



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:02 PM GMT

PDB ID : 4UPO
Title : Structure of rat neuronal nitric oxide synthase heme domain in complex with N-[3-({[(3S,5S)-5-{{(3-{{(Z)-imino(thiophen-2-yl)methyl}amino}benzyl)oxy}methyl}pyrrolidin-3-yl]oxy}methyl)phenyl]thiophene-2-carboximidamide
Authors : Li, H.; Poulos, T.L.
Deposited on : 2014-06-17
Resolution : 1.95 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

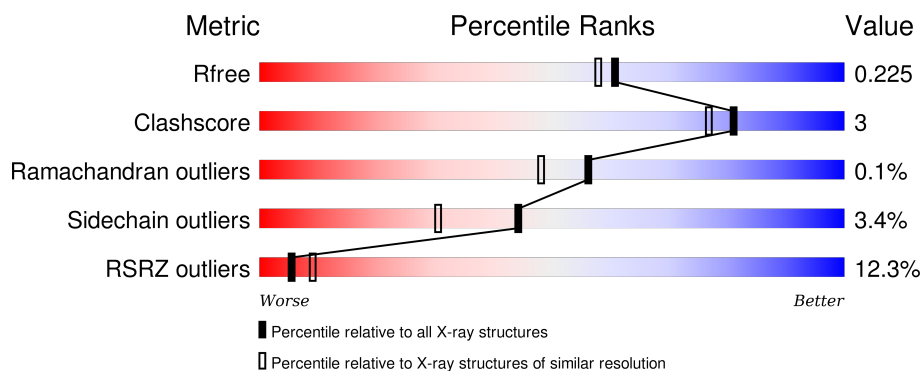
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	1833 (1.96-1.96)
Clashscore	102246	1953 (1.96-1.96)
Ramachandran outliers	100387	1936 (1.96-1.96)
Sidechain outliers	100360	1936 (1.96-1.96)
RSRZ outliers	91569	1835 (1.96-1.96)

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	
1	B	422	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	ACT	A	860	-	-	-	X
5	ACT	B	860	-	-	-	X

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 7173 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	408	Total	C	N	O	S	0	2	1
			3323	2127	567	608	21			
1	B	411	Total	C	N	O	S	0	3	0
			3357	2148	574	614	21			

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



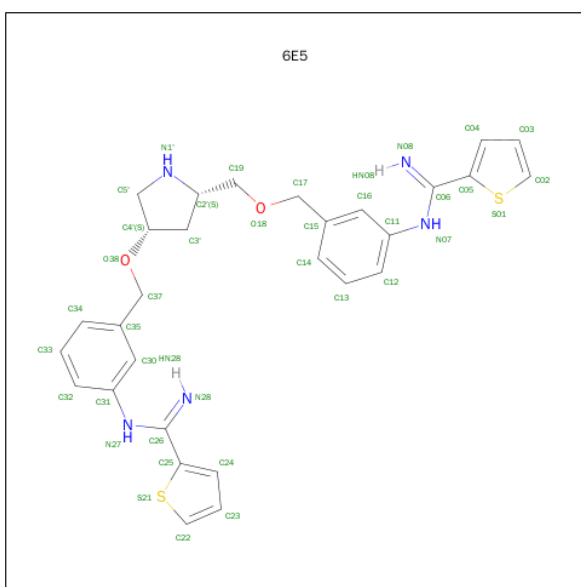
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- 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-[3-({[(3S,5S)-5-{{(3-{{[(Z)-IMINO(THIOPHEN-2-YL)METHYL}AMINO}BENZYL)OXY]METHYL}PYRROLIDIN-3-YL)OXY}METHYL)PHENYL]THIOPHENE-2-CARBOXIMIDAMIDE (three-letter code: 6E5) (formula: C₂₉H₃₁N₅O₂S₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			38	29	5	2	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	1	Total	C	N	O	S	0	0
			38	29	5	2	2		

