



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:03 PM GMT

PDB ID : 1DWX  
Title : MURINE INDUCIBLE NITRIC OXIDE SYNTHASE OXYGENASE DIMER  
N-HYDROXYARGININE AND TETRAHYDROBIOPTERIN  
Authors : Crane, B.R.; Arvai, A.S.; Getzoff, E.D.; Stuehr, D.J.; Tainer, J.A.  
Deposited on : 1999-12-14  
Resolution : 2.60 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) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

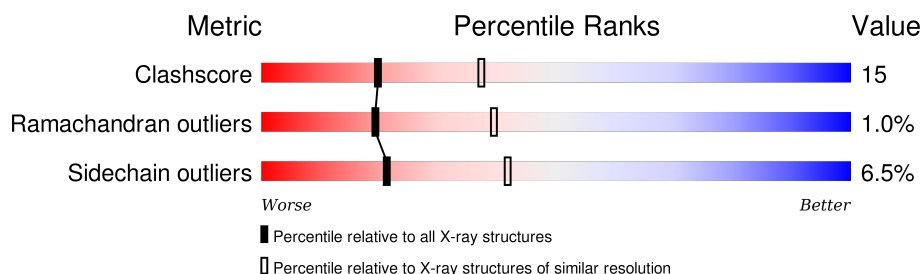
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.60 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))
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	420	
1	B	420	

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 7235 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.

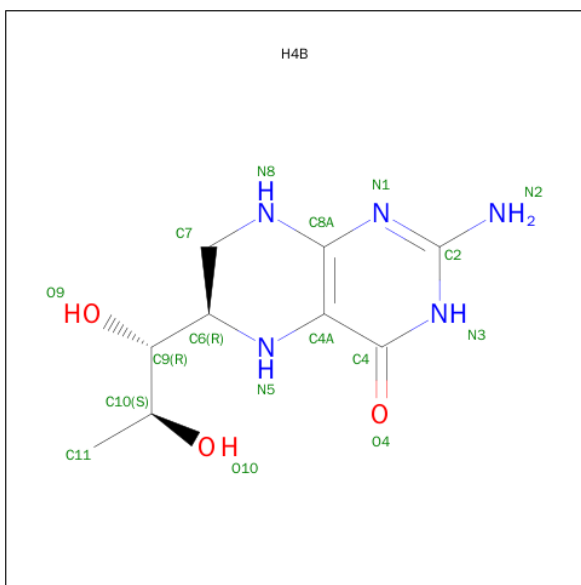
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	414	Total	C	N	O	S	0	0	0
			3376	2163	582	611	20			
1	B	411	Total	C	N	O	S	0	0	0
			3356	2153	579	604	20			

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



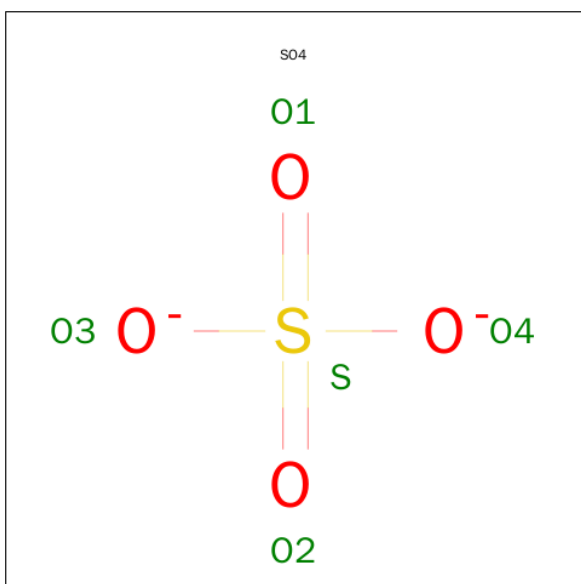
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- 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		

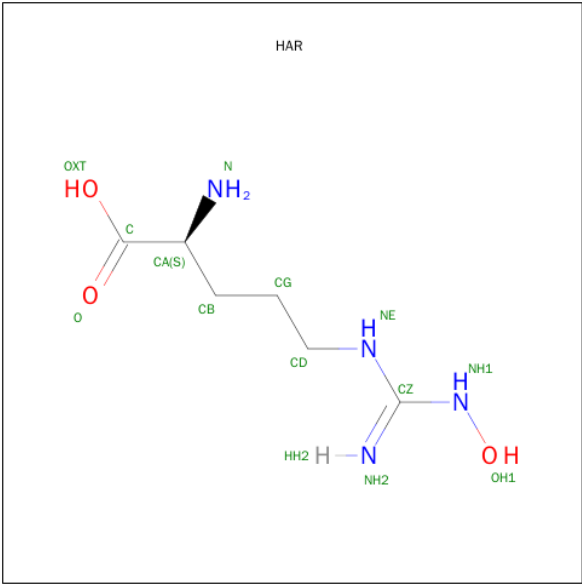
- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is N-OMEGA-HYDROXY-L-ARGININE (three-letter code: HAR) (formula:

