



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 03:58 pm GMT

PDB ID : 1Q9I  
Title : The A251C:S430C double mutant of flavocytochrome c3 from *Shewanella frigidimarina*  
Authors : Rothery, E.L.; Mowat, C.G.; Miles, C.S.; Walkinshaw, M.D.; Reid, G.A.; Chapman, S.K.  
Deposited on : 2003-08-25  
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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

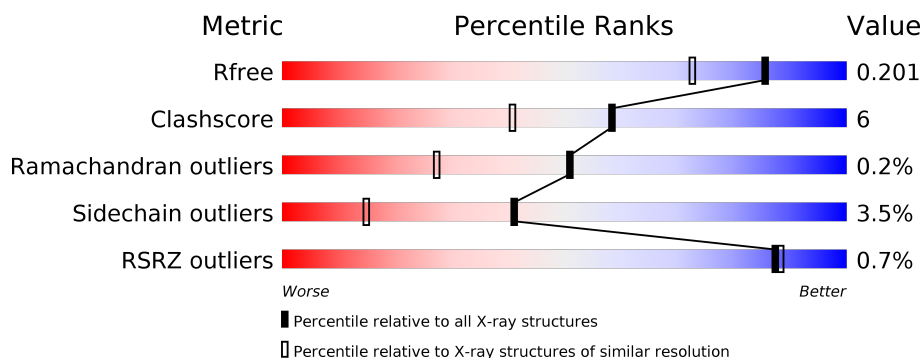
# 1 Overall quality at a glance

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))
$R_{free}$	100719	2696 (1.60-1.60)
Clashscore	112137	2967 (1.60-1.60)
Ramachandran outliers	110173	2887 (1.60-1.60)
Sidechain outliers	110143	2886 (1.60-1.60)
RSRZ outliers	101464	2714 (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.

Mol	Chain	Length	Quality of chain
1	A	571	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: -10px; left: 0; right: 0;">%</div> <div style="position: absolute; top: 10px; left: 0; right: 0;">91% 6% ..</div> </div> </div>

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 5653 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 flavocytochrome c3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	568	Total	C	N	O	S	0	0	0
			4190	2600	739	824	27			

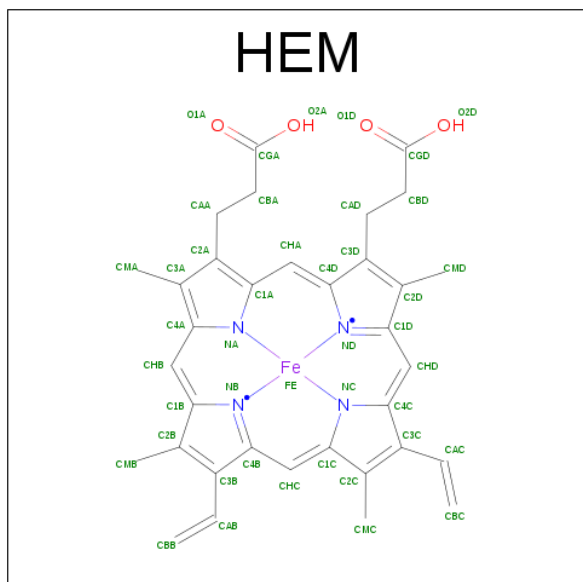
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	251	CYS	ALA	ENGINEERED	UNP Q02469
A	430	CYS	SER	ENGINEERED	UNP Q02469

