



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 03:25 AM GMT

PDB ID : 3E6T  
Title : Structure of murine INOS oxygenase domain with inhibitor AR-C118901  
Authors : Garcin, E.D.; Arvai, A.S.; Rosenfeld, R.J.; Kroeger, M.D.; Crane, B.R.; Andersson, G.; Andrews, G.; Hamley, P.J.; Mallinder, P.R.; Nicholls, D.J.; St-Gallay, S.A.; Tinker, A.C.; Gensmantel, N.P.; Mete, A.; Cheshire, D.R.; Connolly, S.; Stueh, D.J.; Aberg, A.; Wallace, A.V.; Tainer, J.A.; Getzoff, E.D.  
Deposited on : 2008-08-15  
Resolution : 2.50 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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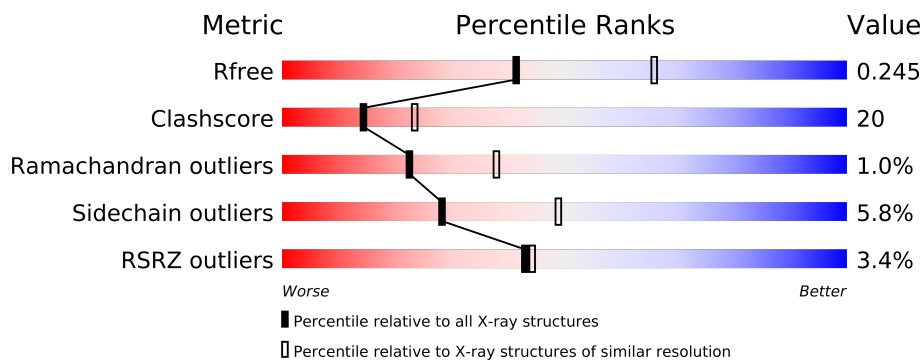
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 2.50 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))
$R_{free}$	66092	2784 (2.50-2.50)
Clashscore	79885	3562 (2.50-2.50)
Ramachandran outliers	78287	3480 (2.50-2.50)
Sidechain outliers	78261	3482 (2.50-2.50)
RSRZ outliers	66119	2785 (2.50-2.50)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	433	
1	B	433	

## 2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 7161 atoms, of which 0 are hydrogen and 0 are deuterium.

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, inducible.

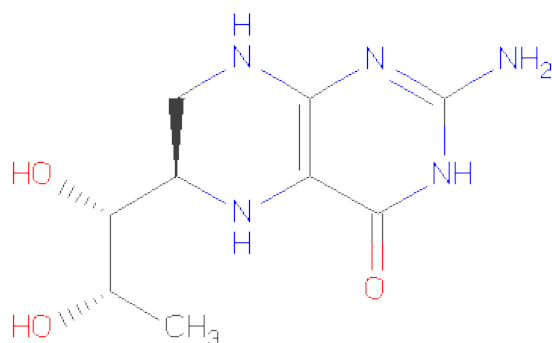
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	415	Total	C	N	O	S	0	0	0
			3385	2171	582	612	20			
1	B	410	Total	C	N	O	S	0	0	0
			3347	2148	577	602	20			

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



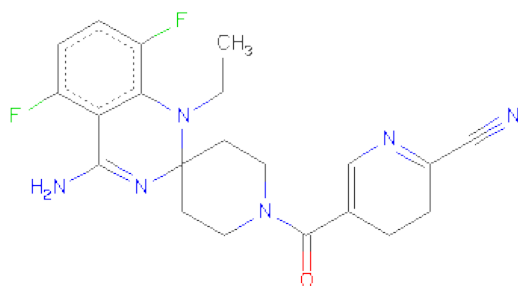
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- 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 5-(4'-AMINO-1'-ETHYL-5',8'-DIFLUORO-1'H-SPIRO[PIPERIDINE-4,2'-QUINAZOLINE]-1-YLCARBONYL)PICOLINONITRILE (three-letter code: 1A2) (formula: C<sub>21</sub>H<sub>22</sub>F<sub>2</sub>N<sub>6</sub>O).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	F	N	O	0	0
			30	21	2	6	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	1	Total	C	F	N	O	0	0
			30	21	2	6	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	112	Total	O	0	0
			112	112		
5	B	137	Total	O	0	0
			137	137		



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	213.99Å 213.99Å 116.82Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.98 – 2.50 29.86 – 2.50	Depositor EDS
% Data completeness (in resolution range)	88.7 (19.98-2.50) 88.8 (29.86-2.50)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.29 (at 2.51Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.236 , 0.267 0.227 , 0.245	Depositor DCC
$R_{free}$ test set	2461 reflections (5.34%)	DCC
Wilson B-factor (Å <sup>2</sup> )	45.3	Xtriage
Anisotropy	0.753	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 36.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 52558 reflections (0.002%)	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	7161	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.37% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 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, H4B, 1A2

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	0.38	0/3484	0.63	1/4737 (0.0%)
1	B	0.37	0/3445	0.63	0/4684
All	All	0.38	0/6929	0.63	1/9421 (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	365	GLY	N-CA-C	-5.28	99.89	113.10

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3385	0	3278	150	0
1	B	3347	0	3248	111	0
2	A	43	0	30	2	0
2	B	43	0	30	3	0
3	A	17	0	14	0	0
3	B	17	0	14	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	30	0	22	3	0
4	B	30	0	22	5	0
5	A	112	0	0	16	0
5	B	137	0	0	4	0
All	All	7161	0	6658	265	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 20.

