



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:07 PM GMT

PDB ID : 1S5Y  
Title : The crystal structure of Trematomus bernacchii hemoglobin oxidized by ferri-cyanide  
Authors : Vitagliano, L.; Bonomi, G.; Riccio, A.; di Prisco, G.; Smulevich, G.; Maz-zarella, L.  
Deposited on : 2004-01-22  
Resolution : 2.50 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

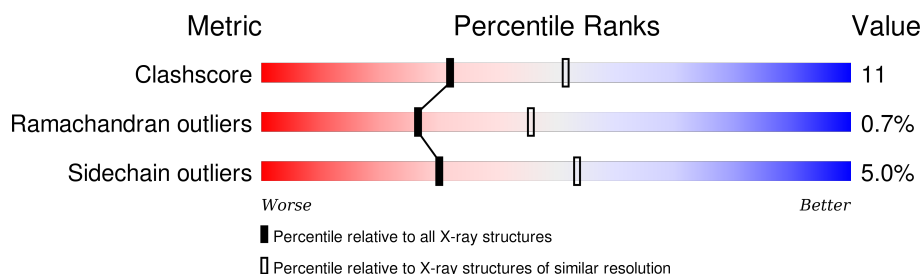
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)

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	143	
1	C	143	
2	B	146	
2	D	146	

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4562 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.

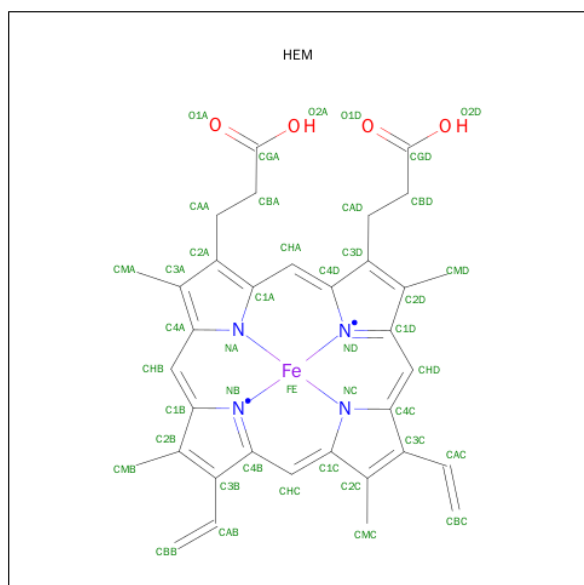
- Molecule 1 is a protein called Hemoglobin alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	143	Total	C	N	O	S	0	0	0
			1104	710	190	199	5			
1	C	143	Total	C	N	O	S	0	0	0
			1104	710	190	199	5			

- Molecule 2 is a protein called Hemoglobin beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	136	Total	C	N	O	S	0	0	0
			1050	669	182	194	5			
2	D	136	Total	C	N	O	S	0	0	0
			1050	669	182	194	5			

- 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	B	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	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 4 is water.

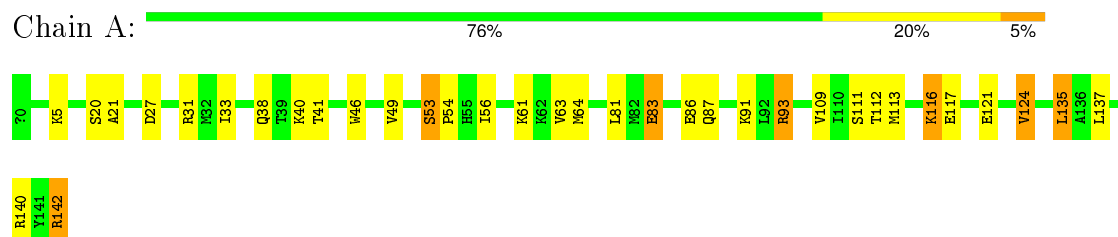
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	29	Total 29	O 29	0	0
4	B	17	Total 17	O 17	0	0
4	C	17	Total 17	O 17	0	0
4	D	19	Total 19	O 19	0	0