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

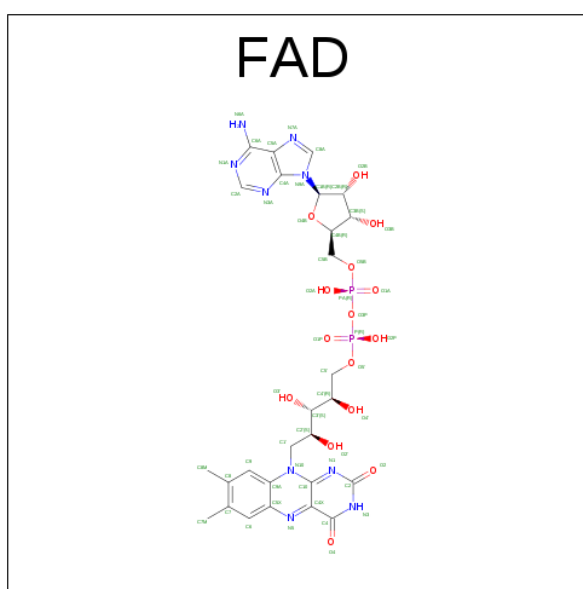
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Na	0	0
			1	1		

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



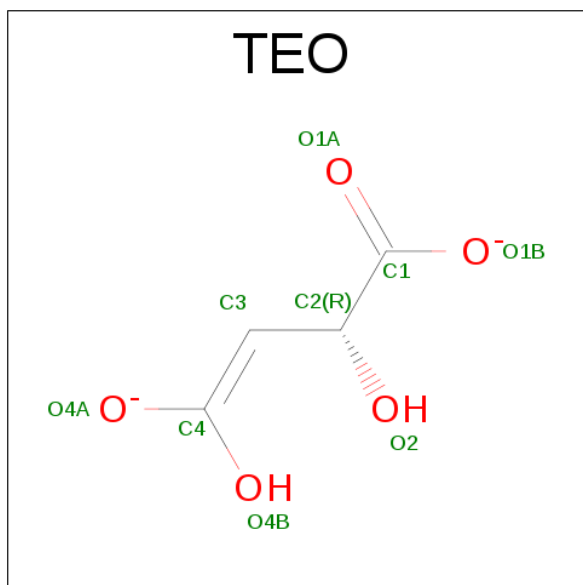
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	Fe	N	O	
			43	34	1	4	4	
3	A	1	Total	C	Fe	N	O	
			43	34	1	4	4	
3	A	1	Total	C	Fe	N	O	
			43	34	1	4	4	
3	A	1	Total	C	Fe	N	O	
			43	34	1	4	4	

- Molecule 4 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula:  $C_{27}H_{33}N_9O_{15}P_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	
			53	27	9	15	2	

- Molecule 5 is MALATE LIKE INTERMEDIATE (three-letter code: TEO) (formula:  $C_4H_4O_5$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			9	4	5		

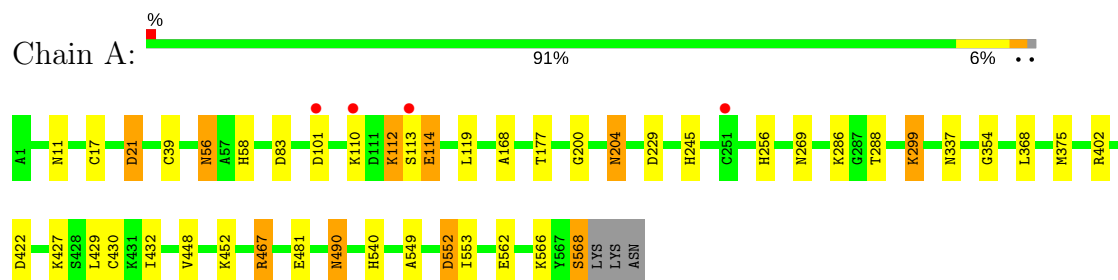
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1228	Total	O	0	0
			1228	1228		

### 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 the various outlier classes 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.

