



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:31 PM GMT

PDB ID : 4IMT
Title : Structure of rat neuronal nitric oxide synthase in complex with 6,6'-((4-(3-aminopropyl)-1,3-phenylene)bis(ethane-2,1-diyl))bis(4-methylpyridin-2-amine)
Authors : Li, H.; Poulos, T.L.
Deposited on : 2013-01-03
Resolution : 2.20 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

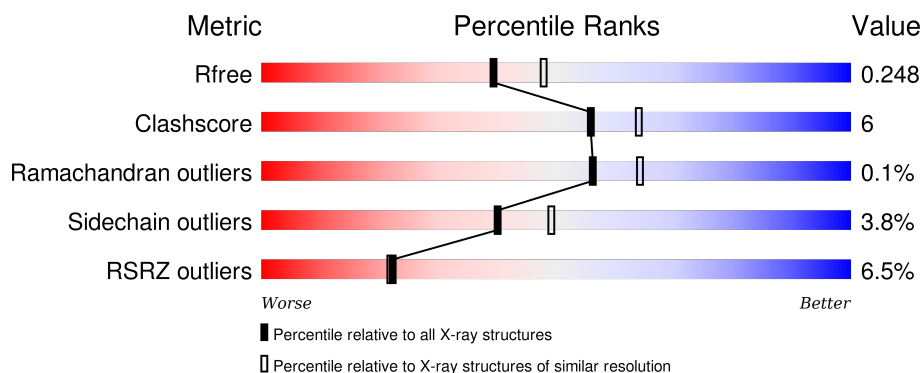
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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>9%</div> <div>80%</div> <div>14%</div> <div>• •</div> </div>
1	B	422	<div> <div>4%</div> <div>85%</div> <div>11%</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	1EW	A	803	-	-	-	X

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 7024 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	1	0
			3319	2125	566	607	21			
1	B	411	Total	C	N	O	S	0	2	0
			3354	2146	574	613	21			

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



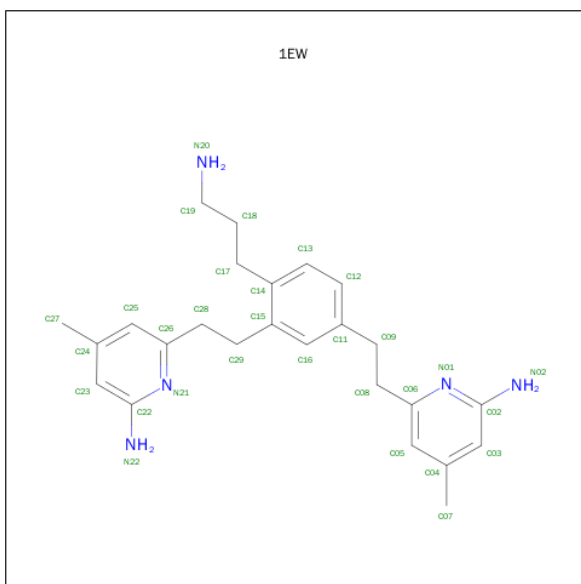
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- 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 6,6'-{[4-(3-AMINOPROPYL)BENZENE-1,3-DIYL]DIETHANE-2,1-DIYL}BIS(4-METHYLPYRIDIN-2-AMINE) (three-letter code: 1EW) (formula: C₂₅H₃₃N₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	N	0	0
			30	25	5		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	N	0	0
			30	25	5		

