



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:52 PM GMT

PDB ID : 4JSH  
Title : Structure of rat neuronal nitric oxide synthase heme domain in complex with  
4-methyl-6-((3-(piperidin-4-ylmethoxy)phenoxy)methyl)pyridin-2-amine  
Authors : Li, H.; Poulos, T.L.  
Deposited on : 2013-03-22  
Resolution : 2.35 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](mailto: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.

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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

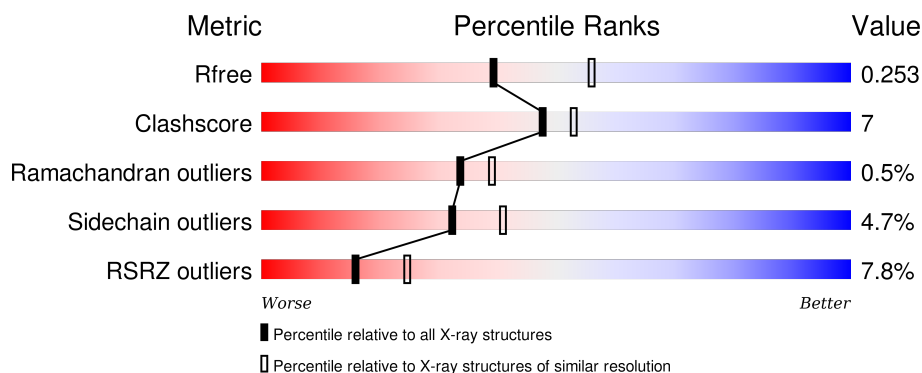
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.35 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	1352 (2.38-2.34)
Clashscore	102246	1456 (2.38-2.34)
Ramachandran outliers	100387	1435 (2.38-2.34)
Sidechain outliers	100360	1436 (2.38-2.34)
RSRZ outliers	91569	1358 (2.38-2.34)

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  $\geq 3$ , 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>10%</div> <div>79%</div> <div>19%</div> <div>..</div> </div>
1	B	422	<div> <div>5%</div> <div>78%</div> <div>20%</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
3	H4B	B	802	-	-	-	X
4	Q15	A	803	-	-	-	X
5	ACT	B	803	-	-	-	X

## 2 Entry composition [i](#)

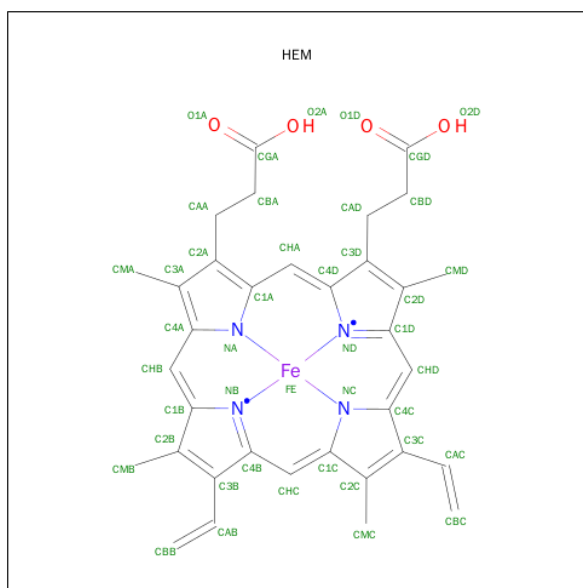
There are 7 unique types of molecules in this entry. The entry contains 7082 atoms, of which 0 are hydrogens and 0 are deuteriums.