All (265) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:195:ILE:HB	5:B:2039:HOH:O	1.32	1.30
1:A:153:ILE:HD12	1:A:153:ILE:H	1.22	1.00
2:A:901:HEM:HBA2	4:A:905:1A2:H82	1.42	1.00
1:A:99:THR:HG22	1:A:100:SER:H	1.31	0.93
1:B:301:LEU:HD13	1:B:315:ILE:HD11	1.51	0.90
2:B:901:HEM:HBA2	4:B:905:1A2:H82	1.56	0.87
1:B:252:ARG:HH21	1:B:489:PRO:HD3	1.44	0.81
1:B:141:ILE:HD11	1:B:163:VAL:HG21	1.66	0.78
1:A:99:THR:HG22	1:A:100:SER:N	1.98	0.77
1:A:134:LEU:O	1:A:138:ILE:HG12	1.83	0.77
1:A:251:PHE:O	1:A:252:ARG:HG2	1.84	0.77
1:A:221:ILE:HG21	1:A:301:LEU:HD21	1.66	0.77
1:A:195:ILE:HG21	1:A:368:MET:HE3	1.66	0.77
1:A:132:GLU:O	1:A:135:PRO:HD2	1.87	0.74
1:A:224:HIS:ND1	1:A:239:THR:HG22	2.03	0.72
1:A:239:THR:HG23	1:A:362:PRO:HG2	1.72	0.71
1:B:152:LYS:HD2	1:B:155:GLU:OE2	1.90	0.71
1:A:124:ARG:HD3	5:A:1018:HOH:O	1.90	0.71
1:A:336:LEU:HB3	1:A:338:LEU:HD22	1.73	0.70
1:A:141:ILE:CD1	1:A:163:VAL:HG21	2.22	0.70
1:A:153:ILE:HD12	1:A:153:ILE:N	2.02	0.69
1:A:217:MET:HE2	1:A:305:ALA:HB2	1.75	0.68
1:B:343:LEU:O	5:B:2073:HOH:O	2.12	0.68
1:B:195:ILE:CG2	1:B:437:PHE:HB2	2.24	0.68
1:B:285:ILE:HD11	1:B:291:PRO:HB3	1.76	0.67
1:A:405:LYS:O	1:A:409:VAL:HG23	1.95	0.66
1:B:186:MET:HE1	1:B:189:ARG:HH11	1.60	0.66
1:A:89:ILE:HD12	1:A:89:ILE:N	2.11	0.66
1:A:290:LYS:CD	1:A:290:LYS:H	2.08	0.65
1:B:417:LEU:O	1:B:421:GLN:HG3	1.95	0.65
1:B:215:GLN:O	1:B:219:GLN:HG3	1.97	0.64