### 3 Residue-property plots

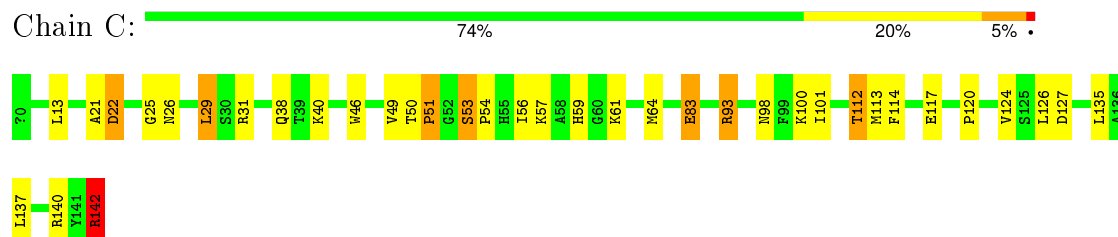
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

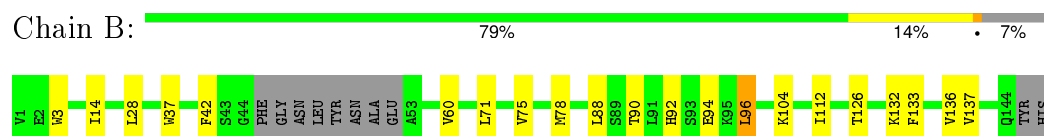
- Molecule 1: Hemoglobin alpha chain



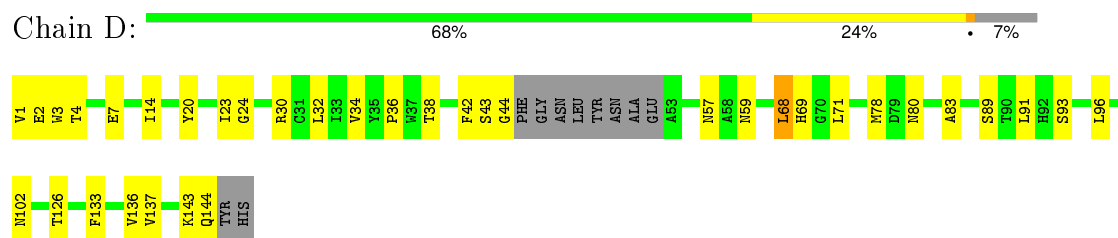
- Molecule 1: Hemoglobin alpha chain



- Molecule 2: Hemoglobin beta chain



- Molecule 2: Hemoglobin beta chain



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	55.99 Å 62.98 Å 63.50 Å 77.06° 69.79° 84.17°	Depositor
Resolution (Å)	25.00 – 2.50	Depositor
% Data completeness (in resolution range)	(Not available) (25.00-2.50)	Depositor
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	CNS	Depositor
R, $R_{free}$	0.199 , 0.247	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4562	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

## 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, ACE

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	1.22	1/1127 (0.1%)	0.64	0/1523
1	C	0.47	1/1127 (0.1%)	0.63	0/1523
2	B	0.38	0/1071	0.57	0/1449
2	D	0.39	0/1071	0.59	0/1449
All	All	0.71	2/4396 (0.0%)	0.61	0/5944

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	142	ARG	C-OXT	-38.14	0.50	1.23
1	C	142	ARG	C-OXT	6.90	1.36	1.23