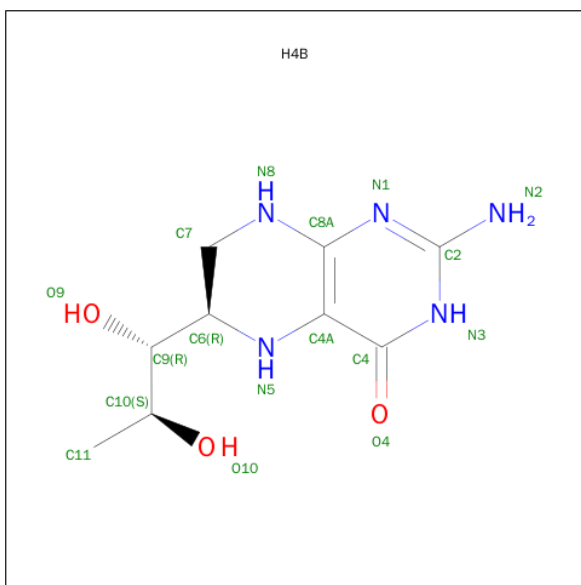
In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Nitric oxide synthase, brain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	418	Total	C	N	O	S	0	0	0
			3407	2176	587	623	21			
1	B	420	Total	C	N	O	S	0	1	0
			3424	2186	590	626	22			

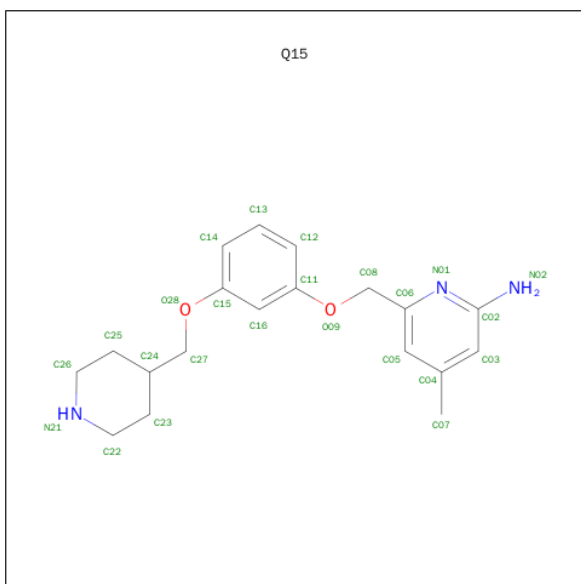
- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula:  $C_{34}H_{32}FeN_4O_4$ ).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
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-METHYL-6-{{3-(PIPERIDIN-4-YLMETHOXY)PHENOXY}METHYL}PYRIDIN-2-AMINE (three-letter code: Q15) (formula: C<sub>19</sub>H<sub>25</sub>N<sub>3</sub>O<sub>2</sub>).



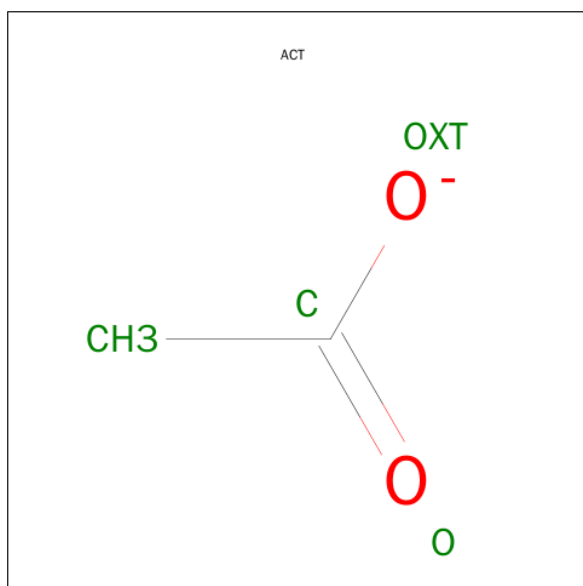
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			24	19	3	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	N	O	0	0
			24	19	3	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		