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:249:HIS:C	1:B:306:ASP:O	2.36	0.64
1:A:290:LYS:CE	1:A:290:LYS:H	2.10	0.64
1:A:215:GLN:O	1:A:219:GLN:HG3	1.97	0.64
1:B:188:TRP:CE3	1:B:200:TRP:HA	2.33	0.64
1:B:221:ILE:HG21	1:B:301:LEU:HD21	1.80	0.63
1:A:195:ILE:HD13	1:A:368:MET:HE1	1.80	0.63
1:B:385:ILE:O	1:B:389:VAL:HG23	1.99	0.63
1:A:266:MET:SD	1:A:272:ARG:HD3	2.39	0.63
1:A:163:VAL:O	1:A:167:ILE:HG13	1.99	0.63
1:A:283:LEU:O	1:A:287:LEU:HG	1.99	0.63
1:A:159:ARG:O	1:A:163:VAL:HG23	1.98	0.63
1:A:176:THR:OG1	1:A:179:GLU:HG3	1.99	0.63
1:A:407:ARG:HD3	5:A:1115:HOH:O	1.97	0.62
2:B:901:HEM:HBA2	4:B:905:1A2:C8	2.28	0.62
1:B:177:LEU:O	1:B:181:ILE:HD13	1.99	0.62
1:B:78:TYR:CZ	1:B:91:HIS:HD2	2.18	0.61
1:A:217:MET:CE	1:A:303:LEU:HB3	2.29	0.61
1:A:467:THR:CG2	1:A:469:VAL:HG22	2.31	0.61
1:A:153:ILE:H	1:A:153:ILE:CD1	1.92	0.61
1:A:438:MET:CE	1:A:469:VAL:HG12	2.30	0.61
1:B:195:ILE:HG22	1:B:437:PHE:HB2	1.81	0.60
1:B:239:THR:O	1:B:361:CYS:HA	2.01	0.60
1:B:303:LEU:HD23	1:B:313:PHE:CD2	2.36	0.60
1:A:290:LYS:H	1:A:290:LYS:HD2	1.66	0.60
1:A:304:GLN:O	1:A:304:GLN:HG3	2.02	0.60
1:A:141:ILE:HD13	1:A:163:VAL:HG21	1.83	0.60
1:A:281:THR:O	1:A:285:ILE:HG12	2.02	0.60
1:B:301:LEU:HB3	1:B:303:LEU:HD21	1.84	0.59
1:B:387:GLU:HG3	1:B:397:THR:HG21	1.85	0.59
1:A:141:ILE:HD11	1:A:163:VAL:HG21	1.85	0.59
1:B:195:ILE:HG22	1:B:195:ILE:O	2.03	0.58
1:A:290:LYS:HE3	1:A:290:LYS:H	1.67	0.58
1:B:303:LEU:O	1:B:310:PRO:HA	2.03	0.58
1:B:141:ILE:CD1	1:B:163:VAL:HG21	2.34	0.58
1:A:272:ARG:HG2	1:A:272:ARG:HH11	1.68	0.58
1:A:445:TYR:HA	1:A:450:GLY:H	1.68	0.58
1:B:175:LEU:HD13	1:B:356:LEU:CD1	2.34	0.58
1:B:467:THR:CG2	1:B:469:VAL:HG22	2.34	0.58
1:B:303:LEU:HD22	1:B:303:LEU:N	2.19	0.57
1:A:411:GLU:O	1:A:414:VAL:HG22	2.03	0.57
1:A:410:THR:O	1:A:414:VAL:HG13	2.04	0.57
1:B:186:MET:HB3	1:B:481:PRO:HG2	1.86	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:241:PHE:HB3	1:B:242:PRO:CD	2.35	0.57
1:B:190:ASN:O	1:B:192:PRO:HD3	2.05	0.57
1:B:459:VAL:HG22	1:B:469:VAL:HG23	1.86	0.57
1:B:149:LYS:HG2	1:B:150:GLU:N	2.20	0.56
2:A:901:HEM:HBA2	4:A:905:1A2:C8	2.28	0.56
1:B:167:ILE:HG23	1:B:171:GLY:O	2.05	0.56
1:A:438:MET:HG3	1:A:468:PRO:HB2	1.86	0.56
1:B:77:GLN:HE21	1:B:77:GLN:HA	1.71	0.56
1:A:134:LEU:HB3	1:A:135:PRO:HD3	1.87	0.56
1:B:271:ILE:HD13	1:B:278:LEU:HD11	1.88	0.56
1:A:241:PHE:HB3	1:A:242:PRO:CD	2.35	0.56
1:A:333:PHE:HB3	5:A:1097:HOH:O	2.04	0.56
1:B:289:TRP:NE1	1:B:314:GLU:OE1	2.37	0.55
1:A:153:ILE:O	1:A:157:LEU:HD13	2.06	0.55
1:B:492:THR:OG1	1:B:492:THR:O	2.22	0.55
1:B:331:GLU:HA	1:B:331:GLU:OE1	2.06	0.55
1:B:149:LYS:HG2	1:B:150:GLU:HG3	1.89	0.54
1:B:134:LEU:HB3	1:B:135:PRO:HD3	1.89	0.54
1:A:350:LEU:HD23	1:A:351:LEU:N	2.22	0.54
1:A:217:MET:HE1	1:A:303:LEU:HB3	1.88	0.54
1:A:238:ILE:HG13	1:A:363:PHE:HB3	1.89	0.54
1:A:188:TRP:CE3	1:A:200:TRP:HA	2.43	0.54
1:B:301:LEU:CD1	1:B:315:ILE:HD11	2.33	0.54
1:A:138:ILE:HG22	1:A:142:ASN:ND2	2.22	0.54
1:A:466:ILE:O	1:A:466:ILE:HG22	2.08	0.54
1:B:133:LEU:C	1:B:133:LEU:HD13	2.28	0.53
1:B:186:MET:HE1	1:B:189:ARG:NH1	2.22	0.53
1:A:465:SER:O	1:A:471:HIS:HE1	1.90	0.53
1:A:264:TYR:HB2	1:A:266:MET:HE2	1.89	0.53
1:A:215:GLN:NE2	5:A:1058:HOH:O	2.36	0.52
1:B:129:PRO:HG2	1:B:132:GLU:HG2	1.91	0.52
1:B:175:LEU:HD13	1:B:356:LEU:HD12	1.91	0.52
1:B:195:ILE:HD11	1:B:458:LEU:O	2.09	0.52
1:A:249:HIS:C	1:A:306:ASP:O	2.48	0.52
1:A:274:ASP:OD2	5:A:1079:HOH:O	2.18	0.52
1:A:149:LYS:HG2	1:A:150:GLU:N	2.24	0.52
1:B:285:ILE:HD12	1:B:291:PRO:HD3	1.92	0.52
1:B:130:LEU:HD21	1:B:167:ILE:HG22	1.92	0.52
1:B:194:CYS:O	1:B:197:ARG:NH1	2.42	0.52
1:A:262:ALA:HB2	1:A:299:LEU:CD2	2.39	0.52
1:B:195:ILE:HG23	1:B:437:PHE:HB2	1.91	0.52
1:A:350:LEU:HD23	1:A:350:LEU:C	2.31	0.51

