



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 09:35 PM GMT

PDB ID : 4JSL
Title : Structure of bovine endothelial nitric oxide synthase heme domain in complex with 6,6'-(heptane-1,7-diyl)bis(4-methylpyridin-2-amine)
Authors : Li, H.; Poulos, T.L.
Deposited on : 2013-03-22
Resolution : 2.04 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

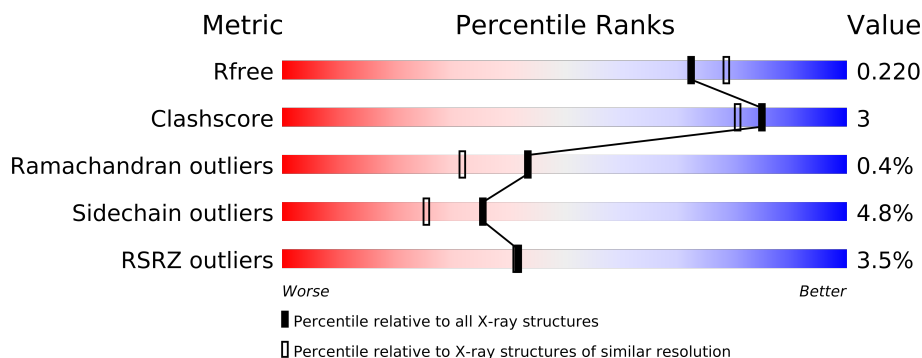
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.04 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	1380 (2.06-2.02)
Clashscore	79885	1577 (2.06-2.02)
Ramachandran outliers	78287	1565 (2.06-2.02)
Sidechain outliers	78261	1565 (2.06-2.02)
RSRZ outliers	66119	1381 (2.06-2.02)

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. 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	443	
1	B	443	

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

Mol	Type	Chain	Res	Geometry	Electron density
5	ACT	B	504	-	X

2 Entry composition i

There are 8 unique types of molecules in this entry. The entry contains 7104 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	As	C	N	O	S	0	2	0
			3234	1	2056	569	592	16			
1	B	403	Total	As	C	N	O	S	0	1	0
			3218	1	2046	567	588	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	100	ARG	CYS	CONFLICT	UNP P29473
B	100	ARG	CYS	CONFLICT	UNP P29473

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



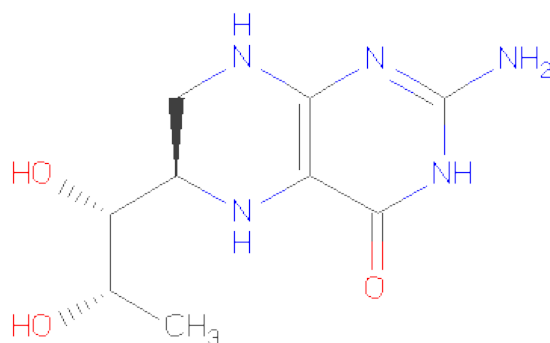
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

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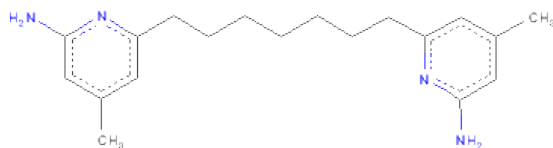
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

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



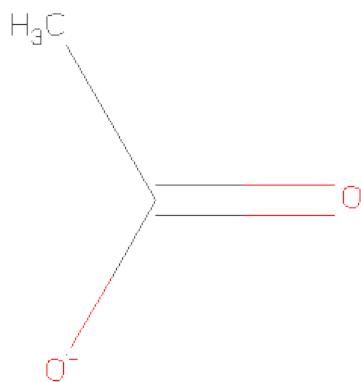
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			17	9	5	3		
3	B	1	Total	C	N	O	0	0
			17	9	5	3		

- Molecule 4 is 6,6'-HEPTANE-1,7-DIYLBIS(4-METHYLPYRIDIN-2-AMINE) (three-letter code: QJ4) (formula: $C_{19}H_{28}N_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	N	0	0
			23	19	4		
4	B	1	Total	C	N	0	0
			23	19	4		

- Molecule 5 is ACETATE ION (three-letter code: ACT) (formula: C₂H₃O₂).



