



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:28 PM GMT

PDB ID : 1FS8  
Title : CYTOCHROME C NITRITE REDUCTASE FROM WOLINELLA  
SUCCINOGENES-SULFATE COMPLEX  
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Deposited on : 2000-09-08  
Resolution : 1.60 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

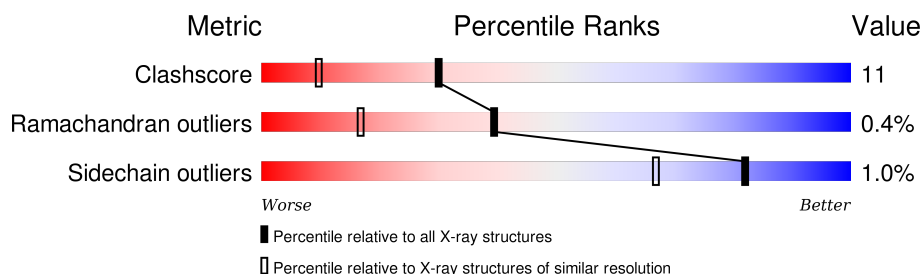
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.60 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	2732 (1.60-1.60)
Ramachandran outliers	100387	2654 (1.60-1.60)
Sidechain outliers	100360	2653 (1.60-1.60)

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	485	

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 4661 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 CYTOCHROME C NITRITE REDUCTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	471	Total	C	N	O	S	7	0	0
			3772	2387	648	716	21			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	45	GLU	LYS	CONFLICT	UNP Q9S1E5

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Ca	0	0
			1	1		

- Molecule 3 is YTTRIUM ION (three-letter code: Y1) (formula: Y).

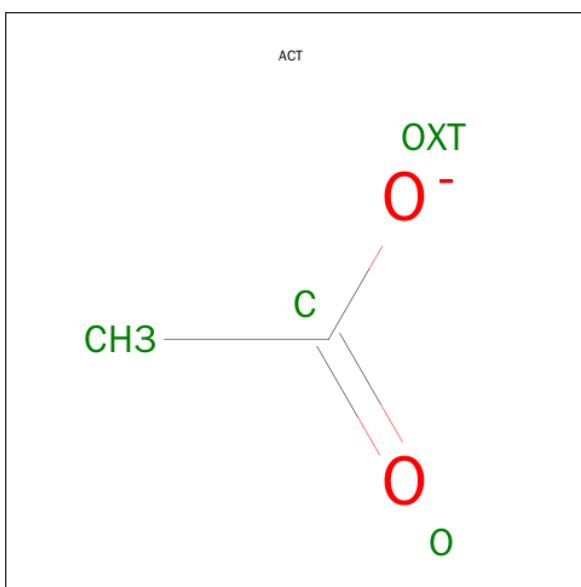
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	4	Total	Y	0	0
			4	4		

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



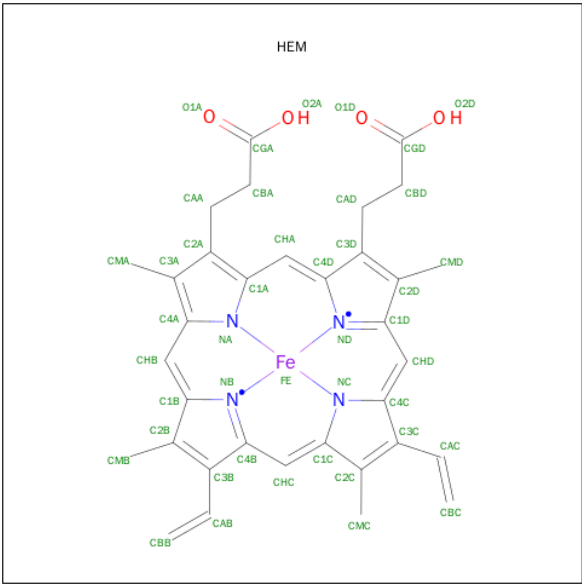
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is ACETATE ION (three-letter code: ACT) (formula: C<sub>2</sub>H<sub>3</sub>O<sub>2</sub>).



