



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:22 PM GMT

PDB ID : 4CAP
Title : Structure of rat neuronal nitric oxide synthase heme domain in complex with
7-(2-(3-(3-Fluorophenyl(propylamino)methyl))quinolin-2- amine
Authors : Li, H.; Poulos, T.L.
Deposited on : 2013-10-08
Resolution : 2.06 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

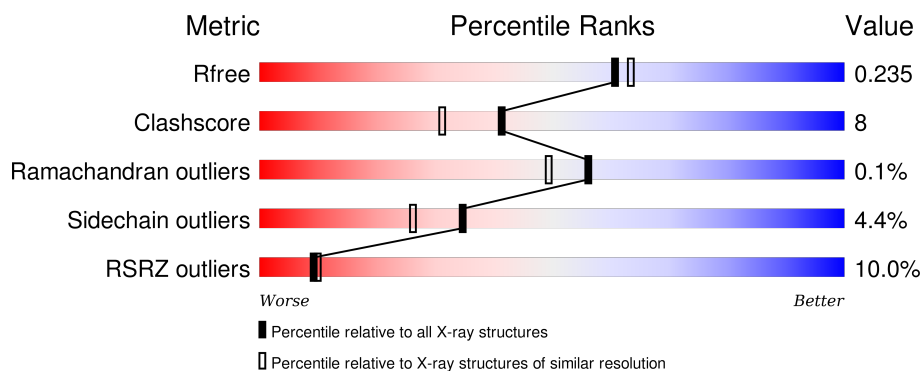
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.06 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	1799 (2.08-2.04)
Clashscore	102246	1910 (2.08-2.04)
Ramachandran outliers	100387	1893 (2.08-2.04)
Sidechain outliers	100360	1893 (2.08-2.04)
RSRZ outliers	91569	1802 (2.08-2.04)

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	
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
5	ACT	A	860	-	-	-	X
5	ACT	B	860	-	-	-	X

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 7067 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
			3326	2130	566	609	21			
1	B	411	Total	C	N	O	S	0	1	0
			3348	2142	574	611	21			

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



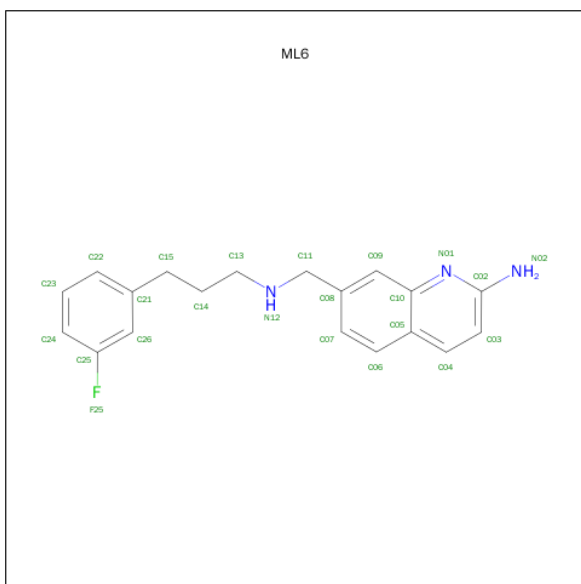
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- 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 7-([3-(3-FLUOROPHENYL)PROPYL]AMINO)METHYL)QUINOLIN-2-AMINE (three-letter code: ML6) (formula: C₁₉H₂₀FN₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	F	N	0	0
			23	19	1	3		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	F	N	0	0
			23	19	1	3		

