



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 11:08 AM GMT

PDB ID : 3O72
Title : Crystal structure of EfeB in complex with heme
Authors : Liu, X.; Du, Q.; Wang, Z.; Zhu, D.; Huang, Y.; Li, N.; Xu, S.; Gu, L.
Deposited on : 2010-07-30
Resolution : 1.95 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

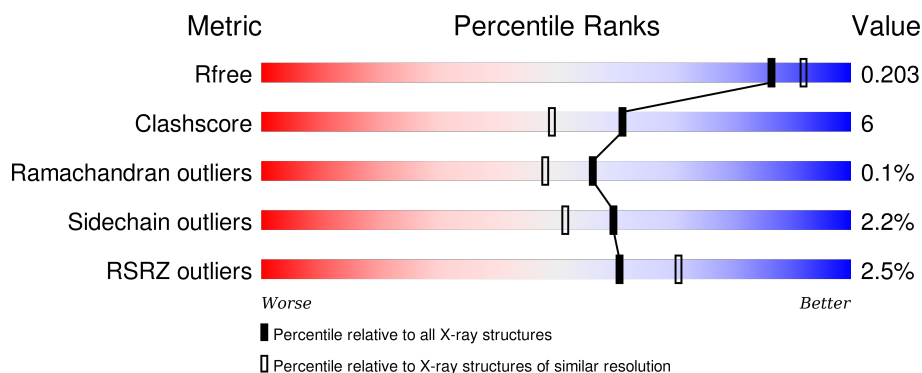
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

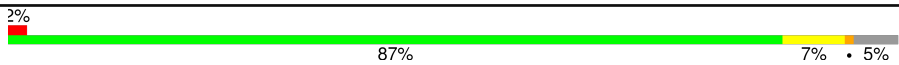



The reported resolution of this entry is 1.95 Å.

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	1833 (1.96-1.96)
Clashscore	102246	1953 (1.96-1.96)
Ramachandran outliers	100387	1936 (1.96-1.96)
Sidechain outliers	100360	1936 (1.96-1.96)
RSRZ outliers	91569	1835 (1.96-1.96)

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	396	
1	B	396	
1	C	396	
1	D	396	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 13094 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 Redox component of a tripartite ferrous iron transporter.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	375	Total	C	N	O	S	Se	0	1	0
			2939	1861	517	550	3	8			
1	B	375	Total	C	N	O	S	Se	0	1	0
			2939	1861	517	550	3	8			
1	C	375	Total	C	N	O	S	Se	0	1	0
			2939	1861	517	550	3	8			
1	D	375	Total	C	N	O	S	Se	0	1	0
			2939	1862	517	549	3	8			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	424	LEU	-	EXPRESSION TAG	UNP C6UPP3
A	425	GLU	-	EXPRESSION TAG	UNP C6UPP3
A	426	HIS	-	EXPRESSION TAG	UNP C6UPP3
A	427	HIS	-	EXPRESSION TAG	UNP C6UPP3
A	428	HIS	-	EXPRESSION TAG	UNP C6UPP3
A	429	HIS	-	EXPRESSION TAG	UNP C6UPP3
A	430	HIS	-	EXPRESSION TAG	UNP C6UPP3
A	431	HIS	-	EXPRESSION TAG	UNP C6UPP3
B	424	LEU	-	EXPRESSION TAG	UNP C6UPP3
B	425	GLU	-	EXPRESSION TAG	UNP C6UPP3
B	426	HIS	-	EXPRESSION TAG	UNP C6UPP3
B	427	HIS	-	EXPRESSION TAG	UNP C6UPP3
B	428	HIS	-	EXPRESSION TAG	UNP C6UPP3
B	429	HIS	-	EXPRESSION TAG	UNP C6UPP3
B	430	HIS	-	EXPRESSION TAG	UNP C6UPP3
B	431	HIS	-	EXPRESSION TAG	UNP C6UPP3
C	424	LEU	-	EXPRESSION TAG	UNP C6UPP3
C	425	GLU	-	EXPRESSION TAG	UNP C6UPP3
C	426	HIS	-	EXPRESSION TAG	UNP C6UPP3
C	427	HIS	-	EXPRESSION TAG	UNP C6UPP3
C	428	HIS	-	EXPRESSION TAG	UNP C6UPP3