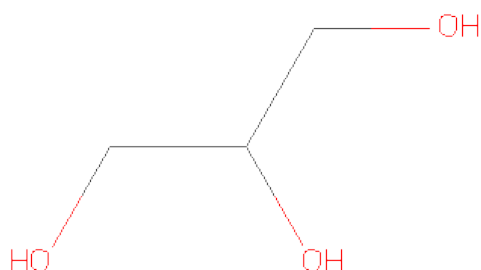
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		

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

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 7 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 8 is water.

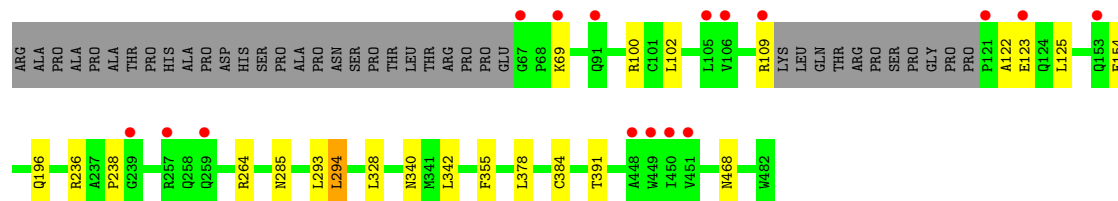
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	228	Total	O	0	0
			228	228		
8	B	229	Total	O	0	0
			229	229		

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.

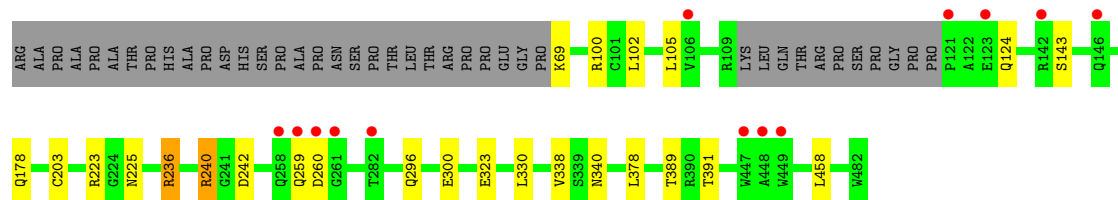
- Molecule 1: Nitric oxide synthase, endothelial

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, α , β , γ	57.86Å 106.58Å 157.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.97 – 2.04 46.97 – 2.04	Depositor EDS
% Data completeness (in resolution range)	99.3 (46.97-2.04) 99.3 (46.97-2.04)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.37 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.169 , 0.208 0.178 , 0.220	Depositor DCC
R_{free} test set	3088 reflections (5.22%)	DCC
Wilson B-factor (Å ²)	35.1	Xtriage
Anisotropy	0.053	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 33.8	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 62232 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7104	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.29% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities 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, CAS, ACT, HEM, QJ4

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	0.69	0/3320	0.70	1/4520 (0.0%)
1	B	0.69	1/3300 (0.0%)	0.69	1/4492 (0.0%)
All	All	0.69	1/6620 (0.0%)	0.69	2/9012 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	203	CYS	CB-SG	-6.61	1.71	1.82

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	240	ARG	NE-CZ-NH2	-5.95	117.32	120.30
1	A	294	LEU	CB-CG-CD1	5.31	120.02	111.00

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	3234	0	0	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3218	0	0	3	0
2	A	43	0	30	6	0
2	B	43	0	30	3	0
3	A	17	0	15	1	0
3	B	17	0	15	0	0
4	A	23	0	28	3	0
4	B	23	0	28	3	0
5	A	8	0	6	0	0
5	B	8	0	6	0	0
6	A	6	0	8	1	0
6	B	6	0	8	1	0
7	A	1	0	0	0	0
8	A	228	0	0	1	0
8	B	229	0	0	1	0
All	All	7104	0	174	19	0

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 3.