- 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		

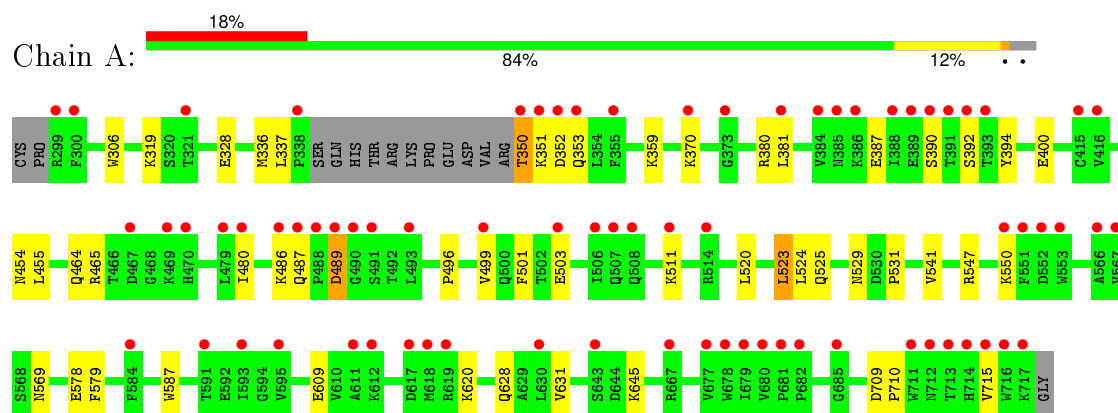
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	122	Total	O	0	0
			122	122		
7	B	166	Total	O	0	0
			166	166		

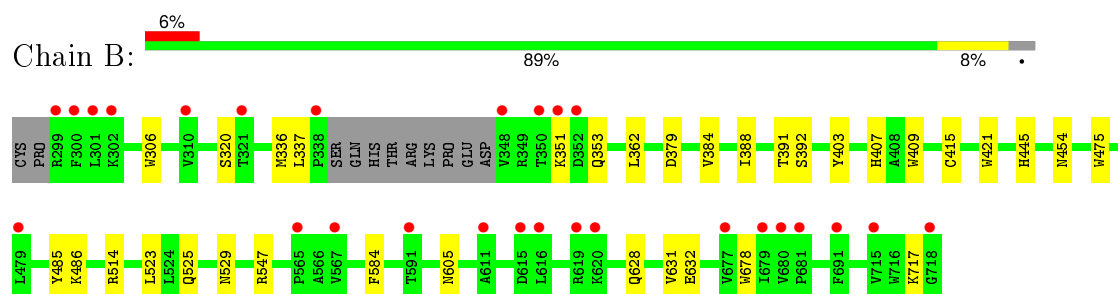
3 Residue-property plots [i](#)

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 ($RSRZ > 2$). 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	52.16Å 110.97Å 164.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.03 – 1.95 38.00 – 1.95	Depositor EDS
% Data completeness (in resolution range)	96.6 (38.03-1.95) 96.5 (38.00-1.95)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.14 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.184 , 0.224 0.184 , 0.225	Depositor DCC
R_{free} test set	3383 reflections (5.21%)	DCC
Wilson B-factor (Å ²)	32.2	Xtriage
Anisotropy	0.781	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 47.2	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	1 of 68356 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7173	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.94% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, ZN, 6E5, H4B, 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	0.56	0/3422	0.68	0/4643
1	B	0.64	0/3459	0.72	1/4689 (0.0%)
All	All	0.60	0/6881	0.70	1/9332 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	379	ASP	CB-CG-OD1	5.33	123.09	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3323	0	3232	26	0
1	B	3357	0	3275	18	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	38	0	29	1	0
4	B	38	0	29	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	4	0	3	0	0
5	B	4	0	3	0	0
6	A	1	0	0	0	0
7	A	122	0	0	2	0
7	B	166	0	0	1	0
All	All	7173	0	6661	44	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 3.