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Chain	Residue	Modelled	Actual	Comment	Reference
C	429	HIS	-	EXPRESSION TAG	UNP C6UPP3
C	430	HIS	-	EXPRESSION TAG	UNP C6UPP3
C	431	HIS	-	EXPRESSION TAG	UNP C6UPP3
D	424	LEU	-	EXPRESSION TAG	UNP C6UPP3
D	425	GLU	-	EXPRESSION TAG	UNP C6UPP3
D	426	HIS	-	EXPRESSION TAG	UNP C6UPP3
D	427	HIS	-	EXPRESSION TAG	UNP C6UPP3
D	428	HIS	-	EXPRESSION TAG	UNP C6UPP3
D	429	HIS	-	EXPRESSION TAG	UNP C6UPP3
D	430	HIS	-	EXPRESSION TAG	UNP C6UPP3
D	431	HIS	-	EXPRESSION TAG	UNP C6UPP3

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- Chemical structure of HEM (Heme) showing a central iron atom coordinated by four nitrogen atoms in a porphyrin-like ring. The structure includes various side chains and a central iron atom labeled 'FE'.

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

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

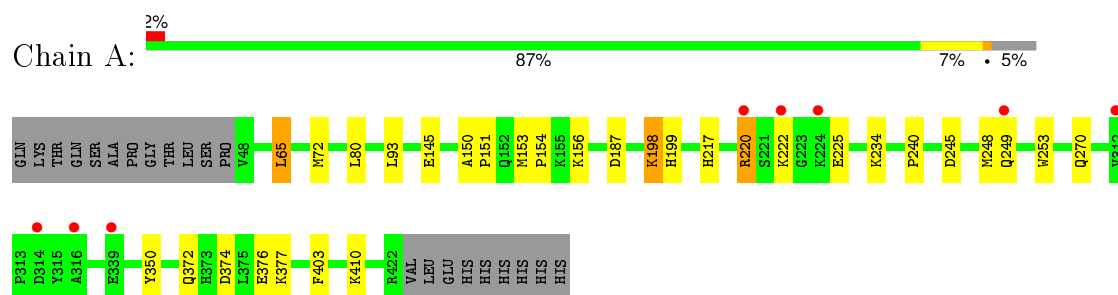
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	341	Total O 341 341	0	0
4	B	316	Total O 316 316	0	0
4	C	269	Total O 269 269	0	0
4	D	232	Total O 232 232	0	0

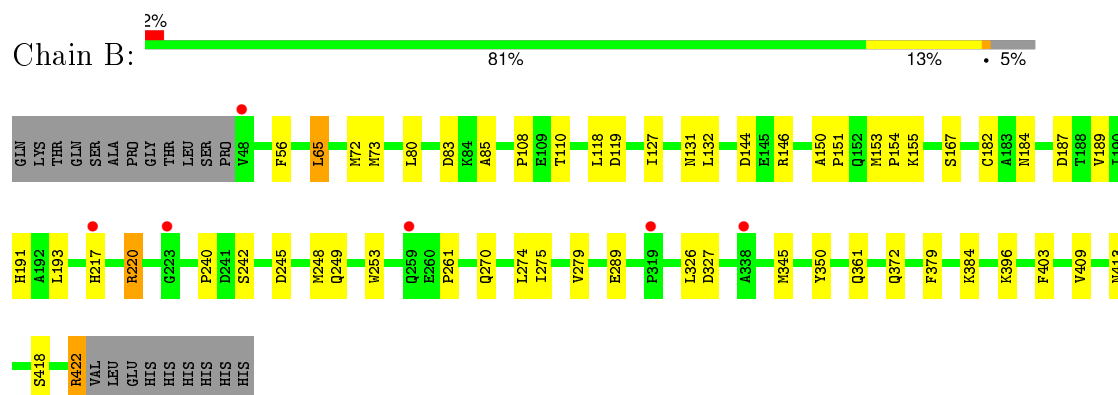
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

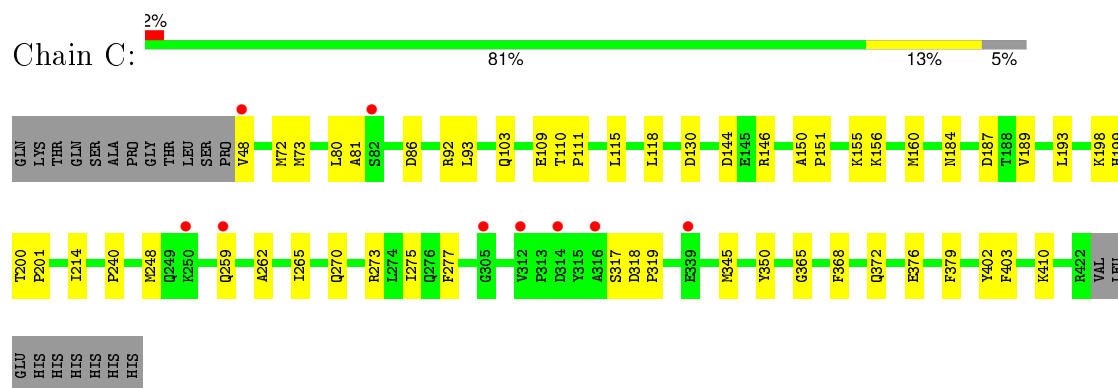
- Molecule 1: Redox component of a tripartite ferrous iron transporter