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:346:VAL:HG22	4:B:905:1A2:H133	1.91	0.51
1:B:243:GLN:HB3	1:B:358:PHE:CE2	2.44	0.51
1:B:80:ARG:NH2	1:B:89:ILE:HD13	2.26	0.51
1:B:258:LEU:HB3	1:B:259:ILE:HD12	1.92	0.51
1:A:80:ARG:NH1	1:A:89:ILE:HG21	2.26	0.51
1:B:259:ILE:HD12	1:B:259:ILE:N	2.26	0.51
1:B:209:ARG:O	1:B:242:PRO:HG3	2.11	0.50
1:A:328:PRO:O	1:A:329:LYS:HD2	2.11	0.50
1:A:488:GLU:OE1	1:A:491:LYS:HE2	2.11	0.50
1:B:77:GLN:NE2	1:B:77:GLN:HA	2.27	0.50
1:A:189:ARG:O	5:A:1036:HOH:O	2.18	0.50
1:A:480:SER:HA	1:A:481:PRO:C	2.32	0.50
1:A:186:MET:HE1	1:A:189:ARG:HH11	1.77	0.50
1:B:258:LEU:HB2	1:B:345:ALA:HB3	1.94	0.49
1:A:346:VAL:HG22	4:A:905:1A2:H133	1.94	0.49
1:B:283:LEU:O	1:B:287:LEU:HG	2.11	0.49
1:A:271:ILE:HD13	1:A:278:LEU:HD11	1.93	0.49
1:B:397:THR:HG22	1:B:397:THR:O	2.12	0.49
1:A:190:ASN:O	1:A:192:PRO:HD3	2.12	0.49
1:A:138:ILE:HG22	1:A:142:ASN:HD21	1.77	0.49
1:A:251:PHE:C	1:A:252:ARG:HG2	2.32	0.49
1:A:368:MET:HE1	1:A:433:ALA:HB1	1.94	0.49
1:B:186:MET:HE3	1:B:189:ARG:HD3	1.94	0.49
1:B:266:MET:HB3	1:B:267:PRO:HD2	1.93	0.49
1:B:438:MET:HE3	1:B:469:VAL:HA	1.95	0.49
1:A:252:ARG:HH21	1:A:489:PRO:HD3	1.76	0.49
1:A:239:THR:CG2	1:A:362:PRO:HG2	2.41	0.48
1:B:78:TYR:CD1	1:B:78:TYR:C	2.86	0.48
1:B:407:ARG:HG3	5:B:2109:HOH:O	2.11	0.48
1:B:327:HIS:CG	1:B:328:PRO:HD2	2.48	0.48
1:A:290:LYS:N	1:A:290:LYS:HD2	2.28	0.48
1:A:303:LEU:O	1:A:310:PRO:HA	2.13	0.48
1:A:254:TRP:CZ3	1:A:490:TRP:HB3	2.48	0.48
1:B:186:MET:CE	1:B:189:ARG:HD3	2.43	0.48
1:A:438:MET:HE3	1:A:469:VAL:HG12	1.96	0.48
1:A:149:LYS:HG2	1:A:150:GLU:HG3	1.95	0.48
1:A:442:GLN:HG3	1:A:443:ASN:N	2.29	0.48
1:B:213:THR:OG1	1:B:216:GLU:HG3	2.14	0.47
1:B:346:VAL:CG2	4:B:905:1A2:H133	2.44	0.47
1:B:445:TYR:HA	1:B:450:GLY:H	1.80	0.47
1:B:371:GLU:OE1	4:B:905:1A2:N6	2.48	0.47
1:A:301:LEU:HB3	1:A:303:LEU:CD1	2.44	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:129:PRO:HG3	5:A:1019:HOH:O	2.14	0.47
1:B:305:ALA:O	1:B:307:GLY:N	2.47	0.47
1:A:148:PHE:CE2	1:A:152:LYS:HE2	2.50	0.47
1:A:144:TYR:O	1:A:147:SER:HB3	2.15	0.47
1:A:264:TYR:HB2	1:A:266:MET:CE	2.45	0.47
1:A:186:MET:CE	1:A:189:ARG:HH11	2.28	0.47
1:B:195:ILE:HG23	1:B:437:PHE:CB	2.45	0.47
1:A:438:MET:HE2	1:A:469:VAL:HG12	1.94	0.47
1:B:124:ARG:HH21	1:B:128:THR:HB	1.79	0.47
1:A:417:LEU:O	1:A:421:GLN:HG3	2.16	0.46
1:B:445:TYR:O	1:B:449:GLY:HA2	2.14	0.46
1:B:453:ALA:HB1	1:B:458:LEU:CD1	2.45	0.46
1:A:301:LEU:HD13	1:A:315:ILE:HD11	1.97	0.46
1:B:129:PRO:HB2	1:B:131:GLU:CD	2.36	0.46
1:A:488:GLU:HG3	5:A:1122:HOH:O	2.14	0.46
1:A:246:ASP:CG	1:A:248:LYS:H	2.18	0.46
1:B:327:HIS:ND1	1:B:328:PRO:HD2	2.31	0.46
1:B:332:TRP:O	1:B:335:GLU:HB2	2.15	0.46
1:B:252:ARG:NH2	1:B:489:PRO:HD3	2.23	0.46
1:A:99:THR:CG2	1:A:100:SER:H	2.03	0.46
1:A:395:LEU:HD23	1:A:404:TRP:HB2	1.98	0.46
1:A:259:ILE:N	1:A:259:ILE:HD12	2.31	0.46
1:A:252:ARG:NH2	1:A:489:PRO:HD3	2.30	0.45
1:B:130:LEU:CD2	1:B:167:ILE:HG22	2.46	0.45
1:A:428:MET:HB3	5:A:1048:HOH:O	2.16	0.45
1:A:78:TYR:CD1	1:A:78:TYR:C	2.90	0.45
1:B:285:ILE:CD1	1:B:291:PRO:HB3	2.45	0.45
1:A:262:ALA:HB2	1:A:299:LEU:HG	1.99	0.45
1:A:332:TRP:O	1:A:335:GLU:HB2	2.16	0.45
1:A:368:MET:CE	1:A:433:ALA:HB1	2.46	0.45
1:B:266:MET:CE	1:B:272:ARG:HE	2.30	0.45
1:A:301:LEU:HB3	1:A:303:LEU:HD13	1.98	0.45
1:A:124:ARG:CD	5:A:1018:HOH:O	2.56	0.44
1:B:84:TRP:NE1	1:B:114:MET:HG3	2.32	0.44
1:A:252:ARG:HD3	1:A:359:PRO:HB2	1.98	0.44