C<sub>6</sub>H<sub>14</sub>N<sub>4</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			13	6	4	3		
5	B	1	Total	C	N	O	0	0
			13	6	4	3		

- Molecule 6 is water.

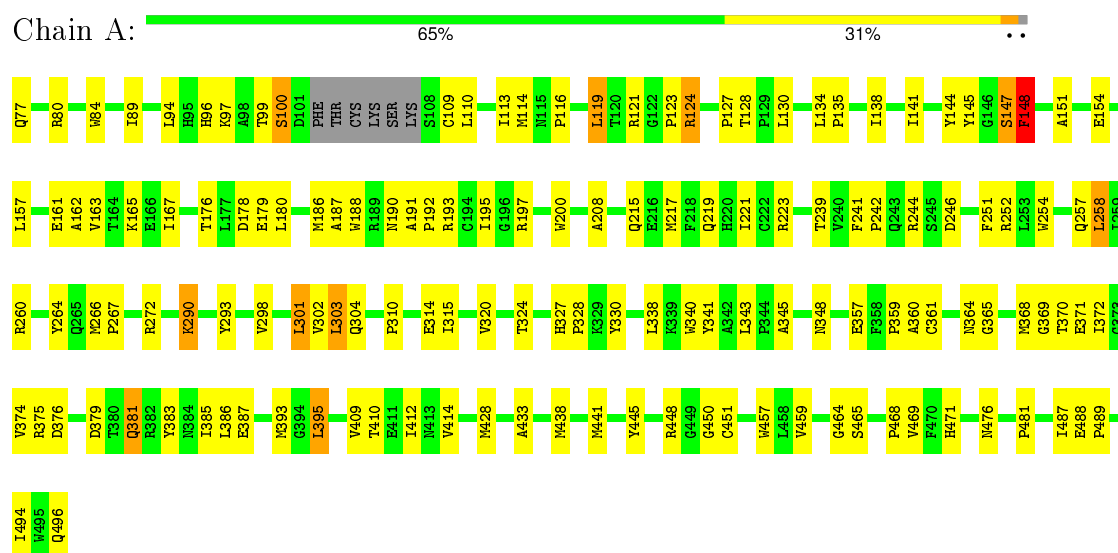
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	164	Total	O	0	0
			164	164		
6	B	188	Total	O	0	0
			188	188		

### 3 Residue-property plots

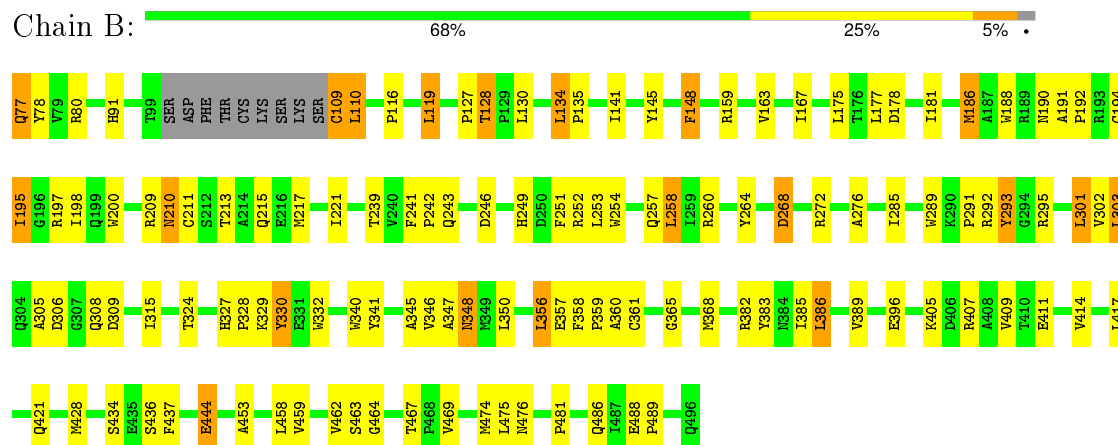
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 ( $\text{RSRZ} > 2$ ). 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.