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

- Molecule 6 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: C<sub>34</sub>H<sub>32</sub>FeN<sub>4</sub>O<sub>4</sub>).



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

- Molecule 7 is water.

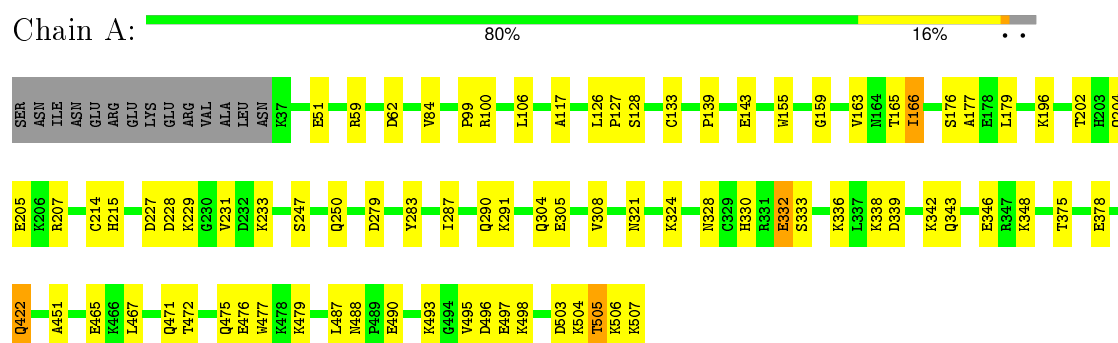
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	650	Total	O		
			650	650	0	0

### 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.

Note EDS was not executed.

#### • Molecule 1: CYTOCHROME C NITRITE REDUCTASE



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	119.01Å 119.01Å 186.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 1.60	Depositor
% Data completeness (in resolution range)	(Not available) (25.00-1.60)	Depositor
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	CNS	Depositor
R, $R_{free}$	0.184 , 0.206	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4661	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	25.0	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: Y1, HEM, CA, SO4, ACT

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	0.29	0/3863	0.56	0/5202

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

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	3772	0	3682	87	0
2	A	1	0	0	0	0
3	A	4	0	0	0	0
4	A	15	0	0	0	0
5	A	4	0	3	0	0
6	A	215	0	150	5	0
7	A	650	0	0	10	0
All	All	4661	0	3835	87	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 11.

