



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:37 AM GMT

PDB ID : 2HX4  
Title : Rat nNOS heme domain complexed with 4-N-(Nw-nitro-L-argininyl)-trans-4-hydroxyamino-L-proline amide  
Authors : Igarashi, J.; Li, H.; Poulos, T.L.  
Deposited on : 2006-08-02  
Resolution : 2.15 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.

---

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

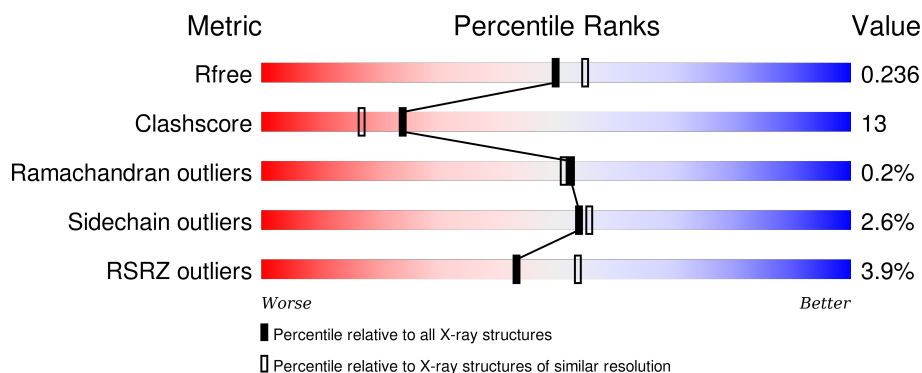
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.15 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	1045 (2.16-2.16)
Clashscore	102246	1152 (2.16-2.16)
Ramachandran outliers	100387	1131 (2.16-2.16)
Sidechain outliers	100360	1131 (2.16-2.16)
RSRZ outliers	91569	1050 (2.16-2.16)

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	
1	B	422	

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

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

## 2 Entry composition [i](#)

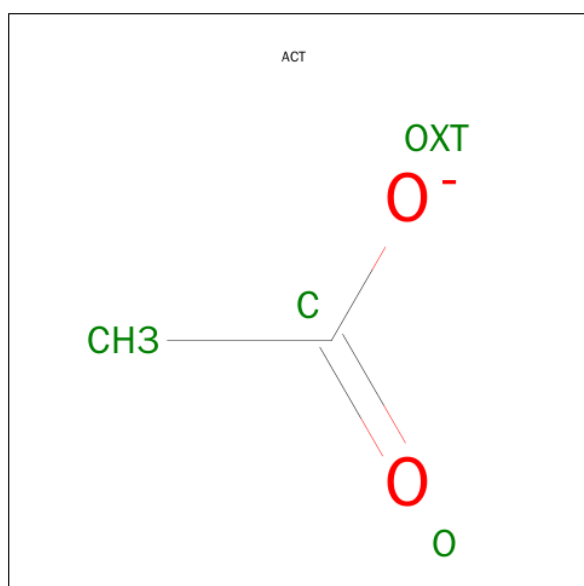
There are 7 unique types of molecules in this entry. The entry contains 7248 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	407	Total	C	N	O	S	0	0	0
			3313	2121	566	605	21			
1	B	411	Total	C	N	O	S	0	0	0
			3345	2140	574	610	21			

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula:  $C_2H_3O_2$ ).

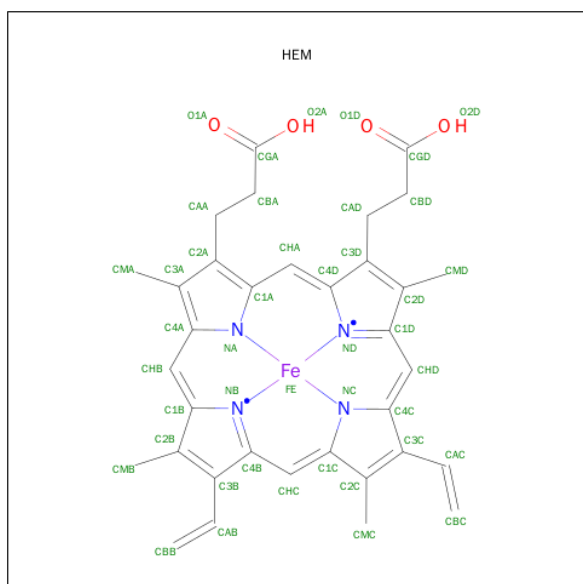


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

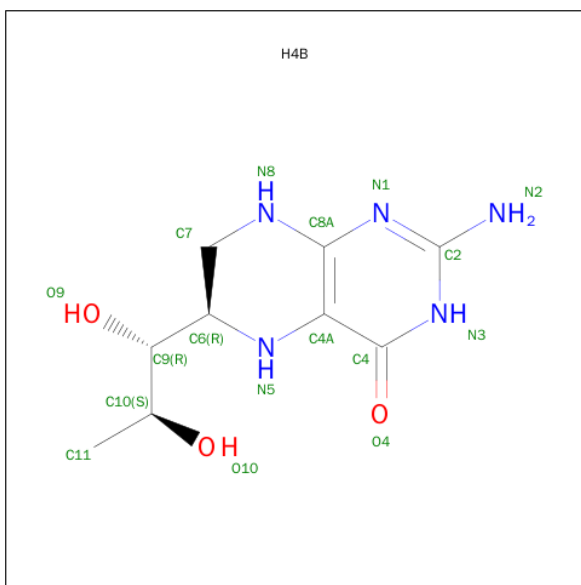
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Zn 1 1	0	0

- Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula:  $C_{34}H_{32}FeN_4O_4$ ).



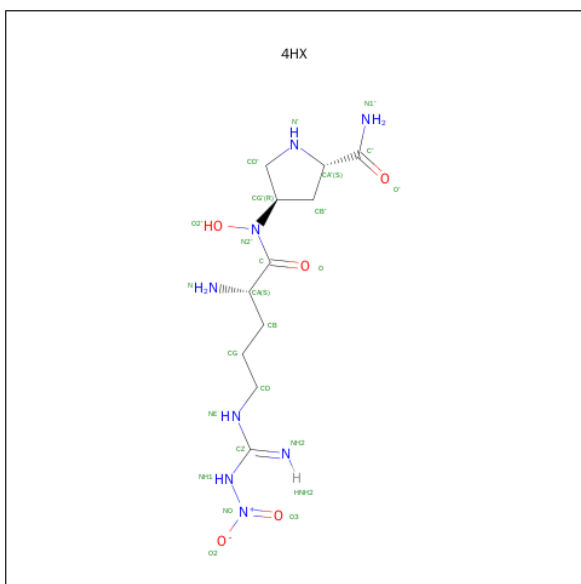
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 5 is 5,6,7,8-TETRAHYDROBIOPTERIN (three-letter code: H4B) (formula:  $\text{C}_9\text{H}_{15}\text{N}_5\text{O}_3$ ).



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

- Molecule 6 is (4R)-4-(HYDROXY{N 5 -[IMINO(NITROAMINO)METHYL]-L-ORNITHYL}AMINO)-L-PROLINAMIDE (three-letter code: 4HX) (formula: C<sub>11</sub>H<sub>22</sub>N<sub>8</sub>O<sub>5</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			24	11	8	5		

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	N	O	0	0
			24	11	8	5		

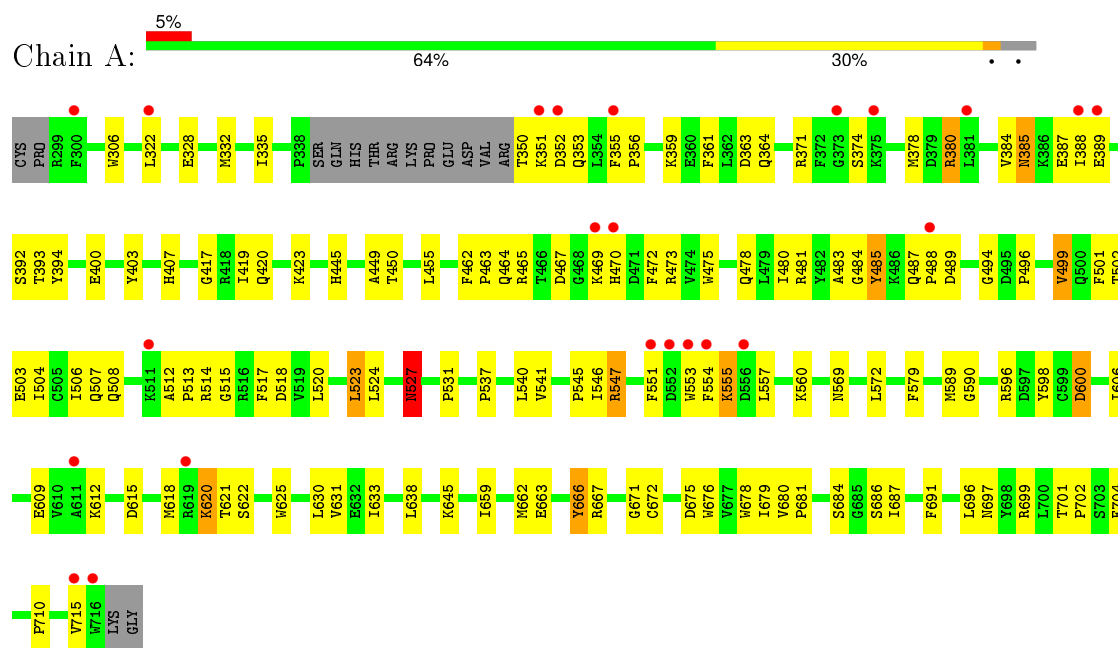
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	171	Total	O	0	0
			171	171		
7	B	242	Total	O	0	0
			242	242		