Note EDS was not executed.

#### • Molecule 1: NITRIC OXIDE SYNTHASE



#### • Molecule 1: NITRIC OXIDE SYNTHASE



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	212.98Å 212.98Å 114.20Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.60	Depositor
% Data completeness (in resolution range)	97.9 (30.00-2.60)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.06	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, $R_{free}$	0.238 , 0.278	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7235	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	51.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, HAR, H4B, SO4

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	0.35	0/3474	0.63	1/4723 (0.0%)
1	B	0.36	0/3454	0.65	2/4696 (0.0%)
All	All	0.36	0/6928	0.64	3/9419 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	309	ASP	CB-CG-OD1	7.18	124.77	118.30
1	A	365	GLY	N-CA-C	-6.14	97.75	113.10
1	B	365	GLY	N-CA-C	-5.66	98.96	113.10

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	3376	0	3269	102	1
1	B	3356	0	3255	100	1
2	A	43	0	30	1	0
2	B	43	0	30	0	0
3	A	17	0	14	1	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	5	0	0	0	0
5	A	13	0	12	0	0
5	B	13	0	12	0	0
6	A	164	0	0	14	0
6	B	188	0	0	20	0
All	All	7235	0	6636	202	2

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 15.

All (202) 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:266:MET:SD	1:A:272:ARG:HD3	2.10	0.92
1:A:221:ILE:HG21	1:A:301:LEU:HD21	1.56	0.87
1:A:99:THR:HG22	1:A:100:SER:H	1.40	0.86
1:A:381:GLN:HE21	1:A:381:GLN:H	1.17	0.86
1:A:195:ILE:HG21	1:A:368:MET:HE1	1.62	0.82
1:B:77:GLN:HE21	1:B:77:GLN:HA	1.46	0.81
1:B:175:LEU:HD21	1:B:356:LEU:HD11	1.66	0.78
1:B:407:ARG:HA	6:B:2143:HOH:O	1.84	0.77
1:B:301:LEU:HB3	1:B:303:LEU:HD11	1.66	0.76
1:B:467:THR:CG2	1:B:469:VAL:HG22	2.18	0.74
1:A:381:GLN:NE2	1:A:381:GLN:H	1.86	0.73
1:B:467:THR:HG22	1:B:469:VAL:HG22	1.71	0.71
1:A:290:LYS:H	1:A:290:LYS:HD3	1.55	0.70
1:A:290:LYS:N	1:A:290:LYS:HD3	2.11	0.66
1:A:134:LEU:O	1:A:138:ILE:HG12	1.96	0.66
1:B:221:ILE:HD12	1:B:303:LEU:HD21	1.77	0.65
1:A:393:MET:HB2	1:A:395:LEU:HD22	1.79	0.65
1:A:298:VAL:HG21	1:A:320:VAL:HG11	1.80	0.64
1:A:144:TYR:O	1:A:147:SER:HB2	1.98	0.64
1:A:188:TRP:CZ3	1:A:200:TRP:HA	2.32	0.64
1:B:215:GLN:HG2	6:B:2066:HOH:O	1.98	0.63
1:A:223:ARG:HD3	6:A:2053:HOH:O	1.98	0.63
1:B:464:GLY:O	1:B:467:THR:HB	1.98	0.63
1:A:387:GLU:HB3	6:A:2048:HOH:O	1.99	0.62
1:A:410:THR:O	1:A:414:VAL:HG23	2.00	0.62
1:B:272:ARG:NH1	1:B:295:ARG:HG3	2.15	0.62
1:B:330:TYR:HD2	1:B:332:TRP:CZ2	2.19	0.61