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	1104	0	1136	35	0
1	C	1104	0	1136	27	0
2	B	1050	0	1047	15	0
2	D	1050	0	1047	25	0
3	A	43	0	30	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	43	0	30	0	0
3	C	43	0	30	0	0
3	D	43	0	30	0	0
4	A	29	0	0	1	0
4	B	17	0	0	1	0
4	C	17	0	0	0	0
4	D	19	0	0	0	0
All	All	4562	0	4486	95	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 (95) 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:93:ARG:HH11	1:A:93:ARG:HB3	1.14	1.10
2:D:68:LEU:HD12	2:D:71:LEU:HD12	1.58	0.84
1:C:93:ARG:HB3	1:C:93:ARG:HH11	1.43	0.83
1:A:135:LEU:HD21	1:C:135:LEU:HD21	1.62	0.80
1:A:86:GLU:OE1	1:A:140:ARG:HD3	1.85	0.76
1:C:64:MET:HE2	1:C:64:MET:HA	1.68	0.75
1:A:93:ARG:NH1	1:A:93:ARG:HB3	1.98	0.71
1:C:29:LEU:HD11	1:C:59:HIS:HD2	1.56	0.69
2:B:92:HIS:HA	2:B:96:LEU:HB2	1.76	0.67
1:A:93:ARG:HH11	1:A:93:ARG:CB	2.00	0.67
2:D:143:LYS:HE3	2:D:144:GLN:HE21	1.59	0.66
2:D:143:LYS:HE3	2:D:144:GLN:NE2	2.11	0.66
2:D:38:THR:HG22	2:D:102:ASN:OD1	1.99	0.63
1:C:21:ALA:O	1:C:64:MET:HG3	1.99	0.62
1:C:114:PHE:HB3	1:C:117:GLU:HB2	1.82	0.62
1:A:33:ILE:HG23	1:A:40:LYS:HG2	1.81	0.62
1:A:27:ASP:OD1	1:A:31:ARG:NE	2.32	0.60
1:A:87:GLN:HG3	1:A:91:LYS:HD3	1.82	0.60
1:A:49:VAL:O	1:A:49:VAL:HG12	2.02	0.58
1:A:87:GLN:HE21	1:A:91:LYS:HD3	1.68	0.58
1:C:124:VAL:HG22	2:D:34:VAL:HA	1.85	0.58
1:C:98:ASN:HD22	1:C:101:ILE:HD12	1.69	0.57
1:A:116:LYS:NZ	1:A:116:LYS:HB2	2.20	0.57
2:D:91:LEU:O	2:D:96:LEU:HD23	2.05	0.57
2:D:133:PHE:O	2:D:137:VAL:HG23	2.04	0.57
2:B:104:LYS:HD3	2:B:104:LYS:C	2.26	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:49:VAL:HG12	1:C:49:VAL:O	2.06	0.56
1:C:120:PRO:HD3	2:D:30:ARG:NH2	2.22	0.55
2:B:104:LYS:HD3	2:B:104:LYS:O	2.06	0.54
1:A:63:VAL:HG12	1:A:64:MET:CE	2.39	0.53
1:A:87:GLN:HE21	1:A:91:LYS:CD	2.21	0.52
2:B:133:PHE:O	2:B:137:VAL:HG23	2.09	0.52
2:D:89:SER:OG	2:D:144:GLN:HB2	2.09	0.52
2:B:42:PHE:HA	4:B:1041:HOH:O	2.10	0.52
1:C:22:ASP:OD2	1:C:61:LYS:HG3	2.10	0.51
1:A:121:GLU:O	1:A:124:VAL:HG23	2.09	0.51
2:D:23:ILE:HG22	2:D:68:LEU:HD21	1.93	0.51
2:B:28:LEU:HD23	2:B:60:VAL:O	2.11	0.50
1:A:21:ALA:O	1:A:64:MET:HG3	2.12	0.50
2:B:37:TRP:HZ3	1:C:142:ARG:HG2	1.75	0.50
1:A:46:TRP:HB2	1:A:49:VAL:CG2	2.42	0.50
1:C:26:ASN:HD22	1:C:57:LYS:HA	1.78	0.49
1:A:81:LEU:HA	1:A:83:GLU:OE2	2.12	0.49
1:A:46:TRP:HZ3	1:A:56:ILE:HD13	1.77	0.48
1:C:46:TRP:HZ3	1:C:56:ILE:HD13	1.77	0.48