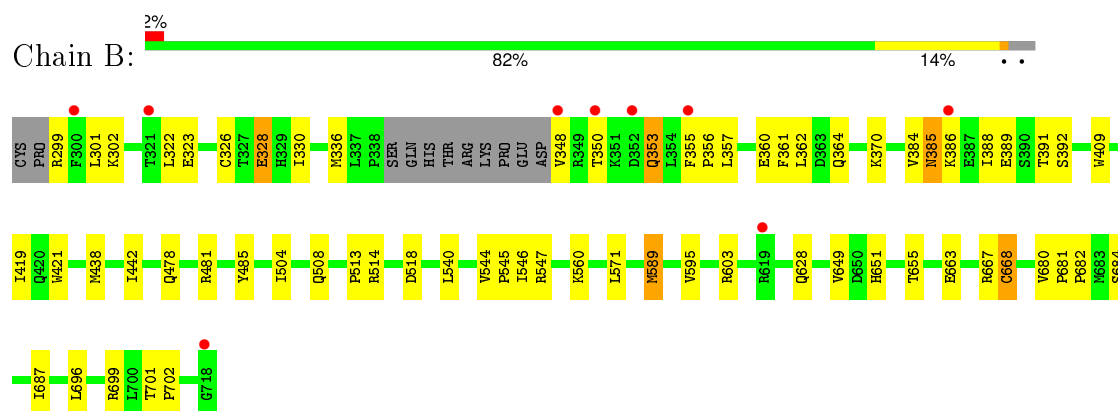
### 3 Residue-property plots [i](#)

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



#### • Molecule 1: Nitric-oxide synthase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	51.92Å 111.00Å 164.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.51 – 2.15 49.51 – 2.15	Depositor EDS
% Data completeness (in resolution range)	98.2 (49.51-2.15) 98.3 (49.51-2.15)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.77 (at 2.16Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.203 , 0.245 0.198 , 0.236	Depositor DCC
$R_{free}$ test set	2603 reflections (5.31%)	DCC
Wilson B-factor (Å <sup>2</sup> )	29.9	Xtriage
Anisotropy	0.659	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 51.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 51616 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7248	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.70% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, ZN, ACT, H4B, 4HX

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	0.44	0/3406	0.66	1/4621 (0.0%)
1	B	0.45	0/3438	0.66	2/4661 (0.0%)
All	All	0.44	0/6844	0.66	3/9282 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	B	326	CYS	CA-CB-SG	6.27	125.29	114.00
1	B	589	MET	N-CA-C	-5.84	95.23	111.00
1	A	589	MET	N-CA-C	-5.65	95.75	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	666	TYR	Sidechain

## 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	3313	0	3221	127	0
1	B	3345	0	3259	55	0
2	A	4	0	3	2	0
2	B	4	0	3	1	0
3	A	1	0	0	0	0
4	A	43	0	30	2	0
4	B	43	0	30	0	0
5	A	17	0	15	1	0
5	B	17	0	15	0	0
6	A	24	0	21	0	0
6	B	24	0	21	1	0
7	A	171	0	0	14	0
7	B	242	0	0	4	0
All	All	7248	0	6618	176	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 13.

All (176) 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:630:LEU:HD12	7:A:1016:HOH:O	1.75	0.86
1:A:523:LEU:HD22	1:A:531:PRO:HB2	1.57	0.85
1:A:467:ASP:OD2	1:A:469:LYS:HB2	1.77	0.85
1:B:478:GLN:HB2	1:B:481:ARG:HG3	1.66	0.77
1:B:299:ARG:HD2	7:B:1060:HOH:O	1.85	0.75
1:A:555:LYS:HB3	1:A:555:LYS:NZ	2.02	0.74
1:A:662:MET:HE3	7:A:1035:HOH:O	1.89	0.72
1:A:473:ARG:NH2	1:A:710:PRO:HD3	2.05	0.71
1:A:672:CYS:HB3	7:A:1035:HOH:O	1.91	0.71
1:A:470:HIS:HB3	1:A:527:ASN:ND2	2.07	0.69
1:A:508:GLN:NE2	1:A:508:GLN:HA	2.10	0.67
4:A:750:HEM:HMC2	4:A:750:HEM:HBC2	1.77	0.67
1:A:508:GLN:HA	1:A:508:GLN:HE21	1.61	0.66
1:A:517:PHE:HB2	1:A:560:LYS:HE2	1.78	0.65