- 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		

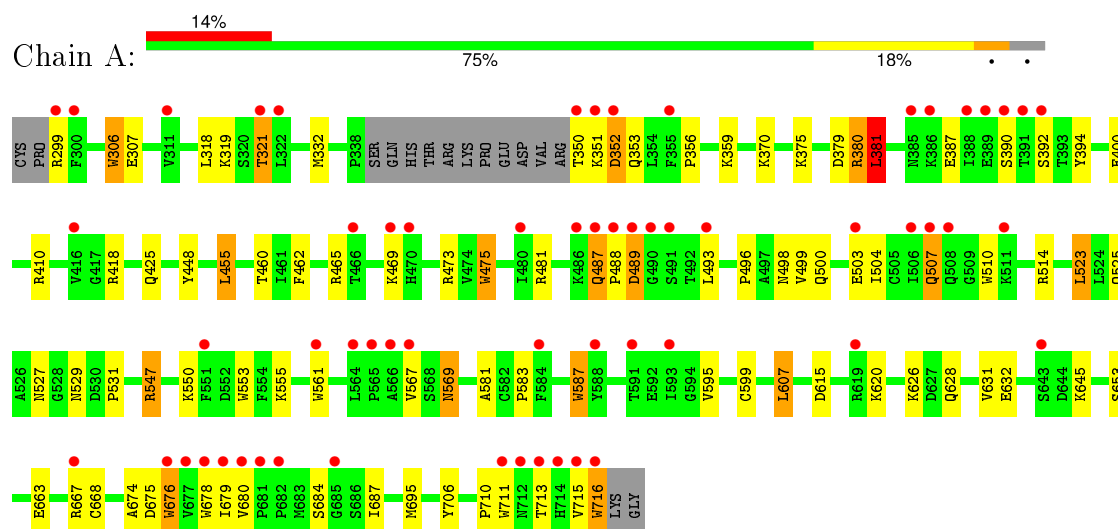
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	105	Total	O	0	0
			105	105		
7	B	113	Total	O	0	0
			113	113		

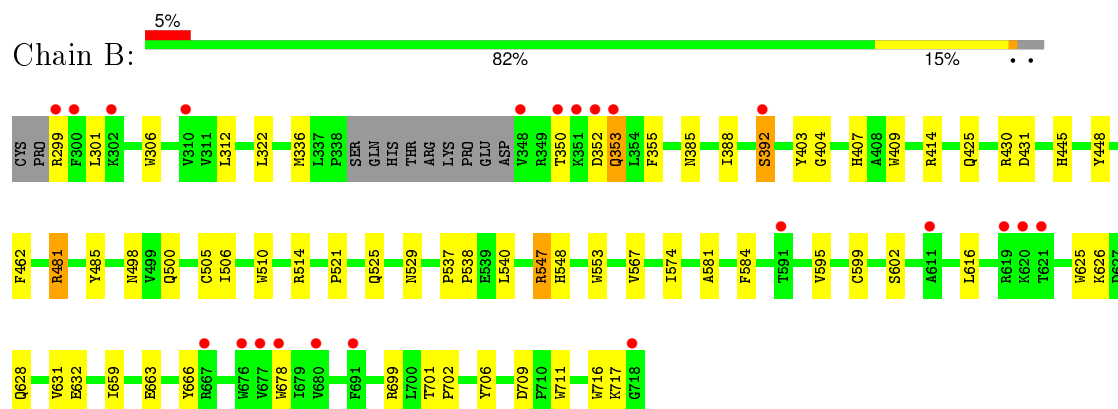
3 Residue-property plots

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, α , β , γ	51.77Å 110.83Å 165.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	92.19 – 2.06 39.06 – 2.06	Depositor EDS
% Data completeness (in resolution range)	98.7 (92.19-2.06) 98.8 (39.06-2.06)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.74 (at 2.06Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.185 , 0.235 0.183 , 0.235	Depositor DCC
R_{free} test set	2919 reflections (5.20%)	DCC
Wilson B-factor (Å ²)	40.6	Xtriage
Anisotropy	0.649	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 48.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	2 of 59136 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	7067	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 5.23% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² 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: HEM, ZN, ML6, 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	0.91	10/3428 (0.3%)	0.88	7/4652 (0.2%)
1	B	0.98	5/3444 (0.1%)	0.93	3/4669 (0.1%)
All	All	0.95	15/6872 (0.2%)	0.91	10/9321 (0.1%)