1:B:258:LEU:HB2	1:B:345:ALA:HB3	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:901:HEM:HHC	2:A:901:HEM:HBB2	1.82	0.60
1:B:251:PHE:O	1:B:360:ALA:HB2	2.01	0.60
1:A:188:TRP:CE3	1:A:200:TRP:HA	2.37	0.60
1:B:459:VAL:HG22	1:B:469:VAL:HG23	1.84	0.60
1:B:243:GLN:HB3	1:B:358:PHE:CE2	2.37	0.59
1:A:360:ALA:HA	6:A:2100:HOH:O	2.02	0.59
1:A:141:ILE:HD11	1:A:163:VAL:HG21	1.85	0.59
1:A:84:TRP:NE1	1:A:114:MET:HG3	2.18	0.59
1:B:386:LEU:HB2	6:B:2135:HOH:O	2.03	0.58
1:B:252:ARG:HD3	1:B:359:PRO:HB2	1.85	0.58
1:B:141:ILE:CD1	1:B:163:VAL:HG21	2.34	0.58
1:A:459:VAL:HG22	1:A:469:VAL:HG23	1.85	0.58
1:B:209:ARG:O	1:B:242:PRO:HG3	2.03	0.58
1:B:329:LYS:HB3	6:B:2111:HOH:O	2.04	0.58
1:B:239:THR:O	1:B:361:CYS:HA	2.03	0.57
1:B:257:GLN:HB2	1:B:345:ALA:O	2.04	0.57
1:B:221:ILE:HG21	1:B:301:LEU:HD21	1.86	0.57
1:A:386:LEU:HB2	6:A:2117:HOH:O	2.03	0.56
1:B:217:MET:HB3	1:B:303:LEU:HD23	1.87	0.56
1:A:191:ALA:O	1:A:197:ARG:NH1	2.39	0.56
1:B:127:PRO:HG3	1:B:246:ASP:HA	1.86	0.56
1:A:195:ILE:HG21	1:A:368:MET:CE	2.33	0.56
1:B:436:SER:HB2	6:B:2158:HOH:O	2.04	0.55
1:B:195:ILE:HD13	1:B:437:PHE:HB2	1.87	0.55
1:A:116:PRO:HG2	1:A:119:LEU:HB2	1.88	0.55
1:A:252:ARG:HB2	1:A:304:GLN:HG2	1.87	0.55
1:A:80:ARG:NH2	6:A:2002:HOH:O	2.39	0.54
1:A:110:LEU:HB3	6:A:2010:HOH:O	2.07	0.54
1:A:258:LEU:HB2	1:A:345:ALA:HB3	1.90	0.54
1:B:301:LEU:HD13	1:B:315:ILE:HD11	1.90	0.54
1:B:80:ARG:NH2	6:B:2003:HOH:O	2.40	0.54
1:B:368:MET:HA	1:B:428:MET:O	2.09	0.53
1:A:99:THR:HG22	1:A:100:SER:N	2.17	0.53
1:B:303:LEU:HD12	1:B:303:LEU:N	2.23	0.53
1:A:340:TRP:HZ3	1:A:383:TYR:CE1	2.27	0.53
1:A:178:ASP:HB3	6:A:2020:HOH:O	2.09	0.53
1:A:304:GLN:O	1:A:304:GLN:HG3	2.07	0.52
1:A:197:ARG:NH2	1:A:451:CYS:SG	2.82	0.52
1:B:141:ILE:HD13	1:B:163:VAL:HG21	1.91	0.52
1:B:188:TRP:CE3	1:B:200:TRP:HA	2.45	0.52
1:B:175:LEU:HD21	1:B:356:LEU:CD1	2.38	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:385:ILE:O	1:B:389:VAL:HG23	2.10	0.52
1:B:285:ILE:HD11	1:B:291:PRO:HB3	1.92	0.52
1:A:163:VAL:O	1:A:167:ILE:HG13	2.10	0.51
1:A:266:MET:HB3	1:A:267:PRO:HD2	1.93	0.51
1:B:348:ASN:HB2	6:B:2120:HOH:O	2.11	0.51
1:A:290:LYS:CD	1:A:290:LYS:H	2.17	0.51
1:B:253:LEU:HD22	1:B:347:ALA:HB1	1.93	0.51
1:A:186:MET:HE1	1:A:448:ARG:NH1	2.26	0.51
1:A:123:PRO:O	1:A:124:ARG:HG3	2.11	0.50
1:A:303:LEU:O	1:A:310:PRO:HA	2.12	0.50
1:B:396:GLU:HB3	6:B:2140:HOH:O	2.11	0.50
1:B:405:LYS:O	1:B:409:VAL:HG12	2.12	0.50
1:A:176:THR:OG1	1:A:179:GLU:HG3	2.12	0.50
1:B:346:VAL:HG23	6:B:2117:HOH:O	2.12	0.50
1:A:371:GLU:O	1:A:375:ARG:HB2	2.11	0.50
1:B:258:LEU:HD13	1:B:347:ALA:HB2	1.94	0.50
1:A:251:PHE:O	1:A:360:ALA:HB2	2.12	0.49