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:371:ARG:HG3	1:A:371:ARG:HH21	1.62	0.64
1:A:696:LEU:HD22	1:B:330:ILE:HD11	1.78	0.63
1:A:507:GLN:O	1:A:507:GLN:HG2	1.99	0.63
1:A:322:LEU:HB2	1:A:699:ARG:HB2	1.80	0.63
1:A:332:MET:HE1	1:B:301:LEU:HD22	1.81	0.62
1:B:668:CYS:HB3	7:B:964:HOH:O	1.99	0.62
1:A:350:THR:HG22	1:A:352:ASP:H	1.65	0.62
1:A:630:LEU:HA	7:A:1016:HOH:O	1.99	0.61
1:B:419:ILE:HG22	2:B:860:ACT:H2	1.81	0.61
1:A:506:ILE:C	1:A:508:GLN:H	2.03	0.61
1:B:322:LEU:HD13	1:B:699:ARG:NH2	2.15	0.61
1:A:378:MET:HA	1:A:378:MET:CE	2.31	0.61
1:A:546:ILE:HG12	1:A:560:LYS:HA	1.81	0.60
1:A:496:PRO:HA	1:A:499:VAL:HG23	1.83	0.60
1:A:351:LYS:HE2	1:A:392:SER:HB3	1.84	0.60
1:B:355:PHE:N	1:B:356:PRO:HD2	2.17	0.59
1:B:386:LYS:O	1:B:389:GLU:HG2	2.01	0.59
4:A:750:HEM:HBB2	4:A:750:HEM:HHC	1.84	0.59
1:A:551:PHE:HB3	1:A:553:TRP:NE1	2.18	0.58
1:A:465:ARG:HG3	7:A:1064:HOH:O	2.04	0.58
1:B:513:PRO:HG2	1:B:518:ASP:CG	2.24	0.58
1:A:662:MET:HG2	7:A:1035:HOH:O	2.02	0.58
1:A:485:TYR:CE2	1:A:512:ALA:HB1	2.38	0.58
1:B:684:SER:HB3	1:B:687:ILE:HD11	1.85	0.57
1:A:555:LYS:HB3	1:A:555:LYS:HZ3	1.66	0.57
1:A:484:GLY:O	1:A:499:VAL:HA	2.04	0.57
1:A:417:GLY:HA2	2:A:860:ACT:H2	1.84	0.57
1:A:332:MET:HE2	1:B:696:LEU:HD21	1.86	0.56
1:A:485:TYR:HE2	1:A:512:ALA:HB1	1.69	0.56
1:A:551:PHE:HB3	1:A:553:TRP:CE2	2.41	0.55
1:A:499:VAL:O	1:A:503:GLU:HG3	2.07	0.55
1:A:378:MET:HA	1:A:378:MET:HE2	1.88	0.55
1:A:596:ARG:O	1:A:600:ASP:HB2	2.05	0.55
1:A:332:MET:CE	1:B:696:LEU:HD21	2.37	0.55
1:A:696:LEU:HD22	1:B:330:ILE:CD1	2.37	0.54
1:A:463:PRO:HB2	1:A:472:PHE:CE1	2.43	0.54
1:B:701:THR:HA	1:B:702:PRO:C	2.28	0.54
1:B:323:GLU:O	1:B:699:ARG:HD3	2.08	0.54
1:A:501:PHE:HD2	1:A:520:LEU:HD13	1.72	0.54
1:B:684:SER:HB3	1:B:687:ILE:CG1	2.38	0.54
1:B:391:THR:O	1:B:392:SER:HB2	2.07	0.54