All (44) 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:523:LEU:HD22	1:A:531:PRO:HB2	1.54	0.89
2:A:750:HEM:HHC	2:A:750:HEM:HBB2	1.72	0.71
1:A:609:GLU:HG3	7:A:2091:HOH:O	1.95	0.66
1:A:350:THR:N	1:A:353:GLN:HE21	1.97	0.61
1:A:631:VAL:HG11	1:B:628:GLN:HG3	1.84	0.59
2:B:750:HEM:HBB2	2:B:750:HEM:HHC	1.82	0.59
1:A:350:THR:N	1:A:353:GLN:NE2	2.50	0.59
1:A:359:LYS:HE2	1:A:381:LEU:HD21	1.83	0.59
1:A:380:ARG:HD3	1:A:400:GLU:OE2	2.03	0.58
1:A:499:VAL:O	1:A:503:GLU:HG3	2.05	0.57
1:A:306:TRP:CD2	1:B:336:MET:HE3	2.40	0.56
1:B:605:ASN:ND2	7:B:2143:HOH:O	2.34	0.55
1:A:455:LEU:HD12	1:A:587:TRP:HB3	1.90	0.54
1:A:524:LEU:O	1:A:531:PRO:HA	2.10	0.52
1:A:496:PRO:HA	1:A:499:VAL:HG23	1.94	0.49
1:A:480:ILE:HD13	1:A:541:VAL:HG13	1.94	0.49
1:A:620:LYS:HB2	1:A:620:LYS:NZ	2.28	0.49
1:B:391:THR:O	1:B:392:SER:HB2	2.12	0.48
1:B:403:TYR:CE1	1:B:407:HIS:CE1	3.01	0.48
1:A:387:GLU:OE2	1:A:394:TYR:HA	2.14	0.47
1:B:475:TRP:HB2	1:B:523:LEU:HB3	1.96	0.47
1:A:628:GLN:HG2	1:B:631:VAL:HG11	1.97	0.47
2:B:750:HEM:HHC	2:B:750:HEM:CBB	2.45	0.46
1:B:409:TRP:CE3	1:B:421:TRP:HA	2.52	0.45
1:B:678:TRP:HH2	4:B:800:6E5:H5'A	1.81	0.45
1:A:328:GLU:CD	1:A:328:GLU:H	2.20	0.45
1:B:485:TYR:CE2	1:B:514:ARG:HA	2.51	0.45
1:A:501:PHE:HD2	1:A:520:LEU:HD13	1.81	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:336:MET:HG3	1:B:306:TRP:NE1	2.33	0.43
4:A:800:6E5:H19	7:A:2076:HOH:O	2.18	0.43
2:A:750:HEM:CMC	2:A:750:HEM:HBC2	2.49	0.43
1:B:584:PHE:CD1	2:B:750:HEM:CAC	3.02	0.42
1:A:628:GLN:NE2	1:B:632:GLU:OE2	2.52	0.42
1:B:445:HIS:C	1:B:445:HIS:CD2	2.93	0.42
1:A:465:ARG:HD2	1:A:578:GLU:OE1	2.19	0.42
1:A:455:LEU:HD12	1:A:587:TRP:CB	2.50	0.42
1:A:709:ASP:HA	1:A:710:PRO:HD3	1.91	0.42
1:B:415:CYS:HB2	2:B:750:HEM:ND	2.34	0.41
1:B:362:LEU:HD11	1:B:384:VAL:HG21	2.02	0.41
1:A:351:LYS:HE3	1:A:392:SER:HB3	2.02	0.41
1:B:525:GLN:HG3	1:B:529:ASN:O	2.20	0.41
1:B:388:ILE:O	1:B:392:SER:N	2.49	0.41
1:A:525:GLN:HG3	1:A:529:ASN:O	2.21	0.40
1:A:464:GLN:HB3	1:A:579:PHE:CE1	2.57	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	406/422 (96%)	392 (97%)	13 (3%)	1 (0%)	52	43
1	B	410/422 (97%)	398 (97%)	12 (3%)	0	100	100
All	All	816/844 (97%)	790 (97%)	25 (3%)	1 (0%)	56	48

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	489	ASP

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%)	348 (95%)	17 (5%)	32	16
1	B	369/377 (98%)	361 (98%)	8 (2%)	60	51
All	All	734/754 (97%)	709 (97%)	25 (3%)	44	30

