



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 11:22 PM GMT

PDB ID : 1WNX  
Title : D136E mutant of Heme Oxygenase from *Corynebacterium diphtheriae* (HmuO)  
Authors : Unno, M.; Matsui, T.; Ikeda-Saito, M.  
Deposited on : 2004-08-10  
Resolution : 1.85 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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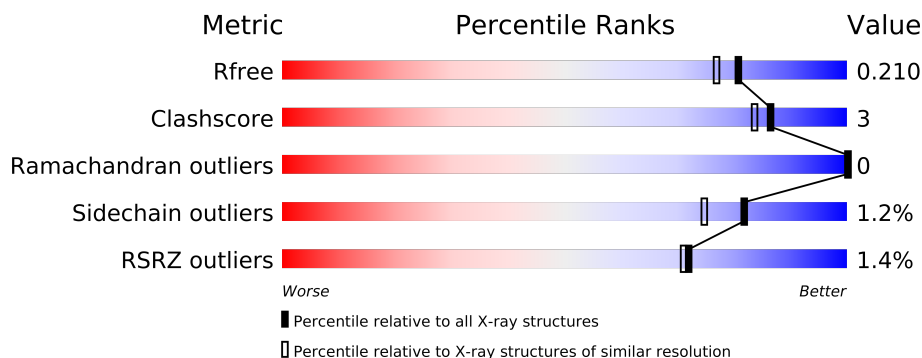
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.85 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	1269 (1.86-1.86)
Clashscore	79885	1470 (1.86-1.86)
Ramachandran outliers	78287	1451 (1.86-1.86)
Sidechain outliers	78261	1451 (1.86-1.86)
RSRZ outliers	66119	1269 (1.86-1.86)

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	215	
1	B	215	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	SO4	A	1001	-	X
2	SO4	A	1003	-	X
2	SO4	A	1004	-	X
2	SO4	A	1005	-	X
2	SO4	B	1002	-	X
2	SO4	B	1007	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
2	SO4	B	1008	-	X
2	SO4	B	1009	-	X
2	SO4	B	1010	-	X
2	SO4	B	1011	-	X
3	NA	A	3001	-	X
4	HEM	A	901	-	X
4	HEM	B	902	-	X

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3835 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 Heme oxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	207	Total	C	N	O	S	0	0	0
			1653	1043	293	314	3			
1	B	210	Total	C	N	O	S	0	1	0
			1673	1054	296	320	3			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	34	LYS	GLU	SEE REMARK 999	UNP P71119
A	60	VAL	ALA	SEE REMARK 999	UNP P71119
A	92	GLY	ASP	SEE REMARK 999	UNP P71119
A	93	SER	GLY	SEE REMARK 999	UNP P71119
A	136	GLU	ASP	ENGINEERED	UNP P71119
A	192	HIS	ASN	SEE REMARK 999	UNP P71119
B	34	LYS	GLU	SEE REMARK 999	UNP P71119
B	60	VAL	ALA	SEE REMARK 999	UNP P71119
B	92	GLY	ASP	SEE REMARK 999	UNP P71119
B	93	SER	GLY	SEE REMARK 999	UNP P71119
B	136	GLU	ASP	ENGINEERED	UNP P71119
B	192	HIS	ASN	SEE REMARK 999	UNP P71119

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).

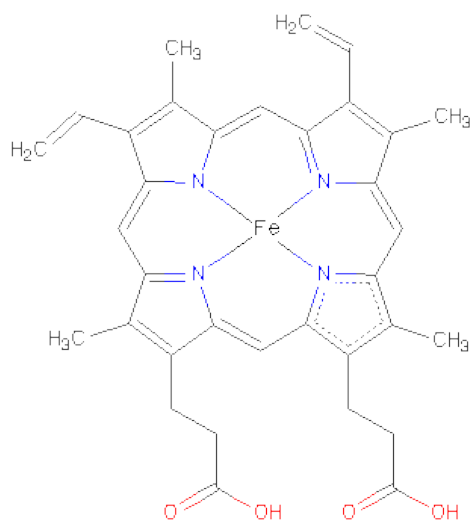


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total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

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Na	0	0
			1	1		

- Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula:  $C_{34}H_{32}FeN_4O_4$ ).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	182	Total	O	0	0
			182	182		
5	B	185	Total	O	0	0
			185	185		