All (19) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:384:CAS:SG	1:A:384:CAS:AS	2.58	1.22
3:A:502:H4B:O4	6:A:506:GOL:H32	1.86	0.75
1:B:236:ARG:NH1	1:B:242:ASP:OD1	2.41	0.54
2:A:501:HEM:C1C	4:A:503:QJ4:H28	2.46	0.51
2:A:501:HEM:HHC	2:A:501:HEM:HBB2	1.96	0.48
6:B:506:GOL:H2	8:B:649:HOH:O	2.15	0.47
1:A:384:CAS:CB	1:A:384:CAS:AS	3.22	0.47
2:B:501:HEM:O2D	4:B:503:QJ4:H11	2.16	0.46
1:B:296:GLN:NE2	1:B:300:GLU:O	2.48	0.46
1:A:264:ARG:NE	1:A:285:ASN:O	2.50	0.45
2:A:501:HEM:O2D	4:A:503:QJ4:N21	2.50	0.45
1:A:196:GLN:OE1	8:A:817:HOH:O	2.21	0.45
2:A:501:HEM:HBA2	4:A:503:QJ4:H18	2.00	0.44
2:B:501:HEM:HBC2	2:B:501:HEM:CMC	2.49	0.43
2:A:501:HEM:HHC	2:A:501:HEM:CBB	2.49	0.43
1:A:355:PHE:CD1	2:A:501:HEM:CAC	3.03	0.42
1:B:338:VAL:CG2	4:B:503:QJ4:H18	2.51	0.41
2:B:501:HEM:C1C	4:B:503:QJ4:H28	2.56	0.41
1:A:236:ARG:NH1	1:A:342:LEU:CD1	2.84	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	402/443 (91%)	390 (97%)	10 (2%)	2 (0%)	38	25
1	B	399/443 (90%)	388 (97%)	10 (2%)	1 (0%)	50	39
All	All	801/886 (90%)	778 (97%)	20 (2%)	3 (0%)	43	32

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	122	ALA
1	B	260	ASP
1	A	238	PRO

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	345/375 (92%)	331 (96%)	14 (4%)	41	33
1	B	343/375 (92%)	324 (94%)	19 (6%)	30	20
All	All	688/750 (92%)	655 (95%)	33 (5%)	35	25

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	69	LYS
1	A	100	ARG
1	A	102	LEU
1	A	109	ARG
1	A	123	GLU
1	A	125	LEU

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Mol	Chain	Res	Type
1	A	154	GLU
1	A	293	LEU
1	A	294	LEU
1	A	328	LEU
1	A	340	ASN
1	A	378	LEU
1	A	391	THR
1	A	468	ASN
1	B	69	LYS
1	B	100	ARG
1	B	102	LEU
1	B	105	LEU
1	B	124	GLN
1	B	143	SER
1	B	178	GLN
1	B	223	ARG
1	B	225	ASN
1	B	236	ARG
1	B	240	ARG
1	B	259	GLN
1	B	323	GLU
1	B	330	LEU
1	B	340	ASN
1	B	378	LEU
1	B	389	THR
1	B	391	THR
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 ⓘ

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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
1	CAS	A	384	1	8,8,9	18.33	3 (37%)	7,9,11	5.28	4 (57%)
1	CAS	B	384	1	8,8,9	8.98	3 (37%)	7,9,11	5.71	3 (42%)

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
1	CAS	A	384	1	-	0/2/7/9	0/0/0/0
1	CAS	B	384	1	-	0/2/7/9	0/0/0/0

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	384	CAS	AS-SG	49.49	2.58	2.26
1	B	384	CAS	AS-SG	19.07	2.38	2.26
1	B	384	CAS	O-C	16.39	1.22	1.11
1	A	384	CAS	O-C	15.12	1.21	1.11
1	B	384	CAS	CA-C	2.75	1.53	1.48
1	A	384	CAS	CA-C	2.64	1.53	1.48

