



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:05 PM GMT

PDB ID : 1IRD
Title : Crystal Structure of Human Carbonmonoxy-Haemoglobin at 1.25 Å Resolution
Authors : Park, S.-Y.; Tame, J.R.H.
Deposited on : 2001-09-27
Resolution : 1.25 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
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

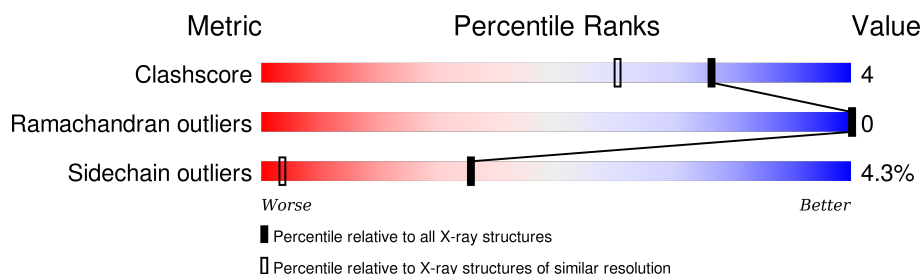
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.25 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	1530 (1.30-1.22)
Ramachandran outliers	100387	1467 (1.30-1.22)
Sidechain outliers	100360	1465 (1.30-1.22)

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 ≥ 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	141	 89% 9% ..
2	B	146	 92% 6% .

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2607 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	141	Total	C	N	O	S	0	0	0
			1069	685	187	194	3			

- Molecule 2 is a protein called Hemoglobin beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	146	Total	C	N	O	S	0	0	0
			1123	724	195	201	3			

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

- Molecule 4 is CARBON MONOXIDE (three-letter code: CMO) (formula: CO).



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

- Molecule 5 is water.


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	171	Total	O	0	0
			171	171		
5	B	154	Total	O	0	0
			154	154		

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 ($\text{RSRZ} > 2$). 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.

Note EDS was not executed.

- Molecule 1: Hemoglobin alpha chain

Chain A:  89% 9% ..



- Molecule 2: Hemoglobin beta chain

Chain B:  92% 6% .



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	53.27 Å 53.27 Å 190.78 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 1.25	Depositor
% Data completeness (in resolution range)	94.5 (25.00-1.25)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.163 , 0.214	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2607	wwPDB-VP
Average B, all atoms (Å ²)	22.0	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: 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	4.08	1/1097 (0.1%)	1.14	10/1491 (0.7%)
2	B	0.58	0/1153	1.12	4/1566 (0.3%)
All	All	2.88	1/2250 (0.0%)	1.13	14/3057 (0.5%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	141	ARG	C-OXT	133.92	3.77	1.23

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	92	ARG	NE-CZ-NH2	-8.95	115.83	120.30
1	A	140	TYR	CB-CG-CD1	-6.60	117.04	121.00
2	B	240	ARG	CD-NE-CZ	6.49	132.68	123.60
2	B	201	VAL	C-N-CA	6.46	137.85	121.70
2	B	202	HIS	CA-C-N	6.25	130.94	117.20
1	A	46	PHE	CB-CG-CD1	-6.18	116.47	120.80
1	A	74	ASP	CB-CG-OD1	5.87	123.58	118.30
1	A	74	ASP	CB-CG-OD2	-5.83	113.05	118.30
1	A	24	TYR	CB-CG-CD1	-5.59	117.65	121.00
1	A	92	ARG	CD-NE-CZ	5.48	131.27	123.60
2	B	330	TYR	CB-CG-CD1	-5.39	117.77	121.00
1	A	141	ARG	NE-CZ-NH2	-5.36	117.62	120.30
1	A	46	PHE	CB-CG-CD2	5.30	124.51	120.80
1	A	75	ASP	CB-CG-OD1	-5.16	113.66	118.30

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	1069	0	1073	7	0
2	B	1123	0	1115	5	0
3	A	43	0	30	5	0
3	B	43	0	30	0	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
5	A	171	0	0	1	0
5	B	154	0	0	1	0
All	All	2607	0	2248	17	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 4.

All (17) 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)
3:A:142:HEM:CBA	3:A:142:HEM:CGA	1.82	1.55
3:A:142:HEM:CGD	3:A:142:HEM:O2D	1.66	1.41
3:A:142:HEM:O2A	3:A:142:HEM:CBA	2.33	0.75
3:A:142:HEM:CAA	3:A:142:HEM:CGA	2.66	0.69
2:B:317:HIS:HE1	5:B:2073:HOH:O	1.85	0.60
3:A:142:HEM:CBD	3:A:142:HEM:O2D	2.48	0.59
1:A:63:ALA:O	1:A:67:THR:HG23	2.04	0.57
1:A:14:TRP:HE1	1:A:67:THR:HG22	1.73	0.54
2:B:208:LYS:O	2:B:212:THR:HG23	2.09	0.53
1:A:74:ASP:HB2	5:A:2172:HOH:O	2.11	0.51
1:A:14:TRP:NE1	1:A:67:THR:HG22	2.29	0.48
1:A:92:ARG:HD2	1:A:92:ARG:HA	1.49	0.47
1:A:80:LEU:O	1:A:84:SER:OG	2.30	0.46
2:B:344:LYS:HE2	2:B:344:LYS:HB3	1.84	0.44
2:B:295:LYS:HE3	2:B:295:LYS:HB2	1.67	0.44
2:B:275:LEU:HA	2:B:275:LEU:HD23	1.86	0.41
1:A:123:ALA:O	1:A:127:LYS:HG3	2.20	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	139/141 (99%)	138 (99%)	1 (1%)	0	100	100
2	B	144/146 (99%)	141 (98%)	3 (2%)	0	100	100
All	All	283/287 (99%)	279 (99%)	4 (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	113/113 (100%)	108 (96%)	5 (4%)	35	3
2	B	118/118 (100%)	113 (96%)	5 (4%)	36	4
All	All	231/231 (100%)	221 (96%)	10 (4%)	35	3

