



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 28, 2017 – 04:40 PM EDT

PDB ID : 5VUT
Title : Structure of rat neuronal nitric oxide synthase heme domain in complex with
4-(2-(((2-Amino-4-methylquinolin-7-yl)methyl)amino)ethyl)benzonitrile
Authors : Li, H.; Poulos, T.L.
Deposited on : unknown
Resolution : 2.00 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029824
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : rb-20029824

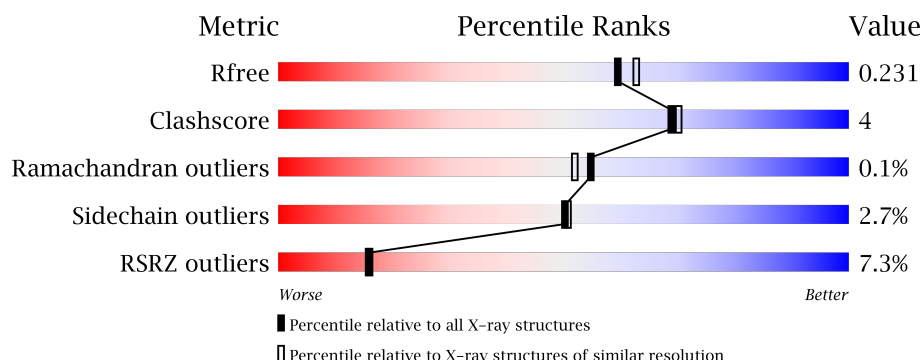
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	100719	6609 (2.00-2.00)
Clashscore	112137	7775 (2.00-2.00)
Ramachandran outliers	110173	7679 (2.00-2.00)
Sidechain outliers	110143	7678 (2.00-2.00)
RSRZ outliers	101464	6696 (2.00-2.00)

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>11%</div> <div> <div></div> <div>82%</div> <div>14%</div> <div></div> </div> <div></div> </div>
1	B	422	<div> <div>3%</div> <div> <div></div> <div>86%</div> <div>11%</div> <div></div> </div> <div></div> </div>

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
4	9OG	A	803	-	-	-	X
4	9OG	B	803	-	-	-	X
5	ACT	A	804	-	-	-	X
5	ACT	B	804	-	-	-	X

2 Entry composition [i](#)

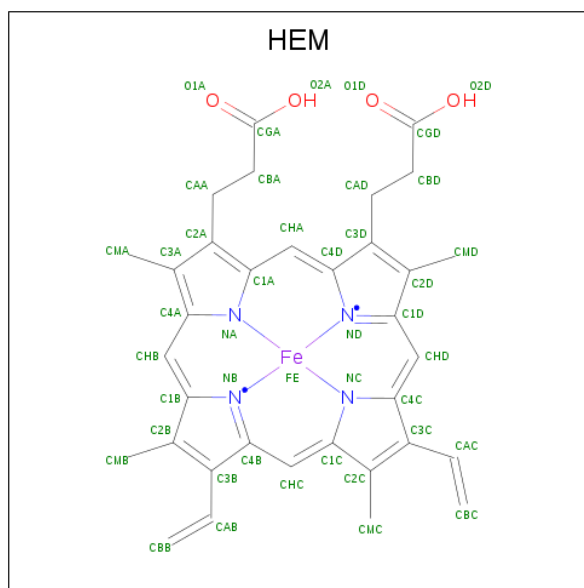
There are 7 unique types of molecules in this entry. The entry contains 7224 atoms, of which 0 are hydrogens and 0 are deuteriums.

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	3	0
			3322	2127	566	607	22			
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: $C_{34}H_{32}FeN_4O_4$).