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	384	CAS	CE2-AS-CE1	13.85	118.37	96.77
1	A	384	CAS	CE2-AS-CE1	9.44	111.49	96.77
1	A	384	CAS	CE1-AS-SG	7.65	124.97	96.44
1	A	384	CAS	AS-SG-CB	-5.89	92.42	102.24
1	B	384	CAS	AS-SG-CB	-4.49	94.76	102.24
1	B	384	CAS	CE1-AS-SG	3.26	108.60	96.44
1	A	384	CAS	C-CA-N	-3.17	110.66	113.83

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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	501	1	49,50,50	2.71	15 (30%)	46,82,82	1.99	10 (21%)
3	H4B	A	502	-	18,18,18	1.02	1 (5%)	24,26,26	1.53	2 (8%)
4	QJ4	A	503	-	24,24,24	0.57	0	31,31,31	1.84	8 (25%)
5	ACT	A	504	-	1,3,3	1.27	0	0,3,3	0.00	-
5	ACT	A	505	-	1,3,3	2.42	1 (100%)	0,3,3	0.00	-
6	GOL	A	506	-	5,5,5	0.34	0	5,5,5	0.72	0
2	HEM	B	501	1	49,50,50	2.51	17 (34%)	46,82,82	2.08	11 (23%)
3	H4B	B	502	-	18,18,18	0.91	1 (5%)	24,26,26	1.62	9 (37%)
4	QJ4	B	503	-	24,24,24	0.76	1 (4%)	31,31,31	2.50	10 (32%)
5	ACT	B	504	-	1,3,3	1.46	0	0,3,3	0.00	-
5	ACT	B	505	-	1,3,3	2.00	0	0,3,3	0.00	-
6	GOL	B	506	-	5,5,5	0.58	0	5,5,5	0.56	0

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	501	1	-	0/14/114/114	0/0/8/8
3	H4B	A	502	-	-	0/8/17/17	0/0/2/2
4	QJ4	A	503	-	-	0/10/10/10	0/2/2/2
5	ACT	A	504	-	-	0/0/0/0	0/0/0/0
5	ACT	A	505	-	-	0/0/0/0	0/0/0/0
6	GOL	A	506	-	-	0/4/4/4	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	B	501	1	-	0/14/114/114	0/0/8/8
3	H4B	B	502	-	-	0/8/17/17	0/0/2/2
4	QJ4	B	503	-	-	0/10/10/10	0/2/2/2
5	ACT	B	504	-	-	0/0/0/0	0/0/0/0
5	ACT	B	505	-	-	0/0/0/0	0/0/0/0
6	GOL	B	506	-	-	0/4/4/4	0/0/0/0

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	HEM	C2D-C1D	9.71	1.47	1.44
2	B	501	HEM	C2B-C1B	6.79	1.46	1.44
2	A	501	HEM	C2B-C1B	6.76	1.46	1.44
2	B	501	HEM	C3D-C4D	5.65	1.46	1.44
2	B	501	HEM	C3B-C2B	-5.54	1.34	1.43
2	A	501	HEM	C3B-C2B	-5.54	1.34	1.43
2	B	501	HEM	C3C-C2C	-5.34	1.34	1.43
2	B	501	HEM	C3B-CAB	4.82	1.55	1.40
2	B	501	HEM	C3D-C2D	4.77	1.52	1.43
2	A	501	HEM	C3C-CAC	4.76	1.55	1.40
2	A	501	HEM	C3C-C2C	-4.62	1.35	1.43
2	A	501	HEM	C3D-C2D	4.60	1.51	1.43
2	A	501	HEM	C3B-CAB	4.43	1.54	1.40
2	B	501	HEM	C3C-CAC	4.43	1.54	1.40
2	B	501	HEM	C4A-C3A	4.05	1.45	1.40
2	A	501	HEM	C4A-C3A	3.95	1.45	1.40
2	A	501	HEM	FE-NA	3.53	2.07	1.92
2	B	501	HEM	FE-NA	3.34	2.06	1.92
2	B	501	HEM	FE-NB	2.96	2.08	1.97
2	A	501	HEM	CMC-C2C	2.79	1.56	1.47
2	B	501	HEM	CMD-C2D	2.66	1.55	1.47
2	A	501	HEM	CAA-C2A	2.66	1.56	1.52
2	B	501	HEM	FE-ND	2.58	2.07	1.97
5	A	505	ACT	CH3-C	2.42	1.52	1.48
2	B	501	HEM	CAA-C2A	2.25	1.56	1.52
2	A	501	HEM	CMD-C2D	2.24	1.54	1.47
2	A	501	HEM	FE-NB	2.23	2.06	1.97
2	B	501	HEM	FE-NC	2.21	2.06	1.97
2	A	501	HEM	CMB-C2B	2.20	1.54	1.47
3	B	502	H4B	C2-N2	2.12	1.35	1.32
2	B	501	HEM	CHB-C1B	2.08	1.38	1.35
2	B	501	HEM	CMB-C2B	2.07	1.53	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502	H4B	C7-C6	2.05	1.54	1.52
2	B	501	HEM	C2D-C1D	2.03	1.45	1.44
4	B	503	QJ4	C26-N21	2.01	1.38	1.34
2	A	501	HEM	C4A-NA	2.01	1.40	1.36