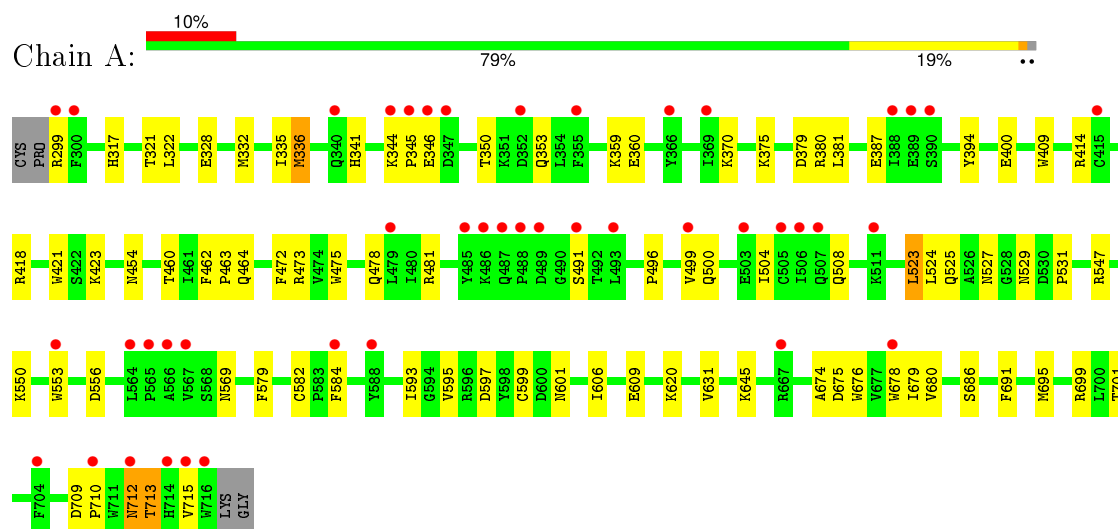
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	30	Total	O	0	0
			30	30		
7	B	44	Total	O	0	0
			44	44		

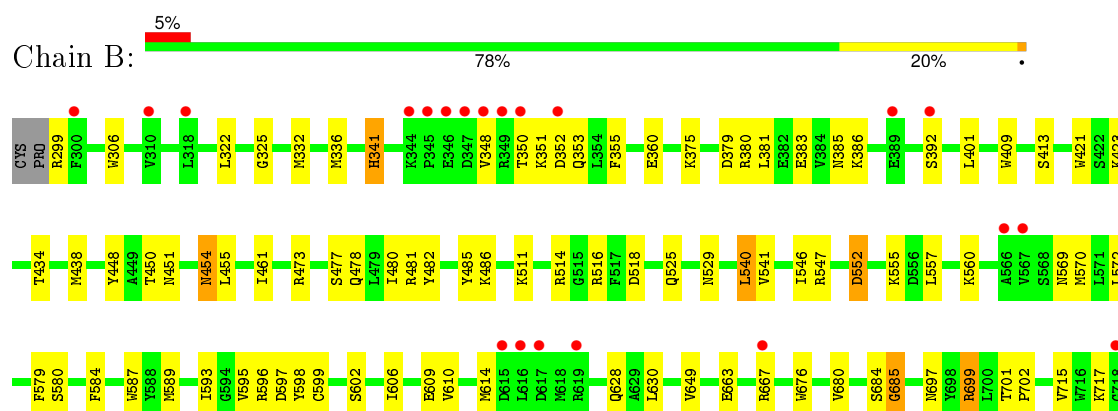
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Nitric oxide synthase, brain



- Molecule 1: Nitric oxide synthase, brain



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	51.73Å 111.56Å 164.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.78 – 2.35 40.75 – 2.33	Depositor EDS
% Data completeness (in resolution range)	99.1 (40.78-2.35) 98.1 (40.75-2.33)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.98 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, $R_{free}$	0.193 , 0.253 0.193 , 0.253	Depositor DCC
$R_{free}$ test set	1988 reflections (5.20%)	DCC
Wilson B-factor (Å <sup>2</sup> )	45.5	Xtriage
Anisotropy	0.939	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 43.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 40836 reflections (0.002%)	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	7082	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.27% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical 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: Q15, 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	0.55	0/3503	0.70	1/4753 (0.0%)
1	B	0.63	0/3523	0.74	0/4777
All	All	0.59	0/7026	0.72	1/9530 (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	A	418	ARG	NE-CZ-NH1	-5.15	117.72	120.30

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	3407	0	3314	40	0
1	B	3424	0	3335	53	0
2	A	43	0	30	8	0
2	B	43	0	30	6	0
3	A	17	0	15	0	0
3	B	17	0	15	1	0
4	A	24	0	25	3	0
4	B	24	0	25	6	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	30	0	0	0	0
7	B	44	0	0	2	0
All	All	7082	0	6795	102	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 7.

