



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 04:46 pm GMT

PDB ID : 1CH4  
Title : MODULE-SUBSTITUTED CHIMERA HEMOGLOBIN BETA-ALPHA (F133V)  
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Deposited on : 1998-06-11  
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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

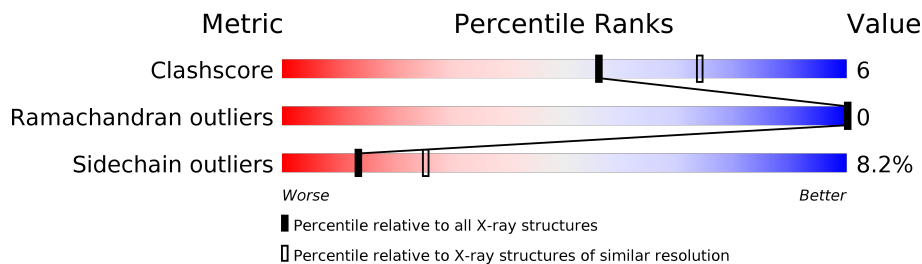
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 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	112137	4554 (2.50-2.50)
Ramachandran outliers	110173	4463 (2.50-2.50)
Sidechain outliers	110143	4465 (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	146	
1	B	146	
1	C	146	
1	D	146	

## 2 Entry composition [i](#)

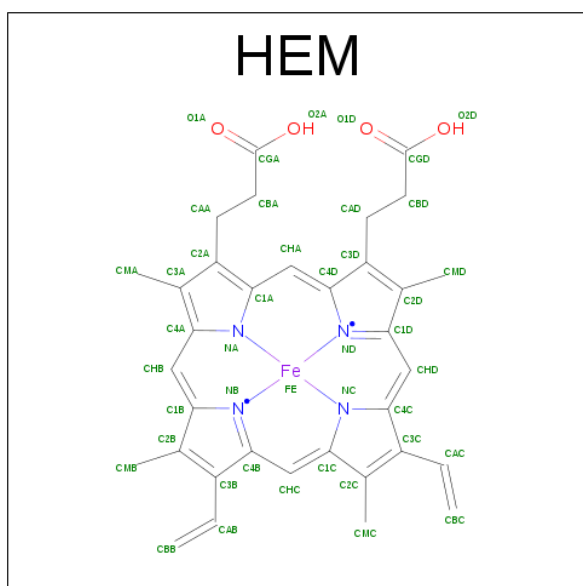
There are 4 unique types of molecules in this entry. The entry contains 4750 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 MODULE-SUBSTITUTED CHIMERA HEMOGLOBIN BETA-ALPHA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	146	Total	C	N	O	S	0	3	0
			1132	727	198	204	3			
1	B	146	Total	C	N	O	S	0	3	0
			1132	727	198	204	3			
1	C	146	Total	C	N	O	S	0	3	0
			1132	727	198	204	3			
1	D	146	Total	C	N	O	S	0	3	0
			1132	727	198	204	3			

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula:  $C_{34}H_{32}FeN_4O_4$ ).



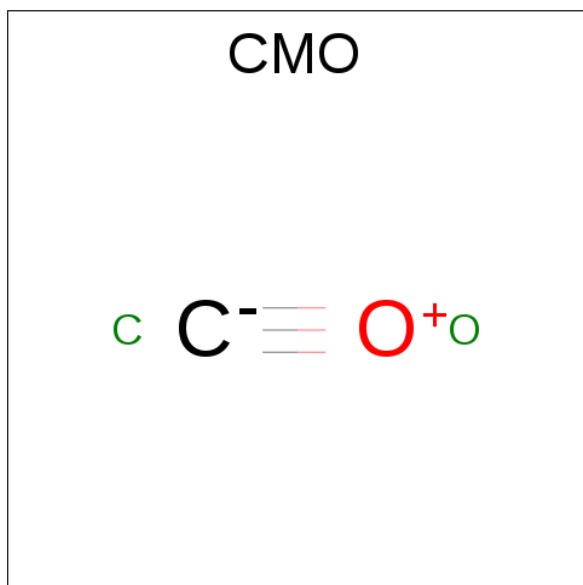
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

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

- Molecule 3 is CARBON MONOXIDE (three-letter code: CMO) (formula: CO).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	8	Total	O		
			8	8	0	0
4	B	8	Total	O		
			8	8	0	0