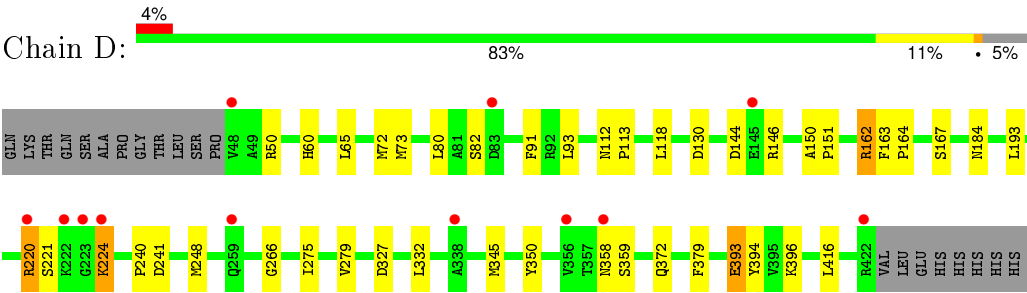
- Molecule 1: Redox component of a tripartite ferrous iron transporter



- Molecule 1: Redox component of a tripartite ferrous iron transporter



- Molecule 1: Redox component of a tripartite ferrous iron transporter



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	50.64Å 155.57Å 97.37Å 90.00° 94.70° 90.00°	Depositor
Resolution (Å)	50.00 – 1.95 42.34 – 1.95	Depositor EDS
% Data completeness (in resolution range)	92.6 (50.00-1.95) 92.6 (42.34-1.95)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.33 (at 1.95Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, R_{free}	0.170 , 0.206 0.168 , 0.203	Depositor DCC
R_{free} test set	1927 reflections (1.91%)	DCC
Wilson B-factor (Å ²)	20.8	Xtriage
Anisotropy	0.233	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 46.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 104476 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13094	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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, OXY

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	0.42	0/3000	0.55	0/4051
1	B	0.39	0/3000	0.56	0/4051
1	C	0.36	0/3000	0.53	0/4051
1	D	0.35	0/3000	0.51	0/4050
All	All	0.38	0/12000	0.54	0/16203

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-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 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	2939	0	2912	26	0
1	B	2939	0	2912	35	0
1	C	2939	0	2912	32	0
1	D	2939	0	2917	34	0
2	A	43	0	30	0	0
2	B	43	0	30	2	0
2	C	43	0	30	5	0
2	D	43	0	30	3	0
3	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
4	A	341	0	0	4	0
4	B	316	0	0	3	0
4	C	269	0	0	2	0
4	D	232	0	0	2	0
All	All	13094	0	11773	131	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 6.