1:A:217:MET:HE3	1:A:303:LEU:HB3	1.98	0.44
1:B:188:TRP:CZ3	1:B:200:TRP:HA	2.52	0.44
1:B:453:ALA:O	1:B:476:ASN:HB2	2.16	0.44
1:B:294:GLY:N	1:B:297:ASP:OD2	2.50	0.44
1:A:151:ALA:HB1	5:A:1022:HOH:O	2.17	0.44
1:A:141:ILE:HD11	1:A:163:VAL:HG11	1.99	0.44
1:A:266:MET:HB2	1:A:270:THR:O	2.17	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:350:LEU:HD23	1:B:350:LEU:C	2.38	0.44
1:B:375:ARG:O	1:B:379:ASP:HB2	2.18	0.44
1:A:241:PHE:HB3	1:A:242:PRO:HD2	2.00	0.44
1:B:438:MET:HE2	1:B:469:VAL:HG12	1.99	0.44
1:A:84:TRP:CE2	1:A:114:MET:HG3	2.53	0.44
1:A:371:GLU:O	1:A:375:ARG:HB2	2.18	0.44
1:A:133:LEU:C	1:A:133:LEU:HD23	2.39	0.44
1:B:441:MET:HE1	1:B:472:GLN:HG2	2.01	0.43
1:B:434:SER:HB3	1:B:468:PRO:HD2	2.00	0.43
1:A:334:GLN:HA	5:A:1096:HOH:O	2.17	0.43
1:A:434:SER:OG	1:A:468:PRO:HD2	2.19	0.43
1:A:333:PHE:O	1:A:336:LEU:HB2	2.18	0.43
1:A:172:THR:OG1	1:A:173:TYR:N	2.50	0.43
1:B:186:MET:HB3	1:B:481:PRO:CG	2.49	0.43
1:B:241:PHE:HB3	1:B:242:PRO:HD2	2.01	0.43
1:B:138:ILE:HG22	1:B:142:ASN:ND2	2.33	0.43
1:A:238:ILE:CG1	1:A:363:PHE:HB3	2.48	0.43
1:B:226:LEU:O	1:B:226:LEU:HD23	2.18	0.43
1:A:290:LYS:N	1:A:290:LYS:HE3	2.34	0.43
1:A:465:SER:C	1:A:467:THR:H	2.21	0.43
1:B:438:MET:CE	1:B:469:VAL:HG12	2.48	0.43
1:A:164:THR:O	1:A:168:GLU:HG2	2.19	0.43
1:A:445:TYR:O	1:A:449:GLY:HA2	2.19	0.43
1:A:397:THR:O	1:A:397:THR:HG22	2.18	0.43
1:B:194:CYS:HB2	2:B:901:HEM:ND	2.34	0.42
1:B:124:ARG:HD2	1:B:244:ARG:HD3	2.01	0.42
1:A:332:TRP:CE3	1:A:392:ARG:HD2	2.54	0.42
1:A:253:LEU:HD12	1:A:253:LEU:N	2.34	0.42
1:A:309:ASP:OD1	1:A:495:TRP:HA	2.19	0.42
1:A:223:ARG:HD2	5:A:1059:HOH:O	2.19	0.42
1:A:272:ARG:HG2	1:A:272:ARG:NH1	2.32	0.42
1:B:348:ASN:ND2	1:B:348:ASN:H	2.18	0.42
1:B:322:GLU:OE2	1:B:339:LYS:HE3	2.19	0.42
1:A:453:ALA:HB3	1:A:474:MET:HB2	2.01	0.42
1:A:195:ILE:HG21	1:A:368:MET:CE	2.43	0.42
1:B:330:TYR:HB3	1:B:332:TRP:CE2	2.55	0.42
1:A:84:TRP:NE1	1:A:114:MET:HG3	2.34	0.42
1:A:123:PRO:O	1:A:124:ARG:HG3	2.20	0.41
1:A:467:THR:HG21	1:A:469:VAL:HG22	2.01	0.41
1:B:351:LEU:HB3	1:B:358:PHE:HB2	2.02	0.41
1:B:210:ASN:N	1:B:210:ASN:HD22	2.17	0.41
1:A:306:ASP:HB3	1:A:307:GLY:H	1.72	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:149:LYS:HE2	1:A:150:GLU:HG3	2.03	0.41
1:B:348:ASN:H	1:B:348:ASN:HD22	1.68	0.41
1:B:78:TYR:HD1	1:B:79:VAL:N	2.19	0.41
1:A:434:SER:OG	1:A:467:THR:HG23	2.20	0.41
1:A:368:MET:HE1	1:A:433:ALA:CB	2.51	0.41
1:A:445:TYR:O	1:A:449:GLY:N	2.54	0.41
1:A:303:LEU:N	1:A:303:LEU:CD1	2.83	0.41
1:B:80:ARG:NH2	1:B:89:ILE:CD1	2.84	0.41
1:A:252:ARG:HH12	1:A:486:GLN:HB3	1.86	0.41
1:A:360:ALA:HA	5:A:1073:HOH:O	2.20	0.41
1:A:258:LEU:HB3	1:A:259:ILE:HD12	2.01	0.41
1:B:153:ILE:HG13	1:B:153:ILE:H	1.67	0.41
1:A:384:ASN:HA	5:A:1104:HOH:O	2.21	0.41
1:A:251:PHE:O	1:A:360:ALA:HB2	2.21	0.41
1:A:301:LEU:HD12	1:A:301:LEU:HA	1.89	0.40
1:B:195:ILE:HG13	5:B:2038:HOH:O	2.22	0.40
1:A:368:MET:CE	1:A:433:ALA:CB	2.99	0.40
1:A:89:ILE:CD1	1:A:89:ILE:N	2.81	0.40
1:B:264:TYR:CE1	1:B:293:TYR:HA	2.55	0.40
1:A:256:SER:HB2	1:A:257:GLN:OE1	2.22	0.40
1:A:242:PRO:HB2	1:A:251:PHE:CE1	2.56	0.40
1:A:249:HIS:C	5:A:1068:HOH:O	2.60	0.40
1:A:379:ASP:HB3	1:A:381:GLN:OE1	2.20	0.40
1:A:138:ILE:O	1:A:142:ASN:ND2	2.54	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	411/433 (95%)	370 (90%)	38 (9%)	3 (1%)	30	50
1	B	406/433 (94%)	362 (89%)	39 (10%)	5 (1%)	19	32
All	All	817/866 (94%)	732 (90%)	77 (9%)	8 (1%)	22	38