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	319	LYS
1	A	337	LEU
1	A	350	THR
1	A	352	ASP
1	A	370	LYS
1	A	390	SER
1	A	454	ASN
1	A	486	LYS
1	A	487	GLN
1	A	489	ASP
1	A	511	LYS
1	A	523	LEU
1	A	547	ARG
1	A	550	LYS
1	A	569	ASN
1	A	645	LYS
1	A	715	VAL
1	B	320	SER
1	B	337	LEU
1	B	351	LYS
1	B	353	GLN
1	B	454	ASN
1	B	486	LYS
1	B	547	ARG
1	B	717	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	454	ASN
1	A	527	ASN
1	A	569	ASN
1	A	628	GLN
1	A	697	ASN
1	B	364	GLN
1	B	385	ASN
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 [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	2.18	8 (26%)	24,82,82	2.58	12 (50%)
3	H4B	A	760	-	13,18,18	1.37	2 (15%)	11,26,26	2.82	6 (54%)
4	6E5	A	800	-	40,42,42	1.31	6 (15%)	35,56,56	1.76	8 (22%)
5	ACT	A	860	-	1,3,3	1.91	0	0,3,3	0.00	-
2	HEM	B	750	1	30,50,50	2.17	7 (23%)	24,82,82	2.44	10 (41%)
3	H4B	B	760	-	13,18,18	1.01	1 (7%)	11,26,26	2.05	4 (36%)
4	6E5	B	800	-	40,42,42	1.26	5 (12%)	35,56,56	1.86	6 (17%)
5	ACT	B	860	-	1,3,3	1.58	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	6E5	A	800	-	-	1/19/36/36	0/5/5/5
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	6E5	B	800	-	-	0/19/36/36	0/5/5/5
5	ACT	B	860	-	-	0/0/0/0	0/0/0/0

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	750	HEM	C2C-C1C	-6.71	1.39	1.52
2	A	750	HEM	C2D-C3D	-6.44	1.35	1.54
2	B	750	HEM	C2D-C3D	-6.28	1.35	1.54
2	A	750	HEM	C2C-C1C	-6.02	1.41	1.52
2	A	750	HEM	C3B-C4B	-3.24	1.49	1.51
2	A	750	HEM	C3D-C4D	-3.23	1.47	1.51
2	B	750	HEM	C3D-C4D	-2.96	1.47	1.51
4	B	800	6E5	C26-N27	-2.93	1.33	1.39
4	B	800	6E5	C31-N27	-2.76	1.36	1.41
4	A	800	6E5	C31-N27	-2.68	1.36	1.41
2	B	750	HEM	C2B-C1B	-2.61	1.43	1.51
2	B	750	HEM	C3B-C4B	-2.57	1.49	1.51
4	A	800	6E5	C26-N27	-2.54	1.34	1.39
2	A	750	HEM	C2B-C1B	-2.39	1.44	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	800	6E5	C06-N07	-2.32	1.34	1.39
3	B	760	H4B	C4-N3	2.21	1.37	1.33
4	A	800	6E5	C05-S01	2.22	1.75	1.72
4	B	800	6E5	C13-C14	2.23	1.43	1.38
2	A	750	HEM	CHC-C1C	2.27	1.41	1.36
2	B	750	HEM	FE-NC	2.40	2.05	1.95
4	B	800	6E5	C05-C06	2.47	1.49	1.45
2	A	750	HEM	FE-NB	2.57	2.11	1.97
2	B	750	HEM	C1C-NC	2.65	1.39	1.36
3	A	760	H4B	C4-N3	3.02	1.38	1.33
2	A	750	HEM	FE-NC	3.07	2.07	1.95
4	B	800	6E5	C25-C26	3.16	1.50	1.45
3	A	760	H4B	C2-N2	3.17	1.40	1.34
4	A	800	6E5	C05-C06	3.20	1.50	1.45
4	A	800	6E5	C25-C26	3.81	1.50	1.45