1:B:188:TRP:O	1:B:197:ARG:HD2	2.12	0.49
1:A:387:GLU:HB2	6:A:2118:HOH:O	2.11	0.49
1:A:141:ILE:CD1	1:A:163:VAL:HG21	2.42	0.49
1:B:159:ARG:O	1:B:163:VAL:HG23	2.12	0.49
1:B:434:SER:CB	1:B:467:THR:HG23	2.42	0.49
1:B:195:ILE:HD13	1:B:437:PHE:CB	2.43	0.49
1:A:145:TYR:HA	1:A:148:PHE:CZ	2.46	0.49
1:B:411:GLU:HA	1:B:414:VAL:HG22	1.95	0.49
1:A:123:PRO:HD3	1:A:487:ILE:HD12	1.95	0.49
1:B:78:TYR:CE1	1:B:91:HIS:HD2	2.31	0.49
1:B:327:HIS:ND1	1:B:328:PRO:HD2	2.28	0.48
1:A:77:GLN:HG2	6:A:2001:HOH:O	2.13	0.48
1:B:260:ARG:HD2	1:B:341:TYR:OH	2.14	0.48
1:B:289:TRP:O	1:B:291:PRO:HD3	2.12	0.48
1:B:330:TYR:N	1:B:330:TYR:CD1	2.81	0.48
1:B:330:TYR:HD1	1:B:330:TYR:N	2.12	0.47
1:B:444:GLU:HG3	6:B:2164:HOH:O	2.12	0.47
1:B:210:ASN:C	1:B:210:ASN:HD22	2.18	0.47
1:B:210:ASN:HD22	1:B:211:CYS:N	2.12	0.47
1:A:465:SER:O	1:A:471:HIS:HE1	1.98	0.47
1:A:386:LEU:HD21	1:A:409:VAL:HG23	1.96	0.47
1:B:453:ALA:HB3	1:B:474:MET:HB2	1.97	0.47
1:A:379:ASP:HB3	1:A:381:GLN:NE2	2.31	0.46
1:A:193:ARG:HD2	6:A:2158:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:LEU:HD23	1:A:97:LYS:HD2	1.98	0.46
1:A:80:ARG:NH2	1:A:89:ILE:HD12	2.29	0.46
1:A:154:GLU:H	1:A:154:GLU:CD	2.19	0.46
1:A:162:ALA:HB2	6:A:2024:HOH:O	2.14	0.46
1:A:438:MET:HG3	1:A:468:PRO:HB2	1.98	0.46
1:A:127:PRO:HG3	1:A:246:ASP:HA	1.96	0.46
1:A:239:THR:O	1:A:361:CYS:HA	2.15	0.46
1:B:361:CYS:SG	1:B:361:CYS:O	2.74	0.45
1:B:110:LEU:N	1:B:110:LEU:HD22	2.31	0.45
1:A:385:ILE:HD11	1:A:412:ILE:HD13	1.97	0.45
1:A:258:LEU:HD22	1:A:345:ALA:HB1	1.98	0.45
1:B:128:THR:HG22	6:B:2062:HOH:O	2.16	0.45
1:B:276:ALA:HB2	1:B:382:ARG:CD	2.46	0.45
1:B:163:VAL:O	1:B:167:ILE:HG13	2.17	0.45
1:B:198:ILE:CG2	6:B:2158:HOH:O	2.65	0.45
1:B:444:GLU:HA	1:B:444:GLU:OE1	2.16	0.45
1:A:488:GLU:HG2	6:A:2151:HOH:O	2.16	0.45
1:B:141:ILE:HD11	1:B:163:VAL:HG21	1.98	0.45
1:A:438:MET:CE	1:A:469:VAL:HG12	2.47	0.45
1:A:488:GLU:HA	1:A:489:PRO:HD2	1.82	0.45
1:A:165:LYS:HB2	1:A:165:LYS:HE3	1.84	0.45
1:B:213:THR:HB	6:B:2065:HOH:O	2.16	0.45
1:A:217:MET:HG2	1:A:241:PHE:CE1	2.52	0.44
1:A:77:GLN:O	1:A:96:HIS:HE1	2.00	0.44
1:A:301:LEU:HD13	1:A:315:ILE:HD11	1.99	0.44
1:B:134:LEU:HB3	1:B:135:PRO:HD3	2.00	0.44
1:B:186:MET:HB3	1:B:481:PRO:HG2	1.98	0.44
1:B:190:ASN:O	1:B:192:PRO:HD3	2.17	0.44
1:B:295:ARG:HG2	6:B:2099:HOH:O	2.15	0.44
1:B:350:LEU:HD21	1:B:357:GLU:HB2	2.00	0.44
1:A:121:ARG:HD3	1:A:121:ARG:HA	1.83	0.44
1:B:194:CYS:O	1:B:197:ARG:HG3	2.18	0.43
1:A:190:ASN:O	1:A:192:PRO:HD3	2.19	0.43
1:B:252:ARG:NH2	1:B:489:PRO:HD3	2.34	0.43
1:A:494:ILE:N	1:A:494:ILE:HD12	2.32	0.43
1:A:208:ALA:O	1:A:242:PRO:HD3	2.19	0.43