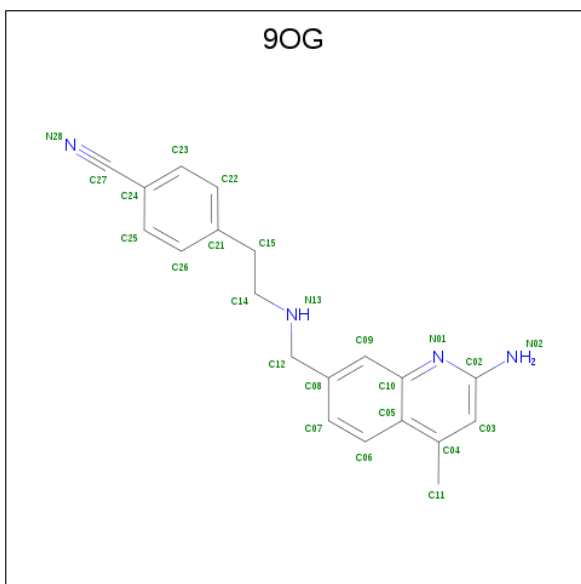
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- 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 4-(2-{{(2-amino-4-methylquinolin-7-yl)methyl}amino}ethyl)benzonitrile (three-letter code: 9OG) (formula: C₂₀H₂₀N₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N		0	0
			24	20	4			

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	N	0	0
			24	20	4		

- 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		

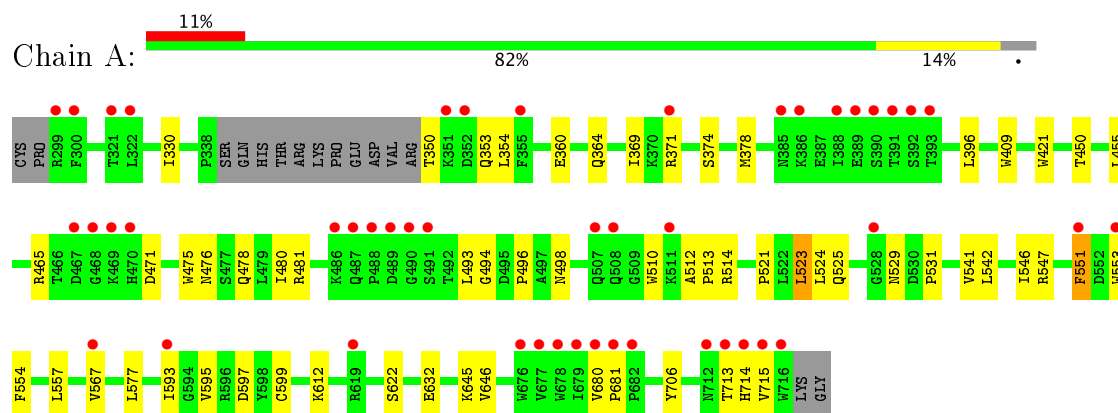
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	162	Total	O	0	0
			162	162		
7	B	206	Total	O	0	0
			206	206		