2:D:24:GLY:HA2	2:D:68:LEU:HD22	1.96	0.47
1:A:111:SER:OG	2:B:112:ILE:HG23	2.13	0.47
2:B:3:TRP:CE2	2:B:78:MET:HE2	2.50	0.47
2:D:80:ASN:ND2	2:D:83:ALA:HB3	2.29	0.47
2:D:32:LEU:HD23	2:D:38:THR:OG1	2.15	0.46
1:A:46:TRP:CZ3	1:A:56:ILE:HD13	2.51	0.45
1:C:61:LYS:HE3	1:C:61:LYS:HB2	1.60	0.45
1:A:83:GLU:HG3	1:A:83:GLU:H	1.52	0.45
1:C:25:GLY:O	1:C:29:LEU:HB2	2.16	0.45
2:D:23:ILE:CG2	2:D:68:LEU:HD21	2.46	0.45
2:B:90:THR:O	2:B:94:GLU:HB2	2.18	0.44
1:A:53:SER:HA	1:A:54:PRO:HD3	1.78	0.44
1:A:142:ARG:NH2	1:C:127:ASP:OD2	2.29	0.44
1:A:5:LYS:HB2	1:A:5:LYS:HE3	1.83	0.44
2:B:71:LEU:O	2:B:75:VAL:HG23	2.17	0.44
1:C:31:ARG:HG2	1:C:31:ARG:HH11	1.82	0.44
2:D:96:LEU:N	2:D:96:LEU:HD22	2.32	0.44
2:D:42:PHE:C	2:D:44:GLY:H	2.20	0.43
1:A:49:VAL:HG13	1:A:56:ILE:HD11	1.99	0.43
2:D:3:TRP:CE2	2:D:78:MET:HE2	2.54	0.43
1:A:116:LYS:HZ3	1:A:116:LYS:HB2	1.84	0.43
1:C:29:LEU:HD11	1:C:59:HIS:CD2	2.45	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:VAL:HG12	1:A:64:MET:HE2	2.01	0.42
1:A:135:LEU:HD21	1:C:135:LEU:CD2	2.41	0.42
2:D:2:GLU:HA	2:D:2:GLU:OE2	2.19	0.42
2:D:1:VAL:HG11	2:D:136:VAL:HG22	2.01	0.42
2:D:20:TYR:CD1	2:D:69:HIS:CD2	3.08	0.42
2:B:132:LYS:O	2:B:136:VAL:HG23	2.19	0.42
1:C:112:THR:HG22	1:C:113:MET:HG3	2.02	0.42
1:A:61:LYS:HB2	1:A:61:LYS:HE3	1.68	0.42
1:C:93:ARG:CB	1:C:93:ARG:HH11	2.24	0.41
2:D:34:VAL:C	2:D:36:PRO:HD3	2.41	0.41
2:D:4:THR:OG1	2:D:7:GLU:HG3	2.20	0.41
2:B:88:LEU:HD23	2:B:88:LEU:HA	1.93	0.41
1:C:53:SER:HA	1:C:54:PRO:HD3	1.84	0.41
2:B:14:ILE:HD11	2:B:126:THR:CG2	2.50	0.41
2:B:14:ILE:HD11	2:B:126:THR:HG22	2.01	0.41
1:A:63:VAL:HG12	1:A:64:MET:HE3	2.02	0.41
1:C:50:THR:O	1:C:53:SER:HB3	2.20	0.41
1:A:20:SER:HA	4:A:1057:HOH:O	2.20	0.41
1:A:109:VAL:CG1	1:A:113:MET:CE	2.99	0.41
1:C:83:GLU:N	1:C:83:GLU:OE2	2.53	0.41
1:A:117:GLU:N	1:A:117:GLU:OE1	2.53	0.41
1:A:87:GLN:HE21	1:A:91:LYS:CE	2.34	0.41
2:D:14:ILE:HD11	2:D:126:THR:CG2	2.50	0.41
2:D:14:ILE:HD11	2:D:126:THR:HG23	2.03	0.41
1:C:50:THR:HB	1:C:51:PRO:HD2	2.02	0.40
2:D:57:ASN:OD1	2:D:59:ASN:N	2.55	0.40
1:C:13:LEU:CD2	1:C:126:LEU:HD22	2.51	0.40
1:A:87:GLN:HE21	1:A:91:LYS:HE2	1.86	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	141/143 (99%)	132 (94%)	8 (6%)	1 (1%)	26	46
1	C	141/143 (99%)	133 (94%)	7 (5%)	1 (1%)	26	46
2	B	132/146 (90%)	127 (96%)	5 (4%)	0	100	100
2	D	132/146 (90%)	122 (92%)	8 (6%)	2 (2%)	13	22
All	All	546/578 (94%)	514 (94%)	28 (5%)	4 (1%)	26	46