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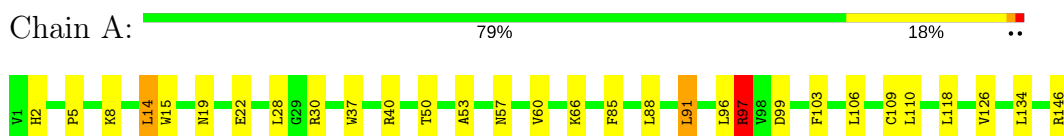
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	13	Total 13	O 13	0	0
4	D	13	Total 13	O 13	0	0

### 3 Residue-property plots [i](#)

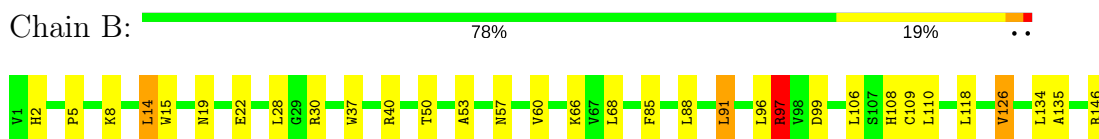
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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 ( $RSRZ > 2$ ). 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.

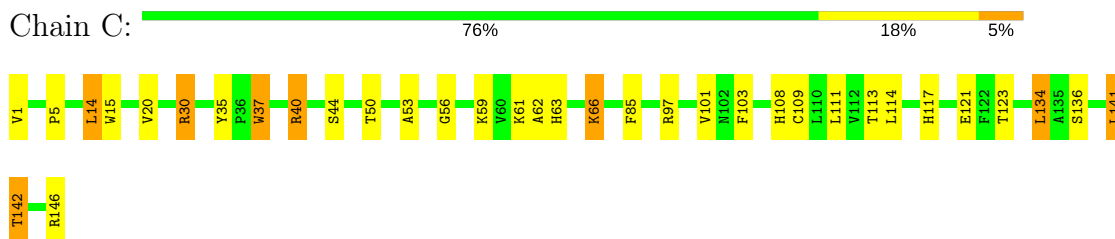
- Molecule 1: MODULE-SUBSTITUTED CHIMERA HEMOGLOBIN BETA-ALPHA



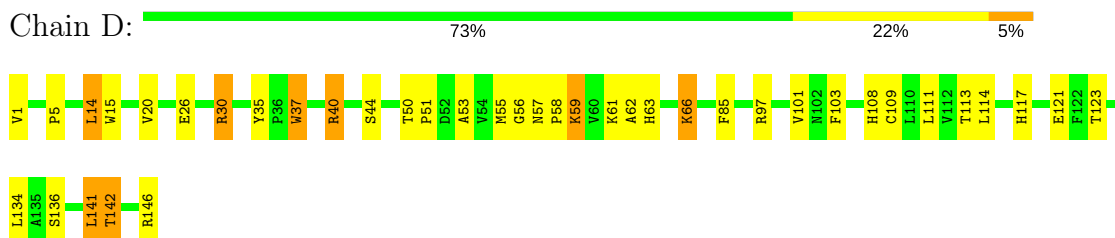
- Molecule 1: MODULE-SUBSTITUTED CHIMERA HEMOGLOBIN BETA-ALPHA



- Molecule 1: MODULE-SUBSTITUTED CHIMERA HEMOGLOBIN BETA-ALPHA



- Molecule 1: MODULE-SUBSTITUTED CHIMERA HEMOGLOBIN BETA-ALPHA



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	62.87Å 81.31Å 55.11Å 90.00° 90.96° 90.00°	Depositor
Resolution (Å)	8.00 – 2.50	Depositor
% Data completeness (in resolution range)	81.4 (8.00-2.50)	Depositor
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	X-PLOR 3.1	Depositor
R, $R_{free}$	0.188 , 0.244	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4750	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	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: CMO, HEM

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	0.82	0/1174	1.42	13/1594 (0.8%)
1	B	0.82	0/1174	1.42	13/1594 (0.8%)
1	C	0.84	0/1174	1.54	17/1594 (1.1%)
1	D	0.84	0/1174	1.54	17/1594 (1.1%)
All	All	0.83	0/4696	1.48	60/6376 (0.9%)

There are no bond length outliers.

