



# Full wwPDB X-ray Structure Validation Report ⓘ

Apr 2, 2014 – 11:30 AM EDT

PDB ID : 4N4N  
Title : Nitrosomonas europea HAO  
Authors : Maalcke, W.J.; Dietl, A.; Marritt, S.J.; Butt, J.N.; Jetten, M.S.M.; Keltjens, J.T.; Barends, T.R.M.B.; Kartal, B.  
Deposited on : 2013-10-08  
Resolution : 2.20 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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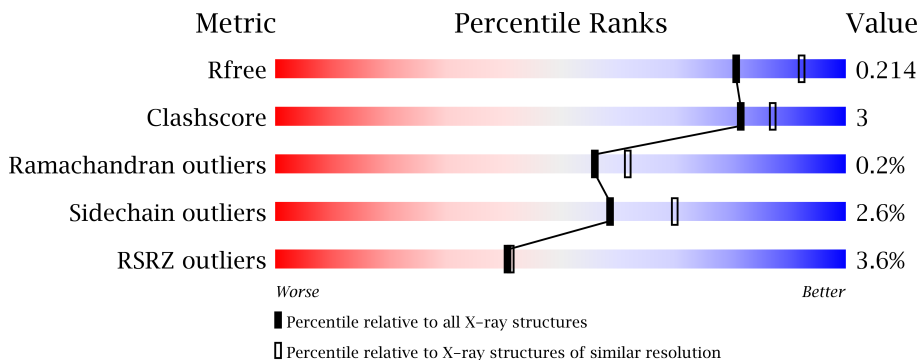
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : **FAILED**  
Xtriage (Phenix) : dev-1439  
EDS : stable22978  
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) : stable22978

# 1 Overall quality at a glance

The reported resolution of this entry is 2.20 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	2938 (2.20-2.20)
Clashscore	79885	3751 (2.20-2.20)
Ramachandran outliers	78287	3681 (2.20-2.20)
Sidechain outliers	78261	3682 (2.20-2.20)
RSRZ outliers	66119	2939 (2.20-2.20)

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	546	
1	C	546	
1	E	546	
2	B	57	
2	D	57	
2	F	57	

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 15450 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.

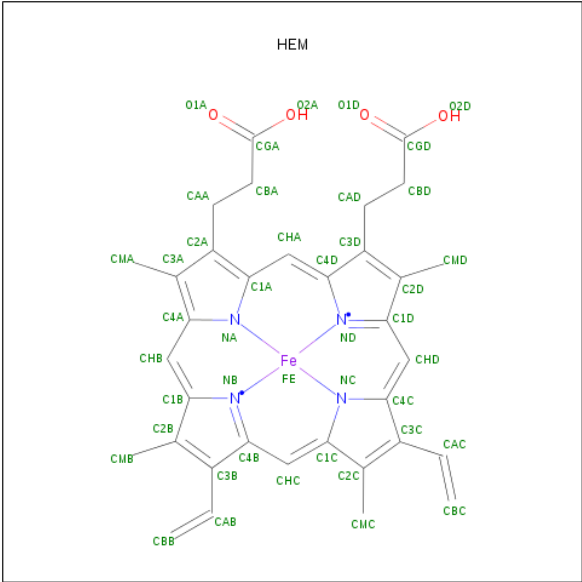
- Molecule 1 is a protein called Hydroxylamine oxidoreductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	504	Total	C	N	O	S	0	0	0
			4018	2499	713	774	32			
1	C	502	Total	C	N	O	S	0	0	0
			4005	2491	710	772	32			
1	E	503	Total	C	N	O	S	0	1	0
			4018	2499	713	774	32			

- Molecule 2 is a protein called hydroxylamine oxidoreductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	56	Total	C	N	O	S	0	0	0
			423	263	75	82	3			
2	D	57	Total	C	N	O	S	0	0	0
			429	266	76	84	3			
2	F	56	Total	C	N	O	S	0	0	0
			425	264	75	83	3			

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula:  $C_{34}H_{32}FeN_4O_4$ ).