### 3 Residue-property plots

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

- Molecule 1: Heme oxygenase

Chain A: 



- Molecule 1: Heme oxygenase

Chain B: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	106.15Å 63.72Å 79.30Å 90.00° 130.28° 90.00°	Depositor
Resolution (Å)	40.00 – 1.85 49.50 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.9 (40.00-1.85) 99.9 (49.50-1.85)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.15 (at 1.86Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, $R_{free}$	0.150 , 0.201 0.166 , 0.210	Depositor DCC
$R_{free}$ test set	3443 reflections (11.08%)	DCC
Wilson B-factor (Å <sup>2</sup> )	19.1	Xtriage
Anisotropy	0.394	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 39.1	EDS
Estimated twinning fraction	0.016 for h,-k,-h-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 34531 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	3835	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.30% 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, SO4, NA

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	1.12	3/1685 (0.2%)	1.11	15/2275 (0.7%)
1	B	1.06	2/1709 (0.1%)	0.99	3/2310 (0.1%)
All	All	1.09	5/3394 (0.1%)	1.05	18/4585 (0.4%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	41	GLU	CD-OE1	6.13	1.32	1.25
1	B	136	GLU	CD-OE2	5.81	1.32	1.25
1	A	109	TYR	CE2-CZ	-5.62	1.31	1.38
1	A	196	GLU	CD-OE1	-5.53	1.19	1.25
1	B	22	LYS	CE-NZ	5.38	1.62	1.49

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	44	ARG	NE-CZ-NH2	-10.39	115.11	120.30
1	A	29	MET	CG-SD-CE	-9.49	85.02	100.20
1	A	44	ARG	NE-CZ-NH1	8.72	124.66	120.30
1	B	29	MET	CG-SD-CE	-7.07	88.89	100.20
1	A	109	TYR	CB-CG-CD2	-7.06	116.77	121.00
1	A	85	ARG	NE-CZ-NH2	-7.00	116.80	120.30
1	A	154	ASP	CB-CG-OD2	6.70	124.33	118.30
1	A	108	ASP	CB-CG-OD2	6.48	124.13	118.30
1	A	109	TYR	CB-CG-CD1	6.28	124.77	121.00
1	A	74	ASP	CB-CG-OD1	6.19	123.87	118.30
1	A	132	ARG	NE-CZ-NH1	6.12	123.36	120.30
1	A	61	ASP	CB-CG-OD1	5.92	123.63	118.30
1	A	97	ARG	NE-CZ-NH1	5.87	123.23	120.30
1	A	132	ARG	NE-CZ-NH2	-5.65	117.48	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	118	ASP	CB-CG-OD1	5.58	123.32	118.30
1	A	97	ARG	NE-CZ-NH2	-5.32	117.64	120.30
1	B	86	ASP	CB-CG-OD2	5.03	122.83	118.30
1	B	210	ASP	CB-CG-OD1	5.01	122.81	118.30

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	1653	0	12	4	0
1	B	1673	0	23	5	0
2	A	25	0	0	0	0
2	B	30	0	0	0	0
3	A	1	0	0	0	0
4	A	43	0	0	1	0
4	B	43	0	0	0	0
5	A	182	0	0	2	0
5	B	185	0	0	3	0
All	All	3835	0	35	9	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 (9) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:141:GLN:CG	5:A:3161:HOH:O	2.49	0.60
1:B:132:ARG:NH1	1:B:204:ASN:OD1	2.41	0.54
1:B:162:HIS:CE1	5:B:1162:HOH:O	2.65	0.49
1:B:97:ARG:NE	5:B:1109:HOH:O	2.48	0.45
1:B:13:LYS:NZ	5:B:1158:HOH:O	2.52	0.43
1:A:208:PHE:CZ	4:A:901:HEM:CBB	3.03	0.41
1:A:112:ARG:NH2	1:A:196:GLU:OE1	2.53	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:58:GLN:OE1	1:B:117:ARG:NH1	2.54	0.41
1:A:141:GLN:NE2	5:A:3161:HOH:O	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	205/215 (95%)	203 (99%)	2 (1%)	0	100	100
1	B	209/215 (97%)	205 (98%)	4 (2%)	0	100	100
All	All	414/430 (96%)	408 (99%)	6 (1%)	0	100	100

There are no Ramachandran outliers to report.