All (131) 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:220:ARG:HB2	1:A:220:ARG:HH11	1.18	1.09
1:A:217:HIS:HA	1:A:220:ARG:HH12	1.29	0.96
1:D:162:ARG:HG3	1:D:162:ARG:HH11	1.37	0.90
1:A:220:ARG:HH11	1:A:220:ARG:CB	1.90	0.85
1:A:217:HIS:HA	1:A:220:ARG:NH1	1.96	0.81
1:D:220:ARG:HG2	1:D:221:SER:N	1.97	0.80
1:B:289:GLU:HG2	4:B:868:HOH:O	1.81	0.80
1:D:72:MSE:HG2	1:D:350:TYR:CG	2.17	0.79
1:A:156:LYS:HD2	1:A:376:GLU:HG3	1.69	0.75
1:C:72:MSE:HG2	1:C:350:TYR:CG	2.22	0.74
1:A:374:ASP:CG	1:A:377:LYS:HE3	2.07	0.74
1:B:144:ASP:OD1	1:B:146:ARG:HD3	1.88	0.73
1:B:361:GLN:HE22	1:C:92:ARG:HH22	1.40	0.69
1:C:144:ASP:OD1	1:C:146:ARG:HD3	1.92	0.69
1:C:262:ALA:O	1:C:265:ILE:HG12	1.93	0.69
1:A:220:ARG:NH1	1:A:220:ARG:HB2	2.02	0.67
1:A:245:ASP:O	1:A:249[A]:GLN:HG3	1.94	0.66
1:A:72:MSE:HG2	1:A:350:TYR:CG	2.30	0.66
1:A:374:ASP:OD1	1:A:377:LYS:HE3	1.96	0.66
1:B:73:MSE:HE3	1:B:189:VAL:HG12	1.78	0.65
1:B:72:MSE:HG2	1:B:350:TYR:CG	2.32	0.64
1:D:162:ARG:CG	1:D:162:ARG:HH11	2.11	0.63
1:D:144:ASP:OD1	1:D:146:ARG:HD3	1.98	0.62
1:D:217:HIS:HB2	1:D:220:ARG:CZ	2.30	0.62
1:D:72:MSE:HG2	1:D:350:TYR:CD1	2.35	0.61
1:D:162:ARG:NH1	1:D:162:ARG:HG3	2.14	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:73:MSE:HE3	1:C:189:VAL:HG12	1.84	0.58
1:B:220:ARG:HB3	1:B:220:ARG:CZ	2.33	0.57
1:B:73:MSE:SE	1:B:193:LEU:HD22	2.54	0.57
1:C:345:MSE:HE1	1:C:379:PHE:HA	1.87	0.56
1:C:72:MSE:HG2	1:C:350:TYR:CD1	2.41	0.55
1:D:217:HIS:HA	1:D:220:ARG:HD3	1.87	0.55
1:D:221:SER:HB2	1:D:224:LYS:HB2	1.88	0.55
1:A:153:MSE:HG3	1:A:154:PRO:HD2	1.88	0.54
1:D:217:HIS:HB2	1:D:220:ARG:NH1	2.21	0.54
1:C:130:ASP:OD1	1:C:184:ASN:HB2	2.06	0.54
1:B:167:SER:OG	1:B:396:LYS:HD2	2.07	0.54
1:A:410:LYS:HG2	4:A:711:HOH:O	2.07	0.54
1:D:93:LEU:HD21	1:D:199:HIS:O	2.08	0.54
1:C:146:ARG:HD2	4:C:783:HOH:O	2.06	0.54
1:B:108:PRO:HD2	1:B:119:ASP:HB2	1.89	0.54
2:B:500:HEM:HBB2	2:B:500:HEM:HHC	1.90	0.54
1:B:275:ILE:HG21	2:B:500:HEM:HBB1	1.88	0.54
1:C:275:ILE:HG21	2:C:500:HEM:HBB1	1.88	0.54
1:B:384:LYS:HE3	1:C:109:GLU:OE1	2.08	0.54
1:D:358:ASN:OD1	1:D:359:SER:N	2.38	0.53
1:A:145:GLU:CD	1:A:145:GLU:H	2.12	0.53
1:D:150:ALA:HB3	1:D:151:PRO:HD3	1.90	0.53
1:D:162:ARG:NE	4:D:563:HOH:O	2.37	0.53
1:D:118:LEU:HA	1:D:198[B]:LYS:HE2	1.91	0.52
1:A:198:LYS:NZ	4:A:688:HOH:O	2.39	0.52
1:D:72:MSE:HG2	1:D:350:TYR:CD2	2.44	0.52
4:B:1081:HOH:O	1:C:410:LYS:HD2	2.10	0.52
1:A:220:ARG:CG	1:A:220:ARG:HH11	2.23	0.52
1:D:240:PRO:HB2	1:D:248:MSE:HE1	1.91	0.52
2:C:500:HEM:CMC	2:C:500:HEM:HBC2	2.40	0.51
1:C:318:ASP:N	1:C:319:PRO:HD3	2.26	0.51
1:C:317:SER:C	1:C:319:PRO:HD3	2.31	0.51
1:B:72:MSE:HG2	1:B:350:TYR:CD2	2.46	0.51
1:C:118:LEU:HA	1:C:198:LYS:HE2	1.93	0.50
2:C:500:HEM:HBC2	2:C:500:HEM:HMC1	1.93	0.50
1:D:130:ASP:OD1	1:D:184:ASN:HB2	2.11	0.50
1:C:270:GLN:HE21	1:C:403:PHE:HB2	1.76	0.50
1:A:72:MSE:HG2	1:A:350:TYR:CD2	2.46	0.50
2:C:500:HEM:HHC	2:C:500:HEM:HBB2	1.94	0.49