- 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		

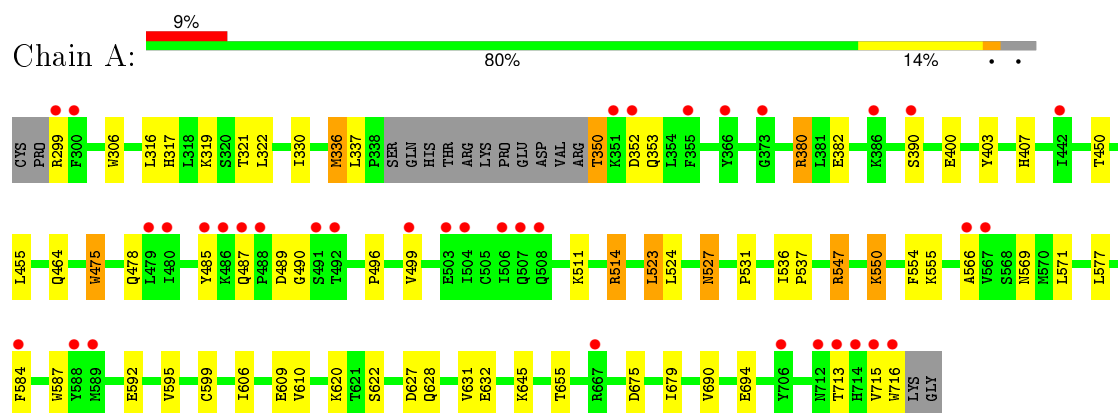
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	69	Total	O	0	0
			69	69		
7	B	93	Total	O	0	0
			93	93		

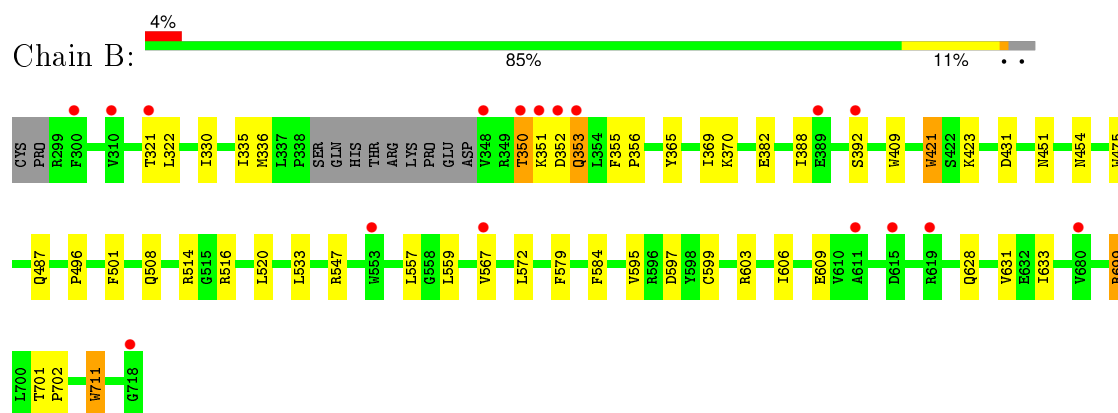
3 Residue-property plots [i](#)

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	51.66Å 110.72Å 163.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.20 40.64 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.4 (50.00-2.20) 99.1 (40.64-2.20)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.199 , 0.248 0.201 , 0.248	Depositor DCC
R_{free} test set	2431 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	39.1	Xtriage
Anisotropy	0.833	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 38.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	5 of 48413 reflections (0.010%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7024	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.24% 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: HEM, ZN, H4B, 1EW, 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	0.73	2/3415 (0.1%)	0.75	2/4633 (0.0%)
1	B	0.80	2/3453 (0.1%)	0.79	0/4681
All	All	0.77	4/6868 (0.1%)	0.77	2/9314 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	711	TRP	CD2-CE2	5.59	1.48	1.41
1	B	421	TRP	CD2-CE2	5.47	1.48	1.41
1	A	716	TRP	CD2-CE2	5.35	1.47	1.41
1	A	475	TRP	CD2-CE2	5.30	1.47	1.41

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	547	ARG	NE-CZ-NH2	5.58	123.09	120.30
1	A	547	ARG	NE-CZ-NH1	-5.57	117.51	120.30

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	3319	0	3227	46	0
1	B	3354	0	3270	35	0
2	A	43	0	30	3	0
2	B	43	0	30	6	0
3	A	17	0	15	0	0
3	B	17	0	15	0	0
4	A	30	0	33	5	0
4	B	30	0	33	1	0
5	A	4	0	3	0	0
5	B	4	0	3	0	0
6	A	1	0	0	0	0
7	A	69	0	0	1	0
7	B	93	0	0	3	0
All	All	7024	0	6659	84	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 6.