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	53	SER
2	D	43	SER
1	A	53	SER
2	D	93	SER

### 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	119/119 (100%)	110 (92%)	9 (8%)	16	30
1	C	119/119 (100%)	107 (90%)	12 (10%)	9	17
2	B	112/120 (93%)	111 (99%)	1 (1%)	84	95
2	D	112/120 (93%)	111 (99%)	1 (1%)	84	95
All	All	462/478 (97%)	439 (95%)	23 (5%)	30	53

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	GLN
1	A	41	THR
1	A	83	GLU
1	A	93	ARG
1	A	112	THR
1	A	116	LYS
1	A	124	VAL

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Mol	Chain	Res	Type
1	A	135	LEU
1	A	137	LEU
2	B	96	LEU
1	C	22	ASP
1	C	29	LEU
1	C	38	GLN
1	C	40	LYS
1	C	51	PRO
1	C	83	GLU
1	C	93	ARG
1	C	100	LYS
1	C	112	THR
1	C	137	LEU
1	C	140	ARG
1	C	142	ARG
2	D	68	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	ASN
1	A	87	GLN
1	A	98	ASN
2	B	17	HIS
2	B	69	HIS
1	C	26	ASN
1	C	87	GLN
1	C	98	ASN
2	D	41	HIS
2	D	69	HIS
2	D	80	ASN
2	D	144	GLN

### 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 [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

4 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
3	HEM	A	200	1,4	30,50,50	2.17	7 (23%)	24,82,82	3.24	16 (66%)
3	HEM	B	400	2	30,50,50	2.59	10 (33%)	24,82,82	3.42	10 (41%)
3	HEM	C	600	1,4	30,50,50	2.24	7 (23%)	24,82,82	3.62	15 (62%)
3	HEM	D	800	2	30,50,50	2.38	7 (23%)	24,82,82	3.44	13 (54%)

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	200	1,4	-	0/10/54/54	0/0/8/8
3	HEM	B	400	2	-	0/10/54/54	0/0/8/8
3	HEM	C	600	1,4	-	0/10/54/54	0/0/8/8
3	HEM	D	800	2	-	0/10/54/54	0/0/8/8

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	400	HEM	C3B-C4B	-7.35	1.45	1.51
3	D	800	HEM	C2D-C3D	-6.22	1.35	1.54
3	C	600	HEM	C2D-C3D	-6.16	1.36	1.54
3	A	200	HEM	C2D-C3D	-5.90	1.36	1.54
3	B	400	HEM	C2D-C3D	-5.72	1.37	1.54
3	D	800	HEM	C3B-C4B	-5.01	1.47	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	400	HEM	C2C-C1C	-3.77	1.45	1.52
3	A	200	HEM	C2C-C1C	-3.05	1.46	1.52
3	D	800	HEM	C2C-C1C	-3.02	1.46	1.52
3	D	800	HEM	C3D-C4D	-3.00	1.47	1.51
3	A	200	HEM	C3B-C4B	-2.99	1.49	1.51
3	B	400	HEM	C3D-C4D	-2.90	1.47	1.51
3	C	600	HEM	C3B-C4B	-2.69	1.49	1.51
3	C	600	HEM	C2D-C1D	-2.20	1.44	1.51
3	C	600	HEM	FE-NC	2.18	2.04	1.95
3	B	400	HEM	FE-NC	2.18	2.04	1.95
3	B	400	HEM	CHD-C4C	2.42	1.42	1.36
3	B	400	HEM	C4C-NC	2.52	1.39	1.36
3	A	200	HEM	C1C-NC	3.08	1.39	1.36
3	A	200	HEM	C3C-CAC	3.39	1.57	1.51
3	B	400	HEM	CBB-CAB	3.96	1.52	1.29
3	C	600	HEM	CBB-CAB	4.16	1.53	1.29
3	D	800	HEM	C1C-NC	4.22	1.41	1.36
3	D	800	HEM	CBB-CAB	4.35	1.54	1.29
3	C	600	HEM	CBC-CAC	4.59	1.55	1.29
3	A	200	HEM	CBC-CAC	4.60	1.55	1.29
3	A	200	HEM	CBB-CAB	4.62	1.56	1.29
3	D	800	HEM	CBC-CAC	4.63	1.56	1.29
3	B	400	HEM	CBC-CAC	4.64	1.56	1.29
3	B	400	HEM	C1C-NC	4.88	1.42	1.36
3	C	600	HEM	C1C-NC	5.85	1.43	1.36