1:C:214:ILE:HG23	1:C:350:TYR:CD1	2.47	0.49
1:C:103:GLN:HB2	4:C:518:HOH:O	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:MSE:HG2	1:A:350:TYR:CD1	2.48	0.49
1:B:72:MSE:HG2	1:B:350:TYR:CD1	2.48	0.49
1:C:73:MSE:SE	1:C:193:LEU:HD22	2.62	0.49
1:D:393:GLU:HG2	1:D:394:TYR:CD2	2.48	0.49
1:B:150:ALA:HB3	1:B:151:PRO:HD3	1.95	0.49
1:A:234:LYS:NZ	4:A:985:HOH:O	2.43	0.48
1:D:345:MSE:HE1	1:D:379:PHE:HA	1.95	0.48
1:D:241:ASP:O	1:D:248:MSE:HE3	2.14	0.48
1:A:65:LEU:HD22	1:A:253:TRP:HZ2	1.79	0.47
1:D:167:SER:HB2	1:D:396:LYS:HE3	1.96	0.47
1:B:144:ASP:OD2	1:B:146:ARG:HD2	2.14	0.47
1:B:217:HIS:CD2	1:B:217:HIS:H	2.33	0.47
1:D:358:ASN:ND2	4:D:798:HOH:O	2.48	0.47
1:B:240:PRO:HB2	1:B:248:MSE:HE1	1.96	0.47
1:A:270:GLN:HE21	1:A:403:PHE:HB2	1.79	0.47
1:B:242:SER:HA	1:B:248:MSE:SE	2.65	0.47
1:C:240:PRO:HB2	1:C:248:MSE:HE1	1.97	0.47
1:D:393:GLU:HG2	1:D:394:TYR:CE2	2.51	0.46
1:B:270:GLN:HE21	1:B:403:PHE:HB2	1.81	0.46
1:C:273:ARG:HB2	1:C:368:PHE:HB3	1.98	0.46
1:D:162:ARG:CG	1:D:162:ARG:NH1	2.72	0.46
1:D:163:PHE:HB3	1:D:164:PRO:HD2	1.98	0.46
1:B:56:PHE:CE2	1:B:409:VAL:HG11	2.51	0.46
1:C:150:ALA:HB3	1:C:151:PRO:HD3	1.97	0.45
1:C:156:LYS:HD2	1:C:376:GLU:HG3	1.99	0.45
1:C:160:MSE:HA	1:C:402:TYR:CE2	2.52	0.45
1:D:91:PHE:HB3	1:D:416:LEU:O	2.16	0.45
1:B:245:ASP:O	1:B:249[A]:GLN:HG3	2.16	0.44
1:A:222:LYS:CE	4:A:848:HOH:O	2.64	0.44
1:B:131:ASN:HB2	1:B:184:ASN:ND2	2.33	0.44
1:B:110:THR:HA	1:B:118:LEU:HD23	1.99	0.44
1:A:217:HIS:CE1	1:A:225:GLU:OE2	2.70	0.44
1:B:153:MSE:HG3	1:B:154:PRO:HD2	2.00	0.44
1:A:240:PRO:HB2	1:A:248:MSE:HE1	1.99	0.44
1:D:60:HIS:HD2	1:D:266:GLY:O	2.00	0.44
1:B:65:LEU:HD22	1:B:253:TRP:HZ2	1.83	0.44
1:D:275:ILE:HG21	2:D:500:HEM:HBB1	2.00	0.44
1:D:73:MSE:SE	1:D:193:LEU:HD22	2.68	0.43
1:D:167:SER:CB	1:D:396:LYS:HE3	2.49	0.43
1:B:345:MSE:HE1	1:B:379:PHE:HA	2.01	0.43
1:C:110:THR:HA	1:C:111:PRO:HD3	1.81	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:327:ASP:HA	1:D:332:LEU:HD11	2.00	0.43
1:B:418:SER:O	1:B:422:ARG:HG2	2.18	0.43
1:C:93:LEU:HD21	1:C:199:HIS:O	2.19	0.43
1:D:112:ASN:HA	1:D:113:PRO:HD3	1.88	0.43
1:C:155:LYS:HG3	1:C:259:GLN:O	2.19	0.43
1:C:277:PHE:CG	1:C:365:GLY:HA2	2.52	0.43
1:C:72:MSE:HG2	1:C:350:TYR:CD2	2.54	0.42
2:D:500:HEM:HHC	2:D:500:HEM:HBB2	2.02	0.42
1:B:83:ASP:HB3	1:B:85:ALA:H	1.84	0.42
1:B:326:LEU:HB2	1:C:115:LEU:HD21	2.02	0.42
1:A:93:LEU:HD21	1:A:199:HIS:O	2.20	0.42
1:B:327:ASP:HB2	4:B:718:HOH:O	2.19	0.42
1:B:155:LYS:NZ	1:B:261:PRO:HD3	2.35	0.41
1:A:220:ARG:NH1	1:A:220:ARG:CG	2.84	0.41
2:D:500:HEM:CMC	2:D:500:HEM:HBC2	2.51	0.41
1:B:155:LYS:HZ2	1:B:261:PRO:HD3	1.85	0.41
1:C:200:THR:N	1:C:201:PRO:CD	2.84	0.41
1:B:127:ILE:HD12	1:B:191:HIS:CE1	2.56	0.40
1:B:132:LEU:HA	1:B:182:CYS:O	2.21	0.40
1:B:422:ARG:HB3	1:B:422:ARG:HH11	1.85	0.40
2:C:500:HEM:CBB	2:C:500:HEM:HHC	2.52	0.40
1:C:81:ALA:HB1	1:C:86:ASP:HB3	2.03	0.40
1:A:150:ALA:HB3	1:A:151:PRO:HD3	2.03	0.40