All (84) 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:350:THR:N	1:A:353:GLN:HE21	1.47	1.12
1:A:350:THR:N	1:A:353:GLN:NE2	2.18	0.92
1:A:523:LEU:HD22	1:A:531:PRO:HB2	1.61	0.81
1:B:352:ASP:HB2	7:B:928:HOH:O	1.84	0.78
1:A:694:GLU:HB3	1:B:335:ILE:HD13	1.74	0.69
1:A:306:TRP:CD2	1:B:336:MET:HE2	2.30	0.67
1:A:321:THR:HG23	1:A:322:LEU:HG	1.79	0.64
2:A:801:HEM:HMC2	2:A:801:HEM:HBC2	1.77	0.64
2:B:801:HEM:HBB2	2:B:801:HEM:HHC	1.80	0.61
1:A:380:ARG:HD3	1:A:400:GLU:OE1	2.00	0.61
1:B:409:TRP:CE3	1:B:421:TRP:HA	2.39	0.57
2:B:801:HEM:HBC2	2:B:801:HEM:CMC	2.35	0.57
1:A:306:TRP:CG	1:B:336:MET:CE	2.88	0.56
1:A:306:TRP:CG	1:B:336:MET:HE1	2.42	0.55
1:B:322:LEU:HD12	1:B:699:ARG:HG2	1.90	0.54
2:B:801:HEM:HBC2	2:B:801:HEM:HMC1	1.90	0.54
1:A:306:TRP:CE2	1:B:336:MET:HE2	2.43	0.53
2:B:801:HEM:O1D	4:B:804:IEW:N21	2.41	0.53
1:A:403:TYR:CE1	1:A:407:HIS:CE1	2.97	0.52
1:B:557:LEU:HD22	1:B:609:GLU:OE2	2.09	0.52
1:B:699:ARG:HH11	1:B:699:ARG:HB2	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:803:1EW:H12	4:A:803:1EW:H9	1.91	0.51
1:A:524:LEU:O	1:A:531:PRO:HA	2.09	0.51
2:A:801:HEM:CMC	2:A:801:HEM:HBC2	2.41	0.51
1:A:306:TRP:CD1	1:B:336:MET:HE1	2.47	0.50
1:A:606:ILE:HA	1:A:609:GLU:OE1	2.11	0.50
1:A:299:ARG:O	1:A:317:HIS:CE1	2.65	0.49
1:B:431:ASP:OD2	7:B:945:HOH:O	2.20	0.49
1:A:337:LEU:HD21	4:A:803:1EW:H3	1.96	0.48
1:B:352:ASP:CB	7:B:928:HOH:O	2.51	0.47
1:B:501:PHE:HD2	1:B:520:LEU:HD13	1.78	0.47
1:A:478:GLN:HA	1:A:566:ALA:O	2.15	0.47
4:A:803:1EW:C17	4:A:803:1EW:H9	2.44	0.47
1:B:388:ILE:O	1:B:392:SER:N	2.39	0.47
1:A:628:GLN:HG2	1:B:631:VAL:HG11	1.96	0.46
1:B:701:THR:HA	1:B:702:PRO:C	2.35	0.46
1:A:475:TRP:HB2	1:A:523:LEU:HB3	1.98	0.46
1:A:592:GLU:OE1	4:A:803:1EW:N01	2.49	0.46
1:A:455:LEU:HD12	1:A:587:TRP:HB3	1.99	0.45
1:A:321:THR:HG23	1:A:322:LEU:H	1.80	0.45
2:A:801:HEM:HHC	2:A:801:HEM:HBB2	1.99	0.45
1:A:485:TYR:HB3	1:A:514:ARG:NH1	2.32	0.44
1:B:475:TRP:CZ3	1:B:711:TRP:HB3	2.53	0.44