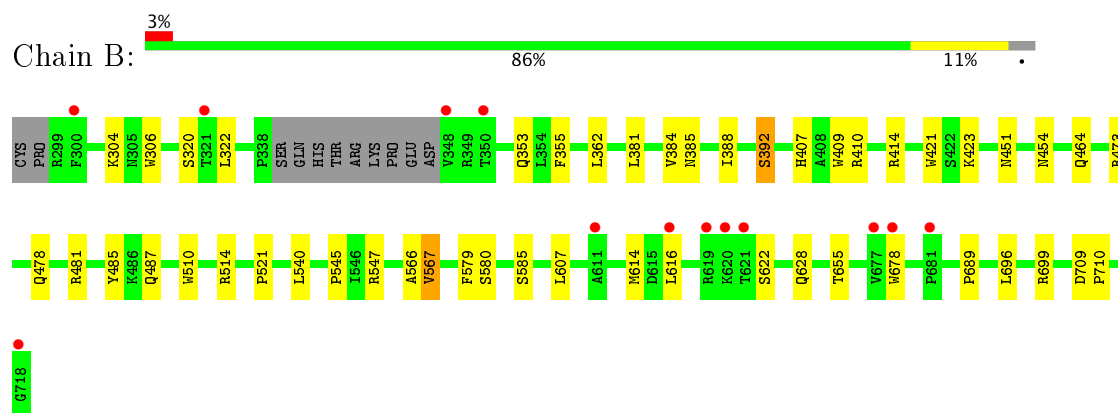
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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	52.17Å 111.04Å 165.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.08 – 2.00 39.08 – 2.00	Depositor EDS
% Data completeness (in resolution range)	97.8 (39.08-2.00) 98.0 (39.08-2.00)	Depositor EDS
R_{merge}	0.26	Depositor
R_{sym}	0.26	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.38 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.8.2_1309	Depositor
R, R_{free}	0.200 , 0.238 0.192 , 0.231	Depositor DCC
R_{free} test set	3182 reflections (4.94%)	DCC
Wilson B-factor (Å ²)	31.9	Xtriage
Anisotropy	0.560	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 54.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7224	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.19% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: 9OG, HEM, ZN, 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	0.39	0/3421	0.53	0/4641
1	B	0.40	0/3459	0.52	1/4689 (0.0%)
All	All	0.40	0/6880	0.53	1/9330 (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	567	VAL	CB-CA-C	-5.55	100.86	111.40