All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	100	SER
1	B	306	ASP
1	A	123	PRO
1	A	200	TRP
1	B	200	TRP
1	B	268	ASP
1	B	344	PRO
1	B	266	MET

### 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	363/381 (95%)	341 (94%)	22 (6%)	26	46
1	B	358/381 (94%)	338 (94%)	20 (6%)	30	51
All	All	721/762 (95%)	679 (94%)	42 (6%)	28	49

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	119	LEU
1	A	148	PHE
1	A	153	ILE
1	A	161	GLU
1	A	180	LEU
1	A	223	ARG
1	A	239	THR
1	A	252	ARG
1	A	257	GLN
1	A	258	LEU
1	A	290	LYS
1	A	301	LEU
1	A	303	LEU
1	A	336	LEU
1	A	338	LEU
1	A	348	ASN

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Mol	Chain	Res	Type
1	A	386	LEU
1	A	387	GLU
1	A	395	LEU
1	A	417	LEU
1	A	467	THR
1	A	475	LEU
1	B	78	TYR
1	B	119	LEU
1	B	128	THR
1	B	130	LEU
1	B	134	LEU
1	B	210	ASN
1	B	226	LEU
1	B	233	ASN
1	B	257	GLN
1	B	258	LEU
1	B	264	TYR
1	B	292	ARG
1	B	301	LEU
1	B	303	LEU
1	B	348	ASN
1	B	417	LEU
1	B	444	GLU
1	B	467	THR
1	B	486	GLN
1	B	492	THR

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	77	GLN
1	A	91	HIS
1	A	142	ASN
1	A	204	GLN
1	A	215	GLN
1	A	219	GLN
1	A	231	ASN
1	A	233	ASN
1	A	348	ASN
1	A	471	HIS
1	B	77	GLN
1	B	91	HIS

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Mol	Chain	Res	Type
1	B	210	ASN
1	B	231	ASN
1	B	233	ASN
1	B	348	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

6 ligands are modelled in this entry.

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	901	-	49,50,50	2.20	14 (28%)	46,82,82	1.41	6 (13%)
3	H4B	A	902	-	18,18,18	1.95	3 (16%)	24,26,26	2.18	6 (25%)
4	1A2	A	905	-	33,33,33	3.76	18 (54%)	49,49,49	2.34	14 (28%)
2	HEM	B	901	-	49,50,50	2.58	14 (28%)	46,82,82	1.30	5 (10%)
3	H4B	B	902	-	18,18,18	1.98	3 (16%)	24,26,26	2.20	6 (25%)
4	1A2	B	905	-	33,33,33	3.92	20 (60%)	49,49,49	2.38	17 (34%)

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	901	-	-	0/14/114/114	0/0/8/8
3	H4B	A	902	-	1/1/3/5	0/8/17/17	0/0/2/2
4	1A2	A	905	-	-	0/10/53/53	0/2/4/4
2	HEM	B	901	-	-	0/14/114/114	0/0/8/8
3	H4B	B	902	-	-	0/8/17/17	0/0/2/2
4	1A2	B	905	-	-	0/10/53/53	0/2/4/4