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	97[A]	ARG	NE-CZ-NH2	-9.28	115.66	120.30
1	C	97[B]	ARG	NE-CZ-NH2	-9.28	115.66	120.30
1	D	97[A]	ARG	NE-CZ-NH2	-9.15	115.72	120.30
1	D	97[B]	ARG	NE-CZ-NH2	-9.15	115.72	120.30
1	D	15	TRP	CD1-CG-CD2	8.49	113.10	106.30
1	C	15	TRP	CD1-CG-CD2	8.47	113.08	106.30
1	B	15	TRP	CD1-CG-CD2	8.33	112.96	106.30
1	A	15	TRP	CD1-CG-CD2	8.27	112.92	106.30
1	A	37	TRP	CD1-CG-CD2	8.13	112.81	106.30
1	B	37	TRP	CD1-CG-CD2	8.09	112.77	106.30
1	D	37	TRP	CD1-CG-CD2	8.06	112.75	106.30
1	C	37	TRP	CD1-CG-CD2	8.02	112.72	106.30
1	C	15	TRP	CE2-CD2-CG	-7.95	100.94	107.30
1	D	15	TRP	CE2-CD2-CG	-7.95	100.94	107.30
1	C	97[A]	ARG	NE-CZ-NH1	7.73	124.16	120.30
1	C	97[B]	ARG	NE-CZ-NH1	7.73	124.16	120.30
1	D	97[A]	ARG	NE-CZ-NH1	7.66	124.13	120.30
1	D	97[B]	ARG	NE-CZ-NH1	7.66	124.13	120.30
1	A	30	ARG	NE-CZ-NH1	7.52	124.06	120.30
1	B	30	ARG	NE-CZ-NH1	7.51	124.06	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	15	TRP	CE2-CD2-CG	-7.50	101.30	107.30
1	A	15	TRP	CE2-CD2-CG	-7.46	101.33	107.30
1	A	37	TRP	CE2-CD2-CG	-7.41	101.37	107.30
1	B	37	TRP	CE2-CD2-CG	-7.37	101.40	107.30
1	D	37	TRP	CE2-CD2-CG	-7.30	101.46	107.30
1	C	37	TRP	CE2-CD2-CG	-7.22	101.53	107.30
1	A	109	CYS	CA-CB-SG	-7.06	101.29	114.00
1	B	109	CYS	CA-CB-SG	-7.05	101.30	114.00
1	B	40[A]	ARG	NE-CZ-NH1	6.59	123.60	120.30
1	B	40[B]	ARG	NE-CZ-NH1	6.59	123.60	120.30
1	A	40[A]	ARG	NE-CZ-NH1	6.58	123.59	120.30
1	A	40[B]	ARG	NE-CZ-NH1	6.58	123.59	120.30
1	D	30	ARG	NE-CZ-NH1	6.26	123.43	120.30
1	C	30	ARG	NE-CZ-NH1	6.26	123.43	120.30
1	B	14	LEU	CA-CB-CG	6.01	129.11	115.30
1	D	101	VAL	CG1-CB-CG2	-6.00	101.30	110.90
1	A	14	LEU	CA-CB-CG	6.00	129.09	115.30
1	C	101	VAL	CG1-CB-CG2	-5.98	101.33	110.90
1	D	35	TYR	CB-CG-CD2	-5.90	117.46	121.00
1	C	35	TYR	CB-CG-CD2	-5.88	117.47	121.00
1	D	20	VAL	CA-CB-CG2	-5.58	102.53	110.90
1	C	20	VAL	CA-CB-CG2	-5.57	102.54	110.90
1	D	15	TRP	CG-CD2-CE3	5.55	138.89	133.90
1	C	15	TRP	CG-CD2-CE3	5.51	138.86	133.90
1	D	40[A]	ARG	NE-CZ-NH1	5.44	123.02	120.30
1	D	40[B]	ARG	NE-CZ-NH1	5.44	123.02	120.30
1	C	40[A]	ARG	NE-CZ-NH1	5.41	123.00	120.30
1	C	40[B]	ARG	NE-CZ-NH1	5.41	123.00	120.30
1	B	97[A]	ARG	NE-CZ-NH2	-5.29	117.65	120.30
1	B	97[B]	ARG	NE-CZ-NH2	-5.29	117.65	120.30
1	D	15	TRP	CB-CG-CD1	-5.23	120.20	127.00
1	A	97[A]	ARG	NE-CZ-NH2	-5.23	117.69	120.30
1	A	97[B]	ARG	NE-CZ-NH2	-5.23	117.69	120.30
1	C	15	TRP	CB-CG-CD1	-5.22	120.21	127.00
1	B	15	TRP	CG-CD1-NE1	-5.21	104.89	110.10
1	A	15	TRP	CG-CD1-NE1	-5.18	104.92	110.10
1	B	126	VAL	CG1-CB-CG2	-5.15	102.66	110.90
1	A	126	VAL	CG1-CB-CG2	-5.15	102.67	110.90
1	C	63	HIS	CA-CB-CG	-5.11	104.91	113.60
1	D	63	HIS	CA-CB-CG	-5.10	104.94	113.60