All (87) 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:498:LYS:HE3	1:A:506:LYS:HA	1.26	1.09
1:A:100:ARG:HH22	1:A:507:LYS:HD2	1.33	0.94
1:A:100:ARG:NH2	1:A:507:LYS:HD2	1.94	0.82
1:A:333:SER:HB3	1:A:336:LYS:HG3	1.62	0.79
1:A:498:LYS:O	1:A:507:LYS:HE3	1.83	0.79
1:A:488:ASN:HB3	7:A:1298:HOH:O	1.85	0.77
1:A:117:ALA:HB3	1:A:465:GLU:HG2	1.69	0.74
1:A:100:ARG:HH22	1:A:507:LYS:CD	2.00	0.73
1:A:422:GLN:HE21	1:A:422:GLN:HA	1.54	0.73
1:A:497:GLU:HG3	1:A:507:LYS:HA	1.72	0.72
1:A:139:PRO:O	1:A:143:GLU:HG3	1.90	0.70
1:A:467:LEU:O	1:A:471:GLN:HG3	1.94	0.68
1:A:498:LYS:H	1:A:507:LYS:CE	2.07	0.67
1:A:51:GLU:HB2	7:A:1035:HOH:O	1.96	0.66
1:A:497:GLU:H	1:A:507:LYS:HG2	1.60	0.66
1:A:117:ALA:CB	1:A:465:GLU:HG2	2.27	0.65
1:A:496:ASP:HA	1:A:507:LYS:HE2	1.79	0.64
1:A:498:LYS:H	1:A:507:LYS:HE3	1.63	0.64
1:A:106:LEU:HD21	1:A:163:VAL:HA	1.82	0.62
1:A:497:GLU:HB2	1:A:507:LYS:HG2	1.81	0.62
1:A:497:GLU:N	1:A:507:LYS:HE2	2.15	0.62
1:A:228:ASP:OD1	1:A:229:LYS:HG3	1.99	0.62
1:A:204:GLN:HA	1:A:204:GLN:HE21	1.66	0.61
1:A:503:ASP:C	1:A:505:THR:H	2.04	0.59
1:A:62:ASP:HB3	7:A:1272:HOH:O	2.02	0.59
1:A:100:ARG:HH22	1:A:507:LYS:CE	2.17	0.58
1:A:330:HIS:HB3	1:A:332:GLU:OE2	2.05	0.55
1:A:497:GLU:HB2	1:A:507:LYS:HA	1.88	0.55
1:A:328:ASN:HB2	7:A:1122:HOH:O	2.07	0.55
1:A:100:ARG:CZ	1:A:507:LYS:HD2	2.37	0.54
1:A:228:ASP:HB3	7:A:1304:HOH:O	2.07	0.54
1:A:497:GLU:CG	1:A:507:LYS:HA	2.38	0.54
1:A:100:ARG:NH1	1:A:507:LYS:HD2	2.24	0.53
1:A:495:VAL:O	1:A:507:LYS:NZ	2.42	0.53
1:A:215:HIS:HB3	1:A:279:ASP:HB2	1.91	0.53
1:A:490:GLU:HB2	7:A:1298:HOH:O	2.09	0.53
1:A:497:GLU:HA	1:A:504:LYS:HD2	1.91	0.52
1:A:100:ARG:HH12	1:A:507:LYS:HD2	1.74	0.52
1:A:159:GLY:HA3	1:A:477:TRP:CE2	2.45	0.52
1:A:202:THR:OG1	1:A:205:GLU:HG3	2.10	0.51
1:A:179:LEU:HD11	6:A:509:HEM:HBD1	1.92	0.51
1:A:496:ASP:CA	1:A:507:LYS:HE2	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:476:GLU:HG2	7:A:1264:HOH:O	2.11	0.51
1:A:422:GLN:NE2	1:A:422:GLN:HA	2.26	0.50
1:A:504:LYS:O	1:A:505:THR:CB	2.59	0.50
1:A:342:LYS:O	1:A:346:GLU:HG3	2.12	0.50
1:A:196:LYS:NZ	1:A:196:LYS:HB2	2.27	0.50
1:A:290:GLN:NE2	7:A:1219:HOH:O	2.45	0.50
1:A:498:LYS:H	1:A:507:LYS:CG	2.24	0.49
1:A:99:PRO:HG3	6:A:508:HEM:CGD	2.41	0.49
1:A:497:GLU:N	1:A:507:LYS:HG2	2.27	0.49
1:A:497:GLU:CB	1:A:507:LYS:HG2	2.42	0.49
1:A:59:ARG:HD2	1:A:305:GLU:OE2	2.12	0.48
1:A:287:ILE:HA	1:A:290:GLN:HE21	1.78	0.48
1:A:291:LYS:HG2	6:A:512:HEM:CMD	2.43	0.48
1:A:204:GLN:NE2	1:A:204:GLN:HA	2.29	0.48
1:A:207:ARG:NH1	6:A:509:HEM:O1D	2.46	0.48
1:A:472:THR:O	1:A:476:GLU:HG3	2.13	0.48
1:A:227:ASP:OD2	1:A:233:LYS:HE3	2.15	0.47
1:A:503:ASP:C	1:A:505:THR:N	2.68	0.47
1:A:488:ASN:ND2	7:A:1035:HOH:O	2.47	0.47
1:A:321:ASN:OD1	1:A:324:LYS:HG2	2.15	0.46
1:A:247:SER:OG	1:A:250:GLN:HG3	2.16	0.45
1:A:475:GLN:O	1:A:479:LYS:HG3	2.17	0.45
1:A:497:GLU:HB2	1:A:507:LYS:CA	2.46	0.45
1:A:338:LYS:HB2	1:A:338:LYS:HZ2	1.81	0.45
1:A:165:THR:O	1:A:166:ILE:C	2.55	0.45
1:A:84:VAL:HB	1:A:451:ALA:HB1	1.99	0.44
1:A:504:LYS:O	1:A:505:THR:HB	2.18	0.44
1:A:498:LYS:N	1:A:507:LYS:HG3	2.33	0.44
1:A:496:ASP:HA	1:A:507:LYS:CE	2.48	0.43
1:A:133:CYS:HB3	1:A:214:CYS:HB3	1.99	0.43
1:A:328:ASN:ND2	7:A:1127:HOH:O	2.52	0.42
1:A:128:SER:HA	1:A:155:TRP:HA	2.01	0.42
1:A:176:SER:O	1:A:177:ALA:HB3	2.19	0.42
1:A:498:LYS:HG2	1:A:503:ASP:OD1	2.19	0.42
1:A:490:GLU:O	1:A:493:LYS:HG3	2.19	0.42
1:A:375:THR:OG1	1:A:378:GLU:HG3	2.20	0.42
1:A:196:LYS:HB2	1:A:196:LYS:HZ3	1.84	0.42
1:A:100:ARG:HH22	1:A:507:LYS:NZ	2.19	0.41
1:A:227:ASP:OD2	1:A:231:VAL:HB	2.21	0.41
1:A:304:GLN:HA	1:A:308:VAL:O	2.21	0.41
1:A:133:CYS:HB3	1:A:214:CYS:SG	2.61	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:LEU:HB3	1:A:127:PRO:HD2	2.03	0.40
1:A:339:ASP:O	1:A:343:GLN:HG3	2.22	0.40
1:A:291:LYS:HG2	6:A:512:HEM:HMD1	2.03	0.40
1:A:163:VAL:HG21	1:A:487:LEU:HB2	2.04	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	469/485 (97%)	453 (97%)	14 (3%)	2 (0%)	39 17