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:478:GLN:HB2	1:A:481:ARG:HG3	1.90	0.53
1:B:328:GLU:H	1:B:328:GLU:CD	2.11	0.53
1:B:357:LEU:O	1:B:360:GLU:HG3	2.09	0.53
1:A:322:LEU:HD13	1:A:699:ARG:NH2	2.24	0.53
1:A:355:PHE:N	1:A:356:PRO:HD2	2.24	0.53
1:A:384:VAL:O	1:A:388:ILE:HG13	2.09	0.52
1:A:350:THR:HG22	1:A:352:ASP:N	2.24	0.52
1:A:350:THR:HB	1:A:353:GLN:CD	2.30	0.52
1:B:364:GLN:NE2	7:B:1083:HOH:O	2.42	0.52
1:A:351:LYS:HE2	1:A:392:SER:CB	2.40	0.51
1:A:322:LEU:HD13	1:A:699:ARG:HH21	1.76	0.51
1:A:475:TRP:HB2	1:A:523:LEU:HB3	1.92	0.51
1:A:631:VAL:HG11	1:B:628:GLN:CG	2.41	0.51
1:A:598:TYR:O	1:A:606:ILE:HG12	2.11	0.51
1:A:699:ARG:HD2	7:A:1008:HOH:O	2.11	0.50
1:A:501:PHE:CD2	1:A:520:LEU:HD13	2.47	0.50
1:A:667:ARG:HG3	1:A:667:ARG:HH11	1.76	0.50
1:B:355:PHE:CE1	1:B:385:ASN:HB2	2.47	0.50
1:A:489:ASP:O	1:A:489:ASP:OD2	2.29	0.50
1:A:684:SER:HB3	1:A:687:ILE:HD11	1.94	0.50
1:A:387:GLU:OE1	1:A:394:TYR:HA	2.12	0.50
1:A:385:ASN:HD22	1:A:385:ASN:C	2.16	0.49
1:A:361:PHE:O	1:A:364:GLN:HG2	2.13	0.49
1:A:701:THR:HA	1:A:702:PRO:C	2.33	0.49
1:B:362:LEU:HD11	1:B:384:VAL:HG21	1.94	0.49
1:A:506:ILE:C	1:A:508:GLN:N	2.65	0.49
1:A:508:GLN:CA	1:A:508:GLN:HE21	2.20	0.49
1:A:545:PRO:HG2	1:A:547:ARG:NH2	2.28	0.49
1:A:715:VAL:O	1:A:715:VAL:HG23	2.12	0.49
1:A:609:GLU:HG3	7:A:1027:HOH:O	2.13	0.48
1:B:438:MET:O	1:B:442:ILE:HG13	2.12	0.48
6:B:799:4HX:HG'	7:B:943:HOH:O	2.14	0.48
1:A:380:ARG:HD3	1:A:400:GLU:OE1	2.13	0.48
1:A:554:PHE:HB3	7:A:1065:HOH:O	2.13	0.48
1:A:393:THR:OG1	1:A:394:TYR:N	2.47	0.48
1:A:371:ARG:CG	1:A:371:ARG:HH21	2.27	0.47
1:A:353:GLN:O	1:A:356:PRO:HG2	2.15	0.47
1:A:620:LYS:HE3	1:A:622:SER:OG	2.14	0.47
1:B:299:ARG:HB3	1:B:299:ARG:HH11	1.79	0.47
1:A:374:SER:O	1:A:378:MET:HG2	2.14	0.47
1:A:631:VAL:HG11	1:B:628:GLN:HG2	1.96	0.47

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:684:SER:HB3	1:A:687:ILE:HG12	1.96	0.47
1:A:554:PHE:O	1:A:557:LEU:HB2	2.15	0.47
1:A:551:PHE:CD2	1:A:551:PHE:N	2.81	0.47
1:A:659:ILE:O	1:A:663:GLU:HG3	2.15	0.47
1:B:328:GLU:N	1:B:328:GLU:OE1	2.48	0.47
1:B:589:MET:HA	1:B:649:VAL:O	2.15	0.47
1:A:618:MET:HG2	1:A:625:TRP:CD2	2.50	0.46
1:A:524:LEU:O	1:A:531:PRO:HA	2.16	0.46
1:B:388:ILE:O	1:B:392:SER:N	2.44	0.46
1:A:475:TRP:CE2	1:A:710:PRO:HB2	2.51	0.46
1:A:388:ILE:O	1:A:392:SER:HA	2.15	0.46
1:A:675:ASP:O	1:A:679:ILE:HG12	2.16	0.46
1:B:409:TRP:CE3	1:B:421:TRP:HA	2.51	0.46
1:B:546:ILE:HG12	1:B:560:LYS:HA	1.97	0.46
1:A:555:LYS:HB3	1:A:555:LYS:HZ2	1.79	0.46
1:A:504:ILE:O	1:A:508:GLN:HG2	2.15	0.46
1:A:332:MET:HB3	1:A:335:ILE:HG13	1.97	0.45
1:A:609:GLU:HG3	7:A:1034:HOH:O	2.16	0.45
1:B:504:ILE:O	1:B:508:GLN:HG2	2.17	0.45
1:A:464:GLN:HB3	1:A:579:PHE:CE2	2.51	0.45
1:B:684:SER:HB3	1:B:687:ILE:CD1	2.46	0.45
1:A:678:TRP:HA	5:A:760:H4B:N1	2.31	0.45
1:A:483:ALA:HB1	1:A:502:THR:CG2	2.47	0.44
1:B:651:HIS:O	1:B:655:THR:HG23	2.17	0.44
1:B:571:LEU:HD23	1:B:571:LEU:C	2.37	0.44
1:A:494:GLY:O	1:A:496:PRO:HD3	2.18	0.44
1:A:462:PHE:HB3	1:A:463:PRO:CD	2.47	0.44
1:B:302:LYS:HE3	1:B:302:LYS:HB2	1.67	0.44
1:A:419:ILE:HG22	2:A:860:ACT:H1	2.00	0.44
1:A:470:HIS:HB3	1:A:527:ASN:HD21	1.82	0.44
1:B:370:LYS:HA	1:B:370:LYS:HD3	1.89	0.44
1:A:633:ILE:HD12	7:A:1016:HOH:O	2.18	0.43
1:A:508:GLN:NE2	7:A:1002:HOH:O	2.51	0.43
1:A:306:TRP:CD1	1:B:336:MET:HE2	2.53	0.43
1:A:485:TYR:HE2	1:A:512:ALA:CB	2.32	0.43
1:A:485:TYR:CZ	1:A:514:ARG:HA	2.53	0.43
1:A:676:TRP:CZ2	1:A:680:VAL:HG21	2.53	0.43
1:A:684:SER:HB3	1:A:687:ILE:CG1	2.48	0.43
1:B:603:ARG:HH11	1:B:603:ARG:HG3	1.83	0.43
1:B:348:VAL:HG22	1:B:348:VAL:O	2.19	0.43
1:B:595:VAL:HG11	1:B:682:PRO:HB2	2.00	0.43