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	1132	0	1155	12	0
1	B	1132	0	1155	15	0
1	C	1132	0	1156	16	0
1	D	1132	0	1156	18	0
2	A	43	0	30	1	0
2	B	43	0	30	0	0
2	C	43	0	30	0	0
2	D	43	0	30	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
4	A	8	0	0	1	0
4	B	8	0	0	1	0
4	C	13	0	0	0	0
4	D	13	0	0	0	0
All	All	4750	0	4742	55	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 (55) 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:14:LEU:HD22	1:A:118:LEU:HD23	1.80	0.63
1:B:14:LEU:HD22	1:B:118:LEU:HD23	1.80	0.63
1:B:97[B]:ARG:HG3	1:D:37:TRP:HA	1.86	0.57
1:C:1:VAL:HG13	1:C:136:SER:OG	2.04	0.57
1:A:97[B]:ARG:HG3	1:C:37:TRP:HA	1.86	0.57
1:D:1:VAL:HG13	1:D:136:SER:OG	2.04	0.57
1:D:113:THR:O	1:D:117:HIS:HD2	1.89	0.56
1:C:113:THR:O	1:C:117:HIS:HD2	1.89	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:85:PHE:HD1	1:B:88:LEU:HD12	1.75	0.52
1:D:14:LEU:HD12	1:D:121:GLU:HB3	1.92	0.51
1:C:14:LEU:HD12	1:C:121:GLU:HB3	1.92	0.51
1:C:62:ALA:O	1:C:66:LYS:HD2	2.11	0.51
1:A:85:PHE:HD1	1:A:88:LEU:HD12	1.75	0.50
1:D:62:ALA:O	1:D:66:LYS:HD2	2.11	0.50
1:B:97[B]:ARG:HB3	1:D:40[B]:ARG:HD3	1.93	0.50
1:A:97[B]:ARG:HB3	1:C:40[B]:ARG:HD3	1.93	0.49
1:D:109:CYS:O	1:D:113:THR:HG23	2.13	0.49
1:C:109:CYS:O	1:C:113:THR:HG23	2.13	0.48
1:C:85:PHE:HB3	1:C:141:LEU:CD1	2.44	0.47
1:C:142:THR:HG22	1:C:146:ARG:HH12	1.78	0.47
1:D:85:PHE:HB3	1:D:141:LEU:CD1	2.44	0.47
1:D:142:THR:HG22	1:D:146:ARG:HH12	1.78	0.47
1:D:26:GLU:HA	1:D:55:MET:HE1	1.97	0.47
1:B:5:PRO:HA	1:B:8:LYS:HE3	1.97	0.45
1:A:5:PRO:HA	1:A:8:LYS:HE3	1.98	0.45
1:A:91:LEU:HD22	1:A:96:LEU:HG	1.98	0.45
1:B:91:LEU:HD22	1:B:96:LEU:HG	1.99	0.45
1:C:50:THR:H	1:C:53:ALA:HB3	1.83	0.43
1:B:19:ASN:HB3	1:B:22:GLU:HG2	2.00	0.43