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	711	TRP	CD2-CE2	6.42	1.49	1.41
1	A	306	TRP	CD2-CE2	5.93	1.48	1.41
1	A	510	TRP	CD2-CE2	5.92	1.48	1.41
1	B	716	TRP	CD2-CE2	5.90	1.48	1.41
1	A	678	TRP	CD2-CE2	5.84	1.48	1.41
1	A	587	TRP	CD2-CE2	5.70	1.48	1.41
1	B	711	TRP	CD2-CE2	5.55	1.48	1.41
1	A	475	TRP	CD2-CE2	5.54	1.48	1.41
1	A	676	TRP	CD2-CE2	5.52	1.48	1.41
1	B	409	TRP	CD2-CE2	5.27	1.47	1.41
1	A	561	TRP	CD2-CE2	5.26	1.47	1.41
1	B	625	TRP	CD2-CE2	5.26	1.47	1.41
1	B	553	TRP	CD2-CE2	5.23	1.47	1.41
1	A	553	TRP	CD2-CE2	5.02	1.47	1.41
1	A	716	TRP	CD2-CE2	5.00	1.47	1.41

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	430	ARG	NE-CZ-NH1	10.47	125.53	120.30
1	A	410	ARG	NE-CZ-NH1	6.64	123.62	120.30
1	A	410	ARG	NE-CZ-NH2	-6.16	117.22	120.30
1	A	418	ARG	NE-CZ-NH1	-5.88	117.36	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	709	ASP	CB-CG-OD1	5.85	123.57	118.30
1	A	547	ARG	NE-CZ-NH2	-5.74	117.43	120.30
1	A	607	LEU	CA-CB-CG	5.54	128.03	115.30
1	A	455	LEU	CB-CG-CD2	5.12	119.71	111.00
1	B	431	ASP	CB-CG-OD1	5.08	122.87	118.30
1	A	381	LEU	CA-CB-CG	5.07	126.96	115.30

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	3326	0	3239	65	0
1	B	3348	0	3264	51	0
2	A	43	0	30	5	0
2	B	43	0	30	9	0
3	A	17	0	15	0	0
3	B	17	0	15	0	0
4	A	23	0	20	4	0
4	B	23	0	20	6	0
5	A	4	0	3	0	0
5	B	4	0	3	1	0
6	A	1	0	0	0	0
7	A	105	0	0	3	0
7	B	113	0	0	1	0
All	All	7067	0	6639	114	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 8.

All (114) 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)
5:B:860:ACT:H3	7:B:2027:HOH:O	1.37	1.22