- Molecule 1: flavocytochrome c3



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	45.29Å 91.89Å 78.31Å 90.00° 91.50° 90.00°	Depositor
Resolution (Å)	24.00 – 1.60 22.64 – 1.60	Depositor EDS
% Data completeness (in resolution range)	92.2 (24.00-1.60) 92.2 (22.64-1.60)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.00 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, $R_{free}$	0.155 , 0.197 0.158 , 0.201	Depositor DCC
$R_{free}$ test set	3884 reflections (5.25%)	DCC
Wilson B-factor (Å <sup>2</sup> )	14.9	Xtriage
Anisotropy	0.034	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 48.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.023 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	5653	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.55% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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, NA, FAD, TEO

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	0.59	0/4260	0.81	5/5764 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	552	ASP	CB-CG-OD1	7.06	124.65	118.30
1	A	101	ASP	CB-CG-OD2	5.82	123.54	118.30
1	A	422	ASP	CB-CG-OD2	5.70	123.43	118.30
1	A	83	ASP	CB-CG-OD1	5.45	123.21	118.30
1	A	21	ASP	CB-CG-OD2	5.21	122.99	118.30

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	4190	0	4097	47	1
2	A	1	0	0	0	0
3	A	172	0	120	9	0
4	A	53	0	30	4	0
5	A	9	0	2	1	0
6	A	1228	0	0	14	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5653	0	4249	47	1

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 (47) 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:39:CYS:SG	3:A:802:HEM:CAC	2.16	1.34
1:A:17:CYS:SG	3:A:801:HEM:HAC	1.66	1.31
1:A:39:CYS:SG	3:A:802:HEM:HAC	1.81	1.18
1:A:39:CYS:HG	3:A:802:HEM:CAC	1.62	1.06
1:A:368:LEU:HB2	6:A:1417:HOH:O	1.60	0.99
1:A:229:ASP:H	1:A:256:HIS:HE1	1.04	0.94
1:A:204:ASN:H	1:A:204:ASN:HD22	1.18	0.88
1:A:229:ASP:H	1:A:256:HIS:CE1	1.92	0.86
1:A:481:GLU:CG	6:A:1385:HOH:O	2.28	0.81
1:A:299:LYS:HB2	6:A:1432:HOH:O	1.81	0.80
1:A:200:GLY:HA3	1:A:204:ASN:HD21	1.49	0.78
1:A:402:ARG:HH22	5:A:9806:TEO:C3	1.97	0.77
1:A:17:CYS:HG	3:A:801:HEM:HAC	1.51	0.75
1:A:229:ASP:N	1:A:256:HIS:HE1	1.85	0.71
1:A:39:CYS:SG	3:A:802:HEM:C3C	2.84	0.70
1:A:17:CYS:SG	3:A:801:HEM:C3C	2.85	0.69
1:A:204:ASN:ND2	1:A:204:ASN:H	1.91	0.66
1:A:177:THR:OG1	1:A:245:HIS:HE1	1.78	0.66
1:A:375:MET:HE2	4:A:9805:FAD:H6	1.83	0.60
1:A:17:CYS:SG	3:A:801:HEM:CBC	2.85	0.58
1:A:566:LYS:HG3	6:A:1357:HOH:O	2.05	0.55
1:A:568:SER:HB2	6:A:1136:HOH:O	2.06	0.55
1:A:540:HIS:HE1	1:A:552:ASP:OD2	1.91	0.53
1:A:56:ASN:HD22	1:A:58:HIS:H	1.57	0.51
1:A:354:GLY:N	6:A:1182:HOH:O	2.43	0.51
1:A:540:HIS:HD2	6:A:9822:HOH:O	1.94	0.50
1:A:448:VAL:HG12	6:A:1751:HOH:O	2.11	0.49
1:A:113:SER:O	1:A:114:GLU:HB2	2.12	0.49
1:A:39:CYS:SG	3:A:802:HEM:CBC	2.94	0.49
1:A:168:ALA:HA	4:A:9805:FAD:N5	2.29	0.48
1:A:204:ASN:ND2	1:A:204:ASN:N	2.62	0.47
1:A:467:ARG:NH2	6:A:1894:HOH:O	2.49	0.46
1:A:375:MET:HE2	4:A:9805:FAD:C6	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:427:LYS:NZ	6:A:1292:HOH:O	2.36	0.46
1:A:119:LEU:HD13	6:A:2039:HOH:O	2.16	0.45
1:A:56:ASN:ND2	1:A:58:HIS:H	2.14	0.45
1:A:490:ASN:C	1:A:490:ASN:HD22	2.21	0.44
1:A:549:ALA:O	1:A:553:ILE:HG23	2.18	0.43
1:A:562:GLU:O	1:A:566:LYS:HG3	2.19	0.42
1:A:427:LYS:CD	6:A:1292:HOH:O	2.67	0.42
1:A:113:SER:O	1:A:114:GLU:CB	2.68	0.41
1:A:429:LEU:HD23	1:A:432:ILE:HG13	2.02	0.41
1:A:427:LYS:HD3	6:A:1292:HOH:O	2.20	0.41
1:A:168:ALA:HA	4:A:9805:FAD:C5X	2.51	0.40
1:A:177:THR:OG1	1:A:245:HIS:CE1	2.67	0.40
1:A:568:SER:HB3	6:A:1248:HOH:O	2.20	0.40
1:A:112:LYS:HZ3	1:A:112:LYS:HA	1.87	0.40