All (102) 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	1.51	0.90
1:B:350:THR:H	1:B:353:GLN:HE21	1.18	0.86
1:B:699:ARG:HH11	1:B:699:ARG:HG2	1.41	0.85
1:A:341:HIS:HA	1:A:699:ARG:NH1	2.06	0.70
1:B:593:ILE:HA	1:B:597:ASP:HB2	1.75	0.69
1:A:341:HIS:HA	1:A:699:ARG:HH12	1.59	0.67
1:B:595:VAL:O	1:B:599:CYS:HB2	1.95	0.66
1:A:321:THR:HG21	1:A:701:THR:HG22	1.79	0.64
2:A:801:HEM:HMD2	2:A:801:HEM:HBD2	1.81	0.61
1:A:595:VAL:O	1:A:599:CYS:HB2	2.02	0.60
2:B:801:HEM:HBA1	4:B:804:Q15:H8	1.85	0.59
1:B:587:TRP:H	2:B:801:HEM:HAB	1.68	0.58
1:A:345:PRO:HG3	1:A:713:THR:HG21	1.85	0.58
1:B:557:LEU:HD22	1:B:609:GLU:OE1	2.02	0.58
1:B:351:LYS:HE3	1:B:392:SER:OG	2.04	0.58
1:B:477:SER:HB2	1:B:569:ASN:ND2	2.19	0.57
1:A:332:MET:HE3	1:A:335:ILE:HG13	1.86	0.57
2:A:801:HEM:HBD1	4:A:803:Q15:C13	2.35	0.57
1:B:380:ARG:NH1	1:B:383:GLU:OE2	2.37	0.56
1:B:409:TRP:CE3	1:B:421:TRP:HA	2.43	0.54
2:A:801:HEM:CMD	2:A:801:HEM:HBD2	2.36	0.54
1:B:663:GLU:HB3	1:B:667:ARG:NH1	2.23	0.54
1:A:332:MET:CE	1:A:335:ILE:HG13	2.37	0.54
1:B:336:MET:CE	4:B:804:Q15:H17	2.39	0.53
1:B:485:TYR:CZ	1:B:514:ARG:HA	2.44	0.53
1:A:322:LEU:HB3	1:A:699:ARG:HH21	1.74	0.53
1:B:525:GLN:HG3	1:B:529:ASN:O	2.09	0.53
1:A:524:LEU:O	1:A:531:PRO:HA	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:801:HEM:HH C	2:A:801:HEM:HBB2	1.90	0.52
2:A:801:HEM:CM C	2:A:801:HEM:HBC2	2.40	0.52
1:A:478:GLN:HB2	1:A:481:ARG:HG3	1.92	0.52
1:B:584:PHE:CD1	2:B:801:HEM:CAC	2.93	0.52
1:A:464:GLN:HB3	1:A:579:PHE:CE2	2.44	0.51
1:B:336:MET:HE1	4:B:804:Q15:H17	1.93	0.51
1:B:699:ARG:HH11	1:B:699:ARG:CG	2.16	0.51
1:B:701:THR:HA	1:B:702:PRO:C	2.30	0.51
1:A:593:ILE:O	1:A:597:ASP:HB2	2.11	0.51
1:A:500:GLN:O	1:A:504:ILE:HG13	2.11	0.50
1:B:478:GLN:HB2	1:B:481:ARG:HG3	1.92	0.50
2:B:801:HEM:HBC2	2:B:801:HEM:CM C	2.41	0.50
3:B:802:H4B:O4	4:B:804:Q15:H20	2.12	0.50
1:A:525:GLN:HG3	1:A:529:ASN:O	2.11	0.50
1:A:712:ASN:OD1	1:A:712:ASN:N	2.43	0.50
1:B:699:ARG:NH1	1:B:699:ARG:HG2	2.19	0.50
1:B:413:SER:OG	1:B:697:ASN:ND2	2.45	0.49
1:B:480:ILE:HA	7:B:933:HOH:O	2.11	0.49
1:A:675:ASP:O	1:A:679:ILE:HG12	2.12	0.49
1:B:434:THR:O	1:B:438:MET:HG3	2.13	0.49
1:A:380:ARG:HD3	1:A:400:GLU:OE1	2.12	0.49
1:A:460:THR:O	1:A:582:CYS:HA	2.13	0.49
1:A:299:ARG:O	1:A:317:HIS:CE1	2.66	0.48
2:A:801:HEM:HBD1	4:A:803:Q15:H12	1.95	0.48
1:A:631:VAL:HG11	1:B:628:GLN:HG3	1.96	0.48
