



# Full wwPDB X-ray Structure Validation Report i

Oct 21, 2014 – 07:17 PM EDT

PDB ID : 1GDJ  
Title : CRYSTAL STRUCTURE OF FERRIC COMPLEXES OF THE YELLOW LUPIN LEGHEMOGLOBIN WITH ISOQUINOLINE AT 1.8 ANGSTROMS RESOLUTION (RUSSIAN)  
Authors : Harutyunyan, E.; Safonova, T.; Kuranova, I.  
Deposited on : 1994-09-14  
Resolution : 1.70 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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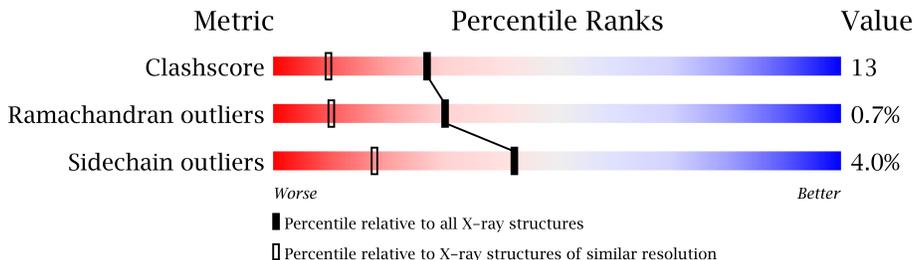
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.16 November 2013  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 21963  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable24103

# 1 Overall quality at a glance

The reported resolution of this entry is 1.70 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	79885	2929 (1.70-1.70)
Ramachandran outliers	78287	2878 (1.70-1.70)
Sidechain outliers	78261	2878 (1.70-1.70)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	153	

## 2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 1470 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

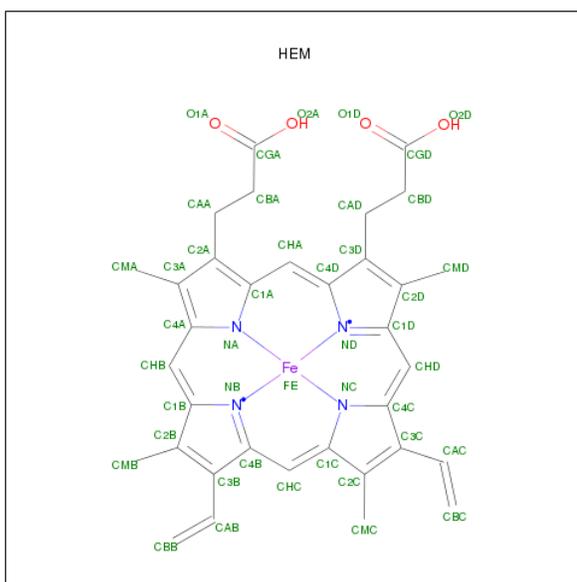
- Molecule 1 is a protein called LEGHEMOGLOBIN (DEOXY).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	153	1184	764	195	224	1	0	3	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	79	GLU	GLN	CONFLICT	UNP P02240
A	150	ASP	ASN	CONFLICT	UNP P02240

- 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
			Total	C	Fe	N	O		
2	A	1	43	34	1	4	4	0	0

- Molecule 3 is water.

<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
3	A	243	Total 243	O 243	0	0

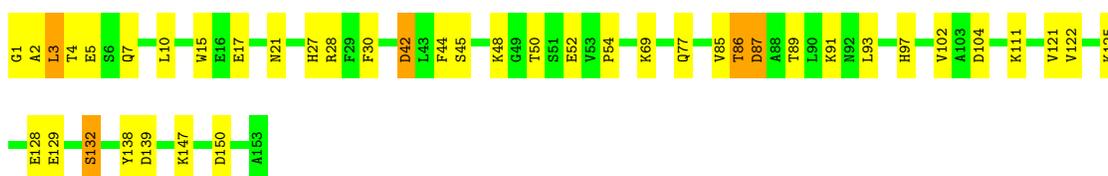
### 3 Residue-property plots i

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: LEGHEMOGLOBIN (DEOXY)