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:567:VAL:HG21	4:A:800:ML6:C07	2.06	0.86
2:B:750:HEM:HHC	2:B:750:HEM:HBB2	1.63	0.80
1:A:306:TRP:CD1	1:B:336:MET:HE2	2.18	0.79
1:B:706:TYR:OH	2:B:750:HEM:O2D	2.04	0.73
2:B:750:HEM:HBC2	2:B:750:HEM:HMC2	1.73	0.71
1:B:336:MET:HG2	4:B:800:ML6:H24	1.74	0.69
1:A:306:TRP:CD2	1:B:336:MET:HE3	2.28	0.68
1:A:351:LYS:HE2	1:A:392:SER:HB3	1.73	0.68
1:B:567:VAL:HG21	4:B:800:ML6:C07	2.29	0.63
1:B:706:TYR:OH	2:B:750:HEM:CGD	2.47	0.63
1:A:487:GLN:HE21	1:A:488:PRO:HD2	1.65	0.61
2:A:750:HEM:HMC2	2:A:750:HEM:HBC2	1.83	0.60
1:A:306:TRP:CG	1:B:336:MET:CE	2.83	0.60
1:A:675:ASP:O	1:A:679:ILE:HG12	2.01	0.60
1:A:473:ARG:NH2	1:A:710:PRO:HD3	2.17	0.60
1:A:350:THR:HB	1:A:353:GLN:NE2	2.15	0.60
1:B:336:MET:HG2	4:B:800:ML6:C24	2.31	0.60
1:A:706:TYR:OH	2:A:750:HEM:O2D	2.21	0.58
1:A:306:TRP:CG	1:B:336:MET:HE2	2.38	0.58
1:A:550:LYS:HB2	1:A:550:LYS:NZ	2.16	0.58
1:A:674:ALA:HB3	1:A:695:MET:HB3	1.84	0.58
1:B:355:PHE:CE1	1:B:385:ASN:HB2	2.38	0.58
1:A:387:GLU:OE2	1:A:394:TYR:HA	2.04	0.58
1:A:352:ASP:O	1:A:356:PRO:HD2	2.04	0.57
1:B:481:ARG:HD3	1:B:498:ASN:ND2	2.20	0.57
1:A:375:LYS:NZ	1:A:379:ASP:OD2	2.36	0.56
1:A:380:ARG:HD3	1:A:400:GLU:OE2	2.05	0.56
1:A:306:TRP:CD2	1:B:336:MET:CE	2.88	0.56
1:A:706:TYR:OH	2:A:750:HEM:CGD	2.55	0.55
1:A:499:VAL:O	1:A:503:GLU:HG3	2.07	0.55
1:A:525:GLN:HG3	1:A:529:ASN:O	2.07	0.54
1:A:455:LEU:HD12	1:A:587:TRP:HB3	1.89	0.54
1:B:425:GLN:HG2	1:B:448:TYR:CZ	2.43	0.54
1:B:352:ASP:HB3	1:B:353:GLN:HE21	1.73	0.54
1:B:659:ILE:O	1:B:663:GLU:HG3	2.08	0.53
1:A:569:ASN:H	1:A:569:ASN:HD22	1.57	0.53
1:B:388:ILE:O	1:B:392:SER:N	2.40	0.53
1:A:359:LYS:HG3	1:A:381:LEU:HD11	1.91	0.52
1:A:307[B]:GLU:OE2	1:B:602:SER:N	2.33	0.52
1:A:306:TRP:CE2	1:B:336:MET:HE3	2.45	0.52
1:A:493:LEU:HA	7:A:2062:HOH:O	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:462:PHE:HB2	1:B:581:ALA:HB3	1.93	0.51
1:A:687:ILE:HD12	1:B:626:LYS:HB3	1.92	0.51
1:A:628:GLN:HG3	1:B:631:VAL:HG11	1.94	0.50
2:B:750:HEM:C4B	4:B:800:ML6:H03	2.46	0.49
1:B:481:ARG:HD3	1:B:498:ASN:HD21	1.77	0.49
1:A:321:THR:HG22	7:A:2004:HOH:O	2.12	0.49
2:A:750:HEM:C4B	4:A:800:ML6:H03	2.48	0.49
2:B:750:HEM:CHC	2:B:750:HEM:HBB2	2.36	0.49
1:B:352:ASP:HB3	1:B:353:GLN:NE2	2.28	0.48
2:B:750:HEM:HHC	2:B:750:HEM:CBB	2.37	0.48
1:A:631:VAL:HG11	1:B:628:GLN:HG3	1.94	0.48
1:A:523:LEU:HD22	1:A:531:PRO:HB2	1.96	0.48
1:A:496:PRO:O	1:A:499:VAL:HG23	2.13	0.48
1:A:684:SER:HB3	1:A:687:ILE:HG12	1.95	0.48
1:B:353:GLN:H	1:B:353:GLN:HE21	1.62	0.48
1:A:350:THR:HB	1:A:353:GLN:CD	2.35	0.47
1:A:350:THR:N	7:A:2010:HOH:O	2.46	0.47