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

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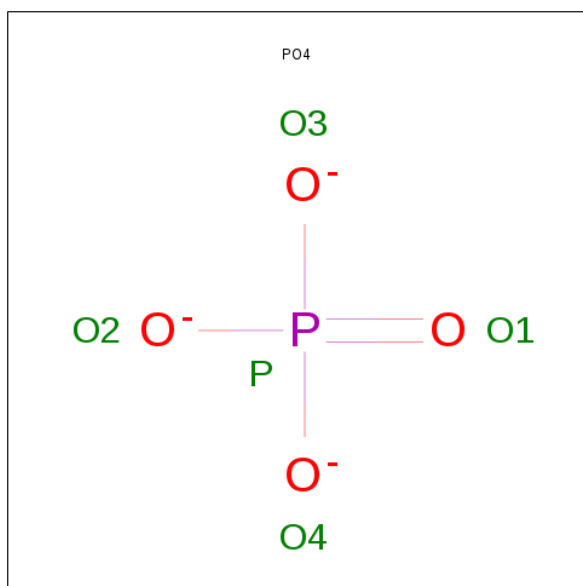
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- ORTEP diagram of the HEC molecule. The structure shows a central iron atom (Fe) coordinated by four nitrogen atoms (N1, N2, N3, N4) in a porphyrin-like ring. The structure is labeled with atom names (C1A, C2A, C3A, C4A, C1B, C2B, C3B, C4B, C1C, C2C, C3C, C4C, C1D, C2D, C3D, C4D) and includes thermal ellipsoids at the 50% probability level.

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

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- WORLD WIDE  
PDB  
PROTEIN DATA BANK

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total K 1 1	0	0

- Molecule 6 is PHOSPHATE ION (three-letter code: PO4) (formula: O<sub>4</sub>P).

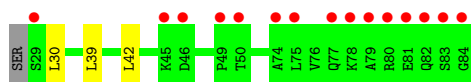


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	E	1	Total O P 5 4 1	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	330	Total O 330 330	0	0
7	B	42	Total O 42 42	0	0
7	C	328	Total O 328 328	0	0
7	D	37	Total O 37 37	0	0
7	E	314	Total O 314 314	0	0
7	F	43	Total O 43 43	0	0





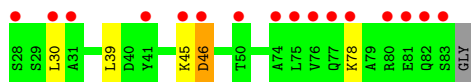
- Molecule 2: hydroxylamine oxidoreductase

Chain D: 



- Molecule 2: hydroxylamine oxidoreductase

Chain F: 





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	140.72Å 141.97Å 106.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.40 – 2.20 47.07 – 2.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (50.40-2.20) 99.6 (47.07-2.20)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.22 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, $R_{free}$	0.179 , 0.213 0.181 , 0.214	Depositor DCC
$R_{free}$ test set	10263 reflections (10.46%)	DCC
Wilson B-factor (Å <sup>2</sup> )	28.1	Xtriage
Anisotropy	0.039	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 35.1	EDS
Estimated twinning fraction	0.022 for k,h,-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 108386 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	15450	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.75% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, K, PO4, HEC

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	0/4121	0.50	0/5583
1	C	0.35	0/4108	0.51	0/5567
1	E	0.35	0/4121	0.51	0/5583
2	B	0.33	0/426	0.53	0/571
2	D	0.33	0/432	0.57	0/579
2	F	0.32	0/428	0.53	0/574
All	All	0.35	0/13636	0.51	0/18457

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 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	4018	0	3821	20	0
1	C	4005	0	3805	22	0
1	E	4018	0	3820	38	0
2	B	423	0	444	0	0
2	D	429	0	449	2	0
2	F	425	0	446	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	301	0	210	7	0
3	C	301	0	210	5	0
3	E	301	0	210	7	0
4	A	43	0	29	2	0
4	C	43	0	29	3	0
4	E	43	0	29	2	0
5	A	1	0	0	0	0
6	E	5	0	0	0	0
7	A	330	0	0	1	0
7	B	42	0	0	0	0
7	C	328	0	0	0	0
7	D	37	0	0	0	0
7	E	314	0	0	1	0
7	F	43	0	0	0	0
All	All	15450	0	13502	90	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 3.