1:D:50:THR:H	1:D:53:ALA:HB3	1.83	0.43
1:A:57:ASN:O	1:A:60:VAL:HB	2.19	0.43
1:D:57:ASN:HA	1:D:58:PRO:HD3	1.78	0.43
1:A:50:THR:H	1:A:53:ALA:HB3	1.84	0.43
1:B:108[B]:HIS:ND1	1:C:108[B]:HIS:CD2	2.87	0.42
1:C:103:PHE:CE1	1:C:141:LEU:HD23	2.54	0.42
1:B:50:THR:H	1:B:53:ALA:HB3	1.84	0.42
1:B:57:ASN:O	1:B:60:VAL:HB	2.19	0.42
1:D:30:ARG:HD2	1:D:113:THR:HG22	2.02	0.42
1:D:103:PHE:CE1	1:D:141:LEU:HD23	2.54	0.42
1:A:19:ASN:HB3	1:A:22:GLU:HG2	2.00	0.41
1:B:146:ARG:HE	1:B:146:ARG:HA	1.86	0.41
1:B:68:LEU:HD23	1:B:68:LEU:HA	1.88	0.41
1:C:56:GLY:HA2	1:C:61:LYS:NZ	2.36	0.41
1:C:30:ARG:HD2	1:C:113:THR:HG22	2.02	0.41
1:D:51:PRO:O	1:D:55:MET:HG2	2.21	0.41
1:D:56:GLY:HA2	1:D:61:LYS:NZ	2.36	0.41
1:A:103:PHE:HD1	2:A:147:HEM:HBB2	1.85	0.41
4:A:241:HOH:O	1:D:108[B]:HIS:CE1	2.73	0.41
1:A:146:ARG:HD3	1:B:135:ALA:HB3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:220:HOH:O	1:C:108[B]:HIS:CE1	2.73	0.40
1:B:14:LEU:HG	1:B:126:VAL:HG11	2.03	0.40
1:C:134:LEU:HD12	1:C:134:LEU:HA	1.93	0.40
1:D:59:LYS:HZ2	1:D:59:LYS:HB2	1.86	0.40
1:B:97[A]:ARG:HD3	1:B:97[A]:ARG:HH11	1.78	0.40
1:A:146:ARG:HA	1:A:146:ARG:HE	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	147/146 (101%)	143 (97%)	4 (3%)	0	100	100
1	B	147/146 (101%)	143 (97%)	4 (3%)	0	100	100
1	C	147/146 (101%)	140 (95%)	7 (5%)	0	100	100
1	D	147/146 (101%)	141 (96%)	6 (4%)	0	100	100
All	All	588/584 (101%)	567 (96%)	21 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 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	125/122 (102%)	115 (92%)	10 (8%)	14	27
1	B	125/122 (102%)	115 (92%)	10 (8%)	14	27
1	C	125/122 (102%)	114 (91%)	11 (9%)	12	22
1	D	125/122 (102%)	114 (91%)	11 (9%)	12	22
All	All	500/488 (102%)	458 (92%)	42 (8%)	13	24