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	505	THR
1	A	166	ILE

### 5.3.2 Protein sidechains [i](#)

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	397/410 (97%)	393 (99%)	4 (1%)	82 67

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

Mol	Chain	Res	Type
1	A	283	TYR
1	A	332	GLU
1	A	348	LYS
1	A	422	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	ASN
1	A	204	GLN
1	A	290	GLN
1	A	328	ASN
1	A	335	GLN
1	A	422	GLN
1	A	480	GLN

### 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](#)

Of 14 ligands modelled in this entry, 5 are monoatomic - leaving 9 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
6	HEM	A	508	1,4	30,50,50	2.66	12 (40%)	24,82,82	2.18	7 (29%)
6	HEM	A	509	1	30,50,50	2.70	11 (36%)	24,82,82	2.14	8 (33%)
6	HEM	A	510	1	30,50,50	2.64	11 (36%)	24,82,82	2.08	7 (29%)
6	HEM	A	511	1	30,50,50	2.76	9 (30%)	24,82,82	2.19	8 (33%)
6	HEM	A	512	1	30,50,50	2.75	10 (33%)	24,82,82	2.36	8 (33%)
4	SO4	A	656	6	4,4,4	0.25	0	6,6,6	0.11	0
4	SO4	A	657	-	4,4,4	0.22	0	6,6,6	0.08	0
4	SO4	A	658	-	4,4,4	0.23	0	6,6,6	0.13	0
5	ACT	A	659	-	1,3,3	3.89	1 (100%)	0,3,3	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
6	HEM	A	508	1,4	-	0/10/54/54	0/0/8/8
6	HEM	A	509	1	-	0/10/54/54	0/0/8/8
6	HEM	A	510	1	-	0/10/54/54	0/0/8/8
6	HEM	A	511	1	-	0/10/54/54	0/0/8/8
6	HEM	A	512	1	-	0/10/54/54	0/0/8/8
4	SO4	A	656	6	-	0/0/0/0	0/0/0/0
4	SO4	A	657	-	-	0/0/0/0	0/0/0/0
4	SO4	A	658	-	-	0/0/0/0	0/0/0/0
5	ACT	A	659	-	-	0/0/0/0	0/0/0/0