All (90) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:179[B]:LYS:HD3	1:E:179[B]:LYS:N	1.62	1.10
1:E:179[B]:LYS:CD	1:E:179[B]:LYS:H	1.69	1.03
1:E:179[B]:LYS:HD3	1:E:179[B]:LYS:H	0.81	0.94
2:F:45:LYS:O	2:F:46:ASP:HB2	1.82	0.78
1:C:37:LEU:HD11	1:C:53:LEU:HD13	1.66	0.78
1:A:477:MET:SD	1:A:498:MET:CE	2.76	0.72
1:C:153:ARG:HD2	1:C:160:GLU:HA	1.70	0.72
1:C:149:GLU:O	1:C:153:ARG:HG3	1.91	0.70
1:E:207:ARG:HH11	1:E:229:ALA:HA	1.55	0.70
1:A:477:MET:SD	1:A:498:MET:HE2	2.33	0.69
1:A:37:LEU:HD11	1:A:53:LEU:HD13	1.75	0.69
4:A:608:HEC:HMB1	4:A:608:HEC:HBB3	1.76	0.68
1:E:207:ARG:NH1	1:E:229:ALA:HA	2.13	0.63
1:E:318:VAL:HG13	1:E:322:ASP:HB2	1.80	0.62
1:C:253:CYS:HA	4:E:601:HEC:HMC3	1.81	0.61
1:C:120:ARG:HH21	1:E:330:ASN:HB3	1.66	0.60
4:C:601:HEC:HMB1	4:C:601:HEC:HBB3	1.83	0.59
1:E:240:VAL:HG13	1:E:449:THR:HG23	1.84	0.59
4:E:601:HEC:HMB1	4:E:601:HEC:HBB3	1.86	0.58
1:A:477:MET:SD	1:A:498:MET:HE1	2.44	0.58
1:E:207:ARG:NH1	1:E:208:GLU:OE2	2.37	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:97:VAL:HG12	1:E:181:LYS:HE2	1.86	0.57
1:E:477:MET:SD	1:E:498:MET:HE1	2.45	0.57
1:A:311:ARG:NH2	7:A:924:HOH:O	2.35	0.56
1:E:179[B]:LYS:CD	1:E:179[B]:LYS:N	2.42	0.56
1:A:526:GLU:OE2	1:C:521:ARG:HD2	2.06	0.56
4:A:608:HEC:HMC3	1:E:253:CYS:HA	1.87	0.55
1:C:153:ARG:HD3	1:C:159:GLY:O	2.06	0.55
1:E:116:ARG:O	1:E:120:ARG:HG3	2.07	0.54
1:E:477:MET:SD	1:E:498:MET:CE	2.96	0.54
1:A:357:ASN:O	1:A:359:PRO:HD3	2.09	0.53
1:A:120:ARG:HH21	1:C:330:ASN:HB3	1.74	0.52
3:A:605:HEM:HMA1	3:A:606:HEM:HBA2	1.91	0.51
1:C:114:TRP:CE3	1:C:114:TRP:HA	2.46	0.51
1:A:114:TRP:HA	1:A:114:TRP:CE3	2.47	0.50
1:E:477:MET:HB2	1:E:494:GLY:HA2	1.92	0.50
1:A:385:THR:HA	1:A:388:HIS:O	2.12	0.49
1:E:483:HIS:HD2	3:E:606:HEM:C1C	2.30	0.49
3:E:606:HEM:HMA1	3:E:607:HEM:HBA2	1.94	0.49
1:E:354:ARG:NH2	1:E:401:ASP:OD1	2.45	0.49
1:A:253:CYS:HA	4:C:601:HEC:HMC3	1.93	0.49
1:E:240:VAL:HG22	7:E:802:HOH:O	2.12	0.49
1:E:26:ILE:HD11	1:E:201:LEU:HD13	1.96	0.48
1:C:318:VAL:HG13	1:C:322:ASP:HB2	1.95	0.48
1:E:114:TRP:HA	1:E:114:TRP:CE3	2.49	0.48