Chain A:



## 4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	93.48Å 51.60Å 38.33Å 90.00° 98.80° 90.00°	Depositor
Resolution (Å)	8.00 – 1.70	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-1.70)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, $R_{free}$	0.165 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1470	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	16.0	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

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	1/1224 (0.1%)	1.77	15/1663 (0.9%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	132	SER	CA-CB	5.31	1.60	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	2	ALA	C-N-CA	22.62	178.25	121.70
1	A	139	ASP	CB-CG-OD1	9.30	126.67	118.30
1	A	132	SER	O-C-N	-7.30	111.01	122.70
1	A	104	ASP	CB-CG-OD1	7.15	124.74	118.30
1	A	2	ALA	O-C-N	-7.09	111.36	122.70
1	A	28	ARG	CD-NE-CZ	6.91	133.27	123.60
1	A	86	THR	N-CA-CB	5.77	121.27	110.30
1	A	87	ASP	CB-CG-OD1	5.38	123.14	118.30
1	A	132	SER	CA-CB-OG	-5.31	96.86	111.20
1	A	42	ASP	CB-CG-OD1	-5.30	113.53	118.30
1	A	50	THR	CA-CB-CG2	5.25	119.75	112.40
1	A	5	GLU	CA-CB-CG	5.21	124.87	113.40
1	A	28	ARG	NE-CZ-NH1	-5.18	117.71	120.30
1	A	3	LEU	CA-C-O	5.17	130.95	120.10
1	A	30	PHE	CB-CG-CD1	-5.06	117.26	120.80

There are no chirality outliers.

There are no planarity outliers.

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1184	0	1200	32	0
2	A	43	0	30	0	0
3	A	243	0	0	9	1
All	All	1470	0	1230	32	1

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 13.

All (32) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:54:PRO:HB3	3:A:372:HOH:O	1.80	0.81
1:A:122:VAL:HG12	3:A:364:HOH:O	1.85	0.76
1:A:44:PHE:O	1:A:48:LYS:HG3	1.93	0.68
1:A:17:GLU:HB3	3:A:364:HOH:O	1.94	0.66
1:A:10:LEU:HD23	3:A:404:HOH:O	1.95	0.66
1:A:1:GLY:N	1:A:7:GLN:HE22	2.01	0.59
1:A:97[B]:HIS:CD2	1:A:102:VAL:HG21	2.39	0.57
1:A:93:LEU:HD23	1:A:97[B]:HIS:ND1	2.20	0.57
1:A:77:GLN:NE2	1:A:85:VAL:H	2.02	0.57
1:A:97[B]:HIS:HD2	1:A:102:VAL:HG21	1.72	0.55
1:A:93:LEU:HD23	1:A:97[B]:HIS:CE1	2.42	0.54
1:A:77:GLN:NE2	3:A:187:HOH:O	2.41	0.54
1:A:121:VAL:HG12	3:A:364:HOH:O	2.08	0.53
1:A:147:LYS:HE3	3:A:443:HOH:O	2.08	0.53
1:A:93:LEU:HD23	1:A:97[B]:HIS:HD1	1.71	0.53
1:A:93:LEU:CD2	1:A:97[B]:HIS:CE1	2.91	0.53
1:A:42:ASP:O	1:A:48:LYS:HE2	2.08	0.52
1:A:97[B]:HIS:HD2	1:A:102:VAL:HG11	1.75	0.51
1:A:15:TRP:CZ2	1:A:69:LYS:HD2	2.47	0.49
1:A:85:VAL:HG12	1:A:86:THR:O	2.13	0.49
1:A:86:THR:HG22	1:A:87:ASP:H	1.79	0.48
1:A:111:LYS:HB2	1:A:138:TYR:CE1	2.49	0.47
1:A:128:GLU:HG2	3:A:310:HOH:O	2.14	0.47
1:A:27:HIS:HE1	3:A:198:HOH:O	1.98	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:89:THR:O	1:A:93:LEU:HB2	2.17	0.45
1:A:1:GLY:H3	1:A:7:GLN:HE22	1.65	0.43
1:A:86:THR:HG22	1:A:87:ASP:N	2.34	0.43
1:A:45:SER:HA	1:A:48:LYS:HE3	2.00	0.43
1:A:17:GLU:CD	1:A:125:LYS:HZ1	2.23	0.42
1:A:1:GLY:HA3	1:A:4:THR:HG23	2.01	0.42
1:A:87:ASP:O	1:A:91:LYS:HG3	2.20	0.42
1:A:1:GLY:H1	1:A:7:GLN:HE22	1.68	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	Distance(Å)	Clash(Å)
3:A:385:HOH:O	3:A:385:HOH:O[2_657]	2.06	0.14