1:A:550:LYS:NZ	1:A:550:LYS:CB	2.77	0.47
1:A:460:THR:O	1:A:583:PRO:HD2	2.15	0.47
1:A:332:MET:CE	1:B:301:LEU:HD22	2.45	0.47
1:A:465:ARG:HH11	1:A:465:ARG:HB2	1.80	0.47
1:A:595:VAL:O	1:A:599:CYS:HB2	2.14	0.47
1:A:531:PRO:HD2	1:A:716:TRP:HZ3	1.79	0.46
1:A:632:GLU:OE2	1:B:628:GLN:NE2	2.48	0.46
1:A:462:PHE:HB2	1:A:581:ALA:HB3	1.97	0.46
1:A:507:GLN:O	1:A:507:GLN:HG2	2.16	0.46
1:A:465:ARG:HB2	1:A:465:ARG:NH1	2.30	0.46
1:B:505:CYS:O	1:B:506:ILE:C	2.50	0.46
1:B:425:GLN:HG2	1:B:448:TYR:CE1	2.51	0.45
1:B:322:LEU:HD13	1:B:699:ARG:HH21	1.81	0.45
1:B:414:ARG:HD3	1:B:678:TRP:CD2	2.51	0.45
1:A:299:ARG:HG3	1:A:318:LEU:HD11	1.98	0.45
1:B:701:THR:HA	1:B:702:PRO:C	2.37	0.45
1:A:500:GLN:O	1:A:504:ILE:HG13	2.17	0.44
1:B:510:TRP:CE2	1:B:521:PRO:HD3	2.53	0.44
1:A:481:ARG:HD3	1:A:498:ASN:HD21	1.82	0.44
1:B:353:GLN:N	1:B:353:GLN:HE21	2.17	0.43
1:B:403:TYR:CE2	1:B:407:HIS:CE1	3.06	0.43
1:B:404:GLY:HA3	1:B:574:ILE:HD13	2.01	0.43
1:A:332:MET:HE3	1:B:301:LEU:HD22	2.01	0.42
1:A:350:THR:HB	1:A:353:GLN:HG3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:676:TRP:CZ2	1:A:680:VAL:HG21	2.55	0.42
1:A:455:LEU:HD12	1:A:587:TRP:CB	2.49	0.42
1:B:525:GLN:HG3	1:B:529:ASN:O	2.19	0.42
1:A:607:LEU:HD13	1:A:626:LYS:HG2	2.01	0.42
1:A:496:PRO:HA	1:A:499:VAL:HG23	2.02	0.42
1:B:595:VAL:O	1:B:599:CYS:HB2	2.20	0.42
1:B:445:HIS:C	1:B:445:HIS:CD2	2.92	0.42
1:B:537:PRO:HA	1:B:538:PRO:HD3	1.97	0.42
1:A:351:LYS:HE2	1:A:392:SER:CB	2.44	0.41
1:B:388:ILE:O	1:B:392:SER:HA	2.19	0.41
2:B:750:HEM:NC	4:B:800:ML6:H04	2.35	0.41
1:B:312:LEU:HB3	1:B:666:TYR:CD2	2.54	0.41
1:B:299:ARG:NH1	1:B:299:ARG:HB3	2.35	0.41
4:A:800:ML6:H23	1:B:306:TRP:CZ3	2.56	0.41
1:B:584:PHE:CD2	2:B:750:HEM:CAC	3.03	0.41
1:A:350:THR:HB	1:A:353:GLN:CG	2.50	0.41
1:A:503:GLU:O	1:A:507:GLN:HB3	2.21	0.41
1:A:531:PRO:HD2	1:A:716:TRP:CZ3	2.55	0.41
1:A:667:ARG:NH1	1:A:668:CYS:SG	2.94	0.41
1:B:355:PHE:CE1	1:B:385:ASN:CB	3.04	0.41
4:B:800:ML6:H26	4:B:800:ML6:H14	1.91	0.41
1:A:475:TRP:CE2	1:A:531:PRO:HG3	2.55	0.41
1:A:475:TRP:CZ2	1:A:531:PRO:HG3	2.56	0.41
1:B:485:TYR:CZ	1:B:514:ARG:HA	2.56	0.41
1:B:616:LEU:HA	1:B:616:LEU:HD23	1.86	0.41
1:A:567:VAL:HG21	4:A:800:ML6:C06	2.51	0.40
1:B:548:HIS:NE2	1:B:632:GLU:OE1	2.49	0.40
1:A:569:ASN:N	1:A:569:ASN:HD22	2.16	0.40
1:B:547:ARG:N	1:B:547:ARG:HD3	2.35	0.40
2:A:750:HEM:HBD2	2:A:750:HEM:CMD	2.52	0.40
1:A:425:GLN:HG2	1:A:448:TYR:CZ	2.56	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	406/422 (96%)	387 (95%)	18 (4%)	1 (0%)	52	43
1	B	408/422 (97%)	394 (97%)	14 (3%)	0	100	100
All	All	814/844 (96%)	781 (96%)	32 (4%)	1 (0%)	56	49