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	3322	0	3232	32	0
1	B	3357	0	3275	24	0
2	A	43	0	30	3	0
2	B	43	0	30	2	0
3	A	17	0	15	0	0
3	B	17	0	15	0	0
4	A	24	0	0	0	0
4	B	24	0	0	0	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	162	0	0	3	0
7	B	206	0	0	0	0
All	All	7224	0	6603	58	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (58) 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	1.61	0.82
2:A:801:HEM:HHC	2:A:801:HEM:HBB2	1.66	0.77
2:A:801:HEM:HMC2	2:A:801:HEM:HBC2	1.70	0.72
2:B:801:HEM:HBB2	2:B:801:HEM:HHC	1.73	0.71
2:B:801:HEM:HBC2	2:B:801:HEM:HMC2	1.77	0.67
1:A:350:THR:N	1:A:353:GLN:OE1	2.34	0.60
1:B:478:GLN:HB2	1:B:481:ARG:HG3	1.82	0.60
1:A:480:ILE:HD13	1:A:541:VAL:HG13	1.83	0.59
1:A:475:TRP:HB2	1:A:523:LEU:HB3	1.86	0.57
1:A:330:ILE:HD11	1:B:696:LEU:HB3	1.86	0.56
1:A:632:GLU:OE2	1:B:628:GLN:NE2	2.41	0.53
1:A:475:TRP:CZ2	1:A:531:PRO:HG3	2.44	0.52
1:B:322:LEU:HD13	1:B:699:ARG:HH21	1.74	0.52
1:B:510:TRP:CE2	1:B:521:PRO:HD3	2.46	0.50
1:A:481:ARG:NH1	1:A:498:ASN:OD1	2.44	0.49
1:B:362:LEU:HD11	1:B:384:VAL:HG21	1.95	0.49
1:A:706:TYR:OH	2:A:801:HEM:O1D	2.23	0.49
1:B:388:ILE:O	1:B:392:SER:N	2.38	0.48
1:B:409:TRP:CE3	1:B:421:TRP:HA	2.48	0.48
1:A:465:ARG:NH2	1:A:471:ASP:OD2	2.38	0.47
1:B:451:ASN:HB3	1:B:454:ASN:O	2.15	0.47
1:A:369:ILE:HG13	1:A:371:ARG:HG3	1.96	0.46
1:A:360:GLU:O	1:A:364:GLN:HG2	2.16	0.46
1:B:414:ARG:HD3	1:B:678:TRP:CD2	2.51	0.45
1:A:714:HIS:ND1	7:A:902:HOH:O	2.32	0.45
1:A:450:THR:HA	1:A:455:LEU:HD22	1.98	0.44
1:A:374:SER:O	1:A:378:MET:HG2	2.18	0.44
1:B:487:GLN:OE1	1:B:514:ARG:NH2	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:494:GLY:O	1:A:496:PRO:HD3	2.18	0.43
1:B:510:TRP:CG	1:B:521:PRO:HG3	2.53	0.43
1:B:614:MET:HE2	1:B:616:LEU:HD11	1.99	0.43
1:A:525:GLN:HG3	1:A:529:ASN:O	2.18	0.43
1:A:409:TRP:CE3	1:A:421:TRP:HA	2.53	0.43
1:A:478:GLN:HB2	1:A:481:ARG:HG3	2.00	0.43
1:A:553:TRP:HZ3	1:A:557:LEU:HD11	1.83	0.43
1:B:355:PHE:CE1	1:B:385:ASN:HB2	2.54	0.42
1:B:381:LEU:HA	1:B:381:LEU:HD23	1.87	0.42
1:B:566:ALA:HB2	1:B:585:SER:HB3	2.00	0.42
1:A:593:ILE:HA	1:A:597:ASP:HB2	2.01	0.42
1:A:551:PHE:HD1	1:A:553:TRP:CZ2	2.37	0.42
1:B:464:GLN:HB3	1:B:579:PHE:CE2	2.55	0.42
1:A:512:ALA:HA	1:A:513:PRO:HD3	1.94	0.42
1:A:396:LEU:HG	1:A:577:LEU:HD12	2.01	0.42
1:A:546:ILE:HG22	1:A:554:PHE:HE2	1.85	0.41
1:B:304:LYS:HD3	1:B:306:TRP:CE2	2.55	0.41
1:B:407:HIS:CE1	1:B:410:ARG:HH11	2.38	0.41
1:B:709:ASP:HA	1:B:710:PRO:HD3	1.95	0.41
1:A:510:TRP:CD1	1:A:521:PRO:HG3	2.56	0.41
1:B:473:ARG:HD3	1:B:580:SER:HB2	2.03	0.41
1:B:485:TYR:CZ	1:B:514:ARG:HA	2.56	0.41
1:A:542:LEU:HD21	1:A:646:VAL:HG22	2.03	0.41
1:A:554:PHE:HB3	7:A:925:HOH:O	2.21	0.41
1:A:680:VAL:HA	1:A:681:PRO:HD3	1.92	0.41
1:A:595:VAL:O	1:A:599:CYS:HB2	2.21	0.40
1:A:524:LEU:O	1:A:531:PRO:HA	2.21	0.40
1:B:655:THR:OG1	1:B:689:PRO:HD2	2.22	0.40
1:B:607:LEU:HA	1:B:607:LEU:HD23	1.97	0.40
1:A:493:LEU:HA	7:A:1019:HOH:O	2.20	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	405/422 (96%)	385 (95%)	19 (5%)	1 (0%)	51	48
1	B	410/422 (97%)	398 (97%)	12 (3%)	0	100	100
All	All	815/844 (97%)	783 (96%)	31 (4%)	1 (0%)	55	52

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	514	ARG

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%)	354 (97%)	11 (3%)	46	46
1	B	369/377 (98%)	360 (98%)	9 (2%)	54	56
All	All	734/754 (97%)	714 (97%)	20 (3%)	50	51

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	354	LEU
1	A	476	ASN
1	A	523	LEU
1	A	547	ARG
1	A	551	PHE
1	A	567	VAL
1	A	612	LYS
1	A	622	SER
1	A	645	LYS
1	A	713	THR
1	A	715	VAL
1	B	320	SER

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Mol	Chain	Res	Type
1	B	353	GLN
1	B	392	SER
1	B	423	LYS
1	B	540	LEU
1	B	545	PRO
1	B	547	ARG
1	B	567	VAL
1	B	622	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	353	GLN
1	B	407	HIS
1	B	535	GLN