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	HEM	C3B-C4B-NB	-8.38	108.00	114.00
2	B	501	HEM	C3B-C4B-NB	-7.85	108.39	114.00
4	B	503	QJ4	C02-N01-C06	6.43	122.75	118.23
4	B	503	QJ4	C14-C26-N21	6.32	124.27	115.69
4	A	503	QJ4	C02-N01-C06	5.80	122.30	118.23
4	B	503	QJ4	C22-N21-C26	5.79	122.30	118.23
2	B	501	HEM	C4D-ND-C1D	5.13	110.41	105.16
3	A	502	H4B	C4-C4A-C8A	4.40	118.63	114.56
2	B	501	HEM	C3A-C4A-NA	-4.20	106.24	109.41
2	A	501	HEM	C4D-ND-C1D	4.09	109.35	105.16
2	A	501	HEM	C4A-CHB-C1B	-4.01	122.19	127.47
4	B	503	QJ4	C04-C05-C06	-3.96	117.65	120.29
4	B	503	QJ4	C14-C26-C25	-3.59	116.19	121.17
2	A	501	HEM	CMA-C3A-C4A	-3.54	123.18	128.62
2	B	501	HEM	C4A-C3A-C2A	3.51	109.44	107.00
2	A	501	HEM	CHC-C4B-NB	3.39	127.40	124.58
2	B	501	HEM	CBA-CAA-C2A	-3.37	106.76	112.69
2	B	501	HEM	CBD-CAD-C3D	-3.23	107.32	114.37
4	A	503	QJ4	C04-C05-C06	-3.21	118.15	120.29
4	A	503	QJ4	C08-C06-C05	3.13	125.52	121.17
4	B	503	QJ4	C25-C26-N21	-3.10	119.54	122.99
3	B	502	H4B	C4A-C4-N3	2.92	121.54	114.06
3	B	502	H4B	C4-C4A-C8A	2.89	117.23	114.56
4	A	503	QJ4	C14-C26-N21	2.84	119.55	115.69
4	A	503	QJ4	C22-N21-C26	2.79	120.19	118.23
3	A	502	H4B	C6-C7-N8	-2.75	108.11	111.66
2	A	501	HEM	CBA-CAA-C2A	-2.74	107.86	112.69
3	B	502	H4B	C4-C4A-N5	2.70	123.03	119.10
2	B	501	HEM	C1B-NB-C4B	2.69	107.92	105.16
3	B	502	H4B	C4A-N5-C6	-2.65	113.95	121.16
2	B	501	HEM	C4C-NC-C1C	2.59	108.22	105.53
2	A	501	HEM	CMA-C3A-C2A	2.52	129.69	124.94
4	B	503	QJ4	N22-C22-N21	2.48	121.35	116.59
4	B	503	QJ4	C05-C06-N01	-2.46	120.25	122.99