All (72) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	905	1A2	C25-N26	10.84	1.34	1.14
4	A	905	1A2	C25-N26	10.69	1.34	1.14
2	B	901	HEM	C3D-C4D	8.23	1.46	1.44
2	B	901	HEM	C2D-C1D	7.77	1.46	1.44
4	A	905	1A2	C23-C22	-7.54	1.39	1.49
4	B	905	1A2	C23-C22	-7.38	1.39	1.49
4	A	905	1A2	C22-N21	6.46	1.44	1.31
4	B	905	1A2	C15-C7	6.41	1.58	1.52
4	B	905	1A2	C22-N21	6.26	1.44	1.31
4	B	905	1A2	C5-N6	6.13	1.37	1.28
2	B	901	HEM	C3D-C2D	-6.09	1.33	1.43
4	A	905	1A2	C5-N6	5.93	1.37	1.28
2	A	901	HEM	C4A-C3A	5.92	1.47	1.40
3	B	902	H4B	C7-C6	-5.60	1.46	1.52
3	A	902	H4B	C7-C6	-5.60	1.46	1.52
4	B	905	1A2	C20-C19	5.57	1.45	1.35
4	B	905	1A2	C25-C22	-5.49	1.36	1.43
4	A	905	1A2	C11-C7	5.37	1.57	1.52
2	A	901	HEM	C3D-C2D	-5.32	1.34	1.43
4	A	905	1A2	C20-C19	5.24	1.45	1.35
3	B	902	H4B	C7-N8	-5.21	1.38	1.46
4	A	905	1A2	C15-C7	5.19	1.57	1.52
2	A	901	HEM	CHA-C4D	5.13	1.43	1.35
2	B	901	HEM	C3B-C2B	-5.07	1.34	1.43
4	A	905	1A2	C25-C22	-4.97	1.37	1.43
3	A	902	H4B	C7-N8	-4.80	1.38	1.46
2	A	901	HEM	C3B-C2B	-4.80	1.35	1.43
4	B	905	1A2	C11-C7	4.67	1.56	1.52
4	B	905	1A2	C8-N8	4.52	1.51	1.47
2	B	901	HEM	C3B-C4B	4.52	1.49	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	901	HEM	C2D-C1D	4.35	1.45	1.44
4	B	905	1A2	C16-N13	4.07	1.42	1.34
2	B	901	HEM	C3C-C2C	-3.89	1.37	1.43
2	A	901	HEM	C3B-C4B	3.70	1.48	1.44
4	A	905	1A2	C4A-C4	3.68	1.44	1.38
4	B	905	1A2	C24-C19	-3.68	1.43	1.50
2	B	901	HEM	CHA-C4D	3.61	1.41	1.35
4	A	905	1A2	C7-N8	3.59	1.53	1.48
4	B	905	1A2	C7-N8	3.46	1.53	1.48
4	A	905	1A2	C24-C23	-3.34	1.38	1.51
2	B	901	HEM	FE-NB	3.32	2.09	1.97
4	B	905	1A2	C20-N21	3.27	1.44	1.36
4	A	905	1A2	C24-C19	-3.20	1.44	1.50
4	A	905	1A2	C16-N13	3.13	1.40	1.34
2	B	901	HEM	CHB-C1B	3.12	1.40	1.35
2	A	901	HEM	C3C-C2C	-3.05	1.38	1.43
4	A	905	1A2	C16-C19	-2.96	1.45	1.50
2	A	901	HEM	FE-NA	2.86	2.04	1.92
4	B	905	1A2	C24-C23	-2.83	1.40	1.51
2	A	901	HEM	C1A-C2A	2.82	1.48	1.43
4	A	905	1A2	C8-N8	2.79	1.49	1.47
4	B	905	1A2	C16-C19	-2.65	1.45	1.50
4	B	905	1A2	C4A-C4	2.61	1.42	1.38
4	A	905	1A2	F10-C1	-2.61	1.29	1.35
2	B	901	HEM	CHD-C4C	2.60	1.41	1.36
2	A	901	HEM	FE-NC	2.55	2.07	1.97
2	A	901	HEM	C2B-C1B	-2.52	1.43	1.44
4	A	905	1A2	C20-N21	2.51	1.42	1.36
2	A	901	HEM	FE-NB	2.49	2.06	1.97
3	A	902	H4B	C4-N3	-2.47	1.33	1.37
4	B	905	1A2	O17-C16	2.45	1.28	1.23
2	B	901	HEM	FE-NA	2.33	2.02	1.92
4	B	905	1A2	F10-C1	-2.29	1.29	1.35
2	A	901	HEM	CHB-C1B	2.24	1.39	1.35
4	A	905	1A2	O17-C16	2.24	1.28	1.23
2	B	901	HEM	FE-ND	2.24	2.06	1.97
2	A	901	HEM	C1A-NA	2.23	1.40	1.36
3	B	902	H4B	C4-N3	-2.23	1.33	1.37
4	B	905	1A2	C4A-C5	-2.21	1.40	1.47
2	B	901	HEM	CMC-C2C	2.19	1.54	1.47
2	B	901	HEM	FE-NC	2.06	2.05	1.97
4	B	905	1A2	C14-N13	2.01	1.50	1.47