1:A:674:ALA:HB3	1:A:695:MET:HB3	1.96	0.48
1:A:496:PRO:HA	1:A:499:VAL:HG23	1.95	0.48
1:B:572:LEU:HB3	1:B:579:PHE:HB2	1.96	0.47
1:A:387:GLU:OE1	1:A:394:TYR:HA	2.15	0.47
1:A:686:SER:HA	1:A:691:PHE:CG	2.50	0.47
1:B:341:HIS:NE2	2:B:801:HEM:O2D	2.42	0.47
2:B:801:HEM:HBD1	4:B:804:Q15:C13	2.44	0.47
2:A:801:HEM:C1C	4:A:803:Q15:H6	2.50	0.46
1:B:570:MET:SD	4:B:804:Q15:H12	2.56	0.46
1:A:606:ILE:HA	1:A:609:GLU:OE1	2.17	0.45
1:B:480:ILE:HD13	1:B:541:VAL:HG13	1.98	0.45
1:A:414:ARG:HD3	1:A:678:TRP:CD2	2.52	0.45
1:B:598:TYR:HB3	1:B:606:ILE:HG12	1.98	0.45
1:B:325:GLY:O	1:B:332:MET:HG3	2.16	0.45
1:A:676:TRP:CZ2	1:A:680:VAL:HG21	2.52	0.45
1:A:350:THR:OG1	1:A:353:GLN:HG3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:359:LYS:HE2	1:A:381:LEU:HD21	1.99	0.44
1:B:595:VAL:HG13	1:B:630:LEU:HD11	1.99	0.44
1:B:448:TYR:CD2	1:B:448:TYR:C	2.91	0.44
1:B:557:LEU:CD2	1:B:609:GLU:OE1	2.65	0.44
1:A:409:TRP:CE3	1:A:421:TRP:HA	2.53	0.44
1:B:684:SER:O	1:B:685:GLY:C	2.56	0.43
1:B:401:LEU:HD11	1:B:461:ILE:HG21	2.00	0.43
1:A:336:MET:HG3	1:B:306:TRP:NE1	2.35	0.42
1:B:473:ARG:HD3	1:B:580:SER:HB2	2.02	0.42
1:B:299:ARG:CZ	1:B:299:ARG:HB3	2.50	0.42
1:A:475:TRP:HB2	1:A:523:LEU:HB3	2.01	0.42
1:A:553:TRP:O	1:A:556:ASP:HB2	2.20	0.42
1:A:462:PHE:HB3	1:A:463:PRO:CD	2.49	0.42
1:B:546:ILE:HG12	1:B:560:LYS:HA	2.01	0.42
1:B:353:GLN:HG2	1:B:353:GLN:H	1.66	0.41
1:B:451:ASN:HB3	1:B:454:ASN:O	2.20	0.41
1:B:482:TYR:HA	1:B:518:ASP:O	2.20	0.41
1:A:375:LYS:HG2	1:A:379:ASP:OD1	2.20	0.41
1:B:552:ASP:OD1	1:B:555:LYS:NZ	2.53	0.41
1:B:589:MET:HA	1:B:649:VAL:O	2.21	0.41
1:B:540:LEU:HD12	1:B:540:LEU:HA	1.89	0.41
1:B:381:LEU:HA	1:B:381:LEU:HD23	1.92	0.41
1:B:610:VAL:O	1:B:614:MET:HG3	2.21	0.41
1:B:375:LYS:HE2	1:B:379:ASP:OD2	2.21	0.41
1:A:321:THR:HG21	1:A:701:THR:CG2	2.47	0.41
1:B:676:TRP:CZ2	1:B:680:VAL:HG21	2.56	0.41
1:A:472:PHE:HA	1:A:525:GLN:O	2.21	0.40
1:A:709:ASP:HA	1:A:710:PRO:HD2	1.83	0.40
1:B:355:PHE:CE1	1:B:385:ASN:HB2	2.57	0.40
1:B:596:ARG:HD3	7:B:926:HOH:O	2.22	0.40
1:B:715:VAL:O	1:B:717:LYS:HE3	2.21	0.40
1:B:450:THR:O	1:B:455:LEU:HD12	2.21	0.40
1:A:584:PHE:CD1	2:A:801:HEM:CAC	3.05	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	416/422 (99%)	394 (95%)	21 (5%)	1 (0%)	52	63
1	B	419/422 (99%)	400 (96%)	16 (4%)	3 (1%)	26	29
All	All	835/844 (99%)	794 (95%)	37 (4%)	4 (0%)	34	39