### 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	170/175 (97%)	167 (98%)	3 (2%)	71	57
1	B	172/175 (98%)	171 (99%)	1 (1%)	92	89
All	All	342/350 (98%)	338 (99%)	4 (1%)	82	75

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	83	LEU
1	A	89	LYS

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Mol	Chain	Res	Type
1	A	184	GLU
1	B	169	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 ⓘ

Of 14 ligands modelled in this entry, 1 is monoatomic - leaving 13 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	SO4	A	1001	-	4,4,4	0.06	0	6,6,6	0.24	0
2	SO4	A	1003	-	4,4,4	0.30	0	6,6,6	0.59	0
2	SO4	A	1004	-	4,4,4	0.29	0	6,6,6	0.49	0
2	SO4	A	1005	-	4,4,4	0.92	0	6,6,6	0.96	0
2	SO4	A	1006	-	4,4,4	0.16	0	6,6,6	0.77	0
4	HEM	A	901	1,5	49,50,50	2.87	18 (36%)	46,82,82	2.33	16 (34%)
2	SO4	B	1002	-	4,4,4	0.17	0	6,6,6	0.34	0
2	SO4	B	1007	-	4,4,4	0.36	0	6,6,6	0.33	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	SO4	B	1008	-	4,4,4	0.22	0	6,6,6	0.23	0
2	SO4	B	1009	-	4,4,4	0.20	0	6,6,6	0.22	0
2	SO4	B	1010	-	4,4,4	1.26	1 (25%)	6,6,6	0.55	0
2	SO4	B	1011	-	4,4,4	0.17	0	6,6,6	0.28	0
4	HEM	B	902	1,5	49,50,50	3.64	13 (26%)	46,82,82	1.84	7 (15%)

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	SO4	A	1001	-	-	0/0/0/0	0/0/0/0
2	SO4	A	1003	-	-	0/0/0/0	0/0/0/0
2	SO4	A	1004	-	-	0/0/0/0	0/0/0/0
2	SO4	A	1005	-	-	0/0/0/0	0/0/0/0
2	SO4	A	1006	-	-	0/0/0/0	0/0/0/0
4	HEM	A	901	1,5	-	0/14/114/114	0/0/8/8
2	SO4	B	1002	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1007	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1008	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1009	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1010	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1011	-	-	0/0/0/0	0/0/0/0
4	HEM	B	902	1,5	-	0/14/114/114	0/0/8/8

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	902	HEM	C2B-C1B	20.16	1.49	1.44
4	B	902	HEM	C3C-C2C	-7.52	1.30	1.43
4	A	901	HEM	C3D-C4D	7.48	1.46	1.44
4	A	901	HEM	C2D-C1D	7.40	1.46	1.44
4	A	901	HEM	C4A-C3A	6.62	1.48	1.40
4	A	901	HEM	C3B-C2B	-6.00	1.33	1.43
4	A	901	HEM	C3C-C2C	-5.48	1.34	1.43
4	A	901	HEM	C3D-C2D	5.15	1.52	1.43
4	B	902	HEM	C3D-C2D	5.05	1.52	1.43
4	B	902	HEM	C3C-CAC	4.66	1.55	1.40
4	A	901	HEM	C3C-CAC	4.47	1.54	1.40
4	B	902	HEM	C3B-C2B	-4.41	1.36	1.43
4	B	902	HEM	C3B-CAB	4.41	1.54	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	901	HEM	CAA-C2A	4.30	1.59	1.52
4	B	902	HEM	C2D-C1D	-3.85	1.43	1.44
4	B	902	HEM	C4A-C3A	3.73	1.44	1.40
4	A	901	HEM	C3B-CAB	3.60	1.51	1.40
4	B	902	HEM	FE-NC	3.44	2.10	1.97
4	A	901	HEM	FE-NC	3.25	2.10	1.97
4	A	901	HEM	FE-NA	3.21	2.06	1.92
4	A	901	HEM	CMD-C2D	3.10	1.57	1.47
4	A	901	HEM	FE-ND	2.93	2.08	1.97
4	A	901	HEM	C1C-NC	-2.89	1.33	1.38
4	B	902	HEM	CMD-C2D	2.67	1.55	1.47
4	A	901	HEM	C4A-CHB	-2.65	1.32	1.39
4	A	901	HEM	CMC-C2C	2.62	1.55	1.47
4	A	901	HEM	CMB-C2B	2.55	1.55	1.47
4	A	901	HEM	C2C-C1C	2.40	1.50	1.43
2	B	1010	SO4	O1-S	2.19	1.54	1.47
4	B	902	HEM	CMC-C2C	2.08	1.53	1.47
4	B	902	HEM	CMB-C2B	2.02	1.53	1.47
4	B	902	HEM	FE-NA	2.02	2.01	1.92