All (54) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	511	HEM	C3B-C4B	-7.73	1.45	1.51
6	A	512	HEM	C3B-C4B	-6.91	1.45	1.51
6	A	509	HEM	C3B-C4B	-6.61	1.46	1.51
6	A	511	HEM	C2D-C3D	-6.49	1.35	1.54
6	A	510	HEM	C2D-C3D	-6.46	1.35	1.54
6	A	508	HEM	C2D-C3D	-6.41	1.35	1.54
6	A	512	HEM	C2D-C3D	-6.38	1.35	1.54
6	A	509	HEM	C2D-C3D	-6.13	1.36	1.54
6	A	509	HEM	C3B-CAB	-5.76	1.40	1.51
6	A	508	HEM	C3D-C4D	-5.76	1.44	1.51
6	A	512	HEM	C3D-C4D	-5.69	1.44	1.51
6	A	508	HEM	C3B-CAB	-5.65	1.40	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	512	HEM	C3B-CAB	-5.65	1.40	1.51
6	A	508	HEM	C3B-C4B	-5.54	1.46	1.51
6	A	509	HEM	C3C-CAC	-5.50	1.41	1.51
6	A	510	HEM	C3B-CAB	-5.47	1.41	1.51
6	A	510	HEM	C3B-C4B	-5.46	1.47	1.51
6	A	511	HEM	C3C-CAC	-5.46	1.41	1.51
6	A	512	HEM	C3C-CAC	-5.45	1.41	1.51
6	A	510	HEM	C3C-CAC	-5.27	1.41	1.51
6	A	511	HEM	C3B-CAB	-5.20	1.41	1.51
6	A	508	HEM	C3C-CAC	-5.12	1.41	1.51
6	A	511	HEM	C3D-C4D	-5.06	1.45	1.51
6	A	510	HEM	C3D-C4D	-4.87	1.45	1.51
6	A	509	HEM	C3D-C4D	-4.85	1.45	1.51
6	A	510	HEM	C2C-C1C	-4.75	1.43	1.52
6	A	509	HEM	C2C-C1C	-3.93	1.45	1.52
6	A	511	HEM	C2C-C1C	-3.84	1.45	1.52
6	A	512	HEM	C2C-C1C	-3.81	1.45	1.52
6	A	508	HEM	C2C-C1C	-3.66	1.45	1.52
6	A	508	HEM	C2D-C1D	-2.22	1.44	1.51
6	A	511	HEM	C2B-C1B	-2.18	1.44	1.51
6	A	510	HEM	C2D-C1D	-2.15	1.44	1.51
6	A	508	HEM	C2B-C1B	-2.12	1.44	1.51
6	A	509	HEM	C2B-C1B	-2.12	1.44	1.51
6	A	510	HEM	C2B-C1B	-2.11	1.44	1.51
6	A	512	HEM	C1C-NC	2.02	1.38	1.36
6	A	509	HEM	CBB-CAB	2.07	1.41	1.29
6	A	509	HEM	CBC-CAC	2.13	1.41	1.29
6	A	511	HEM	CBC-CAC	2.14	1.41	1.29
6	A	510	HEM	CBB-CAB	2.14	1.41	1.29
6	A	512	HEM	CBC-CAC	2.15	1.41	1.29
6	A	512	HEM	CBB-CAB	2.19	1.42	1.29
6	A	509	HEM	C4C-NC	2.23	1.38	1.36
6	A	512	HEM	C4C-NC	2.24	1.38	1.36
6	A	508	HEM	C4C-NC	2.25	1.38	1.36
6	A	511	HEM	CBB-CAB	2.32	1.42	1.29
6	A	510	HEM	C4C-NC	2.34	1.38	1.36
6	A	508	HEM	CBC-CAC	2.34	1.42	1.29
6	A	510	HEM	CBC-CAC	2.37	1.43	1.29
6	A	508	HEM	CBB-CAB	2.38	1.43	1.29
6	A	508	HEM	C1C-NC	2.44	1.39	1.36
6	A	509	HEM	C1C-NC	3.01	1.39	1.36
5	A	659	ACT	CH3-C	3.89	1.54	1.48