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	352	ASP
1	B	685	GLY
1	B	322	LEU
1	A	491	SER

### 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	374/377 (99%)	353 (94%)	21 (6%)	26	31
1	B	376/377 (100%)	362 (96%)	14 (4%)	41	53
All	All	750/754 (100%)	715 (95%)	35 (5%)	32	41

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	328	GLU
1	A	336	MET

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Mol	Chain	Res	Type
1	A	344	LYS
1	A	346	GLU
1	A	360	GLU
1	A	370	LYS
1	A	423	LYS
1	A	454	ASN
1	A	473	ARG
1	A	508	GLN
1	A	523	LEU
1	A	527	ASN
1	A	547	ARG
1	A	550	LYS
1	A	569	ASN
1	A	601	ASN
1	A	620	LYS
1	A	645	LYS
1	A	712	ASN
1	A	713	THR
1	A	715	VAL
1	B	341	HIS
1	B	348	VAL
1	B	360	GLU
1	B	386	LYS
1	B	423	LYS
1	B	454	ASN
1	B	486	LYS
1	B	511	LYS
1	B	516	ARG
1	B	540	LEU
1	B	547	ARG
1	B	552	ASP
1	B	602	SER
1	B	699	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	454	ASN
1	A	507	GLN
1	A	508	GLN
1	A	527	ASN
1	A	535	GLN

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Mol	Chain	Res	Type
1	A	569	ASN
1	A	605	ASN
1	A	642	GLN
1	A	697	ASN
1	B	353	GLN
1	B	364	GLN
1	B	385	ASN
1	B	425	GLN
1	B	454	ASN
1	B	601	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	801	1	30,50,50	2.10	8 (26%)	24,82,82	2.46	9 (37%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	H4B	A	802	-	13,18,18	0.87	0	11,26,26	2.41	4 (36%)
4	Q15	A	803	-	26,26,26	0.52	0	32,34,34	1.57	4 (12%)
5	ACT	A	804	-	1,3,3	2.09	1 (100%)	0,3,3	0.00	-
2	HEM	B	801	1	30,50,50	1.95	7 (23%)	24,82,82	2.64	12 (50%)
3	H4B	B	802	-	13,18,18	0.91	0	11,26,26	2.31	5 (45%)
5	ACT	B	803	-	1,3,3	2.57	1 (100%)	0,3,3	0.00	-
4	Q15	B	804	-	26,26,26	0.63	0	32,34,34	1.54	4 (12%)

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	Q15	A	803	-	-	0/10/18/18	1/3/3/3
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	Q15	B	804	-	-	0/10/18/18	0/3/3/3