There are no symmetry-related clashes.

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	374/396 (94%)	366 (98%)	8 (2%)	0	100	100
1	B	374/396 (94%)	364 (97%)	10 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	374/396 (94%)	365 (98%)	9 (2%)	0	100	100
1	D	374/396 (94%)	363 (97%)	10 (3%)	1 (0%)	46	35
All	All	1496/1584 (94%)	1458 (98%)	37 (2%)	1 (0%)	56	48

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	224	LYS

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	313/323 (97%)	307 (98%)	6 (2%)	65	58
1	B	313/323 (97%)	304 (97%)	9 (3%)	50	38
1	C	313/323 (97%)	309 (99%)	4 (1%)	76	72
1	D	313/323 (97%)	304 (97%)	9 (3%)	50	38
All	All	1252/1292 (97%)	1224 (98%)	28 (2%)	60	51

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	LEU
1	A	80	LEU
1	A	187	ASP
1	A	198	LYS
1	A	220	ARG
1	A	372	GLN
1	B	65	LEU
1	B	80	LEU
1	B	187	ASP
1	B	220	ARG
1	B	274	LEU

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Mol	Chain	Res	Type
1	B	279	VAL
1	B	372	GLN
1	B	413	ASN
1	B	422	ARG
1	C	48	VAL
1	C	80	LEU
1	C	187	ASP
1	C	372	GLN
1	D	50	ARG
1	D	65	LEU
1	D	80	LEU
1	D	82	SER
1	D	162	ARG
1	D	220	ARG
1	D	279	VAL
1	D	372	GLN
1	D	393	GLU

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	51	ASN
1	A	165	ASN
1	A	259	GLN
1	A	270	GLN
1	A	310	HIS
1	B	259	GLN
1	B	270	GLN
1	B	276	GLN
1	C	51	ASN
1	C	165	ASN
1	C	270	GLN
1	D	60	HIS
1	D	140	HIS
1	D	259	GLN
1	D	270	GLN
1	D	310	HIS

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 ⓘ

8 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	500	1,3	30,50,50	2.21	8 (26%)	24,82,82	2.35	11 (45%)
3	OXY	A	600	2	1,1,1	0.80	0	0,0,0	0.00	-
2	HEM	B	500	1,3	30,50,50	2.05	7 (23%)	24,82,82	2.46	10 (41%)
3	OXY	B	600	2	1,1,1	0.76	0	0,0,0	0.00	-
2	HEM	C	500	1,3	30,50,50	2.20	9 (30%)	24,82,82	2.38	11 (45%)
3	OXY	C	600	2	1,1,1	0.78	0	0,0,0	0.00	-
2	HEM	D	500	1,3	30,50,50	2.14	7 (23%)	24,82,82	2.35	10 (41%)
3	OXY	D	600	2	1,1,1	0.77	0	0,0,0	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	500	1,3	-	0/10/54/54	0/0/8/8
3	OXY	A	600	2	-	0/0/0/0	0/0/0/0
2	HEM	B	500	1,3	-	0/10/54/54	0/0/8/8
3	OXY	B	600	2	-	0/0/0/0	0/0/0/0
2	HEM	C	500	1,3	-	0/10/54/54	0/0/8/8
3	OXY	C	600	2	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	D	500	1,3	-	0/10/54/54	0/0/8/8
3	OXY	D	600	2	-	0/0/0/0	0/0/0/0