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	400	HEM	C3C-CAC-CBC	-10.47	108.39	124.46
3	D	800	HEM	C3C-CAC-CBC	-10.27	108.70	124.46
3	C	600	HEM	CAA-CBA-CGA	-4.67	104.18	112.75
3	C	600	HEM	C3B-CAB-CBB	-3.36	119.30	124.46
3	D	800	HEM	CBD-CAD-C3D	-3.08	104.58	113.55
3	B	400	HEM	CMA-C3A-C4A	-2.94	123.49	128.36
3	A	200	HEM	CMA-C3A-C4A	-2.88	123.61	128.36
3	C	600	HEM	CAA-C2A-C3A	-2.85	120.87	129.00
3	D	800	HEM	CMA-C3A-C4A	-2.70	123.90	128.36
3	A	200	HEM	C3B-CAB-CBB	-2.60	120.46	124.46
3	A	200	HEM	CHC-C4B-NB	-2.26	119.09	124.52
3	C	600	HEM	CMA-C3A-C4A	-2.20	124.72	128.36
3	D	800	HEM	C3B-CAB-CBB	-2.17	121.13	124.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	200	HEM	CHD-C1D-ND	-2.04	119.61	124.52
3	A	200	HEM	CMA-C3A-C2A	2.06	129.54	125.24
3	C	600	HEM	CMA-C3A-C2A	2.20	129.84	125.24
3	C	600	HEM	C2C-C1C-CHC	2.48	127.45	123.68
3	A	200	HEM	C3B-C4B-CHC	2.64	126.89	123.16
3	B	400	HEM	C2D-C3D-C4D	2.88	106.39	101.50
3	C	600	HEM	C2D-C3D-C4D	2.91	106.44	101.50
3	B	400	HEM	CAD-C3D-C4D	3.04	123.20	112.47
3	D	800	HEM	C1D-CHD-C4C	3.05	130.92	125.82
3	C	600	HEM	C1D-CHD-C4C	3.18	131.14	125.82
3	D	800	HEM	C4B-CHC-C1C	3.19	131.15	125.82
3	D	800	HEM	CAD-C3D-C4D	3.20	123.76	112.47
3	A	200	HEM	CAA-CBA-CGA	3.24	118.68	112.75
3	C	600	HEM	CMD-C2D-C3D	3.28	128.88	114.35
3	A	200	HEM	CMD-C2D-C3D	3.40	129.37	114.35
3	B	400	HEM	CMD-C2D-C3D	3.43	129.52	114.35
3	A	200	HEM	C4B-CHC-C1C	3.45	131.59	125.82
3	C	600	HEM	CAD-C3D-C4D	3.46	124.69	112.47
3	A	200	HEM	CAD-C3D-C4D	3.50	124.81	112.47
3	B	400	HEM	C4B-CHC-C1C	3.57	131.79	125.82
3	D	800	HEM	CMD-C2D-C3D	3.66	130.52	114.35
3	D	800	HEM	CMB-C2B-C3B	3.80	126.02	116.53
3	A	200	HEM	C2D-C3D-C4D	3.81	107.95	101.50
3	A	200	HEM	C1D-CHD-C4C	3.85	132.26	125.82
3	B	400	HEM	CMB-C2B-C3B	3.95	126.38	116.53
3	D	800	HEM	C2D-C3D-C4D	3.96	108.21	101.50
3	A	200	HEM	CMB-C2B-C3B	4.01	126.54	116.53
3	D	800	HEM	CAA-CBA-CGA	4.10	120.27	112.75
3	B	400	HEM	CBA-CAA-C2A	4.22	120.09	112.53
3	C	600	HEM	CAA-C2A-C1A	4.30	131.68	127.01
3	C	600	HEM	CMB-C2B-C3B	4.53	127.84	116.53
3	D	800	HEM	CMC-C2C-C3C	4.61	128.04	116.53
3	C	600	HEM	CMC-C2C-C3C	4.82	128.56	116.53
3	A	200	HEM	CAD-C3D-C2D	4.87	127.22	113.22
3	D	800	HEM	CAD-C3D-C2D	5.10	127.88	113.22
3	B	400	HEM	CMC-C2C-C3C	5.42	130.06	116.53
3	C	600	HEM	CAD-C3D-C2D	5.43	128.83	113.22
3	A	200	HEM	CMC-C2C-C3C	5.61	130.54	116.53
3	B	400	HEM	CAD-C3D-C2D	5.97	130.39	113.22
3	A	200	HEM	CBA-CAA-C2A	7.79	126.48	112.53
3	C	600	HEM	CBA-CAA-C2A	10.42	131.20	112.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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 ⓘ

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.