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	901	HEM	C3B-C4B-NB	-7.28	108.79	114.00
4	B	902	HEM	C4D-ND-C1D	6.22	111.52	105.16
4	A	901	HEM	C4C-NC-C1C	5.34	111.09	105.53
4	B	902	HEM	CHD-C1D-ND	5.19	128.90	124.58
4	A	901	HEM	C4D-ND-C1D	4.35	109.61	105.16
4	B	902	HEM	C3B-C4B-NB	-4.13	111.05	114.00
4	A	901	HEM	C3A-C4A-NA	-3.83	106.52	109.41
4	B	902	HEM	C2D-C1D-ND	-3.60	108.67	112.93
4	A	901	HEM	CHB-C4A-NA	3.44	130.33	124.58
4	A	901	HEM	CAD-C3D-C4D	3.42	130.68	124.53
4	A	901	HEM	CHD-C4C-NC	3.40	127.69	124.73
4	A	901	HEM	CMA-C3A-C4A	-3.16	123.76	128.62
4	A	901	HEM	CAA-CBA-CGA	-3.10	103.49	113.47
4	A	901	HEM	C2D-C1D-ND	-2.93	109.47	112.93
4	B	902	HEM	CAD-CBD-CGD	2.78	122.15	113.48
4	A	901	HEM	C4A-CHB-C1B	-2.75	123.85	127.47
4	A	901	HEM	CHB-C1B-NB	-2.74	120.55	124.31
4	B	902	HEM	CHC-C4B-NB	2.69	126.82	124.58
4	A	901	HEM	C4A-NA-C1A	2.64	110.24	106.76

*Continued on next page...*

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	901	HEM	O1A-CGA-CBA	-2.48	114.48	123.03
4	A	901	HEM	CHD-C1D-ND	2.43	126.60	124.58
4	B	902	HEM	CHD-C4C-NC	-2.42	122.63	124.73
4	A	901	HEM	C2A-C1A-NA	-2.00	106.95	109.73

There are no chirality outliers.

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	207/215 (96%)	-0.48	1 (0%) 88 89	12, 20, 33, 43	0
1	B	210/215 (97%)	-0.43	3 (1%) 72 71	13, 21, 37, 49	0
All	All	417/430 (96%)	-0.45	4 (0%) 72 78	12, 20, 37, 49	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	3	THR	3.9
1	B	6	ALA	3.5
1	A	167	ALA	3.2
1	B	4	ALA	2.3

### 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( $\text{\AA}^2$ )	Q<0.9
2	SO4	B	1008	5/5	0.39	37.46	90,91,92,92	0
3	NA	A	3001	1/1	0.31	13.52	26,26,26,26	0
2	SO4	B	1011	5/5	0.29	10.33	83,85,85,85	0
2	SO4	B	1002	5/5	0.24	8.14	70,70,72,73	0
2	SO4	B	1010	5/5	0.22	7.57	43,44,50,54	0
2	SO4	B	1009	5/5	0.44	6.24	87,87,88,89	0
2	SO4	B	1007	5/5	0.20	5.70	71,73,75,75	0
2	SO4	A	1004	5/5	0.26	4.76	68,69,72,72	0
2	SO4	A	1005	5/5	0.27	4.27	41,41,48,51	0
4	HEM	B	902	43/43	0.14	2.91	21,42,72,79	0
4	HEM	A	901	43/43	0.14	2.84	20,32,62,76	0
2	SO4	A	1003	5/5	0.17	2.21	45,49,50,54	0
2	SO4	A	1001	5/5	0.18	2.03	75,75,77,78	0
2	SO4	A	1006	5/5	0.23	0.82	62,65,68,68	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.