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:403:TYR:CE1	1:A:407:HIS:CE1	3.07	0.43
1:A:684:SER:HB3	1:A:687:ILE:CD1	2.48	0.43
1:B:299:ARG:CB	1:B:299:ARG:NH1	2.82	0.42
1:A:545:PRO:HG2	1:A:547:ARG:HH21	1.83	0.42
1:A:480:ILE:HD13	1:A:541:VAL:HG13	1.99	0.42
1:B:485:TYR:CE1	1:B:514:ARG:HA	2.53	0.42
1:B:544:VAL:HA	1:B:545:PRO:HD3	1.88	0.42
1:A:666:TYR:HA	1:A:671:GLY:N	2.34	0.42
1:B:299:ARG:HB3	1:B:299:ARG:NH1	2.34	0.42
1:B:485:TYR:CZ	1:B:514:ARG:HA	2.55	0.42
1:B:680:VAL:HA	1:B:681:PRO:HD3	1.91	0.42
1:A:359:LYS:HG2	1:A:363:ASP:OD2	2.19	0.42
1:A:449:ALA:O	1:A:455:LEU:HA	2.20	0.42
1:A:546:ILE:HG13	1:A:598:TYR:OH	2.19	0.42
1:A:483:ALA:HB1	1:A:502:THR:HG23	2.01	0.42
1:A:667:ARG:HG3	1:A:667:ARG:NH1	2.34	0.42
1:A:572:LEU:HD13	1:A:704:PHE:CE2	2.55	0.41
1:A:686:SER:HA	1:A:691:PHE:CG	2.54	0.41
1:A:537:PRO:HB2	1:A:540:LEU:HG	2.02	0.41
1:A:487:GLN:HB3	1:A:488:PRO:HD2	2.03	0.41
1:A:680:VAL:HA	1:A:681:PRO:HD3	1.91	0.41
1:A:480:ILE:HA	7:A:996:HOH:O	2.20	0.41
1:B:663:GLU:HB3	1:B:667:ARG:NH1	2.35	0.41
1:A:445:HIS:C	1:A:445:HIS:CD2	2.94	0.41
1:A:512:ALA:HA	1:A:513:PRO:HD3	1.91	0.41
1:A:621:THR:HG23	7:A:1030:HOH:O	2.20	0.41
1:A:517:PHE:CB	1:A:560:LYS:HE2	2.47	0.41
1:B:353:GLN:HB3	1:B:353:GLN:HE21	1.55	0.41
1:A:590:GLY:HA3	1:A:638:LEU:HD21	2.02	0.41
1:A:515:GLY:N	1:A:518:ASP:OD2	2.41	0.40
1:A:487:GLN:HB3	1:A:488:PRO:CD	2.52	0.40
1:A:420:GLN:OE1	1:A:423:LYS:HE2	2.21	0.40
1:A:328:GLU:OE2	1:B:328:GLU:OE1	2.39	0.40
1:B:696:LEU:HA	1:B:696:LEU:HD23	1.92	0.40
1:B:360:GLU:HG3	1:B:361:PHE:H	1.87	0.40
1:A:450:THR:HA	1:A:455:LEU:HD22	2.04	0.40
1:B:350:THR:O	1:B:353:GLN:HG2	2.21	0.40
1:A:612:LYS:O	1:A:615:ASP:N	2.44	0.40
1:B:328:GLU:N	1:B:328:GLU:CD	2.75	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	403/422 (96%)	368 (91%)	33 (8%)	2 (0%)	34	26
1	B	407/422 (96%)	392 (96%)	15 (4%)	0	100	100
All	All	810/844 (96%)	760 (94%)	48 (6%)	2 (0%)	52	51

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	527	ASN
1	A	499	VAL

### 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	363/377 (96%)	350 (96%)	13 (4%)	42	40
1	B	366/377 (97%)	360 (98%)	6 (2%)	70	76
All	All	729/754 (97%)	710 (97%)	19 (3%)	54	55

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	380	ARG
1	A	385	ASN
1	A	389	GLU
1	A	485	TYR

