	0
5	ACT	B	860	4/4	0.96	0.07	-1.14	52,53,54,56	0

6.5 Other polymers

There are no such residues in this entry.