All (1) 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:A:286:LYS:NZ	6:A:1145:HOH:O[2_646]	1.36	0.84

## 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	566/571 (99%)	547 (97%)	18 (3%)	1 (0%)	51 27

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	114	GLU

### 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	434/446 (97%)	419 (96%)	15 (4%)	41	15

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	21	ASP
1	A	56	ASN
1	A	110	LYS
1	A	112	LYS
1	A	204	ASN
1	A	269	ASN
1	A	288	THR
1	A	299	LYS
1	A	337	ASN
1	A	430	CYS
1	A	452	LYS
1	A	467	ARG
1	A	490	ASN
1	A	568	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	35	GLN
1	A	56	ASN
1	A	91	ASN
1	A	116	GLN
1	A	204	ASN
1	A	245	HIS
1	A	256	HIS
1	A	490	ASN
1	A	540	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 ⓘ

Of 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 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$
3	HEM	A	801	1	28,50,50	2.08	8 (28%)	17,82,82	1.61	5 (29%)
3	HEM	A	802	1	28,50,50	2.15	6 (21%)	17,82,82	1.46	4 (23%)
3	HEM	A	803	1	28,50,50	2.16	9 (32%)	17,82,82	2.28	7 (41%)
3	HEM	A	804	1	28,50,50	2.34	10 (35%)	17,82,82	1.58	4 (23%)
4	FAD	A	9805	-	51,58,58	2.45	21 (41%)	54,89,89	2.69	23 (42%)
5	TEO	A	9806	-	1,8,8	1.54	0	0,10,10	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
3	HEM	A	801	1	-	0/6/54/54	0/0/8/8
3	HEM	A	802	1	-	0/6/54/54	0/0/8/8
3	HEM	A	803	1	-	0/6/54/54	0/0/8/8

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	HEM	A	804	1	-	0/6/54/54	0/0/8/8
4	FAD	A	9805	-	-	0/28/50/50	0/6/6/6
5	TEO	A	9806	-	-	0/1/8/8	0/0/0/0