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	801	1	28,50,50	2.22	7 (25%)	17,82,82	2.61	5 (29%)
3	H4B	A	802	-	14,18,18	0.87	0	12,26,26	2.41	4 (33%)
4	9OG	A	803	-	26,26,26	1.56	2 (7%)	32,35,35	1.36	3 (9%)
5	ACT	A	804	-	1,3,3	1.02	0	0,3,3	0.00	-
2	HEM	B	801	1	28,50,50	2.18	8 (28%)	17,82,82	2.01	4 (23%)
3	H4B	B	802	-	14,18,18	0.84	0	12,26,26	2.25	4 (33%)
4	9OG	B	803	-	26,26,26	1.58	2 (7%)	32,35,35	1.16	3 (9%)
5	ACT	B	804	-	1,3,3	0.92	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	801	1	-	0/6/54/54	0/0/8/8
3	H4B	A	802	-	-	0/8/17/17	0/2/2/2
4	9OG	A	803	-	-	0/9/9/9	0/3/3/3
5	ACT	A	804	-	-	0/0/0/0	0/0/0/0
2	HEM	B	801	1	-	0/6/54/54	0/0/8/8
3	H4B	B	802	-	-	0/8/17/17	0/2/2/2
4	9OG	B	803	-	-	0/9/9/9	0/3/3/3
5	ACT	B	804	-	-	0/0/0/0	0/0/0/0

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	803	9OG	C24-C27	-6.42	1.29	1.44
4	B	803	9OG	C24-C27	-6.28	1.29	1.44
2	A	801	HEM	C3B-C2B	-4.39	1.34	1.40
2	B	801	HEM	C3B-C2B	-4.05	1.35	1.40
2	A	801	HEM	C3C-C2C	-3.81	1.35	1.40
2	B	801	HEM	C3C-C2C	-3.48	1.35	1.40
4	B	803	9OG	C05-C10	-2.13	1.39	1.42
2	A	801	HEM	C1C-NC	2.17	1.39	1.36
2	B	801	HEM	C1B-NB	2.20	1.39	1.36
4	A	803	9OG	C02-N01	2.31	1.35	1.33
2	B	801	HEM	C4D-ND	2.41	1.39	1.36
2	B	801	HEM	C1C-NC	2.45	1.39	1.36
2	A	801	HEM	C4D-ND	2.87	1.40	1.36
2	B	801	HEM	C3B-CAB	3.59	1.55	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	HEM	C3C-CAC	3.77	1.55	1.47
2	A	801	HEM	C3B-CAB	3.95	1.55	1.47
2	B	801	HEM	C3C-CAC	4.02	1.55	1.47
2	A	801	HEM	C3D-C2D	5.26	1.53	1.37
2	B	801	HEM	C3D-C2D	5.38	1.53	1.37

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	HEM	CBA-CAA-C2A	-6.64	99.79	112.48
2	A	801	HEM	CBD-CAD-C3D	-5.35	102.27	112.47
2	B	801	HEM	C1D-C2D-C3D	-4.47	103.89	107.00
2	B	801	HEM	CBA-CAA-C2A	-4.06	104.71	112.48
2	A	801	HEM	C1D-C2D-C3D	-3.58	104.50	107.00
4	A	803	9OG	C05-C10-N01	-3.35	119.38	122.84
3	A	802	H4B	N3-C2-N1	-3.32	120.07	125.45
2	A	801	HEM	CAD-CBD-CGD	-3.31	107.01	112.66
3	B	802	H4B	N3-C2-N1	-2.97	120.63	125.45
2	B	801	HEM	CAD-CBD-CGD	-2.73	107.99	112.66
2	B	801	HEM	CBD-CAD-C3D	-2.72	107.28	112.47
4	B	803	9OG	C05-C10-N01	-2.63	120.12	122.84
4	B	803	9OG	C12-C08-C09	-2.39	118.00	121.69
2	A	801	HEM	C4C-C3C-C2C	2.09	108.36	106.90
4	A	803	9OG	N02-C02-N01	2.20	119.42	118.06
3	B	802	H4B	C2-N1-C8A	2.52	120.19	114.51
3	A	802	H4B	C2-N1-C8A	2.95	121.17	114.51
3	B	802	H4B	C4-N3-C2	3.35	120.87	116.06
3	A	802	H4B	C4-N3-C2	3.48	121.06	116.06
4	B	803	9OG	C04-C05-C10	3.49	119.96	117.99
4	A	803	9OG	C04-C05-C10	3.86	120.17	117.99
3	B	802	H4B	C4-C4A-C8A	5.00	119.09	114.56
3	A	802	H4B	C4-C4A-C8A	5.11	119.19	114.56