1:A:675:ASP:O	1:A:679:ILE:HG12	2.17	0.44
1:A:487:GLN:O	1:A:490:GLY:N	2.51	0.44
1:A:527:ASN:HD22	1:A:527:ASN:HA	1.62	0.44
1:B:597:ASP:OD1	1:B:603:ARG:NH1	2.50	0.43
4:A:803:1EW:H12	4:A:803:1EW:C28	2.48	0.43
1:A:450:THR:HA	1:A:455:LEU:HD22	1.99	0.43
1:A:316:LEU:HD12	1:A:319:LYS:HE2	2.00	0.43
1:A:496:PRO:O	1:A:499:VAL:HG23	2.19	0.43
1:B:501:PHE:CD2	1:B:520:LEU:HD13	2.53	0.43
1:A:571:LEU:C	1:A:571:LEU:HD23	2.39	0.43
1:A:350:THR:OG1	1:A:353:GLN:NE2	2.52	0.43
1:A:536:ILE:O	1:A:537:PRO:C	2.57	0.42
1:A:554:PHE:HB3	7:A:910:HOH:O	2.18	0.42
1:A:321:THR:HG23	1:A:322:LEU:N	2.34	0.42
1:B:487:GLN:OE1	1:B:514:ARG:NH2	2.53	0.42
1:B:599:CYS:SG	1:B:633:ILE:HD13	2.59	0.42
1:B:451:ASN:HB3	1:B:454:ASN:O	2.20	0.42
1:B:595:VAL:O	1:B:599:CYS:HB2	2.19	0.42
2:B:801:HEM:CBB	2:B:801:HEM:HHC	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:632:GLU:OE2	1:B:628:GLN:NE2	2.49	0.42
1:B:508:GLN:HB2	1:B:533:LEU:HD11	2.02	0.41
1:A:336:MET:HB3	1:A:336:MET:HE2	1.93	0.41
1:B:369:ILE:O	1:B:370:LYS:HB2	2.20	0.41
1:A:330:ILE:HG23	1:A:330:ILE:O	2.20	0.41
1:B:350:THR:OG1	1:B:353:GLN:NE2	2.53	0.41
1:B:365:TYR:CE2	1:B:369:ILE:HD13	2.55	0.41
1:A:595:VAL:O	1:A:599:CYS:HB2	2.21	0.41
1:A:550:LYS:HB2	1:A:550:LYS:NZ	2.35	0.41
1:A:382:GLU:OE1	1:A:382:GLU:HA	2.20	0.41
1:A:523:LEU:HD22	1:A:531:PRO:CB	2.41	0.41
1:A:627:ASP:O	1:A:631:VAL:HG23	2.20	0.41
1:A:464:GLN:NE2	1:A:577:LEU:HD22	2.36	0.41
1:B:409:TRP:CH2	2:B:801:HEM:HMC3	2.56	0.40
1:A:655:THR:HB	1:A:690:VAL:HG22	2.02	0.40
1:B:572:LEU:HB3	1:B:579:PHE:HB2	2.02	0.40
1:B:567:VAL:HB	1:B:584:PHE:CZ	2.57	0.40
1:A:606:ILE:O	1:A:610:VAL:HG23	2.21	0.40
1:A:566:ALA:HA	1:A:584:PHE:O	2.22	0.40
1:B:559:LEU:HD11	1:B:606:ILE:CD1	2.52	0.40
1:B:355:PHE:N	1:B:356:PRO:HD2	2.36	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	404/422 (96%)	391 (97%)	12 (3%)	1 (0%)	52	59
1	B	409/422 (97%)	393 (96%)	16 (4%)	0	100	100
All	All	813/844 (96%)	784 (96%)	28 (3%)	1 (0%)	56	64