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	HIS
1	A	28	LEU
1	A	66	LYS
1	A	91	LEU
1	A	97[A]	ARG
1	A	97[B]	ARG
1	A	99	ASP
1	A	106	LEU
1	A	110	LEU
1	A	134	LEU
1	B	2	HIS
1	B	28	LEU
1	B	66	LYS
1	B	91	LEU
1	B	97[A]	ARG
1	B	97[B]	ARG
1	B	99	ASP
1	B	106	LEU
1	B	110	LEU
1	B	134	LEU
1	C	5	PRO
1	C	14	LEU
1	C	44	SER
1	C	59	LYS
1	C	66	LYS
1	C	111	LEU
1	C	114	LEU
1	C	123	THR
1	C	134	LEU
1	C	141	LEU
1	C	142	THR
1	D	5	PRO
1	D	14	LEU

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Mol	Chain	Res	Type
1	D	44	SER
1	D	59	LYS
1	D	66	LYS
1	D	111	LEU
1	D	114	LEU
1	D	123	THR
1	D	134	LEU
1	D	141	LEU
1	D	142	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	117	HIS
1	D	117	HIS

### 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](#)

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	147	1,3	28,50,50	1.74	6 (21%)	17,82,82	1.44	2 (11%)
3	CMO	A	148	2	0,1,1	0.00	-	0,0,0	0.00	-
2	HEM	B	147	1,3	28,50,50	1.74	6 (21%)	17,82,82	1.44	2 (11%)
3	CMO	B	148	2	0,1,1	0.00	-	0,0,0	0.00	-
2	HEM	C	147	1,3	28,50,50	1.69	6 (21%)	17,82,82	1.78	6 (35%)
3	CMO	C	148	2	0,1,1	0.00	-	0,0,0	0.00	-
2	HEM	D	147	1,3	28,50,50	1.69	6 (21%)	17,82,82	1.78	6 (35%)
3	CMO	D	148	2	0,1,1	0.00	-	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	147	1,3	-	0/6/54/54	0/0/8/8
3	CMO	A	148	2	-	0/0/0/0	0/0/0/0
2	HEM	B	147	1,3	-	0/6/54/54	0/0/8/8
3	CMO	B	148	2	-	0/0/0/0	0/0/0/0
2	HEM	C	147	1,3	-	0/6/54/54	0/0/8/8
3	CMO	C	148	2	-	0/0/0/0	0/0/0/0
2	HEM	D	147	1,3	-	0/6/54/54	0/0/8/8
3	CMO	D	148	2	-	0/0/0/0	0/0/0/0

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	147	HEM	C3B-CAB	-4.15	1.39	1.47
2	C	147	HEM	C3B-CAB	-4.12	1.39	1.47
2	A	147	HEM	C3C-CAC	-4.10	1.39	1.47
2	B	147	HEM	C3C-CAC	-4.08	1.39	1.47
2	D	147	HEM	C3C-CAC	-4.04	1.39	1.47
2	C	147	HEM	C3C-CAC	-4.03	1.39	1.47
2	B	147	HEM	C3B-CAB	-3.74	1.40	1.47
2	A	147	HEM	C3B-CAB	-3.72	1.40	1.47
2	B	147	HEM	C3B-C2B	-3.59	1.35	1.40
2	A	147	HEM	C3B-C2B	-3.55	1.35	1.40
2	A	147	HEM	C3C-C2C	-3.44	1.35	1.40
2	B	147	HEM	C3C-C2C	-3.38	1.35	1.40
2	C	147	HEM	C3C-C2C	-3.36	1.35	1.40
2	D	147	HEM	C3C-C2C	-3.34	1.35	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	147	HEM	C3B-C2B	-2.92	1.36	1.40
2	D	147	HEM	C3B-C2B	-2.92	1.36	1.40
2	D	147	HEM	CBC-CAC	2.27	1.44	1.28
2	C	147	HEM	CBC-CAC	2.28	1.44	1.28
2	C	147	HEM	CBB-CAB	2.38	1.45	1.28
2	B	147	HEM	CBB-CAB	2.38	1.45	1.28
2	A	147	HEM	CBB-CAB	2.38	1.45	1.28
2	D	147	HEM	CBB-CAB	2.39	1.45	1.28
2	B	147	HEM	CBC-CAC	2.42	1.45	1.28
2	A	147	HEM	CBC-CAC	2.42	1.45	1.28

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	147	HEM	CMD-C2D-C1D	-2.64	124.41	128.46
2	C	147	HEM	CMD-C2D-C1D	-2.64	124.41	128.46
2	D	147	HEM	CAA-CBA-CGA	-2.38	108.59	112.66
2	C	147	HEM	CAA-CBA-CGA	-2.37	108.62	112.66
2	D	147	HEM	C3B-C4B-NB	2.01	111.80	109.21
2	C	147	HEM	C3B-C4B-NB	2.01	111.81	109.21
2	C	147	HEM	CMD-C2D-C3D	2.06	128.84	124.94
2	D	147	HEM	CMD-C2D-C3D	2.07	128.84	124.94
2	D	147	HEM	CAD-CBD-CGD	2.08	116.21	112.66
2	C	147	HEM	CAD-CBD-CGD	2.08	116.22	112.66
2	A	147	HEM	CBA-CAA-C2A	2.41	117.10	112.48
2	B	147	HEM	CBA-CAA-C2A	2.42	117.11	112.48
2	B	147	HEM	CAD-CBD-CGD	2.77	117.40	112.66
2	A	147	HEM	CAD-CBD-CGD	2.78	117.40	112.66
2	D	147	HEM	CBA-CAA-C2A	3.51	119.19	112.48
2	C	147	HEM	CBA-CAA-C2A	3.53	119.24	112.48

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	147	HEM	1	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 is therefore empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section is therefore empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section is therefore empty.