1:C:114:TRP:HE3	1:C:114:TRP:HA	1.79	0.47
1:C:491:TYR:HD2	4:C:601:HEC:HMC2	1.79	0.47
1:E:107:HIS:HA	1:E:110:GLU:HB3	1.96	0.47
1:A:483:HIS:HD2	3:A:605:HEM:C1C	2.33	0.47
1:A:114:TRP:HE3	1:A:114:TRP:HA	1.79	0.46
3:E:603:HEM:CHC	3:E:605:HEM:HBC1	2.45	0.46
1:E:411:TYR:CE1	1:E:470:ALA:HB2	2.50	0.46
1:E:153:ARG:HD3	1:E:160:GLU:HA	1.98	0.46
1:A:227:SER:O	3:A:606:HEM:HBD1	2.16	0.46
1:C:354:ARG:NH2	1:C:401:ASP:OD1	2.47	0.46
1:C:477:MET:HB2	1:C:494:GLY:HA2	1.99	0.45
3:A:601:HEM:HBC1	3:A:606:HEM:C4B	2.53	0.44
1:A:123:HIS:HB3	1:A:167:VAL:HB	1.99	0.44
1:C:381:VAL:O	1:C:385:THR:HG23	2.17	0.44
1:E:227:SER:O	3:E:607:HEM:HBD1	2.18	0.44
1:A:303:HIS:CD2	3:A:607:HEM:ND	2.86	0.44
1:A:272:PHE:O	3:A:602:HEM:HBA1	2.18	0.43
1:E:114:TRP:HA	1:E:114:TRP:HE3	1.84	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:269:ARG:HA	1:E:270:HIS:HA	1.80	0.43
1:A:29:VAL:HA	1:A:30:PRO:HD3	1.91	0.43
1:E:254:THR:O	1:E:258:THR:HG23	2.18	0.42
1:A:130:ARG:NH2	1:A:163:THR:HB	2.34	0.42
1:E:483:HIS:CD2	3:E:606:HEM:C1C	3.06	0.42
1:C:90:PRO:HB3	1:C:188:ILE:HB	2.01	0.42
3:C:606:HEM:HMA1	3:C:607:HEM:HBA2	2.00	0.42
1:E:499:ASN:O	1:E:503:VAL:HG23	2.19	0.42
1:E:84:PRO:HB3	1:E:190:MET:HB2	2.00	0.42
1:C:526:GLU:OE2	1:E:521:ARG:HD2	2.19	0.42
1:E:139:TYR:CE1	1:E:142:LYS:HD3	2.55	0.42
1:C:155:MET:CE	1:C:198:THR:HG23	2.50	0.42
1:C:406:GLU:HG3	2:D:39:LEU:HD21	2.02	0.42
1:E:98:ALA:HB1	1:E:102:ASP:HB2	2.01	0.42
1:E:295:TRP:CZ2	1:E:299:THR:HG21	2.55	0.42
1:C:84:PRO:HB3	1:C:190:MET:HB2	2.01	0.42
1:C:190:MET:HA	1:C:191:PRO:HD3	1.92	0.41
3:C:604:HEM:HAA1	3:C:605:HEM:HMA1	2.02	0.41
3:E:602:HEM:HBC1	3:E:607:HEM:C4B	2.56	0.41
1:A:257:HIS:CE1	3:A:606:HEM:HMD2	2.55	0.41
1:C:483:HIS:HD2	3:C:606:HEM:C1C	2.39	0.41
3:C:602:HEM:HBC1	3:C:607:HEM:C4B	2.56	0.41
2:D:31:ALA:HA	2:D:32:PRO:HD3	1.95	0.41
1:E:349:ILE:HD13	3:E:607:HEM:C2A	2.56	0.40
1:E:279:LYS:HA	1:E:280:PRO:HD3	1.97	0.40
1:E:149:GLU:O	1:E:153:ARG:HG3	2.21	0.40
3:C:603:HEM:CHC	3:C:605:HEM:HBC1	2.52	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	502/546 (92%)	487 (97%)	14 (3%)	1 (0%)	56 62