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	364/377 (97%)	346 (95%)	18 (5%)	31	36
1	B	368/377 (98%)	358 (97%)	10 (3%)	52	64
All	All	732/754 (97%)	704 (96%)	28 (4%)	40	49

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	336	MET
1	A	350	THR
1	A	352	ASP
1	A	380	ARG
1	A	390	SER
1	A	511	LYS
1	A	514	ARG
1	A	523	LEU
1	A	527	ASN
1	A	547	ARG
1	A	550	LYS
1	A	555	LYS
1	A	569	ASN
1	A	620	LYS
1	A	622	SER
1	A	645	LYS
1	A	713	THR
1	A	715	VAL
1	B	321	THR
1	B	330	ILE
1	B	350	THR
1	B	351	LYS

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Mol	Chain	Res	Type
1	B	353	GLN
1	B	382	GLU
1	B	423	LYS
1	B	516	ARG
1	B	547	ARG
1	B	699	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	353	GLN
1	A	454	ASN
1	A	527	ASN
1	A	569	ASN
1	A	605	ASN
1	A	697	ASN
1	B	353	GLN
1	B	364	GLN
1	B	454	ASN
1	B	478	GLN
1	B	535	GLN
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	801	1	30,50,50	2.70	14 (46%)	24,82,82	2.95	11 (45%)
3	H4B	A	802	-	13,18,18	0.69	0	11,26,26	2.57	6 (54%)
4	1EW	A	803	-	32,32,32	0.81	2 (6%)	43,43,43	2.05	11 (25%)
5	ACT	A	804	-	1,3,3	1.24	0	0,3,3	0.00	-
2	HEM	B	801	1	30,50,50	2.54	12 (40%)	24,82,82	3.12	9 (37%)
3	H4B	B	802	-	13,18,18	1.11	1 (7%)	11,26,26	2.71	5 (45%)
5	ACT	B	803	-	1,3,3	1.70	0	0,3,3	0.00	-
4	1EW	B	804	-	32,32,32	0.61	0	43,43,43	1.95	8 (18%)

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	-	0/10/54/54	0/0/8/8
3	H4B	A	802	-	-	0/8/17/17	0/2/2/2
4	1EW	A	803	-	-	0/14/14/14	0/3/3/3
5	ACT	A	804	-	-	0/0/0/0	0/0/0/0
2	HEM	B	801	1	-	0/10/54/54	0/0/8/8
3	H4B	B	802	-	-	0/8/17/17	0/2/2/2
5	ACT	B	803	-	-	0/0/0/0	0/0/0/0
4	1EW	B	804	-	-	0/14/14/14	0/3/3/3

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	801	HEM	C3B-C4B	-4.73	1.47	1.51
2	A	801	HEM	CAD-C3D	-4.08	1.45	1.54
2	A	801	HEM	C3D-C4D	-3.95	1.46	1.51
2	A	801	HEM	C3B-C4B	-3.89	1.48	1.51
2	A	801	HEM	C2D-C3D	-3.35	1.44	1.54
2	B	801	HEM	C3D-C4D	-3.30	1.47	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	801	HEM	C2C-C1C	-3.24	1.46	1.52
2	B	801	HEM	C2D-C3D	-2.95	1.45	1.54
2	A	801	HEM	C3B-CAB	-2.70	1.46	1.51
2	B	801	HEM	C3B-CAB	-2.44	1.46	1.51
4	A	803	1EW	C05-C04	-2.05	1.35	1.38
4	A	803	1EW	C26-N21	2.01	1.38	1.34
2	B	801	HEM	FE-NB	2.06	2.08	1.97
2	A	801	HEM	CHD-C1D	2.12	1.44	1.38
2	B	801	HEM	CHC-C4B	2.14	1.44	1.38
2	A	801	HEM	C4A-CHB	2.20	1.45	1.39
3	B	802	H4B	C4-N3	2.51	1.37	1.33
2	A	801	HEM	CHC-C4B	2.67	1.46	1.38
2	A	801	HEM	FE-NB	2.75	2.12	1.97
2	A	801	HEM	C2A-C3A	2.96	1.46	1.37
2	A	801	HEM	CHC-C1C	3.15	1.43	1.36
2	B	801	HEM	C2A-C3A	3.39	1.47	1.37
2	B	801	HEM	CHC-C1C	3.59	1.44	1.36
2	B	801	HEM	CHD-C4C	4.32	1.46	1.36
2	A	801	HEM	CHD-C4C	4.65	1.47	1.36
2	B	801	HEM	C4C-NC	4.87	1.42	1.36
2	B	801	HEM	C1C-NC	5.66	1.43	1.36
2	A	801	HEM	C1C-NC	5.67	1.43	1.36
2	A	801	HEM	C4C-NC	6.10	1.43	1.36