1:B:191:ALA:O	1:B:197:ARG:HD3	2.18	0.43
1:B:145:TYR:HE2	1:B:159:ARG:HG2	1.84	0.43
1:B:78:TYR:CE1	1:B:91:HIS:CD2	3.06	0.43
1:A:187:ALA:HB2	1:A:481:PRO:HB2	2.01	0.43
1:A:257:GLN:HB2	1:A:345:ALA:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:486:GLN:HG3	6:B:2120:HOH:O	2.18	0.43
1:B:148:PHE:HB3	6:B:2029:HOH:O	2.18	0.43
1:A:301:LEU:HG	1:A:303:LEU:HD11	2.01	0.43
1:A:241:PHE:HB3	1:A:242:PRO:HD2	2.01	0.42
1:B:116:PRO:HG2	1:B:119:LEU:HB2	2.00	0.42
1:A:372:ILE:HA	1:A:376:ASP:HB2	2.01	0.42
1:B:305:ALA:O	1:B:308:GLN:HG2	2.19	0.42
1:B:254:TRP:HB2	1:B:302:VAL:HB	2.01	0.42
1:A:221:ILE:CG2	1:A:301:LEU:HD21	2.40	0.42
1:B:434:SER:HB3	1:B:467:THR:HG23	2.02	0.42
1:A:445:TYR:HA	1:A:450:GLY:H	1.84	0.42
1:B:467:THR:HG21	1:B:469:VAL:HG22	1.96	0.42
1:A:267:PRO:HB3	1:B:293:TYR:CD2	2.54	0.41
1:B:127:PRO:HA	6:B:2086:HOH:O	2.19	0.41
1:B:453:ALA:HB1	1:B:458:LEU:CD1	2.50	0.41
1:A:260:ARG:HD2	1:A:341:TYR:OH	2.20	0.41
1:A:244:ARG:HD2	1:A:357:GLU:OE2	2.20	0.41
1:A:241:PHE:HB3	1:A:242:PRO:CD	2.51	0.41
1:A:330:TYR:CD1	1:A:330:TYR:N	2.88	0.41
1:B:417:LEU:O	1:B:421:GLN:HB2	2.20	0.41
1:A:110:LEU:HD12	1:A:113:ILE:HG13	2.01	0.41
1:A:254:TRP:HB2	1:A:302:VAL:HB	2.02	0.41
1:B:141:ILE:HD11	1:B:163:VAL:HG11	2.02	0.41
1:A:457:TRP:HA	3:A:902:H4B:N1	2.35	0.41
1:B:488:GLU:HA	1:B:489:PRO:HD2	1.88	0.41
1:B:241:PHE:HB3	1:B:242:PRO:CD	2.51	0.41
1:B:437:PHE:CZ	1:B:458:LEU:HD13	2.54	0.41
1:A:465:SER:O	1:A:471:HIS:CE1	2.73	0.41
1:B:340:TRP:HZ3	1:B:383:TYR:CE1	2.39	0.41
1:B:177:LEU:O	1:B:181:ILE:HD13	2.20	0.41
1:A:121:ARG:HD2	6:A:2104:HOH:O	2.20	0.41
1:B:198:ILE:HG21	6:B:2158:HOH:O	2.20	0.41
1:A:314:GLU:HG3	6:A:2083:HOH:O	2.20	0.41
1:A:217:MET:HE1	1:A:303:LEU:HB3	2.01	0.41
1:A:252:ARG:HD3	1:A:359:PRO:HB2	2.01	0.41
1:A:215:GLN:HG3	1:A:219:GLN:NE2	2.36	0.41
1:A:134:LEU:HB3	1:A:135:PRO:HD3	2.03	0.41
1:A:124:ARG:NH1	1:A:128:THR:OG1	2.50	0.41
1:A:192:PRO:HB2	1:A:476:ASN:ND2	2.36	0.41
1:A:368:MET:HA	1:A:428:MET:O	2.22	0.40
1:B:241:PHE:HB3	1:B:242:PRO:HD2	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:343:LEU:HD11	1:A:364:ASN:ND2	2.36	0.40
1:B:453:ALA:HB1	1:B:458:LEU:HD12	2.03	0.40
1:A:368:MET:HE2	1:A:433:ALA:HB3	2.04	0.40
1:A:188:TRP:O	1:A:197:ARG:HD2	2.21	0.40
1:B:382:ARG:NH2	6:B:2133:HOH:O	2.54	0.40
1:A:327:HIS:ND1	1:A:328:PRO:HD2	2.37	0.40
1:B:78:TYR:CD1	1:B:78:TYR:C	2.95	0.40
1:B:249:HIS:HB3	1:B:306:ASP:OD1	2.20	0.40
1:A:370:THR:HA	1:A:374:VAL:HG23	2.03	0.40
1:B:463:SER:N	6:B:2172:HOH:O	2.48	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109:CYS:SG	1:B:109:CYS:SG[9_766]	1.72	0.48
1:A:109:CYS:SG	1:A:109:CYS:SG[11_655]	1.80	0.40