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	366/377 (97%)	342 (93%)	24 (7%)	21	11
1	B	367/377 (97%)	359 (98%)	8 (2%)	60	55
All	All	733/754 (97%)	701 (96%)	32 (4%)	35	26

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	319	LYS
1	A	321	THR
1	A	352	ASP
1	A	370	LYS
1	A	380	ARG
1	A	381	LEU
1	A	390	SER
1	A	469	LYS
1	A	487	GLN
1	A	489	ASP
1	A	507	GLN
1	A	514	ARG

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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	615	ASP
1	A	620	LYS
1	A	645	LYS
1	A	653	SER
1	A	663	GLU
1	A	713	THR
1	A	715	VAL
1	B	350	THR
1	B	353	GLN
1	B	392	SER
1	B	481	ARG
1	B	500	GLN
1	B	540	LEU
1	B	547	ARG
1	B	717	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	353	GLN
1	A	454	ASN
1	A	487	GLN
1	A	507	GLN
1	A	508	GLN
1	A	569	ASN
1	A	642	GLN
1	A	697	ASN
1	B	353	GLN
1	B	364	GLN
1	B	454	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$
2	HEM	A	750	1	30,50,50	2.88	13 (43%)	24,82,82	3.21	12 (50%)
3	H4B	A	760	-	13,18,18	1.21	2 (15%)	11,26,26	2.79	5 (45%)
4	ML6	A	800	-	25,25,25	1.68	6 (24%)	33,33,33	1.34	3 (9%)
5	ACT	A	860	-	1,3,3	2.94	1 (100%)	0,3,3	0.00	-
2	HEM	B	750	1	30,50,50	2.81	14 (46%)	24,82,82	3.02	10 (41%)
3	H4B	B	760	-	13,18,18	1.26	1 (7%)	11,26,26	2.26	3 (27%)
4	ML6	B	800	-	25,25,25	1.02	2 (8%)	33,33,33	1.73	6 (18%)
5	ACT	B	860	-	1,3,3	0.50	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	750	1	-	0/10/54/54	0/0/8/8

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	H4B	A	760	-	-	0/8/17/17	0/2/2/2
4	ML6	A	800	-	-	0/8/8/8	0/3/3/3
5	ACT	A	860	-	-	0/0/0/0	0/0/0/0
2	HEM	B	750	1	-	0/10/54/54	0/0/8/8
3	H4B	B	760	-	-	0/8/17/17	0/2/2/2
4	ML6	B	800	-	-	0/8/8/8	0/3/3/3
5	ACT	B	860	-	-	0/0/0/0	0/0/0/0

All (39) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	750	HEM	C3D-C4D	-6.19	1.43	1.51
2	B	750	HEM	C2D-C3D	-2.94	1.45	1.54
2	A	750	HEM	C3B-C4B	-2.79	1.49	1.51
2	B	750	HEM	C3D-C4D	-2.55	1.48	1.51
2	A	750	HEM	CAD-C3D	-2.42	1.49	1.54
2	A	750	HEM	C2D-C3D	-2.35	1.47	1.54
2	B	750	HEM	C2C-C1C	-2.24	1.48	1.52
2	B	750	HEM	C3B-C4B	-2.06	1.50	1.51
2	B	750	HEM	C4A-CHB	2.02	1.45	1.39
3	A	760	H4B	C4-N3	2.03	1.36	1.33
2	B	750	HEM	FE-NB	2.05	2.08	1.97
4	A	800	ML6	C24-C25	2.09	1.41	1.37
4	B	800	ML6	C04-C03	2.15	1.41	1.36
2	B	750	HEM	FE-ND	2.27	2.09	1.97
3	B	760	H4B	C7-N8	2.32	1.49	1.46
2	A	750	HEM	C4A-CHB	2.36	1.46	1.39
2	A	750	HEM	FE-NB	2.64	2.11	1.97
2	B	750	HEM	C1A-CHA	2.79	1.47	1.39
2	A	750	HEM	C2A-C3A	2.79	1.46	1.37
4	A	800	ML6	C06-C07	2.80	1.42	1.36
3	A	760	H4B	C2-N2	2.81	1.39	1.34
4	A	800	ML6	C10-N01	2.85	1.42	1.37
2	B	750	HEM	CHC-C4B	2.87	1.46	1.38
4	A	800	ML6	C04-C03	2.91	1.42	1.36
4	A	800	ML6	C09-C08	2.93	1.44	1.37
5	A	860	ACT	CH3-C	2.94	1.52	1.48
4	B	800	ML6	C02-N01	2.98	1.36	1.33
2	A	750	HEM	CHC-C4B	3.06	1.47	1.38
2	A	750	HEM	CHD-C1D	3.23	1.47	1.38
2	B	750	HEM	C2A-C3A	3.47	1.47	1.37
2	B	750	HEM	CHC-C1C	3.49	1.44	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	750	HEM	CHD-C4C	4.50	1.46	1.36
4	A	800	ML6	C02-N01	4.54	1.38	1.33
2	A	750	HEM	CHC-C1C	4.75	1.47	1.36
2	B	750	HEM	CHD-C4C	5.00	1.48	1.36
2	A	750	HEM	C1C-NC	6.39	1.43	1.36
2	A	750	HEM	C4C-NC	6.91	1.44	1.36
2	B	750	HEM	C4C-NC	7.09	1.44	1.36
2	B	750	HEM	C1C-NC	8.39	1.46	1.36