All (46) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	800	6E5	C03-C02-S01	-6.26	106.30	113.23
4	A	800	6E5	C23-C22-S21	-5.63	107.00	113.23
4	B	800	6E5	C23-C22-S21	-5.26	107.41	113.23
2	A	750	HEM	C3C-CAC-CBC	-5.18	116.52	124.46
4	A	800	6E5	C03-C02-S01	-4.48	108.28	113.23
3	A	760	H4B	N3-C2-N1	-3.78	119.34	125.53
2	B	750	HEM	CBA-CAA-C2A	-3.51	106.24	112.53
2	A	750	HEM	CBA-CAA-C2A	-3.36	106.50	112.53
2	A	750	HEM	C1D-CHD-C4C	-3.10	120.64	125.82
2	B	750	HEM	C1D-CHD-C4C	-2.96	120.88	125.82
4	A	800	6E5	O18-C19-C2'	-2.83	101.05	109.72
3	B	760	H4B	N3-C2-N1	-2.79	120.97	125.53
2	B	750	HEM	C3C-CAC-CBC	-2.51	120.61	124.46
2	B	750	HEM	C3B-C4B-NB	-2.50	106.85	111.63
4	B	800	6E5	C31-N27-C26	-2.37	122.48	128.32
4	A	800	6E5	C31-N27-C26	-2.30	122.67	128.32
4	B	800	6E5	C12-C11-C16	-2.18	117.08	119.69
4	A	800	6E5	C11-N07-C06	-2.17	122.99	128.32
2	A	750	HEM	CAA-C2A-C3A	-2.11	122.97	129.00
2	A	750	HEM	C3B-C4B-NB	-2.09	107.63	111.63
4	A	800	6E5	C14-C15-C16	2.09	121.64	118.55
4	B	800	6E5	C14-C15-C16	2.13	121.70	118.55
4	A	800	6E5	C25-C26-N27	2.24	121.47	115.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	750	HEM	C2D-C3D-C4D	2.37	105.51	101.50
2	A	750	HEM	C2D-C3D-C4D	2.44	105.63	101.50
4	B	800	6E5	C13-C12-C11	2.51	122.90	119.72
3	B	760	H4B	C2-N1-C8A	2.52	120.19	114.54
2	A	750	HEM	CMD-C2D-C3D	2.61	125.91	114.35
3	A	760	H4B	N2-C2-N3	2.66	121.61	117.20
2	B	750	HEM	CMD-C2D-C3D	2.73	126.41	114.35
2	A	750	HEM	CAA-C2A-C1A	2.92	130.18	127.01
2	A	750	HEM	CMB-C2B-C3B	2.96	123.91	116.53
3	B	760	H4B	C4-N3-C2	2.97	120.06	115.94
4	A	800	6E5	C3'-C4'-C5'	2.99	106.38	103.59
3	A	760	H4B	C4A-C8A-N8	3.62	122.69	118.43
3	B	760	H4B	C4-C4A-C8A	3.84	118.03	114.56
3	A	760	H4B	C2-N1-C8A	3.85	123.19	114.54
2	B	750	HEM	CAD-C3D-C4D	3.89	126.20	112.47
3	A	760	H4B	C4-N3-C2	3.96	121.44	115.94
2	A	750	HEM	CAD-C3D-C4D	4.25	127.45	112.47
2	B	750	HEM	CMC-C2C-C3C	4.25	127.14	116.53
3	A	760	H4B	C4-C4A-C8A	4.28	118.44	114.56
2	B	750	HEM	CMB-C2B-C3B	4.59	127.99	116.53
2	A	750	HEM	CMC-C2C-C3C	4.61	128.03	116.53
2	A	750	HEM	CAD-C3D-C2D	4.76	126.89	113.22
2	B	750	HEM	CAD-C3D-C2D	5.24	128.28	113.22

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	800	6E5	N28-C26-N27-C31

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	750	HEM	2	0
4	A	800	6E5	1	0
2	B	750	HEM	4	0
4	B	800	6E5	1	0

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, 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	408/422 (96%)	0.87	74 (18%) 2 2	26, 53, 95, 131	0
1	B	411/422 (97%)	0.29	27 (6%) 22 31	27, 41, 70, 97	0
All	All	819/844 (97%)	0.58	101 (12%) 5 9	26, 46, 89, 131	0