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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	410/420 (98%)	359 (88%)	45 (11%)	6 (2%)	13	26
1	B	407/420 (97%)	371 (91%)	34 (8%)	2 (0%)	34	60
All	All	817/840 (97%)	730 (89%)	79 (10%)	8 (1%)	19	39

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	100	SER
1	A	148	PHE

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Mol	Chain	Res	Type
1	A	151	ALA
1	B	293	TYR
1	A	293	TYR
1	B	268	ASP
1	A	369	GLY
1	A	464	GLY

### 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	362/368 (98%)	342 (94%)	20 (6%)	27	51
1	B	359/368 (98%)	332 (92%)	27 (8%)	17	33
All	All	721/736 (98%)	674 (94%)	47 (6%)	21	42

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	119	LEU
1	A	124	ARG
1	A	130	LEU
1	A	147	SER
1	A	148	PHE
1	A	157	LEU
1	A	161	GLU
1	A	180	LEU
1	A	258	LEU
1	A	264	TYR
1	A	290	LYS
1	A	301	LEU
1	A	303	LEU
1	A	324	THR
1	A	338	LEU
1	A	348	ASN
1	A	381	GLN

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Mol	Chain	Res	Type
1	A	395	LEU
1	A	441	MET
1	A	496	GLN
1	B	77	GLN
1	B	109	CYS
1	B	110	LEU
1	B	119	LEU
1	B	128	THR
1	B	130	LEU
1	B	134	LEU
1	B	148	PHE
1	B	178	ASP
1	B	186	MET
1	B	195	ILE
1	B	210	ASN
1	B	258	LEU
1	B	264	TYR
1	B	268	ASP
1	B	292	ARG
1	B	301	LEU
1	B	303	LEU
1	B	324	THR
1	B	330	TYR
1	B	348	ASN
1	B	356	LEU
1	B	386	LEU
1	B	444	GLU
1	B	462	VAL
1	B	475	LEU
1	B	476	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	HIS
1	A	96	HIS
1	A	219	GLN
1	A	231	ASN
1	A	308	GLN
1	A	348	ASN
1	A	364	ASN
1	A	381	GLN

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Mol	Chain	Res	Type
1	A	442	GLN
1	A	471	HIS
1	A	472	GLN
1	A	476	ASN
1	B	77	GLN
1	B	91	HIS
1	B	210	ASN
1	B	231	ASN
1	B	348	ASN
1	B	442	GLN
1	B	476	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](#)

7 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	1	30,50,50	2.69	9 (30%)	24,82,82	2.47	9 (37%)
3	H4B	A	902	-	13,18,18	1.66	1 (7%)	11,26,26	2.16	5 (45%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	SO4	A	903	-	4,4,4	0.68	0	6,6,6	0.18	0
5	HAR	A	906	-	5,12,12	0.36	0	3,14,14	0.27	0
2	HEM	B	901	1	30,50,50	2.74	9 (30%)	24,82,82	2.43	8 (33%)
3	H4B	B	902	-	13,18,18	1.71	2 (15%)	11,26,26	2.03	5 (45%)
5	HAR	B	906	-	5,12,12	0.39	0	3,14,14	0.44	0