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	750	HEM	CBA-CAA-C2A	-8.00	98.20	112.53
2	B	750	HEM	CBA-CAA-C2A	-7.16	99.70	112.53
3	A	760	H4B	N3-C2-N1	-4.92	117.47	125.53
4	B	800	ML6	C11-C08-C09	-4.23	115.26	121.98
4	B	800	ML6	C03-C02-N01	-4.05	117.77	122.20
2	B	750	HEM	C3C-CAC-CBC	-3.28	119.43	124.46
2	A	750	HEM	CAD-CBD-CGD	-3.13	100.25	113.02
2	A	750	HEM	C3C-CAC-CBC	-3.06	119.76	124.46
2	A	750	HEM	C3B-CAB-CBB	-2.90	120.01	124.46
2	B	750	HEM	C3B-CAB-CBB	-2.82	120.13	124.46
3	B	760	H4B	N3-C2-N1	-2.65	121.19	125.53
4	B	800	ML6	C24-C25-C26	-2.43	120.21	123.35
2	A	750	HEM	CBD-CAD-C3D	-2.32	106.80	113.55
2	B	750	HEM	CBD-CAD-C3D	-2.29	106.89	113.55
2	A	750	HEM	C2D-C3D-C4D	2.14	105.12	101.50
4	A	800	ML6	C09-C10-N01	2.17	122.30	118.78
3	A	760	H4B	N2-C2-N3	2.20	120.85	117.20
2	A	750	HEM	C2C-C1C-NC	2.57	114.53	110.21
4	A	800	ML6	C02-N01-C10	2.65	120.94	117.78
4	B	800	ML6	N02-C02-N01	2.67	119.77	118.03
3	A	760	H4B	N2-C2-N1	2.70	121.67	117.20
2	B	750	HEM	CMD-C2D-C3D	2.93	127.29	114.35
2	A	750	HEM	CMD-C2D-C3D	2.98	127.51	114.35
2	A	750	HEM	CAD-C3D-C4D	3.03	123.16	112.47
3	A	760	H4B	C2-N1-C8A	3.05	121.39	114.54
2	B	750	HEM	C2C-C1C-NC	3.05	115.36	110.21
4	B	800	ML6	C08-C11-N12	3.23	120.89	112.88
3	B	760	H4B	C4-N3-C2	3.72	121.10	115.94
4	B	800	ML6	C02-N01-C10	3.86	122.39	117.78
4	A	800	ML6	C08-C11-N12	4.12	123.10	112.88

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	760	H4B	C4-C4A-C8A	4.37	118.52	114.56
2	B	750	HEM	CAD-C3D-C2D	4.67	126.65	113.22
2	B	750	HEM	CAD-C3D-C4D	5.05	130.28	112.47
2	B	750	HEM	CMC-C2C-C3C	5.16	129.41	116.53
3	A	760	H4B	C4-N3-C2	5.39	123.42	115.94
2	A	750	HEM	CMC-C2C-C3C	6.01	131.53	116.53
2	A	750	HEM	CMB-C2B-C3B	6.06	131.65	116.53
2	A	750	HEM	CAD-C3D-C2D	6.42	131.66	113.22
2	B	750	HEM	CMB-C2B-C3B	6.46	132.66	116.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 22 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	750	HEM	5	0
4	A	800	ML6	4	0
2	B	750	HEM	9	0
4	B	800	ML6	6	0
5	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, 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%)	0.56	60 (14%) 3 3	32, 58, 101, 140	0
1	B	411/422 (97%)	0.08	22 (5%) 29 32	30, 47, 78, 101	0
All	All	818/844 (96%)	0.32	82 (10%) 9 10	30, 52, 93, 140	0