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	HEM	C3B-C4B	-7.25	1.45	1.51
2	D	500	HEM	C3B-C4B	-7.24	1.45	1.51
2	C	500	HEM	C3B-C4B	-7.21	1.45	1.51
2	A	500	HEM	C3D-C4D	-5.72	1.44	1.51
2	B	500	HEM	C3B-C4B	-5.58	1.46	1.51
2	B	500	HEM	C3D-C4D	-5.15	1.45	1.51
2	C	500	HEM	C3D-C4D	-5.02	1.45	1.51
2	D	500	HEM	C3D-C4D	-4.80	1.45	1.51
2	C	500	HEM	C2C-C1C	-4.23	1.44	1.52
2	B	500	HEM	C2C-C1C	-3.58	1.45	1.52
2	D	500	HEM	C2C-C1C	-3.52	1.45	1.52
2	A	500	HEM	C2C-C1C	-3.38	1.46	1.52
2	C	500	HEM	C2D-C1D	-2.24	1.44	1.51
2	D	500	HEM	C2D-C1D	-2.21	1.44	1.51
2	B	500	HEM	C2B-C1B	-2.19	1.44	1.51
2	A	500	HEM	C2B-C1B	-2.13	1.44	1.51
2	C	500	HEM	C2B-C1B	-2.11	1.45	1.51
2	D	500	HEM	C3C-CAC	2.05	1.55	1.51
2	C	500	HEM	FE-ND	2.07	2.08	1.97
2	A	500	HEM	FE-ND	2.12	2.08	1.97
2	B	500	HEM	FE-NC	2.29	2.04	1.95
2	C	500	HEM	CAA-C2A	2.29	1.55	1.52
2	C	500	HEM	C1C-NC	2.32	1.38	1.36
2	A	500	HEM	CAA-C2A	2.34	1.56	1.52
2	D	500	HEM	FE-NB	2.34	2.09	1.97
2	A	500	HEM	C4C-NC	2.42	1.39	1.36
2	C	500	HEM	FE-NC	2.47	2.05	1.95
2	B	500	HEM	C1C-NC	2.49	1.39	1.36
2	A	500	HEM	C1C-NC	2.63	1.39	1.36
2	B	500	HEM	FE-ND	2.66	2.11	1.97
2	D	500	HEM	FE-NC	2.72	2.06	1.95