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	1	-	0/10/54/54	0/0/8/8
3	H4B	A	902	-	-	0/8/17/17	0/2/2/2
4	SO4	A	903	-	-	0/0/0/0	0/0/0/0
5	HAR	A	906	-	-	0/5/13/13	0/0/0/0
2	HEM	B	901	1	-	0/10/54/54	0/0/8/8
3	H4B	B	902	-	-	0/8/17/17	0/2/2/2
5	HAR	B	906	-	-	0/5/13/13	0/0/0/0

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	901	HEM	C2D-C3D	-7.12	1.33	1.54
2	B	901	HEM	C2D-C3D	-6.87	1.33	1.54
2	A	901	HEM	C3B-CAB	-6.41	1.39	1.51
2	A	901	HEM	C3C-CAC	-6.37	1.39	1.51
2	B	901	HEM	C3B-CAB	-6.25	1.39	1.51
2	B	901	HEM	C3C-CAC	-6.11	1.39	1.51
2	B	901	HEM	C3B-C4B	-5.74	1.46	1.51
3	B	902	H4B	C7-N8	-5.60	1.38	1.46
3	A	902	H4B	C7-N8	-5.44	1.39	1.46
2	A	901	HEM	C3D-C4D	-4.25	1.46	1.51
2	A	901	HEM	C3B-C4B	-4.02	1.48	1.51
2	B	901	HEM	C3D-C4D	-3.98	1.46	1.51
2	A	901	HEM	C2C-C1C	-3.29	1.46	1.52
2	B	901	HEM	C2C-C1C	-3.05	1.46	1.52
2	B	901	HEM	C2B-C1B	-2.81	1.42	1.51
2	B	901	HEM	C2D-C1D	-2.28	1.44	1.51
2	A	901	HEM	C2B-C1B	-2.18	1.44	1.51
3	B	902	H4B	C4A-N5	-2.10	1.33	1.38
2	A	901	HEM	FE-NC	2.00	2.03	1.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	901	HEM	C4C-NC	2.06	1.38	1.36
2	B	901	HEM	C4C-NC	3.12	1.39	1.36

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	901	HEM	CBA-CAA-C2A	-3.72	105.85	112.53
2	A	901	HEM	CBA-CAA-C2A	-3.66	105.98	112.53
3	B	902	H4B	N3-C2-N1	-2.43	121.55	125.53
3	A	902	H4B	N3-C2-N1	-2.35	121.68	125.53
2	A	901	HEM	C3B-C4B-CHC	2.06	126.06	123.16
3	B	902	H4B	C7-C6-N5	2.29	115.21	110.45
2	B	901	HEM	CMD-C2D-C3D	2.37	124.84	114.35
2	A	901	HEM	CMD-C2D-C3D	2.42	125.05	114.35
3	B	902	H4B	C4-C4A-C8A	2.50	116.83	114.56
3	B	902	H4B	C2-N1-C8A	2.75	120.72	114.54
3	A	902	H4B	C2-N1-C8A	2.87	120.99	114.54
3	A	902	H4B	C4-C4A-C8A	2.92	117.21	114.56
3	A	902	H4B	C7-C6-N5	2.92	116.52	110.45
2	A	901	HEM	C3B-CAB-CBB	2.94	128.96	124.46
2	B	901	HEM	C3B-C4B-CHC	2.95	127.31	123.16
2	B	901	HEM	C2D-C3D-C4D	2.96	106.52	101.50
3	B	902	H4B	C4-N3-C2	3.39	120.64	115.94
3	A	902	H4B	C4-N3-C2	3.44	120.71	115.94
2	A	901	HEM	CAD-C3D-C2D	3.51	123.32	113.22
2	B	901	HEM	CAD-C3D-C2D	3.90	124.44	113.22
2	A	901	HEM	C2D-C3D-C4D	3.96	108.21	101.50
2	A	901	HEM	CAD-C3D-C4D	4.48	128.26	112.47
2	B	901	HEM	CAD-C3D-C4D	4.67	128.96	112.47
2	A	901	HEM	CMB-C2B-C3B	4.83	128.58	116.53
2	B	901	HEM	CMC-C2C-C3C	5.18	129.45	116.53
2	B	901	HEM	CMB-C2B-C3B	5.28	129.71	116.53
2	A	901	HEM	CMC-C2C-C3C	5.78	130.96	116.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	901	HEM	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	902	H4B	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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.