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	500/546 (92%)	485 (97%)	14 (3%)	1 (0%)	56	62
1	E	502/546 (92%)	488 (97%)	14 (3%)	0	100	100
2	B	54/57 (95%)	53 (98%)	1 (2%)	0	100	100
2	D	55/57 (96%)	54 (98%)	1 (2%)	0	100	100
2	F	54/57 (95%)	52 (96%)	1 (2%)	1 (2%)	12	7
All	All	1667/1809 (92%)	1619 (97%)	45 (3%)	3 (0%)	56	62

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	46	ASP
1	C	292	HIS
1	A	292	HIS

### 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	430/459 (94%)	421 (98%)	9 (2%)	66	78
1	C	429/459 (94%)	419 (98%)	10 (2%)	63	74
1	E	430/459 (94%)	421 (98%)	9 (2%)	66	78
2	B	48/49 (98%)	45 (94%)	3 (6%)	25	27
2	D	49/49 (100%)	45 (92%)	4 (8%)	17	15
2	F	49/49 (100%)	46 (94%)	3 (6%)	26	28
All	All	1435/1524 (94%)	1397 (97%)	38 (3%)	59	70

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	53	LEU
1	A	95	LYS
1	A	101	LYS
1	A	114	TRP

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Mol	Chain	Res	Type
1	A	201	LEU
1	A	382	LEU
1	A	464	LEU
1	A	467	LEU
1	A	474	LEU
2	B	30	LEU
2	B	39	LEU
2	B	42	LEU
1	C	53	LEU
1	C	114	TRP
1	C	201	LEU
1	C	267	HIS
1	C	290	VAL
1	C	312	ASP
1	C	318	VAL
1	C	320	LEU
1	C	464	LEU
1	C	474	LEU
2	D	30	LEU
2	D	39	LEU
2	D	66	ARG
2	D	77	GLN
1	E	114	TRP
1	E	201	LEU
1	E	240	VAL
1	E	267	HIS
1	E	318	VAL
1	E	432	THR
1	E	464	LEU
1	E	467	LEU
1	E	474	LEU
2	F	30	LEU
2	F	39	LEU
2	F	78	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	386	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

Mogul failed to run properly - this section will therefore be empty.

## 5.5 Carbohydrates ⓘ

Mogul failed to run properly - this section will therefore be empty.

## 5.6 Ligand geometry ⓘ

Mogul failed to run properly - this section will therefore be empty.

## 5.7 Other polymers ⓘ

Mogul failed to run properly - this section will therefore be empty.

## 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	504/546 (92%)	-0.28	6 (1%) 75 76	18, 26, 38, 61	0
1	C	502/546 (91%)	-0.24	9 (1%) 65 66	18, 26, 40, 56	0
1	E	503/546 (92%)	-0.29	5 (0%) 79 80	18, 26, 39, 54	0
2	B	56/57 (98%)	0.97	15 (26%) 1 1	28, 40, 67, 68	0
2	D	57/57 (100%)	0.43	10 (17%) 2 2	26, 32, 63, 65	0
2	F	56/57 (98%)	0.96	16 (28%) 1 1	29, 41, 69, 70	0
All	All	1678/1809 (92%)	-0.17	61 (3%) 41 41	18, 27, 44, 70	0