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	512	HEM	CAA-C2A-C1A	-4.98	121.60	127.01
6	A	508	HEM	C3B-CAB-CBB	-3.53	119.05	124.46
6	A	511	HEM	C3B-CAB-CBB	-2.96	119.92	124.46
6	A	509	HEM	C3B-CAB-CBB	-2.79	120.17	124.46
6	A	509	HEM	C3C-CAC-CBC	-2.65	120.39	124.46
6	A	512	HEM	CBA-CAA-C2A	-2.34	108.33	112.53
6	A	510	HEM	C3C-CAC-CBC	-2.10	121.23	124.46
6	A	511	HEM	C3C-CAC-CBC	-2.07	121.28	124.46
6	A	510	HEM	CMD-C2D-C3D	2.49	125.35	114.35
6	A	511	HEM	CMD-C2D-C3D	2.50	125.42	114.35
6	A	511	HEM	C2D-C3D-C4D	2.83	106.29	101.50
6	A	508	HEM	CMD-C2D-C3D	2.83	126.89	114.35
6	A	512	HEM	CMD-C2D-C3D	2.88	127.08	114.35
6	A	509	HEM	CMD-C2D-C3D	2.90	127.17	114.35
6	A	510	HEM	C2D-C3D-C4D	3.02	106.61	101.50
6	A	512	HEM	CAD-C3D-C4D	3.07	123.30	112.47
6	A	512	HEM	C2D-C3D-C4D	3.16	106.85	101.50
6	A	509	HEM	C2D-C3D-C4D	3.23	106.98	101.50
6	A	509	HEM	CAD-C3D-C4D	3.51	124.85	112.47
6	A	508	HEM	C2D-C3D-C4D	3.63	107.66	101.50
6	A	509	HEM	CMB-C2B-C3B	3.71	125.79	116.53
6	A	508	HEM	CAD-C3D-C2D	3.77	124.05	113.22
6	A	508	HEM	CMB-C2B-C3B	3.94	126.36	116.53
6	A	510	HEM	CAD-C3D-C2D	4.01	124.76	113.22
6	A	512	HEM	CMC-C2C-C3C	4.02	126.57	116.53
6	A	509	HEM	CMC-C2C-C3C	4.03	126.58	116.53
6	A	512	HEM	CMB-C2B-C3B	4.08	126.71	116.53
6	A	510	HEM	CMC-C2C-C3C	4.22	127.06	116.53
6	A	511	HEM	CAD-C3D-C2D	4.34	125.68	113.22
6	A	510	HEM	CMB-C2B-C3B	4.34	127.38	116.53
6	A	508	HEM	CAD-C3D-C4D	4.36	127.86	112.47
6	A	511	HEM	CAD-C3D-C4D	4.39	127.95	112.47
6	A	511	HEM	CMC-C2C-C3C	4.41	127.53	116.53
6	A	508	HEM	CMC-C2C-C3C	4.44	127.61	116.53
6	A	510	HEM	CAD-C3D-C4D	4.57	128.60	112.47
6	A	511	HEM	CMB-C2B-C3B	4.60	128.02	116.53
6	A	509	HEM	CAD-C3D-C2D	5.19	128.13	113.22
6	A	512	HEM	CAD-C3D-C2D	5.78	129.83	113.22

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	508	HEM	1	0
6	A	509	HEM	2	0
6	A	512	HEM	2	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.