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	HEM	CBA-CAA-C2A	-7.48	99.13	112.53
2	A	801	HEM	CBA-CAA-C2A	-6.28	101.27	112.53
4	A	803	1EW	C28-C26-C25	-4.80	114.39	121.13
4	B	804	1EW	C28-C29-C15	-4.26	105.81	112.76
4	B	804	1EW	C05-C06-N01	-3.53	118.97	122.96
3	A	802	H4B	N3-C2-N1	-3.31	120.10	125.53
4	A	803	1EW	C05-C06-N01	-3.27	119.27	122.96
4	A	803	1EW	C29-C15-C16	-3.10	112.83	119.31
2	A	801	HEM	CBD-CAD-C3D	-2.76	105.51	113.55
4	B	804	1EW	C09-C11-C12	-2.73	114.35	121.25
3	B	802	H4B	N3-C2-N1	-2.65	121.19	125.53
2	B	801	HEM	C3C-CAC-CBC	-2.56	120.52	124.46
2	A	801	HEM	CAD-CBD-CGD	-2.38	103.30	113.02
2	A	801	HEM	C3B-CAB-CBB	-2.28	120.95	124.46
4	B	804	1EW	C18-C17-C14	-2.27	106.77	113.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	HEM	C2C-C1C-CHC	-2.09	120.50	123.68
4	A	803	1EW	N02-C02-N01	2.11	120.34	116.50
4	A	803	1EW	C29-C15-C14	2.50	127.53	122.70
3	B	802	H4B	C4A-C8A-N8	2.71	121.63	118.43
4	A	803	1EW	C02-N01-C06	2.83	120.24	118.23
4	B	804	1EW	C08-C06-N01	2.88	119.94	115.69
3	A	802	H4B	N2-C2-N3	3.11	122.34	117.20
2	A	801	HEM	C2C-C1C-NC	3.12	115.46	110.21
2	B	801	HEM	C2C-C1C-NC	3.13	115.49	110.21
3	B	802	H4B	C2-N1-C8A	3.18	121.69	114.54
3	A	802	H4B	C2-N1-C8A	3.29	121.94	114.54
3	A	802	H4B	C4A-C8A-N8	3.46	122.51	118.43
4	A	803	1EW	N22-C22-N21	3.47	122.82	116.50
2	A	801	HEM	CMD-C2D-C3D	3.49	129.78	114.35
4	A	803	1EW	C08-C06-N01	3.50	120.86	115.69
4	B	804	1EW	C09-C11-C16	3.52	126.47	120.56
4	A	803	1EW	C28-C29-C15	3.53	118.51	112.76
3	A	802	H4B	C4-N3-C2	3.56	120.88	115.94
2	B	801	HEM	CMD-C2D-C3D	3.66	130.52	114.35
2	A	801	HEM	CAA-CBA-CGA	3.82	119.75	112.75
3	A	802	H4B	C4-C4A-C8A	4.03	118.21	114.56
2	A	801	HEM	CAD-C3D-C4D	4.20	127.27	112.47
4	A	803	1EW	C22-N21-C26	4.31	121.29	118.23
3	B	802	H4B	N2-C2-N3	4.54	124.72	117.20
2	A	801	HEM	CMB-C2B-C3B	4.69	128.23	116.53
2	B	801	HEM	CAD-C3D-C4D	4.76	129.25	112.47
2	B	801	HEM	CAD-C3D-C2D	5.07	127.80	113.22
3	B	802	H4B	C4-C4A-C8A	5.09	119.17	114.56
2	A	801	HEM	CAD-C3D-C2D	5.21	128.20	113.22
4	B	804	1EW	C22-N21-C26	5.24	121.95	118.23
2	B	801	HEM	CMC-C2C-C3C	5.64	130.60	116.53