All (54) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	802	HEM	C3B-C2B	-5.06	1.33	1.40
3	A	804	HEM	C3C-C2C	-4.78	1.34	1.40
4	A	9805	FAD	C10-N1	-4.55	1.26	1.33
3	A	801	HEM	C3B-C2B	-4.43	1.34	1.40
4	A	9805	FAD	C1'-N10	-4.42	1.43	1.48
3	A	803	HEM	C3C-C2C	-4.21	1.34	1.40
3	A	804	HEM	C3B-C2B	-4.10	1.34	1.40
3	A	803	HEM	C3B-C2B	-4.08	1.35	1.40
3	A	801	HEM	C3C-C2C	-3.93	1.35	1.40
3	A	802	HEM	C3C-C2C	-3.70	1.35	1.40
4	A	9805	FAD	O4-C4	-3.38	1.16	1.24
4	A	9805	FAD	O4B-C1B	-2.42	1.37	1.41
4	A	9805	FAD	PA-O2A	-2.32	1.43	1.55
4	A	9805	FAD	P-O1P	-2.28	1.42	1.50
4	A	9805	FAD	C5'-C4'	-2.22	1.48	1.51
4	A	9805	FAD	O3'-C3'	2.00	1.47	1.43
3	A	801	HEM	CMA-C3A	2.02	1.55	1.51
3	A	804	HEM	C1D-ND	2.10	1.40	1.36
3	A	803	HEM	C1D-ND	2.11	1.40	1.36
3	A	804	HEM	CAA-C2A	2.11	1.55	1.52
3	A	804	HEM	C4D-ND	2.12	1.39	1.36
3	A	801	HEM	C4D-ND	2.13	1.39	1.36
3	A	804	HEM	C1B-NB	2.16	1.39	1.36
3	A	802	HEM	CMB-C2B	2.16	1.56	1.51
4	A	9805	FAD	C2'-C3'	2.25	1.57	1.53
3	A	801	HEM	C1C-NC	2.28	1.39	1.36
4	A	9805	FAD	C4-N3	2.73	1.38	1.33
3	A	804	HEM	C4C-NC	2.75	1.40	1.36
4	A	9805	FAD	C4'-C3'	2.77	1.58	1.53
3	A	803	HEM	CAA-C2A	2.78	1.56	1.52
4	A	9805	FAD	C2-N1	2.88	1.43	1.38
4	A	9805	FAD	O2'-C2'	2.96	1.49	1.43
4	A	9805	FAD	O3B-C3B	2.98	1.49	1.43
3	A	803	HEM	C3B-CAB	2.99	1.53	1.47
3	A	803	HEM	C1B-NB	3.06	1.40	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	803	HEM	C4D-ND	3.12	1.40	1.36
3	A	801	HEM	C3B-CAB	3.17	1.54	1.47
4	A	9805	FAD	C8-C7	3.21	1.49	1.41
3	A	802	HEM	C3B-CAB	3.27	1.54	1.47
3	A	803	HEM	C3C-CAC	3.32	1.54	1.47
4	A	9805	FAD	C2-N3	3.53	1.45	1.38
3	A	804	HEM	C3C-CAC	3.99	1.55	1.47
3	A	801	HEM	C3D-C2D	4.12	1.49	1.37
3	A	802	HEM	C3C-CAC	4.14	1.56	1.47
3	A	804	HEM	C3B-CAB	4.25	1.56	1.47
3	A	801	HEM	C3C-CAC	4.28	1.56	1.47
4	A	9805	FAD	C9A-N10	4.63	1.44	1.38
3	A	803	HEM	C3D-C2D	4.68	1.51	1.37
4	A	9805	FAD	C5X-N5	4.77	1.42	1.35
3	A	804	HEM	C3D-C2D	4.77	1.51	1.37
3	A	802	HEM	C3D-C2D	4.81	1.51	1.37
4	A	9805	FAD	C2A-N3A	5.11	1.40	1.32
4	A	9805	FAD	C6-C5X	5.14	1.49	1.41
4	A	9805	FAD	C4-C4X	6.24	1.53	1.41