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type
1	A	523	LEU
1	A	527	ASN
1	A	547	ARG
1	A	555	LYS
1	A	569	ASN
1	A	600	ASP
1	A	620	LYS
1	A	645	LYS
1	A	697	ASN
1	B	328	GLU
1	B	353	GLN
1	B	385	ASN
1	B	540	LEU
1	B	547	ARG
1	B	668	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	353	GLN
1	A	385	ASN
1	A	425	GLN
1	A	454	ASN
1	A	507	GLN
1	A	508	GLN
1	A	527	ASN
1	A	569	ASN
1	A	605	ASN
1	A	642	GLN
1	A	697	ASN
1	B	364	GLN
1	B	385	ASN
1	B	454	ASN
1	B	507	GLN
1	B	527	ASN
1	B	535	GLN
1	B	601	ASN
1	B	605	ASN
1	B	642	GLN
1	B	697	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

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	$\# Z  > 2$
4	HEM	A	750	1	30,50,50	2.85	10 (33%)	24,82,82	2.42	10 (41%)
5	H4B	A	760	-	13,18,18	2.35	3 (23%)	11,26,26	4.10	6 (54%)
6	4HX	A	799	-	16,24,24	0.87	0	15,32,32	2.15	4 (26%)
2	ACT	A	860	-	1,3,3	2.30	1 (100%)	0,3,3	0.00	-
4	HEM	B	750	1	30,50,50	2.57	8 (26%)	24,82,82	2.17	8 (33%)
5	H4B	B	760	-	13,18,18	2.27	3 (23%)	11,26,26	4.15	6 (54%)
6	4HX	B	799	-	16,24,24	0.74	0	15,32,32	2.09	5 (33%)
2	ACT	B	860	-	1,3,3	2.54	1 (100%)	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
4	HEM	A	750	1	-	0/10/54/54	0/0/8/8
5	H4B	A	760	-	-	0/8/17/17	0/2/2/2
6	4HX	A	799	-	-	0/16/36/36	0/1/1/1
2	ACT	A	860	-	-	0/0/0/0	0/0/0/0
4	HEM	B	750	1	-	0/10/54/54	0/0/8/8
5	H4B	B	760	-	-	0/8/17/17	0/2/2/2
6	4HX	B	799	-	-	0/16/36/36	0/1/1/1
2	ACT	B	860	-	-	0/0/0/0	0/0/0/0

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	750	HEM	C2D-C3D	-6.41	1.35	1.54
4	A	750	HEM	C3D-C4D	-6.36	1.43	1.51
4	B	750	HEM	C2D-C3D	-6.17	1.36	1.54
4	A	750	HEM	C3C-CAC	-6.16	1.39	1.51
4	B	750	HEM	C3D-C4D	-6.04	1.43	1.51
4	B	750	HEM	C3B-CAB	-5.56	1.40	1.51
4	A	750	HEM	C3B-CAB	-5.31	1.41	1.51
4	B	750	HEM	C3C-CAC	-5.24	1.41	1.51
4	A	750	HEM	C2C-C1C	-4.93	1.43	1.52
4	A	750	HEM	C3B-C4B	-4.38	1.48	1.51
4	B	750	HEM	C3B-C4B	-4.26	1.48	1.51
4	B	750	HEM	C2C-C1C	-3.87	1.45	1.52
4	A	750	HEM	C2B-C1B	-2.41	1.44	1.51
4	A	750	HEM	C2D-C1D	-2.11	1.45	1.51
4	B	750	HEM	CHD-C4C	2.14	1.41	1.36
2	A	860	ACT	CH3-C	2.30	1.52	1.48
4	B	750	HEM	C1C-NC	2.45	1.39	1.36
2	B	860	ACT	CH3-C	2.54	1.52	1.48
4	A	750	HEM	CHD-C4C	2.59	1.42	1.36
4	A	750	HEM	C1C-NC	2.68	1.39	1.36
5	B	760	H4B	C4A-N5	3.81	1.46	1.38
5	A	760	H4B	C4A-N5	4.29	1.47	1.38
5	A	760	H4B	C6-N5	4.32	1.54	1.45
5	B	760	H4B	C6-N5	4.34	1.54	1.45
5	B	760	H4B	C4-N3	4.99	1.42	1.33
5	A	760	H4B	C4-N3	5.34	1.43	1.33

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	760	H4B	N3-C2-N1	-4.81	117.65	125.53