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	716	TRP	6.6
1	A	715	VAL	6.4
1	B	348	VAL	6.3
1	B	718	GLY	6.0
1	A	488	PRO	5.9
1	B	300	PHE	5.7
1	B	619	ARG	5.0
1	A	486	LYS	4.8
1	A	355	PHE	4.7
1	A	352	ASP	4.7
1	A	713	THR	4.6
1	A	712	ASN	4.2
1	A	392	SER	3.9
1	A	487	GLN	3.8
1	B	350	THR	3.8
1	A	507	GLN	3.7
1	A	391	THR	3.7
1	A	506	ILE	3.7
1	A	351	LYS	3.5
1	A	714	HIS	3.3
1	A	491	SER	3.3
1	A	390	SER	3.3
1	A	480	ILE	3.3
1	A	677	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	679	ILE	3.2
1	A	490	GLY	3.2
1	B	310	VAL	3.2
1	A	300	PHE	3.2
1	A	489	ASP	3.1
1	A	619	ARG	3.1
1	A	388	ILE	3.0
1	A	350	THR	3.0
1	B	677	VAL	3.0
1	A	680	VAL	3.0
1	B	299	ARG	2.9
1	A	681	PRO	2.9
1	A	567	VAL	2.8
1	A	561	TRP	2.7
1	A	469	LYS	2.7
1	A	470	HIS	2.7
1	A	503	GLU	2.7
1	A	711	TRP	2.6
1	A	385	ASN	2.6
1	B	620	LYS	2.6
1	B	621	THR	2.6
1	A	593	ILE	2.6
1	A	389	GLU	2.5
1	B	676	TRP	2.5
1	A	466	THR	2.5
1	A	311	VAL	2.4
1	A	416	VAL	2.4
1	A	566	ALA	2.4
1	B	667	ARG	2.4
1	A	565	PRO	2.4
1	B	352	ASP	2.4
1	A	322	LEU	2.4
1	A	299	ARG	2.4
1	A	588	TYR	2.3
1	A	321	THR	2.3
1	B	591	THR	2.3
1	B	351	LYS	2.3
1	A	511	LYS	2.3
1	A	591	THR	2.2
1	A	667	ARG	2.2
1	A	685	GLY	2.2
1	B	392	SER	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	386	LYS	2.2
1	B	691	PHE	2.2
1	B	302	LYS	2.2
1	A	584	PHE	2.2
1	A	682	PRO	2.2
1	A	678	TRP	2.2
1	B	611	ALA	2.2
1	A	551	PHE	2.1
1	A	676	TRP	2.1
1	A	643	SER	2.1
1	A	493	LEU	2.1
1	A	564	LEU	2.1
1	B	353	GLN	2.1
1	B	678	TRP	2.0
1	B	680	VAL	2.0
1	A	508	GLN	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, 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
5	ACT	B	860	4/4	0.91	0.15	2.92	58,67,69,70	0
5	ACT	A	860	4/4	0.97	0.17	2.36	67,67,78,78	0
4	ML6	A	800	23/23	0.88	0.26	1.32	35,42,70,77	0
4	ML6	B	800	23/23	0.95	0.20	1.11	35,47,70,71	0
2	HEM	B	750	43/43	0.98	0.18	1.08	34,39,49,59	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	HEM	A	750	43/43	0.98	0.21	0.92	33,38,47,60	0
3	H4B	B	760	17/17	0.96	0.20	0.90	36,38,44,48	0
3	H4B	A	760	17/17	0.96	0.19	0.39	36,41,45,48	0
6	ZN	A	1717	1/1	0.99	0.10	0.16	44,44,44,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.