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	HEM	C2D-C3D	-6.23	1.35	1.54
2	A	801	HEM	C2C-C1C	-6.21	1.40	1.52
2	B	801	HEM	C2C-C1C	-5.93	1.41	1.52
2	B	801	HEM	C2D-C3D	-5.70	1.37	1.54
2	A	801	HEM	C2B-C1B	-2.75	1.42	1.51
2	B	801	HEM	C3D-C4D	-2.57	1.48	1.51
2	B	801	HEM	C2B-C1B	-2.38	1.44	1.51
2	A	801	HEM	C3D-C4D	-2.37	1.48	1.51
2	A	801	HEM	C3B-C4B	-2.34	1.49	1.51
2	B	801	HEM	C3B-CAB	-2.00	1.47	1.51
2	A	801	HEM	C3C-CAC	-2.00	1.47	1.51
2	B	801	HEM	C1C-NC	2.04	1.38	1.36
5	A	804	ACT	CH3-C	2.09	1.51	1.48
2	A	801	HEM	FE-NB	2.34	2.09	1.97
5	B	803	ACT	CH3-C	2.57	1.52	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	801	HEM	FE-NC	2.82	2.06	1.95
2	A	801	HEM	FE-NC	3.13	2.08	1.95

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	HEM	C3C-CAC-CBC	-4.28	117.89	124.46
3	B	802	H4B	N3-C2-N1	-3.44	119.90	125.53
2	A	801	HEM	C1D-CHD-C4C	-3.41	120.12	125.82
4	A	803	Q15	C04-C05-C06	-3.39	118.15	120.28
2	B	801	HEM	C1D-CHD-C4C	-3.11	120.62	125.82
2	B	801	HEM	C3B-CAB-CBB	-3.09	119.72	124.46
3	A	802	H4B	N3-C2-N1	-2.95	120.69	125.53
4	B	804	Q15	C05-C06-N01	-2.51	120.12	122.96
2	B	801	HEM	C3B-C4B-NB	-2.42	107.00	111.63
2	B	801	HEM	CBA-CAA-C2A	-2.42	108.20	112.53
2	B	801	HEM	C3C-CAC-CBC	-2.39	120.79	124.46
2	A	801	HEM	C3B-C4B-NB	-2.24	107.35	111.63
2	B	801	HEM	CAA-CBA-CGA	2.15	116.68	112.75
2	A	801	HEM	C2D-C3D-C4D	2.23	105.28	101.50
2	A	801	HEM	CMD-C2D-C3D	2.45	125.19	114.35
3	A	802	H4B	C2-N1-C8A	2.46	120.06	114.54
4	B	804	Q15	C22-N21-C26	2.49	118.60	110.33
3	B	802	H4B	C2-N1-C8A	2.70	120.61	114.54
4	A	803	Q15	C22-N21-C26	2.96	120.16	110.33
3	B	802	H4B	N2-C2-N1	3.00	122.16	117.20
4	B	804	Q15	O28-C27-C24	3.20	115.35	107.97
2	A	801	HEM	CAD-C3D-C4D	3.24	123.90	112.47
2	B	801	HEM	CMD-C2D-C3D	3.40	129.37	114.35
3	B	802	H4B	C4-N3-C2	3.51	120.80	115.94
2	B	801	HEM	CAD-C3D-C4D	3.57	125.05	112.47
2	A	801	HEM	CMB-C2B-C3B	3.66	125.67	116.53
3	A	802	H4B	C4-N3-C2	3.69	121.06	115.94
2	B	801	HEM	CBD-CAD-C3D	3.87	124.81	113.55
4	A	803	Q15	C02-N01-C06	3.92	121.01	118.23
2	B	801	HEM	CMC-C2C-C3C	3.94	126.36	116.53
3	B	802	H4B	C4-C4A-C8A	4.04	118.22	114.56
4	A	803	Q15	C27-O28-C15	4.09	126.69	117.89
2	A	801	HEM	CMC-C2C-C3C	4.12	126.82	116.53
2	B	801	HEM	CMB-C2B-C3B	4.17	126.93	116.53
4	B	804	Q15	C02-N01-C06	4.75	121.61	118.23
3	A	802	H4B	C4-C4A-C8A	5.31	119.37	114.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	HEM	CAD-C3D-C2D	6.12	130.82	113.22
2	B	801	HEM	CAD-C3D-C2D	6.37	131.54	113.22

There are no chirality outliers.