## 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	154/153 (101%)	150 (97%)	3 (2%)	1 (1%)	33 13

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	3	LEU

### 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	128/125 (102%)	123 (96%)	5 (4%)	43 18

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	ASN
1	A	52	GLU
1	A	129	GLU
1	A	132	SER
1	A	150	ASP

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	21	ASN
1	A	25	HIS
1	A	27	HIS
1	A	56	ASN
1	A	61	GLN
1	A	63	HIS
1	A	77	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

1 ligand is 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	154	1	42,50,50	3.81	18 (42%)	27,82,82	1.95	10 (37%)

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	154	1	-	0/14/114/114	0/0/8/8

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	154	HEM	C3C-C2C	-12.29	1.35	1.45
2	A	154	HEM	C3B-C2B	-11.92	1.34	1.45
2	A	154	HEM	CMD-C2D	6.95	1.56	1.45
2	A	154	HEM	CMB-C2B	6.95	1.56	1.45
2	A	154	HEM	CMC-C2C	5.25	1.54	1.45
2	A	154	HEM	C3D-C2D	-5.07	1.30	1.43
2	A	154	HEM	C3C-CAC	4.71	1.55	1.40
2	A	154	HEM	FE-NB	4.45	2.11	1.95
2	A	154	HEM	C1A-NA	3.97	1.43	1.36
2	A	154	HEM	C3B-CAB	3.97	1.53	1.40
2	A	154	HEM	FE-NC	3.64	2.10	1.95
2	A	154	HEM	FE-ND	3.27	2.08	1.95
2	A	154	HEM	C4D-ND	3.08	1.41	1.33
2	A	154	HEM	C4A-C3A	3.02	1.48	1.43
2	A	154	HEM	CAA-C2A	2.63	1.56	1.52
2	A	154	HEM	C1C-NC	2.53	1.40	1.33
2	A	154	HEM	C4C-NC	2.35	1.39	1.33
2	A	154	HEM	C1B-C2B	2.28	1.47	1.45

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	154	HEM	CAD-C3D-C4D	-3.85	120.17	125.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	154	HEM	CMB-C2B-C3B	3.25	133.67	126.21
2	A	154	HEM	CMA-C3A-C4A	-3.23	123.50	128.46
2	A	154	HEM	C4A-NA-C1A	-3.06	104.12	107.93
2	A	154	HEM	CMA-C3A-C2A	3.03	130.65	124.94
2	A	154	HEM	O2D-CGD-O1D	2.95	130.76	123.31
2	A	154	HEM	CBD-CAD-C3D	-2.83	108.37	114.51
2	A	154	HEM	C3A-C4A-NA	2.24	110.99	109.50
2	A	154	HEM	O1A-CGA-CBA	-2.06	116.09	123.06
2	A	154	HEM	O2A-CGA-CBA	2.01	121.14	114.19

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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.