4	B	804	1EW	C02-N01-C06	5.83	122.37	118.23
2	A	801	HEM	CMC-C2C-C3C	5.85	131.12	116.53
4	A	803	1EW	C28-C26-N21	6.02	124.58	115.69
2	B	801	HEM	CMB-C2B-C3B	6.54	132.86	116.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	HEM	3	0
4	A	803	1EW	5	0
2	B	801	HEM	6	0
4	B	804	1EW	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	407/422 (96%)	0.50	36 (8%) 12 11	31, 58, 96, 127	0
1	B	411/422 (97%)	0.04	17 (4%) 41 39	31, 47, 74, 101	0
All	All	818/844 (96%)	0.27	53 (6%) 22 22	31, 52, 91, 127	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	488	PRO	7.2
1	A	715	VAL	6.0
1	A	716	TRP	5.7
1	B	300	PHE	5.4
1	A	713	THR	4.2
1	B	619	ARG	4.1
1	A	507	GLN	4.0
1	A	566	ALA	3.4
1	A	485	TYR	3.2
1	B	353	GLN	3.2
1	A	352	ASP	3.2
1	A	355	PHE	3.2
1	B	310	VAL	3.0
1	A	503	GLU	3.0
1	A	300	PHE	3.0
1	B	351	LYS	2.9
1	A	487	GLN	2.9
1	B	389	GLU	2.9
1	A	506	ILE	2.9
1	A	479	LEU	2.9
1	A	390	SER	2.8
1	A	504	ILE	2.8
1	A	486	LYS	2.8
1	A	714	HIS	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	567	VAL	2.7
1	A	712	ASN	2.7
1	A	480	ILE	2.7
1	A	667	ARG	2.7
1	A	508	GLN	2.6
1	A	706	TYR	2.6
1	B	350	THR	2.5
1	B	615	ASP	2.5
1	B	611	ALA	2.5
1	A	492	THR	2.5
1	B	321	THR	2.5
1	A	299	ARG	2.5
1	B	392	SER	2.4
1	B	718	GLY	2.4
1	A	588	TYR	2.3
1	A	373	GLY	2.3
1	A	584	PHE	2.3
1	A	366	TYR	2.2
1	A	499	VAL	2.2
1	A	386	LYS	2.2
1	A	589	MET	2.1
1	B	567	VAL	2.1
1	B	553	TRP	2.1
1	B	680	VAL	2.1
1	A	491	SER	2.1
1	B	352	ASP	2.1
1	A	351	LYS	2.1
1	A	442	ILE	2.0
1	B	348	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 ⓘ

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(Å ²)	Q<0.9
4	1EW	A	803	30/30	0.86	0.35	2.75	40,77,94,97	0
2	HEM	B	801	43/43	0.98	0.18	1.01	34,37,53,66	0
4	1EW	B	804	30/30	0.90	0.20	0.80	29,55,69,76	0
5	ACT	B	803	4/4	0.95	0.13	0.76	60,60,61,63	0
2	HEM	A	801	43/43	0.97	0.22	0.76	32,39,53,59	0
3	H4B	B	802	17/17	0.91	0.17	0.20	46,49,51,53	0
3	H4B	A	802	17/17	0.93	0.15	-0.16	49,53,55,57	0
5	ACT	A	804	4/4	0.97	0.11	-0.89	59,61,63,67	0
6	ZN	A	805	1/1	0.99	0.07	-1.15	44,44,44,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.