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	760	H4B	N3-C2-N1	-4.58	118.03	125.53
6	A	799	4HX	O'-C'-N1'	-4.44	116.50	123.08
6	B	799	4HX	O'-C'-N1'	-4.37	116.60	123.08
4	B	750	HEM	CBA-CAA-C2A	-2.91	107.31	112.53
4	A	750	HEM	CBD-CAD-C3D	-2.65	105.84	113.55
4	B	750	HEM	CAA-C2A-C1A	-2.48	124.32	127.01
4	A	750	HEM	CAA-C2A-C1A	-2.20	124.62	127.01
6	B	799	4HX	CB-CA-N	-2.03	105.55	109.81
6	B	799	4HX	CB'-CA'-C'	2.09	114.58	111.15
6	A	799	4HX	CB'-CA'-C'	2.51	115.25	111.15
4	A	750	HEM	CMD-C2D-C3D	2.52	125.51	114.35
4	A	750	HEM	C3B-CAB-CBB	2.53	128.33	124.46
4	B	750	HEM	CAD-C3D-C4D	2.59	121.59	112.47
4	B	750	HEM	CMB-C2B-C3B	2.65	123.16	116.53
6	A	799	4HX	CA'-C'-N1'	2.84	121.19	116.60
5	A	760	H4B	N2-C2-N1	2.87	121.95	117.20
4	A	750	HEM	C3B-C4B-CHC	2.91	127.26	123.16
4	B	750	HEM	C2D-C3D-C4D	2.92	106.45	101.50
4	A	750	HEM	CAD-C3D-C4D	3.17	123.67	112.47
5	B	760	H4B	N2-C2-N1	3.20	122.50	117.20
6	B	799	4HX	CA'-C'-N1'	3.27	121.88	116.60
5	B	760	H4B	C2-N1-C8A	3.36	122.09	114.54
4	B	750	HEM	CMD-C2D-C3D	3.63	130.40	114.35
4	B	750	HEM	CMC-C2C-C3C	3.65	125.64	116.53
5	B	760	H4B	C4A-C8A-N8	3.68	122.77	118.43
5	A	760	H4B	C4A-C8A-N8	3.78	122.89	118.43
5	A	760	H4B	C2-N1-C8A	3.87	123.25	114.54
6	B	799	4HX	CD'-CG'-N2'	3.99	121.79	113.54
4	A	750	HEM	CMB-C2B-C3B	4.02	126.56	116.53
6	A	799	4HX	CD'-CG'-N2'	4.46	122.77	113.54
4	A	750	HEM	CAD-C3D-C2D	4.62	126.50	113.22
4	A	750	HEM	C2D-C3D-C4D	4.83	109.68	101.50
4	A	750	HEM	CMC-C2C-C3C	5.02	129.06	116.53
5	B	760	H4B	C4-N3-C2	6.12	124.44	115.94
5	A	760	H4B	C4-N3-C2	6.14	124.46	115.94
4	B	750	HEM	CAD-C3D-C2D	6.27	131.26	113.22
5	A	760	H4B	C4-C4A-C8A	8.96	122.67	114.56
5	B	760	H4B	C4-C4A-C8A	9.51	123.17	114.56

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	750	HEM	2	0
5	A	760	H4B	1	0
2	A	860	ACT	2	0
6	B	799	4HX	1	0
2	B	860	ACT	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, 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	407/422 (96%)	0.26	23 (5%) 27 37	18, 40, 70, 85	0
1	B	411/422 (97%)	-0.07	9 (2%) 65 73	19, 32, 55, 76	0
All	All	818/844 (96%)	0.09	32 (3%) 43 54	18, 35, 66, 85	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	321	THR	4.2
1	B	348	VAL	4.1
1	A	352	ASP	3.7
1	A	355	PHE	3.7
1	B	352	ASP	3.6
1	A	488	PRO	3.4
1	B	300	PHE	3.4
1	A	551	PHE	3.3
1	B	718	GLY	2.8
1	B	619	ARG	2.8
1	A	511	LYS	2.8
1	A	716	TRP	2.7
1	A	375	LYS	2.7
1	A	611	ALA	2.7
1	A	619	ARG	2.7
1	A	553	TRP	2.6
1	B	386	LYS	2.6
1	A	389	GLU	2.6
1	B	350	THR	2.5
1	A	300	PHE	2.5
1	A	554	PHE	2.5
1	A	373	GLY	2.4
1	A	552	ASP	2.4
1	A	322	LEU	2.4

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	B	355	PHE	2.3
1	A	388	ILE	2.2
1	A	381	LEU	2.2
1	A	470	HIS	2.2
1	A	715	VAL	2.1
1	A	351	LYS	2.1
1	A	556	ASP	2.0
1	A	469	LYS	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
2	ACT	A	860	4/4	0.94	0.20	4.93	47,50,50,51	0
2	ACT	B	860	4/4	0.92	0.21	4.70	49,49,50,50	0
6	4HX	B	799	24/24	0.95	0.14	1.57	25,32,46,51	0
6	4HX	A	799	24/24	0.94	0.13	1.46	27,37,53,55	0
3	ZN	A	900	1/1	1.00	0.11	0.37	30,30,30,30	0
4	HEM	B	750	43/43	0.98	0.12	0.33	19,22,27,33	0
4	HEM	A	750	43/43	0.98	0.10	-0.45	19,24,30,31	0
5	H4B	A	760	17/17	0.96	0.11	-0.45	25,27,30,31	0
5	H4B	B	760	17/17	0.97	0.11	-0.79	16,21,26,27	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.