There are no torsion outliers.

All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	803	Q15	C22-C23-C24-C25-C26-N21

5 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	HEM	8	0
4	A	803	Q15	3	0
2	B	801	HEM	6	0
3	B	802	H4B	1	0
4	B	804	Q15	6	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	418/422 (99%)	0.66	44 (10%) 8 14	39, 69, 121, 155	0
1	B	420/422 (99%)	0.26	21 (5%) 32 47	34, 57, 90, 131	0
All	All	838/844 (99%)	0.46	65 (7%) 16 25	34, 61, 111, 155	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	345	PRO	6.5
1	A	716	TRP	5.8
1	A	715	VAL	5.7
1	B	300	PHE	4.9
1	A	488	PRO	4.6
1	A	486	LYS	4.6
1	B	718	GLY	4.5
1	B	345	PRO	4.4
1	A	344	LYS	4.3
1	A	347	ASP	4.3
1	A	507	GLN	4.0
1	B	619	ARG	4.0
1	A	567	VAL	3.9
1	A	355	PHE	3.9
1	A	390	SER	3.8
1	B	347	ASP	3.6
1	A	493	LEU	3.5
1	B	346	GLU	3.4
1	A	487	GLN	3.4
1	A	389	GLU	3.4
1	A	506	ILE	3.3
1	A	503	GLU	3.2
1	A	584	PHE	3.2
1	A	714	HIS	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	566	ALA	3.0
1	B	348	VAL	2.9
1	A	352	ASP	2.9
1	B	352	ASP	2.9
1	B	567	VAL	2.8
1	A	489	ASP	2.7
1	A	565	PRO	2.7
1	A	678	TRP	2.7
1	A	340	GLN	2.6
1	B	667	ARG	2.6
1	A	366	TYR	2.6
1	A	588	TYR	2.6
1	A	388	ILE	2.6
1	A	712	ASN	2.6
1	A	491	SER	2.5
1	A	479	LEU	2.5
1	A	346	GLU	2.5
1	A	299	ARG	2.5
1	B	615	ASP	2.4
1	A	485	TYR	2.4
1	A	505	CYS	2.4
1	A	710	PRO	2.4
1	B	350	THR	2.4
1	B	310	VAL	2.3
1	A	300	PHE	2.3
1	B	392	SER	2.3
1	B	344	LYS	2.2
1	A	499	VAL	2.2
1	B	616	LEU	2.2
1	A	369	ILE	2.2
1	B	389	GLU	2.2
1	B	566	ALA	2.2
1	A	511	LYS	2.1
1	A	667	ARG	2.1
1	B	349	ARG	2.1
1	B	617	ASP	2.1
1	A	415	CYS	2.1
1	A	553	TRP	2.1
1	A	704	PHE	2.0
1	A	564	LEU	2.0
1	B	318	LEU	2.0

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

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

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
5	ACT	B	803	4/4	0.74	0.35	8.42	64,71,72,74	0
4	Q15	A	803	24/24	0.93	0.38	2.80	49,60,105,111	0
3	H4B	B	802	17/17	0.94	0.23	2.02	52,57,63,64	0
3	H4B	A	802	17/17	0.97	0.20	0.84	48,55,62,64	0
2	HEM	A	801	43/43	0.97	0.23	0.57	37,44,58,68	0
2	HEM	B	801	43/43	0.98	0.17	0.41	37,45,64,72	0
5	ACT	A	804	4/4	0.94	0.14	-0.05	56,61,64,65	0
6	ZN	A	805	1/1	0.99	0.12	-0.44	55,55,55,55	0
4	Q15	B	804	24/24	0.96	0.17	-0.47	30,51,82,82	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.