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	HEM	C3B-CAB-CBB	-4.01	118.31	124.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	HEM	C3B-CAB-CBB	-3.51	119.07	124.46
2	C	500	HEM	C3B-CAB-CBB	-3.33	119.34	124.46
2	D	500	HEM	C3B-CAB-CBB	-3.20	119.55	124.46
2	C	500	HEM	C3C-CAC-CBC	-2.66	120.37	124.46
2	A	500	HEM	C3B-C4B-NB	-2.53	106.80	111.63
2	C	500	HEM	CBD-CAD-C3D	-2.32	106.79	113.55
2	B	500	HEM	C3B-C4B-NB	-2.30	107.23	111.63
2	B	500	HEM	CBD-CAD-C3D	-2.29	106.90	113.55
2	A	500	HEM	CBD-CAD-C3D	-2.22	107.09	113.55
2	C	500	HEM	C3B-C4B-NB	-2.16	107.49	111.63
2	D	500	HEM	C3B-C4B-NB	-2.15	107.51	111.63
2	D	500	HEM	CMA-C3A-C4A	-2.09	124.91	128.36
2	A	500	HEM	C3B-C4B-CHC	2.02	126.00	123.16
2	C	500	HEM	C3B-C4B-CHC	2.05	126.05	123.16
2	A	500	HEM	C2C-C1C-CHC	2.21	127.04	123.68
2	C	500	HEM	C2D-C3D-C4D	2.42	105.61	101.50
2	A	500	HEM	C2D-C3D-C4D	2.61	105.92	101.50
2	B	500	HEM	C2D-C3D-C4D	2.61	105.92	101.50
2	D	500	HEM	C3B-C4B-CHC	2.68	126.94	123.16
2	B	500	HEM	CMD-C2D-C3D	2.70	126.29	114.35
2	A	500	HEM	CMD-C2D-C3D	2.76	126.56	114.35
2	D	500	HEM	C2D-C3D-C4D	2.81	106.26	101.50
2	D	500	HEM	CMD-C2D-C3D	2.90	127.17	114.35
2	D	500	HEM	CMB-C2B-C3B	3.06	124.18	116.53
2	C	500	HEM	CMD-C2D-C3D	3.08	127.98	114.35
2	B	500	HEM	C3B-C4B-CHC	3.34	127.86	123.16
2	C	500	HEM	CMB-C2B-C3B	3.68	125.72	116.53
2	A	500	HEM	CMC-C2C-C3C	3.90	126.27	116.53
2	B	500	HEM	CMB-C2B-C3B	4.01	126.55	116.53
2	A	500	HEM	CMB-C2B-C3B	4.20	127.01	116.53
2	D	500	HEM	CAD-C3D-C4D	4.26	127.50	112.47
2	B	500	HEM	CAD-C3D-C4D	4.29	127.59	112.47
2	C	500	HEM	CAD-C3D-C4D	4.35	127.82	112.47
2	A	500	HEM	CAD-C3D-C4D	4.37	127.90	112.47
2	A	500	HEM	CAD-C3D-C2D	4.51	126.18	113.22
2	D	500	HEM	CAD-C3D-C2D	4.52	126.22	113.22
2	C	500	HEM	CMC-C2C-C3C	4.56	127.92	116.53
2	B	500	HEM	CAD-C3D-C2D	4.62	126.49	113.22
2	C	500	HEM	CAD-C3D-C2D	4.64	126.56	113.22
2	D	500	HEM	CMC-C2C-C3C	4.92	128.82	116.53
2	B	500	HEM	CMC-C2C-C3C	4.98	128.97	116.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	500	HEM	2	0
2	C	500	HEM	5	0
2	D	500	HEM	3	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	367/396 (92%)	-0.04	8 (2%) 65 74	11, 18, 39, 52	1 (0%)
1	B	367/396 (92%)	-0.08	6 (1%) 74 83	11, 20, 41, 54	1 (0%)
1	C	367/396 (92%)	0.10	9 (2%) 61 71	13, 23, 42, 64	1 (0%)
1	D	367/396 (92%)	0.22	14 (3%) 44 56	15, 25, 45, 63	1 (0%)
All	All	1468/1584 (92%)	0.05	37 (2%) 61 71	11, 22, 42, 64	4 (0%)

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	48	VAL	10.8
1	D	223	GLY	5.9
1	D	259	GLN	5.3
1	C	48	VAL	4.1
1	A	220	ARG	3.9
1	D	358	ASN	3.7
1	D	224	LYS	3.6
1	B	259	GLN	3.5
1	D	145	GLU	3.4
1	D	222	LYS	3.3
1	C	316	ALA	3.1
1	B	217	HIS	3.0
1	D	422	ARG	2.9
1	C	259	GLN	2.9
1	B	223	GLY	2.9
1	C	339	GLU	2.7
1	B	338	ALA	2.6
1	D	220	ARG	2.5
1	D	83	ASP	2.5
1	C	312	VAL	2.4
1	A	222	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	316	ALA	2.3
1	A	249[A]	GLN	2.3
1	C	250	LYS	2.3
1	A	339	GLU	2.3
1	A	224	LYS	2.3
1	C	314	ASP	2.2
1	C	305	GLY	2.2
1	A	312	VAL	2.2
1	C	82	SER	2.1
1	A	314	ASP	2.1
1	D	198[A]	LYS	2.1
1	D	338	ALA	2.1
1	B	48	VAL	2.1
1	D	217	HIS	2.1
1	B	319	PRO	2.1
1	D	356	VAL	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(\AA^2)	Q<0.9
2	HEM	B	500	43/43	0.99	0.09	-0.27	11,14,17,18	0
2	HEM	C	500	43/43	0.97	0.10	-0.28	16,19,25,26	0
2	HEM	A	500	43/43	0.98	0.09	-0.30	9,13,17,18	0
2	HEM	D	500	43/43	0.98	0.08	-1.26	17,19,23,24	0
3	OXY	A	600	2/2	0.95	0.15	-	18,18,18,24	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	OXY	B	600	2/2	0.97	0.08	-	19,19,19,22	0
3	OXY	C	600	2/2	0.95	0.15	-	26,26,26,29	0
3	OXY	D	600	2/2	0.97	0.07	-	24,24,24,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.