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	905	1A2	C14-C15-C7	-6.89	106.88	113.16
4	A	905	1A2	C14-C15-C7	-6.83	106.94	113.16
3	B	902	H4B	C7-C6-C9	6.38	123.50	113.66
3	A	902	H4B	C7-C6-C9	5.99	122.90	113.66
3	A	902	H4B	C6-C7-N8	5.73	119.06	111.66
3	B	902	H4B	C6-C7-N8	5.32	118.52	111.66
4	B	905	1A2	C7-N6-C5	5.23	125.86	116.59
2	B	901	HEM	C3B-C4B-NB	-5.16	110.31	114.00
4	B	905	1A2	C14-N13-C12	-5.10	102.79	112.52
2	A	901	HEM	C3B-C4B-NB	-5.06	110.38	114.00
4	B	905	1A2	C15-C14-N13	-5.00	103.82	110.94
4	A	905	1A2	C15-C14-N13	-4.92	103.94	110.94
4	A	905	1A2	C7-N6-C5	4.69	124.90	116.59
4	A	905	1A2	C14-N13-C12	-4.44	104.05	112.52
4	A	905	1A2	C19-C20-N21	-4.40	116.46	125.14
4	B	905	1A2	C15-C7-N8	4.38	116.53	110.48
4	A	905	1A2	C15-C7-N8	4.27	116.39	110.48
4	A	905	1A2	C12-C11-C7	-4.24	109.29	113.16
4	B	905	1A2	C4A-C8A-C1	-3.99	115.79	118.72
2	A	901	HEM	CBD-CAD-C3D	-3.92	105.82	114.37
4	B	905	1A2	C19-C20-N21	-3.90	117.44	125.14
4	A	905	1A2	C11-C12-N13	-3.71	105.66	110.94
4	A	905	1A2	C4A-C8A-C1	-3.51	116.14	118.72
2	B	901	HEM	CHC-C4B-NB	-3.23	121.90	124.58
4	B	905	1A2	C4A-C5-N6	-3.20	118.09	122.86
4	B	905	1A2	C12-C11-C7	-2.99	110.43	113.16
4	B	905	1A2	C11-C12-N13	-3.00	106.67	110.94
3	A	902	H4B	N8-C8A-N1	2.99	120.21	115.82
3	B	902	H4B	N8-C8A-N1	2.93	120.12	115.82
4	A	905	1A2	C4A-C5-N6	-2.90	118.54	122.86
3	A	902	H4B	C4-C4A-C8A	2.84	117.19	114.56
3	B	902	H4B	C4-C4A-C8A	2.81	117.17	114.56
4	B	905	1A2	N8-C7-N6	-2.76	102.54	110.19
4	B	905	1A2	C24-C23-C22	2.74	119.83	112.16
4	A	905	1A2	C23-C24-C19	2.57	121.79	114.41
4	B	905	1A2	C23-C24-C19	2.56	121.77	114.41
2	A	901	HEM	C4A-C3A-C2A	2.54	108.76	107.00
4	A	905	1A2	C24-C23-C22	2.49	119.13	112.16
3	A	902	H4B	C7-C6-N5	2.46	115.83	109.85
4	A	905	1A2	N8-C7-N6	-2.46	103.37	110.19
4	A	905	1A2	F9-C4-C4A	2.42	121.94	117.85
3	B	902	H4B	C7-C6-N5	2.40	115.67	109.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	905	1A2	C19-C16-N13	-2.37	117.91	119.83
4	B	905	1A2	C25-C22-N21	2.32	121.90	116.56
2	B	901	HEM	CAD-C3D-C4D	2.26	128.59	124.53
4	B	905	1A2	C23-C22-N21	-2.26	120.62	123.79
2	A	901	HEM	C1A-C2A-C3A	-2.23	104.61	106.92
2	B	901	HEM	C2D-C1D-ND	-2.23	110.30	112.93
3	A	902	H4B	C4A-C4-N3	2.22	119.76	114.06
3	B	902	H4B	C4A-C4-N3	2.20	119.70	114.06
2	A	901	HEM	C2D-C1D-ND	-2.12	110.43	112.93
2	A	901	HEM	CMA-C3A-C4A	-2.09	125.40	128.62
4	B	905	1A2	F9-C4-C4A	2.07	121.36	117.85
2	B	901	HEM	CAA-C2A-C3A	-2.06	123.11	129.00

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	902	H4B	C6

There are no torsion outliers.

There are no ring outliers.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	415/433 (95%)	0.12	16 (3%) 37 38	29, 51, 85, 123	0
1	B	410/433 (94%)	0.03	12 (2%) 49 51	30, 50, 77, 95	0
All	All	825/866 (95%)	0.08	28 (3%) 43 44	29, 50, 81, 123	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	102	PHE	10.1
1	A	103	THR	8.1
1	B	494	ILE	5.3
1	A	101	ASP	5.2
1	A	100	SER	4.0
1	A	495	TRP	3.9
1	B	110	LEU	3.7
1	A	108	SER	3.5
1	B	495	TRP	3.4
1	A	494	ILE	3.2
1	A	110	LEU	3.1
1	B	77	GLN	3.1
1	B	148	PHE	3.0
1	A	157	LEU	2.6
1	A	111	GLY	2.5
1	A	78	TYR	2.4
1	A	148	PHE	2.3
1	A	446	ARG	2.3
1	B	150	GLU	2.3
1	A	152	LYS	2.3
1	B	171	GLY	2.2
1	A	149	LYS	2.2
1	B	111	GLY	2.2
1	B	87	GLY	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	492	THR	2.1
1	B	267	PRO	2.1
1	B	492	THR	2.1
1	B	152	LYS	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 ⓘ

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	1A2	A	905	30/30	0.21	1.72	45,50,58,59	0
4	1A2	B	905	30/30	0.21	1.33	48,52,55,60	0
2	HEM	A	901	43/43	0.18	0.64	27,30,41,44	0
2	HEM	B	901	43/43	0.17	0.27	29,31,40,47	0
3	H4B	B	902	17/17	0.14	-0.34	40,41,46,47	0
3	H4B	A	902	17/17	0.11	-1.01	46,48,49,49	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.