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	502	H4B	C6-C7-N8	-2.42	108.53	111.66
2	B	501	HEM	C4A-CHB-C1B	-2.36	124.36	127.47
3	B	502	H4B	N8-C8A-N1	2.35	119.27	115.82
4	A	503	QJ4	C12-C13-C14	-2.33	103.86	113.96
2	B	501	HEM	C2D-C1D-ND	-2.32	110.19	112.93
4	A	503	QJ4	C05-C06-N01	-2.28	120.45	122.99
2	A	501	HEM	CAD-C3D-C4D	2.27	128.61	124.53
3	B	502	H4B	C2-N1-C8A	2.26	120.82	117.61
2	A	501	HEM	C2D-C1D-ND	-2.25	110.28	112.93
3	B	502	H4B	N2-C2-N3	2.24	120.32	117.86
4	B	503	QJ4	C05-C04-C03	2.23	121.09	118.10
4	B	503	QJ4	C09-C08-C06	2.14	120.73	113.16
3	B	502	H4B	C7-C6-C9	-2.12	110.38	113.66
2	B	501	HEM	CAD-C3D-C4D	2.11	128.31	124.53
2	A	501	HEM	CBD-CAD-C3D	-2.07	109.85	114.37
4	A	503	QJ4	C10-C09-C08	-2.01	105.23	113.96

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, 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	405/443 (91%)	0.01	16 (3%) 36 36	24, 36, 58, 77	0
1	B	403/443 (90%)	0.02	13 (3%) 45 45	23, 38, 62, 84	0
All	All	808/886 (91%)	0.02	29 (3%) 42 40	23, 37, 60, 84	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	259	GLN	6.5
1	A	67	GLY	5.0
1	A	259	GLN	4.4
1	B	121	PRO	3.7
1	A	239	GLY	3.6
1	B	123	GLU	3.2
1	B	448	ALA	3.2
1	A	109	ARG	3.1
1	A	123	GLU	3.0
1	B	260	ASP	3.0
1	A	69	LYS	3.0
1	A	450	ILE	2.7
1	B	449	TRP	2.7
1	A	257	ARG	2.6
1	A	448	ALA	2.6
1	B	146	GLN	2.5
1	B	447	TRP	2.5
1	A	449	TRP	2.4
1	A	91	GLN	2.4
1	B	261	GLY	2.4
1	A	451	VAL	2.4
1	B	106	VAL	2.3
1	B	258	GLN	2.3
1	A	105	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	282	THR	2.2
1	A	121	PRO	2.1
1	A	106	VAL	2.1
1	A	153	GLN	2.1
1	B	142	ARG	2.0

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

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	CAS	A	384	9/10	0.11	0.76	37,37,66,66	0
1	CAS	B	384	9/10	0.10	-0.53	46,47,63,65	0

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, 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
5	ACT	B	504	4/4	0.20	5.29	44,45,46,46	0
6	GOL	A	506	6/6	0.24	1.70	55,57,57,60	0
6	GOL	B	506	6/6	0.23	1.22	51,52,54,54	0
2	HEM	A	501	43/43	0.19	0.89	24,27,36,42	0
4	QJ4	B	503	23/23	0.19	0.85	28,38,49,49	0
2	HEM	B	501	43/43	0.16	0.65	25,27,35,38	0
3	H4B	B	502	17/17	0.20	0.28	27,29,31,33	0
3	H4B	A	502	17/17	0.20	0.27	25,29,31,32	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	QJ4	A	503	23/23	0.16	0.26	24,33,45,46	0
5	ACT	B	505	4/4	0.09	0.08	32,34,34,35	0
5	ACT	A	505	4/4	0.09	-0.03	29,29,29,31	0
5	ACT	A	504	4/4	0.12	-0.09	43,44,45,46	0
7	ZN	A	507	1/1	0.10	-1.71	30,30,30,30	0

6.5 Other polymers ⓘ

There are no such residues in this entry.