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	9805	FAD	C4X-C4-N3	-5.97	114.99	123.48
3	A	803	HEM	CBD-CAD-C3D	-5.31	102.33	112.47
4	A	9805	FAD	O5'-C5'-C4'	-5.31	95.18	109.36
4	A	9805	FAD	N3A-C2A-N1A	-4.24	125.16	128.86
4	A	9805	FAD	O4'-C4'-C3'	-4.15	98.79	109.09
4	A	9805	FAD	C9A-C5X-N5	-3.98	116.32	122.24
4	A	9805	FAD	C8M-C8-C9	-3.54	111.47	120.34
3	A	804	HEM	CMA-C3A-C4A	-3.49	123.09	128.46
4	A	9805	FAD	C4X-C10-N10	-3.36	118.19	120.52
4	A	9805	FAD	O3'-C3'-C2'	-3.23	100.81	108.82
3	A	803	HEM	CMA-C3A-C4A	-2.98	123.88	128.46
4	A	9805	FAD	O3B-C3B-C2B	-2.93	102.46	111.83
3	A	803	HEM	C1D-C2D-C3D	-2.91	104.97	107.00
4	A	9805	FAD	C4A-C5A-N7A	-2.75	106.76	109.41
3	A	801	HEM	CBD-CAD-C3D	-2.53	107.64	112.47
3	A	802	HEM	CMA-C3A-C4A	-2.53	124.58	128.46
3	A	801	HEM	CBA-CAA-C2A	-2.52	107.66	112.48
3	A	802	HEM	C3B-C4B-NB	-2.38	106.13	109.21
3	A	804	HEM	CBA-CAA-C2A	-2.35	107.99	112.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	801	HEM	CMA-C3A-C4A	-2.35	124.85	128.46
3	A	802	HEM	CBA-CAA-C2A	-2.29	108.11	112.48
3	A	803	HEM	C3C-C4C-NC	-2.25	106.69	110.94
4	A	9805	FAD	C6-C7-C8	-2.14	116.14	119.95
4	A	9805	FAD	C10-C4X-N5	2.08	122.98	120.59
3	A	803	HEM	CMC-C2C-C3C	2.22	129.01	124.89
3	A	801	HEM	C4A-C3A-C2A	2.25	108.56	107.00
4	A	9805	FAD	N6A-C6A-N1A	2.28	123.29	118.77
3	A	801	HEM	CMB-C2B-C3B	2.33	129.21	124.89
3	A	804	HEM	CMB-C2B-C3B	2.34	129.24	124.89
3	A	804	HEM	C4A-C3A-C2A	2.34	108.63	107.00
4	A	9805	FAD	O2'-C2'-C1'	2.37	115.27	109.79
3	A	802	HEM	CMC-C2C-C3C	2.46	129.46	124.89
3	A	803	HEM	C4C-C3C-C2C	2.74	108.81	106.90
4	A	9805	FAD	C5X-C9A-N10	2.95	119.85	117.66
4	A	9805	FAD	C1B-N9A-C4A	3.04	131.89	126.64
4	A	9805	FAD	C9-C8-C7	3.10	125.49	119.95
4	A	9805	FAD	C2A-N1A-C6A	3.13	124.24	118.77
4	A	9805	FAD	C6-C5X-C9A	3.22	123.18	119.00
3	A	803	HEM	C4A-C3A-C2A	3.59	109.50	107.00
4	A	9805	FAD	O2B-C2B-C3B	4.32	125.66	111.83
4	A	9805	FAD	C4X-N5-C5X	5.21	122.27	116.76
4	A	9805	FAD	O4'-C4'-C5'	5.24	121.67	110.00
4	A	9805	FAD	C4-N3-C2	7.11	121.38	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	801	HEM	4	0
3	A	802	HEM	5	0
4	A	9805	FAD	4	0
5	A	9806	TEO	1	0

## 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	568/571 (99%)	-0.34	4 (0%) 87 88	8, 15, 25, 37	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	113	SER	3.8
1	A	101	ASP	3.2
1	A	110	LYS	2.9
1	A	251	CYS	2.1

### 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, 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	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
5	TEO	A	9806	9/9	0.94	0.10	1.45	13,17,22,25	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
3	HEM	A	802	43/43	0.95	0.10	1.09	13,18,27,32	0
3	HEM	A	803	43/43	0.97	0.10	0.44	11,16,29,36	0
3	HEM	A	801	43/43	0.97	0.09	0.33	11,16,21,23	0
3	HEM	A	804	43/43	0.98	0.08	-0.00	9,11,17,19	0
4	FAD	A	9805	53/53	0.98	0.06	-0.67	7,10,15,16	0
2	NA	A	9810	1/1	1.00	0.04	-4.72	9,9,9,9	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.