



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 04:01 PM GMT

PDB ID : 4E44  
Title : Crystal structure of the hMHF1/hMHF2 Histone-Fold Tetramer  
Authors : Fox III, D.; Zhao, Y.; Yang, W.; Weidong, W.  
Deposited on : 2012-03-12  
Resolution : 2.10 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) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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) : trunk26865

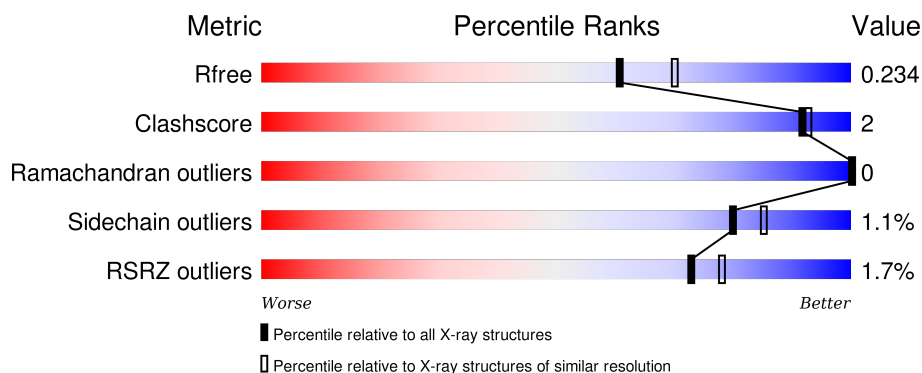
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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))
$R_{free}$	91344	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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	112	<div> <div>88%</div> <div>7%</div> </div>
1	C	112	<div> <div>83%</div> <div>13%</div> </div>
2	B	83	<div> <div>80%</div> <div>10%</div> <div>10%</div> </div>
2	D	83	<div> <div>87%</div> <div>10%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2783 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 Centromere protein S.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	104	Total	C	N	O	S	0	0	0
			810	503	138	163	6			
1	C	97	Total	C	N	O	S	0	1	0
			744	465	135	139	5			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	EXPRESSION TAG	UNP Q8N2Z9
A	0	SER	-	EXPRESSION TAG	UNP Q8N2Z9
C	-1	GLY	-	EXPRESSION TAG	UNP Q8N2Z9
C	0	SER	-	EXPRESSION TAG	UNP Q8N2Z9

- Molecule 2 is a protein called Centromere protein X.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	75	Total	C	N	O	S	0	0	0
			563	362	99	101	1			
2	D	75	Total	C	N	O	S	0	0	0
			569	366	97	105	1			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	GLY	-	EXPRESSION TAG	UNP A8MT69
B	0	SER	-	EXPRESSION TAG	UNP A8MT69
D	-1	GLY	-	EXPRESSION TAG	UNP A8MT69
D	0	SER	-	EXPRESSION TAG	UNP A8MT69

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	23	Total 23	O 23	0	0
3	C	32	Total 32	O 32	0	0
3	B	15	Total 15	O 15	0	0
3	D	27	Total 27	O 27	0	0



- Molecule 1: Centromere protein S



- GLY SER MET GLU GLU GLU ALA THR GLU E9 A86 S89 L92 Y95 T103 A104 Q105 ILE ASN LEU GLU ARG

- GLY SER MET GLU GLY ALA GLY ALA **G7** V15 V39 L42 **K43** V44 L63 R64 L70 V73 L74 F81

- GLY  
SER  
MET  
GLU  
GLY  
ALA  
GLY  
ALA  
G7  
R17  
V44  
D61  
F81

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	40.78Å 91.93Å 61.43Å 90.00° 96.30° 90.00°	Depositor
Resolution (Å)	45.97 – 2.10 30.53 – 2.10	Depositor EDS
% Data completeness (in resolution range)	96.6 (45.97-2.10) 96.6 (30.53-2.10)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.91 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, $R_{free}$	0.200 , 0.231 0.202 , 0.234	Depositor DCC
$R_{free}$ test set	1334 reflections (5.53%)	DCC
Wilson B-factor (Å <sup>2</sup> )	32.6	Xtriage
Anisotropy	0.037	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 41.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 25419 reflections	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	2783	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.45% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

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	0.73	0/819	0.77	1/1101 (0.1%)
1	C	0.80	0/756	0.81	0/1020
2	B	0.65	0/569	0.72	0/769
2	D	0.81	0/575	0.86	2/776 (0.3%)
All	All	0.75	0/2719	0.79	3/3666 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	61	ASP	CB-CG-OD1	6.09	123.78	118.30
2	D	17	ARG	NE-CZ-NH1	6.08	123.34	120.30
1	A	12	ARG	NE-CZ-NH2	5.01	122.81	120.30

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	810	0	773	3	0
1	C	744	0	711	4	0
2	B	563	0	567	7	0
2	D	569	0	575	3	0
3	A	23	0	0	0	0
3	B	15	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	32	0	0	0	0
3	D	27	0	0	0	0
All	All	2783	0	2626	11	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 2.

All (11) 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:95:TYR:CE2	2:B:44:VAL:HG21	2.14	0.82
1:C:86:ALA:HB1	1:C:92:LEU:HG	1.89	0.53
2:B:63:LEU:HD12	2:B:64:ARG:N	2.26	0.51
1:A:58:CYS:SG	2:B:42:LEU:HD11	2.54	0.48
1:C:95:TYR:CE2	2:D:44:VAL:HG21	2.50	0.47
1:A:53:LEU:CD2	2:B:74:LEU:HD22	2.45	0.47
1:C:92:LEU:CD1	2:D:44:VAL:HG11	2.44	0.47
1:C:95:TYR:CE2	2:D:44:VAL:CG2	2.98	0.46
2:B:70:LEU:O	2:B:73:VAL:HG22	2.20	0.41
2:B:15:VAL:HG21	2:B:39:VAL:HG22	2.02	0.41
2:B:63:LEU:HD12	2:B:63:LEU:C	2.40	0.41

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	102/112 (91%)	101 (99%)	1 (1%)	0	100	100
1	C	96/112 (86%)	94 (98%)	2 (2%)	0	100	100
2	B	73/83 (88%)	73 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	73/83 (88%)	73 (100%)	0	0	100	100
All	All	344/390 (88%)	341 (99%)	3 (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	84/99 (85%)	83 (99%)	1 (1%)	78	84
1	C	74/99 (75%)	73 (99%)	1 (1%)	74	80
2	B	57/68 (84%)	56 (98%)	1 (2%)	66	72
2	D	59/68 (87%)	59 (100%)	0	100	100
All	All	274/334 (82%)	271 (99%)	3 (1%)	80	85

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
1	C	89	SER
2	B	63	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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](#)

There are no ligands in this entry.

## 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 [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	104/112 (92%)	0.06	1 (0%) 84 87	27, 37, 54, 80	0
1	C	97/112 (86%)	0.08	2 (2%) 67 72	19, 32, 66, 82	0
2	B	75/83 (90%)	-0.01	1 (1%) 79 84	26, 38, 54, 61	0
2	D	75/83 (90%)	-0.03	2 (2%) 58 65	20, 29, 46, 61	0
All	All	351/390 (90%)	0.03	6 (1%) 73 78	19, 34, 58, 82	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	103	ILE	3.6
1	A	95	TYR	3.4
2	B	7	GLY	3.1
1	C	104	ALA	2.7
2	D	7	GLY	2.7
2	D	81	PHE	2.4

### 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](#)

There are no ligands in this entry.

## 6.5 Other polymers

There are no such residues in this entry.