All (101) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	300	PHE	8.1
1	B	350	THR	6.7
1	A	715	VAL	6.4
1	B	348	VAL	6.3
1	B	718	GLY	6.2
1	A	716	TRP	6.1
1	A	355	PHE	5.9
1	A	488	PRO	5.4
1	A	351	LYS	5.1
1	A	352	ASP	5.1
1	A	350	THR	4.7
1	B	351	LYS	4.7
1	A	388	ILE	4.5
1	A	551	PHE	4.5
1	A	392	SER	4.4
1	A	300	PHE	4.4
1	A	390	SER	4.3
1	B	352	ASP	4.3
1	B	619	ARG	4.2
1	A	490	GLY	4.1
1	A	506	ILE	4.0
1	A	507	GLN	3.9
1	A	489	ASP	3.9
1	A	717	LYS	3.8

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Mol	Chain	Res	Type	RSRZ
1	A	386	LYS	3.8
1	A	619	ARG	3.8
1	A	491	SER	3.7
1	A	713	THR	3.7
1	A	391	THR	3.7
1	A	385	ASN	3.6
1	A	503	GLU	3.5
1	A	470	HIS	3.5
1	A	714	HIS	3.3
1	A	487	GLN	3.3
1	A	511	LYS	3.3
1	A	389	GLU	3.2
1	B	299	ARG	3.1
1	A	682	PRO	3.1
1	A	416	VAL	3.0
1	A	469	LYS	3.0
1	A	381	LEU	3.0
1	A	480	ILE	2.9
1	B	567	VAL	2.9
1	A	678	TRP	2.9
1	B	302	LYS	2.8
1	A	321	THR	2.8
1	A	679	ILE	2.8
1	B	620	LYS	2.8
1	A	681	PRO	2.8
1	A	393	THR	2.8
1	B	310	VAL	2.8
1	B	616	LEU	2.7
1	A	567	VAL	2.7
1	B	715	VAL	2.7
1	A	552	ASP	2.7
1	B	677	VAL	2.7
1	B	591	THR	2.7
1	A	667	ARG	2.7
1	A	514	ARG	2.7
1	A	415	CYS	2.6
1	A	486	LYS	2.6
1	B	691	PHE	2.6
1	A	508	GLN	2.6
1	A	467	ASP	2.6
1	A	591	THR	2.6
1	A	373	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	353	GLN	2.5
1	A	677	VAL	2.5
1	A	680	VAL	2.5
1	A	712	ASN	2.5
1	A	370	LYS	2.5
1	A	595	VAL	2.5
1	A	630	LEU	2.5
1	B	615	ASP	2.4
1	A	584	PHE	2.4
1	A	299	ARG	2.4
1	A	618	MET	2.4
1	B	301	LEU	2.4
1	A	711	TRP	2.3
1	A	499	VAL	2.3
1	A	553	TRP	2.3
1	B	321	THR	2.3
1	A	593	ILE	2.3
1	A	685	GLY	2.3
1	A	493	LEU	2.3
1	A	611	ALA	2.3
1	A	550	LYS	2.2
1	A	338	PRO	2.2
1	B	611	ALA	2.2
1	B	338	PRO	2.2
1	B	680	VAL	2.2
1	B	679	ILE	2.2
1	B	565	PRO	2.2
1	A	617	ASP	2.1
1	A	384	VAL	2.1
1	B	479	LEU	2.1
1	A	643	SER	2.1
1	B	681	PRO	2.1
1	A	479	LEU	2.0
1	A	566	ALA	2.0
1	A	612	LYS	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 [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, 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(\AA^2)	Q<0.9
5	ACT	B	860	4/4	0.87	0.20	7.28	54,59,61,63	0
5	ACT	A	860	4/4	0.93	0.18	5.67	64,66,68,70	0
4	6E5	B	800	38/38	0.92	0.21	1.38	25,75,118,120	0
4	6E5	A	800	38/38	0.87	0.21	1.33	18,56,93,94	0
2	HEM	B	750	43/43	0.98	0.16	0.85	25,29,35,40	0
2	HEM	A	750	43/43	0.97	0.19	0.66	28,31,35,37	0
3	H4B	B	760	17/17	0.96	0.17	0.62	28,30,34,34	0
3	H4B	A	760	17/17	0.97	0.18	0.58	28,31,33,33	0
6	ZN	A	1717	1/1	0.99	0.08	-0.95	38,38,38,38	0

6.5 Other polymers [i](#)

There are no such residues in this entry.