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	HEM	3	0
2	B	801	HEM	2	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	407/422 (96%)	0.51	47 (11%) 5 5	23, 50, 92, 126	0
1	B	411/422 (97%)	-0.01	13 (3%) 48 48	23, 39, 71, 95	0
All	All	818/844 (96%)	0.25	60 (7%) 16 16	23, 43, 86, 126	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	715	VAL	6.6
1	B	300	PHE	5.6
1	A	488	PRO	5.4
1	A	716	TRP	5.3
1	A	388	ILE	5.0
1	A	355	PHE	4.4
1	A	713	THR	4.3
1	A	712	ASN	4.1
1	A	390	SER	3.9
1	A	352	ASP	3.9
1	A	490	GLY	3.8
1	A	486	LYS	3.6
1	B	620	LYS	3.6
1	A	507	GLN	3.6
1	A	551	PHE	3.6
1	B	611	ALA	3.6
1	A	351	LYS	3.4
1	B	619	ARG	3.3
1	A	389	GLU	3.3
1	A	619	ARG	3.2
1	B	616	LEU	3.2
1	A	391	THR	3.1
1	A	470	HIS	3.0
1	B	348	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	489	ASP	3.0
1	A	469	LYS	3.0
1	B	677	VAL	3.0
1	A	392	SER	2.9
1	A	681	PRO	2.9
1	A	299	ARG	2.9
1	A	385	ASN	2.9
1	A	528	GLY	2.8
1	A	678	TRP	2.8
1	A	679	ILE	2.8
1	A	393	THR	2.7
1	A	487	GLN	2.7
1	A	677	VAL	2.6
1	A	491	SER	2.6
1	A	508	GLN	2.6
1	A	386	LYS	2.6
1	B	718	GLY	2.5
1	A	300	PHE	2.5
1	A	371	ARG	2.5
1	B	681	PRO	2.5
1	A	714	HIS	2.4
1	A	468	GLY	2.4
1	A	680	VAL	2.4
1	A	682	PRO	2.4
1	B	678	TRP	2.3
1	A	322	LEU	2.2
1	A	676	TRP	2.2
1	B	350	THR	2.2
1	A	593	ILE	2.1
1	A	553	TRP	2.1
1	A	511	LYS	2.1
1	A	321	THR	2.1
1	B	621	THR	2.1
1	B	321	THR	2.1
1	A	467	ASP	2.1
1	A	567	VAL	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	804	4/4	0.95	0.17	3.70	46,46,49,52	0
5	ACT	A	804	4/4	0.94	0.14	2.31	62,63,64,64	0
4	9OG	A	803	24/24	0.90	0.28	2.16	24,47,81,86	0
4	9OG	B	803	24/24	0.92	0.24	2.12	22,45,69,70	0
2	HEM	A	801	43/43	0.97	0.20	1.21	25,33,48,51	0
2	HEM	B	801	43/43	0.98	0.17	0.93	19,31,42,49	0
3	H4B	B	802	17/17	0.96	0.18	0.24	24,27,32,34	0
3	H4B	A	802	17/17	0.96	0.16	0.16	22,31,36,41	0
6	ZN	A	805	1/1	1.00	0.10	-0.35	36,36,36,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.