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LEU
1	A	84	SER
1	A	92	ARG
1	A	95	PRO
1	A	141	ARG
2	B	201	VAL
2	B	202	HIS
2	B	243	GLU

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Mol	Chain	Res	Type
2	B	295	LYS
2	B	320	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
2	B	239	GLN
2	B	317	HIS
2	B	343	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](#)

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	142	1,4	30,50,50	2.69	10 (33%)	24,82,82	2.46	9 (37%)
4	CMO	A	143	3	0,1,1	0.00	-	0,0,0	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	HEM	B	347	2,4	30,50,50	2.70	12 (40%)	24,82,82	4.06	12 (50%)
4	CMO	B	348	3	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
3	HEM	A	142	1,4	-	0/10/54/54	0/0/8/8
4	CMO	A	143	3	-	0/0/0/0	0/0/0/0
3	HEM	B	347	2,4	-	0/10/54/54	0/0/8/8
4	CMO	B	348	3	-	0/0/0/0	0/0/0/0

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	142	HEM	C2D-C3D	-7.82	1.31	1.54
3	B	347	HEM	C2D-C3D	-7.05	1.33	1.54
3	A	142	HEM	C2C-C1C	-6.69	1.39	1.52
3	A	142	HEM	CAA-C2A	-4.61	1.44	1.52
3	B	347	HEM	C2A-C3A	-4.38	1.24	1.37
3	B	347	HEM	C3D-C4D	-4.35	1.46	1.51
3	A	142	HEM	C3B-C4B	-3.99	1.48	1.51
3	A	142	HEM	C2B-C1B	-3.63	1.40	1.51
3	B	347	HEM	C3B-C4B	-3.23	1.49	1.51
3	B	347	HEM	CAD-C3D	-3.12	1.47	1.54
3	B	347	HEM	CBB-CAB	-2.97	1.12	1.29
3	B	347	HEM	CMB-C2B	-2.94	1.46	1.53
3	A	142	HEM	C1A-CHA	-2.81	1.32	1.39
3	A	142	HEM	C3C-CAC	-2.32	1.47	1.51
3	B	347	HEM	CHC-C1C	-2.24	1.31	1.36
3	A	142	HEM	CMD-C2D	2.01	1.57	1.53
3	B	347	HEM	CMA-C3A	2.19	1.56	1.51
3	A	142	HEM	FE-ND	2.19	2.09	1.97
3	B	347	HEM	CAA-C2A	2.65	1.56	1.52
3	A	142	HEM	CHD-C4C	3.21	1.43	1.36
3	B	347	HEM	C4C-NC	5.17	1.42	1.36
3	B	347	HEM	C1C-NC	5.20	1.42	1.36

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	347	HEM	CAA-C2A-C1A	-9.35	116.86	127.01
3	B	347	HEM	CMA-C3A-C4A	-4.98	120.13	128.36
3	B	347	HEM	C2C-C1C-NC	-4.41	102.77	110.21
3	A	142	HEM	CBA-CAA-C2A	-3.97	105.41	112.53
3	B	347	HEM	C3C-CAC-CBC	-3.33	119.35	124.46
3	A	142	HEM	CAA-CBA-CGA	-2.53	108.10	112.75
3	A	142	HEM	CMA-C3A-C4A	-2.42	124.36	128.36
3	A	142	HEM	C2D-C3D-C4D	2.01	104.91	101.50
3	B	347	HEM	CMA-C3A-C2A	2.75	131.00	125.24
3	B	347	HEM	CMD-C2D-C3D	3.13	128.18	114.35
3	B	347	HEM	CAD-C3D-C4D	3.65	125.33	112.47
3	A	142	HEM	CAD-C3D-C4D	3.69	125.47	112.47
3	A	142	HEM	CMD-C2D-C3D	3.84	131.33	114.35
3	A	142	HEM	CMB-C2B-C3B	4.36	127.42	116.53
3	A	142	HEM	CMC-C2C-C3C	4.69	128.23	116.53
3	B	347	HEM	C2C-C1C-CHC	4.81	131.00	123.68
3	A	142	HEM	CAD-C3D-C2D	5.67	129.52	113.22
3	B	347	HEM	CAD-C3D-C2D	6.25	131.20	113.22
3	B	347	HEM	CMC-C2C-C3C	6.70	133.25	116.53
3	B	347	HEM	CMB-C2B-C3B	6.81	133.54	116.53
3	B	347	HEM	C3B-CAB-CBB	7.65	136.20	124.46

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	142	HEM	5	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](#)

EDS was not executed - this section will therefore be empty.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.