All (61) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	84	GLY	7.7
2	D	83	SER	5.9
2	B	79	ALA	5.5
1	A	527	GLY	4.8
2	F	28	SER	4.5
1	E	527	GLY	4.4
2	D	82	GLN	4.3
2	F	81	GLU	4.2
2	B	77	GLN	4.0
1	C	524	LYS	4.0
2	B	83	SER	3.8
2	B	46	ASP	3.8
2	D	84	GLY	3.7
2	B	81	GLU	3.7
2	B	75	LEU	3.5
2	D	81	GLU	3.5
2	D	46	ASP	3.4
2	B	78	LYS	3.3
1	C	523	ASN	3.3

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Mol	Chain	Res	Type	RSRZ
2	B	82	GLN	3.1
2	D	28	SER	3.1
2	B	29	SER	3.0
2	F	45	LYS	3.0
2	F	46	ASP	3.0
2	F	31	ALA	2.9
2	F	77	GLN	2.8
2	D	77	GLN	2.7
2	F	82	GLN	2.7
2	B	49	PRO	2.6
2	F	76	VAL	2.6
1	A	38	LYS	2.6
2	F	50	THR	2.6
1	C	133	LYS	2.6
2	B	45	LYS	2.5
1	E	515	LEU	2.5
2	D	30	LEU	2.5
1	E	161	LYS	2.5
2	D	79	ALA	2.5
1	A	525	LEU	2.5
2	F	41	TYR	2.4
1	E	135	ASP	2.4
2	F	30	LEU	2.4
2	D	29	SER	2.4
1	C	518	LEU	2.4
1	A	528	LYS	2.3
2	F	75	LEU	2.3
2	F	80	ARG	2.3
2	F	74	ALA	2.3
1	C	522	VAL	2.2
2	B	80	ARG	2.2
2	B	74	ALA	2.2
2	F	78	LYS	2.2
1	A	524	LYS	2.1
1	C	42	GLY	2.1
2	F	83	SER	2.1
2	B	50	THR	2.1
1	C	520	ALA	2.1
1	E	297	ALA	2.1
1	C	521	ARG	2.0
1	C	135	ASP	2.0
1	A	487	GLY	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
6	PO4	E	609	5/5	0.17	1.90	53,53,53,54	0
3	HEM	C	607	43/43	0.14	1.22	17,20,24,26	0
3	HEM	A	606	43/43	0.14	1.06	18,19,23,24	0
4	HEC	C	601	43/43	0.14	0.69	21,22,24,25	0
3	HEM	E	605	43/43	0.11	0.64	22,24,26,27	0
3	HEM	A	601	43/43	0.12	0.58	19,21,23,23	0
3	HEM	E	602	43/43	0.12	0.54	19,21,22,23	0
3	HEM	C	605	43/43	0.10	0.52	21,22,24,24	0
3	HEM	C	603	43/43	0.10	0.47	19,22,24,24	0
3	HEM	C	608	43/43	0.14	0.45	18,19,22,25	0
3	HEM	E	607	43/43	0.12	0.33	18,19,22,24	0
3	HEM	C	606	43/43	0.14	0.28	17,19,22,24	0
3	HEM	E	608	43/43	0.11	0.22	19,21,22,23	0
3	HEM	E	606	43/43	0.12	0.22	17,18,19,20	0
3	HEM	A	605	43/43	0.12	0.06	17,19,22,23	0
3	HEM	E	604	43/43	0.11	-0.11	26,31,34,34	0
3	HEM	A	603	43/43	0.10	-0.13	22,24,25,26	0
3	HEM	C	602	43/43	0.10	-0.16	19,21,23,23	0
3	HEM	E	603	43/43	0.10	-0.20	22,24,24,25	0
4	HEC	A	608	43/43	0.12	-0.28	20,22,23,24	0
3	HEM	A	604	43/43	0.10	-0.28	16,21,22,23	0
4	HEC	E	601	43/43	0.11	-0.31	21,23,25,26	0
3	HEM	A	607	43/43	0.11	-0.31	19,21,22,24	0
3	HEM	A	602	43/43	0.10	-0.32	22,23,24,24	0
3	HEM	C	604	43/43	0.10	-0.36	23,28,29,31	0
5	K	A	609	1/1	0.15	-0.83	38,38,38,38	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.